Full wwPDB X-ray Structure Validation Report ⓘ

Mar 14, 2026 – 09:22 PM UTC

PDB ID	:	4V8K / pdb_00004v8k
Title	:	Crystal structure of the LH1-RC complex from Thermochromatium tepidum in P21 form
Authors	:	Niwa, S.; Takeda, K.; Wang-Otomo, Z.-Y.; Miki, K.
Deposited on	:	2013-11-22
Resolution	:	3.01 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

The following versions of software and data (see references ⓘ) were used in the production of this report:

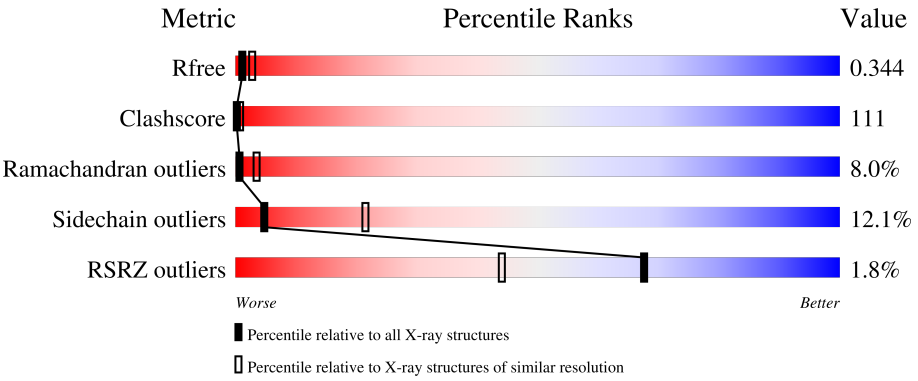
MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 3.01 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



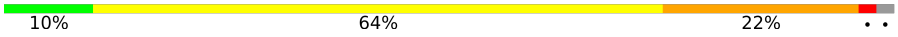
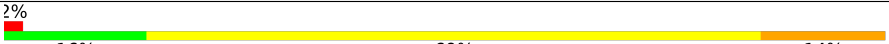
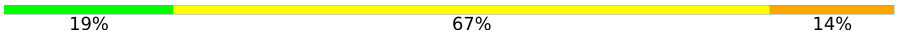
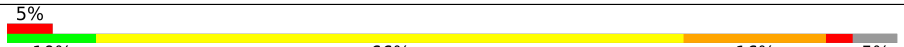
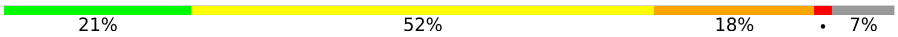
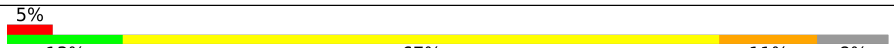
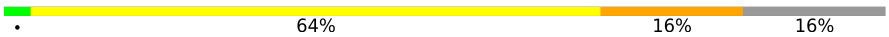
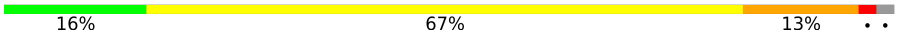
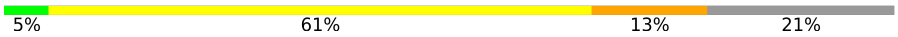
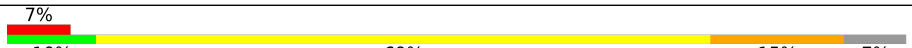
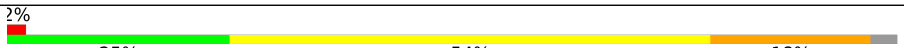
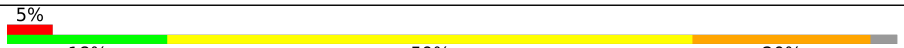
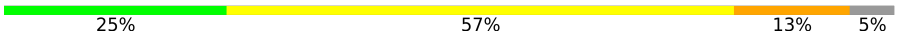
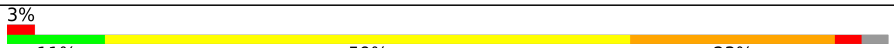
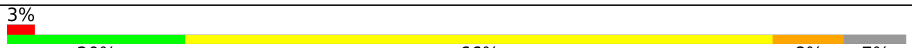
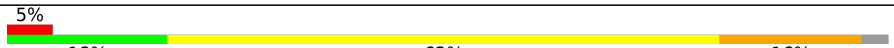
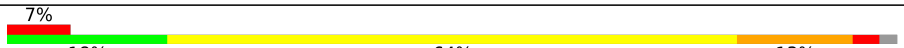
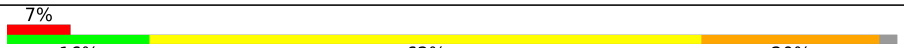
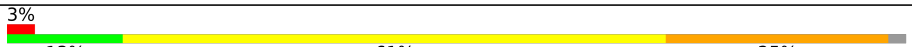
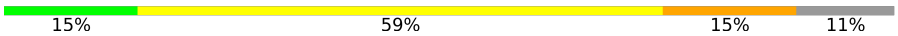
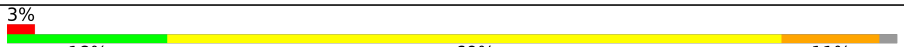
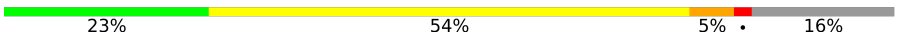
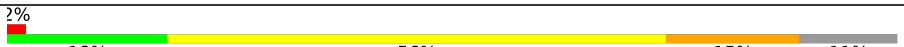
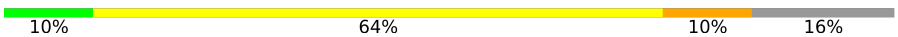
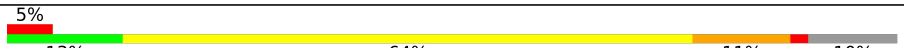
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	2672 (3.00-3.00)
Clashscore	190562	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)
RSRZ outliers	180081	2671 (3.00-3.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=5% The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	AC	404	11% 48% 17% 22%
1	BC	404	13% 49% 15% 22%
2	AL	281	8% 63% 26%
2	BL	281	9% 73% 17%
3	AM	325	8% 60% 27%

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Mol	Chain	Length	Quality of chain
3	BM	325	
4	AH	259	
4	BH	259	
5	A1	61	
5	A3	61	
5	A5	61	
5	A7	61	
5	A9	61	
5	AA	61	
5	AD	61	
5	AF	61	
5	AI	61	
5	AK	61	
5	AO	61	
5	AQ	61	
5	AS	61	
5	AU	61	
5	AW	61	
5	AY	61	
5	B1	61	
5	B3	61	
5	B5	61	
5	B7	61	
5	B9	61	
5	BA	61	

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Mol	Chain	Length	Quality of chain
5	BD	61	5% 59% 13% 26%
5	BF	61	7% 13% 66% 13% 8%
5	BI	61	2% 8% 62% 10% 18%
5	BK	61	3% 21% 61% 16%
5	BO	61	2% 16% 64% 13%
5	BQ	61	3% 26% 62% 8%
5	BS	61	3% 21% 69% 5%
5	BU	61	2% 13% 66% 16% 5%
5	BW	61	2% 13% 62% 20% 5%
5	BY	61	13% 56% 20% 11%
6	A0	47	11% 45% 28% 15%
6	A2	47	21% 55% 9% 15%
6	A4	47	4% 17% 55% 13% 15%
6	A6	47	2% 30% 47% 9% 15%
6	A8	47	6% 9% 60% 17% 15%
6	AB	47	2% 17% 55% 13% 15%
6	AE	47	4% 23% 57% 15%
6	AG	47	9% 57% 19% 15%
6	AJ	47	2% 15% 62% 9% 15%
6	AN	47	2% 13% 57% 15% 15%
6	AP	47	2% 6% 62% 17% 15%
6	AR	47	4% 17% 55% 13% 15%
6	AT	47	17% 45% 23% 15%
6	AV	47	2% 19% 51% 15% 15%
6	AX	47	9% 72% 15%

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Mol	Chain	Length	Quality of chain
6	AZ	47	15% 53% 13% 15%
6	B0	47	4% 13% 55% 15% 15%
6	B2	47	15% 60% 11% 15%
6	B4	47	4% 26% 49% 11% 15%
6	B6	47	26% 51% 9% 15%
6	B8	47	9% 60% 15% 15%
6	BB	47	15% 57% 13% 15%
6	BE	47	17% 64% 15%
6	BG	47	9% 13% 66% 15%
6	BJ	47	11% 68% 6% 15%
6	BN	47	17% 55% 13% 15%
6	BP	47	2% 9% 66% 11% 15%
6	BR	47	9% 28% 47% 11% 15%
6	BT	47	13% 53% 19% 15%
6	BV	47	2% 19% 57% 9% 15%
6	BX	47	21% 53% 11% 15%
6	BZ	47	23% 53% 6% 15%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
14	CRT	A0	101	-	-	X	-
14	CRT	A1	103	-	-	X	-
14	CRT	A2	102	-	-	X	-
14	CRT	A5	103	-	-	X	-
14	CRT	A7	102	-	-	X	-
14	CRT	AA	102	-	-	X	-
14	CRT	AB	102	-	-	X	-
14	CRT	AJ	102	-	-	X	-

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
14	CRT	AP	102	-	-	X	-
14	CRT	AS	104	-	-	X	-
14	CRT	AW	102	-	-	X	-
14	CRT	AX	102	-	-	X	-
14	CRT	B0	101	-	-	X	-
14	CRT	B1	103	-	-	X	-
14	CRT	B2	102	-	-	X	-
14	CRT	B5	103	-	-	X	-
14	CRT	B7	102	-	-	X	-
14	CRT	BA	102	-	-	X	-
14	CRT	BB	102	-	-	X	-
14	CRT	BP	102	-	-	X	-
14	CRT	BU	103	-	-	X	-
14	CRT	BV	102	-	-	X	-
15	PEF	AM	407	-	X	-	-
15	PEF	AS	101	-	-	X	-
9	BCL	A0	102	-	-	X	-
9	BCL	A1	102	-	-	X	-
9	BCL	A2	101	-	-	X	-
9	BCL	A3	103	-	-	X	-
9	BCL	A3	104	-	-	X	-
9	BCL	A5	102	-	-	X	-
9	BCL	A6	101	-	-	X	-
9	BCL	A7	103	-	-	X	-
9	BCL	A8	101	-	-	X	-
9	BCL	A9	102	-	-	X	-
9	BCL	AA	101	-	-	X	-
9	BCL	AB	101	-	-	X	-
9	BCL	AD	102	-	-	X	-
9	BCL	AE	101	-	-	X	-
9	BCL	AF	102	-	-	X	-
9	BCL	AG	101	-	-	X	-
9	BCL	AI	102	-	-	X	-
9	BCL	AJ	101	-	-	X	-
9	BCL	AK	102	-	-	X	-
9	BCL	AL	301	-	-	X	-
9	BCL	AL	303	-	-	X	-
9	BCL	AM	402	-	-	X	-
9	BCL	AN	101	-	-	X	-
9	BCL	AO	102	-	-	X	-
9	BCL	AP	101	-	-	X	-
9	BCL	AQ	102	-	-	X	-

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
9	BCL	AR	101	-	-	X	-
9	BCL	AS	103	-	-	X	-
9	BCL	AT	101	-	-	X	-
9	BCL	AU	102	-	-	X	-
9	BCL	AV	102	-	-	X	-
9	BCL	AW	101	-	-	X	-
9	BCL	AX	101	-	-	X	-
9	BCL	AY	102	-	-	X	-
9	BCL	AZ	101	-	-	X	-
9	BCL	B0	102	-	-	X	-
9	BCL	B1	102	-	-	X	-
9	BCL	B2	101	-	-	X	-
9	BCL	B3	102	-	-	X	-
9	BCL	B6	101	-	-	X	-
9	BCL	B7	103	-	-	X	-
9	BCL	B8	101	-	-	X	-
9	BCL	B9	102	-	-	X	-
9	BCL	BA	101	-	-	X	-
9	BCL	BB	101	-	-	X	-
9	BCL	BD	102	-	-	X	-
9	BCL	BE	101	-	-	X	-
9	BCL	BF	102	-	-	X	-
9	BCL	BG	101	-	-	X	-
9	BCL	BI	102	-	-	X	-
9	BCL	BJ	101	-	-	X	-
9	BCL	BK	102	-	-	X	-
9	BCL	BL	301	-	-	X	-
9	BCL	BM	402	-	-	X	-
9	BCL	BN	101	-	-	X	-
9	BCL	BO	102	-	-	X	-
9	BCL	BP	101	-	-	X	-
9	BCL	BQ	103	-	-	X	-
9	BCL	BQ	104	-	-	X	-
9	BCL	BS	102	-	-	X	-
9	BCL	BT	101	-	-	X	-
9	BCL	BU	102	-	-	X	-
9	BCL	BV	101	-	-	X	-
9	BCL	BW	102	-	-	X	-
9	BCL	BX	101	-	-	X	-
9	BCL	BY	102	-	-	X	-
9	BCL	BZ	101	-	-	X	-

2 Entry composition (i)

There are 17 unique types of molecules in this entry. The entry contains 50862 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Photosynthetic reaction center cytochrome c subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	AC	317	Total	C	N	O	S	0	0	0
			2458	1551	430	460	17			
1	BC	317	Total	C	N	O	S	0	0	0
			2458	1551	430	460	17			

- Molecule 2 is a protein called Photosynthetic reaction center L subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	AL	280	Total	C	N	O	S	0	0	0
			2231	1501	359	361	10			
2	BL	280	Total	C	N	O	S	0	0	0
			2231	1501	359	361	10			

- Molecule 3 is a protein called Photosynthetic reaction center M subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	AM	319	Total	C	N	O	S	0	0	0
			2551	1713	417	410	11			
3	BM	319	Total	C	N	O	S	0	0	0
			2551	1713	417	410	11			

- Molecule 4 is a protein called Photosynthetic reaction center H subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	AH	258	Total	C	N	O	S	0	0	0
			1982	1275	339	363	5			
4	BH	258	Total	C	N	O	S	0	0	0
			1982	1275	339	363	5			

- Molecule 5 is a protein called LH1 alpha polypeptide.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	AA	48	Total	C	N	O	S	0	0	0
			392	265	62	64	1			
5	AD	57	Total	C	N	O	S	0	0	0
			447	295	74	77	1			
5	AF	59	Total	C	N	O	S	0	0	0
			462	304	76	80	2			
5	AI	59	Total	C	N	O	S	0	0	0
			462	304	76	80	2			
5	AK	58	Total	C	N	O	S	0	0	0
			455	300	75	78	2			
5	AO	59	Total	C	N	O	S	0	0	0
			462	304	76	80	2			
5	AQ	57	Total	C	N	O	S	0	0	0
			447	295	74	77	1			
5	AS	59	Total	C	N	O	S	0	0	0
			462	304	76	80	2			
5	AU	60	Total	C	N	O	S	0	0	0
			473	313	77	81	2			
5	AW	60	Total	C	N	O	S	0	0	0
			473	313	77	81	2			
5	AY	60	Total	C	N	O	S	0	0	0
			473	313	77	81	2			
5	A1	58	Total	C	N	O	S	0	0	0
			455	300	75	78	2			
5	A3	57	Total	C	N	O	S	0	0	0
			447	295	74	77	1			
5	A5	56	Total	C	N	O	S	0	0	0
			444	294	73	75	2			
5	A7	51	Total	C	N	O	S	0	0	0
			417	279	67	69	2			
5	A9	60	Total	C	N	O	S	0	0	0
			473	313	77	81	2			
5	BA	55	Total	C	N	O	S	0	0	0
			448	299	72	75	2			
5	BD	45	Total	C	N	O	S	0	0	0
			370	250	59	60	1			
5	BF	56	Total	C	N	O	S	0	0	0
			444	294	73	75	2			
5	BI	50	Total	C	N	O	S	0	0	0
			409	274	66	68	1			
5	BK	60	Total	C	N	O	S	0	0	0
			473	313	77	81	2			
5	BO	59	Total	C	N	O	S	0	0	0
			462	304	76	80	2			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	BQ	59	Total	C	N	O	S	0	0	0
			467	310	76	79	2			
5	BS	59	Total	C	N	O	S	0	0	0
			462	304	76	80	2			
5	BU	58	Total	C	N	O	S	0	0	0
			462	307	75	78	2			
5	BW	58	Total	C	N	O	S	0	0	0
			455	300	75	78	2			
5	BY	54	Total	C	N	O	S	0	0	0
			426	284	69	72	1			
5	B1	54	Total	C	N	O	S	0	0	0
			426	284	69	72	1			
5	B3	60	Total	C	N	O	S	0	0	0
			473	313	77	81	2			
5	B5	51	Total	C	N	O	S	0	0	0
			417	279	67	69	2			
5	B7	54	Total	C	N	O	S	0	0	0
			426	284	69	72	1			
5	B9	51	Total	C	N	O	S	0	0	0
			417	279	67	69	2			

- Molecule 6 is a protein called LH1 beta polypeptide.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
6	AB	40	Total	C	N	O	S	0	0	0
			337	228	52	55	2			
6	AE	40	Total	C	N	O	S	0	0	0
			337	228	52	55	2			
6	AG	40	Total	C	N	O	S	0	0	0
			337	228	52	55	2			
6	AJ	40	Total	C	N	O	S	0	0	0
			337	228	52	55	2			
6	AN	40	Total	C	N	O	S	0	0	0
			337	228	52	55	2			
6	AP	40	Total	C	N	O	S	0	0	0
			337	228	52	55	2			
6	AR	40	Total	C	N	O	S	0	0	0
			337	228	52	55	2			
6	AT	40	Total	C	N	O	S	0	0	0
			337	228	52	55	2			
6	AV	40	Total	C	N	O	S	0	0	0
			337	228	52	55	2			

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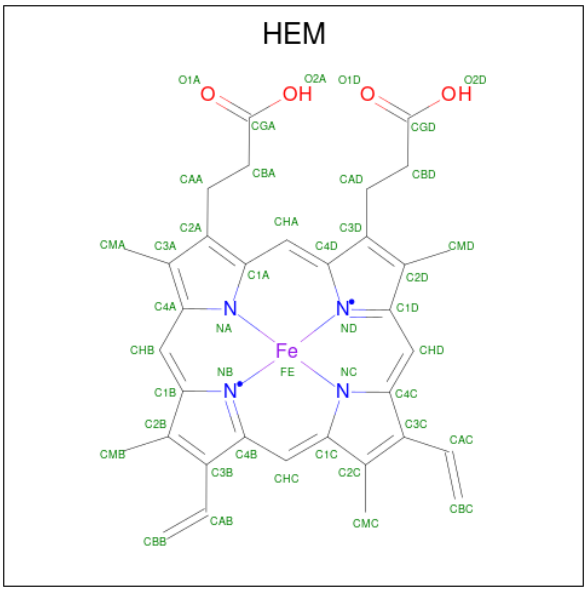
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
6	AX	40	Total	C	N	O	S	0	0	0
			337	228	52	55	2			
6	AZ	40	Total	C	N	O	S	0	0	0
			337	228	52	55	2			
6	A2	40	Total	C	N	O	S	0	0	0
			337	228	52	55	2			
6	A4	40	Total	C	N	O	S	0	0	0
			337	228	52	55	2			
6	A6	40	Total	C	N	O	S	0	0	0
			337	228	52	55	2			
6	A8	40	Total	C	N	O	S	0	0	0
			337	228	52	55	2			
6	A0	40	Total	C	N	O	S	0	0	0
			337	228	52	55	2			
6	BB	40	Total	C	N	O	S	0	0	0
			337	228	52	55	2			
6	BE	40	Total	C	N	O	S	0	0	0
			337	228	52	55	2			
6	BG	40	Total	C	N	O	S	0	0	0
			337	228	52	55	2			
6	BJ	40	Total	C	N	O	S	0	0	0
			337	228	52	55	2			
6	BN	40	Total	C	N	O	S	0	0	0
			337	228	52	55	2			
6	BP	40	Total	C	N	O	S	0	0	0
			337	228	52	55	2			
6	BR	40	Total	C	N	O	S	0	0	0
			337	228	52	55	2			
6	BT	40	Total	C	N	O	S	0	0	0
			337	228	52	55	2			
6	BV	40	Total	C	N	O	S	0	0	0
			337	228	52	55	2			
6	BX	40	Total	C	N	O	S	0	0	0
			337	228	52	55	2			
6	BZ	40	Total	C	N	O	S	0	0	0
			337	228	52	55	2			
6	B2	40	Total	C	N	O	S	0	0	0
			337	228	52	55	2			
6	B4	40	Total	C	N	O	S	0	0	0
			337	228	52	55	2			
6	B6	40	Total	C	N	O	S	0	0	0
			337	228	52	55	2			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
6	B8	40	Total	C	N	O	S	0	0	0
			337	228	52	55	2			
6	B0	40	Total	C	N	O	S	0	0	0
			337	228	52	55	2			

- Molecule 7 is PROTOPORPHYRIN IX CONTAINING FE (CCD ID: HEM) (formula: $C_{34}H_{32}FeN_4O_4$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
7	AC	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
7	AC	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
7	AC	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
7	AC	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
7	BC	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
7	BC	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
7	BC	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
7	BC	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		

- Molecule 8 is CALCIUM ION (CCD ID: CA) (formula: Ca).

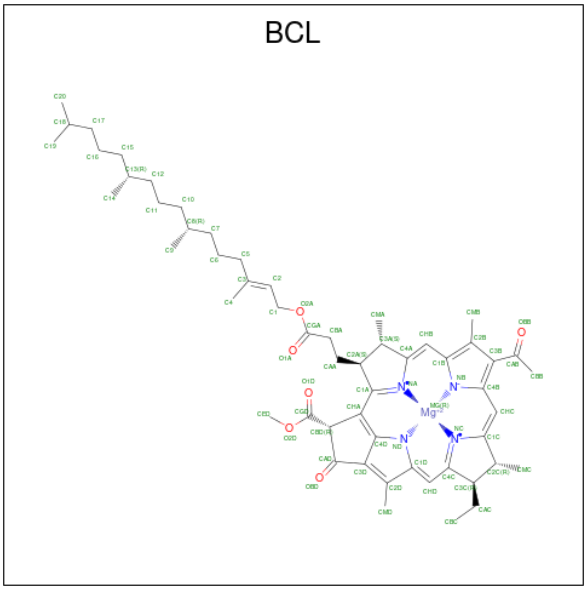
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	AC	1	Total Ca 1 1	0	0
8	AA	1	Total Ca 1 1	0	0
8	AD	1	Total Ca 1 1	0	0
8	AF	1	Total Ca 1 1	0	0
8	AI	1	Total Ca 1 1	0	0
8	AK	1	Total Ca 1 1	0	0
8	AO	1	Total Ca 1 1	0	0
8	AQ	1	Total Ca 1 1	0	0
8	AS	1	Total Ca 1 1	0	0
8	AU	1	Total Ca 1 1	0	0
8	AV	1	Total Ca 1 1	0	0
8	AY	1	Total Ca 1 1	0	0
8	A1	1	Total Ca 1 1	0	0
8	A3	1	Total Ca 1 1	0	0
8	A5	1	Total Ca 1 1	0	0
8	A7	1	Total Ca 1 1	0	0
8	A9	1	Total Ca 1 1	0	0
8	BC	1	Total Ca 1 1	0	0
8	BA	1	Total Ca 1 1	0	0
8	BD	1	Total Ca 1 1	0	0
8	BF	1	Total Ca 1 1	0	0
8	BI	1	Total Ca 1 1	0	0

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	BK	1	Total Ca 1 1	0	0
8	BO	1	Total Ca 1 1	0	0
8	BQ	1	Total Ca 1 1	0	0
8	BS	1	Total Ca 1 1	0	0
8	BU	1	Total Ca 1 1	0	0
8	BW	1	Total Ca 1 1	0	0
8	BY	1	Total Ca 1 1	0	0
8	B1	1	Total Ca 1 1	0	0
8	B3	1	Total Ca 1 1	0	0
8	B5	1	Total Ca 1 1	0	0
8	B7	1	Total Ca 1 1	0	0
8	B9	1	Total Ca 1 1	0	0

- Molecule 9 is BACTERIOCHLOROPHYLL A (CCD ID: BCL) (formula: $C_{55}H_{74}MgN_4O_6$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
9	AL	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
9	AL	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
9	AM	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
9	AM	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
9	AA	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
9	AB	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
9	AD	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
9	AE	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
9	AF	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
9	AG	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
9	AI	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
9	AJ	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
9	AK	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
9	AN	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
9	AO	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
9	AP	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
9	AQ	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
9	AR	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
9	AS	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
9	AT	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
9	AU	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
9	AV	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
9	AW	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
9	AX	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
9	AY	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
9	AZ	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
9	A1	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
9	A2	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
9	A3	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
9	A3	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
9	A5	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
9	A6	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
9	A7	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
9	A8	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
9	A9	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
9	A0	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
9	BL	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
9	BL	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
9	BM	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
9	BM	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
9	BA	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
9	BB	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
9	BD	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		

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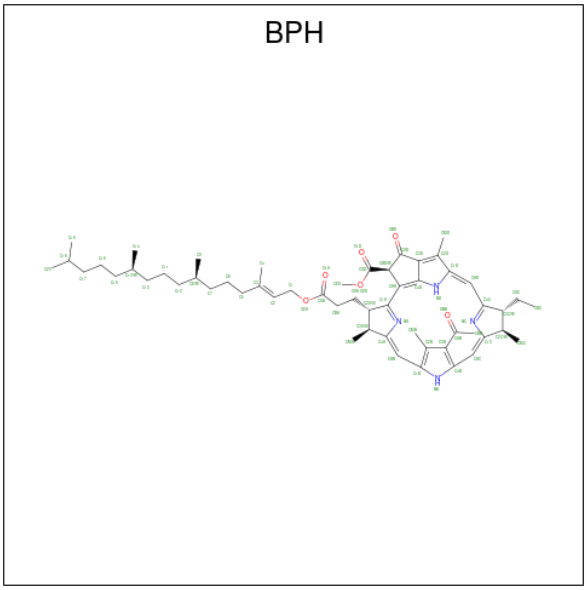
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
9	BE	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
9	BF	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
9	BG	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
9	BI	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
9	BJ	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
9	BK	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
9	BN	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
9	BO	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
9	BP	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
9	BQ	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
9	BQ	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
9	BS	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
9	BT	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
9	BU	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
9	BV	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
9	BW	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
9	BX	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
9	BY	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
9	BZ	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
9	B1	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
9	B2	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
9	B3	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
9	B4	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
9	B5	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
9	B6	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
9	B7	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
9	B8	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
9	B9	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
9	B0	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		

- Molecule 10 is BACTERIOPHEOPHYTIN A (CCD ID: BPH) (formula: $C_{55}H_{76}N_4O_6$).



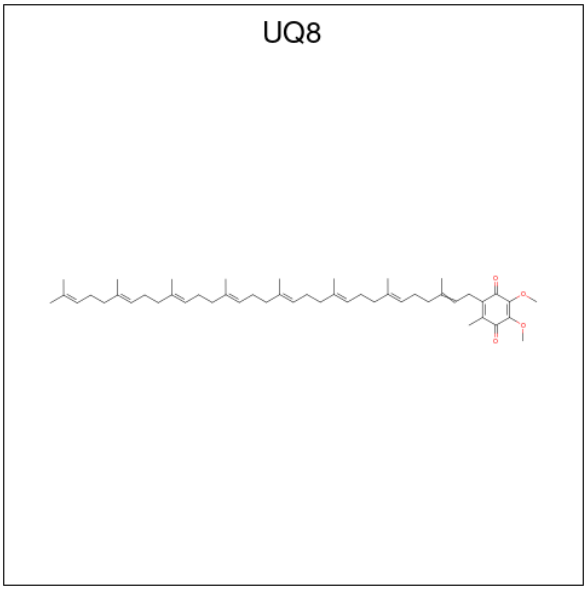
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
10	AL	1	Total	C	N	O	0	0
			65	55	4	6		
10	AM	1	Total	C	N	O	0	0
			65	55	4	6		
10	BL	1	Total	C	N	O	0	0
			65	55	4	6		

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
10	BM	1	Total C N O 65 55 4 6	0	0

- Molecule 11 is Ubiquinone-8 (CCD ID: UQ8) (formula: $C_{49}H_{74}O_4$).

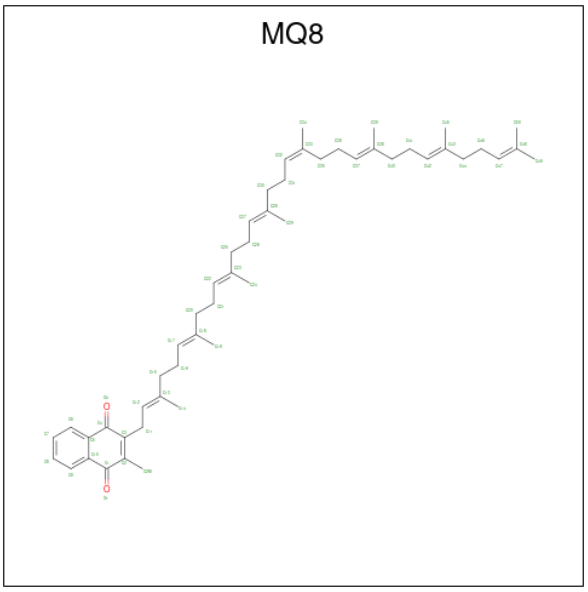


Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
11	AL	1	Total C O 53 49 4	0	0
11	BL	1	Total C O 53 49 4	0	0

- Molecule 12 is FE (III) ION (CCD ID: FE) (formula: Fe).

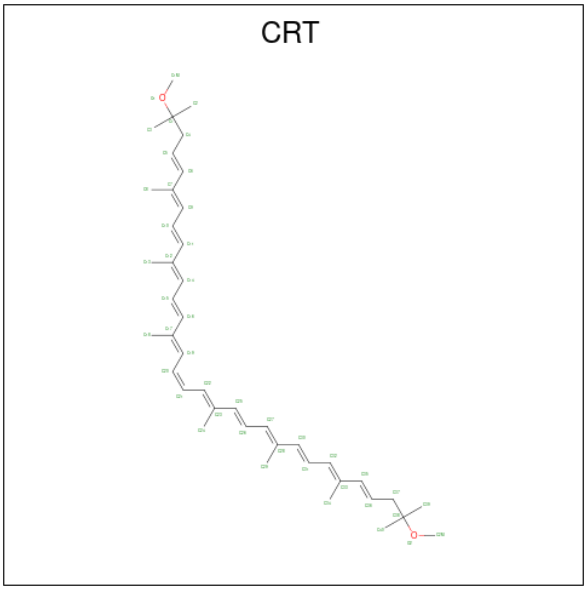
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
12	AM	1	Total Fe 1 1	0	0
12	BM	1	Total Fe 1 1	0	0

- Molecule 13 is MENAQUINONE 8 (CCD ID: MQ8) (formula: $C_{51}H_{72}O_2$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
13	AM	1	Total C O 53 51 2	0	0
13	BM	1	Total C O 53 51 2	0	0

- Molecule 14 is SPIRILLOXANTHIN (CCD ID: CRT) (formula: $C_{42}H_{60}O_2$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
14	AM	1	Total C O 44 42 2	0	0
14	AA	1	Total C O 44 42 2	0	0

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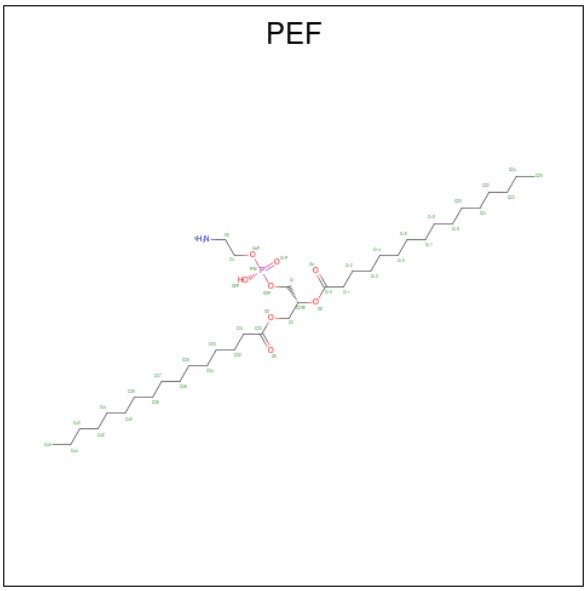
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
14	AB	1	Total C O 44 42 2	0	0
14	AG	1	Total C O 44 42 2	0	0
14	AJ	1	Total C O 44 42 2	0	0
14	AN	1	Total C O 44 42 2	0	0
14	AP	1	Total C O 44 42 2	0	0
14	AR	1	Total C O 44 42 2	0	0
14	AS	1	Total C O 44 42 2	0	0
14	AT	1	Total C O 44 42 2	0	0
14	AW	1	Total C O 44 42 2	0	0
14	AX	1	Total C O 44 42 2	0	0
14	A1	1	Total C O 44 42 2	0	0
14	A2	1	Total C O 44 42 2	0	0
14	A5	1	Total C O 44 42 2	0	0
14	A7	1	Total C O 44 42 2	0	0
14	A0	1	Total C O 44 42 2	0	0
14	BM	1	Total C O 44 42 2	0	0
14	BA	1	Total C O 44 42 2	0	0
14	BB	1	Total C O 44 42 2	0	0
14	BF	1	Total C O 44 42 2	0	0
14	BG	1	Total C O 44 42 2	0	0
14	BN	1	Total C O 44 42 2	0	0

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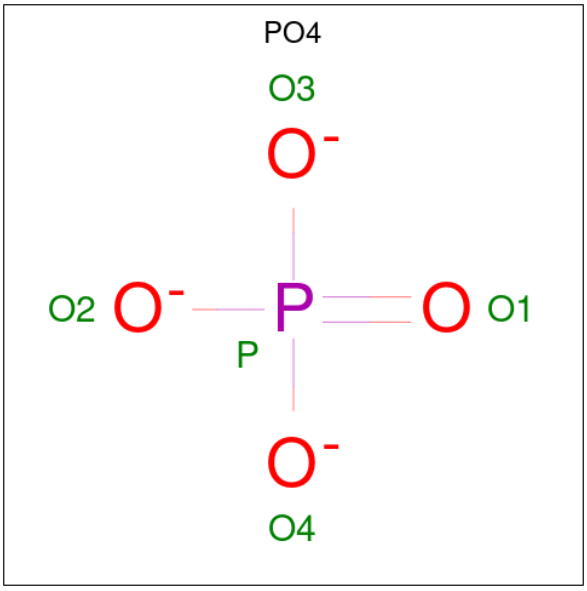
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
14	BO	1	Total C O 44 42 2	0	0
14	BP	1	Total C O 44 42 2	0	0
14	BS	1	Total C O 44 42 2	0	0
14	BU	1	Total C O 44 42 2	0	0
14	BV	1	Total C O 44 42 2	0	0
14	BW	1	Total C O 44 42 2	0	0
14	B1	1	Total C O 44 42 2	0	0
14	B2	1	Total C O 44 42 2	0	0
14	B5	1	Total C O 44 42 2	0	0
14	B7	1	Total C O 44 42 2	0	0
14	B0	1	Total C O 44 42 2	0	0

- Molecule 15 is DI-PALMITOYL-3-SN-PHOSPHATIDYLETHANOLAMINE (CCD ID: PEF) (formula: $C_{37}H_{74}NO_8P$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
15	AM	1	Total C N O P 19 9 1 8 1	0	0
15	AM	1	Total C N O P 14 6 1 6 1	0	0
15	AM	1	Total C N O P 47 37 1 8 1	0	0
15	AH	1	Total C N O P 19 9 1 8 1	0	0
15	AS	1	Total C N O P 47 37 1 8 1	0	0
15	BM	1	Total C N O P 19 9 1 8 1	0	0
15	BQ	1	Total C N O P 47 37 1 8 1	0	0

- Molecule 16 is PHOSPHATE ION (CCD ID: PO4) (formula: O_4P).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
16	AM	1	Total O P 5 4 1	0	0
16	AH	1	Total O P 5 4 1	0	0
16	A3	1	Total O P 5 4 1	0	0
16	BH	1	Total O P 5 4 1	0	0

- Molecule 17 is water.

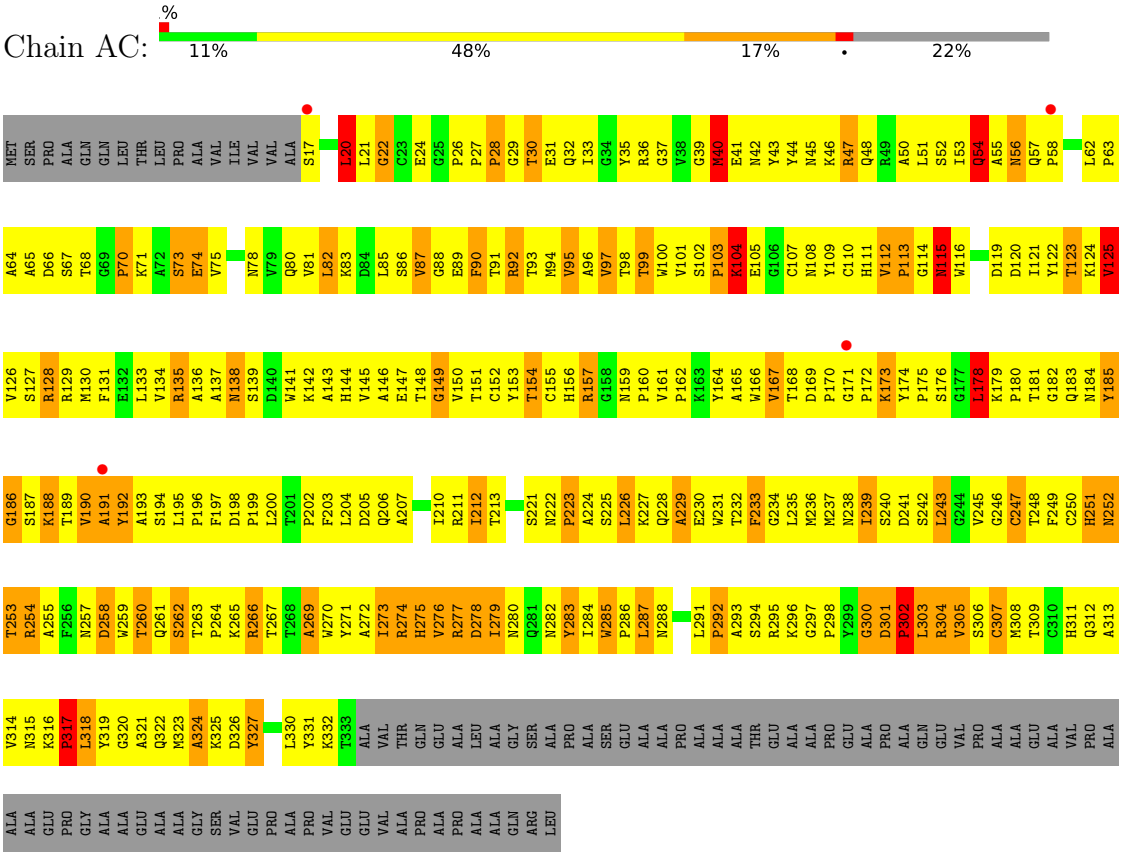
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
17	AC	1	Total O 1 1	0	0
17	AL	3	Total O 3 3	0	0
17	AM	3	Total O 3 3	0	0
17	AH	2	Total O 2 2	0	0
17	AA	1	Total O 1 1	0	0
17	AI	1	Total O 1 1	0	0
17	AW	1	Total O 1 1	0	0
17	BC	1	Total O 1 1	0	0
17	BL	3	Total O 3 3	0	0
17	BM	3	Total O 3 3	0	0
17	BH	1	Total O 1 1	0	0
17	B1	1	Total O 1 1	0	0

3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density (RSRZ > 2). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

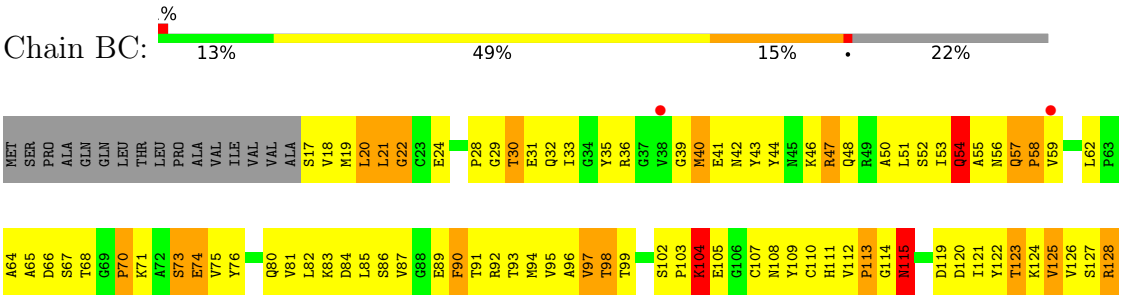
• Molecule 1: Photosynthetic reaction center cytochrome c subunit

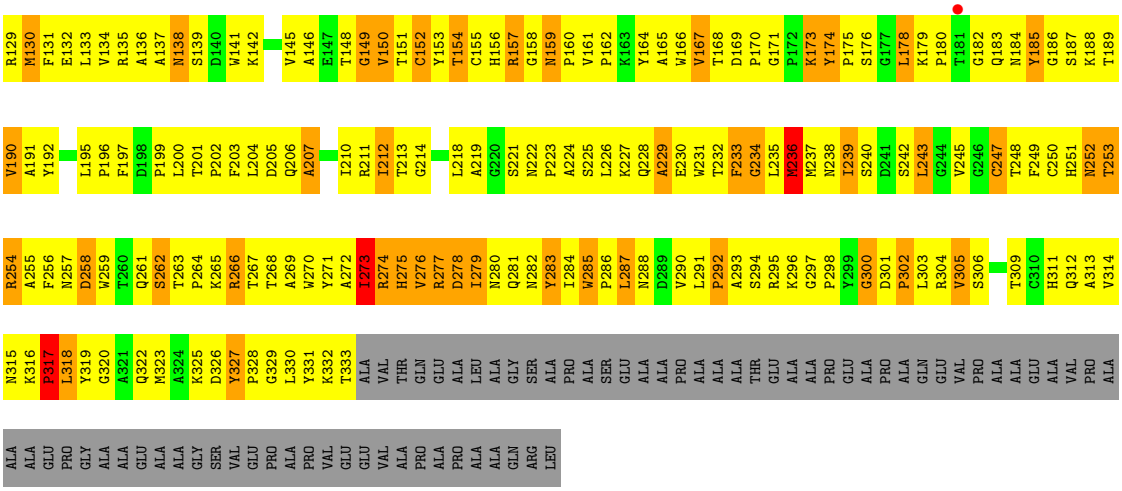
Chain AC:



• Molecule 1: Photosynthetic reaction center cytochrome c subunit

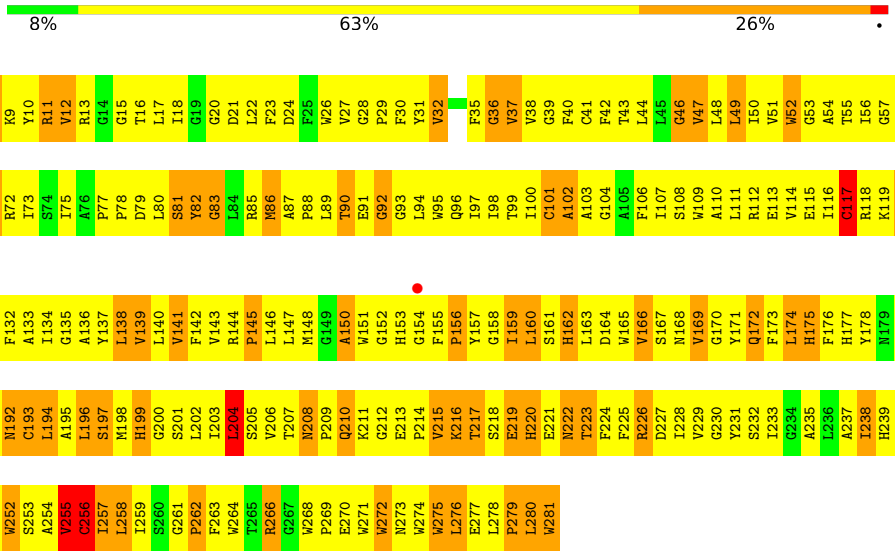
Chain BC:





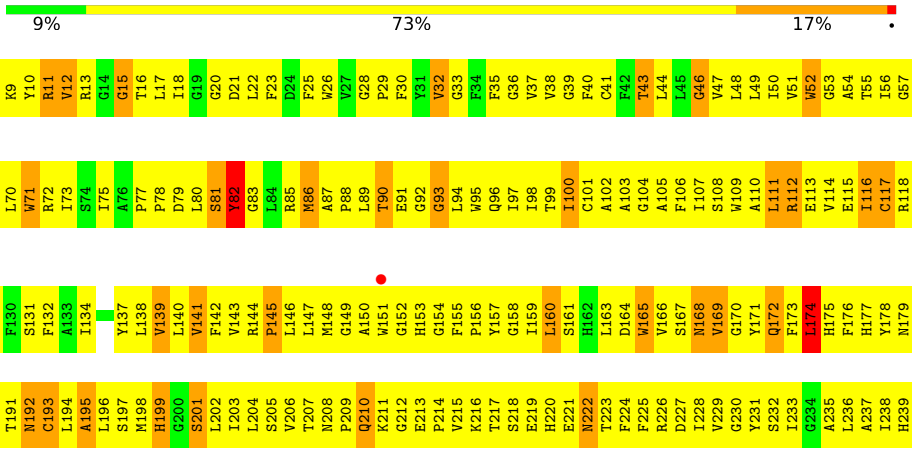
• Molecule 2: Photosynthetic reaction center L subunit

Chain AL:



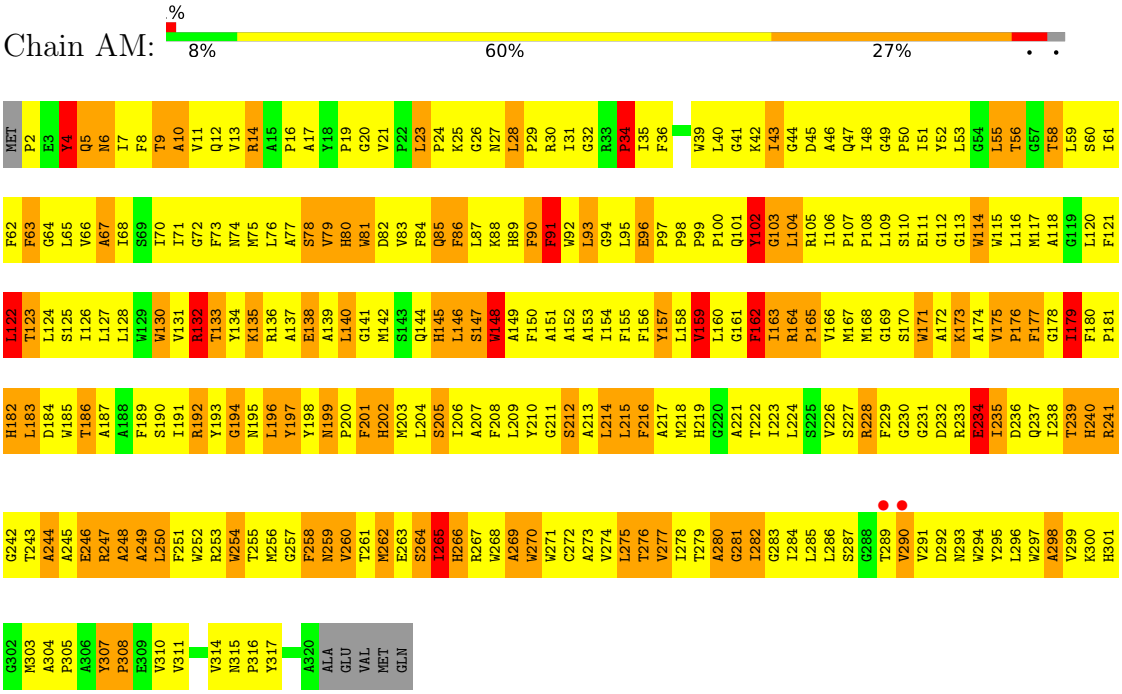
• Molecule 2: Photosynthetic reaction center L subunit

Chain BL:

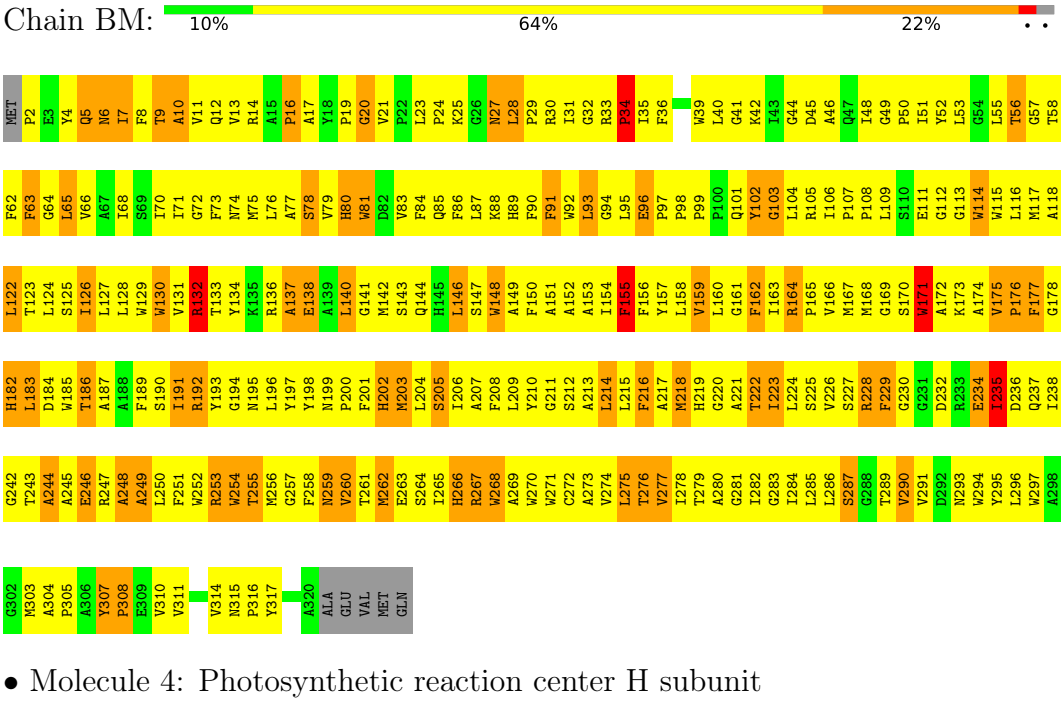




• Molecule 3: Photosynthetic reaction center M subunit

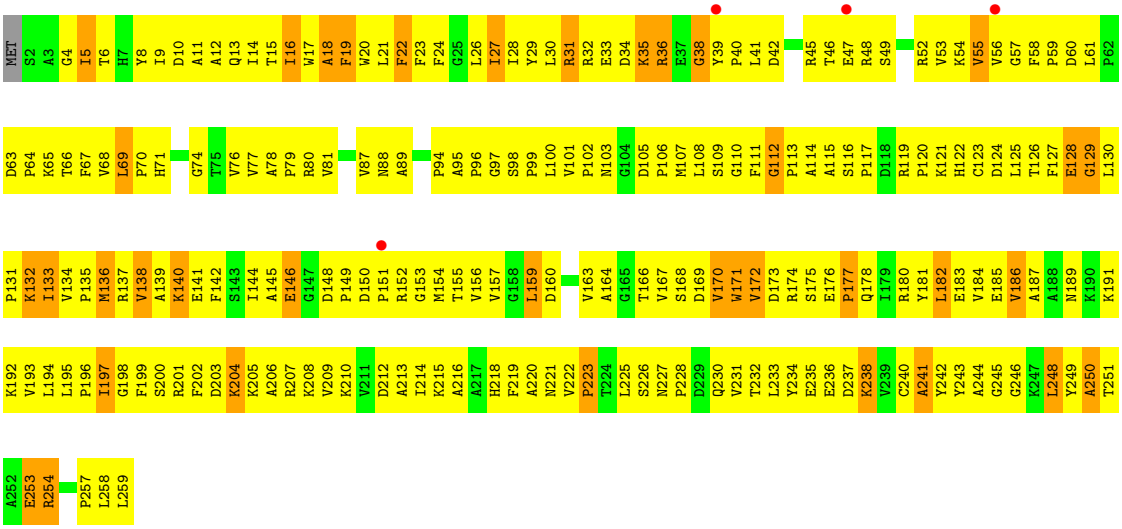


• Molecule 3: Photosynthetic reaction center M subunit



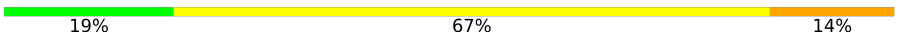
• Molecule 4: Photosynthetic reaction center H subunit

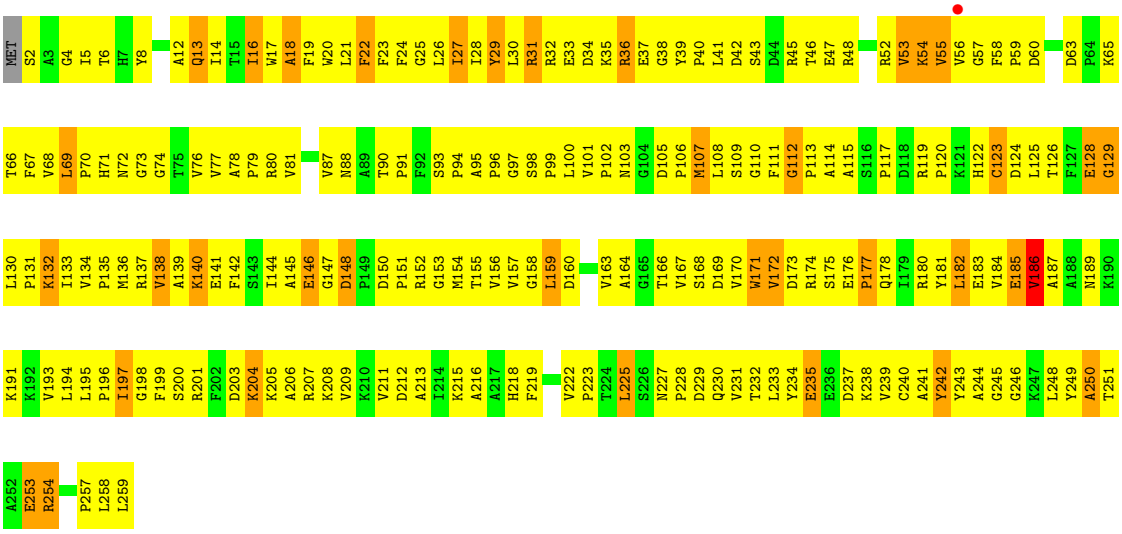




• Molecule 4: Photosynthetic reaction center H subunit

Chain BH:

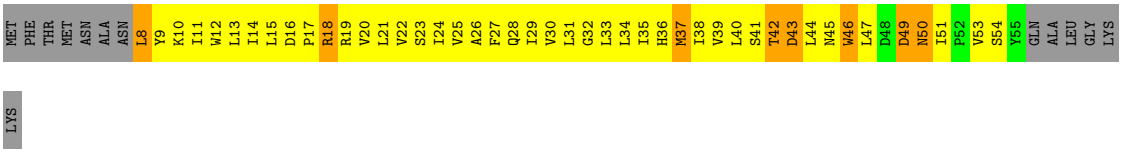




• Molecule 5: LH1 alpha polypeptide

Chain AA:

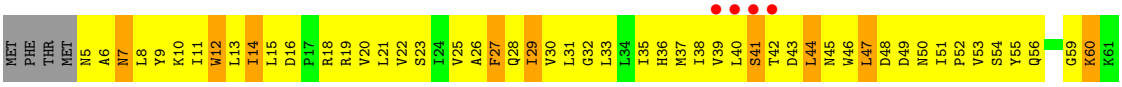




• Molecule 5: LH1 alpha polypeptide

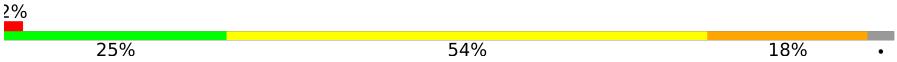
Chain AD:

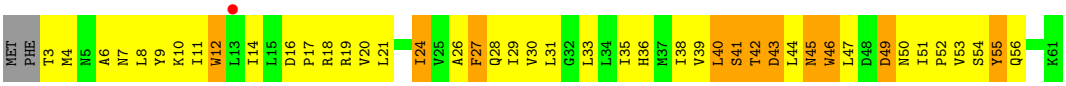




• Molecule 5: LH1 alpha polypeptide

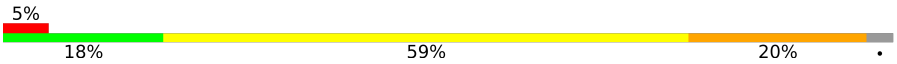
Chain AF:

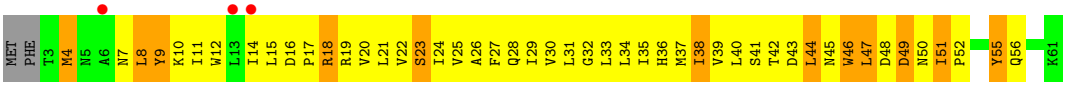




• Molecule 5: LH1 alpha polypeptide

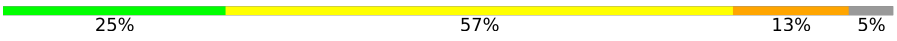
Chain AI:

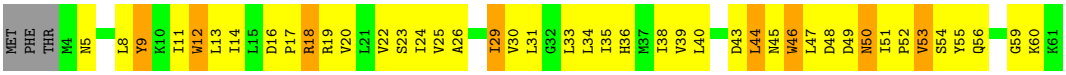




• Molecule 5: LH1 alpha polypeptide

Chain AK:

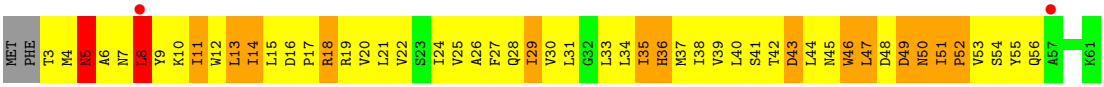




• Molecule 5: LH1 alpha polypeptide

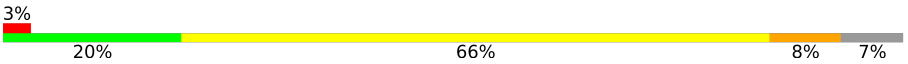
Chain AO:

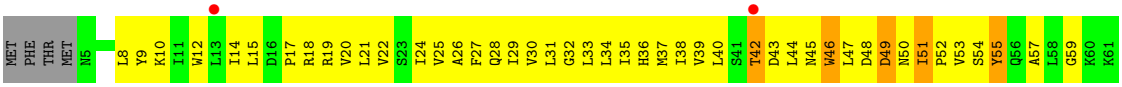




• Molecule 5: LH1 alpha polypeptide

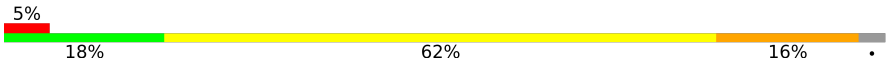
Chain AQ:

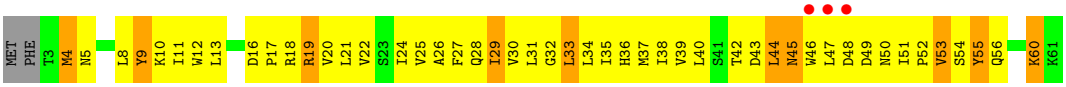




• Molecule 5: LH1 alpha polypeptide

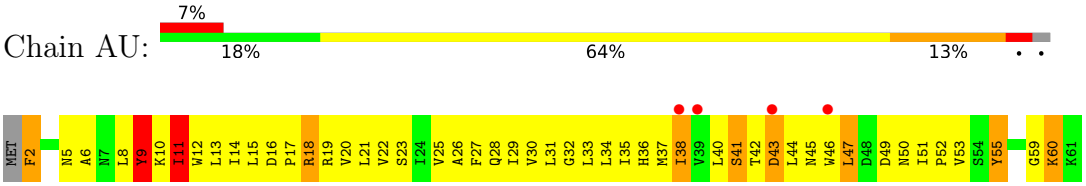
Chain AS:





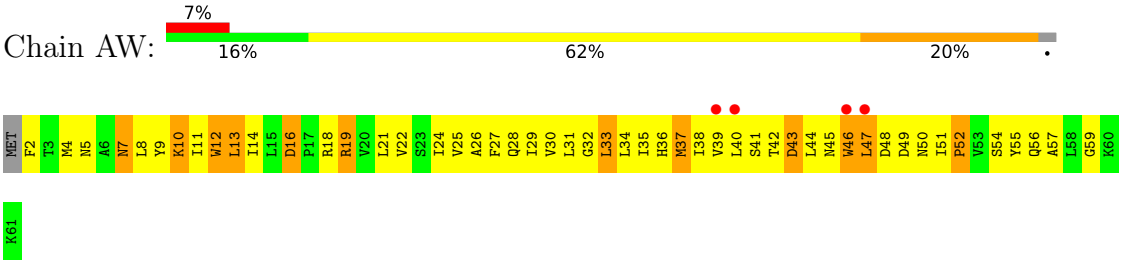
• Molecule 5: LH1 alpha polypeptide

Chain AU:



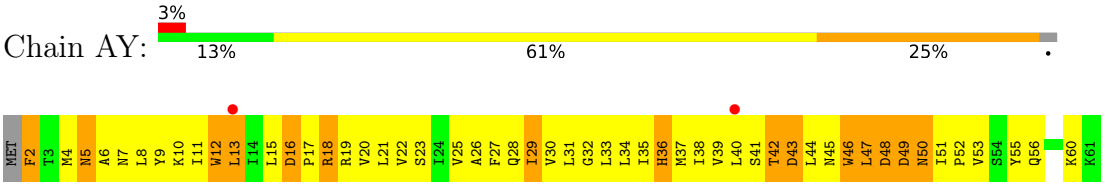
• Molecule 5: LH1 alpha polypeptide

Chain AW:



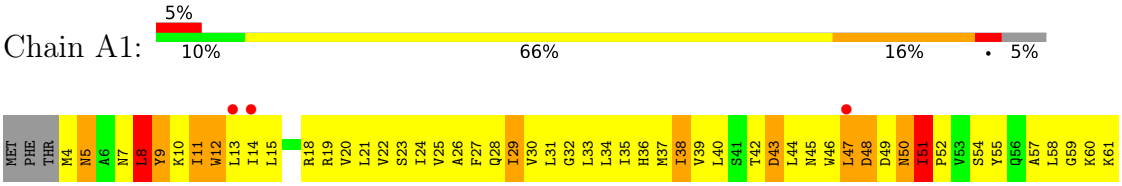
• Molecule 5: LH1 alpha polypeptide

Chain AY:



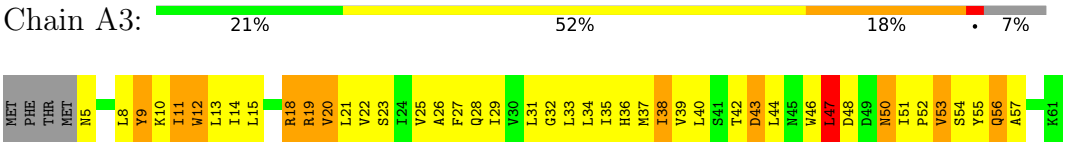
• Molecule 5: LH1 alpha polypeptide

Chain A1:



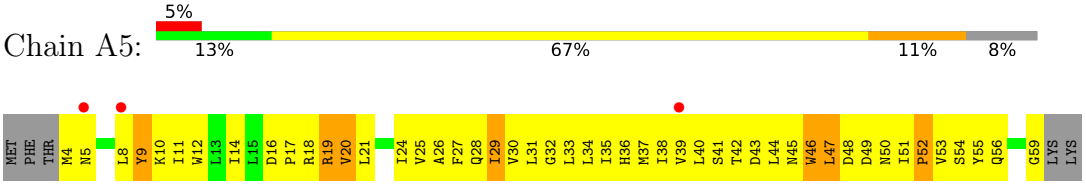
• Molecule 5: LH1 alpha polypeptide

Chain A3:



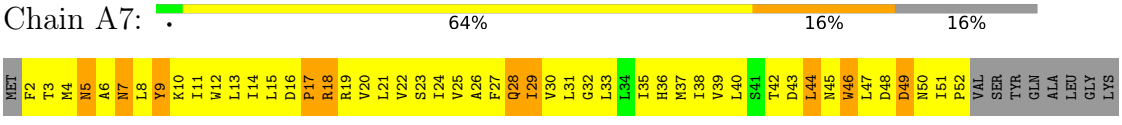
• Molecule 5: LH1 alpha polypeptide

Chain A5:



• Molecule 5: LH1 alpha polypeptide

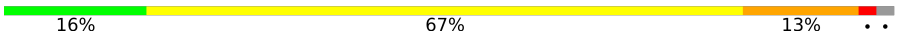
Chain A7:

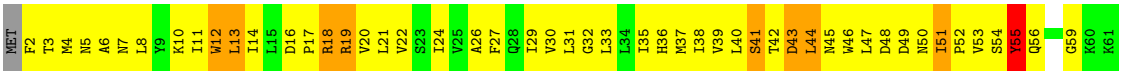


LYS

• Molecule 5: LH1 alpha polypeptide

Chain A9:

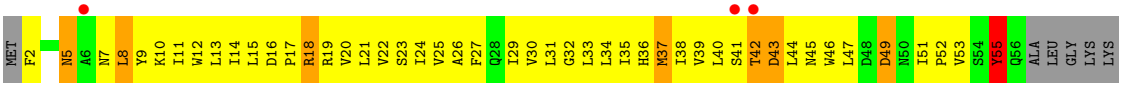




• Molecule 5: LH1 alpha polypeptide

Chain BA:





• Molecule 5: LH1 alpha polypeptide

Chain BD:





LYS

• Molecule 5: LH1 alpha polypeptide

Chain BF:



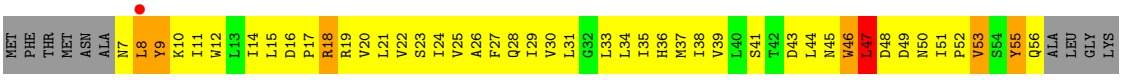


LYS

• Molecule 5: LH1 alpha polypeptide

Chain BI:



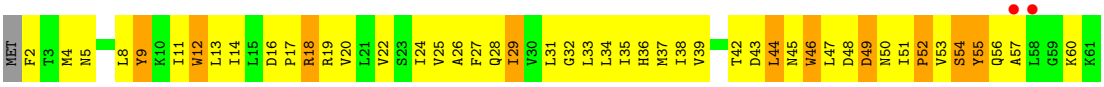


LYS

• Molecule 5: LH1 alpha polypeptide

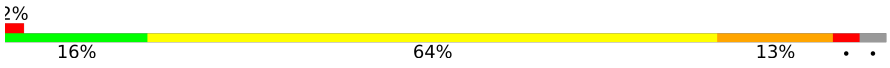
Chain BK:

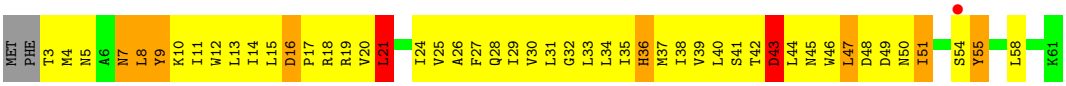




• Molecule 5: LH1 alpha polypeptide

Chain BO:

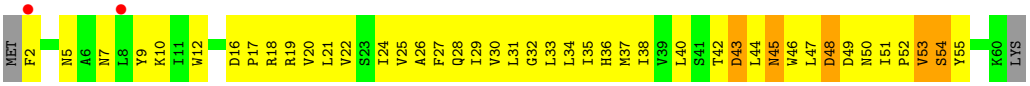




• Molecule 5: LH1 alpha polypeptide

Chain BQ:





• Molecule 5: LH1 alpha polypeptide

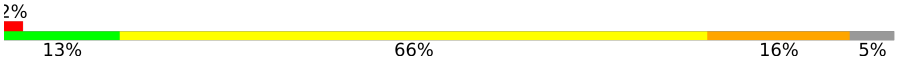
Chain BS:





• Molecule 5: LH1 alpha polypeptide

Chain BU:

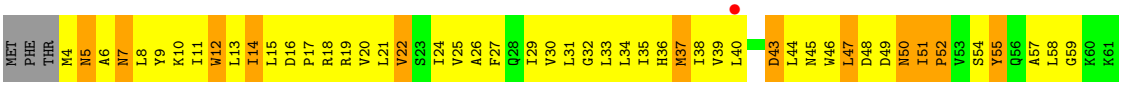




• Molecule 5: LH1 alpha polypeptide

Chain BW:

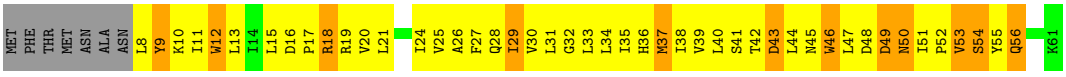




• Molecule 5: LH1 alpha polypeptide

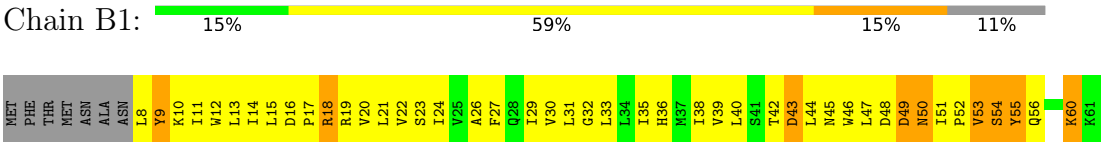
Chain BY:





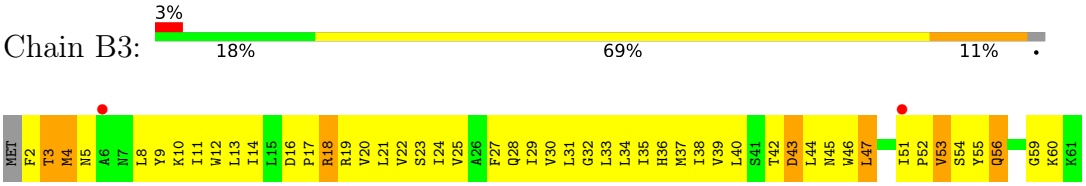
• Molecule 5: LH1 alpha polypeptide

Chain B1:



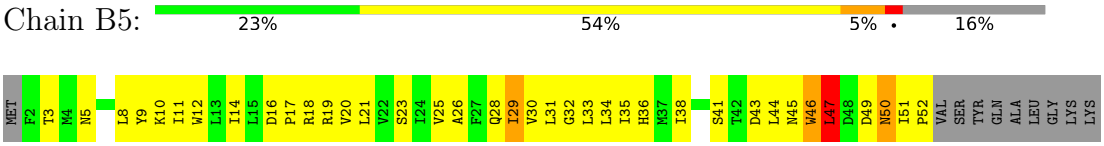
• Molecule 5: LH1 alpha polypeptide

Chain B3:



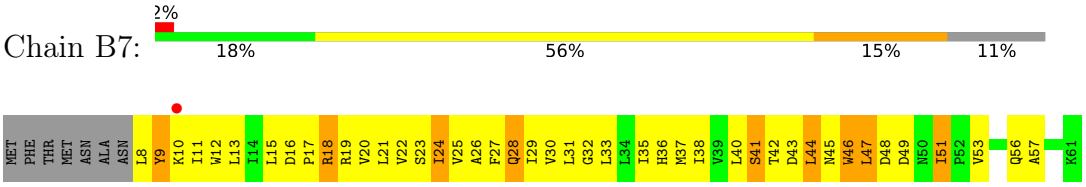
• Molecule 5: LH1 alpha polypeptide

Chain B5:



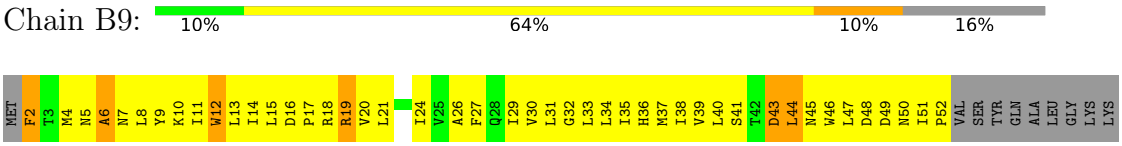
• Molecule 5: LH1 alpha polypeptide

Chain B7:



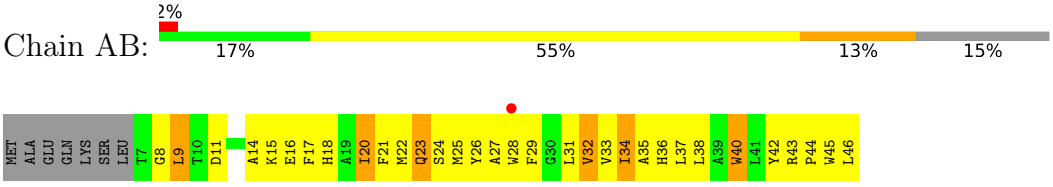
• Molecule 5: LH1 alpha polypeptide

Chain B9:



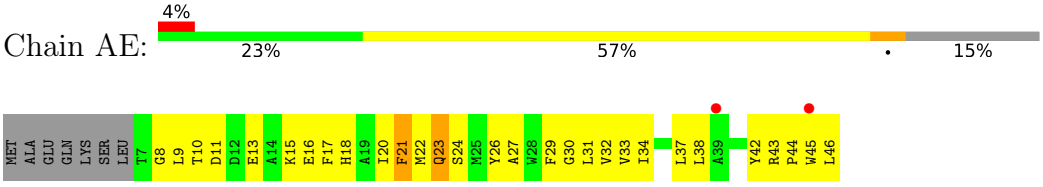
• Molecule 6: LH1 beta polypeptide

Chain AB:



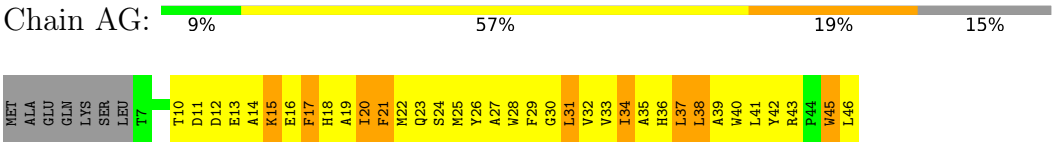
• Molecule 6: LH1 beta polypeptide

Chain AE:



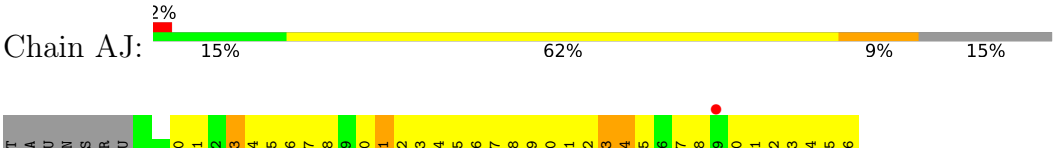
• Molecule 6: LH1 beta polypeptide

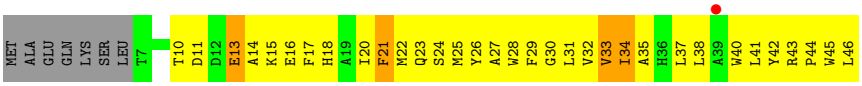
Chain AG:



• Molecule 6: LH1 beta polypeptide

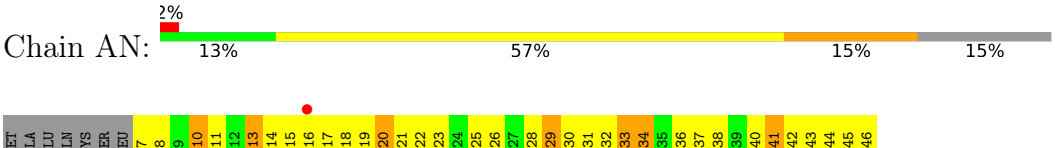
Chain AJ:

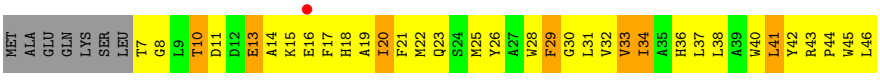




• Molecule 6: LH1 beta polypeptide

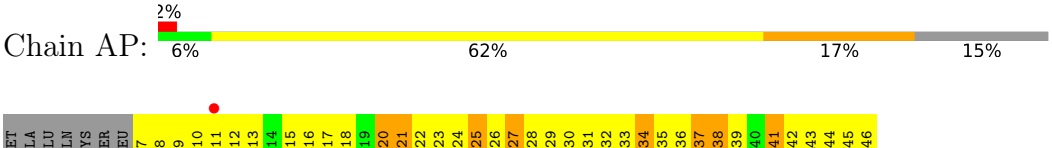
Chain AN:

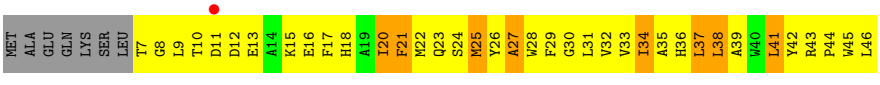




• Molecule 6: LH1 beta polypeptide

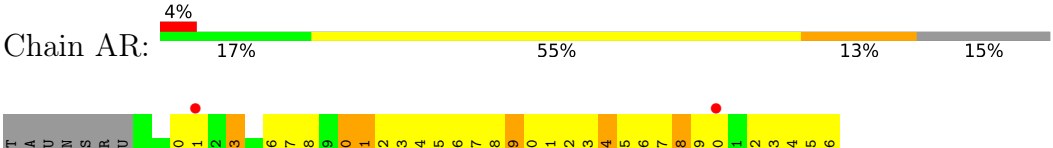
Chain AP:

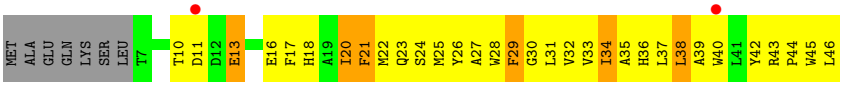




• Molecule 6: LH1 beta polypeptide

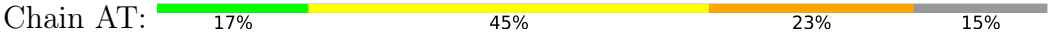
Chain AR:





• Molecule 6: LH1 beta polypeptide

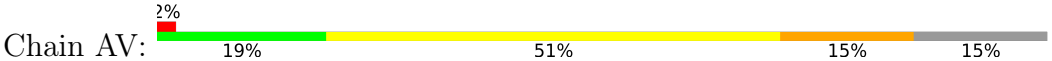
Chain AT:

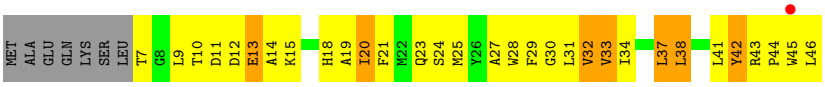




• Molecule 6: LH1 beta polypeptide

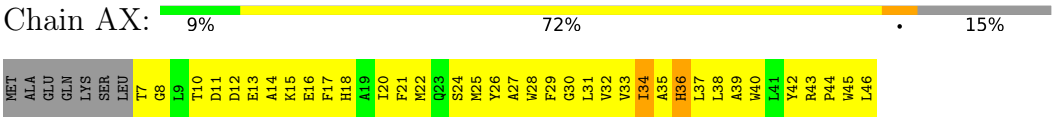
Chain AV:





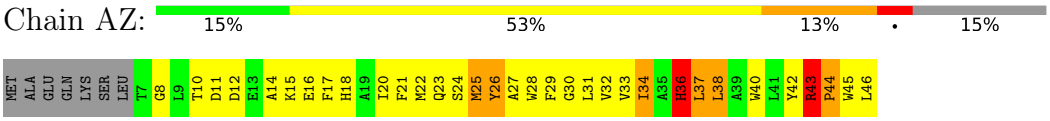
• Molecule 6: LH1 beta polypeptide

Chain AX:



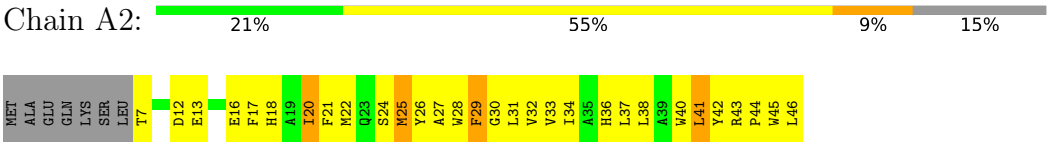
• Molecule 6: LH1 beta polypeptide

Chain AZ:



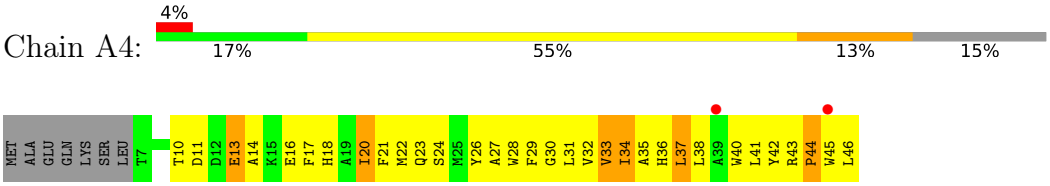
• Molecule 6: LH1 beta polypeptide

Chain A2:



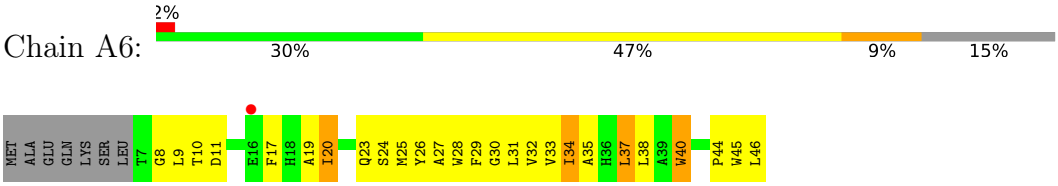
• Molecule 6: LH1 beta polypeptide

Chain A4:



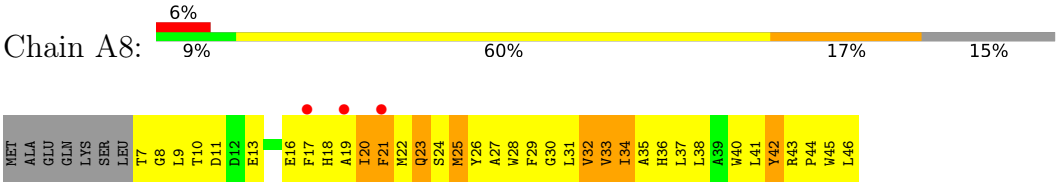
• Molecule 6: LH1 beta polypeptide

Chain A6:



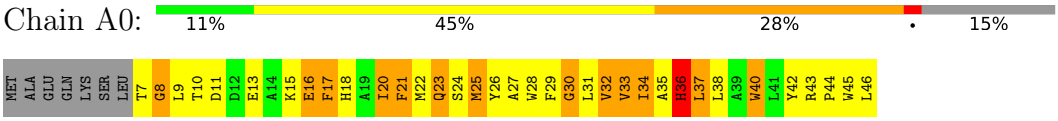
• Molecule 6: LH1 beta polypeptide

Chain A8:



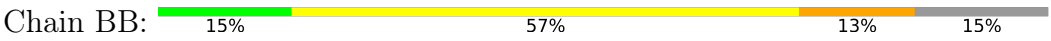
• Molecule 6: LH1 beta polypeptide

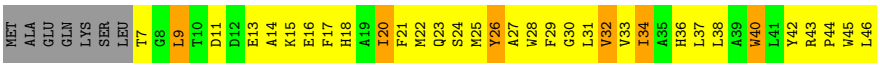
Chain A0:



• Molecule 6: LH1 beta polypeptide

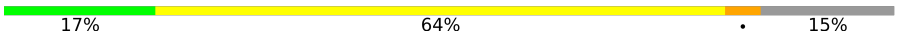
Chain BB:

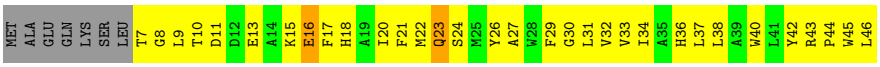




• Molecule 6: LH1 beta polypeptide

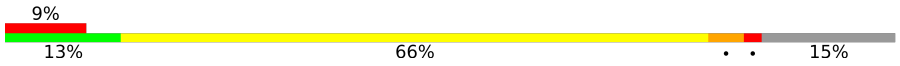
Chain BE:

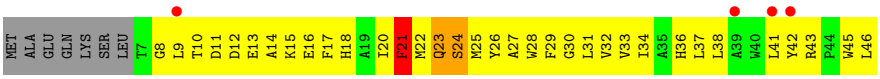




• Molecule 6: LH1 beta polypeptide

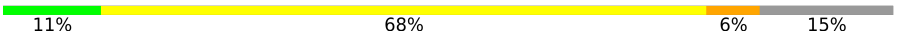
Chain BG:

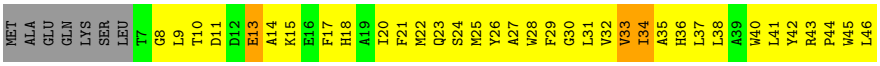




• Molecule 6: LH1 beta polypeptide

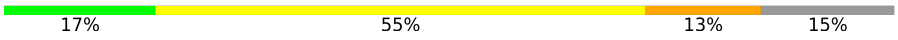
Chain BJ:

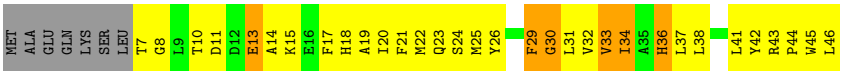




• Molecule 6: LH1 beta polypeptide

Chain BN:

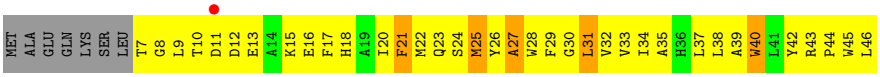




• Molecule 6: LH1 beta polypeptide

Chain BP:





• Molecule 6: LH1 beta polypeptide

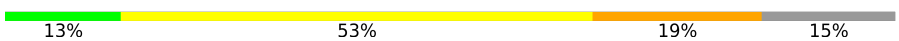
Chain BR:





• Molecule 6: LH1 beta polypeptide

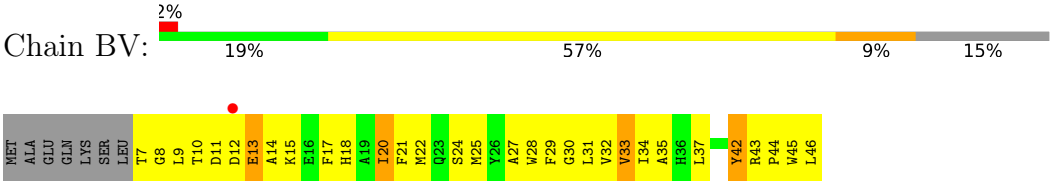
Chain BT:





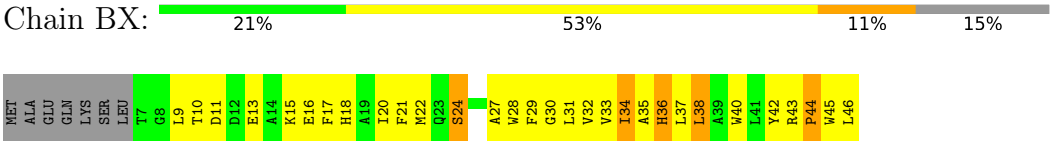
• Molecule 6: LH1 beta polypeptide

Chain BV:



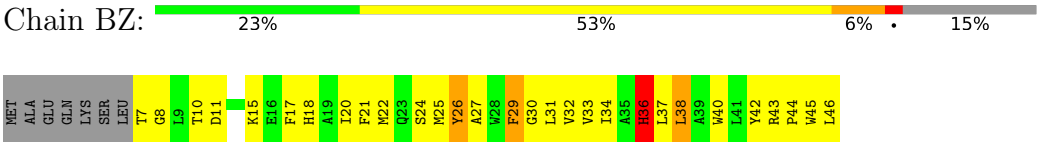
• Molecule 6: LH1 beta polypeptide

Chain BX:



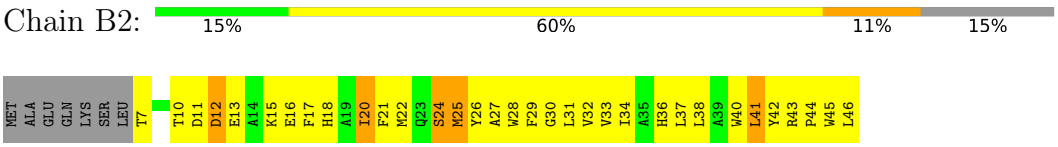
• Molecule 6: LH1 beta polypeptide

Chain BZ:



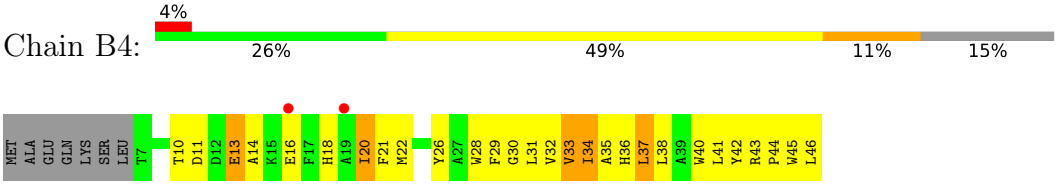
• Molecule 6: LH1 beta polypeptide

Chain B2:



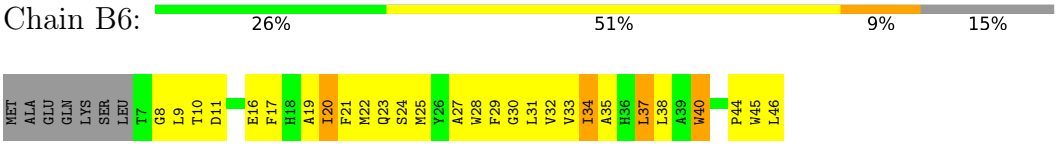
• Molecule 6: LH1 beta polypeptide

Chain B4:



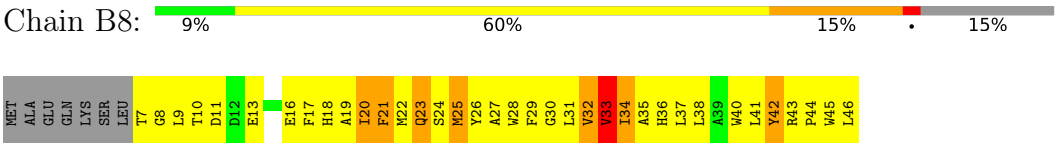
• Molecule 6: LH1 beta polypeptide

Chain B6:



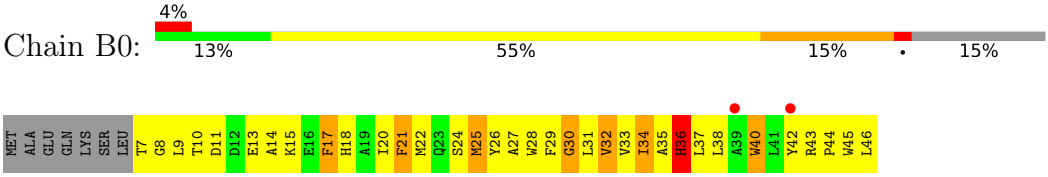
• Molecule 6: LH1 beta polypeptide

Chain B8:



• Molecule 6: LH1 beta polypeptide

Chain B0:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	167.16Å 145.43Å 210.53Å 90.00° 108.50° 90.00°	Depositor
Resolution (Å)	43.79 – 3.01 43.79 – 3.01	Depositor EDS
% Data completeness (in resolution range)	(Not available) (43.79-3.01) 69.3 (43.79-3.01)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
$< I/\sigma(I) >$ [1]	1.67 (at 3.01Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.1_1168)	Depositor
R, R_{free}	0.335 , 0.356 0.342 , 0.344	Depositor DCC
R_{free} test set	8241 reflections (4.92%)	wwPDB-VP
Wilson B-factor (Å^2)	55.8	Xtriage
Anisotropy	0.472	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.17 , 97.2	EDS
L-test for twinning[2]	$< \|L\| > = 0.35$, $< L^2 > = 0.18$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.88	EDS
Total number of atoms	50862	wwPDB-VP
Average B, all atoms (Å^2)	134.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.99% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality (i)

5.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: FE, BCL, PEF, PO4, MQ8, HEM, BPH, CRT, CA, UQ8

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	#$\|Z\|>5$	RMSZ	#$\|Z\|>5$
1	AC	0.75	1/2528 (0.0%)	1.20	27/3451 (0.8%)
1	BC	0.74	1/2528 (0.0%)	1.19	23/3451 (0.7%)
2	AL	0.43	0/2318	0.95	10/3167 (0.3%)
2	BL	0.38	0/2318	0.90	9/3167 (0.3%)
3	AM	0.42	0/2651	0.98	22/3628 (0.6%)
3	BM	0.41	0/2651	0.98	18/3628 (0.5%)
4	AH	0.40	0/2037	0.90	6/2776 (0.2%)
4	BH	0.38	0/2037	0.89	5/2776 (0.2%)
5	A1	0.48	0/464	1.21	6/634 (0.9%)
5	A3	0.40	0/456	1.07	5/624 (0.8%)
5	A5	0.47	0/453	0.98	2/620 (0.3%)
5	A7	0.38	0/426	0.89	2/583 (0.3%)
5	A9	0.40	0/483	1.11	7/660 (1.1%)
5	AA	0.41	0/401	0.89	0/550
5	AD	0.39	0/456	0.93	2/624 (0.3%)
5	AF	0.38	0/471	1.01	4/644 (0.6%)
5	AI	0.54	0/471	1.12	3/644 (0.5%)
5	AK	0.36	0/464	0.87	1/634 (0.2%)
5	AO	0.56	0/471	1.33	9/644 (1.4%)
5	AQ	0.38	0/456	0.94	1/624 (0.2%)
5	AS	0.42	0/471	1.06	3/644 (0.5%)
5	AU	0.46	0/483	0.96	3/660 (0.5%)
5	AW	0.41	0/483	0.93	1/660 (0.2%)
5	AY	0.47	0/483	1.06	5/660 (0.8%)
5	B1	0.42	0/435	0.88	0/595
5	B3	0.40	0/483	0.92	1/660 (0.2%)
5	B5	0.38	0/426	0.95	0/583
5	B7	0.41	0/435	0.88	1/595 (0.2%)
5	B9	0.50	0/426	0.94	2/583 (0.3%)
5	BA	0.40	0/458	0.90	1/627 (0.2%)
5	BD	0.63	1/378 (0.3%)	1.13	4/518 (0.8%)
5	BF	0.54	1/453 (0.2%)	1.04	3/620 (0.5%)

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	#|Z| >5	RMSZ	#|Z| >5
5	BI	0.44	0/418	0.90	1/573 (0.2%)
5	BK	0.38	0/483	0.82	1/660 (0.2%)
5	BO	0.53	0/471	1.18	6/644 (0.9%)
5	BQ	0.38	0/477	0.89	1/653 (0.2%)
5	BS	0.38	0/471	0.97	2/644 (0.3%)
5	BU	0.53	2/472 (0.4%)	0.94	3/646 (0.5%)
5	BW	0.44	0/464	0.94	2/634 (0.3%)
5	BY	0.42	0/435	0.79	1/595 (0.2%)
6	A0	0.56	0/350	0.86	2/476 (0.4%)
6	A2	0.35	0/350	0.76	0/476
6	A4	0.55	0/350	0.87	0/476
6	A6	0.39	0/350	0.92	0/476
6	A8	0.55	0/350	0.93	0/476
6	AB	0.46	0/350	0.79	0/476
6	AE	0.42	0/350	0.83	0/476
6	AG	0.55	0/350	0.97	0/476
6	AJ	0.46	0/350	0.84	0/476
6	AN	0.48	0/350	0.93	0/476
6	AP	0.51	0/350	0.89	1/476 (0.2%)
6	AR	0.43	0/350	0.85	0/476
6	AT	0.41	0/350	0.94	2/476 (0.4%)
6	AV	0.47	0/350	1.09	2/476 (0.4%)
6	AX	0.43	0/350	0.85	0/476
6	AZ	0.58	1/350 (0.3%)	1.06	6/476 (1.3%)
6	B0	0.55	0/350	0.87	2/476 (0.4%)
6	B2	0.46	0/350	0.90	1/476 (0.2%)
6	B4	0.54	0/350	0.92	0/476
6	B6	0.40	0/350	0.92	0/476
6	B8	0.55	0/350	0.93	1/476 (0.2%)
6	BB	0.49	0/350	0.86	0/476
6	BE	0.42	0/350	0.87	0/476
6	BG	0.54	0/350	1.09	1/476 (0.2%)
6	BJ	0.48	0/350	0.85	0/476
6	BN	0.55	0/350	1.05	2/476 (0.4%)
6	BP	0.50	0/350	0.85	1/476 (0.2%)
6	BR	0.45	0/350	0.87	0/476
6	BT	0.54	0/350	1.02	2/476 (0.4%)
6	BV	0.48	0/350	1.19	4/476 (0.8%)
6	BX	0.39	0/350	0.89	0/476
6	BZ	0.49	0/350	0.90	4/476 (0.8%)
All	All	0.49	7/44845 (0.0%)	0.98	234/61215 (0.4%)

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	BC	236	MET	SD-CE	-7.05	1.61	1.79
5	BD	37	MET	SD-CE	5.62	1.93	1.79
1	AC	40	MET	SD-CE	5.57	1.93	1.79
5	BU	17	PRO	N-CD	5.32	1.55	1.47
5	BF	37	MET	SD-CE	5.25	1.92	1.79
6	AZ	44	PRO	N-CD	5.23	1.55	1.47
5	BU	16	ASP	C-N	5.23	1.41	1.34

All (234) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	AC	301	ASP	N-CA-C	10.60	117.50	108.07
5	BO	21	LEU	N-CA-C	9.99	123.20	111.02
5	AY	49	ASP	N-CA-C	9.36	121.09	111.07
5	A1	8	LEU	N-CA-C	9.18	122.44	111.33
3	AM	122	LEU	N-CA-C	-9.07	99.96	112.30
1	AC	178	LEU	N-CA-C	9.05	126.58	107.67
6	BV	42	TYR	N-CA-C	8.94	120.95	111.03
1	AC	305	VAL	N-CA-C	8.84	122.10	107.24
5	BD	41	SER	N-CA-C	8.84	120.53	111.07
5	AO	8	LEU	N-CA-C	8.68	120.74	111.28
5	AO	6	ALA	N-CA-C	8.64	120.31	111.07
1	AC	275	HIS	N-CA-C	-8.48	102.19	112.54
5	AI	51	ILE	N-CA-C	8.38	117.59	107.61
1	BC	275	HIS	N-CA-C	-8.38	102.32	112.54
5	AO	5	ASN	N-CA-C	8.36	120.39	111.28
3	AM	146	LEU	N-CA-C	-8.19	98.35	111.04
5	AO	36	HIS	N-CA-C	-8.09	102.41	111.71
5	A1	11	ILE	N-CA-C	8.08	118.86	110.62
1	BC	305	VAL	N-CA-C	8.03	120.73	107.24
1	BC	73	SER	N-CA-C	-7.92	102.60	111.71
1	AC	73	SER	N-CA-C	-7.88	102.67	111.82
6	BN	30	GLY	N-CA-C	-7.84	103.40	112.50
1	AC	178	LEU	CB-CA-C	-7.77	100.98	112.09
1	BC	327	TYR	CA-C-N	7.76	127.81	119.28
1	BC	327	TYR	C-N-CA	7.76	127.81	119.28
1	AC	226	LEU	N-CA-C	-7.65	101.93	111.11
5	BS	45	ASN	N-CA-C	-7.63	100.25	110.55
3	BM	122	LEU	N-CA-C	-7.56	102.28	112.94
2	AL	160	LEU	N-CA-C	-7.52	105.28	114.75
3	AM	163	ILE	N-CA-C	7.40	120.34	111.09
5	BO	16	ASP	CA-C-N	7.31	126.58	118.97
5	BO	16	ASP	C-N-CA	7.31	126.58	118.97

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
5	AO	35	ILE	N-CA-C	-7.29	104.75	111.67
3	BM	130	TRP	N-CA-C	-7.16	103.21	112.23
6	AP	27	ALA	N-CA-C	-7.16	103.48	111.28
5	BD	35	ILE	N-CA-C	7.04	117.13	110.30
1	BC	159	ASN	CA-C-N	7.03	126.55	119.24
1	BC	159	ASN	C-N-CA	7.03	126.55	119.24
3	BM	140	LEU	N-CA-C	-7.02	103.83	112.88
5	AF	24	ILE	N-CA-C	-7.00	105.02	111.67
4	AH	35	LYS	N-CA-C	6.97	121.04	112.54
1	BC	178	LEU	N-CA-C	6.93	122.19	107.70
5	BQ	48	ASP	N-CA-C	6.93	122.34	109.24
3	BM	268	TRP	N-CA-C	-6.86	103.03	111.33
3	BM	155	PHE	N-CA-C	-6.80	103.10	111.40
5	AF	39	VAL	N-CA-C	6.74	116.86	110.53
5	A7	6	ALA	N-CA-C	6.74	119.19	111.11
3	AM	123	THR	N-CA-C	-6.70	104.98	113.43
4	AH	160	ASP	N-CA-C	-6.70	105.11	113.28
3	AM	30	ARG	N-CA-C	-6.70	102.17	110.41
1	AC	327	TYR	CA-C-N	6.64	126.59	119.28
1	AC	327	TYR	C-N-CA	6.64	126.59	119.28
3	BM	146	LEU	N-CA-C	-6.63	102.68	111.24
3	BM	202	HIS	N-CA-C	-6.57	104.19	112.93
6	AV	12	ASP	N-CA-C	6.57	118.09	111.07
3	AM	202	HIS	N-CA-C	-6.53	104.25	112.93
4	BH	43	SER	N-CA-C	6.51	120.16	109.94
5	A3	50	ASN	N-CA-C	6.51	116.58	108.19
1	AC	269	ALA	N-CA-C	-6.49	105.89	113.88
6	BP	27	ALA	N-CA-C	-6.48	103.49	111.33
5	AU	41	SER	N-CA-C	6.46	118.11	111.14
5	BF	5	ASN	N-CA-C	-6.45	105.23	113.23
1	AC	104	LYS	N-CA-C	-6.45	105.43	113.55
3	AM	228	ARG	N-CA-C	-6.45	104.33	111.36
6	BV	12	ASP	N-CA-C	6.44	117.97	111.07
1	BC	123	THR	N-CA-C	-6.44	104.18	111.14
5	A9	12	TRP	N-CA-C	6.44	120.59	112.87
2	BL	123	GLY	N-CA-C	-6.44	102.79	113.02
3	BM	23	LEU	CA-C-N	6.42	127.87	119.84
3	BM	23	LEU	C-N-CA	6.42	127.87	119.84
3	AM	130	TRP	N-CA-C	-6.39	104.75	112.54
3	AM	140	LEU	N-CA-C	-6.38	104.04	113.61
1	BC	277	ARG	N-CA-C	-6.37	103.46	111.11
5	BW	50	ASN	N-CA-C	6.37	119.59	109.65

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
5	A1	9	TYR	N-CA-C	6.37	120.15	112.38
4	AH	238	LYS	N-CA-C	6.29	117.80	111.07
6	AZ	43	ARG	CA-C-N	-6.28	111.99	119.84
6	AZ	43	ARG	C-N-CA	-6.28	111.99	119.84
3	AM	23	LEU	CA-C-N	6.25	127.65	119.84
3	AM	23	LEU	C-N-CA	6.25	127.65	119.84
5	B9	12	TRP	N-CA-C	6.25	119.75	111.75
5	BO	36	HIS	N-CA-C	-6.20	104.58	111.71
3	BM	30	ARG	N-CA-C	-6.20	102.35	110.53
1	AC	123	THR	N-CA-C	-6.18	104.45	111.07
1	BC	152	CYS	N-CA-C	-6.18	105.22	112.89
5	AO	11	ILE	N-CA-C	6.15	116.90	110.62
5	AY	36	HIS	N-CA-C	-6.14	103.89	111.33
5	BU	20	VAL	N-CA-C	6.09	116.26	110.53
6	B2	12	ASP	N-CA-C	6.09	117.72	111.14
5	A9	13	LEU	N-CA-C	-6.06	104.23	111.69
5	BF	42	THR	N-CA-C	6.06	117.68	111.14
1	BC	104	LYS	N-CA-C	-6.04	105.95	113.55
6	BV	33	VAL	N-CA-C	-6.03	104.47	110.62
5	A3	12	TRP	N-CA-C	6.00	120.33	112.89
6	BT	27	ALA	N-CA-C	-6.00	104.82	111.36
2	BL	100	ILE	N-CA-C	-6.00	104.48	110.30
2	BL	15	GLY	N-CA-C	-5.99	106.37	115.00
2	AL	226	ARG	N-CA-C	-5.98	106.07	113.19
5	A9	55	TYR	N-CA-C	5.96	117.78	111.28
3	AM	56	THR	N-CA-C	-5.96	104.78	111.28
5	AS	9	TYR	N-CA-C	-5.95	104.73	112.23
3	BM	228	ARG	N-CA-C	-5.95	104.37	111.69
4	BH	138	VAL	N-CA-C	-5.94	104.24	112.50
2	AL	255	VAL	N-CA-C	-5.94	104.54	110.30
6	AV	33	VAL	N-CA-C	-5.94	104.56	110.62
5	AS	4	MET	N-CA-C	5.92	117.41	111.07
1	BC	229	ALA	N-CA-C	-5.90	104.93	111.71
5	A7	5	ASN	N-CA-C	5.88	118.44	111.33
5	BS	55	TYR	N-CA-C	-5.88	104.95	111.71
4	BH	160	ASP	N-CA-C	-5.87	106.08	113.18
6	AT	27	ALA	N-CA-C	-5.85	104.98	111.36
4	BH	159	LEU	N-CA-C	-5.84	105.42	112.54
5	AI	41	SER	N-CA-C	5.84	118.11	111.11
5	AS	45	ASN	N-CA-C	-5.84	102.78	110.43
1	AC	307	CYS	N-CA-C	-5.83	105.05	111.82
5	BD	45	ASN	N-CA-C	-5.82	102.67	110.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
5	BF	20	VAL	N-CA-C	5.82	116.00	110.53
3	AM	159	VAL	N-CA-C	-5.81	105.72	112.98
1	AC	92	ARG	N-CA-C	-5.80	104.15	111.11
5	AY	42	THR	N-CA-C	5.79	118.03	108.48
3	BM	137	ALA	N-CA-C	5.76	117.24	110.97
5	BD	29	ILE	CB-CG1-CD1	-5.76	101.71	113.80
1	BC	276	VAL	N-CA-C	-5.73	106.12	111.45
3	AM	141	GLY	N-CA-C	-5.72	107.35	115.30
5	AQ	42	THR	N-CA-C	5.70	119.76	112.12
5	AO	43	ASP	N-CA-C	5.67	117.92	111.11
1	AC	318	LEU	N-CA-C	-5.67	105.66	112.58
5	BW	47	LEU	N-CA-C	5.66	117.31	111.03
1	BC	57	GLN	CA-C-N	5.64	126.12	120.14
1	BC	57	GLN	C-N-CA	5.64	126.12	120.14
2	BL	195	ALA	N-CA-C	-5.64	104.14	111.02
1	BC	119	ASP	N-CA-C	-5.61	103.70	110.44
6	BG	21	PHE	N-CA-CB	-5.60	101.85	109.98
1	AC	101	VAL	CA-C-N	-5.60	116.54	123.21
1	AC	101	VAL	C-N-CA	-5.60	116.54	123.21
1	AC	229	ALA	N-CA-C	-5.59	105.34	111.82
5	AW	13	LEU	N-CA-C	5.57	117.03	111.07
5	A1	51	ILE	N-CA-C	5.55	120.88	108.88
5	A5	42	THR	N-CA-C	5.55	119.56	112.12
5	AF	42	THR	N-CA-C	5.54	117.13	111.14
2	AL	36	GLY	N-CA-C	-5.52	106.01	113.24
6	AZ	38	LEU	N-CA-C	-5.51	105.81	112.54
2	BL	199	HIS	N-CA-C	-5.50	105.18	111.07
5	BO	8	LEU	N-CA-C	5.49	119.46	112.87
3	AM	67	ALA	N-CA-C	-5.48	104.69	111.33
3	AM	55	LEU	N-CA-C	-5.47	104.24	111.96
1	BC	178	LEU	CB-CA-C	-5.47	102.75	111.17
6	BT	26	TYR	N-CA-C	-5.46	105.87	112.54
5	AY	50	ASN	N-CA-C	5.46	122.43	110.80
3	BM	96	GLU	CA-C-N	5.46	126.00	120.38
3	BM	96	GLU	C-N-CA	5.46	126.00	120.38
6	AZ	8	GLY	N-CA-C	-5.45	107.49	115.72
5	AO	14	ILE	N-CA-C	5.45	117.86	111.05
3	BM	159	VAL	N-CA-C	-5.44	105.93	113.00
5	AU	6	ALA	N-CA-C	-5.43	107.30	114.31
1	AC	207	ALA	N-CA-C	5.42	116.81	110.41
5	A3	37	MET	N-CA-C	5.42	117.00	111.14
2	BL	54	ALA	N-CA-C	-5.42	104.82	111.75

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
6	BZ	38	LEU	N-CA-C	-5.42	105.93	112.54
1	BC	186	GLY	N-CA-C	-5.41	100.35	113.18
2	AL	258	LEU	N-CA-C	-5.41	105.81	112.90
5	B9	19	ARG	N-CA-C	5.40	116.85	111.07
6	B0	36	HIS	N-CA-C	-5.40	105.43	112.23
5	B3	18	ARG	N-CA-C	5.39	116.97	111.14
6	BZ	8	GLY	N-CA-C	-5.39	107.58	115.72
5	AF	55	TYR	N-CA-C	-5.38	106.56	113.23
5	AO	11	ILE	CB-CA-C	-5.38	104.88	112.14
5	A1	5	ASN	N-CA-C	5.38	118.94	112.38
6	BZ	36	HIS	N-CA-C	-5.37	105.42	111.28
5	BU	16	ASP	CA-C-N	-5.37	113.09	119.05
5	BU	16	ASP	C-N-CA	-5.37	113.09	119.05
5	AD	27	PHE	N-CA-C	5.36	117.12	111.28
3	AM	212	SER	N-CA-C	-5.33	105.36	111.07
6	AZ	36	HIS	N-CA-C	-5.33	105.47	111.28
6	AT	26	TYR	N-CA-C	-5.32	106.05	112.54
3	AM	86	PHE	N-CA-C	-5.32	104.89	111.33
5	A9	41	SER	N-CA-C	5.32	117.08	111.28
1	AC	75	VAL	N-CA-C	5.32	116.68	110.62
5	A5	19	ARG	N-CA-C	5.31	116.76	110.97
1	AC	119	ASP	N-CA-C	-5.29	103.74	111.30
5	BY	12	TRP	N-CA-C	5.28	119.35	113.01
5	A3	20	VAL	N-CA-C	5.28	115.49	110.53
2	AL	49	LEU	N-CA-C	-5.27	105.54	111.28
4	AH	138	VAL	N-CA-C	-5.27	104.38	111.44
2	AL	220	HIS	N-CA-C	-5.27	105.06	112.12
5	BO	43	ASP	N-CA-C	5.27	117.43	111.11
2	AL	150	ALA	N-CA-C	5.26	116.99	108.26
6	A0	36	HIS	N-CA-C	-5.25	105.61	112.23
1	BC	207	ALA	N-CA-C	5.25	116.60	110.41
2	AL	194	LEU	N-CA-C	-5.25	105.09	112.12
3	BM	253	ARG	N-CA-C	-5.24	105.25	111.69
2	AL	248	SER	N-CA-C	-5.23	104.59	111.96
5	AU	11	ILE	N-CA-C	5.21	120.18	109.34
5	A9	42	THR	N-CA-C	5.21	119.10	112.12
5	A1	48	ASP	N-CA-C	5.20	118.94	112.59
3	AM	145	HIS	N-CA-C	-5.19	108.21	114.75
3	AM	96	GLU	CA-C-N	5.18	125.72	120.38
3	AM	96	GLU	C-N-CA	5.18	125.72	120.38
3	AM	14	ARG	N-CA-C	5.18	117.12	108.99
5	AD	47	LEU	N-CA-C	5.17	117.33	111.02

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	AC	78	ASN	N-CA-C	5.17	115.87	107.23
1	BC	75	VAL	N-CA-C	5.17	115.89	110.62
5	BK	49	ASP	N-CA-C	5.17	116.60	111.07
5	A3	18	ARG	N-CA-C	5.17	116.72	111.14
1	BC	318	LEU	N-CA-C	-5.16	106.28	112.58
1	AC	251	HIS	N-CA-C	5.16	118.14	110.14
5	AY	12	TRP	N-CA-C	5.15	119.05	112.87
1	AC	277	ARG	N-CA-C	-5.15	102.87	111.37
6	BZ	26	TYR	N-CA-C	-5.15	105.79	111.71
5	B7	51	ILE	N-CA-C	5.12	114.54	108.96
1	BC	252	ASN	N-CA-C	-5.10	99.94	110.80
3	BM	203	MET	N-CA-C	-5.10	106.75	112.87
2	BL	160	LEU	N-CA-C	-5.10	106.89	114.64
1	AC	186	GLY	N-CA-C	-5.09	101.11	113.18
1	AC	20	LEU	N-CA-C	5.08	116.92	109.24
4	AH	159	LEU	N-CA-C	-5.08	106.91	113.01
2	BL	64	ASP	N-CA-C	-5.08	106.91	113.01
5	AK	50	ASN	N-CA-C	5.08	114.74	108.19
5	AI	23	SER	N-CA-C	-5.08	105.44	110.97
5	BA	49	ASP	N-CA-C	5.07	116.49	111.07
6	A0	30	GLY	N-CA-C	-5.07	106.51	112.64
2	BL	112	ARG	N-CA-C	-5.05	105.05	111.11
6	BV	42	TYR	CB-CA-C	-5.05	103.12	110.95
6	B8	33	VAL	N-CA-C	-5.04	105.07	110.36
3	BM	56	THR	N-CA-C	-5.04	105.92	111.71
5	A9	6	ALA	N-CA-C	5.03	119.67	111.37
5	BI	47	LEU	N-CA-C	5.03	116.57	111.14
6	BN	36	HIS	N-CA-C	-5.03	106.98	113.01
6	B0	30	GLY	N-CA-C	-5.02	106.56	112.64
5	A9	51	ILE	N-CA-C	5.02	114.68	109.01
4	BH	42	ASP	N-CA-C	5.01	120.29	110.56
1	AC	252	ASN	N-CA-C	-5.01	104.14	111.30
4	AH	23	PHE	N-CA-C	-5.01	105.47	111.03
6	AZ	26	TYR	N-CA-C	-5.01	105.95	111.71

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (i)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen

atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	AC	2458	0	2377	508	0
1	BC	2458	0	2377	514	0
2	AL	2231	0	2192	677	0
2	BL	2231	0	2192	598	0
3	AM	2551	0	2526	779	0
3	BM	2551	0	2526	699	0
4	AH	1982	0	1981	421	0
4	BH	1982	0	1981	392	0
5	A1	455	0	460	169	0
5	A3	447	0	451	136	0
5	A5	444	0	456	141	0
5	A7	417	0	441	163	0
5	A9	473	0	476	130	0
5	AA	392	0	412	144	0
5	AD	447	0	451	129	0
5	AF	462	0	467	160	0
5	AI	462	0	467	154	0
5	AK	455	0	460	94	0
5	AO	462	0	467	160	0
5	AQ	447	0	451	123	0
5	AS	462	0	467	190	0
5	AU	473	0	476	155	0
5	AW	473	0	476	189	0
5	AY	473	0	476	170	0
5	B1	426	0	434	144	0
5	B3	473	0	476	139	0
5	B5	417	0	441	95	0
5	B7	426	0	434	136	0
5	B9	417	0	441	146	0
5	BA	448	0	462	159	0
5	BD	370	0	399	139	0
5	BF	444	0	456	150	0
5	BI	409	0	426	112	0
5	BK	473	0	476	112	0
5	BO	462	0	467	155	0
5	BQ	467	0	474	124	0
5	BS	462	0	467	118	0
5	BU	462	0	472	180	0
5	BW	455	0	460	163	0
5	BY	426	0	434	154	0
6	A0	337	0	323	115	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	A2	337	0	323	107	0
6	A4	337	0	323	88	0
6	A6	337	0	323	63	0
6	A8	337	0	323	111	0
6	AB	337	0	321	94	0
6	AE	337	0	323	72	0
6	AG	337	0	323	92	0
6	AJ	337	0	323	97	0
6	AN	337	0	323	87	0
6	AP	337	0	323	110	0
6	AR	337	0	323	78	0
6	AT	337	0	323	81	0
6	AV	337	0	323	101	0
6	AX	337	0	323	88	0
6	AZ	337	0	323	102	0
6	B0	337	0	323	128	0
6	B2	337	0	323	135	0
6	B4	337	0	323	73	0
6	B6	337	0	323	56	0
6	B8	337	0	323	79	0
6	BB	337	0	323	105	0
6	BE	337	0	323	75	0
6	BG	337	0	323	81	0
6	BJ	337	0	323	97	0
6	BN	337	0	323	70	0
6	BP	337	0	323	97	0
6	BR	337	0	323	73	0
6	BT	337	0	323	66	0
6	BV	337	0	323	110	0
6	BX	337	0	323	76	0
6	BZ	337	0	323	86	0
7	AC	172	0	120	37	0
7	BC	172	0	120	34	0
8	A1	1	0	0	0	0
8	A3	1	0	0	0	0
8	A5	1	0	0	0	0
8	A7	1	0	0	0	0
8	A9	1	0	0	0	0
8	AA	1	0	0	0	0
8	AC	1	0	0	0	0
8	AD	1	0	0	0	0
8	AF	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
8	AI	1	0	0	0	0
8	AK	1	0	0	0	0
8	AO	1	0	0	0	0
8	AQ	1	0	0	0	0
8	AS	1	0	0	0	0
8	AU	1	0	0	0	0
8	AV	1	0	0	0	0
8	AY	1	0	0	0	0
8	B1	1	0	0	0	0
8	B3	1	0	0	0	0
8	B5	1	0	0	0	0
8	B7	1	0	0	0	0
8	B9	1	0	0	0	0
8	BA	1	0	0	0	0
8	BC	1	0	0	0	0
8	BD	1	0	0	0	0
8	BF	1	0	0	0	0
8	BI	1	0	0	0	0
8	BK	1	0	0	0	0
8	BO	1	0	0	0	0
8	BQ	1	0	0	0	0
8	BS	1	0	0	0	0
8	BU	1	0	0	0	0
8	BW	1	0	0	0	0
8	BY	1	0	0	0	0
9	A0	66	0	72	69	0
9	A1	66	0	74	52	0
9	A2	66	0	74	28	0
9	A3	132	0	148	60	0
9	A5	66	0	74	34	0
9	A6	66	0	74	27	0
9	A7	66	0	74	52	0
9	A8	66	0	74	46	0
9	A9	66	0	74	36	0
9	AA	66	0	74	42	0
9	AB	66	0	74	35	0
9	AD	66	0	74	27	0
9	AE	66	0	74	32	0
9	AF	66	0	74	38	0
9	AG	66	0	74	36	0
9	AI	66	0	74	36	0
9	AJ	66	0	74	43	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
9	AK	66	0	74	58	0
9	AL	132	0	148	55	0
9	AM	132	0	148	41	0
9	AN	66	0	74	53	0
9	AO	66	0	74	50	0
9	AP	66	0	74	37	0
9	AQ	66	0	74	25	0
9	AR	66	0	74	34	0
9	AS	66	0	74	32	0
9	AT	66	0	74	25	0
9	AU	66	0	73	46	0
9	AV	66	0	74	30	0
9	AW	66	0	72	40	0
9	AX	66	0	74	33	0
9	AY	66	0	74	48	0
9	AZ	66	0	72	41	0
9	B0	66	0	74	57	0
9	B1	66	0	74	37	0
9	B2	66	0	74	43	0
9	B3	66	0	74	47	0
9	B4	66	0	74	20	0
9	B5	66	0	74	20	0
9	B6	66	0	74	24	0
9	B7	66	0	74	39	0
9	B8	66	0	74	36	0
9	B9	66	0	74	38	0
9	BA	66	0	74	38	0
9	BB	66	0	74	47	0
9	BD	66	0	74	34	0
9	BE	66	0	74	46	0
9	BF	66	0	74	39	0
9	BG	66	0	74	46	0
9	BI	66	0	74	50	0
9	BJ	66	0	74	26	0
9	BK	66	0	74	26	0
9	BL	132	0	148	42	0
9	BM	132	0	148	44	0
9	BN	66	0	74	30	0
9	BO	66	0	74	48	0
9	BP	66	0	74	35	0
9	BQ	132	0	148	51	0
9	BS	66	0	74	28	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
9	BT	66	0	74	22	0
9	BU	66	0	74	38	0
9	BV	66	0	74	23	0
9	BW	66	0	74	42	0
9	BX	66	0	74	33	0
9	BY	66	0	74	35	0
9	BZ	66	0	74	31	0
10	AL	65	0	76	17	0
10	AM	65	0	76	12	0
10	BL	65	0	76	11	0
10	BM	65	0	76	7	0
11	AL	53	0	74	8	0
11	BL	53	0	74	12	0
12	AM	1	0	0	0	0
12	BM	1	0	0	0	0
13	AM	53	0	72	10	0
13	BM	53	0	72	8	0
14	A0	44	0	60	28	0
14	A1	44	0	60	29	0
14	A2	44	0	60	51	0
14	A5	44	0	60	29	0
14	A7	44	0	60	46	0
14	AA	44	0	60	24	0
14	AB	44	0	60	46	0
14	AG	44	0	60	11	0
14	AJ	44	0	60	21	0
14	AM	44	0	60	13	0
14	AN	44	0	60	15	0
14	AP	44	0	60	24	0
14	AR	44	0	60	20	0
14	AS	44	0	60	77	0
14	AT	44	0	60	18	0
14	AW	44	0	60	32	0
14	AX	44	0	60	47	0
14	B0	44	0	60	44	0
14	B1	44	0	60	41	0
14	B2	44	0	60	75	0
14	B5	44	0	60	22	0
14	B7	44	0	60	39	0
14	BA	44	0	60	26	0
14	BB	44	0	60	27	0
14	BF	44	0	60	19	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
14	BG	44	0	60	13	0
14	BM	44	0	60	8	0
14	BN	44	0	60	15	0
14	BO	44	0	60	13	0
14	BP	44	0	60	29	0
14	BS	44	0	60	10	0
14	BU	44	0	60	61	0
14	BV	44	0	60	59	0
14	BW	44	0	60	20	0
15	AH	19	0	11	8	0
15	AM	80	0	92	23	0
15	AS	47	0	73	37	0
15	BM	19	0	11	2	0
15	BQ	47	0	73	12	0
16	A3	5	0	0	0	0
16	AH	5	0	0	1	0
16	AM	5	0	0	0	0
16	BH	5	0	0	1	0
17	AA	1	0	0	0	0
17	AC	1	0	0	0	0
17	AH	2	0	0	0	0
17	AI	1	0	0	0	0
17	AL	3	0	0	2	0
17	AM	3	0	0	2	0
17	AW	1	0	0	1	0
17	B1	1	0	0	0	0
17	BC	1	0	0	0	0
17	BH	1	0	0	0	0
17	BL	3	0	0	3	0
17	BM	3	0	0	0	0
All	All	50862	0	51516	11377	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 111.

All (11377) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
14:A7:102:CRT:C22	14:A7:102:CRT:C21	1.82	1.56
14:AS:104:CRT:C9	6:AV:20:ILE:HD12	1.38	1.54
5:AS:30:VAL:HG22	15:AS:101:PEF:C41	1.47	1.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:AS:30:VAL:CG2	15:AS:101:PEF:H412	1.46	1.43
9:AW:101:BCL:O2A	9:AW:101:BCL:C1	1.65	1.42
5:BU:11:ILE:HG23	14:BU:103:CRT:C8	1.47	1.41
14:BV:102:CRT:H393	5:BW:36:HIS:CB	1.52	1.39
5:B1:13:LEU:HD12	14:B1:103:CRT:C2	1.53	1.38
14:AB:102:CRT:C2	5:A9:13:LEU:HD12	1.51	1.37
14:AB:102:CRT:C3	5:A9:10:LYS:HB3	1.55	1.34
5:BU:11:ILE:CG2	14:BU:103:CRT:H83	1.57	1.34
9:A7:103:BCL:O2A	9:A7:103:BCL:C1	1.75	1.33
5:A5:25:VAL:HG11	9:A5:102:BCL:C19	1.60	1.32
9:A0:102:BCL:O2A	9:A0:102:BCL:C1	1.76	1.31
5:BU:14:ILE:CG1	14:BU:103:CRT:H33	1.59	1.30
6:B2:21:PHE:CE1	14:B2:102:CRT:H16	1.66	1.30
5:BU:10:LYS:O	14:BU:103:CRT:H23	1.28	1.30
6:AP:38:LEU:O	6:AP:41:LEU:HD23	1.32	1.29
9:AX:101:BCL:O2A	9:AX:101:BCL:C1	1.77	1.29
9:A6:101:BCL:O2A	9:A6:101:BCL:C1	1.79	1.29
5:BU:14:ILE:HG13	14:BU:103:CRT:C3	1.63	1.28
5:B1:13:LEU:CD1	14:B1:103:CRT:H23	1.63	1.26
14:AB:102:CRT:H33	5:A9:10:LYS:CB	1.68	1.24
14:BV:102:CRT:C39	5:BW:36:HIS:HB3	1.65	1.24
14:A2:102:CRT:C2M	5:A3:40:LEU:HD11	1.64	1.24
6:BZ:46:LEU:HD22	6:B2:42:TYR:OH	1.34	1.24
5:BA:36:HIS:CB	14:B0:101:CRT:H391	1.66	1.23
6:B8:27:ALA:O	6:B8:31:LEU:HG	1.35	1.23
6:BV:21:PHE:CD1	14:BV:102:CRT:H14	1.74	1.23
5:AS:30:VAL:HG21	15:AS:101:PEF:C39	1.68	1.22
5:BO:43:ASP:HA	5:BQ:48:ASP:CB	1.67	1.22
5:B1:11:ILE:HA	14:B1:103:CRT:C8	1.69	1.22
14:BU:103:CRT:H2M1	5:BY:37:MET:N	1.53	1.21
5:A7:36:HIS:CB	14:A7:102:CRT:H2M3	1.71	1.20
6:A8:27:ALA:O	6:A8:31:LEU:HG	1.35	1.20
9:A1:102:BCL:O2A	9:A1:102:BCL:C1	1.89	1.19
6:AV:27:ALA:O	6:AV:31:LEU:HG	1.42	1.19
5:AF:28:GLN:HB3	9:AF:102:BCL:C1	1.74	1.18
6:B2:21:PHE:CD1	14:B2:102:CRT:H14	1.78	1.18
6:A0:45:TRP:HE1	9:A0:102:BCL:C19	1.56	1.17
6:B2:17:PHE:HB2	14:B2:102:CRT:H41	1.24	1.17
9:A0:102:BCL:H162	9:A0:102:BCL:HBB2	1.25	1.17
5:A1:8:LEU:HD23	5:A1:9:TYR:H	1.08	1.16
5:A5:25:VAL:HG11	9:A5:102:BCL:H191	1.17	1.16

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:B2:29:PHE:CE1	9:B2:101:BCL:H11	1.79	1.16
6:A2:17:PHE:CD1	14:A2:102:CRT:H6	1.81	1.15
5:B3:43:ASP:HB2	5:B5:47:LEU:HD13	1.19	1.15
2:AL:89:LEU:HA	2:AL:94:LEU:H	1.09	1.15
14:AS:104:CRT:C18	9:AU:102:BCL:H92	1.76	1.15
14:BP:102:CRT:H342	9:BQ:103:BCL:HBA1	1.17	1.15
14:B7:102:CRT:C34	9:B7:103:BCL:HBA1	1.75	1.15
2:AL:266:ARG:HB2	2:AL:266:ARG:HH11	1.09	1.15
6:AZ:46:LEU:CB	5:A1:52:PRO:HD3	1.76	1.15
6:AG:32:VAL:HG11	9:AG:101:BCL:HBA2	1.24	1.15
14:BV:102:CRT:H393	5:BW:36:HIS:HB2	1.25	1.14
6:B0:17:PHE:HB2	14:B0:101:CRT:H6	1.26	1.14
3:AM:241:ARG:O	4:AH:119:ARG:HD3	1.46	1.14
5:AI:19:ARG:O	5:AI:23:SER:HB3	1.47	1.14
14:AP:102:CRT:H342	9:AQ:102:BCL:HBA1	1.20	1.14
6:AP:32:VAL:HG11	9:AP:101:BCL:HBA2	1.15	1.14
14:AJ:102:CRT:H342	9:AK:102:BCL:HBA1	1.26	1.14
5:A3:10:LYS:HB3	14:A7:102:CRT:H22A	1.29	1.14
6:B2:21:PHE:HD1	14:B2:102:CRT:C14	1.59	1.14
6:B6:32:VAL:HG21	9:B6:101:BCL:HBA2	1.15	1.14
6:AV:7:THR:HG22	14:AX:102:CRT:H1M1	1.28	1.13
5:A7:37:MET:H	14:A7:102:CRT:H2M1	1.13	1.13
14:A5:103:CRT:H293	9:A9:102:BCL:H42	1.25	1.13
5:BA:36:HIS:HB2	14:B0:101:CRT:C39	1.79	1.13
5:BU:43:ASP:HA	5:BW:47:LEU:O	1.49	1.13
6:AR:46:LEU:HB3	6:AT:42:TYR:CZ	1.84	1.12
5:AO:8:LEU:O	5:AO:11:ILE:HG13	1.48	1.12
6:AT:27:ALA:O	6:AT:31:LEU:HG	1.50	1.12
6:A0:17:PHE:HD1	6:A0:18:HIS:N	1.47	1.12
6:B2:13:GLU:O	14:B2:102:CRT:H32A	1.48	1.12
4:BH:6:THR:HB	5:BF:41:SER:HB3	1.25	1.12
5:AS:30:VAL:HG21	15:AS:101:PEF:H391	1.28	1.11
5:AW:9:TYR:HA	6:AX:18:HIS:ND1	1.63	1.11
14:A2:102:CRT:H2M1	5:A3:40:LEU:HD11	1.22	1.11
6:AG:21:PHE:HD1	6:AG:22:MET:N	1.47	1.11
5:A5:4:MET:HG2	6:A8:24:SER:HA	1.18	1.11
6:AZ:46:LEU:HB2	5:A1:52:PRO:HD3	1.22	1.11
5:BA:36:HIS:CB	14:B0:101:CRT:C39	2.27	1.11
5:AF:44:LEU:HB2	6:AG:43:ARG:HH11	1.16	1.11
5:A9:2:PHE:N	5:A9:5:ASN:HD22	1.46	1.11
14:BG:102:CRT:H342	9:BI:102:BCL:HBA1	1.29	1.11

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
14:AS:104:CRT:C9	6:AV:20:ILE:CD1	2.28	1.11
5:BF:4:MET:HG2	6:BJ:23:GLN:HG3	1.11	1.10
6:B0:17:PHE:CD1	14:B0:101:CRT:H9	1.85	1.10
4:BH:5:ILE:HG23	4:BH:6:THR:H	1.05	1.10
6:A6:44:PRO:HG2	5:A7:52:PRO:HG2	1.26	1.10
5:B3:11:ILE:HG12	14:B7:102:CRT:H81	1.26	1.10
4:AH:136:MET:HA	4:AH:139:ALA:HB3	1.31	1.10
5:AI:9:TYR:HA	6:AJ:18:HIS:CE1	1.86	1.10
5:BW:26:ALA:O	5:BW:29:ILE:HG22	1.50	1.10
6:B2:13:GLU:HB3	14:B2:102:CRT:H33	1.32	1.10
6:B2:20:ILE:HG21	14:B2:102:CRT:C7	1.80	1.10
14:B2:102:CRT:H2M1	5:B3:36:HIS:HB3	1.31	1.10
5:BO:43:ASP:CA	5:BQ:48:ASP:HB3	1.82	1.10
14:B2:102:CRT:C2M	5:B3:36:HIS:HB3	1.82	1.10
14:AB:102:CRT:H23	5:A9:13:LEU:CD1	1.82	1.09
5:AU:12:TRP:HE1	6:AV:18:HIS:HA	1.13	1.09
5:A3:13:LEU:HB2	14:A7:102:CRT:H1M1	1.31	1.09
1:AC:153:TYR:HB3	1:AC:323:MET:HE1	1.34	1.09
1:AC:165:ALA:HB1	1:AC:303:LEU:HB3	1.30	1.09
5:AO:12:TRP:NE1	6:AP:18:HIS:HA	1.67	1.09
5:AS:50:ASN:HA	5:AU:60:LYS:HA	1.29	1.09
3:BM:79:VAL:HG21	3:BM:85:GLN:HB3	1.33	1.09
2:AL:203:ILE:HG21	3:AM:266:HIS:HD1	0.96	1.09
5:AF:19:ARG:NH2	5:AI:18:ARG:HH21	1.50	1.09
5:AO:12:TRP:HE1	6:AP:18:HIS:HA	0.97	1.09
5:AS:10:LYS:O	14:AS:104:CRT:H33	1.50	1.09
6:BV:21:PHE:HD1	14:BV:102:CRT:H14	1.01	1.09
5:A5:4:MET:HE3	6:A8:24:SER:HB3	1.14	1.08
6:BZ:46:LEU:HD22	6:B2:42:TYR:CZ	1.87	1.08
14:BS:103:CRT:H342	9:BU:102:BCL:HBA1	1.15	1.08
4:BH:227:ASN:HD22	4:BH:228:PRO:HD2	1.13	1.08
5:BA:27:PHE:CZ	5:BD:29:ILE:HD11	1.86	1.08
5:B7:29:ILE:HA	9:B7:103:BCL:H11	1.33	1.08
14:B7:102:CRT:H342	9:B7:103:BCL:HBA1	1.10	1.08
6:BV:17:PHE:CD1	14:BV:102:CRT:H6	1.89	1.08
5:B5:5:ASN:HA	5:B5:8:LEU:HG	1.35	1.08
5:AF:4:MET:HB2	6:AJ:23:GLN:HG3	1.33	1.07
5:A1:5:ASN:HA	5:A1:8:LEU:HB3	1.34	1.07
5:A1:9:TYR:HA	6:A2:18:HIS:ND1	1.69	1.07
14:BW:103:CRT:H342	9:B1:102:BCL:HBA1	1.35	1.07
5:B7:43:ASP:HB2	5:B9:47:LEU:HD12	1.34	1.07

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:BF:13:LEU:HD12	14:BF:103:CRT:H1M1	1.33	1.07
5:BW:16:ASP:HB2	5:BW:19:ARG:NE	1.70	1.07
5:A3:19:ARG:O	5:A3:23:SER:HB2	1.53	1.06
4:BH:125:LEU:HA	4:BH:131:PRO:HA	1.35	1.06
6:BE:33:VAL:HG23	9:BE:101:BCL:H141	1.31	1.06
5:BU:12:TRP:HE1	6:BV:18:HIS:HA	0.93	1.06
5:BF:45:ASN:HB3	5:BF:49:ASP:HB3	1.36	1.06
6:A6:32:VAL:CG2	9:A6:101:BCL:HBA2	1.85	1.06
5:A7:29:ILE:HB	9:A7:103:BCL:H43	1.32	1.06
6:BX:32:VAL:HG11	9:BX:101:BCL:HBA2	1.35	1.06
14:AW:102:CRT:H183	9:AY:102:BCL:C9	1.84	1.06
5:BU:21:LEU:O	5:BU:25:VAL:HG23	1.53	1.06
5:B7:46:TRP:CH2	9:B7:103:BCL:HBC3	1.91	1.06
5:A3:11:ILE:HG12	14:A7:102:CRT:H81	1.34	1.06
3:BM:179:ILE:HD13	3:BM:179:ILE:H	1.14	1.06
5:AW:49:ASP:HB2	5:AY:56:GLN:HB2	1.36	1.05
6:B2:20:ILE:HG21	14:B2:102:CRT:C8	1.86	1.05
6:AR:27:ALA:O	6:AR:31:LEU:HG	1.56	1.05
6:A6:32:VAL:HG21	9:A6:101:BCL:CBA	1.84	1.05
5:B1:11:ILE:CA	14:B1:103:CRT:C8	2.34	1.05
5:AK:5:ASN:ND2	6:AN:22:MET:HE3	1.69	1.05
14:B5:103:CRT:H342	9:B9:102:BCL:HBA1	1.32	1.05
5:AW:2:PHE:HA	5:AW:5:ASN:HD22	1.13	1.05
5:BW:16:ASP:HB2	5:BW:19:ARG:HE	1.19	1.05
5:B1:11:ILE:HA	14:B1:103:CRT:H82	1.32	1.05
5:AO:8:LEU:HA	6:AR:20:ILE:HD11	1.32	1.05
5:A3:43:ASP:HB2	5:A5:47:LEU:HD13	1.06	1.05
6:B2:21:PHE:HD1	14:B2:102:CRT:H14	0.94	1.05
6:AP:46:LEU:HB2	5:AQ:51:ILE:HG21	1.34	1.04
14:BG:102:CRT:H2M2	5:BI:37:MET:HE2	1.28	1.04
1:AC:183:GLN:HE22	1:AC:230:GLU:HG2	1.20	1.04
1:BC:20:LEU:HD22	1:BC:21:LEU:H	1.21	1.04
3:BM:63:PHE:HB3	3:BM:125:SER:HB2	1.39	1.04
14:BB:102:CRT:H2M3	5:BD:36:HIS:CB	1.87	1.04
6:BG:32:VAL:HG11	9:BG:101:BCL:HBA2	1.40	1.04
6:BV:21:PHE:HE1	14:BV:102:CRT:H16	1.19	1.03
5:AA:43:ASP:HA	5:AD:48:ASP:HB3	1.38	1.03
1:AC:280:ASN:OD1	1:AC:304:ARG:HB3	1.57	1.03
3:AM:79:VAL:HG21	3:AM:85:GLN:HB3	1.35	1.03
6:A0:17:PHE:CD1	6:A0:17:PHE:C	2.35	1.03
5:BK:27:PHE:CE2	5:BO:29:ILE:HD11	1.92	1.03

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:BS:28:GLN:O	9:BS:102:BCL:H11	1.56	1.03
2:AL:203:ILE:HG21	3:AM:266:HIS:ND1	1.74	1.03
5:AW:51:ILE:HB	5:AW:52:PRO:HA	1.40	1.03
5:A1:12:TRP:HZ2	6:A2:21:PHE:CD2	1.77	1.03
6:BZ:46:LEU:HD21	6:B2:42:TYR:CE2	1.94	1.03
4:AH:31:ARG:HA	4:AH:31:ARG:HE	1.23	1.03
6:AB:23:GLN:HG3	5:A9:4:MET:HE1	1.34	1.03
1:BC:135:ARG:HH12	1:BC:332:LYS:HA	1.21	1.03
3:AM:105:ARG:HA	5:AO:42:THR:HG22	1.38	1.02
6:A0:32:VAL:HG21	9:A0:102:BCL:HBA2	1.41	1.02
2:BL:239:HIS:CD2	3:BM:223:ILE:HG13	1.94	1.02
5:BA:36:HIS:CG	14:B0:101:CRT:H392	1.93	1.02
4:AH:125:LEU:HA	4:AH:131:PRO:HA	1.37	1.02
2:BL:120:LEU:HD21	3:BM:250:LEU:HD23	1.40	1.02
14:B1:103:CRT:H342	9:B5:102:BCL:HBA1	1.39	1.02
6:AB:32:VAL:HG21	9:AB:101:BCL:HBA2	1.38	1.02
5:AF:42:THR:O	5:AI:48:ASP:HB3	1.59	1.02
5:AS:51:ILE:HB	5:AS:52:PRO:HA	1.40	1.02
5:AY:8:LEU:HB3	6:AZ:18:HIS:CE1	1.94	1.02
5:A7:36:HIS:HB3	14:A7:102:CRT:H2M3	1.39	1.02
6:AB:29:PHE:HE1	9:AB:101:BCL:H11	1.21	1.02
5:A1:7:ASN:HB3	5:A1:10:LYS:HE3	1.39	1.02
5:A1:44:LEU:HD13	6:A2:43:ARG:HD2	1.38	1.02
5:A7:37:MET:N	14:A7:102:CRT:H2M1	1.74	1.02
14:BP:102:CRT:C34	9:BQ:103:BCL:HBA1	1.88	1.02
3:BM:104:LEU:HD11	3:BM:169:GLY:HA2	1.39	1.02
5:BY:8:LEU:HA	6:B2:20:ILE:HD11	1.41	1.02
6:AG:28:TRP:NE1	6:AG:32:VAL:HG21	1.74	1.01
6:AN:30:GLY:O	6:AN:34:ILE:HG22	1.59	1.01
5:A5:4:MET:HG3	6:A8:27:ALA:CB	1.89	1.01
3:BM:187:ALA:HA	9:BM:402:BCL:HBC1	1.42	1.01
5:BU:43:ASP:HA	5:BW:47:LEU:C	1.84	1.01
14:AS:104:CRT:C18	9:AU:102:BCL:C9	2.37	1.01
6:B2:20:ILE:HG21	14:B2:102:CRT:H83	1.42	1.01
6:B2:21:PHE:HB2	14:B2:102:CRT:H11	1.42	1.01
1:AC:45:ASN:HD22	1:AC:48:GLN:HB2	1.25	1.01
4:AH:227:ASN:HD22	4:AH:228:PRO:HD2	1.20	1.01
5:AO:8:LEU:HD12	6:AP:18:HIS:HE1	1.20	1.01
6:A8:33:VAL:HG23	9:A8:101:BCL:H142	1.43	1.01
3:BM:140:LEU:HD23	3:BM:142:MET:HG3	1.42	1.01
9:BQ:103:BCL:HBC1	9:BQ:104:BCL:HBC3	1.37	1.01

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:AB:20:ILE:HD12	14:AB:102:CRT:H10	1.41	1.01
4:BH:31:ARG:HA	4:BH:31:ARG:HE	1.25	1.01
14:BA:102:CRT:H342	9:BF:102:BCL:HBA1	1.41	1.01
5:BK:27:PHE:HE2	5:BO:29:ILE:HD11	1.22	1.01
6:BZ:46:LEU:HB2	5:B1:52:PRO:HD3	1.42	1.01
9:AE:101:BCL:C1B	9:AF:102:BCL:HMB2	1.92	1.00
10:BM:403:BPH:HMA1	15:BQ:101:PEF:H411	1.41	1.00
14:BB:102:CRT:C2M	5:BD:36:HIS:HB2	1.90	1.00
5:BO:29:ILE:HA	9:BO:102:BCL:H11	1.39	1.00
6:BV:17:PHE:HD1	14:BV:102:CRT:H6	1.21	1.00
2:AL:22:LEU:HB2	5:A7:19:ARG:CB	1.90	1.00
5:BU:12:TRP:NE1	6:BV:18:HIS:HA	1.74	1.00
5:B1:11:ILE:N	14:B1:103:CRT:C8	2.25	1.00
5:BA:27:PHE:CE1	5:BD:29:ILE:HD11	1.95	1.00
5:BQ:36:HIS:CE1	9:BQ:104:BCL:HMD3	1.96	1.00
6:BT:9:LEU:HD22	6:BT:13:GLU:HG3	1.43	1.00
5:BU:14:ILE:CG1	14:BU:103:CRT:C3	2.26	1.00
5:BU:18:ARG:O	5:BU:22:VAL:HG12	1.61	1.00
2:BL:177:HIS:HB3	3:BM:183:LEU:HD22	1.43	1.00
5:B1:13:LEU:HB3	14:B1:103:CRT:H1M3	1.38	1.00
6:B2:21:PHE:CD1	14:B2:102:CRT:H16	1.95	1.00
4:AH:159:LEU:HD22	4:AH:254:ARG:NH2	1.77	1.00
5:AY:28:GLN:HB3	9:AY:102:BCL:C2	1.92	1.00
1:BC:165:ALA:HB1	1:BC:303:LEU:HB3	1.42	1.00
5:BO:4:MET:HB2	6:BR:23:GLN:HB3	1.42	1.00
5:AF:28:GLN:HB3	9:AF:102:BCL:H12	1.36	1.00
2:AL:160:LEU:HD11	3:AM:305:PRO:HD3	1.43	0.99
6:B0:21:PHE:HB2	14:B0:101:CRT:C12	1.91	0.99
1:AC:280:ASN:HB3	1:AC:304:ARG:HD2	1.40	0.99
1:BC:280:ASN:OD1	1:BC:304:ARG:HB3	1.60	0.99
4:BH:5:ILE:CG2	5:BD:42:THR:HG21	1.93	0.99
14:B2:102:CRT:H342	9:B3:102:BCL:HBA1	1.42	0.99
9:AO:102:BCL:ND	9:AP:101:BCL:HMD3	1.76	0.99
5:BQ:51:ILE:HG13	5:BQ:52:PRO:HA	1.44	0.99
5:BU:14:ILE:CD1	14:BU:103:CRT:C3	2.41	0.99
5:B7:44:LEU:CD2	5:B7:46:TRP:HE3	1.75	0.99
6:AG:46:LEU:HB3	6:AJ:42:TYR:CZ	1.97	0.99
6:AP:21:PHE:HE1	6:AP:25:MET:HB2	1.22	0.99
5:A5:4:MET:HG2	6:A8:24:SER:CA	1.92	0.99
1:BC:39:GLY:HA3	2:BL:168:ASN:ND2	1.77	0.99
5:BF:50:ASN:CG	6:BG:43:ARG:HH22	1.71	0.99

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:BO:4:MET:HB2	6:BR:23:GLN:CB	1.92	0.99
6:B2:20:ILE:CG2	14:B2:102:CRT:C7	2.39	0.99
6:B0:21:PHE:HB2	14:B0:101:CRT:C14	1.91	0.99
14:AW:102:CRT:H183	9:AY:102:BCL:H91	1.43	0.99
1:BC:267:THR:HG21	3:BM:314:VAL:HB	1.41	0.99
14:AW:102:CRT:C18	9:AY:102:BCL:C9	2.39	0.99
5:A3:43:ASP:HB2	5:A5:47:LEU:CD1	1.92	0.99
5:BA:46:TRP:CB	6:BB:43:ARG:HH22	1.75	0.99
6:B2:17:PHE:CD1	14:B2:102:CRT:H9	1.97	0.99
14:AB:102:CRT:H31A	5:A9:10:LYS:O	1.63	0.98
3:BM:250:LEU:HG	3:BM:254:TRP:HE1	1.24	0.98
14:BN:102:CRT:H342	9:BO:102:BCL:HBA1	1.43	0.98
4:BH:136:MET:HA	4:BH:139:ALA:HB3	1.42	0.98
14:A7:102:CRT:C22	14:A7:102:CRT:C20	2.42	0.98
4:BH:159:LEU:HD22	4:BH:254:ARG:NH2	1.77	0.98
14:BU:103:CRT:C2M	5:BY:37:MET:N	2.26	0.98
6:AN:41:LEU:HD23	6:AN:42:TYR:N	1.77	0.98
14:BV:102:CRT:C39	5:BW:36:HIS:CB	2.30	0.98
5:AA:18:ARG:HG3	5:A9:14:ILE:HG23	1.46	0.98
5:B3:44:LEU:HD21	9:B4:101:BCL:HBC3	1.45	0.98
1:AC:236:MET:SD	7:AC:503:HEM:NB	2.35	0.98
2:AL:175:HIS:HA	17:AL:402:HOH:O	1.64	0.98
5:A1:8:LEU:HD23	5:A1:9:TYR:N	1.77	0.98
4:AH:55:VAL:HG13	4:AH:56:VAL:H	1.26	0.98
5:AW:5:ASN:HA	5:AW:8:LEU:HD12	1.46	0.98
9:AW:101:BCL:C1	9:AW:101:BCL:CGA	2.42	0.98
5:A7:43:ASP:HB2	5:A9:47:LEU:HD12	1.44	0.98
5:AF:8:LEU:HD21	6:AJ:24:SER:OG	1.62	0.98
5:BW:9:TYR:HA	6:BX:18:HIS:CG	1.98	0.98
6:A6:45:TRP:CD1	6:A6:46:LEU:H	1.82	0.97
5:BD:46:TRP:CH2	9:BD:102:BCL:HBC3	1.99	0.97
6:AP:32:VAL:HG11	9:AP:101:BCL:CBA	1.94	0.97
6:AZ:45:TRP:CE3	9:AZ:101:BCL:HAC2	2.00	0.97
4:BH:55:VAL:HG13	4:BH:56:VAL:H	1.28	0.97
5:BW:24:ILE:HD11	9:BY:102:BCL:H18	1.45	0.97
2:AL:273:ASN:HA	2:AL:276:LEU:HD23	1.45	0.97
6:AG:45:TRP:HA	5:AI:52:PRO:HD2	1.44	0.97
5:AS:9:TYR:HB2	6:AT:15:LYS:HA	1.46	0.97
14:A5:103:CRT:C29	9:A9:102:BCL:H42	1.93	0.97
6:BV:21:PHE:CE1	14:BV:102:CRT:H16	1.98	0.97
3:BM:25:LYS:HD2	6:BP:8:GLY:HA3	1.45	0.97

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:BA:43:ASP:HA	5:BD:48:ASP:HB3	1.45	0.97
14:BB:102:CRT:H2M3	5:BD:36:HIS:HB2	0.98	0.97
14:A1:103:CRT:H342	9:A5:102:BCL:H3A	1.46	0.97
6:AR:46:LEU:HB3	6:AT:42:TYR:CE2	2.00	0.97
6:BE:45:TRP:O	6:BE:46:LEU:HG	1.65	0.97
14:AB:102:CRT:H2M3	5:AD:36:HIS:HB2	1.47	0.97
5:BF:9:TYR:CE1	5:BF:10:LYS:HD3	2.00	0.97
5:AI:42:THR:HB	5:AK:48:ASP:OD1	1.64	0.97
5:BA:36:HIS:HB2	14:B0:101:CRT:H391	0.98	0.97
14:BV:102:CRT:H393	5:BW:36:HIS:HB3	1.23	0.96
9:BW:102:BCL:ND	9:BX:101:BCL:HMD3	1.80	0.96
14:AB:102:CRT:H23	5:A9:13:LEU:HD12	0.99	0.96
5:AO:12:TRP:HE1	6:AP:18:HIS:CA	1.77	0.96
5:AS:46:TRP:CZ3	9:AS:103:BCL:HBC3	1.99	0.96
5:A7:36:HIS:HB2	14:A7:102:CRT:C2M	1.94	0.96
5:AY:11:ILE:HD13	9:A1:102:BCL:H151	1.46	0.96
1:BC:165:ALA:CB	1:BC:303:LEU:HB3	1.94	0.96
2:AL:78:PRO:HB3	2:AL:92:GLY:HA3	1.47	0.96
14:AP:102:CRT:C34	9:AQ:102:BCL:HBA1	1.95	0.96
14:AS:104:CRT:C10	6:AV:20:ILE:CD1	2.43	0.96
5:B5:10:LYS:HB3	14:B5:103:CRT:H5	1.46	0.96
4:AH:231:VAL:HG23	4:AH:235:GLU:HG3	1.47	0.96
5:AS:37:MET:HG2	15:AS:101:PEF:H452	1.47	0.96
14:AS:104:CRT:C10	6:AV:20:ILE:HD12	1.95	0.96
14:AS:104:CRT:H36	5:AW:33:LEU:HA	1.47	0.96
6:BN:10:THR:HG22	6:BN:11:ASP:H	1.27	0.96
5:BY:12:TRP:HE1	6:BZ:18:HIS:HA	1.31	0.96
2:BL:188:PHE:HB3	2:BL:249:ALA:HB2	1.45	0.96
4:BH:130:LEU:HG	4:BH:131:PRO:HD2	1.44	0.96
5:BF:11:ILE:HB	14:BF:103:CRT:H82	1.44	0.96
9:BV:101:BCL:HMA1	9:BW:102:BCL:HHB	1.48	0.96
2:AL:126:VAL:HB	2:AL:127:PRO:HD3	1.47	0.96
6:A0:20:ILE:HD12	14:A0:101:CRT:H133	1.47	0.96
4:BH:5:ILE:HG23	4:BH:6:THR:N	1.81	0.96
5:B3:28:GLN:HG3	9:B3:102:BCL:H12	1.47	0.96
1:AC:175:PRO:HD2	1:AC:179:LYS:HB2	1.46	0.96
5:BF:29:ILE:HA	9:BF:102:BCL:H11	1.46	0.96
5:BO:50:ASN:CG	6:BP:43:ARG:HH22	1.73	0.96
14:BU:103:CRT:H14	6:BX:24:SER:HB3	1.46	0.96
5:BA:46:TRP:HB2	6:BB:43:ARG:HH22	1.24	0.95
6:AB:18:HIS:HE1	6:AB:22:MET:HE2	1.30	0.95

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:AB:29:PHE:CE1	9:AB:101:BCL:H11	2.00	0.95
14:AS:104:CRT:C7	6:AV:20:ILE:HD12	1.94	0.95
6:BB:32:VAL:HG21	9:BB:101:BCL:HBA2	1.48	0.95
9:BO:102:BCL:ND	9:BP:101:BCL:HMD3	1.80	0.95
6:AG:21:PHE:CD1	6:AG:21:PHE:C	2.45	0.95
5:AI:9:TYR:HA	6:AJ:18:HIS:ND1	1.79	0.95
5:A7:37:MET:N	14:A7:102:CRT:C2M	2.28	0.95
5:B7:33:LEU:O	14:B7:102:CRT:H2M1	1.66	0.95
6:B0:32:VAL:HG21	9:B0:102:BCL:HBA2	1.47	0.95
5:AS:11:ILE:HA	14:AS:104:CRT:H82	1.47	0.95
6:B2:17:PHE:CD1	14:B2:102:CRT:H6	2.01	0.95
4:AH:195:LEU:HD12	4:AH:196:PRO:HD2	1.49	0.95
6:AB:40:TRP:HZ3	6:AB:45:TRP:H	1.05	0.95
5:BA:55:TYR:HE1	5:B9:44:LEU:HB3	1.28	0.95
4:AH:123:CYS:HA	4:AH:232:THR:HA	1.48	0.95
6:AG:46:LEU:HB3	6:AJ:42:TYR:CE2	2.02	0.95
5:AQ:51:ILE:HG13	5:AQ:52:PRO:HA	1.45	0.95
5:B3:11:ILE:N	14:B7:102:CRT:H82	1.81	0.95
5:AO:13:LEU:O	6:AP:7:THR:HA	1.66	0.95
6:AP:30:GLY:O	6:AP:34:ILE:HG22	1.67	0.95
14:A5:103:CRT:H342	9:A9:102:BCL:HBA1	1.46	0.95
1:BC:153:TYR:HB3	1:BC:323:MET:HE1	1.49	0.95
2:BL:183:MET:HA	17:BL:401:HOH:O	1.63	0.95
6:B2:17:PHE:HB2	14:B2:102:CRT:C4	1.96	0.95
9:A3:104:BCL:H61	6:A4:29:PHE:CE1	2.01	0.95
6:A0:17:PHE:HD1	6:A0:17:PHE:C	1.74	0.95
9:BW:102:BCL:C1D	9:BX:101:BCL:CMD	2.45	0.95
5:AF:27:PHE:CE1	5:AI:29:ILE:HD11	2.02	0.94
5:A5:16:ASP:HB2	5:A5:19:ARG:HG2	1.47	0.94
9:A5:102:BCL:HBC2	9:A6:101:BCL:HMD1	1.48	0.94
2:BL:230:GLY:HA2	3:BM:51:ILE:HB	1.49	0.94
9:BA:101:BCL:HBA1	14:B0:101:CRT:H342	1.45	0.94
14:BF:103:CRT:H342	9:BK:102:BCL:HBA1	1.46	0.94
6:BZ:46:LEU:CB	5:B1:52:PRO:HD3	1.96	0.94
10:AM:403:BPH:H9C3	15:AM:409:PEF:H222	1.48	0.94
5:BU:12:TRP:HE1	6:BV:18:HIS:CA	1.80	0.94
9:AL:301:BCL:HBA2	9:AL:303:BCL:HBC1	1.47	0.94
6:AV:7:THR:HG22	14:AX:102:CRT:C1M	1.97	0.94
3:BM:275:LEU:HA	3:BM:278:ILE:HD12	1.48	0.94
4:BH:227:ASN:ND2	4:BH:228:PRO:HD2	1.80	0.94
5:AF:19:ARG:HH22	5:AI:18:ARG:NH2	1.65	0.94

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:AG:21:PHE:CD1	6:AG:22:MET:N	2.35	0.94
6:B2:29:PHE:HE1	9:B2:101:BCL:H11	1.26	0.94
4:BH:176:GLU:HG3	4:BH:178:GLN:HG2	1.50	0.94
5:BF:4:MET:CG	6:BJ:23:GLN:HG3	1.96	0.94
5:A5:25:VAL:HG11	9:A5:102:BCL:H192	1.50	0.94
5:A5:43:ASP:HB2	5:A7:47:LEU:HB3	1.45	0.94
1:AC:165:ALA:CB	1:AC:303:LEU:HB3	1.98	0.94
5:AS:13:LEU:HB2	14:AS:104:CRT:C3	1.96	0.94
5:BO:43:ASP:HA	5:BQ:48:ASP:HB3	0.96	0.94
6:B4:13:GLU:HA	6:B4:16:GLU:HG2	1.48	0.94
6:B0:40:TRP:HH2	6:B0:46:LEU:HG	1.32	0.94
5:AW:10:LYS:HD2	6:AZ:20:ILE:HD12	1.48	0.94
6:AZ:10:THR:HG22	6:AZ:11:ASP:H	1.30	0.94
5:A5:51:ILE:HB	5:A5:52:PRO:HA	1.50	0.94
1:AC:291:LEU:HD23	1:AC:292:PRO:HD2	1.48	0.94
4:BH:5:ILE:CG2	4:BH:6:THR:H	1.80	0.94
6:BG:27:ALA:O	6:BG:31:LEU:HG	1.68	0.94
9:BQ:103:BCL:OBD	6:BR:32:VAL:HG13	1.65	0.94
6:B2:21:PHE:CE1	14:B2:102:CRT:C16	2.50	0.94
1:AC:39:GLY:HA3	2:AL:168:ASN:ND2	1.83	0.93
9:AL:301:BCL:H2C	9:AM:402:BCL:H2C	1.49	0.93
14:AW:102:CRT:H182	9:AY:102:BCL:H8	1.49	0.93
5:BO:4:MET:HB2	6:BR:23:GLN:CG	1.99	0.93
2:AL:196:LEU:HD23	3:AM:216:PHE:HB2	1.47	0.93
6:AR:46:LEU:HB3	6:AT:42:TYR:OH	1.69	0.93
5:A1:12:TRP:HE1	6:A2:18:HIS:HA	1.32	0.93
4:AH:55:VAL:HA	5:AA:19:ARG:HH12	1.31	0.93
14:AS:104:CRT:H183	9:AU:102:BCL:C9	1.97	0.93
6:A6:29:PHE:CE1	9:A6:101:BCL:H11	2.02	0.93
6:BP:10:THR:HG22	6:BP:11:ASP:H	1.31	0.93
4:BH:6:THR:CB	5:BF:41:SER:HB3	1.99	0.93
5:BU:12:TRP:CE2	6:BV:17:PHE:CE2	2.55	0.93
5:AU:12:TRP:NE1	6:AV:18:HIS:HA	1.82	0.93
1:BC:203:PHE:CE1	1:BC:210:ILE:HG12	2.01	0.93
5:BS:45:ASN:HB3	5:BS:49:ASP:HB3	1.48	0.93
5:B9:51:ILE:HB	5:B9:52:PRO:HA	1.47	0.93
2:AL:158:GLY:HA3	2:AL:161:SER:HB3	1.49	0.93
5:BU:16:ASP:HB3	5:BU:18:ARG:HH11	1.34	0.93
5:AS:30:VAL:CG2	15:AS:101:PEF:C39	2.46	0.93
5:A7:36:HIS:HB2	14:A7:102:CRT:H2M3	1.45	0.93
6:BX:46:LEU:HB2	5:BY:52:PRO:HD3	1.48	0.93

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:A2:17:PHE:HE1	14:A2:102:CRT:H9	1.33	0.93
5:BA:47:LEU:HD12	5:B9:43:ASP:HB2	1.49	0.93
4:AH:5:ILE:HG23	4:AH:6:THR:H	1.33	0.93
5:B7:43:ASP:CA	5:B9:48:ASP:HB3	1.99	0.93
4:AH:234:TYR:O	4:AH:238:LYS:HG2	1.69	0.92
14:AS:104:CRT:H183	9:AU:102:BCL:H92	1.51	0.92
5:AY:12:TRP:HE1	6:AZ:18:HIS:HA	1.34	0.92
3:BM:105:ARG:HA	5:BO:42:THR:HG22	1.51	0.92
6:B4:30:GLY:O	6:B4:33:VAL:HG12	1.70	0.92
6:B8:32:VAL:HG11	9:B8:101:BCL:HBA2	1.50	0.92
3:AM:200:PRO:HA	3:AM:203:MET:HG2	1.51	0.92
5:AF:9:TYR:CE1	5:AF:10:LYS:HD3	2.03	0.92
14:AS:104:CRT:H342	9:AW:101:BCL:HBA1	1.51	0.92
5:A1:15:LEU:HA	5:A3:18:ARG:HH12	1.34	0.92
9:A1:102:BCL:H92	14:A2:102:CRT:C18	1.98	0.92
5:A5:4:MET:SD	6:A8:27:ALA:HB3	2.10	0.92
5:BA:46:TRP:HB2	6:B0:46:LEU:OXT	1.69	0.92
5:B1:11:ILE:N	14:B1:103:CRT:H83	1.82	0.92
2:AL:87:ALA:H	2:AL:96:GLN:HE22	0.93	0.92
4:AH:227:ASN:ND2	4:AH:228:PRO:HD2	1.83	0.92
1:BC:97:VAL:HG21	1:BC:131:PHE:HZ	1.31	0.92
2:BL:196:LEU:HD11	3:BM:269:ALA:HB1	1.48	0.92
5:BU:13:LEU:O	6:BV:7:THR:HA	1.69	0.92
9:B1:102:BCL:CBC	9:B2:101:BCL:HBC3	2.00	0.92
5:B3:11:ILE:CA	14:B7:102:CRT:H82	1.99	0.92
5:AF:28:GLN:HA	9:AG:101:BCL:HED1	1.51	0.92
6:AV:46:LEU:HB3	6:AX:42:TYR:OH	1.68	0.92
1:BC:280:ASN:HB3	1:BC:304:ARG:HD2	1.52	0.92
3:BM:299:VAL:HB	3:BM:304:ALA:HB3	1.51	0.92
5:BU:26:ALA:O	5:BU:29:ILE:HG22	1.69	0.92
5:B3:11:ILE:HG12	14:B7:102:CRT:C8	2.00	0.92
14:A2:102:CRT:H2M2	5:A3:40:LEU:HD11	1.47	0.92
3:AM:104:LEU:HD11	3:AM:169:GLY:HA2	1.51	0.92
5:AU:26:ALA:O	5:AU:29:ILE:HG22	1.69	0.92
5:AS:11:ILE:HG12	14:AS:104:CRT:H81	1.49	0.92
6:A2:25:MET:HE2	9:A3:103:BCL:H171	1.51	0.92
5:BK:11:ILE:HG12	14:BP:102:CRT:H81	1.49	0.92
14:BV:102:CRT:H342	9:BW:102:BCL:HBA1	1.49	0.92
5:BY:10:LYS:HB2	14:B2:102:CRT:H82	1.52	0.92
6:B2:21:PHE:HE1	14:B2:102:CRT:H16	1.15	0.92
1:AC:32:GLN:HB2	2:AL:80:LEU:HD12	1.49	0.92

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:A4:30:GLY:O	6:A4:33:VAL:HG12	1.70	0.92
5:BF:44:LEU:HB2	6:BG:43:ARG:HH11	1.30	0.92
5:AD:39:VAL:O	5:AD:43:ASP:HB3	1.70	0.92
6:AP:27:ALA:O	6:AP:31:LEU:HG	1.70	0.92
6:BV:46:LEU:HB3	6:BX:42:TYR:OH	1.71	0.91
5:A1:52:PRO:HD2	5:A1:55:TYR:OH	1.70	0.91
9:A0:102:BCL:HBB2	9:A0:102:BCL:C16	2.00	0.91
6:BV:17:PHE:HD1	14:BV:102:CRT:C6	1.83	0.91
5:AF:11:ILE:N	14:AJ:102:CRT:H82	1.83	0.91
5:A7:4:MET:SD	6:A0:24:SER:OG	2.28	0.91
1:BC:175:PRO:HD2	1:BC:179:LYS:HB2	1.52	0.91
2:BL:89:LEU:HA	2:BL:94:LEU:H	1.33	0.91
3:BM:175:VAL:HA	3:BM:185:TRP:CD1	2.05	0.91
14:AB:102:CRT:C2	5:A9:13:LEU:CD1	2.45	0.91
1:BC:39:GLY:HA3	2:BL:168:ASN:HD22	1.33	0.91
2:BL:196:LEU:HD13	3:BM:216:PHE:HB2	1.51	0.91
4:BH:5:ILE:HG22	5:BD:42:THR:HG21	1.50	0.91
5:BF:27:PHE:CE1	5:BI:29:ILE:HD11	2.06	0.91
5:B7:43:ASP:HA	5:B9:48:ASP:HB3	1.49	0.91
5:AS:8:LEU:HD22	5:AS:11:ILE:HD11	1.53	0.91
1:BC:135:ARG:NH1	1:BC:332:LYS:HA	1.84	0.91
6:BV:20:ILE:HG21	14:BV:102:CRT:C7	2.01	0.91
9:B8:101:BCL:HMC2	9:B9:102:BCL:HBB1	1.52	0.91
3:AM:136:ARG:NH1	3:AM:136:ARG:HA	1.86	0.91
5:A7:29:ILE:HA	9:A7:103:BCL:H11	1.51	0.91
5:BD:15:LEU:HB3	5:BD:20:VAL:HG21	1.50	0.91
5:B1:11:ILE:HA	14:B1:103:CRT:H81	1.49	0.91
9:AK:102:BCL:HMD2	9:AN:101:BCL:CHD	2.01	0.91
6:BN:29:PHE:O	6:BN:33:VAL:HB	1.71	0.91
3:AM:235:ILE:H	3:AM:235:ILE:HD12	1.35	0.91
6:AN:41:LEU:HD23	6:AN:41:LEU:C	1.95	0.91
5:AO:8:LEU:HD12	6:AP:18:HIS:CE1	2.05	0.91
1:BC:81:VAL:HG11	1:BC:131:PHE:HB3	1.52	0.91
5:BK:46:TRP:HA	5:BK:49:ASP:OD1	1.69	0.91
5:BO:13:LEU:O	6:BP:7:THR:HA	1.70	0.91
5:BS:36:HIS:CE1	9:BT:101:BCL:HMD3	2.06	0.91
9:AK:102:BCL:C1D	9:AN:101:BCL:HMD1	2.00	0.91
9:AO:102:BCL:ND	9:AP:101:BCL:CMD	2.32	0.91
6:BZ:46:LEU:CD2	6:B2:42:TYR:CE2	2.54	0.91
14:AB:102:CRT:H21A	5:A9:13:LEU:HD12	1.51	0.90
1:AC:39:GLY:HA3	2:AL:168:ASN:HD22	1.33	0.90

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:A8:101:BCL:H203	9:A8:101:BCL:H152	1.49	0.90
6:A0:16:GLU:OE2	14:A0:101:CRT:H21A	1.70	0.90
5:BQ:35:ILE:HA	5:BQ:38:ILE:HG22	1.52	0.90
5:B3:5:ASN:HA	5:B3:8:LEU:HD12	1.53	0.90
2:AL:59:THR:HB	2:AL:63:SER:HB3	1.53	0.90
3:AM:63:PHE:HZ	5:AQ:33:LEU:HD23	1.34	0.90
3:AM:279:THR:HA	3:AM:282:ILE:HD12	1.53	0.90
6:AP:32:VAL:CG1	9:AP:101:BCL:HBA2	2.00	0.90
5:BA:36:HIS:CG	14:B0:101:CRT:C39	2.55	0.90
5:AF:4:MET:CB	6:AJ:23:GLN:HG3	2.00	0.90
2:BL:4:LEU:HD12	3:BM:250:LEU:HD12	1.54	0.90
5:BO:4:MET:HE2	6:BR:23:GLN:HB2	1.52	0.90
6:BV:21:PHE:HD1	14:BV:102:CRT:C14	1.84	0.90
1:AC:166:TRP:HE1	1:AC:305:VAL:C	1.79	0.90
6:A8:33:VAL:HG23	9:A8:101:BCL:C14	2.01	0.90
9:BW:102:BCL:HBC1	9:BX:101:BCL:HBC3	1.51	0.90
3:AM:278:ILE:O	3:AM:282:ILE:HG13	1.71	0.90
5:AF:19:ARG:HH12	5:AI:18:ARG:NH2	1.70	0.90
5:AO:43:ASP:HA	5:AQ:48:ASP:HB3	1.53	0.90
6:BT:10:THR:HG22	6:BT:11:ASP:H	1.36	0.90
1:AC:267:THR:HG21	3:AM:314:VAL:HB	1.53	0.90
5:BI:26:ALA:O	5:BI:29:ILE:HG22	1.70	0.90
14:BO:103:CRT:H342	9:BS:102:BCL:HBA1	1.53	0.90
9:BQ:104:BCL:HBA2	6:BR:32:VAL:HG11	1.54	0.90
2:AL:276:LEU:H	2:AL:276:LEU:HD22	1.37	0.90
6:BP:13:GLU:HA	6:BP:16:GLU:CD	1.97	0.90
5:A5:4:MET:CG	6:A8:27:ALA:HB3	2.02	0.90
4:BH:45:ARG:HD3	4:BH:97:GLY:H	1.36	0.90
6:B2:17:PHE:CB	14:B2:102:CRT:H41	2.01	0.90
5:B1:32:GLY:HA3	9:B1:102:BCL:O1A	1.72	0.90
6:B2:32:VAL:HG11	9:B2:101:BCL:HBA2	1.53	0.90
6:AE:29:PHE:CE1	9:AE:101:BCL:H11	2.07	0.89
6:AP:21:PHE:CE1	6:AP:25:MET:HB2	2.07	0.89
2:BL:197:SER:HB3	3:BM:273:ALA:HB1	1.54	0.89
6:A6:32:VAL:HG21	9:A6:101:BCL:HBA2	0.92	0.89
5:BD:26:ALA:O	5:BD:29:ILE:HG22	1.72	0.89
6:B0:40:TRP:CH2	6:B0:46:LEU:HG	2.07	0.89
1:AC:183:GLN:NE2	1:AC:230:GLU:HG2	1.86	0.89
6:AP:46:LEU:HB2	5:AQ:51:ILE:CG2	2.00	0.89
6:BV:29:PHE:CE1	9:BV:101:BCL:H11	2.07	0.89
2:AL:89:LEU:HA	2:AL:94:LEU:N	1.87	0.89

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:AH:35:LYS:HZ3	4:AH:57:GLY:HA3	1.34	0.89
14:AA:102:CRT:H342	9:AF:102:BCL:HBA1	1.52	0.89
6:AT:10:THR:HG22	6:AT:11:ASP:H	1.38	0.89
5:AY:13:LEU:HD21	6:AZ:14:ALA:CB	2.01	0.89
1:BC:17:SER:HB3	3:BM:91:PHE:HZ	1.36	0.89
14:BU:103:CRT:C34	9:BY:102:BCL:HBA1	2.01	0.89
6:B2:17:PHE:CE1	14:B2:102:CRT:H9	2.08	0.89
6:B6:29:PHE:CE1	9:B6:101:BCL:H11	2.07	0.89
1:AC:285:TRP:CZ3	1:AC:302:PRO:HD3	2.06	0.89
6:AZ:46:LEU:HB2	5:A1:52:PRO:CD	2.02	0.89
1:BC:22:GLY:HA3	2:BL:263:PHE:HB3	1.55	0.89
6:B6:32:VAL:HG21	9:B6:101:BCL:CBA	2.01	0.89
6:AZ:45:TRP:CE3	9:AZ:101:BCL:H2C	2.07	0.89
14:A2:102:CRT:H2M1	5:A3:40:LEU:CD1	2.03	0.89
5:A7:2:PHE:N	5:A7:5:ASN:HB3	1.87	0.89
5:A9:31:LEU:HD21	9:A0:102:BCL:HMA1	1.53	0.89
2:BL:217:THR:H	2:BL:220:HIS:CE1	1.90	0.89
9:BM:402:BCL:H143	15:BQ:101:PEF:H442	1.54	0.89
6:BG:31:LEU:O	6:BG:34:ILE:HG13	1.73	0.89
5:BQ:50:ASN:HD22	5:BS:56:GLN:HA	1.37	0.89
1:AC:280:ASN:O	1:AC:285:TRP:HB2	1.73	0.89
9:BL:301:BCL:HBA2	9:BL:303:BCL:HBC1	1.54	0.89
6:B0:17:PHE:CE1	14:B0:101:CRT:H11	2.07	0.89
5:AF:19:ARG:HH22	5:AI:18:ARG:HH21	0.94	0.89
1:AC:148:THR:HG23	1:AC:322:GLN:HA	1.55	0.89
2:AL:89:LEU:HG	2:AL:97:ILE:HD11	1.53	0.89
6:AT:17:PHE:HD1	14:AT:102:CRT:H6	1.36	0.89
6:BB:42:TYR:OH	6:B0:46:LEU:HB3	1.72	0.89
2:AL:87:ALA:H	2:AL:96:GLN:NE2	1.71	0.89
5:AS:30:VAL:HG21	15:AS:101:PEF:H392	1.52	0.89
6:A2:17:PHE:CE1	14:A2:102:CRT:H6	2.08	0.89
2:BL:97:ILE:HA	2:BL:100:ILE:HD12	1.54	0.89
4:BH:234:TYR:O	4:BH:238:LYS:HG2	1.72	0.89
2:AL:22:LEU:HB2	5:A7:19:ARG:HB3	1.53	0.88
3:AM:109:LEU:HB2	5:AQ:42:THR:HB	1.54	0.88
5:AA:47:LEU:HD12	5:A9:43:ASP:HB2	1.55	0.88
6:AJ:32:VAL:HG11	9:AJ:101:BCL:HBA2	1.54	0.88
5:A5:25:VAL:CG1	9:A5:102:BCL:H191	2.03	0.88
5:B1:11:ILE:CA	14:B1:103:CRT:H82	2.00	0.88
2:AL:52:TRP:NE1	5:A9:38:ILE:HA	1.87	0.88
5:AF:28:GLN:HB3	9:AF:102:BCL:H11	1.52	0.88

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:AF:102:BCL:HBA2	9:AG:101:BCL:OBD	1.73	0.88
3:AM:175:VAL:HA	3:AM:185:TRP:CD1	2.08	0.88
5:AS:46:TRP:HZ3	9:AS:103:BCL:HBC3	1.37	0.88
3:BM:156:PHE:HZ	9:BM:402:BCL:HBD	1.37	0.88
6:BV:43:ARG:NH1	5:BW:55:TYR:HB3	1.87	0.88
9:AQ:102:BCL:OBD	6:AR:32:VAL:HG13	1.73	0.88
9:AU:102:BCL:H2A	9:AU:102:BCL:O1D	1.73	0.88
6:BJ:10:THR:HG22	6:BJ:11:ASP:H	1.38	0.88
6:BZ:10:THR:HG22	6:BZ:11:ASP:H	1.38	0.88
5:B1:44:LEU:HD13	6:B2:43:ARG:HD2	1.54	0.88
14:AA:102:CRT:H83	6:AE:20:ILE:HD13	1.54	0.88
4:AH:176:GLU:HG3	4:AH:178:GLN:HG2	1.55	0.88
5:AO:13:LEU:HD23	5:AO:14:ILE:H	1.36	0.88
5:AS:10:LYS:HB3	14:AS:104:CRT:C1M	2.03	0.88
5:AW:21:LEU:O	5:AW:25:VAL:HG23	1.74	0.88
6:A8:44:PRO:O	5:A9:52:PRO:HD2	1.72	0.88
5:B5:36:HIS:CE1	9:B6:101:BCL:HMD3	2.07	0.88
5:AD:40:LEU:HD11	5:AD:47:LEU:HD23	1.53	0.88
6:A0:40:TRP:HH2	6:A0:46:LEU:HG	1.37	0.88
14:AG:102:CRT:H342	9:AI:102:BCL:HBA1	1.54	0.88
9:AZ:101:BCL:H203	6:A2:38:LEU:HD21	1.54	0.88
5:A1:11:ILE:HG22	14:A1:103:CRT:H10	1.54	0.88
5:BU:12:TRP:CD1	6:BV:17:PHE:HD2	1.92	0.88
6:B4:13:GLU:HA	6:B4:16:GLU:CG	2.03	0.88
2:AL:87:ALA:N	2:AL:96:GLN:HE22	1.72	0.88
2:AL:150:ALA:HB3	2:AL:153:HIS:HB2	1.53	0.88
9:AF:102:BCL:CED	6:AG:31:LEU:HD22	2.04	0.88
5:AS:34:LEU:HA	15:AS:101:PEF:C45	2.03	0.88
5:A1:8:LEU:HG	6:A2:18:HIS:NE2	1.89	0.88
5:B1:13:LEU:HD12	14:B1:103:CRT:H23	0.88	0.88
6:AT:29:PHE:CD1	9:AT:101:BCL:H11	2.08	0.87
14:AX:102:CRT:H342	9:AY:102:BCL:H3A	1.55	0.87
5:AY:51:ILE:HB	5:AY:52:PRO:HA	1.56	0.87
6:AZ:46:LEU:HD22	6:A2:42:TYR:OH	1.73	0.87
5:BF:8:LEU:HD21	6:BJ:24:SER:OG	1.74	0.87
5:BO:50:ASN:ND2	6:BP:43:ARG:HH22	1.71	0.87
14:BU:103:CRT:H343	9:BY:102:BCL:HBA1	1.54	0.87
1:AC:153:TYR:CD1	1:AC:157:ARG:HA	2.10	0.87
5:A1:8:LEU:HG	6:A2:18:HIS:CE1	2.08	0.87
5:BS:4:MET:SD	6:BV:24:SER:HB3	2.14	0.87
2:AL:68:TYR:HA	2:AL:73:ILE:HD11	1.56	0.87

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:AT:18:HIS:O	6:AT:22:MET:HG2	1.73	0.87
5:A3:50:ASN:HB2	5:A5:59:GLY:O	1.73	0.87
5:A7:2:PHE:N	5:A7:5:ASN:CB	2.37	0.87
4:BH:35:LYS:NZ	4:BH:57:GLY:HA3	1.88	0.87
2:AL:182:HIS:N	2:AL:256:CYS:SG	2.47	0.87
3:BM:59:LEU:HD11	5:BQ:29:ILE:HG21	1.57	0.87
14:BU:103:CRT:C14	6:BX:24:SER:HB3	2.04	0.87
5:BY:55:TYR:HD1	5:BY:56:GLN:H	1.20	0.87
2:AL:177:HIS:HB3	3:AM:183:LEU:HD22	1.56	0.87
5:BQ:43:ASP:OD1	5:BQ:44:LEU:HG	1.73	0.87
4:AH:35:LYS:NZ	4:AH:57:GLY:HA3	1.89	0.87
14:AN:102:CRT:H342	9:AO:102:BCL:HBA1	1.57	0.87
6:BZ:46:LEU:CD2	6:B2:42:TYR:CZ	2.57	0.87
2:AL:120:LEU:HD21	3:AM:250:LEU:HD23	1.54	0.87
6:AB:24:SER:OG	5:A9:4:MET:HE2	1.74	0.87
5:AS:50:ASN:CA	5:AU:60:LYS:HA	2.05	0.87
6:AT:17:PHE:CD1	14:AT:102:CRT:H6	2.10	0.87
5:A1:50:ASN:CG	5:A1:51:ILE:H	1.81	0.87
1:BC:291:LEU:O	1:BC:296:LYS:HE3	1.74	0.87
5:AO:13:LEU:O	6:AP:7:THR:CA	2.22	0.87
5:AW:12:TRP:HZ2	6:AX:21:PHE:CG	1.93	0.87
5:A5:21:LEU:HD11	9:A5:102:BCL:C14	2.05	0.87
9:A6:101:BCL:C1	9:A6:101:BCL:CGA	2.53	0.87
9:A9:102:BCL:HBC2	9:A0:102:BCL:HMD1	1.56	0.87
4:BH:94:PRO:HG2	6:B0:8:GLY:HA3	1.54	0.87
5:B1:13:LEU:CB	14:B1:103:CRT:H1M3	2.04	0.87
14:AR:102:CRT:H342	9:AS:103:BCL:HBA1	1.57	0.87
5:A5:4:MET:CG	6:A8:24:SER:HA	2.04	0.87
5:A7:36:HIS:CB	14:A7:102:CRT:C2M	2.53	0.87
4:BH:113:PRO:HG2	4:BH:248:LEU:HD22	1.57	0.87
5:BK:36:HIS:CE1	9:BN:101:BCL:HMD3	2.09	0.87
5:BI:35:ILE:HA	5:BI:38:ILE:HG22	1.57	0.86
6:BP:32:VAL:HG11	9:BP:101:BCL:HBA2	1.57	0.86
5:BW:49:ASP:CG	5:BW:50:ASN:H	1.83	0.86
5:B7:12:TRP:HZ3	5:B7:17:PRO:HB3	1.40	0.86
5:A7:44:LEU:CD2	5:A7:46:TRP:HE3	1.87	0.86
5:A7:46:TRP:CH2	9:A7:103:BCL:HBC3	2.10	0.86
2:BL:202:LEU:HD21	2:BL:221:GLU:HB3	1.54	0.86
3:BM:41:GLY:HA3	3:BM:46:ALA:HB2	1.57	0.86
5:BF:38:ILE:HD11	5:BI:37:MET:HE1	1.55	0.86
6:BT:18:HIS:O	6:BT:22:MET:HG2	1.74	0.86

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:BZ:46:LEU:CD2	6:B2:42:TYR:OH	2.20	0.86
5:A1:4:MET:HE2	6:A4:27:ALA:HB3	1.55	0.86
6:A0:17:PHE:CD1	6:A0:18:HIS:N	2.37	0.86
4:BH:54:LYS:HA	4:BH:54:LYS:HE3	1.56	0.86
3:AM:70:ILE:HG21	3:AM:118:ALA:HB2	1.58	0.86
4:AH:124:ASP:HB2	4:AH:233:LEU:HD21	1.55	0.86
6:AZ:45:TRP:CZ3	9:AZ:101:BCL:HAC2	2.10	0.86
1:BC:167:VAL:HG21	1:BC:298:PRO:HD2	1.57	0.86
6:B2:20:ILE:HG12	14:B2:102:CRT:H81	1.56	0.86
5:AA:18:ARG:HD2	5:AA:18:ARG:H	1.41	0.86
6:AB:20:ILE:HD12	14:AB:102:CRT:C10	2.05	0.86
6:A0:45:TRP:NE1	9:A0:102:BCL:C19	2.37	0.86
3:AM:63:PHE:CE2	3:AM:124:LEU:HD12	2.11	0.86
5:AW:9:TYR:HA	6:AX:18:HIS:CE1	2.11	0.86
5:AW:26:ALA:O	5:AW:29:ILE:HG22	1.74	0.86
1:BC:275:HIS:O	1:BC:279:ILE:HG13	1.75	0.86
6:BE:33:VAL:HG23	9:BE:101:BCL:C14	2.05	0.86
14:BS:103:CRT:C34	9:BU:102:BCL:HBA1	2.05	0.86
5:A1:12:TRP:CZ2	6:A2:21:PHE:CD2	2.64	0.86
6:A0:45:TRP:HE1	9:A0:102:BCL:H191	1.38	0.86
4:BH:35:LYS:HZ3	4:BH:57:GLY:HA3	1.39	0.86
4:BH:123:CYS:HA	4:BH:232:THR:HA	1.55	0.86
1:AC:135:ARG:HH12	1:AC:332:LYS:HA	1.39	0.86
5:AF:44:LEU:HB2	6:AG:43:ARG:NH1	1.89	0.86
5:AF:44:LEU:HD22	6:AG:43:ARG:HD2	1.57	0.86
5:AI:7:ASN:HD22	6:AN:20:ILE:HG13	1.40	0.86
5:A5:43:ASP:CB	5:A7:47:LEU:HB3	2.04	0.86
2:BL:150:ALA:HB3	2:BL:153:HIS:HB2	1.58	0.86
3:BM:235:ILE:H	3:BM:235:ILE:HD12	1.37	0.86
5:BA:15:LEU:HD21	5:BD:21:LEU:HD23	1.56	0.86
5:BU:12:TRP:CD2	6:BV:17:PHE:CE2	2.64	0.86
2:AL:97:ILE:HA	2:AL:100:ILE:HD12	1.58	0.85
5:AO:50:ASN:CG	5:AO:51:ILE:H	1.82	0.85
5:AQ:35:ILE:HA	5:AQ:38:ILE:HG22	1.56	0.85
5:AU:18:ARG:O	5:AU:22:VAL:HG12	1.75	0.85
6:A4:13:GLU:HA	6:A4:16:GLU:HG2	1.56	0.85
6:A0:10:THR:HG22	6:A0:11:ASP:H	1.40	0.85
5:BY:43:ASP:HA	5:B1:48:ASP:HB3	1.58	0.85
6:BZ:46:LEU:HD22	6:B2:42:TYR:HH	1.39	0.85
9:BZ:101:BCL:HBB1	9:B1:102:BCL:HMC1	1.58	0.85
2:BL:10:TYR:HA	4:BH:112:GLY:HA2	1.56	0.85

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:BH:151:PRO:HA	4:BH:154:MET:SD	2.16	0.85
5:B1:50:ASN:HB3	5:B3:60:LYS:HA	1.57	0.85
5:B7:36:HIS:CB	14:B7:102:CRT:H2M3	2.06	0.85
1:AC:167:VAL:HG23	1:AC:301:ASP:OD2	1.77	0.85
5:AI:27:PHE:HE2	5:AK:29:ILE:HD11	1.40	0.85
1:BC:97:VAL:HG21	1:BC:131:PHE:CZ	2.11	0.85
1:BC:285:TRP:CZ3	1:BC:302:PRO:HD3	2.11	0.85
3:BM:60:SER:HA	3:BM:128:LEU:HD23	1.56	0.85
5:B5:30:VAL:HG13	5:B5:31:LEU:H	1.40	0.85
5:AO:50:ASN:CG	6:AP:43:ARG:HH22	1.84	0.85
6:AP:10:THR:HG22	6:AP:11:ASP:H	1.41	0.85
5:AS:34:LEU:HA	15:AS:101:PEF:H453	1.56	0.85
5:AW:7:ASN:H	5:AW:7:ASN:HD22	1.25	0.85
1:BC:167:VAL:HG23	1:BC:301:ASP:OD2	1.76	0.85
1:AC:311:HIS:HA	1:AC:317:PRO:HG3	1.59	0.85
3:AM:241:ARG:HG2	3:AM:242:GLY:H	1.42	0.85
14:A2:102:CRT:C2M	5:A3:40:LEU:CD1	2.52	0.85
6:A4:31:LEU:O	6:A4:34:ILE:HG22	1.77	0.85
6:B0:10:THR:HG22	6:B0:11:ASP:H	1.40	0.85
4:AH:114:ALA:HB2	4:AH:245:GLY:HA3	1.58	0.85
5:AF:50:ASN:ND2	5:AF:51:ILE:HG12	1.90	0.85
5:A5:16:ASP:HB2	5:A5:19:ARG:CG	2.07	0.85
2:AL:29:PRO:O	3:AM:254:TRP:HA	1.76	0.85
2:AL:239:HIS:CD2	3:AM:223:ILE:HG13	2.12	0.85
4:AH:151:PRO:HA	4:AH:154:MET:SD	2.17	0.85
2:BL:52:TRP:NE1	5:B9:38:ILE:HA	1.91	0.85
6:B2:13:GLU:HB3	14:B2:102:CRT:C3	2.07	0.85
2:AL:253:SER:HB2	9:AL:301:BCL:H2A	1.59	0.85
5:AK:5:ASN:HD21	6:AN:22:MET:HE3	1.37	0.85
5:AO:29:ILE:HA	9:AO:102:BCL:H12	1.58	0.85
14:AT:102:CRT:H342	9:AU:102:BCL:CBA	2.07	0.85
2:BL:59:THR:HB	2:BL:63:SER:HB3	1.57	0.85
4:BH:168:SER:HB3	4:BH:183:GLU:HB3	1.59	0.85
5:A3:53:VAL:HA	5:A3:55:TYR:CE2	2.12	0.85
1:BC:153:TYR:CD1	1:BC:157:ARG:HA	2.11	0.85
4:BH:171:TRP:HB2	4:BH:181:TYR:HB2	1.59	0.85
1:AC:53:ILE:HG12	1:AC:319:TYR:CE1	2.12	0.84
4:AH:77:VAL:O	4:AH:80:ARG:HD3	1.76	0.84
6:AB:17:PHE:HE1	14:AB:102:CRT:H9	1.42	0.84
6:AP:38:LEU:O	6:AP:41:LEU:CD2	2.22	0.84
5:AS:10:LYS:HB3	14:AS:104:CRT:O1	1.77	0.84

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:A1:50:ASN:ND2	5:A1:51:ILE:HG23	1.92	0.84
1:BC:280:ASN:O	1:BC:285:TRP:HB2	1.76	0.84
3:BM:253:ARG:HA	3:BM:257:GLY:O	1.77	0.84
5:B7:27:PHE:CE1	5:B9:29:ILE:HD11	2.12	0.84
2:AL:140:LEU:HD12	2:AL:257:ILE:HG21	1.59	0.84
2:AL:206:VAL:O	2:AL:209:PRO:HD3	1.77	0.84
9:A5:102:BCL:H143	14:A7:102:CRT:H132	1.58	0.84
5:BU:51:ILE:HB	5:BU:52:PRO:HA	1.59	0.84
1:AC:164:TYR:HB3	1:AC:309:THR:HA	1.57	0.84
6:A2:17:PHE:CD1	14:A2:102:CRT:C6	2.60	0.84
6:A2:41:LEU:HD23	6:A2:42:TYR:N	1.91	0.84
6:BB:40:TRP:HZ3	6:BB:45:TRP:H	1.24	0.84
6:B0:33:VAL:HG12	6:B0:37:LEU:CD1	2.07	0.84
5:AY:18:ARG:O	5:AY:22:VAL:HG12	1.76	0.84
3:BM:136:ARG:NH1	3:BM:136:ARG:HA	1.92	0.84
5:BO:38:ILE:HG13	5:BO:39:VAL:N	1.93	0.84
6:BZ:44:PRO:HG2	5:B1:55:TYR:OH	1.78	0.84
5:B1:13:LEU:HB3	14:B1:103:CRT:C1M	2.07	0.84
9:AK:102:BCL:C4D	9:AN:101:BCL:HMD3	2.08	0.84
5:A1:12:TRP:NE1	6:A2:18:HIS:HA	1.93	0.84
1:BC:263:THR:HG22	3:BM:311:VAL:HB	1.60	0.84
5:B7:36:HIS:HB2	14:B7:102:CRT:H2M3	1.59	0.84
2:AL:29:PRO:HB2	3:AM:253:ARG:HD2	1.60	0.84
5:A7:12:TRP:CZ3	5:A7:17:PRO:HB3	2.11	0.84
9:BN:101:BCL:CHC	9:BO:102:BCL:HBB3	2.08	0.84
9:BO:102:BCL:HBC2	9:BP:101:BCL:HHD	1.59	0.84
14:BP:102:CRT:H2M3	5:BQ:36:HIS:HB2	1.58	0.84
5:B3:43:ASP:CB	5:B5:47:LEU:HD13	2.06	0.84
1:AC:20:LEU:HD22	1:AC:21:LEU:H	1.42	0.84
5:AD:15:LEU:HB3	5:AD:20:VAL:HG21	1.58	0.84
1:BC:20:LEU:HD22	1:BC:21:LEU:N	1.92	0.84
5:BO:43:ASP:HB2	5:BQ:47:LEU:HB3	1.60	0.84
5:B1:40:LEU:HD12	5:B1:45:ASN:HA	1.57	0.84
1:AC:73:SER:HB3	1:AC:83:LYS:HB2	1.60	0.84
2:AL:148:MET:HB3	2:AL:153:HIS:ND1	1.91	0.84
3:AM:215:LEU:HD21	13:AM:405:MQ8:H193	1.59	0.84
4:AH:31:ARG:HB3	4:AH:59:PRO:HG3	1.60	0.84
5:AF:52:PRO:HB2	5:AF:55:TYR:HE1	1.41	0.84
5:A1:11:ILE:HG22	14:A1:103:CRT:H81	1.60	0.84
2:BL:22:LEU:HB2	5:B7:19:ARG:CB	2.08	0.84
14:BF:103:CRT:H132	9:BI:102:BCL:H143	1.59	0.84

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:BJ:32:VAL:HG11	9:BJ:101:BCL:HBA2	1.59	0.84
6:B4:31:LEU:O	6:B4:34:ILE:HG22	1.76	0.84
5:B7:44:LEU:HD22	5:B7:46:TRP:HE3	1.43	0.84
4:AH:171:TRP:HB2	4:AH:181:TYR:HB2	1.60	0.84
2:BL:279:PRO:HG3	5:BY:41:SER:HB2	1.59	0.84
5:BD:46:TRP:CZ3	9:BD:102:BCL:HBC3	2.13	0.84
5:BS:40:LEU:HD11	5:BS:47:LEU:HD23	1.60	0.84
5:B7:36:HIS:CE1	9:B8:101:BCL:HMD3	2.12	0.84
6:AJ:10:THR:HB	6:AJ:13:GLU:OE2	1.76	0.84
5:AY:13:LEU:N	5:AY:13:LEU:HD23	1.93	0.84
5:BO:7:ASN:HB3	6:BR:20:ILE:HD12	1.57	0.84
3:AM:63:PHE:HE2	3:AM:124:LEU:HD12	1.43	0.83
14:AW:102:CRT:C18	9:AY:102:BCL:H8	2.08	0.83
5:A5:14:ILE:HG21	5:A7:18:ARG:HG2	1.61	0.83
2:BL:177:HIS:HB3	3:BM:183:LEU:CD2	2.08	0.83
2:BL:178:TYR:HD2	2:BL:269:PRO:HG3	1.43	0.83
6:BB:22:MET:O	6:BB:26:TYR:CE1	2.31	0.83
5:BY:20:VAL:HB	9:B1:102:BCL:H201	1.60	0.83
2:AL:230:GLY:HA2	3:AM:51:ILE:HB	1.60	0.83
3:AM:63:PHE:CZ	5:AQ:33:LEU:HD23	2.13	0.83
3:AM:202:HIS:O	3:AM:206:ILE:HG13	1.78	0.83
9:BD:102:BCL:C1D	9:BE:101:BCL:CMD	2.56	0.83
5:B1:52:PRO:HD2	5:B1:55:TYR:HE2	1.43	0.83
5:B3:19:ARG:O	5:B3:23:SER:HB2	1.76	0.83
6:B0:17:PHE:HB2	14:B0:101:CRT:C6	2.08	0.83
1:AC:142:LYS:HA	1:AC:145:VAL:HG23	1.58	0.83
5:AY:28:GLN:HB3	9:AY:102:BCL:H2	1.60	0.83
5:A7:43:ASP:HA	5:A9:48:ASP:HB3	1.58	0.83
3:BM:293:ASN:CG	3:BM:296:LEU:HG	2.03	0.83
6:BX:45:TRP:O	6:BX:46:LEU:HG	1.77	0.83
3:AM:159:VAL:HA	3:AM:163:ILE:HG22	1.61	0.83
5:A1:10:LYS:HB2	14:A1:103:CRT:H83	1.58	0.83
14:BG:102:CRT:H2M3	5:BI:36:HIS:HB3	1.60	0.83
5:A3:13:LEU:HD12	14:A7:102:CRT:H1M3	1.58	0.83
3:BM:187:ALA:HA	9:BM:402:BCL:CBC	2.08	0.83
6:BP:21:PHE:CE1	14:BP:102:CRT:H16	2.13	0.83
5:BQ:42:THR:HG23	5:BQ:43:ASP:H	1.40	0.83
3:AM:202:HIS:CE1	3:AM:206:ILE:HD11	2.14	0.83
3:AM:265:ILE:HG22	3:AM:266:HIS:H	1.43	0.83
5:AS:34:LEU:CB	15:AS:101:PEF:H442	2.08	0.83
5:A3:11:ILE:HG12	14:A7:102:CRT:C8	2.07	0.83

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:BM:204:LEU:HD12	3:BM:279:THR:HG21	1.59	0.83
4:BH:204:LYS:H	4:BH:204:LYS:HD2	1.40	0.83
9:BQ:103:BCL:C1D	9:BQ:104:BCL:CMD	2.56	0.83
9:A3:103:BCL:O1D	9:A3:103:BCL:H2A	1.78	0.83
3:BM:202:HIS:O	3:BM:206:ILE:HG13	1.78	0.83
6:BN:20:ILE:H	6:BN:20:ILE:HD12	1.44	0.83
9:AQ:102:BCL:HBC2	9:AR:101:BCL:HMD1	1.59	0.83
5:AW:46:TRP:CH2	9:AW:101:BCL:H2C	2.13	0.83
9:A8:101:BCL:C1C	9:A9:102:BCL:HBB3	2.09	0.83
6:B4:13:GLU:CA	6:B4:16:GLU:HG2	2.09	0.83
3:AM:66:VAL:HG11	3:AM:121:PHE:HD2	1.42	0.83
5:AF:8:LEU:HA	6:AJ:20:ILE:HD11	1.61	0.83
5:AS:10:LYS:HD3	14:AS:104:CRT:H1M1	1.59	0.83
14:A2:102:CRT:H31	9:A3:103:BCL:HBA1	1.60	0.83
3:BM:160:LEU:HD23	3:BM:284:ILE:HG21	1.61	0.83
5:BO:29:ILE:HB	9:BO:102:BCL:H43	1.59	0.83
6:B2:10:THR:HG22	6:B2:13:GLU:OE1	1.79	0.83
2:AL:188:PHE:CE2	2:AL:248:SER:HB3	2.14	0.83
2:AL:275:TRP:O	2:AL:278:LEU:HG	1.78	0.83
3:AM:179:ILE:HG12	3:AM:180:PHE:H	1.41	0.83
6:AX:32:VAL:HG11	9:AX:101:BCL:HBA2	1.58	0.83
5:A1:15:LEU:HD23	5:A3:18:ARG:NH1	1.92	0.83
9:A1:102:BCL:C9	14:A2:102:CRT:H183	2.08	0.83
6:B0:33:VAL:HG22	9:B0:102:BCL:H143	1.59	0.83
5:AI:44:LEU:HD13	6:AJ:43:ARG:HD3	1.61	0.82
9:A3:103:BCL:HBC2	9:A3:104:BCL:HMD1	1.60	0.82
5:A5:4:MET:CE	6:A8:24:SER:HB3	2.04	0.82
9:A7:103:BCL:C1	9:A7:103:BCL:CGA	2.57	0.82
6:BB:20:ILE:HD11	5:B9:8:LEU:HD23	1.61	0.82
6:BJ:37:LEU:HD21	9:BJ:101:BCL:H151	1.61	0.82
5:BY:50:ASN:HD21	6:BZ:43:ARG:NH1	1.77	0.82
1:AC:81:VAL:HG11	1:AC:131:PHE:HB3	1.58	0.82
6:AE:21:PHE:CD1	6:AE:22:MET:N	2.47	0.82
6:BB:18:HIS:CE1	6:BB:22:MET:HB2	2.15	0.82
5:BU:14:ILE:HD12	14:BU:103:CRT:C3	2.06	0.82
9:AF:102:BCL:H2	6:AG:28:TRP:CH2	2.14	0.82
5:AS:40:LEU:HD11	5:AS:47:LEU:HD23	1.59	0.82
9:A1:102:BCL:H92	14:A2:102:CRT:H183	1.59	0.82
6:A2:17:PHE:HD1	14:A2:102:CRT:C6	1.93	0.82
3:BM:84:PHE:CZ	5:BW:37:MET:HG2	2.14	0.82
14:BB:102:CRT:H342	9:BD:102:BCL:HBA1	1.60	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:B9:13:LEU:O	6:B0:7:THR:HB	1.80	0.82
5:BI:9:TYR:HB2	6:BJ:15:LYS:HA	1.58	0.82
5:BQ:43:ASP:HA	5:BS:47:LEU:O	1.78	0.82
1:BC:17:SER:HB3	3:BM:91:PHE:CZ	2.13	0.82
5:BS:4:MET:SD	5:BS:8:LEU:HD21	2.20	0.82
5:B7:44:LEU:HD21	5:B7:46:TRP:HE3	1.43	0.82
1:AC:135:ARG:HG2	1:AC:330:LEU:HA	1.61	0.82
6:AG:33:VAL:O	6:AG:37:LEU:HB2	1.79	0.82
6:AG:45:TRP:O	6:AG:46:LEU:HB2	1.78	0.82
5:AI:39:VAL:HG11	9:AI:102:BCL:HBC1	1.59	0.82
5:AU:18:ARG:H	5:AU:18:ARG:HD2	1.44	0.82
6:A2:46:LEU:HD22	6:A4:42:TYR:HE2	1.41	0.82
6:A8:43:ARG:NH2	5:A9:55:TYR:HB2	1.94	0.82
2:BL:17:LEU:HD11	2:BL:114:VAL:HB	1.60	0.82
4:BH:195:LEU:HD12	4:BH:196:PRO:HD2	1.60	0.82
5:AF:11:ILE:HD12	5:AF:14:ILE:HD11	1.61	0.82
9:A1:102:BCL:C9	14:A2:102:CRT:C18	2.57	0.82
2:BL:158:GLY:HA3	2:BL:161:SER:HB3	1.60	0.82
3:AM:136:ARG:HA	3:AM:136:ARG:HH11	1.41	0.82
3:AM:140:LEU:HD23	3:AM:142:MET:HG3	1.59	0.82
5:AK:12:TRP:NE1	6:AN:17:PHE:HD2	1.78	0.82
5:AY:36:HIS:CE1	9:AZ:101:BCL:HMD3	2.15	0.82
4:BH:128:GLU:H	4:BH:128:GLU:CD	1.88	0.82
6:BB:22:MET:HG3	6:BB:26:TYR:CE1	2.14	0.82
6:BR:10:THR:HG22	6:BR:11:ASP:H	1.45	0.82
5:BU:10:LYS:O	14:BU:103:CRT:C2	2.20	0.82
6:BZ:46:LEU:HD21	6:B2:42:TYR:HE2	1.43	0.82
15:AH:301:PEF:H41	15:AH:301:PEF:H12	1.61	0.82
3:BM:7:ILE:HB	15:BM:407:PEF:N	1.93	0.82
5:BA:17:PRO:O	5:BA:21:LEU:HB2	1.79	0.82
9:BO:102:BCL:ND	9:BP:101:BCL:CMD	2.43	0.82
14:BP:102:CRT:H2M3	5:BQ:36:HIS:CB	2.09	0.82
1:AC:200:LEU:HD11	1:AC:238:ASN:ND2	1.94	0.82
4:AH:130:LEU:HG	4:AH:131:PRO:HD2	1.61	0.82
6:AB:17:PHE:CE1	14:AB:102:CRT:H9	2.15	0.82
5:AF:19:ARG:CZ	5:AI:18:ARG:HH21	1.92	0.82
9:AJ:101:BCL:CHC	9:AK:102:BCL:HBB3	2.10	0.82
14:AW:102:CRT:C18	9:AY:102:BCL:H92	2.08	0.82
14:A5:103:CRT:H14	5:A7:21:LEU:HD22	1.62	0.82
6:BE:44:PRO:HG2	5:BF:55:TYR:OH	1.79	0.82
5:BI:50:ASN:CG	5:BI:51:ILE:H	1.85	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:BO:7:ASN:HD22	6:BR:20:ILE:HD12	1.45	0.82
9:BQ:103:BCL:HBC2	9:BQ:104:BCL:HMD1	1.61	0.82
6:B2:21:PHE:CZ	9:B3:102:BCL:H203	2.14	0.82
9:AZ:101:BCL:C20	6:A2:38:LEU:HD21	2.10	0.81
3:BM:28:LEU:HB3	3:BM:29:PRO:HD2	1.60	0.81
5:BU:12:TRP:CD2	6:BV:17:PHE:HE2	1.98	0.81
5:B9:12:TRP:HE1	6:B0:18:HIS:CB	1.92	0.81
5:AK:16:ASP:HB2	5:AK:19:ARG:HG2	1.60	0.81
5:A5:21:LEU:HD11	9:A5:102:BCL:H141	1.60	0.81
1:BC:98:THR:O	1:BC:103:PRO:HD3	1.80	0.81
3:BM:208:PHE:CZ	3:BM:275:LEU:HD13	2.15	0.81
5:BW:16:ASP:H	5:BW:19:ARG:HE	1.28	0.81
5:BW:49:ASP:CG	5:BW:50:ASN:N	2.37	0.81
5:BY:50:ASN:HB3	5:B1:60:LYS:HA	1.62	0.81
5:B3:46:TRP:CZ3	9:B3:102:BCL:HBC3	2.14	0.81
4:AH:47:GLU:HG3	5:AA:19:ARG:HA	1.61	0.81
6:AG:28:TRP:NE1	6:AG:32:VAL:CG2	2.43	0.81
9:AO:102:BCL:HBC2	9:AP:101:BCL:HHD	1.60	0.81
6:AT:45:TRP:CE3	9:AT:101:BCL:H2C	2.15	0.81
5:AW:2:PHE:HA	5:AW:5:ASN:ND2	1.94	0.81
3:BM:170:SER:C	3:BM:172:ALA:H	1.88	0.81
6:BG:23:GLN:O	6:BG:26:TYR:HB2	1.80	0.81
5:BO:20:VAL:O	5:BO:24:ILE:HG12	1.80	0.81
5:BQ:36:HIS:CE1	9:BQ:104:BCL:CMD	2.63	0.81
9:BW:102:BCL:C1D	9:BX:101:BCL:HMD3	2.10	0.81
9:BW:102:BCL:CHD	9:BX:101:BCL:HMD1	2.10	0.81
5:B1:10:LYS:HD2	6:B4:20:ILE:HG13	1.62	0.81
9:B3:102:BCL:HBC2	9:B4:101:BCL:HMD1	1.63	0.81
6:B6:40:TRP:CZ3	6:B6:44:PRO:HA	2.15	0.81
14:B7:102:CRT:H342	9:B7:103:BCL:CBA	2.04	0.81
3:AM:300:LYS:HE2	3:AM:300:LYS:HA	1.63	0.81
9:AK:102:BCL:HAC2	9:AN:101:BCL:HBC3	1.59	0.81
9:AX:101:BCL:C1	9:AX:101:BCL:CGA	2.58	0.81
6:AE:21:PHE:HD1	6:AE:22:MET:N	1.78	0.81
2:BL:207:THR:HG21	3:BM:238:ILE:HG13	1.60	0.81
5:AA:46:TRP:HB2	6:A0:46:LEU:OXT	1.80	0.81
3:BM:159:VAL:HG11	3:BM:281:GLY:O	1.80	0.81
6:BB:22:MET:O	6:BB:26:TYR:CD1	2.33	0.81
5:BD:9:TYR:CE1	6:BE:11:ASP:HB3	2.16	0.81
5:B3:11:ILE:HA	14:B7:102:CRT:H82	1.61	0.81
5:B9:32:GLY:HA3	9:B9:102:BCL:O1A	1.80	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:B0:21:PHE:HB2	14:B0:101:CRT:C11	2.11	0.81
2:AL:203:ILE:CG2	3:AM:266:HIS:HD1	1.88	0.81
2:AL:207:THR:HG21	3:AM:238:ILE:HG13	1.60	0.81
3:AM:301:HIS:ND1	4:AH:8:TYR:HB3	1.96	0.81
5:BF:4:MET:HG2	6:BJ:23:GLN:CG	2.03	0.81
3:AM:299:VAL:HB	3:AM:304:ALA:HB3	1.62	0.81
4:AH:159:LEU:HD22	4:AH:254:ARG:HH22	1.41	0.81
3:BM:260:VAL:HG12	4:BH:34:ASP:HB3	1.62	0.81
4:BH:124:ASP:HB2	4:BH:233:LEU:HD21	1.62	0.81
5:BU:14:ILE:H	14:BU:103:CRT:H22A	1.44	0.81
6:B6:45:TRP:CD1	6:B6:46:LEU:H	1.99	0.81
5:B9:35:ILE:HG21	9:B0:102:BCL:C4D	2.10	0.81
5:AF:44:LEU:O	5:AF:44:LEU:HD12	1.81	0.81
5:AQ:44:LEU:HD12	5:AQ:46:TRP:HE3	1.46	0.81
5:A5:12:TRP:HZ3	5:A5:17:PRO:HA	1.45	0.81
9:A6:101:BCL:HMC1	9:A7:103:BCL:HBB1	1.61	0.81
1:BC:270:TRP:O	1:BC:273:ILE:HD12	1.80	0.81
3:BM:161:GLY:O	3:BM:165:PRO:HD2	1.81	0.81
5:B3:36:HIS:NE2	9:B4:101:BCL:HMD3	1.95	0.81
3:AM:60:SER:HA	3:AM:128:LEU:HD23	1.63	0.81
5:AD:16:ASP:OD2	5:AD:18:ARG:HG2	1.81	0.81
2:BL:206:VAL:HG12	3:BM:142:MET:HE1	1.63	0.81
3:BM:178:GLY:O	3:BM:182:HIS:HB3	1.81	0.81
4:BH:69:LEU:HB3	4:BH:70:PRO:HD2	1.62	0.81
5:BI:18:ARG:NH1	5:BI:18:ARG:HB3	1.95	0.81
5:BO:13:LEU:O	5:BO:13:LEU:HD23	1.79	0.81
6:BV:20:ILE:HG21	14:BV:102:CRT:C6	2.11	0.81
5:BY:16:ASP:HB3	5:BY:18:ARG:HE	1.44	0.81
6:B2:12:ASP:HA	6:B2:15:LYS:HE2	1.61	0.81
9:B8:101:BCL:C1C	9:B9:102:BCL:HBB3	2.11	0.81
3:AM:260:VAL:HG12	4:AH:34:ASP:HB3	1.61	0.80
5:A1:15:LEU:HD23	5:A3:18:ARG:HH12	1.45	0.80
5:A7:44:LEU:HD21	5:A7:46:TRP:HE3	1.45	0.80
6:BV:20:ILE:CG2	14:BV:102:CRT:C7	2.58	0.80
5:B3:8:LEU:HD23	6:B6:20:ILE:HD11	1.62	0.80
1:AC:242:SER:O	1:AC:313:ALA:HA	1.80	0.80
2:AL:203:ILE:HG13	3:AM:266:HIS:CE1	2.16	0.80
6:AN:29:PHE:O	6:AN:33:VAL:HB	1.79	0.80
9:AO:102:BCL:C1D	9:AP:101:BCL:CMD	2.59	0.80
5:AS:34:LEU:CD1	15:AS:101:PEF:H442	2.12	0.80
5:AA:32:GLY:N	9:AB:101:BCL:HED2	1.95	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:AO:31:LEU:O	5:AO:35:ILE:HG12	1.81	0.80
9:AQ:102:BCL:C1D	9:AR:101:BCL:CMD	2.59	0.80
6:A4:13:GLU:HA	6:A4:16:GLU:CD	2.07	0.80
9:BD:102:BCL:ND	9:BE:101:BCL:HMD3	1.96	0.80
5:BK:45:ASN:C	5:BK:49:ASP:HB3	2.05	0.80
14:B2:102:CRT:C2M	5:B3:40:LEU:HD11	2.12	0.80
6:B4:32:VAL:HG11	9:B4:101:BCL:HBA2	1.61	0.80
2:AL:227:ASP:O	3:AM:51:ILE:HG13	1.81	0.80
6:AR:10:THR:HG22	6:AR:11:ASP:H	1.46	0.80
5:AS:37:MET:HG2	15:AS:101:PEF:C45	2.11	0.80
5:A1:21:LEU:HD11	9:A1:102:BCL:C14	2.12	0.80
5:BU:49:ASP:CG	5:BU:50:ASN:H	1.90	0.80
3:AM:175:VAL:HG13	3:AM:176:PRO:HD2	1.64	0.80
6:A2:17:PHE:HD1	14:A2:102:CRT:H6	1.41	0.80
5:A5:11:ILE:N	14:A5:103:CRT:H82	1.96	0.80
9:BA:101:BCL:HBC2	9:BB:101:BCL:HMD1	1.62	0.80
5:BF:28:GLN:O	9:BG:101:BCL:HED1	1.80	0.80
14:BV:102:CRT:C2M	5:BW:37:MET:HB2	2.11	0.80
6:BX:45:TRP:CE3	9:BX:101:BCL:HAC2	2.17	0.80
2:AL:94:LEU:HA	2:AL:97:ILE:HD12	1.63	0.80
4:BH:54:LYS:HE2	5:BD:23:SER:HA	1.63	0.80
5:B1:19:ARG:HH22	5:B3:18:ARG:NH1	1.80	0.80
6:B2:13:GLU:C	14:B2:102:CRT:H32A	2.06	0.80
2:AL:266:ARG:HH11	2:AL:266:ARG:CB	1.92	0.80
6:AE:23:GLN:HG3	6:AE:24:SER:H	1.46	0.80
5:AW:10:LYS:HB2	14:AW:102:CRT:H83	1.64	0.80
6:BB:22:MET:HG3	6:BB:26:TYR:OH	1.82	0.80
5:BF:29:ILE:HB	9:BF:102:BCL:H43	1.62	0.80
5:B1:18:ARG:HD2	5:B1:19:ARG:H	1.46	0.80
5:AS:34:LEU:HD13	15:AS:101:PEF:H442	1.63	0.80
5:A1:7:ASN:O	5:A1:10:LYS:HG3	1.81	0.80
2:BL:57:GLY:HA3	2:BL:66:GLN:HG2	1.64	0.80
3:BM:164:ARG:HB3	3:BM:165:PRO:HD3	1.63	0.80
3:BM:199:ASN:C	3:BM:203:MET:HE3	2.06	0.80
5:BA:33:LEU:HA	14:B0:101:CRT:H2M3	1.64	0.80
14:BU:103:CRT:H2M1	5:BY:37:MET:CA	2.12	0.80
3:AM:107:PRO:HG2	3:AM:113:GLY:HA2	1.63	0.80
3:AM:215:LEU:HD12	3:AM:218:MET:SD	2.22	0.80
5:AO:7:ASN:HB2	6:AR:20:ILE:HG12	1.62	0.80
5:AO:8:LEU:CD1	6:AP:18:HIS:HE1	1.94	0.80
5:AS:42:THR:HG21	5:AU:47:LEU:HB3	1.63	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:AW:34:LEU:HD21	14:AX:102:CRT:H403	1.62	0.80
6:A4:10:THR:HG22	6:A4:11:ASP:H	1.47	0.80
6:A4:13:GLU:HA	6:A4:16:GLU:CG	2.12	0.80
9:A8:101:BCL:NC	9:A9:102:BCL:HBB3	1.97	0.80
1:BC:73:SER:HB3	1:BC:83:LYS:HB2	1.62	0.80
3:BM:11:VAL:HG13	4:BH:148:ASP:HB3	1.64	0.80
6:BR:27:ALA:O	6:BR:31:LEU:HG	1.82	0.80
6:BR:46:LEU:HB3	6:BT:42:TYR:CE2	2.16	0.80
9:BW:102:BCL:HMD3	6:BX:36:HIS:CD2	2.17	0.80
5:BY:45:ASN:HB3	5:BY:48:ASP:O	1.82	0.80
2:AL:140:LEU:HD23	2:AL:140:LEU:O	1.80	0.80
5:AO:14:ILE:HG23	5:AO:15:LEU:HG	1.64	0.80
5:AS:10:LYS:CD	14:AS:104:CRT:H1M1	2.12	0.80
14:A2:102:CRT:C2M	5:A3:36:HIS:HB3	2.11	0.80
6:A0:40:TRP:CH2	6:A0:46:LEU:HG	2.17	0.80
1:BC:203:PHE:HE1	1:BC:210:ILE:HG12	1.43	0.80
2:BL:87:ALA:H	2:BL:96:GLN:HE22	1.30	0.80
5:BA:35:ILE:HD11	14:BB:102:CRT:H372	1.64	0.80
9:BK:102:BCL:HED1	6:BN:31:LEU:HB3	1.64	0.80
5:BO:18:ARG:NH1	5:BO:18:ARG:HB2	1.94	0.80
5:BW:29:ILE:HA	9:BW:102:BCL:H11	1.64	0.80
1:AC:164:TYR:CB	1:AC:309:THR:HA	2.12	0.79
3:AM:159:VAL:HA	3:AM:163:ILE:CG2	2.11	0.79
5:AQ:43:ASP:HB2	5:AS:47:LEU:HD12	1.63	0.79
6:AT:32:VAL:HG21	9:AT:101:BCL:HBA2	1.63	0.79
9:A3:104:BCL:H61	6:A4:29:PHE:HE1	1.43	0.79
5:A7:44:LEU:HD23	6:A8:43:ARG:HH11	1.47	0.79
6:B2:21:PHE:HE1	14:B2:102:CRT:C16	1.88	0.79
5:B3:13:LEU:HB2	14:B7:102:CRT:C2	2.12	0.79
6:B4:10:THR:HG22	6:B4:11:ASP:H	1.46	0.79
5:B5:19:ARG:O	5:B5:23:SER:HB2	1.83	0.79
5:B7:12:TRP:CZ3	5:B7:17:PRO:HB3	2.17	0.79
5:B7:46:TRP:CD1	5:B7:47:LEU:HD22	2.16	0.79
6:AE:30:GLY:O	6:AE:33:VAL:HG12	1.81	0.79
5:AS:34:LEU:HB2	15:AS:101:PEF:H442	1.64	0.79
6:A6:40:TRP:CZ3	6:A6:44:PRO:HA	2.17	0.79
2:AL:223:THR:HA	2:AL:226:ARG:HB3	1.65	0.79
4:AH:121:LYS:HE3	4:BH:72:ASN:O	1.81	0.79
6:AB:23:GLN:CG	5:A9:4:MET:HE1	2.11	0.79
6:AE:10:THR:HG22	6:AE:11:ASP:H	1.47	0.79
6:AJ:27:ALA:O	6:AJ:31:LEU:HG	1.82	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:AW:36:HIS:CE1	9:AX:101:BCL:HMD3	2.17	0.79
1:BC:311:HIS:HA	1:BC:317:PRO:HG3	1.61	0.79
2:BL:113:GLU:HB3	2:BL:127:PRO:HG3	1.62	0.79
5:BA:44:LEU:HD13	6:BB:43:ARG:NH1	1.97	0.79
9:BA:101:BCL:HBB3	9:B0:102:BCL:CHC	2.13	0.79
5:BD:36:HIS:CE1	9:BE:101:BCL:HMD3	2.16	0.79
6:BP:20:ILE:HD13	14:BP:102:CRT:C5	2.12	0.79
5:BQ:26:ALA:O	5:BQ:29:ILE:HG22	1.82	0.79
6:B2:20:ILE:CG2	14:B2:102:CRT:C8	2.60	0.79
5:AI:26:ALA:O	5:AI:29:ILE:HG22	1.82	0.79
5:AQ:14:ILE:O	5:AS:18:ARG:CZ	2.31	0.79
5:B7:36:HIS:HB3	14:B7:102:CRT:H391	1.62	0.79
9:BG:101:BCL:HBB3	9:BI:102:BCL:C4B	2.12	0.79
6:B2:20:ILE:HG23	14:B2:102:CRT:C9	2.12	0.79
1:AC:42:ASN:HA	2:AL:172:GLN:OE1	1.83	0.79
3:AM:240:HIS:HE1	4:AH:69:LEU:HD11	1.48	0.79
5:AU:19:ARG:NE	5:AW:18:ARG:NH2	2.31	0.79
3:BM:76:LEU:HD23	5:BU:37:MET:HE3	1.65	0.79
9:BA:101:BCL:HBC1	9:BB:101:BCL:HBC3	1.62	0.79
6:BE:10:THR:HG22	6:BE:11:ASP:H	1.45	0.79
14:BP:102:CRT:H342	9:BQ:103:BCL:CBA	2.06	0.79
5:BQ:44:LEU:HD12	5:BQ:46:TRP:HE3	1.46	0.79
9:BW:102:BCL:HMD3	6:BX:36:HIS:HD2	1.45	0.79
5:BY:13:LEU:HD21	6:BZ:10:THR:O	1.82	0.79
9:AL:303:BCL:OBD	3:AM:206:ILE:HD12	1.83	0.79
4:AH:5:ILE:HD11	5:AF:40:LEU:CD1	2.13	0.79
6:AT:29:PHE:CE1	9:AT:101:BCL:H11	2.17	0.79
5:BA:8:LEU:HB3	6:BE:20:ILE:CG2	2.13	0.79
5:BF:44:LEU:HD12	5:BF:44:LEU:O	1.83	0.79
9:BG:101:BCL:HMB3	9:BI:102:BCL:CHB	2.13	0.79
5:B3:51:ILE:HA	5:B3:53:VAL:H	1.47	0.79
1:AC:24:GLU:O	2:AL:263:PHE:HA	1.82	0.79
1:AC:122:TYR:HA	1:AC:125:VAL:HG23	1.65	0.79
5:A3:13:LEU:HD21	6:A4:10:THR:O	1.83	0.79
3:BM:63:PHE:HZ	5:BQ:33:LEU:HD23	1.45	0.79
5:B1:43:ASP:HB2	5:B3:47:LEU:HD12	1.65	0.79
10:AM:403:BPH:C9	15:AM:409:PEF:H222	2.13	0.79
5:AO:51:ILE:HG12	5:AO:52:PRO:HD2	1.63	0.79
5:AS:51:ILE:HB	5:AS:52:PRO:CA	2.12	0.79
5:AU:46:TRP:HA	5:AU:49:ASP:OD1	1.83	0.79
5:A5:43:ASP:OD2	5:A7:47:LEU:HA	1.82	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:A0:32:VAL:HG21	9:A0:102:BCL:CBA	2.12	0.79
1:BC:152:CYS:O	1:BC:156:HIS:HB2	1.83	0.79
5:B7:47:LEU:HD22	5:B7:47:LEU:H	1.48	0.79
1:AC:274:ARG:HA	1:AC:277:ARG:HG2	1.65	0.79
5:AF:49:ASP:CG	5:AF:50:ASN:H	1.90	0.79
6:AP:34:ILE:HD13	6:AP:35:ALA:N	1.98	0.79
5:AU:43:ASP:HA	5:AW:47:LEU:C	2.08	0.79
5:BW:16:ASP:CB	5:BW:19:ARG:HE	1.93	0.79
5:B9:12:TRP:HE1	6:B0:18:HIS:HB2	1.47	0.79
1:AC:263:THR:HB	1:AC:264:PRO:HD2	1.65	0.78
6:AE:13:GLU:OE1	6:AE:13:GLU:N	2.17	0.78
6:AR:32:VAL:HG11	9:AR:101:BCL:HBA2	1.66	0.78
5:AS:13:LEU:CD1	14:AS:104:CRT:H32A	2.12	0.78
5:AW:9:TYR:HA	6:AX:18:HIS:CG	2.17	0.78
5:AY:16:ASP:HB2	5:AY:19:ARG:NH2	1.98	0.78
6:BN:17:PHE:CD1	14:BN:102:CRT:H6	2.18	0.78
5:BO:4:MET:HB2	6:BR:23:GLN:HG3	1.62	0.78
5:BS:24:ILE:HD11	9:BU:102:BCL:H162	1.66	0.78
6:B2:45:TRP:O	6:B2:46:LEU:HG	1.83	0.78
5:B9:35:ILE:HG13	9:B0:102:BCL:O1D	1.83	0.78
3:AM:84:PHE:CZ	5:AW:37:MET:HG2	2.18	0.78
9:BA:101:BCL:ND	9:BB:101:BCL:HMD3	1.98	0.78
6:BP:21:PHE:HE1	14:BP:102:CRT:H19	1.48	0.78
9:BQ:103:BCL:CHD	9:BQ:104:BCL:HMD1	2.12	0.78
1:AC:40:MET:SD	1:AC:252:ASN:HA	2.23	0.78
3:AM:159:VAL:HG11	3:AM:281:GLY:O	1.83	0.78
3:AM:161:GLY:O	3:AM:165:PRO:HD2	1.82	0.78
6:AE:45:TRP:O	6:AE:46:LEU:HG	1.84	0.78
6:AG:29:PHE:O	6:AG:33:VAL:HG23	1.83	0.78
5:A1:5:ASN:CA	5:A1:8:LEU:HB3	2.13	0.78
14:A5:103:CRT:H342	9:A9:102:BCL:CBA	2.13	0.78
9:A6:101:BCL:CHC	9:A7:103:BCL:HBB3	2.13	0.78
14:BU:103:CRT:H2M1	5:BY:37:MET:H	1.41	0.78
5:BY:10:LYS:HB2	14:B2:102:CRT:C8	2.13	0.78
2:AL:12:VAL:HG22	2:AL:13:ARG:N	1.98	0.78
5:AD:10:LYS:HB3	14:AG:102:CRT:H5	1.64	0.78
5:AS:46:TRP:CZ3	9:AS:103:BCL:H2C	2.18	0.78
5:BY:9:TYR:HA	6:BZ:18:HIS:CG	2.18	0.78
1:AC:315:ASN:OD1	1:AC:316:LYS:HG3	1.83	0.78
2:AL:190:PHE:HE1	3:AM:209:LEU:HD21	1.48	0.78
6:AR:21:PHE:HB2	14:AR:102:CRT:H14	1.66	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:A7:50:ASN:CG	5:A7:51:ILE:H	1.89	0.78
1:BC:270:TRP:HA	1:BC:273:ILE:HD12	1.64	0.78
2:BL:8:LYS:HE2	4:BH:87:VAL:HG21	1.64	0.78
6:B4:13:GLU:HA	6:B4:16:GLU:CD	2.09	0.78
5:B5:14:ILE:HG21	5:B7:18:ARG:HG2	1.65	0.78
1:AC:191:ALA:O	1:AC:192:TYR:HB2	1.83	0.78
2:AL:4:LEU:HD12	3:AM:250:LEU:HD12	1.66	0.78
2:AL:266:ARG:HB2	2:AL:266:ARG:NH1	1.94	0.78
3:AM:40:LEU:HD13	3:AM:48:ILE:HD11	1.63	0.78
4:AH:172:VAL:HG23	4:AH:173:ASP:N	1.98	0.78
14:AB:102:CRT:H82	5:A9:10:LYS:HB2	1.65	0.78
6:AX:45:TRP:O	6:AX:46:LEU:HG	1.83	0.78
5:A1:27:PHE:HE2	5:A3:29:ILE:HD11	1.48	0.78
9:A0:102:BCL:H162	9:A0:102:BCL:CBB	2.12	0.78
1:BC:249:PHE:HD1	1:BC:250:CYS:SG	2.06	0.78
2:BL:38:VAL:HG23	2:BL:39:GLY:H	1.48	0.78
6:B4:22:MET:O	6:B4:26:TYR:HB2	1.84	0.78
5:B7:42:THR:HB	5:B9:48:ASP:CG	2.08	0.78
5:AI:30:VAL:O	5:AI:33:LEU:HG	1.84	0.78
5:A3:12:TRP:NE1	6:A4:18:HIS:HB2	1.99	0.78
6:A6:44:PRO:CG	5:A7:52:PRO:HG2	2.11	0.78
1:BC:41:GLU:OE1	2:BL:153:HIS:CD2	2.37	0.78
2:BL:148:MET:HB3	2:BL:153:HIS:ND1	1.98	0.78
5:BI:52:PRO:HG2	5:BI:55:TYR:HE2	1.48	0.78
5:BW:24:ILE:HD11	9:BY:102:BCL:C18	2.13	0.78
5:B5:11:ILE:N	14:B5:103:CRT:H82	1.99	0.78
5:B9:50:ASN:HD22	5:B9:51:ILE:HG12	1.47	0.78
2:AL:210:GLN:HB2	2:AL:213:GLU:HG3	1.65	0.78
3:AM:59:LEU:HD11	5:AQ:29:ILE:HD13	1.64	0.78
6:BE:30:GLY:O	6:BE:33:VAL:HG12	1.84	0.78
9:BI:102:BCL:HBC2	9:BJ:101:BCL:HHD	1.63	0.78
5:BS:24:ILE:HD13	9:BU:102:BCL:H202	1.64	0.78
5:B5:16:ASP:HB2	5:B5:19:ARG:HG2	1.65	0.78
2:AL:12:VAL:HG22	2:AL:13:ARG:H	1.47	0.78
3:AM:250:LEU:HG	3:AM:254:TRP:HE1	1.48	0.78
5:AF:43:ASP:OD1	5:AF:44:LEU:HD23	1.84	0.78
9:AK:102:BCL:HAC2	9:AN:101:BCL:CBC	2.14	0.78
14:AS:104:CRT:H181	9:AU:102:BCL:H92	1.63	0.78
5:A3:13:LEU:HB2	14:A7:102:CRT:C1M	2.13	0.78
5:BU:21:LEU:HD11	6:BV:17:PHE:HE1	1.48	0.78
6:B6:40:TRP:HZ3	6:B6:44:PRO:HA	1.48	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:B0:17:PHE:HE1	14:B0:101:CRT:H11	1.44	0.78
1:AC:153:TYR:CB	1:AC:323:MET:HE1	2.14	0.78
1:AC:274:ARG:HA	1:AC:277:ARG:CG	2.15	0.78
3:AM:70:ILE:CG2	3:AM:118:ALA:HB2	2.14	0.78
9:AK:102:BCL:HED1	6:AN:31:LEU:HB3	1.66	0.78
5:AS:10:LYS:O	14:AS:104:CRT:C3	2.32	0.78
5:AS:31:LEU:HD11	14:AT:102:CRT:H35	1.63	0.78
9:A3:103:BCL:HMD3	6:A4:36:HIS:CE1	2.19	0.78
6:BE:32:VAL:HG21	9:BE:101:BCL:HBA2	1.66	0.78
5:BO:4:MET:CB	6:BR:23:GLN:HB3	2.14	0.78
5:BO:9:TYR:HA	6:BP:18:HIS:CG	2.19	0.78
1:AC:237:MET:SD	2:AL:174:LEU:HD23	2.23	0.77
5:AY:38:ILE:HD12	5:AY:39:VAL:N	1.99	0.77
9:BQ:103:BCL:ND	9:BQ:104:BCL:HMD3	1.98	0.77
5:BU:12:TRP:HZ2	6:BV:21:PHE:HD2	1.28	0.77
1:AC:203:PHE:CE1	1:AC:210:ILE:HG12	2.19	0.77
1:AC:296:LYS:HA	1:AC:301:ASP:O	1.83	0.77
6:AG:45:TRP:CD1	6:AG:46:LEU:N	2.52	0.77
5:A3:33:LEU:HD12	5:A3:34:LEU:N	1.98	0.77
1:BC:122:TYR:HA	1:BC:125:VAL:HG23	1.66	0.77
14:BO:103:CRT:H14	6:BR:21:PHE:HB2	1.67	0.77
5:BY:43:ASP:OD1	5:BY:44:LEU:HD23	1.83	0.77
6:B2:46:LEU:HD22	6:B4:42:TYR:HE2	1.48	0.77
3:AM:276:THR:HG22	3:AM:277:VAL:N	1.98	0.77
9:A9:102:BCL:ND	9:A0:102:BCL:HMD3	2.00	0.77
1:BC:234:GLY:O	1:BC:237:MET:HB2	1.83	0.77
2:BL:17:LEU:HG	2:BL:115:GLU:HG2	1.67	0.77
3:BM:179:ILE:HD13	3:BM:179:ILE:N	1.98	0.77
5:BO:49:ASP:OD1	5:BO:50:ASN:N	2.16	0.77
9:BQ:103:BCL:CBC	9:BQ:104:BCL:HBC3	2.12	0.77
5:BU:11:ILE:HG23	14:BU:103:CRT:H83	0.79	0.77
5:BU:18:ARG:H	5:BU:18:ARG:HD2	1.49	0.77
2:AL:178:TYR:HD2	2:AL:269:PRO:HG3	1.47	0.77
3:AM:159:VAL:HG13	3:AM:285:LEU:HD13	1.66	0.77
6:AB:18:HIS:CE1	6:AB:22:MET:HE2	2.19	0.77
5:AD:36:HIS:CE1	9:AE:101:BCL:HMD3	2.18	0.77
6:AZ:24:SER:O	6:AZ:27:ALA:HB3	1.85	0.77
6:A4:22:MET:O	6:A4:26:TYR:HB2	1.84	0.77
9:BM:401:BCL:HBC1	9:BM:402:BCL:HBD	1.64	0.77
5:BF:26:ALA:O	5:BF:29:ILE:HG22	1.83	0.77
5:BU:11:ILE:CB	14:BU:103:CRT:H83	2.13	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:AO:12:TRP:NE1	6:AP:18:HIS:CA	2.41	0.77
5:AQ:43:ASP:HA	5:AS:47:LEU:C	2.09	0.77
9:AQ:102:BCL:ND	9:AR:101:BCL:HMD3	1.99	0.77
14:AS:104:CRT:H391	5:AW:36:HIS:HB3	1.66	0.77
9:A1:102:BCL:C1D	9:A2:101:BCL:HMD1	2.14	0.77
2:BL:199:HIS:HB2	2:BL:238:ILE:HG12	1.67	0.77
5:BF:51:ILE:HG23	5:BF:52:PRO:HA	1.66	0.77
5:BU:46:TRP:CZ3	9:BU:102:BCL:HAC1	2.19	0.77
5:B5:5:ASN:HA	5:B5:8:LEU:CG	2.13	0.77
5:B7:51:ILE:HD12	5:B7:51:ILE:O	1.84	0.77
2:AL:160:LEU:O	2:AL:163:LEU:HB2	1.85	0.77
3:AM:264:SER:O	3:AM:267:ARG:HB2	1.84	0.77
9:AW:101:BCL:C1	9:AW:101:BCL:O1A	2.33	0.77
9:A8:101:BCL:HMC2	9:A9:102:BCL:HBB1	1.66	0.77
2:BL:37:VAL:HG23	2:BL:38:VAL:H	1.50	0.77
2:BL:276:LEU:HD22	2:BL:276:LEU:H	1.49	0.77
4:BH:231:VAL:HG23	4:BH:235:GLU:HG3	1.66	0.77
14:BA:102:CRT:H23	6:BE:16:GLU:HG3	1.64	0.77
6:BB:17:PHE:HE1	14:BB:102:CRT:H9	1.50	0.77
2:AL:217:THR:H	2:AL:220:HIS:CE1	2.02	0.77
3:AM:293:ASN:CG	3:AM:296:LEU:HG	2.09	0.77
5:AY:9:TYR:HB2	6:AZ:15:LYS:HA	1.65	0.77
1:BC:221:SER:O	1:BC:223:PRO:HD3	1.84	0.77
3:BM:200:PRO:O	3:BM:203:MET:HG2	1.83	0.77
5:AI:35:ILE:O	5:AI:39:VAL:HG23	1.83	0.77
9:AK:102:BCL:C1D	9:AN:101:BCL:CMD	2.62	0.77
6:AR:18:HIS:O	6:AR:22:MET:HB2	1.84	0.77
5:AS:30:VAL:CG2	15:AS:101:PEF:H391	2.09	0.77
5:AW:49:ASP:HB2	5:AY:56:GLN:CB	2.14	0.77
5:AY:36:HIS:NE2	9:AZ:101:BCL:HMD3	2.00	0.77
5:AY:44:LEU:HD22	6:AZ:43:ARG:HD2	1.67	0.77
6:A4:13:GLU:CA	6:A4:16:GLU:HG2	2.14	0.77
1:BC:94:MET:SD	7:BC:501:HEM:ND	2.57	0.77
6:BJ:17:PHE:O	6:BJ:20:ILE:HG22	1.85	0.77
5:B3:2:PHE:HE1	5:B3:5:ASN:ND2	1.82	0.77
2:AL:96:GLN:O	2:AL:100:ILE:HG13	1.85	0.77
6:AG:21:PHE:HD1	6:AG:21:PHE:C	1.87	0.77
9:AK:102:BCL:ND	9:AN:101:BCL:HMD3	2.00	0.77
9:AO:102:BCL:C1D	9:AP:101:BCL:HMD1	2.15	0.77
6:AZ:38:LEU:O	6:AZ:38:LEU:HD23	1.85	0.77
9:BL:303:BCL:OBD	3:BM:206:ILE:HD12	1.85	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:BG:21:PHE:CD1	6:BG:21:PHE:C	2.61	0.77
5:BW:10:LYS:HB2	14:BW:103:CRT:H83	1.65	0.77
5:B7:15:LEU:HB3	5:B7:20:VAL:HG11	1.67	0.77
1:AC:175:PRO:CD	1:AC:179:LYS:HB2	2.14	0.77
9:AF:102:BCL:HED1	6:AG:31:LEU:HD22	1.67	0.77
5:A1:10:LYS:HD3	14:A1:103:CRT:H22A	1.65	0.77
5:A1:21:LEU:HD11	9:A1:102:BCL:H141	1.67	0.77
6:A2:20:ILE:HD12	14:A2:102:CRT:C8	2.14	0.77
5:A7:7:ASN:HB2	5:A7:10:LYS:NZ	1.98	0.77
1:BC:267:THR:HG21	3:BM:314:VAL:CB	2.15	0.77
4:BH:168:SER:HB3	4:BH:183:GLU:CB	2.15	0.77
9:BD:102:BCL:C1D	9:BE:101:BCL:HMD1	2.15	0.77
5:BO:26:ALA:O	5:BO:29:ILE:HG22	1.85	0.77
14:BO:103:CRT:H2M3	5:BS:36:HIS:HB2	1.66	0.77
9:BQ:103:BCL:C1D	9:BQ:104:BCL:HMD3	2.14	0.77
5:AF:11:ILE:H	14:AJ:102:CRT:H82	1.50	0.76
6:AG:38:LEU:HA	6:AG:41:LEU:HD12	1.67	0.76
5:AO:49:ASP:CG	5:AO:50:ASN:H	1.90	0.76
5:AQ:26:ALA:O	5:AQ:29:ILE:HG22	1.85	0.76
5:BA:36:HIS:CB	14:B0:101:CRT:H392	2.12	0.76
6:B6:32:VAL:CG2	9:B6:101:BCL:HBA2	2.07	0.76
4:AH:31:ARG:HA	4:AH:31:ARG:NE	1.96	0.76
5:BF:38:ILE:CD1	5:BI:37:MET:HE1	2.14	0.76
5:BF:44:LEU:HB2	6:BG:43:ARG:NH1	2.00	0.76
1:AC:141:TRP:O	1:AC:145:VAL:HG22	1.84	0.76
1:AC:225:SER:H	1:AC:228:GLN:NE2	1.84	0.76
3:AM:59:LEU:CD1	5:AQ:29:ILE:HG21	2.15	0.76
5:AQ:15:LEU:HD23	5:AS:18:ARG:HD3	1.66	0.76
9:A1:102:BCL:C1	9:A1:102:BCL:CGA	2.63	0.76
5:A3:44:LEU:HD12	5:A3:44:LEU:O	1.85	0.76
3:BM:222:THR:HG21	3:BM:252:TRP:HE1	1.50	0.76
5:BA:18:ARG:H	5:BA:18:ARG:HD2	1.49	0.76
5:BA:36:HIS:NE2	9:BB:101:BCL:HMD3	2.01	0.76
5:BF:44:LEU:HB3	5:BI:55:TYR:OH	1.85	0.76
6:BP:21:PHE:CD1	14:BP:102:CRT:H16	2.20	0.76
4:AH:186:VAL:HG12	4:AH:187:ALA:H	1.50	0.76
5:AK:30:VAL:O	5:AK:33:LEU:HG	1.86	0.76
9:A1:102:BCL:C1D	9:A2:101:BCL:CMD	2.63	0.76
2:BL:188:PHE:C	2:BL:190:PHE:H	1.92	0.76
6:BR:24:SER:O	6:BR:27:ALA:HB3	1.86	0.76
5:BU:18:ARG:H	5:BU:18:ARG:CD	1.96	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:B2:16:GLU:OE2	14:B2:102:CRT:H1M1	1.85	0.76
3:AM:206:ILE:HA	9:AM:402:BCL:HMA3	1.67	0.76
14:AA:102:CRT:H403	5:AD:35:ILE:HD13	1.67	0.76
5:AK:9:TYR:OH	6:AN:11:ASP:HB3	1.85	0.76
6:AP:17:PHE:O	6:AP:20:ILE:HG22	1.86	0.76
6:AP:31:LEU:O	6:AP:34:ILE:HG23	1.86	0.76
3:BM:253:ARG:HH11	3:BM:258:PHE:HA	1.51	0.76
6:BB:21:PHE:HZ	6:BB:25:MET:HE3	1.50	0.76
9:BG:101:BCL:HBB3	9:BI:102:BCL:CHC	2.16	0.76
5:BI:36:HIS:CE1	9:BJ:101:BCL:HMD3	2.21	0.76
6:BZ:38:LEU:HD23	6:BZ:38:LEU:O	1.86	0.76
9:B8:101:BCL:HMC1	9:B9:102:BCL:OBB	1.84	0.76
1:AC:169:ASP:OD1	1:AC:170:PRO:HD2	1.86	0.76
6:AZ:45:TRP:CD2	9:AZ:101:BCL:H2C	2.20	0.76
5:BS:7:ASN:HB3	5:BS:10:LYS:HE3	1.68	0.76
5:BU:12:TRP:CD1	6:BV:17:PHE:CD2	2.74	0.76
5:B7:46:TRP:CZ3	9:B7:103:BCL:HBC3	2.20	0.76
1:AC:280:ASN:HB3	1:AC:304:ARG:CD	2.16	0.76
3:AM:41:GLY:HA3	3:AM:46:ALA:HB2	1.68	0.76
5:AO:20:VAL:O	5:AO:24:ILE:HG12	1.85	0.76
5:BD:49:ASP:HB2	5:BF:56:GLN:HG3	1.66	0.76
5:BI:11:ILE:HA	14:BN:102:CRT:H82	1.67	0.76
1:AC:94:MET:SD	7:AC:501:HEM:ND	2.59	0.76
2:AL:129:ALA:HA	2:AL:247:LEU:HD11	1.66	0.76
3:AM:265:ILE:CG2	3:AM:266:HIS:N	2.47	0.76
4:AH:168:SER:HB3	4:AH:183:GLU:HB3	1.68	0.76
6:AB:20:ILE:CG1	5:A9:7:ASN:HB2	2.16	0.76
9:AN:101:BCL:HBB3	9:AO:102:BCL:C4B	2.15	0.76
14:AR:102:CRT:H2M3	5:AS:36:HIS:HB2	1.66	0.76
9:A6:101:BCL:CMC	9:A7:103:BCL:HBB1	2.16	0.76
2:BL:53:GLY:HA3	2:BL:75:ILE:HD11	1.66	0.76
5:BD:27:PHE:CZ	5:BF:29:ILE:HD11	2.20	0.76
5:B1:52:PRO:HD2	5:B1:55:TYR:CE2	2.21	0.76
3:AM:160:LEU:HD23	3:AM:284:ILE:HG21	1.67	0.76
4:AH:114:ALA:HB2	4:AH:245:GLY:CA	2.16	0.76
6:AV:7:THR:CG2	14:AX:102:CRT:H1M1	2.14	0.76
5:A1:19:ARG:HH21	5:A3:18:ARG:NH2	1.84	0.76
6:A2:21:PHE:HE1	14:A2:102:CRT:H16	1.51	0.76
3:BM:224:LEU:HA	3:BM:227:SER:HB2	1.68	0.76
3:BM:301:HIS:CE1	4:BH:8:TYR:HB3	2.21	0.76
6:BZ:24:SER:O	6:BZ:27:ALA:HB3	1.86	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:B2:21:PHE:CD1	14:B2:102:CRT:C16	2.66	0.76
6:B2:29:PHE:HE1	9:B2:101:BCL:C1	1.99	0.76
5:B3:44:LEU:HD21	9:B4:101:BCL:CBC	2.16	0.76
2:AL:17:LEU:HD11	2:AL:114:VAL:HB	1.68	0.76
2:AL:192:ASN:HD22	2:AL:193:CYS:N	1.83	0.76
5:AD:21:LEU:O	5:AD:25:VAL:HG23	1.86	0.76
6:AP:21:PHE:CD1	6:AP:21:PHE:C	2.63	0.76
6:AP:30:GLY:O	6:AP:33:VAL:HG12	1.85	0.76
6:AR:30:GLY:O	6:AR:33:VAL:HG12	1.86	0.76
5:AU:9:TYR:HA	6:AV:18:HIS:CG	2.21	0.76
5:A7:2:PHE:HD1	5:A7:3:THR:N	1.84	0.76
1:BC:295:ARG:HD2	7:BC:502:HEM:O1D	1.86	0.76
2:BL:126:VAL:HB	2:BL:127:PRO:HD3	1.66	0.76
3:BM:175:VAL:HG13	3:BM:176:PRO:HD2	1.66	0.76
14:BV:102:CRT:C2M	5:BW:37:MET:CB	2.64	0.76
5:B5:45:ASN:C	5:B5:49:ASP:HB3	2.11	0.76
1:AC:97:VAL:HG13	7:AC:502:HEM:HMB2	1.68	0.75
1:AC:205:ASP:HB2	1:AC:304:ARG:HE	1.51	0.75
1:AC:285:TRP:CE3	1:AC:302:PRO:HG3	2.21	0.75
3:AM:79:VAL:CG2	3:AM:85:GLN:HB3	2.16	0.75
9:AA:101:BCL:CHB	9:A0:102:BCL:HMB2	2.16	0.75
14:AB:102:CRT:C3	5:A9:10:LYS:CB	2.42	0.75
5:AQ:10:LYS:HB2	14:AT:102:CRT:H83	1.67	0.75
9:A1:102:BCL:H8	14:A2:102:CRT:H182	1.69	0.75
6:BE:9:LEU:HD22	6:BE:13:GLU:HG3	1.68	0.75
5:B3:8:LEU:O	5:B3:11:ILE:HG13	1.86	0.75
1:AC:36:ARG:HB3	2:AL:79:ASP:OD1	1.86	0.75
1:AC:249:PHE:HD1	1:AC:250:CYS:SG	2.09	0.75
2:AL:177:HIS:HB3	3:AM:183:LEU:CD2	2.16	0.75
5:AI:11:ILE:HG12	9:AK:102:BCL:H141	1.68	0.75
6:AP:13:GLU:HA	6:AP:16:GLU:CD	2.10	0.75
9:AZ:101:BCL:C1B	9:A1:102:BCL:HMB2	2.16	0.75
5:A3:14:ILE:HD13	6:A6:17:PHE:HE2	1.49	0.75
14:A0:101:CRT:H35	9:A0:102:BCL:CMA	2.16	0.75
1:BC:33:ILE:H	1:BC:33:ILE:HD12	1.52	0.75
2:BL:3:MET:SD	2:BL:8:LYS:HA	2.26	0.75
4:BH:31:ARG:HA	4:BH:31:ARG:NE	2.01	0.75
4:BH:159:LEU:HD22	4:BH:254:ARG:HH22	1.51	0.75
5:BD:39:VAL:O	5:BD:43:ASP:HB3	1.87	0.75
5:BI:52:PRO:HG2	5:BI:55:TYR:CE2	2.22	0.75
5:BO:7:ASN:HD22	6:BR:20:ILE:CD1	1.99	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:B2:20:ILE:HG12	14:B2:102:CRT:C8	2.17	0.75
5:AF:27:PHE:HA	5:AF:30:VAL:HG12	1.68	0.75
5:AY:50:ASN:CG	5:AY:51:ILE:H	1.94	0.75
6:A4:13:GLU:C	6:A4:16:GLU:HG2	2.11	0.75
5:A5:52:PRO:O	5:A5:55:TYR:CE2	2.39	0.75
3:BM:76:LEU:CD2	5:BU:37:MET:HE3	2.16	0.75
6:BJ:10:THR:HB	6:BJ:13:GLU:OE2	1.84	0.75
5:BU:12:TRP:CG	6:BV:17:PHE:HD2	2.04	0.75
9:B8:101:BCL:CMC	9:B9:102:BCL:HBB1	2.15	0.75
3:AM:200:PRO:HA	3:AM:203:MET:CG	2.16	0.75
9:AA:101:BCL:HBC2	9:AB:101:BCL:HMD1	1.67	0.75
5:AF:4:MET:HB2	6:AJ:23:GLN:CG	2.13	0.75
5:AS:4:MET:O	5:AS:8:LEU:HG	1.86	0.75
5:A7:37:MET:H	14:A7:102:CRT:C2M	1.92	0.75
2:BL:96:GLN:O	2:BL:100:ILE:HG13	1.85	0.75
6:BE:23:GLN:HG3	6:BE:24:SER:H	1.49	0.75
9:BO:102:BCL:C1D	9:BP:101:BCL:CMD	2.63	0.75
6:BR:18:HIS:O	6:BR:22:MET:HB2	1.86	0.75
2:AL:46:GLY:HA3	10:AL:302:BPH:H9C3	1.68	0.75
6:AJ:17:PHE:O	6:AJ:20:ILE:HG22	1.86	0.75
5:AO:8:LEU:CD1	6:AP:18:HIS:CE1	2.70	0.75
5:AS:10:LYS:CG	14:AS:104:CRT:H1M1	2.17	0.75
1:BC:20:LEU:HG	2:BL:271:TRP:NE1	2.01	0.75
1:BC:242:SER:O	1:BC:313:ALA:HA	1.87	0.75
9:BB:101:BCL:CHC	9:BD:102:BCL:HBB3	2.16	0.75
3:AM:197:TYR:CE1	9:AM:402:BCL:HMC3	2.21	0.75
5:AD:28:GLN:O	5:AD:32:GLY:N	2.20	0.75
5:AF:9:TYR:HA	6:AG:18:HIS:CE1	2.22	0.75
5:AI:11:ILE:N	14:AN:102:CRT:H82	2.01	0.75
5:AW:11:ILE:HG21	9:AY:102:BCL:H162	1.67	0.75
5:A9:2:PHE:N	5:A9:5:ASN:ND2	2.30	0.75
6:BJ:33:VAL:HG23	9:BJ:101:BCL:H143	1.69	0.75
5:BO:43:ASP:HA	5:BQ:48:ASP:CG	2.10	0.75
9:B6:101:BCL:HMC1	9:B7:103:BCL:HBB1	1.68	0.75
3:AM:253:ARG:HA	3:AM:257:GLY:O	1.87	0.75
3:AM:273:ALA:O	3:AM:276:THR:HB	1.87	0.75
6:AN:22:MET:HG3	6:AN:26:TYR:HE2	1.49	0.75
5:AO:36:HIS:O	5:AO:40:LEU:HB2	1.86	0.75
6:AR:24:SER:O	6:AR:27:ALA:HB3	1.87	0.75
6:A2:16:GLU:HB2	14:A2:102:CRT:H1M3	1.68	0.75
5:BD:9:TYR:HB2	6:BE:15:LYS:HA	1.67	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:BR:46:LEU:HB3	6:BT:42:TYR:CZ	2.21	0.75
6:BV:33:VAL:O	6:BV:37:LEU:HD23	1.86	0.75
1:AC:243:LEU:H	1:AC:243:LEU:HD12	1.52	0.75
5:AA:29:ILE:HD11	14:A0:101:CRT:H343	1.69	0.75
9:AA:101:BCL:ND	9:AB:101:BCL:HMD3	2.00	0.75
6:A2:17:PHE:CE1	14:A2:102:CRT:H9	2.19	0.75
4:BH:45:ARG:HA	4:BH:96:PRO:HB3	1.68	0.75
6:BV:17:PHE:HA	14:BV:102:CRT:H41	1.69	0.75
9:BZ:101:BCL:HBB1	9:B1:102:BCL:CMC	2.17	0.75
2:AL:196:LEU:CD2	3:AM:216:PHE:HB2	2.16	0.75
3:AM:208:PHE:HB3	3:AM:276:THR:OG1	1.87	0.75
4:AH:136:MET:HE3	4:AH:172:VAL:HG11	1.69	0.75
5:AO:18:ARG:O	5:AO:22:VAL:HG12	1.87	0.75
5:AS:46:TRP:CE3	9:AS:103:BCL:H2C	2.21	0.75
6:A6:40:TRP:HZ3	6:A6:45:TRP:H	1.32	0.75
1:BC:122:TYR:HA	1:BC:125:VAL:CG2	2.16	0.75
3:BM:56:THR:HG21	3:BM:131:VAL:HG11	1.68	0.75
5:BA:18:ARG:HG3	5:B9:14:ILE:HG23	1.68	0.75
5:AQ:8:LEU:HD23	6:AR:22:MET:HE1	1.69	0.74
5:AW:8:LEU:HD22	5:AW:11:ILE:HD11	1.68	0.74
5:AY:4:MET:HB3	5:AY:8:LEU:HG	1.68	0.74
5:A7:43:ASP:CA	5:A9:48:ASP:HB3	2.16	0.74
6:BG:21:PHE:C	6:BG:21:PHE:HD1	1.95	0.74
6:BP:17:PHE:O	6:BP:20:ILE:HG22	1.86	0.74
14:BU:103:CRT:H11	6:BX:21:PHE:HA	1.69	0.74
9:B1:102:BCL:HBC1	9:B2:101:BCL:HBC3	1.68	0.74
5:B3:36:HIS:CE1	9:B4:101:BCL:HMD3	2.21	0.74
5:B5:46:TRP:HA	5:B5:49:ASP:CG	2.12	0.74
3:AM:179:ILE:HD13	3:AM:179:ILE:N	2.03	0.74
3:AM:243:THR:HG22	4:AH:237:ASP:OD1	1.86	0.74
14:AB:102:CRT:H5	5:A9:10:LYS:HB3	1.69	0.74
5:A5:24:ILE:HG13	9:A7:103:BCL:H201	1.69	0.74
2:BL:129:ALA:HB1	2:BL:247:LEU:HD21	1.68	0.74
2:BL:253:SER:C	9:BL:301:BCL:HED3	2.12	0.74
5:BK:9:TYR:OH	6:BN:11:ASP:HB3	1.87	0.74
5:BO:10:LYS:HB2	14:BO:103:CRT:H5	1.69	0.74
14:BU:103:CRT:H343	9:BY:102:BCL:CBA	2.17	0.74
14:B2:102:CRT:H342	9:B3:102:BCL:CBA	2.16	0.74
1:AC:122:TYR:HA	1:AC:125:VAL:CG2	2.17	0.74
1:AC:235:LEU:HG	1:AC:239:ILE:HD11	1.69	0.74
6:AJ:29:PHE:O	6:AJ:33:VAL:HB	1.88	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:A0:45:TRP:CD1	6:A0:46:LEU:H	2.04	0.74
2:BL:279:PRO:HG2	5:BY:37:MET:SD	2.27	0.74
3:BM:158:LEU:O	3:BM:163:ILE:HG22	1.88	0.74
5:BD:49:ASP:HB2	5:BF:56:GLN:CG	2.18	0.74
5:BO:8:LEU:HG	6:BP:18:HIS:CE1	2.23	0.74
5:BS:36:HIS:CE1	9:BT:101:BCL:CMD	2.70	0.74
3:AM:253:ARG:HB2	3:AM:259:ASN:OD1	1.87	0.74
4:AH:5:ILE:CD1	5:AF:47:LEU:HD12	2.17	0.74
4:AH:133:ILE:HD11	4:AH:171:TRP:HB3	1.67	0.74
5:AF:27:PHE:CZ	5:AI:29:ILE:HD11	2.23	0.74
14:AW:102:CRT:C18	9:AY:102:BCL:C8	2.66	0.74
5:A1:13:LEU:O	6:A2:7:THR:HA	1.86	0.74
5:A5:4:MET:CG	6:A8:27:ALA:CB	2.59	0.74
1:BC:32:GLN:HB2	2:BL:80:LEU:HD12	1.69	0.74
2:BL:273:ASN:HA	2:BL:276:LEU:HD23	1.68	0.74
4:BH:55:VAL:HG13	4:BH:56:VAL:N	2.02	0.74
5:BQ:19:ARG:NH1	15:BQ:101:PEF:H51	2.02	0.74
6:AB:44:PRO:C	5:AD:52:PRO:HG3	2.12	0.74
6:AG:46:LEU:HB3	6:AJ:42:TYR:OH	1.86	0.74
9:A8:101:BCL:O1D	9:A8:101:BCL:H2A	1.87	0.74
3:BM:250:LEU:HG	3:BM:254:TRP:NE1	2.02	0.74
5:BA:19:ARG:NH1	5:BD:22:VAL:HG11	2.02	0.74
5:BA:27:PHE:HA	5:BA:30:VAL:HG12	1.68	0.74
5:BD:44:LEU:HD12	5:BD:44:LEU:O	1.86	0.74
6:BG:30:GLY:O	6:BG:33:VAL:HG12	1.87	0.74
5:BK:12:TRP:NE1	6:BN:17:PHE:HD2	1.84	0.74
5:BQ:44:LEU:HD22	6:BR:43:ARG:HD2	1.69	0.74
5:BS:20:VAL:O	5:BS:24:ILE:HG12	1.88	0.74
9:AL:301:BCL:HBC3	9:AL:301:BCL:HHD	1.68	0.74
3:AM:171:TRP:HA	3:AM:171:TRP:CE3	2.22	0.74
5:AI:10:LYS:C	14:AN:102:CRT:H82	2.13	0.74
6:AN:20:ILE:H	6:AN:20:ILE:HD12	1.53	0.74
5:AW:12:TRP:CZ2	6:AX:21:PHE:CG	2.76	0.74
14:AX:102:CRT:H343	9:AY:102:BCL:HBA1	1.69	0.74
14:A2:102:CRT:H2M1	5:A3:36:HIS:HB3	1.69	0.74
5:A3:43:ASP:CB	5:A5:47:LEU:HD13	2.02	0.74
9:A8:101:BCL:HMC1	9:A9:102:BCL:OBB	1.88	0.74
5:BA:34:LEU:O	5:BA:38:ILE:HG23	1.86	0.74
5:BY:50:ASN:HD21	6:BZ:43:ARG:HH12	1.32	0.74
5:B9:32:GLY:N	9:B0:102:BCL:HED2	2.01	0.74
1:AC:96:ALA:C	1:AC:98:THR:H	1.94	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:AL:103:ALA:O	2:AL:107:ILE:HG13	1.87	0.74
5:AQ:31:LEU:HG	9:AR:101:BCL:HED3	1.67	0.74
5:AS:20:VAL:HB	9:AU:102:BCL:C20	2.18	0.74
6:A8:33:VAL:CG2	9:A8:101:BCL:C14	2.66	0.74
1:BC:212:ILE:HD13	1:BC:212:ILE:N	2.02	0.74
6:BJ:29:PHE:O	6:BJ:33:VAL:HB	1.88	0.74
6:BV:43:ARG:HH11	5:BW:55:TYR:HB3	1.52	0.74
14:BW:103:CRT:H342	9:B1:102:BCL:CBA	2.15	0.74
5:B1:43:ASP:HB2	5:B3:47:LEU:CD1	2.18	0.74
5:B3:2:PHE:HE1	5:B3:5:ASN:CG	1.96	0.74
1:AC:280:ASN:CB	1:AC:304:ARG:HD2	2.17	0.74
15:AM:407:PEF:H32	4:AH:29:TYR:CE2	2.23	0.74
5:AO:21:LEU:O	5:AO:25:VAL:HG23	1.88	0.74
9:AW:101:BCL:O1D	9:AW:101:BCL:H2A	1.87	0.74
6:AZ:32:VAL:HG11	9:AZ:101:BCL:HBA2	1.70	0.74
5:A9:4:MET:O	5:A9:8:LEU:HG	1.86	0.74
1:BC:167:VAL:HG21	1:BC:297:GLY:HA3	1.68	0.74
3:BM:79:VAL:CG2	3:BM:85:GLN:HB3	2.15	0.74
6:BB:22:MET:HG3	6:BB:26:TYR:CZ	2.23	0.74
5:BD:33:LEU:O	5:BD:37:MET:HG3	1.88	0.74
5:B1:19:ARG:NH2	5:B3:18:ARG:NH1	2.36	0.74
6:B2:21:PHE:CD1	14:B2:102:CRT:C14	2.52	0.74
5:B7:37:MET:N	14:B7:102:CRT:C2M	2.50	0.74
2:AL:188:PHE:HB3	2:AL:249:ALA:HB2	1.68	0.74
5:AY:8:LEU:CD1	6:AZ:22:MET:HE1	2.18	0.74
5:A5:28:GLN:O	5:A5:32:GLY:N	2.21	0.74
5:B1:12:TRP:HZ3	5:B1:20:VAL:HG21	1.53	0.74
4:AH:136:MET:SD	4:AH:170:VAL:HG23	2.28	0.74
5:AI:27:PHE:CE2	5:AK:29:ILE:HD11	2.22	0.74
14:AR:102:CRT:C34	9:AS:103:BCL:HBA1	2.17	0.74
1:BC:296:LYS:HA	1:BC:301:ASP:O	1.88	0.74
2:BL:194:LEU:O	2:BL:198:MET:HG3	1.87	0.74
3:BM:34:PRO:HG3	3:BM:50:PRO:N	2.03	0.74
3:BM:301:HIS:ND1	4:BH:8:TYR:HB3	2.03	0.74
6:BB:29:PHE:O	6:BB:32:VAL:HG12	1.87	0.74
6:BT:42:TYR:CD2	6:BT:43:ARG:HG2	2.22	0.74
5:BU:16:ASP:HB3	5:BU:18:ARG:NH1	2.02	0.74
1:AC:45:ASN:ND2	1:AC:48:GLN:HB2	2.01	0.73
3:AM:59:LEU:HG	3:AM:128:LEU:HD21	1.70	0.73
3:AM:178:GLY:O	3:AM:182:HIS:HB3	1.88	0.73
6:AP:38:LEU:C	6:AP:41:LEU:HD23	2.11	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:A3:27:PHE:HE2	5:A5:29:ILE:HG13	1.52	0.73
14:A7:102:CRT:H31	9:A7:103:BCL:HBA1	1.70	0.73
6:A8:46:LEU:HD22	6:A0:42:TYR:CE2	2.22	0.73
1:BC:41:GLU:OE1	2:BL:153:HIS:NE2	2.21	0.73
2:BL:86:MET:HG3	5:B7:37:MET:HG3	1.68	0.73
3:BM:279:THR:HA	3:BM:282:ILE:HD12	1.70	0.73
4:BH:114:ALA:HB2	4:BH:245:GLY:HA3	1.68	0.73
6:B0:33:VAL:O	6:B0:37:LEU:HG	1.88	0.73
3:AM:25:LYS:HD2	6:AP:8:GLY:HA3	1.69	0.73
5:AF:19:ARG:NH1	5:AI:18:ARG:NH2	2.36	0.73
2:BL:52:TRP:O	2:BL:56:ILE:HG12	1.86	0.73
3:BM:202:HIS:CE1	3:BM:206:ILE:HD11	2.23	0.73
5:BA:52:PRO:HG2	6:B0:44:PRO:HG2	1.70	0.73
5:BS:46:TRP:CD1	5:BS:47:LEU:HD22	2.23	0.73
5:B3:44:LEU:HD12	5:B3:44:LEU:O	1.87	0.73
10:AM:403:BPH:H101	15:AS:101:PEF:H182	1.70	0.73
5:AF:44:LEU:CB	6:AG:43:ARG:HH11	1.98	0.73
6:AG:45:TRP:CD1	6:AG:46:LEU:H	2.06	0.73
5:AQ:52:PRO:HG3	5:AQ:55:TYR:OH	1.87	0.73
5:AU:25:VAL:HG13	9:AU:102:BCL:H52	1.70	0.73
5:AY:13:LEU:CD2	6:AZ:14:ALA:CB	2.67	0.73
5:A7:36:HIS:HB3	14:A7:102:CRT:H391	1.69	0.73
6:A8:32:VAL:HG11	9:A8:101:BCL:HBA2	1.69	0.73
1:BC:263:THR:HB	1:BC:264:PRO:HD2	1.71	0.73
1:BC:281:GLN:OE1	1:BC:285:TRP:CD1	2.41	0.73
5:BI:8:LEU:HD12	6:BJ:18:HIS:HE1	1.53	0.73
5:BS:31:LEU:HD11	14:BS:103:CRT:H35	1.69	0.73
5:BU:14:ILE:HD12	14:BU:103:CRT:H31A	1.69	0.73
5:BW:9:TYR:HA	6:BX:18:HIS:CD2	2.23	0.73
5:B1:19:ARG:HH22	5:B3:18:ARG:CZ	2.02	0.73
1:AC:157:ARG:HE	1:AC:312:GLN:CD	1.96	0.73
4:AH:227:ASN:HD22	4:AH:228:PRO:CD	2.00	0.73
9:AW:101:BCL:ND	9:AX:101:BCL:HMD3	2.03	0.73
6:A2:21:PHE:CD1	14:A2:102:CRT:H14	2.23	0.73
3:BM:34:PRO:HD3	3:BM:50:PRO:HB3	1.68	0.73
6:BB:29:PHE:CE1	9:BB:101:BCL:H11	2.23	0.73
6:BP:10:THR:HG22	6:BP:11:ASP:N	2.04	0.73
5:BY:43:ASP:CA	5:B1:48:ASP:HB3	2.17	0.73
5:B3:14:ILE:HD11	14:B7:102:CRT:H42	1.70	0.73
5:B7:16:ASP:O	5:B7:20:VAL:HG22	1.89	0.73
6:B0:33:VAL:CG1	6:B0:37:LEU:HD11	2.19	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:AM:249:ALA:HB2	13:AM:405:MQ8:H61	1.71	0.73
15:AM:408:PEF:H52	4:AH:204:LYS:HE2	1.68	0.73
4:AH:5:ILE:HD11	5:AF:40:LEU:HD12	1.70	0.73
5:AO:7:ASN:ND2	5:AO:7:ASN:H	1.86	0.73
5:AW:51:ILE:HB	5:AW:52:PRO:CA	2.15	0.73
6:A4:13:GLU:O	6:A4:16:GLU:HG2	1.88	0.73
3:BM:55:LEU:HD23	5:BQ:22:VAL:HG23	1.70	0.73
5:BF:27:PHE:CZ	5:BI:29:ILE:HD11	2.23	0.73
14:BV:102:CRT:H2M1	5:BW:37:MET:HB2	1.69	0.73
5:AD:50:ASN:CG	5:AD:51:ILE:H	1.96	0.73
5:AK:44:LEU:CD2	5:AK:46:TRP:HB3	2.18	0.73
9:A5:102:BCL:O1D	9:A5:102:BCL:H2A	1.89	0.73
6:BB:21:PHE:CZ	6:BB:25:MET:HE3	2.23	0.73
5:BI:44:LEU:HD12	5:BI:44:LEU:O	1.89	0.73
5:BU:31:LEU:HD23	9:BV:101:BCL:HED3	1.69	0.73
6:B2:17:PHE:HD1	14:B2:102:CRT:H9	1.53	0.73
5:B7:44:LEU:HD22	5:B7:46:TRP:CE3	2.23	0.73
1:AC:316:LYS:HG2	7:AC:504:HEM:HAD2	1.68	0.73
2:AL:120:LEU:O	2:AL:122:ILE:HG23	1.87	0.73
3:AM:144:GLN:HB3	3:AM:147:SER:OG	1.88	0.73
5:AK:29:ILE:HB	9:AK:102:BCL:H43	1.70	0.73
5:AO:29:ILE:HD12	9:AO:102:BCL:H11	1.69	0.73
9:AW:101:BCL:HBC2	9:AX:101:BCL:HMD1	1.70	0.73
5:AY:28:GLN:C	9:AY:102:BCL:H12	2.14	0.73
1:BC:164:TYR:CE2	1:BC:312:GLN:HG2	2.24	0.73
2:BL:216:LYS:HD2	2:BL:220:HIS:CD2	2.24	0.73
6:BJ:17:PHE:HE1	6:BJ:21:PHE:HB2	1.54	0.73
6:B2:11:ASP:O	6:B2:15:LYS:HG2	1.88	0.73
5:B9:31:LEU:O	5:B9:35:ILE:HG12	1.88	0.73
1:AC:22:GLY:HA3	2:AL:263:PHE:HB3	1.70	0.73
1:AC:291:LEU:O	1:AC:296:LYS:HE3	1.89	0.73
2:AL:4:LEU:HD21	3:AM:253:ARG:HH21	1.53	0.73
5:AS:10:LYS:CG	14:AS:104:CRT:C1M	2.67	0.73
5:A7:12:TRP:HZ3	5:A7:17:PRO:HB3	1.51	0.73
1:BC:166:TRP:HE1	1:BC:305:VAL:C	1.96	0.73
3:BM:290:VAL:HG12	3:BM:291:VAL:H	1.54	0.73
10:BM:403:BPH:HMA1	15:BQ:101:PEF:C41	2.16	0.73
10:BM:403:BPH:H3A	15:BQ:101:PEF:H431	1.71	0.73
5:B1:18:ARG:HG2	5:B1:18:ARG:HH11	1.53	0.73
14:B1:103:CRT:C34	9:B5:102:BCL:HBA1	2.17	0.73
6:AB:20:ILE:HG12	5:A9:7:ASN:HB2	1.68	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:AF:38:ILE:HD11	5:AI:37:MET:HE1	1.70	0.73
5:A3:52:PRO:O	5:A3:55:TYR:CZ	2.41	0.73
6:BE:33:VAL:O	6:BE:37:LEU:HD23	1.89	0.73
6:BP:45:TRP:O	6:BP:46:LEU:HG	1.87	0.73
6:BV:17:PHE:CA	14:BV:102:CRT:H41	2.19	0.73
5:BW:16:ASP:O	5:BW:19:ARG:HG2	1.88	0.73
5:B1:14:ILE:HD12	5:B1:15:LEU:N	2.04	0.73
3:AM:166:VAL:HG22	3:AM:171:TRP:CZ3	2.23	0.73
5:A1:10:LYS:HB3	14:A1:103:CRT:H5	1.70	0.73
5:A5:25:VAL:CG1	9:A5:102:BCL:C19	2.54	0.73
5:A9:36:HIS:NE2	9:A0:102:BCL:HMD3	2.03	0.73
5:BD:32:GLY:N	9:BE:101:BCL:HED2	2.04	0.73
9:BI:102:BCL:OBD	6:BJ:32:VAL:HG13	1.88	0.73
9:BX:101:BCL:CMC	5:BY:47:LEU:HD21	2.19	0.73
2:AL:235:ALA:HA	11:AL:304:UQ8:H3MB	1.69	0.72
4:AH:168:SER:HB3	4:AH:183:GLU:CB	2.19	0.72
14:AA:102:CRT:H9	6:AE:17:PHE:CD1	2.23	0.72
6:AG:23:GLN:O	6:AG:26:TYR:HB2	1.88	0.72
5:AK:52:PRO:HB2	5:AK:55:TYR:CD1	2.23	0.72
6:AR:45:TRP:CD1	6:AR:46:LEU:H	2.06	0.72
14:AT:102:CRT:H342	9:AU:102:BCL:HBA2	1.69	0.72
5:A5:43:ASP:HB2	5:A7:47:LEU:CB	2.18	0.72
14:A5:103:CRT:H293	9:A9:102:BCL:C4	2.14	0.72
1:BC:130:MET:HE1	1:BC:284:ILE:HD11	1.71	0.72
3:BM:57:GLY:O	3:BM:61:ILE:HG13	1.89	0.72
3:AM:200:PRO:N	3:AM:203:MET:HE3	2.04	0.72
6:AG:29:PHE:O	6:AG:33:VAL:CG2	2.37	0.72
9:AI:102:BCL:HBC2	9:AJ:101:BCL:HMD1	1.71	0.72
5:AS:47:LEU:HD22	5:AS:47:LEU:H	1.54	0.72
14:AT:102:CRT:H342	9:AU:102:BCL:HBA1	1.71	0.72
5:AY:33:LEU:HD12	5:AY:34:LEU:N	2.04	0.72
6:AZ:22:MET:HG3	6:AZ:26:TYR:HE1	1.53	0.72
6:AZ:46:LEU:HB3	5:A1:52:PRO:HD3	1.71	0.72
5:A3:12:TRP:HE1	6:A4:18:HIS:HB2	1.54	0.72
6:A4:40:TRP:CZ3	6:A4:44:PRO:HA	2.24	0.72
2:BL:190:PHE:HE1	3:BM:209:LEU:HD21	1.53	0.72
3:BM:222:THR:CG2	3:BM:252:TRP:HE1	2.02	0.72
9:BW:102:BCL:C1D	9:BX:101:BCL:HMD1	2.18	0.72
5:B1:18:ARG:O	5:B1:22:VAL:HG12	1.89	0.72
6:B6:27:ALA:O	6:B6:31:LEU:HG	1.89	0.72
5:B9:31:LEU:HD11	5:B9:35:ILE:HD11	1.71	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:AL:10:TYR:OH	3:AM:246:GLU:HG2	1.89	0.72
2:AL:144:ARG:HB3	2:AL:145:PRO:HD3	1.71	0.72
5:AI:18:ARG:HB3	5:AI:18:ARG:CZ	2.19	0.72
6:AX:17:PHE:CE1	14:AX:102:CRT:H6	2.24	0.72
5:A7:25:VAL:HG13	9:A7:103:BCL:H52	1.72	0.72
9:BO:102:BCL:HBD	9:BP:101:BCL:OBD	1.90	0.72
6:B4:29:PHE:HZ	9:B4:101:BCL:H101	1.52	0.72
6:B0:40:TRP:HH2	6:B0:46:LEU:CG	2.00	0.72
2:AL:48:LEU:HA	2:AL:51:VAL:HG23	1.71	0.72
3:AM:279:THR:HA	3:AM:282:ILE:CD1	2.19	0.72
6:AE:9:LEU:HD22	6:AE:13:GLU:HG3	1.71	0.72
5:AS:13:LEU:HD12	14:AS:104:CRT:H32A	1.68	0.72
1:BC:135:ARG:HG2	1:BC:330:LEU:C	2.14	0.72
3:BM:164:ARG:HA	3:BM:167:MET:HB3	1.71	0.72
6:BB:24:SER:HA	5:B9:4:MET:HE2	1.69	0.72
9:BE:101:BCL:HHB	9:BF:102:BCL:HMA1	1.70	0.72
5:BK:44:LEU:CD2	5:BK:46:TRP:HB3	2.19	0.72
5:B3:5:ASN:HA	5:B3:8:LEU:CD1	2.18	0.72
5:B3:21:LEU:O	5:B3:25:VAL:HG23	1.89	0.72
5:B7:37:MET:H	14:B7:102:CRT:H2M1	1.54	0.72
3:AM:215:LEU:C	3:AM:217:ALA:H	1.95	0.72
6:AE:29:PHE:CZ	9:AE:101:BCL:H42	2.24	0.72
5:AW:54:SER:HA	17:AW:201:HOH:O	1.88	0.72
2:BL:4:LEU:H	2:BL:7:GLU:HB3	1.54	0.72
2:BL:206:VAL:O	2:BL:209:PRO:HD3	1.88	0.72
14:BB:102:CRT:H2M2	5:BD:37:MET:CE	2.19	0.72
5:BU:12:TRP:CE2	6:BV:17:PHE:CD2	2.76	0.72
5:B7:43:ASP:HB2	5:B9:47:LEU:CD1	2.16	0.72
1:AC:301:ASP:HB2	1:AC:302:PRO:HD2	1.72	0.72
3:AM:171:TRP:HA	3:AM:171:TRP:HE3	1.55	0.72
3:AM:214:LEU:O	3:AM:217:ALA:HB3	1.88	0.72
14:AA:102:CRT:H11	6:AE:17:PHE:HE1	1.53	0.72
5:AO:13:LEU:HD23	5:AO:14:ILE:N	2.05	0.72
5:A5:12:TRP:CZ3	5:A5:17:PRO:HA	2.25	0.72
6:A0:24:SER:O	6:A0:27:ALA:HB3	1.90	0.72
6:A0:32:VAL:CG2	9:A0:102:BCL:HBA2	2.18	0.72
3:BM:40:LEU:HD13	3:BM:48:ILE:HD11	1.70	0.72
9:BD:102:BCL:CHD	9:BE:101:BCL:HMD1	2.20	0.72
5:BI:39:VAL:HG11	9:BI:102:BCL:HBC1	1.71	0.72
5:BU:36:HIS:CE1	9:BV:101:BCL:HMD3	2.23	0.72
9:B1:102:BCL:HAC2	9:B2:101:BCL:HBC1	1.70	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:B6:101:BCL:CHC	9:B7:103:BCL:HBB3	2.20	0.72
1:AC:154:THR:HG22	1:AC:155:CYS:N	2.04	0.72
3:AM:58:THR:O	3:AM:61:ILE:HG22	1.88	0.72
4:AH:128:GLU:CD	4:AH:128:GLU:H	1.97	0.72
5:AW:12:TRP:CE3	5:AW:12:TRP:HA	2.23	0.72
9:AW:101:BCL:HMD1	6:AX:36:HIS:HA	1.71	0.72
14:AW:102:CRT:H35	5:AY:31:LEU:HD11	1.71	0.72
9:A1:102:BCL:ND	9:A2:101:BCL:HMD3	2.05	0.72
2:BL:148:MET:SD	2:BL:262:PRO:HG3	2.29	0.72
3:BM:179:ILE:O	3:BM:183:LEU:HB2	1.90	0.72
6:B4:13:GLU:C	6:B4:16:GLU:HG2	2.15	0.72
6:B0:45:TRP:CD1	6:B0:46:LEU:H	2.07	0.72
1:AC:121:ILE:HG23	1:AC:123:THR:HG23	1.72	0.72
2:AL:17:LEU:HG	2:AL:115:GLU:HG2	1.71	0.72
3:AM:215:LEU:O	3:AM:218:MET:N	2.22	0.72
5:AS:11:ILE:HG12	14:AS:104:CRT:C8	2.19	0.72
5:AW:14:ILE:HG21	5:AY:21:LEU:HD12	1.71	0.72
5:A3:52:PRO:O	5:A3:55:TYR:CE1	2.42	0.72
6:A8:29:PHE:HZ	9:A8:101:BCL:C7	2.03	0.72
6:A0:21:PHE:HB2	14:A0:101:CRT:C14	2.19	0.72
1:BC:157:ARG:HH12	1:BC:318:LEU:CG	2.02	0.72
3:BM:200:PRO:HA	3:BM:203:MET:SD	2.30	0.72
6:BN:44:PRO:HD2	5:BO:55:TYR:OH	1.90	0.72
9:BW:102:BCL:CBC	9:BX:101:BCL:HBC3	2.20	0.72
5:BY:10:LYS:HG3	6:B2:20:ILE:HD13	1.71	0.72
5:B7:36:HIS:CG	14:B7:102:CRT:H393	2.25	0.72
5:AA:10:LYS:O	5:AA:13:LEU:HD13	1.90	0.72
5:AS:31:LEU:O	5:AS:35:ILE:HG12	1.90	0.72
6:A6:27:ALA:O	6:A6:31:LEU:HG	1.89	0.72
5:A7:29:ILE:HG23	5:A7:30:VAL:N	2.04	0.72
1:BC:153:TYR:CB	1:BC:323:MET:HE1	2.18	0.72
5:BO:29:ILE:CA	9:BO:102:BCL:H11	2.18	0.72
2:AL:29:PRO:CB	3:AM:253:ARG:HD2	2.20	0.72
3:AM:170:SER:C	3:AM:172:ALA:H	1.98	0.72
4:AH:65:LYS:N	4:AH:78:ALA:O	2.23	0.72
5:AA:28:GLN:HB3	9:AA:101:BCL:H2	1.71	0.72
6:AG:16:GLU:O	6:AG:20:ILE:HG22	1.89	0.72
6:AG:31:LEU:O	6:AG:34:ILE:HG23	1.89	0.72
14:AS:104:CRT:H2M1	5:AW:37:MET:HB2	1.70	0.72
14:BA:102:CRT:H403	5:BD:35:ILE:HD13	1.72	0.72
9:BO:102:BCL:HAC2	9:BP:101:BCL:HAC1	1.71	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:BW:46:TRP:CH2	9:BW:102:BCL:H2C	2.24	0.72
1:AC:20:LEU:HD22	1:AC:21:LEU:N	2.05	0.71
3:AM:73:PHE:HZ	5:AS:38:ILE:HA	1.53	0.71
5:AS:37:MET:CG	15:AS:101:PEF:H452	2.19	0.71
1:BC:254:ARG:HH21	3:BM:295:TYR:HE1	1.36	0.71
3:BM:63:PHE:CE2	3:BM:124:LEU:HB2	2.25	0.71
5:BD:33:LEU:O	5:BD:37:MET:CG	2.38	0.71
5:BU:46:TRP:CH2	9:BU:102:BCL:H2C	2.24	0.71
5:B5:30:VAL:HG13	5:B5:31:LEU:N	2.05	0.71
5:B7:42:THR:HB	5:B9:48:ASP:OD2	1.90	0.71
6:B0:29:PHE:HD1	9:B0:102:BCL:H11	1.54	0.71
2:AL:188:PHE:C	2:AL:190:PHE:H	1.97	0.71
3:AM:199:ASN:HD21	3:AM:283:GLY:HA3	1.55	0.71
5:AI:27:PHE:O	5:AI:30:VAL:HG12	1.89	0.71
5:A7:46:TRP:CD1	5:A7:47:LEU:HD22	2.24	0.71
3:BM:124:LEU:O	3:BM:128:LEU:N	2.23	0.71
4:BH:31:ARG:HB3	4:BH:59:PRO:HG3	1.71	0.71
4:BH:45:ARG:HD3	4:BH:97:GLY:N	2.03	0.71
9:BE:101:BCL:C1B	9:BF:102:BCL:HMB3	2.20	0.71
6:BZ:46:LEU:HB2	5:B1:52:PRO:CD	2.19	0.71
5:B7:17:PRO:O	5:B7:21:LEU:HG	1.89	0.71
5:B9:36:HIS:NE2	9:B0:102:BCL:HMD3	2.04	0.71
3:AM:73:PHE:CZ	5:AS:38:ILE:HA	2.25	0.71
3:AM:208:PHE:CZ	3:AM:275:LEU:HD13	2.26	0.71
9:AE:101:BCL:NB	9:AF:102:BCL:HMB2	2.04	0.71
5:AU:44:LEU:HD22	6:AV:43:ARG:HD3	1.71	0.71
6:A2:20:ILE:HD12	14:A2:102:CRT:H81	1.71	0.71
5:A9:44:LEU:O	5:A9:44:LEU:HD22	1.91	0.71
2:BL:196:LEU:HB2	3:BM:216:PHE:CD2	2.25	0.71
3:BM:228:ARG:H	3:BM:228:ARG:HD2	1.55	0.71
5:BA:36:HIS:CD2	9:BB:101:BCL:HMD2	2.25	0.71
5:BI:17:PRO:O	5:BI:21:LEU:HB3	1.90	0.71
9:BK:102:BCL:HMD3	6:BN:36:HIS:CD2	2.24	0.71
6:BP:44:PRO:O	5:BQ:52:PRO:HG2	1.91	0.71
5:BY:31:LEU:O	5:BY:35:ILE:HG12	1.89	0.71
3:AM:138:GLU:C	3:AM:140:LEU:H	1.97	0.71
4:AH:225:LEU:O	4:AH:225:LEU:HD12	1.90	0.71
5:AI:31:LEU:HB3	9:AJ:101:BCL:HED3	1.73	0.71
9:AR:101:BCL:HMA1	9:AS:103:BCL:HMA1	1.71	0.71
9:BI:102:BCL:C1D	9:BJ:101:BCL:CMD	2.68	0.71
5:BQ:43:ASP:HB2	5:BS:47:LEU:HA	1.71	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:B4:29:PHE:HE1	9:B4:101:BCL:H72	1.55	0.71
5:AO:11:ILE:N	14:AR:102:CRT:H82	2.06	0.71
5:AU:12:TRP:HA	5:AU:12:TRP:CE3	2.23	0.71
5:A7:44:LEU:O	5:A7:44:LEU:HD22	1.90	0.71
9:A9:102:BCL:C1D	9:A0:102:BCL:CMD	2.69	0.71
2:BL:2:ALA:N	4:BH:45:ARG:HB2	2.04	0.71
5:BA:14:ILE:HG13	5:BA:15:LEU:CD2	2.20	0.71
14:BB:102:CRT:C2M	5:BD:37:MET:HE3	2.21	0.71
6:B0:24:SER:O	6:B0:27:ALA:HB3	1.90	0.71
2:AL:43:THR:HA	10:AL:302:BPH:H7C1	1.72	0.71
3:AM:260:VAL:HB	3:AM:264:SER:OG	1.91	0.71
4:AH:189:ASN:HB3	4:AH:191:LYS:HG3	1.71	0.71
3:BM:197:TYR:CE1	9:BM:402:BCL:HMC3	2.26	0.71
9:BK:102:BCL:C1D	9:BN:101:BCL:HMD1	2.20	0.71
5:BW:16:ASP:H	5:BW:19:ARG:NE	1.88	0.71
5:B3:2:PHE:CE1	5:B3:5:ASN:ND2	2.58	0.71
5:B9:12:TRP:NE1	6:B0:18:HIS:HB2	2.04	0.71
2:AL:148:MET:SD	2:AL:262:PRO:HG3	2.30	0.71
2:AL:177:HIS:NE2	9:AL:301:BCL:HMC2	2.06	0.71
4:AH:52:ARG:HB2	4:AH:54:LYS:NZ	2.05	0.71
6:AG:28:TRP:CE2	6:AG:32:VAL:CG2	2.74	0.71
9:AQ:102:BCL:C1D	9:AR:101:BCL:HMD3	2.21	0.71
5:A1:19:ARG:O	5:A1:23:SER:HB3	1.91	0.71
2:BL:12:VAL:HG22	2:BL:13:ARG:N	2.04	0.71
6:BG:25:MET:HE2	9:BI:102:BCL:H203	1.73	0.71
5:BU:19:ARG:NE	5:BW:18:ARG:NH2	2.38	0.71
5:B7:18:ARG:O	5:B7:22:VAL:HG12	1.91	0.71
5:B7:36:HIS:NE2	9:B8:101:BCL:HMD3	2.06	0.71
3:AM:126:ILE:HG12	9:AM:402:BCL:H142	1.72	0.71
5:AA:50:ASN:CG	6:AB:43:ARG:NH2	2.49	0.71
14:AA:102:CRT:C2	6:AE:16:GLU:HG3	2.20	0.71
6:AX:25:MET:HG3	14:AX:102:CRT:H21	1.71	0.71
1:BC:270:TRP:HE3	1:BC:271:TYR:CD1	2.08	0.71
5:BA:55:TYR:CE1	5:B9:44:LEU:HB3	2.18	0.71
6:B2:21:PHE:HA	14:B2:102:CRT:C11	2.21	0.71
5:B5:43:ASP:CG	5:B7:47:LEU:O	2.33	0.71
6:B8:45:TRP:O	6:B8:46:LEU:HG	1.91	0.71
9:B9:102:BCL:HBC2	9:B0:102:BCL:HMD1	1.73	0.71
6:B0:33:VAL:HG12	6:B0:37:LEU:HD11	1.71	0.71
4:AH:69:LEU:HB3	4:AH:70:PRO:HD2	1.72	0.71
5:AY:13:LEU:HD21	6:AZ:14:ALA:HB1	1.71	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:BM:60:SER:CA	3:BM:128:LEU:HD23	2.21	0.71
5:BO:50:ASN:CG	5:BO:51:ILE:H	1.98	0.71
6:BP:38:LEU:O	6:BP:38:LEU:HD23	1.90	0.71
6:BP:46:LEU:HB3	6:BR:42:TYR:OH	1.91	0.71
5:BY:30:VAL:HA	5:BY:33:LEU:HG	1.72	0.71
5:B9:17:PRO:O	5:B9:21:LEU:HB2	1.91	0.71
5:AA:11:ILE:N	14:AA:102:CRT:H82	2.06	0.71
5:AA:27:PHE:HA	5:AA:30:VAL:HG12	1.73	0.71
5:AO:38:ILE:HD13	14:AP:102:CRT:H403	1.73	0.71
5:AS:17:PRO:O	5:AS:21:LEU:HG	1.91	0.71
9:AV:102:BCL:H2A	9:AV:102:BCL:O1D	1.89	0.71
1:BC:62:LEU:HG	1:BC:327:TYR:OH	1.90	0.71
1:BC:276:VAL:HG13	1:BC:277:ARG:N	2.03	0.71
2:BL:243:LEU:HD23	11:BL:304:UQ8:H46B	1.73	0.71
6:BV:27:ALA:O	6:BV:31:LEU:HG	1.91	0.71
5:BY:48:ASP:O	5:BY:49:ASP:HB3	1.89	0.71
5:B7:44:LEU:HD23	6:B8:43:ARG:HH11	1.56	0.71
5:B9:40:LEU:HD13	5:B9:47:LEU:HD23	1.71	0.71
2:AL:71:TRP:HD1	3:AM:303:MET:HG2	1.54	0.70
3:AM:28:LEU:HB3	3:AM:29:PRO:HD2	1.72	0.70
3:AM:71:ILE:HD13	3:AM:177:PHE:CE1	2.26	0.70
3:AM:156:PHE:CD2	9:AM:402:BCL:H52	2.26	0.70
6:A0:36:HIS:HE1	9:A0:102:BCL:C1B	2.04	0.70
2:BL:44:LEU:C	2:BL:46:GLY:H	1.99	0.70
2:BL:206:VAL:HG11	2:BL:221:GLU:HG2	1.73	0.70
5:BU:43:ASP:CA	5:BW:47:LEU:O	2.36	0.70
14:BV:102:CRT:H342	9:BW:102:BCL:CBA	2.20	0.70
6:B8:20:ILE:O	6:B8:23:GLN:HG3	1.90	0.70
1:AC:157:ARG:NH1	1:AC:318:LEU:HD21	2.05	0.70
1:AC:167:VAL:HG21	1:AC:297:GLY:HA3	1.71	0.70
6:AB:40:TRP:HZ3	6:AB:45:TRP:N	1.85	0.70
14:AB:102:CRT:H342	9:AD:102:BCL:HBA1	1.73	0.70
5:AY:38:ILE:HD12	5:AY:39:VAL:H	1.54	0.70
2:BL:140:LEU:O	2:BL:140:LEU:HD23	1.91	0.70
2:BL:144:ARG:HB3	2:BL:145:PRO:HD3	1.73	0.70
3:BM:208:PHE:HZ	3:BM:275:LEU:HD13	1.53	0.70
3:BM:253:ARG:HB2	3:BM:259:ASN:OD1	1.90	0.70
5:BA:33:LEU:HA	14:B0:101:CRT:C2M	2.22	0.70
5:BQ:51:ILE:CG1	5:BQ:52:PRO:HA	2.20	0.70
5:BU:11:ILE:HA	14:BU:103:CRT:H5	1.73	0.70
14:BW:103:CRT:H2M1	5:B1:36:HIS:HB3	1.74	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:B2:41:LEU:C	6:B2:41:LEU:HD12	2.15	0.70
6:B0:32:VAL:CG2	9:B0:102:BCL:HBA2	2.19	0.70
9:AL:303:BCL:O1D	9:AL:303:BCL:H2A	1.91	0.70
4:AH:78:ALA:HA	4:AH:79:PRO:C	2.16	0.70
6:AN:33:VAL:O	6:AN:37:LEU:HD23	1.90	0.70
5:AO:11:ILE:O	5:AO:14:ILE:HG22	1.90	0.70
5:AW:7:ASN:HD22	5:AW:7:ASN:N	1.86	0.70
5:A1:51:ILE:HB	5:A1:52:PRO:HA	1.74	0.70
5:A3:5:ASN:CG	6:A4:22:MET:HE3	2.15	0.70
5:A3:11:ILE:HA	14:A7:102:CRT:H82	1.73	0.70
6:A8:20:ILE:O	6:A8:23:GLN:HG3	1.91	0.70
4:BH:77:VAL:HG23	4:BH:80:ARG:HB3	1.71	0.70
5:BF:7:ASN:CG	6:BJ:20:ILE:HD13	2.16	0.70
5:BU:21:LEU:HD11	6:BV:17:PHE:CE1	2.27	0.70
14:BV:102:CRT:C2M	5:BW:33:LEU:O	2.37	0.70
6:BZ:22:MET:O	6:BZ:25:MET:HB3	1.91	0.70
5:B9:44:LEU:O	5:B9:44:LEU:HD22	1.90	0.70
3:AM:63:PHE:HB3	3:AM:125:SER:HB2	1.72	0.70
5:AA:43:ASP:CA	5:AD:48:ASP:HB3	2.19	0.70
6:AJ:17:PHE:HA	6:AJ:20:ILE:HG22	1.73	0.70
1:BC:280:ASN:HB3	1:BC:304:ARG:CD	2.22	0.70
2:BL:10:TYR:HA	4:BH:112:GLY:CA	2.22	0.70
3:BM:34:PRO:CG	3:BM:50:PRO:HD3	2.22	0.70
5:BA:14:ILE:HG13	5:BA:15:LEU:HD22	1.73	0.70
6:BV:46:LEU:HB3	6:BX:42:TYR:HH	1.54	0.70
5:B5:16:ASP:HB2	5:B5:19:ARG:CG	2.21	0.70
9:B6:101:BCL:CMC	9:B7:103:BCL:HBB1	2.21	0.70
5:B7:31:LEU:O	5:B7:35:ILE:HG13	1.90	0.70
14:AS:104:CRT:C39	5:AW:36:HIS:CB	2.70	0.70
5:AY:44:LEU:HD13	6:AZ:43:ARG:HD2	1.73	0.70
6:AZ:27:ALA:O	6:AZ:31:LEU:HG	1.91	0.70
4:BH:176:GLU:O	4:BH:178:GLN:N	2.23	0.70
14:BB:102:CRT:H2M1	5:BD:37:MET:HG2	1.74	0.70
5:BK:11:ILE:HA	14:BP:102:CRT:H82	1.74	0.70
6:BR:30:GLY:O	6:BR:33:VAL:HG12	1.91	0.70
5:BS:5:ASN:HA	5:BS:8:LEU:HG	1.73	0.70
5:BU:12:TRP:HZ2	6:BV:21:PHE:CD2	2.09	0.70
5:B3:13:LEU:HD12	14:B7:102:CRT:H22A	1.71	0.70
5:B3:14:ILE:HD13	6:B6:17:PHE:HE2	1.57	0.70
6:AE:32:VAL:HG21	9:AE:101:BCL:HBA2	1.74	0.70
5:AF:4:MET:CG	6:AJ:23:GLN:HG3	2.21	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:AK:46:TRP:HA	5:AK:49:ASP:OD1	1.92	0.70
6:AV:42:TYR:CD2	6:AV:43:ARG:HG3	2.27	0.70
6:AX:17:PHE:CE1	14:AX:102:CRT:H42	2.27	0.70
5:AY:31:LEU:O	5:AY:35:ILE:HG12	1.91	0.70
6:A0:45:TRP:O	6:A0:46:LEU:HB2	1.89	0.70
1:BC:270:TRP:HA	1:BC:273:ILE:CD1	2.22	0.70
3:BM:215:LEU:C	3:BM:217:ALA:H	1.97	0.70
3:BM:274:VAL:O	3:BM:278:ILE:HG13	1.92	0.70
4:BH:13:GLN:O	4:BH:16:ILE:HG22	1.91	0.70
14:BF:103:CRT:H342	9:BK:102:BCL:CBA	2.21	0.70
5:BW:5:ASN:HB3	6:BX:22:MET:HE3	1.72	0.70
5:B1:10:LYS:C	14:B1:103:CRT:H83	2.17	0.70
5:B7:36:HIS:HB2	14:B7:102:CRT:C2M	2.21	0.70
1:AC:121:ILE:CG2	1:AC:123:THR:HG23	2.22	0.70
1:AC:254:ARG:HH12	3:AM:307:TYR:HE1	1.38	0.70
2:AL:204:LEU:HD11	3:AM:267:ARG:HD2	1.73	0.70
2:AL:237:ALA:HA	2:AL:240:ARG:HG3	1.72	0.70
5:AF:35:ILE:HD13	14:AG:102:CRT:H403	1.74	0.70
1:BC:292:PRO:O	1:BC:296:LYS:HG3	1.91	0.70
6:BJ:31:LEU:O	6:BJ:34:ILE:HG23	1.92	0.70
1:AC:234:GLY:O	1:AC:237:MET:HB2	1.92	0.70
2:AL:204:LEU:HD21	3:AM:267:ARG:HD2	1.73	0.70
3:AM:83:VAL:HG23	3:AM:84:PHE:HD1	1.57	0.70
4:AH:32:ARG:HH21	4:AH:60:ASP:HB2	1.55	0.70
4:AH:235:GLU:HA	4:AH:238:LYS:HB2	1.73	0.70
9:AA:101:BCL:C1D	9:AB:101:BCL:CMD	2.70	0.70
1:BC:95:VAL:O	1:BC:98:THR:HB	1.91	0.70
1:BC:266:ARG:HG3	7:BC:503:HEM:HMD1	1.74	0.70
3:BM:215:LEU:O	3:BM:218:MET:HG3	1.91	0.70
4:BH:65:LYS:N	4:BH:78:ALA:O	2.25	0.70
1:AC:90:PHE:CD1	1:AC:90:PHE:C	2.70	0.70
1:AC:130:MET:SD	7:AC:502:HEM:ND	2.65	0.70
4:AH:39:TYR:HE1	4:AH:48:ARG:HH12	1.38	0.70
6:AX:22:MET:HG3	6:AX:26:TYR:HE2	1.57	0.70
5:A7:35:ILE:HD12	9:A8:101:BCL:O1D	1.90	0.70
1:BC:191:ALA:O	1:BC:192:TYR:HB2	1.91	0.70
6:BJ:30:GLY:O	6:BJ:34:ILE:HG22	1.90	0.70
6:BR:29:PHE:HD1	6:BR:29:PHE:H	1.38	0.70
6:BT:29:PHE:CD1	9:BT:101:BCL:H11	2.26	0.70
3:AM:131:VAL:C	3:AM:133:THR:H	1.99	0.70
6:AB:42:TYR:OH	6:A0:46:LEU:HB3	1.92	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:AI:31:LEU:HB3	9:AJ:101:BCL:CED	2.21	0.70
9:AY:102:BCL:HBC2	9:AZ:101:BCL:HHD	1.72	0.70
6:A2:45:TRP:CE2	9:A2:101:BCL:H2C	2.26	0.70
5:A3:21:LEU:O	5:A3:25:VAL:HG23	1.92	0.70
9:A3:103:BCL:H71	6:A4:28:TRP:CE3	2.27	0.70
6:A0:20:ILE:CD1	14:A0:101:CRT:H133	2.21	0.70
3:BM:300:LYS:HE2	3:BM:300:LYS:HA	1.73	0.70
5:BD:28:GLN:HB3	9:BD:102:BCL:C1	2.22	0.70
6:BN:43:ARG:HB3	5:BO:55:TYR:CE2	2.27	0.70
6:BN:46:LEU:O	5:BO:51:ILE:HG13	1.90	0.70
5:BQ:31:LEU:CG	9:BQ:104:BCL:HED3	2.22	0.70
5:B1:36:HIS:CE1	9:B2:101:BCL:HMD3	2.27	0.70
5:B3:4:MET:O	5:B3:8:LEU:HG	1.92	0.70
1:AC:298:PRO:C	1:AC:300:GLY:H	1.99	0.69
10:AL:302:BPH:H102	9:AL:303:BCL:H193	1.73	0.69
3:AM:114:TRP:HA	3:AM:114:TRP:CE3	2.27	0.69
3:AM:286:LEU:HD13	4:AH:12:ALA:HB1	1.73	0.69
9:AA:101:BCL:C1B	9:A0:102:BCL:HMB2	2.22	0.69
9:BN:101:BCL:HMC3	9:BO:102:BCL:HBB1	1.73	0.69
14:BP:102:CRT:H2M1	5:BQ:33:LEU:O	1.92	0.69
5:BQ:31:LEU:HG	9:BQ:104:BCL:HED3	1.73	0.69
9:BU:102:BCL:C4D	9:BV:101:BCL:HMD3	2.21	0.69
6:BX:27:ALA:O	6:BX:31:LEU:HG	1.91	0.69
5:B3:28:GLN:HG3	9:B3:102:BCL:C1	2.21	0.69
5:B9:35:ILE:HG21	9:B0:102:BCL:ND	2.07	0.69
1:AC:97:VAL:O	1:AC:97:VAL:HG12	1.90	0.69
1:AC:148:THR:HA	1:AC:322:GLN:HB3	1.74	0.69
1:AC:193:ALA:HB3	1:AC:195:LEU:HD12	1.73	0.69
3:AM:27:ASN:HD21	5:AO:19:ARG:NH1	1.90	0.69
9:AK:102:BCL:ND	9:AN:101:BCL:CMD	2.56	0.69
6:AT:21:PHE:CD2	14:AT:102:CRT:H14	2.27	0.69
5:AW:46:TRP:CZ3	9:AW:101:BCL:H2C	2.28	0.69
6:A2:46:LEU:HD22	6:A4:42:TYR:CE2	2.27	0.69
14:A2:102:CRT:H2M2	5:A3:40:LEU:CD1	2.20	0.69
14:A7:102:CRT:C21	14:A7:102:CRT:C23	2.68	0.69
6:A8:45:TRP:O	6:A8:46:LEU:HG	1.91	0.69
9:BI:102:BCL:ND	9:BJ:101:BCL:HMD3	2.07	0.69
5:BO:25:VAL:CG1	9:BO:102:BCL:H52	2.21	0.69
5:BO:25:VAL:HG13	9:BO:102:BCL:H52	1.72	0.69
9:BO:102:BCL:C1D	9:BP:101:BCL:HMD1	2.21	0.69
5:BW:51:ILE:HB	5:BW:52:PRO:CA	2.22	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AC:99:THR:HA	1:AC:103:PRO:HB3	1.74	0.69
2:AL:82:TYR:HA	2:AL:85:ARG:HE	1.56	0.69
5:AA:8:LEU:HD22	6:AE:20:ILE:HG23	1.75	0.69
5:AA:11:ILE:CD1	5:AA:14:ILE:HD11	2.22	0.69
5:AS:36:HIS:O	5:AS:40:LEU:N	2.21	0.69
14:AS:104:CRT:H183	9:AU:102:BCL:H91	1.74	0.69
5:AU:16:ASP:OD2	5:AU:19:ARG:HD3	1.92	0.69
6:AX:32:VAL:O	6:AX:36:HIS:HB2	1.92	0.69
5:A1:36:HIS:CE1	9:A2:101:BCL:HMD3	2.27	0.69
3:BM:205:SER:HB2	9:BM:402:BCL:HMA1	1.74	0.69
6:BN:33:VAL:O	6:BN:37:LEU:HD23	1.92	0.69
9:BO:102:BCL:HAC2	9:BP:101:BCL:CBC	2.22	0.69
5:BQ:43:ASP:HB2	5:BS:47:LEU:CA	2.21	0.69
5:BU:11:ILE:HA	14:BU:103:CRT:C5	2.21	0.69
5:BY:12:TRP:NE1	6:BZ:18:HIS:HA	2.06	0.69
6:B2:20:ILE:CG2	14:B2:102:CRT:H83	2.20	0.69
1:AC:156:HIS:CE1	1:AC:160:PRO:O	2.45	0.69
1:AC:221:SER:O	1:AC:223:PRO:HD3	1.93	0.69
3:AM:164:ARG:HB3	3:AM:165:PRO:HD3	1.73	0.69
5:AY:32:GLY:HA2	9:AZ:101:BCL:O1D	1.91	0.69
5:A3:11:ILE:CA	14:A7:102:CRT:H82	2.23	0.69
5:A7:15:LEU:HB3	5:A7:20:VAL:HG11	1.73	0.69
2:BL:110:ALA:O	2:BL:113:GLU:HB2	1.93	0.69
2:BL:230:GLY:O	3:BM:49:GLY:HA2	1.90	0.69
3:BM:89:HIS:O	3:BM:93:LEU:HG	1.92	0.69
14:BO:103:CRT:C34	9:BS:102:BCL:HBA1	2.22	0.69
14:BS:103:CRT:H342	9:BU:102:BCL:CBA	2.09	0.69
5:BU:16:ASP:OD2	5:BU:19:ARG:HD3	1.92	0.69
14:BV:102:CRT:C2M	5:BW:37:MET:CG	2.70	0.69
5:BW:46:TRP:HH2	9:BW:102:BCL:HBC3	1.56	0.69
1:AC:282:ASN:C	1:AC:283:TYR:HD1	2.01	0.69
2:AL:3:MET:SD	2:AL:8:LYS:HA	2.32	0.69
2:AL:185:ALA:HB2	2:AL:252:TRP:HB3	1.74	0.69
3:AM:98:PRO:HB3	3:AM:107:PRO:HB3	1.74	0.69
4:AH:202:PHE:HB3	4:AH:204:LYS:HZ2	1.56	0.69
6:AG:45:TRP:HA	5:AI:52:PRO:CD	2.21	0.69
5:AK:36:HIS:CE1	9:AN:101:BCL:HMD3	2.28	0.69
5:AU:16:ASP:HB3	5:AU:18:ARG:HH11	1.57	0.69
2:BL:140:LEU:HD12	2:BL:257:ILE:HG21	1.73	0.69
3:BM:166:VAL:HG22	3:BM:171:TRP:CZ3	2.27	0.69
5:B3:12:TRP:NE1	6:B4:18:HIS:HB2	2.07	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:AH:14:ILE:O	4:AH:17:TRP:HB2	1.93	0.69
4:AH:215:LYS:HE3	4:AH:250:ALA:O	1.93	0.69
14:AA:102:CRT:H9	6:AE:17:PHE:HD1	1.58	0.69
5:AI:55:TYR:HD1	5:AI:56:GLN:N	1.91	0.69
5:AO:46:TRP:HD1	5:AO:47:LEU:HD13	1.58	0.69
5:AS:34:LEU:HB2	15:AS:101:PEF:C43	2.22	0.69
5:AU:30:VAL:HG13	5:AU:31:LEU:H	1.57	0.69
5:AY:5:ASN:CG	6:AZ:18:HIS:HD2	2.01	0.69
5:AY:28:GLN:CB	9:AY:102:BCL:C2	2.68	0.69
14:A1:103:CRT:H342	9:A5:102:BCL:C3A	2.22	0.69
14:A2:102:CRT:O2	5:A3:36:HIS:HB3	1.92	0.69
1:BC:157:ARG:NH1	1:BC:318:LEU:HD21	2.07	0.69
3:BM:27:ASN:N	3:BM:27:ASN:HD22	1.90	0.69
4:BH:52:ARG:HB3	4:BH:52:ARG:NH1	2.06	0.69
4:BH:197:ILE:HA	4:BH:200:SER:OG	1.92	0.69
9:BA:101:BCL:C1D	9:BB:101:BCL:CMD	2.71	0.69
5:BQ:19:ARG:HH12	15:BQ:101:PEF:H51	1.57	0.69
5:BU:11:ILE:CG2	14:BU:103:CRT:C8	2.37	0.69
5:BU:32:GLY:HA3	9:BU:102:BCL:O1A	1.93	0.69
6:BZ:32:VAL:HG11	9:BZ:101:BCL:HBA2	1.74	0.69
1:AC:96:ALA:O	1:AC:98:THR:N	2.25	0.69
1:AC:195:LEU:HB3	1:AC:196:PRO:CD	2.22	0.69
2:AL:102:ALA:HB2	10:AL:302:BPH:H112	1.74	0.69
2:AL:237:ALA:O	2:AL:240:ARG:N	2.25	0.69
3:AM:60:SER:CA	3:AM:128:LEU:HD23	2.23	0.69
6:AX:43:ARG:NH1	5:AY:55:TYR:HB3	2.07	0.69
5:A3:56:GLN:NE2	5:A3:56:GLN:H	1.89	0.69
2:BL:3:MET:HG2	2:BL:11:ARG:CZ	2.23	0.69
2:BL:13:ARG:HD2	4:BH:101:VAL:HG22	1.72	0.69
2:BL:120:LEU:O	2:BL:122:ILE:HG23	1.92	0.69
5:BF:11:ILE:CB	14:BF:103:CRT:H82	2.23	0.69
5:BY:16:ASP:HB3	5:BY:18:ARG:NE	2.07	0.69
3:AM:179:ILE:HG12	3:AM:180:PHE:N	2.06	0.69
6:AE:29:PHE:HZ	9:AE:101:BCL:H42	1.57	0.69
5:AF:8:LEU:CA	6:AJ:20:ILE:HD11	2.22	0.69
5:AI:36:HIS:CE1	9:AJ:101:BCL:HMD3	2.27	0.69
6:AZ:18:HIS:NE2	6:AZ:22:MET:HE2	2.08	0.69
5:A1:7:ASN:HB3	5:A1:10:LYS:CE	2.21	0.69
5:A1:52:PRO:CD	5:A1:55:TYR:OH	2.41	0.69
6:A2:45:TRP:CD2	9:A2:101:BCL:H2C	2.28	0.69
1:BC:90:PHE:C	1:BC:90:PHE:CD1	2.69	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BC:243:LEU:H	1:BC:243:LEU:HD12	1.57	0.69
1:BC:249:PHE:CD1	1:BC:250:CYS:SG	2.85	0.69
1:BC:291:LEU:HD23	1:BC:292:PRO:HD2	1.73	0.69
2:BL:223:THR:HA	2:BL:226:ARG:HB3	1.74	0.69
3:BM:171:TRP:CE3	3:BM:171:TRP:HA	2.27	0.69
3:BM:175:VAL:HG22	3:BM:185:TRP:CD2	2.27	0.69
4:BH:135:PRO:HB3	4:BH:171:TRP:NE1	2.07	0.69
4:BH:231:VAL:CG2	4:BH:235:GLU:HG3	2.22	0.69
5:BI:8:LEU:O	5:BI:11:ILE:HG22	1.91	0.69
9:BK:102:BCL:HMD3	6:BN:36:HIS:HD2	1.57	0.69
9:BY:102:BCL:ND	9:BZ:101:BCL:HMD3	2.07	0.69
5:B1:11:ILE:CA	14:B1:103:CRT:H81	2.16	0.69
5:B1:17:PRO:HB3	6:B2:17:PHE:CE2	2.28	0.69
9:B1:102:BCL:CBC	9:B2:101:BCL:CBC	2.71	0.69
6:B2:20:ILE:HG23	14:B2:102:CRT:C7	2.22	0.69
5:B3:46:TRP:CH2	9:B3:102:BCL:HBC3	2.27	0.69
5:B7:8:LEU:O	5:B7:8:LEU:HD13	1.92	0.69
2:AL:42:PHE:HB3	2:AL:101:CYS:HB3	1.75	0.69
2:AL:216:LYS:HD2	2:AL:220:HIS:CD2	2.27	0.69
4:AH:55:VAL:HG13	4:AH:56:VAL:N	2.05	0.69
5:AF:42:THR:HB	5:AI:47:LEU:HD23	1.75	0.69
5:AY:8:LEU:HD22	5:AY:11:ILE:HD11	1.74	0.69
5:A1:44:LEU:CD1	6:A2:43:ARG:HD2	2.21	0.69
9:A1:102:BCL:HMD3	6:A2:36:HIS:CE1	2.27	0.69
2:BL:97:ILE:HA	2:BL:100:ILE:CD1	2.23	0.69
3:BM:199:ASN:HD21	3:BM:283:GLY:HA3	1.58	0.69
4:BH:227:ASN:HD22	4:BH:228:PRO:CD	1.98	0.69
9:BB:101:BCL:HMC3	9:BD:102:BCL:HBB1	1.74	0.69
5:BF:12:TRP:NE1	6:BG:17:PHE:HD1	1.91	0.69
9:BI:102:BCL:C1D	9:BJ:101:BCL:HMD1	2.23	0.69
6:BP:27:ALA:O	6:BP:31:LEU:HG	1.93	0.69
5:BU:2:PHE:HA	5:BU:5:ASN:HD22	1.57	0.69
5:BW:16:ASP:N	5:BW:19:ARG:HE	1.91	0.69
6:B2:32:VAL:HG11	9:B2:101:BCL:CBA	2.23	0.69
5:B3:46:TRP:CZ3	9:B3:102:BCL:CBC	2.76	0.69
6:AB:46:LEU:OXT	6:AE:43:ARG:NH2	2.26	0.69
5:A1:12:TRP:HA	5:A1:12:TRP:CE3	2.28	0.69
5:A3:12:TRP:HE1	6:A4:18:HIS:CB	2.06	0.69
1:BC:121:ILE:CG2	1:BC:123:THR:HG23	2.22	0.69
1:BC:164:TYR:HB3	1:BC:309:THR:HA	1.75	0.69
5:BF:11:ILE:HD12	5:BF:14:ILE:HD11	1.72	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:BF:31:LEU:O	5:BF:35:ILE:HG12	1.92	0.69
5:BY:49:ASP:HB2	5:B1:56:GLN:OE1	1.93	0.69
5:AO:11:ILE:HG12	14:AR:102:CRT:H81	1.75	0.68
5:AQ:48:ASP:O	5:AQ:49:ASP:HB3	1.93	0.68
1:BC:96:ALA:C	1:BC:98:THR:H	2.01	0.68
1:BC:195:LEU:HB3	1:BC:196:PRO:CD	2.23	0.68
5:BA:46:TRP:HB2	6:BB:43:ARG:NH2	2.04	0.68
5:BU:12:TRP:HA	5:BU:12:TRP:CE3	2.28	0.68
5:BW:19:ARG:HG3	5:BW:20:VAL:N	2.08	0.68
2:AL:219:GLU:HG3	4:AH:127:PHE:HB2	1.75	0.68
3:AM:228:ARG:HD3	4:AH:199:PHE:CE1	2.28	0.68
5:AO:5:ASN:HD22	5:AO:8:LEU:HD21	1.59	0.68
5:AQ:44:LEU:HD12	5:AQ:44:LEU:O	1.93	0.68
5:A1:11:ILE:HA	14:A1:103:CRT:H82	1.75	0.68
6:A2:16:GLU:HB2	14:A2:102:CRT:C1M	2.23	0.68
2:BL:170:GLY:HA2	2:BL:176:PHE:CD2	2.29	0.68
2:BL:252:TRP:HA	2:BL:252:TRP:CE3	2.28	0.68
2:BL:253:SER:HB3	9:BL:301:BCL:HED3	1.73	0.68
6:BP:21:PHE:CE1	14:BP:102:CRT:H19	2.28	0.68
9:BW:102:BCL:CMD	6:BX:36:HIS:HD2	2.07	0.68
5:BY:26:ALA:O	5:BY:29:ILE:HG22	1.92	0.68
6:B8:46:LEU:HD22	6:B0:42:TYR:CE2	2.28	0.68
1:AC:242:SER:HA	1:AC:308:MET:SD	2.32	0.68
2:AL:4:LEU:HB2	2:AL:7:GLU:HB2	1.75	0.68
2:AL:71:TRP:N	2:AL:71:TRP:HE3	1.91	0.68
6:AR:21:PHE:HB2	14:AR:102:CRT:C14	2.23	0.68
6:AT:21:PHE:CE2	14:AT:102:CRT:H16	2.28	0.68
9:AU:102:BCL:HBC2	9:AU:102:BCL:CHD	2.24	0.68
5:A5:24:ILE:HG21	14:A7:102:CRT:C21	2.23	0.68
6:BE:38:LEU:O	6:BE:38:LEU:HD23	1.94	0.68
6:BG:46:LEU:HB3	6:BJ:42:TYR:CE2	2.28	0.68
9:BO:102:BCL:H2A	9:BO:102:BCL:O1D	1.92	0.68
5:B3:19:ARG:O	5:B3:23:SER:CB	2.40	0.68
5:B9:36:HIS:CE1	9:B0:102:BCL:HMD3	2.27	0.68
6:B0:36:HIS:HE1	9:B0:102:BCL:C1B	2.06	0.68
1:AC:274:ARG:HG2	1:AC:274:ARG:HH11	1.59	0.68
2:AL:4:LEU:CD2	3:AM:253:ARG:HH21	2.06	0.68
3:AM:34:PRO:HG3	3:AM:50:PRO:N	2.08	0.68
5:AF:19:ARG:NH2	5:AI:18:ARG:NH2	2.32	0.68
5:AS:13:LEU:CB	14:AS:104:CRT:C3	2.71	0.68
6:A8:33:VAL:CG2	9:A8:101:BCL:H142	2.22	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:A9:26:ALA:O	5:A9:29:ILE:HG22	1.93	0.68
1:BC:199:PRO:O	1:BC:203:PHE:HB2	1.94	0.68
3:BM:159:VAL:HA	3:BM:163:ILE:CG2	2.23	0.68
6:BB:36:HIS:HB3	9:BB:101:BCL:H192	1.74	0.68
6:BE:43:ARG:NH1	5:BF:55:TYR:HD2	1.90	0.68
5:BF:33:LEU:H	5:BF:33:LEU:HD12	1.58	0.68
5:BO:25:VAL:HG12	9:BO:102:BCL:H41	1.75	0.68
2:AL:139:VAL:HG23	2:AL:143:VAL:HB	1.75	0.68
2:AL:178:TYR:HB3	2:AL:272:TRP:HD1	1.58	0.68
3:AM:260:VAL:HG22	13:AM:405:MQ8:H112	1.74	0.68
5:AF:38:ILE:CD1	5:AI:37:MET:HE1	2.23	0.68
6:AN:17:PHE:O	6:AN:21:PHE:HB3	1.94	0.68
5:AY:12:TRP:NE1	6:AZ:18:HIS:HA	2.06	0.68
6:AZ:10:THR:HG22	6:AZ:11:ASP:N	2.07	0.68
5:A7:7:ASN:HB2	5:A7:10:LYS:HZ3	1.59	0.68
2:BL:71:TRP:CD1	3:BM:303:MET:HG2	2.29	0.68
3:BM:215:LEU:HA	3:BM:218:MET:SD	2.34	0.68
4:BH:16:ILE:O	4:BH:16:ILE:HD13	1.93	0.68
4:BH:249:TYR:O	4:BH:251:THR:N	2.27	0.68
4:BH:258:LEU:O	5:B5:19:ARG:HD3	1.93	0.68
5:BI:50:ASN:CG	6:BJ:43:ARG:NH2	2.52	0.68
6:B2:13:GLU:O	14:B2:102:CRT:C3	2.34	0.68
5:B3:18:ARG:O	5:B3:22:VAL:HG12	1.94	0.68
5:AA:32:GLY:CA	9:AB:101:BCL:HED2	2.24	0.68
9:AA:101:BCL:HBB3	9:A0:102:BCL:CHC	2.24	0.68
5:AD:36:HIS:CE1	9:AE:101:BCL:CMD	2.77	0.68
5:AF:16:ASP:HB2	5:AF:19:ARG:HB2	1.74	0.68
6:AG:21:PHE:HD1	6:AG:22:MET:CA	2.06	0.68
9:AK:102:BCL:C3D	9:AN:101:BCL:C2D	2.72	0.68
5:AY:28:GLN:O	9:AY:102:BCL:H12	1.93	0.68
5:A3:27:PHE:CE2	5:A5:29:ILE:HG13	2.28	0.68
5:A7:2:PHE:N	5:A7:5:ASN:HB2	2.07	0.68
6:A0:17:PHE:HD1	6:A0:18:HIS:CA	2.07	0.68
1:BC:66:ASP:OD1	1:BC:67:SER:N	2.26	0.68
4:BH:126:THR:HG23	4:BH:130:LEU:O	1.93	0.68
5:BA:9:TYR:HB2	6:BB:18:HIS:HD2	1.59	0.68
5:BI:11:ILE:CA	14:BN:102:CRT:H82	2.22	0.68
5:BO:9:TYR:CD1	6:BP:15:LYS:HB2	2.29	0.68
9:BT:101:BCL:HMA1	9:BU:102:BCL:HHB	1.74	0.68
6:B8:32:VAL:HG11	9:B8:101:BCL:CBA	2.23	0.68
3:AM:175:VAL:CG1	3:AM:176:PRO:HD2	2.24	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:AS:25:VAL:O	5:AS:29:ILE:HG22	1.92	0.68
5:AS:30:VAL:CG2	15:AS:101:PEF:C41	2.32	0.68
5:AU:43:ASP:HB2	5:AW:47:LEU:HB3	1.76	0.68
1:BC:270:TRP:CA	1:BC:273:ILE:HD12	2.22	0.68
1:BC:274:ARG:HA	1:BC:277:ARG:CG	2.23	0.68
5:BA:43:ASP:CA	5:BD:48:ASP:HB3	2.24	0.68
9:BE:101:BCL:CHC	9:BF:102:BCL:HBB3	2.23	0.68
5:BF:13:LEU:HD12	14:BF:103:CRT:C1M	2.19	0.68
6:BV:10:THR:HG22	6:BV:11:ASP:H	1.59	0.68
5:BW:16:ASP:CB	5:BW:19:ARG:HG2	2.23	0.68
2:AL:196:LEU:CD1	3:AM:273:ALA:HB2	2.23	0.68
3:AM:98:PRO:HD2	3:AM:171:TRP:HB3	1.76	0.68
3:AM:114:TRP:HA	3:AM:114:TRP:HE3	1.59	0.68
3:AM:260:VAL:CG2	3:AM:264:SER:OG	2.41	0.68
5:AA:50:ASN:HA	5:AD:59:GLY:C	2.18	0.68
5:A7:47:LEU:HD22	5:A7:47:LEU:H	1.59	0.68
6:BR:44:PRO:O	5:BS:52:PRO:HG3	1.93	0.68
5:BU:45:ASN:OD1	5:BU:47:LEU:HB2	1.94	0.68
14:BU:103:CRT:H31	9:BY:102:BCL:HBA1	1.74	0.68
1:AC:240:SER:HA	7:AC:503:HEM:HMC3	1.74	0.68
5:AA:31:LEU:O	5:AA:35:ILE:HG12	1.93	0.68
14:AA:102:CRT:H23	6:AE:16:GLU:HG3	1.74	0.68
6:AJ:30:GLY:O	6:AJ:34:ILE:HG22	1.92	0.68
6:AP:30:GLY:O	6:AP:34:ILE:CG2	2.42	0.68
5:AY:35:ILE:HG23	5:AY:38:ILE:HD11	1.76	0.68
5:A1:40:LEU:HD12	5:A1:45:ASN:HA	1.76	0.68
5:A7:8:LEU:O	5:A7:8:LEU:HD13	1.94	0.68
9:A8:101:BCL:CMC	9:A9:102:BCL:HBB1	2.23	0.68
5:BQ:10:LYS:HB2	14:BS:103:CRT:H83	1.76	0.68
9:B2:101:BCL:CHB	9:B3:102:BCL:HMB3	2.23	0.68
3:AM:12:GLN:HB2	4:AH:145:ALA:HB2	1.76	0.68
4:AH:22:PHE:CD1	4:AH:22:PHE:C	2.72	0.68
4:AH:133:ILE:CD1	4:AH:171:TRP:HB3	2.23	0.68
5:AF:8:LEU:HD21	6:AJ:24:SER:HG	1.58	0.68
5:AI:18:ARG:HH11	5:AI:18:ARG:HG3	1.58	0.68
5:AO:28:GLN:NE2	14:AP:102:CRT:H25	2.09	0.68
5:AO:50:ASN:CG	5:AO:51:ILE:N	2.52	0.68
9:AQ:102:BCL:CHD	9:AR:101:BCL:HMD1	2.24	0.68
5:AY:30:VAL:HA	5:AY:33:LEU:HG	1.76	0.68
5:A7:25:VAL:HG13	9:A7:103:BCL:C5	2.24	0.68
6:A0:45:TRP:HE1	9:A0:102:BCL:H192	1.56	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:BM:195:ASN:ND2	3:BM:197:TYR:HB2	2.09	0.68
5:BD:10:LYS:HB3	14:BG:102:CRT:H5	1.76	0.68
5:BO:4:MET:HE2	6:BR:23:GLN:CB	2.24	0.68
5:B1:9:TYR:HA	6:B2:18:HIS:CG	2.29	0.68
6:B4:40:TRP:CZ3	6:B4:44:PRO:HA	2.28	0.68
6:B0:32:VAL:HG21	9:B0:102:BCL:CBA	2.21	0.68
2:AL:203:ILE:C	2:AL:205:SER:H	2.00	0.67
4:AH:176:GLU:O	4:AH:178:GLN:N	2.27	0.67
4:AH:221:ASN:HB2	4:AH:242:TYR:OH	1.94	0.67
9:AO:102:BCL:C4D	9:AP:101:BCL:HMD3	2.25	0.67
5:AQ:44:LEU:CD1	5:AQ:46:TRP:HE3	2.06	0.67
6:AR:33:VAL:O	6:AR:37:LEU:HD23	1.94	0.67
5:AY:8:LEU:HD12	6:AZ:22:MET:HE1	1.76	0.67
5:A1:31:LEU:O	5:A1:35:ILE:HG12	1.95	0.67
5:A3:11:ILE:N	14:A7:102:CRT:H82	2.10	0.67
1:BC:210:ILE:O	1:BC:210:ILE:HG22	1.93	0.67
5:BK:12:TRP:HB2	6:BN:14:ALA:HB1	1.74	0.67
5:BW:46:TRP:HA	5:BW:49:ASP:OD1	1.94	0.67
1:AC:95:VAL:O	1:AC:98:THR:HB	1.95	0.67
2:AL:44:LEU:C	2:AL:46:GLY:H	2.02	0.67
2:AL:89:LEU:H	2:AL:89:LEU:HD12	1.58	0.67
3:AM:106:ILE:HG12	5:AO:42:THR:HG21	1.75	0.67
6:AB:27:ALA:O	6:AB:31:LEU:HG	1.95	0.67
5:AF:35:ILE:O	5:AF:38:ILE:HG22	1.93	0.67
9:AJ:101:BCL:HMC1	9:AK:102:BCL:HBB1	1.75	0.67
6:AN:10:THR:HG22	6:AN:11:ASP:H	1.59	0.67
3:BM:114:TRP:HA	3:BM:114:TRP:CE3	2.27	0.67
3:BM:175:VAL:CG1	3:BM:176:PRO:HD2	2.24	0.67
3:BM:179:ILE:HG12	3:BM:180:PHE:H	1.59	0.67
4:BH:23:PHE:C	4:BH:25:GLY:H	2.02	0.67
4:BH:213:ALA:O	4:BH:246:GLY:HA3	1.95	0.67
5:BI:29:ILE:HA	9:BI:102:BCL:H11	1.76	0.67
5:BO:31:LEU:O	5:BO:35:ILE:HG12	1.93	0.67
14:BU:103:CRT:C2M	5:BY:37:MET:CA	2.71	0.67
6:BV:20:ILE:HG23	14:BV:102:CRT:C9	2.24	0.67
1:AC:53:ILE:HG12	1:AC:319:TYR:CZ	2.29	0.67
3:AM:5:GLN:O	3:AM:7:ILE:HG12	1.94	0.67
5:AI:7:ASN:HD22	6:AN:20:ILE:CG1	2.08	0.67
6:AN:17:PHE:CE1	14:AN:102:CRT:H9	2.29	0.67
5:AO:38:ILE:HD13	14:AP:102:CRT:C40	2.24	0.67
5:AY:43:ASP:OD1	5:AY:44:LEU:HD23	1.93	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:A5:10:LYS:CB	14:A5:103:CRT:H5	2.25	0.67
5:A7:2:PHE:CD1	5:A7:3:THR:N	2.62	0.67
1:BC:85:LEU:HD22	1:BC:89:GLU:HG2	1.75	0.67
3:BM:159:VAL:HG12	3:BM:284:ILE:HG22	1.74	0.67
4:BH:6:THR:HB	5:BF:41:SER:CB	2.14	0.67
5:BA:46:TRP:CB	6:BB:43:ARG:NH2	2.54	0.67
5:BD:36:HIS:CE1	9:BE:101:BCL:CMD	2.77	0.67
5:BF:5:ASN:OD1	6:BG:22:MET:HE3	1.95	0.67
5:BW:51:ILE:HB	5:BW:52:PRO:HA	1.76	0.67
5:BY:43:ASP:HB2	5:B1:47:LEU:HG	1.76	0.67
2:AL:97:ILE:O	2:AL:100:ILE:HB	1.95	0.67
5:AS:10:LYS:CB	14:AS:104:CRT:C1M	2.71	0.67
6:AV:7:THR:CG2	14:AX:102:CRT:C1M	2.72	0.67
5:A1:28:GLN:HB3	9:A1:102:BCL:C2	2.25	0.67
5:A3:28:GLN:HG3	9:A3:103:BCL:H12	1.77	0.67
6:A4:37:LEU:HD23	6:A4:38:LEU:N	2.10	0.67
1:BC:280:ASN:CB	1:BC:304:ARG:HD2	2.24	0.67
2:BL:46:GLY:HA3	10:BL:302:BPH:H9C3	1.77	0.67
2:BL:178:TYR:CD2	2:BL:269:PRO:HG3	2.27	0.67
9:BL:301:BCL:HBB3	9:BM:401:BCL:HMD2	1.77	0.67
4:BH:65:LYS:O	4:BH:77:VAL:HA	1.94	0.67
5:BW:8:LEU:O	5:BW:11:ILE:HG13	1.95	0.67
6:B2:13:GLU:CB	14:B2:102:CRT:H33	2.18	0.67
1:AC:94:MET:SD	7:AC:501:HEM:NC	2.68	0.67
1:AC:97:VAL:HG13	7:AC:502:HEM:CMB	2.25	0.67
1:AC:156:HIS:O	1:AC:159:ASN:N	2.28	0.67
2:AL:37:VAL:HG23	2:AL:38:VAL:H	1.58	0.67
2:AL:93:GLY:O	2:AL:97:ILE:HG13	1.93	0.67
2:AL:184:LEU:CD2	2:AL:252:TRP:HE1	2.07	0.67
2:AL:185:ALA:N	2:AL:252:TRP:HD1	1.92	0.67
5:AF:19:ARG:NH1	5:AI:18:ARG:HH21	1.90	0.67
5:AF:49:ASP:HB2	5:AI:56:GLN:HB2	1.77	0.67
5:AU:12:TRP:HE1	6:AV:18:HIS:CA	2.00	0.67
14:AW:102:CRT:H181	9:AY:102:BCL:H92	1.77	0.67
2:BL:189:PHE:HE2	2:BL:253:SER:HG	1.41	0.67
3:BM:136:ARG:HA	3:BM:136:ARG:HH11	1.59	0.67
4:BH:78:ALA:HA	4:BH:79:PRO:C	2.18	0.67
4:BH:119:ARG:HG3	4:BH:237:ASP:OD2	1.95	0.67
5:BO:18:ARG:HB2	5:BO:18:ARG:HH11	1.57	0.67
5:BO:46:TRP:HD1	5:BO:47:LEU:HD13	1.58	0.67
6:BR:29:PHE:O	6:BR:33:VAL:HB	1.95	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:BW:5:ASN:HA	5:BW:8:LEU:HD12	1.77	0.67
5:B1:13:LEU:HD12	14:B1:103:CRT:H22A	1.70	0.67
5:B9:34:LEU:O	5:B9:38:ILE:HG12	1.95	0.67
1:AC:325:LYS:HA	1:AC:331:TYR:OH	1.94	0.67
3:AM:148:TRP:CE3	3:AM:148:TRP:HA	2.30	0.67
4:AH:235:GLU:HA	4:AH:238:LYS:CG	2.25	0.67
5:AI:30:VAL:HA	5:AI:33:LEU:HD23	1.76	0.67
2:BL:68:TYR:HA	2:BL:73:ILE:HD11	1.77	0.67
2:BL:145:PRO:HG3	17:BL:403:HOH:O	1.94	0.67
2:BL:279:PRO:CG	5:BY:41:SER:HB2	2.25	0.67
3:BM:12:GLN:HB2	4:BH:145:ALA:HB2	1.75	0.67
3:BM:253:ARG:NH1	3:BM:258:PHE:HA	2.09	0.67
4:BH:186:VAL:HG12	4:BH:187:ALA:H	1.58	0.67
5:BA:2:PHE:HA	5:BA:5:ASN:HD21	1.60	0.67
5:BA:19:ARG:HH12	5:BD:22:VAL:HG11	1.60	0.67
6:BN:37:LEU:O	6:BN:41:LEU:HG	1.94	0.67
5:BQ:44:LEU:HD22	6:BR:43:ARG:CD	2.24	0.67
5:BW:26:ALA:C	5:BW:29:ILE:HG22	2.17	0.67
5:BY:40:LEU:HD12	5:BY:45:ASN:HA	1.76	0.67
5:B5:31:LEU:O	5:B5:35:ILE:HG12	1.94	0.67
3:AM:265:ILE:CG2	3:AM:266:HIS:H	2.08	0.67
9:AK:102:BCL:HMD2	9:AN:101:BCL:C1D	2.24	0.67
9:AN:101:BCL:CHC	9:AO:102:BCL:HBB3	2.24	0.67
6:AR:46:LEU:CB	6:AT:42:TYR:OH	2.43	0.67
5:AY:8:LEU:HB3	6:AZ:18:HIS:NE2	2.09	0.67
6:A0:32:VAL:CG1	6:A0:33:VAL:N	2.57	0.67
1:BC:36:ARG:HG3	1:BC:36:ARG:O	1.94	0.67
2:BL:16:THR:OG1	4:BH:257:PRO:HB3	1.95	0.67
2:BL:137:TYR:O	2:BL:141:VAL:HG12	1.95	0.67
5:BA:47:LEU:CD1	5:B9:43:ASP:HB2	2.24	0.67
9:BA:101:BCL:HBB1	9:B0:102:BCL:HMC1	1.75	0.67
6:BN:21:PHE:CD2	14:BN:102:CRT:H14	2.29	0.67
6:BX:46:LEU:HD22	6:BZ:42:TYR:CZ	2.30	0.67
1:AC:105:GLU:N	1:AC:105:GLU:OE2	2.28	0.67
2:AL:164:ASP:HA	2:AL:167:SER:OG	1.94	0.67
9:AL:301:BCL:CBA	9:AL:303:BCL:HBC1	2.23	0.67
3:AM:186:THR:HA	9:AM:402:BCL:HMD2	1.77	0.67
5:AY:10:LYS:O	5:AY:13:LEU:HG	1.93	0.67
5:A3:56:GLN:NE2	5:A3:56:GLN:N	2.42	0.67
5:A5:4:MET:HE3	6:A8:24:SER:CB	2.09	0.67
6:A0:9:LEU:HB3	6:A0:13:GLU:HG3	1.77	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:BL:190:PHE:CE1	3:BM:209:LEU:HD21	2.30	0.67
9:BL:301:BCL:H2C	9:BM:402:BCL:H2C	1.75	0.67
5:BA:46:TRP:HA	6:BB:43:ARG:HH12	1.60	0.67
5:BF:49:ASP:OD2	5:BI:56:GLN:HG2	1.95	0.67
9:B8:101:BCL:H2A	9:B8:101:BCL:O1D	1.95	0.67
1:AC:203:PHE:HE1	1:AC:210:ILE:HG12	1.59	0.67
2:AL:163:LEU:HD23	3:AM:197:TYR:HB3	1.77	0.67
2:AL:182:HIS:CA	2:AL:256:CYS:SG	2.83	0.67
9:AA:101:BCL:HED3	9:A0:102:BCL:H92	1.77	0.67
5:AD:7:ASN:O	5:AD:10:LYS:HD3	1.94	0.67
9:AF:102:BCL:OBD	6:AG:32:VAL:HG13	1.94	0.67
6:AJ:45:TRP:O	6:AJ:46:LEU:HB2	1.93	0.67
6:AN:32:VAL:HG12	9:AN:101:BCL:H141	1.77	0.67
9:AN:101:BCL:C4B	9:AO:102:BCL:HBB3	2.25	0.67
6:AP:38:LEU:HD23	6:AP:39:ALA:N	2.09	0.67
6:AP:45:TRP:O	6:AP:46:LEU:HG	1.95	0.67
14:AS:104:CRT:C7	6:AV:20:ILE:CD1	2.65	0.67
5:AU:9:TYR:HB2	6:AV:15:LYS:CD	2.25	0.67
6:AZ:46:LEU:C	5:A1:51:ILE:O	2.38	0.67
6:A8:27:ALA:O	6:A8:31:LEU:CG	2.29	0.67
1:BC:259:TRP:CD1	1:BC:259:TRP:H	2.13	0.67
2:BL:89:LEU:HA	2:BL:94:LEU:N	2.09	0.67
5:BD:29:ILE:O	5:BD:33:LEU:HB2	1.95	0.67
6:BR:34:ILE:C	6:BR:34:ILE:HD12	2.20	0.67
5:BU:49:ASP:CG	5:BU:50:ASN:N	2.50	0.67
1:AC:213:THR:OG1	1:AC:257:ASN:HB2	1.94	0.67
3:AM:176:PRO:HD3	3:AM:185:TRP:CD1	2.30	0.67
9:AF:102:BCL:C1D	9:AG:101:BCL:HMD1	2.24	0.67
5:AS:46:TRP:HZ3	9:AS:103:BCL:CBC	2.07	0.67
14:AS:104:CRT:H391	5:AW:36:HIS:CB	2.25	0.67
9:AY:102:BCL:H2A	9:AY:102:BCL:O1D	1.95	0.67
9:A1:102:BCL:C1	9:A1:102:BCL:O1A	2.43	0.67
5:A3:13:LEU:CB	14:A7:102:CRT:H1M1	2.19	0.67
5:A5:4:MET:HG2	6:A8:24:SER:CB	2.24	0.67
2:BL:71:TRP:HD1	3:BM:303:MET:HG2	1.59	0.67
2:BL:150:ALA:CB	2:BL:153:HIS:HB2	2.25	0.67
3:BM:278:ILE:O	3:BM:282:ILE:HG13	1.93	0.67
4:BH:153:GLY:H	4:BH:167:VAL:CG2	2.08	0.67
4:BH:241:ALA:O	4:BH:244:ALA:HB3	1.94	0.67
6:BE:10:THR:HG22	6:BE:11:ASP:N	2.10	0.67
5:BI:10:LYS:HB3	14:BN:102:CRT:H5	1.76	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AC:94:MET:SD	7:AC:501:HEM:FE	1.87	0.66
2:AL:8:LYS:HD2	4:AH:42:ASP:OD2	1.95	0.66
2:AL:106:PHE:O	2:AL:110:ALA:HB2	1.95	0.66
5:AI:43:ASP:OD2	5:AK:47:LEU:HD12	1.95	0.66
6:A2:16:GLU:CB	14:A2:102:CRT:C1M	2.73	0.66
1:BC:157:ARG:HH12	1:BC:318:LEU:HG	1.59	0.66
6:BB:40:TRP:HA	6:BB:40:TRP:CE3	2.29	0.66
9:BE:101:BCL:CBB	9:BE:101:BCL:HMB1	2.25	0.66
6:BG:32:VAL:HG11	9:BG:101:BCL:CBA	2.21	0.66
6:BN:38:LEU:HD23	6:BN:38:LEU:O	1.94	0.66
5:BW:18:ARG:O	5:BW:22:VAL:CG2	2.42	0.66
6:BX:17:PHE:O	6:BX:20:ILE:HG22	1.95	0.66
5:BY:32:GLY:N	9:BZ:101:BCL:HED2	2.10	0.66
5:BY:36:HIS:CE1	9:BZ:101:BCL:HMD3	2.29	0.66
5:B1:16:ASP:HB3	5:B1:18:ARG:HE	1.59	0.66
9:B1:102:BCL:ND	9:B2:101:BCL:HMD3	2.10	0.66
5:B3:13:LEU:HD21	6:B4:10:THR:O	1.95	0.66
5:B7:44:LEU:HD21	5:B7:46:TRP:CE3	2.29	0.66
1:AC:157:ARG:HE	1:AC:312:GLN:NE2	1.93	0.66
2:AL:150:ALA:CB	2:AL:153:HIS:HB2	2.25	0.66
4:AH:69:LEU:HD23	4:AH:70:PRO:HD2	1.77	0.66
9:AI:102:BCL:CHD	9:AJ:101:BCL:HMD1	2.25	0.66
2:BL:186:ILE:HD13	9:BM:401:BCL:CMD	2.25	0.66
3:BM:70:ILE:HG21	3:BM:118:ALA:HB2	1.77	0.66
3:BM:104:LEU:CD1	3:BM:169:GLY:HA2	2.22	0.66
4:BH:235:GLU:HA	4:BH:238:LYS:CG	2.25	0.66
5:BU:12:TRP:CE3	6:BV:17:PHE:HE2	2.13	0.66
5:BU:14:ILE:CG1	14:BU:103:CRT:H31A	2.25	0.66
5:BU:42:THR:HB	5:BW:48:ASP:CG	2.21	0.66
5:B7:43:ASP:OD1	5:B7:44:LEU:HD12	1.95	0.66
2:AL:38:VAL:HG23	2:AL:39:GLY:H	1.59	0.66
2:AL:43:THR:O	2:AL:47:VAL:HG23	1.94	0.66
2:AL:155:PHE:HB2	2:AL:156:PRO:HD2	1.75	0.66
2:AL:188:PHE:O	2:AL:190:PHE:N	2.27	0.66
4:AH:123:CYS:H	4:AH:232:THR:HG22	1.60	0.66
9:AD:102:BCL:C1D	9:AE:101:BCL:CMD	2.74	0.66
9:AF:102:BCL:HAC2	9:AG:101:BCL:CBC	2.25	0.66
6:AJ:10:THR:HG22	6:AJ:11:ASP:H	1.61	0.66
6:AX:46:LEU:HD13	6:AZ:42:TYR:OH	1.96	0.66
5:AY:35:ILE:O	5:AY:38:ILE:HG13	1.96	0.66
5:A1:11:ILE:CG2	14:A1:103:CRT:H81	2.24	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:A5:8:LEU:HD21	14:A5:103:CRT:H133	1.77	0.66
5:A9:35:ILE:O	5:A9:39:VAL:HG23	1.96	0.66
2:BL:103:ALA:O	2:BL:107:ILE:HG13	1.95	0.66
3:BM:286:LEU:HD13	4:BH:12:ALA:HB1	1.77	0.66
5:BK:11:ILE:CG1	14:BP:102:CRT:H81	2.25	0.66
6:BN:10:THR:HB	6:BN:13:GLU:OE2	1.94	0.66
6:BP:20:ILE:HD13	14:BP:102:CRT:C6	2.25	0.66
6:BR:33:VAL:O	6:BR:37:LEU:HD23	1.96	0.66
6:BT:33:VAL:O	6:BT:37:LEU:HG	1.95	0.66
5:B7:44:LEU:CD2	5:B7:46:TRP:CE3	2.67	0.66
6:B0:31:LEU:O	6:B0:34:ILE:HG23	1.95	0.66
1:AC:190:VAL:C	1:AC:192:TYR:H	2.02	0.66
4:AH:55:VAL:CA	5:AA:19:ARG:HH12	2.05	0.66
4:AH:197:ILE:HA	4:AH:200:SER:OG	1.94	0.66
4:AH:231:VAL:CG2	4:AH:235:GLU:HG3	2.22	0.66
5:AD:36:HIS:O	5:AD:40:LEU:HB2	1.95	0.66
6:AT:9:LEU:HD22	6:AT:13:GLU:HG3	1.77	0.66
6:A4:40:TRP:HZ3	6:A4:45:TRP:N	1.93	0.66
5:A7:43:ASP:OD1	5:A7:44:LEU:HD12	1.94	0.66
2:BL:148:MET:HE3	2:BL:153:HIS:CE1	2.31	0.66
4:BH:136:MET:HE3	4:BH:172:VAL:HG11	1.77	0.66
4:BH:196:PRO:HG2	4:BH:199:PHE:HB2	1.76	0.66
5:BA:51:ILE:HB	5:BA:52:PRO:HA	1.78	0.66
9:BU:102:BCL:C1D	9:BV:101:BCL:HMD1	2.25	0.66
1:AC:227:LYS:O	1:AC:230:GLU:HB3	1.96	0.66
1:AC:249:PHE:CZ	1:AC:265:LYS:HG2	2.29	0.66
1:AC:259:TRP:C	1:AC:261:GLN:N	2.52	0.66
2:AL:113:GLU:HB3	2:AL:127:PRO:HG3	1.78	0.66
2:AL:126:VAL:HB	2:AL:127:PRO:CD	2.23	0.66
4:AH:159:LEU:HD12	4:AH:159:LEU:C	2.21	0.66
6:AG:34:ILE:HD13	6:AG:35:ALA:N	2.11	0.66
5:AI:18:ARG:O	5:AI:22:VAL:HG12	1.95	0.66
5:AQ:43:ASP:N	5:AS:47:LEU:HB3	2.11	0.66
14:AS:104:CRT:H393	5:AW:36:HIS:HB2	1.77	0.66
6:AT:34:ILE:O	6:AT:34:ILE:HD13	1.95	0.66
6:AX:33:VAL:HG22	6:AX:37:LEU:HD23	1.78	0.66
5:A3:46:TRP:CE3	9:A3:103:BCL:H2C	2.30	0.66
9:A6:101:BCL:C1C	9:A7:103:BCL:HBB3	2.25	0.66
6:A8:29:PHE:CZ	9:A8:101:BCL:H61	2.30	0.66
5:A9:40:LEU:HD12	5:A9:45:ASN:HA	1.78	0.66
2:BL:10:TYR:OH	3:BM:246:GLU:HG2	1.95	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:BL:148:MET:HE2	2:BL:262:PRO:HD3	1.77	0.66
2:BL:195:ALA:HA	2:BL:198:MET:HE2	1.77	0.66
3:BM:179:ILE:H	3:BM:179:ILE:CD1	1.92	0.66
6:BB:17:PHE:HB2	14:BB:102:CRT:H21A	1.77	0.66
5:BI:50:ASN:OD1	6:BJ:43:ARG:NH2	2.29	0.66
6:BX:34:ILE:C	6:BX:34:ILE:HD13	2.21	0.66
5:BY:49:ASP:HB2	5:B1:56:GLN:HE22	1.61	0.66
9:B2:101:BCL:HMA1	9:B3:102:BCL:HHB	1.76	0.66
6:B8:27:ALA:O	6:B8:31:LEU:CG	2.29	0.66
9:B9:102:BCL:CMD	6:B0:36:HIS:HD2	2.08	0.66
6:B0:45:TRP:O	6:B0:46:LEU:HB2	1.94	0.66
3:AM:134:TYR:HA	3:AM:144:GLN:NE2	2.10	0.66
3:AM:164:ARG:HA	3:AM:167:MET:HB3	1.76	0.66
6:AN:36:HIS:ND1	9:AN:101:BCL:H162	2.09	0.66
5:AQ:27:PHE:HE2	5:AS:29:ILE:CD1	2.08	0.66
14:AS:104:CRT:H2M1	5:AW:37:MET:CB	2.26	0.66
6:AV:32:VAL:HG11	9:AV:102:BCL:HBA2	1.78	0.66
9:AW:101:BCL:C1D	9:AX:101:BCL:CMD	2.74	0.66
5:A1:12:TRP:HA	5:A1:12:TRP:HE3	1.61	0.66
5:A7:29:ILE:HG23	5:A7:30:VAL:H	1.60	0.66
1:BC:57:GLN:NE2	1:BC:58:PRO:HD2	2.10	0.66
2:BL:188:PHE:O	2:BL:190:PHE:N	2.29	0.66
5:BQ:43:ASP:OD1	5:BQ:44:LEU:CG	2.42	0.66
6:BV:20:ILE:CG2	14:BV:102:CRT:C6	2.74	0.66
6:B2:21:PHE:HB2	14:B2:102:CRT:C11	2.21	0.66
5:B3:14:ILE:CD1	6:B6:17:PHE:HE2	2.09	0.66
6:B8:33:VAL:HG23	9:B8:101:BCL:H143	1.77	0.66
1:AC:192:TYR:O	2:AL:269:PRO:HB3	1.95	0.66
3:AM:291:VAL:HG11	3:AM:297:TRP:HB2	1.77	0.66
4:AH:65:LYS:O	4:AH:77:VAL:HA	1.95	0.66
5:AA:43:ASP:HA	5:AD:48:ASP:CB	2.21	0.66
14:AB:102:CRT:H31A	5:A9:10:LYS:C	2.20	0.66
6:AN:20:ILE:HD12	6:AN:20:ILE:N	2.09	0.66
5:AS:46:TRP:CZ3	9:AS:103:BCL:CBC	2.78	0.66
5:A1:5:ASN:OD1	5:A1:8:LEU:HD12	1.94	0.66
5:A9:36:HIS:CE1	9:A0:102:BCL:HMD3	2.30	0.66
6:A0:40:TRP:HZ3	6:A0:45:TRP:N	1.94	0.66
2:BL:206:VAL:HG12	3:BM:142:MET:CE	2.26	0.66
2:BL:243:LEU:O	2:BL:247:LEU:HB2	1.96	0.66
2:BL:259:ILE:HA	2:BL:263:PHE:HB2	1.78	0.66
3:BM:268:TRP:CD2	4:BH:30:LEU:HD13	2.30	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:BM:290:VAL:HG11	4:BH:12:ALA:HB2	1.77	0.66
5:BA:17:PRO:O	5:BA:21:LEU:CB	2.43	0.66
6:BB:40:TRP:HA	6:BB:40:TRP:HE3	1.60	0.66
6:BJ:38:LEU:O	6:BJ:38:LEU:HD23	1.96	0.66
5:BK:16:ASP:HB2	5:BK:19:ARG:HG2	1.77	0.66
5:BK:33:LEU:HD12	5:BK:34:LEU:N	2.11	0.66
6:B2:46:LEU:HB2	5:B3:52:PRO:HD2	1.77	0.66
5:B5:32:GLY:CA	9:B5:102:BCL:O1A	2.43	0.66
6:B0:36:HIS:CE1	9:B0:102:BCL:C1B	2.79	0.66
1:AC:36:ARG:HB3	2:AL:79:ASP:CG	2.20	0.66
2:AL:89:LEU:H	2:AL:89:LEU:CD1	2.09	0.66
2:AL:171:TYR:C	2:AL:173:PHE:H	2.03	0.66
6:AB:44:PRO:HG2	5:AD:52:PRO:CB	2.26	0.66
9:AD:102:BCL:ND	9:AE:101:BCL:HMD3	2.11	0.66
5:AF:8:LEU:HD23	6:AJ:20:ILE:HD11	1.77	0.66
5:AI:39:VAL:HG11	9:AI:102:BCL:CBC	2.26	0.66
6:AJ:31:LEU:O	6:AJ:34:ILE:HG23	1.96	0.66
1:BC:99:THR:HA	1:BC:103:PRO:HB3	1.76	0.66
1:BC:187:SER:O	1:BC:189:THR:N	2.28	0.66
1:BC:263:THR:HG22	3:BM:311:VAL:CB	2.26	0.66
3:BM:63:PHE:CD2	3:BM:124:LEU:HB2	2.31	0.66
3:BM:114:TRP:HA	3:BM:114:TRP:HE3	1.59	0.66
4:BH:47:GLU:HG3	5:BA:19:ARG:CB	2.26	0.66
6:BG:46:LEU:HB3	6:BJ:42:TYR:CZ	2.31	0.66
5:BW:19:ARG:HG3	5:BW:20:VAL:H	1.60	0.66
6:B4:37:LEU:HD23	6:B4:38:LEU:N	2.11	0.66
6:B4:46:LEU:HB2	5:B5:52:PRO:HD3	1.76	0.66
5:B5:32:GLY:HA3	9:B5:102:BCL:O1A	1.96	0.66
2:AL:273:ASN:HD22	2:AL:276:LEU:HD23	1.60	0.66
6:AE:18:HIS:CE1	6:AE:22:MET:HE2	2.31	0.66
5:A1:15:LEU:HA	5:A3:18:ARG:NH1	2.07	0.66
6:A6:29:PHE:CE1	9:A6:101:BCL:C1	2.79	0.66
6:A6:29:PHE:O	6:A6:33:VAL:HG23	1.96	0.66
9:A0:102:BCL:H192	9:A0:102:BCL:OBB	1.96	0.66
4:BH:106:PRO:HA	4:BH:109:SER:OG	1.96	0.66
5:BD:35:ILE:HG13	9:BE:101:BCL:O1D	1.96	0.66
9:BG:101:BCL:CMB	9:BI:102:BCL:C1B	2.74	0.66
5:BO:9:TYR:CD1	5:BO:9:TYR:C	2.74	0.66
14:B2:102:CRT:H2M3	5:B3:40:LEU:HD11	1.78	0.66
2:AL:82:TYR:CB	2:AL:85:ARG:HE	2.09	0.66
5:AS:8:LEU:HB2	6:AT:18:HIS:NE2	2.11	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:AU:30:VAL:HG13	5:AU:31:LEU:N	2.11	0.66
5:A7:9:TYR:HA	6:A8:18:HIS:ND1	2.10	0.66
5:A9:46:TRP:CH2	9:A9:102:BCL:HBC3	2.31	0.66
6:A0:38:LEU:O	6:A0:38:LEU:HD23	1.96	0.66
2:BL:70:LEU:HB3	2:BL:159:ILE:HG12	1.78	0.66
4:BH:69:LEU:CB	4:BH:70:PRO:HD2	2.25	0.66
9:BF:102:BCL:C1D	9:BG:101:BCL:HMD1	2.26	0.66
5:BK:32:GLY:HA3	9:BK:102:BCL:O1A	1.96	0.66
9:BO:102:BCL:C4D	9:BP:101:BCL:HMD3	2.26	0.66
9:B1:102:BCL:HBC1	9:B2:101:BCL:CBC	2.25	0.66
5:B3:10:LYS:HB3	14:B7:102:CRT:H5	1.76	0.66
5:B3:27:PHE:HE1	5:B3:31:LEU:HD22	1.61	0.66
1:AC:182:GLY:O	1:AC:196:PRO:HA	1.96	0.65
1:AC:253:THR:HG21	2:AL:168:ASN:HA	1.78	0.65
2:AL:167:SER:HA	9:AL:301:BCL:HBC1	1.78	0.65
2:AL:190:PHE:CE1	3:AM:209:LEU:HD21	2.31	0.65
3:AM:104:LEU:HD21	3:AM:169:GLY:CA	2.26	0.65
3:AM:152:ALA:CB	3:AM:274:VAL:HG13	2.26	0.65
6:AB:45:TRP:O	6:AB:46:LEU:HB2	1.96	0.65
5:AD:44:LEU:HD12	5:AD:44:LEU:O	1.95	0.65
5:AF:27:PHE:HE1	5:AI:29:ILE:HD11	1.55	0.65
6:A6:20:ILE:O	6:A6:20:ILE:HD13	1.96	0.65
6:A0:31:LEU:O	6:A0:34:ILE:HG23	1.95	0.65
1:BC:104:LYS:HB3	1:BC:105:GLU:OE2	1.97	0.65
2:BL:253:SER:OG	9:BL:301:BCL:HAA2	1.96	0.65
3:BM:27:ASN:HD21	5:BO:19:ARG:HH11	1.44	0.65
3:BM:84:PHE:HA	5:BW:37:MET:HE1	1.77	0.65
3:BM:148:TRP:HA	3:BM:148:TRP:CE3	2.32	0.65
3:BM:165:PRO:HB3	3:BM:174:ALA:HB2	1.78	0.65
3:BM:215:LEU:HD12	3:BM:218:MET:SD	2.37	0.65
5:BA:47:LEU:HD12	5:B9:43:ASP:CB	2.24	0.65
5:BU:14:ILE:HG13	14:BU:103:CRT:H33	0.74	0.65
9:BU:102:BCL:CHD	9:BU:102:BCL:HBC2	2.26	0.65
5:B1:10:LYS:C	14:B1:103:CRT:C8	2.69	0.65
14:B5:103:CRT:H342	9:B9:102:BCL:CBA	2.18	0.65
14:B0:101:CRT:H32	9:B0:102:BCL:CMA	2.27	0.65
1:AC:126:VAL:O	1:AC:127:SER:C	2.39	0.65
2:AL:253:SER:C	9:AL:301:BCL:HED3	2.21	0.65
3:AM:268:TRP:CD2	4:AH:30:LEU:HD13	2.30	0.65
4:AH:119:ARG:HG3	4:AH:237:ASP:OD2	1.96	0.65
4:AH:259:LEU:HD21	5:A5:19:ARG:C	2.21	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:AA:50:ASN:ND2	5:AA:51:ILE:HG12	2.10	0.65
5:AY:26:ALA:O	5:AY:29:ILE:HG22	1.96	0.65
6:AZ:33:VAL:HG22	6:AZ:37:LEU:HD12	1.77	0.65
5:A3:31:LEU:O	5:A3:35:ILE:HG12	1.96	0.65
5:A3:36:HIS:NE2	9:A3:104:BCL:HMD3	2.11	0.65
5:A5:36:HIS:CE1	9:A6:101:BCL:HMD3	2.31	0.65
6:A0:21:PHE:C	6:A0:21:PHE:CD1	2.75	0.65
1:BC:36:ARG:HB3	2:BL:79:ASP:OD1	1.97	0.65
1:BC:179:LYS:N	1:BC:180:PRO:HD3	2.11	0.65
3:BM:159:VAL:HG13	3:BM:285:LEU:HD13	1.79	0.65
4:BH:52:ARG:O	4:BH:54:LYS:HG2	1.95	0.65
6:BE:45:TRP:HA	5:BF:52:PRO:HG2	1.78	0.65
9:BE:101:BCL:HMB1	9:BE:101:BCL:HBB3	1.77	0.65
6:BX:18:HIS:NE2	6:BX:22:MET:HE2	2.11	0.65
6:BZ:46:LEU:HB3	5:B1:52:PRO:HD3	1.76	0.65
5:B1:10:LYS:HD2	6:B4:20:ILE:CB	2.26	0.65
2:AL:211:LYS:HD3	2:AL:212:GLY:N	2.11	0.65
5:AF:44:LEU:HB3	5:AI:55:TYR:OH	1.96	0.65
5:AI:44:LEU:HD23	5:AK:55:TYR:HE2	1.61	0.65
6:AV:33:VAL:O	6:AV:37:LEU:HD23	1.96	0.65
5:A1:50:ASN:CG	5:A1:51:ILE:N	2.55	0.65
1:BC:85:LEU:HD11	1:BC:329:GLY:HA3	1.77	0.65
3:BM:171:TRP:HA	3:BM:171:TRP:HE3	1.60	0.65
9:BN:101:BCL:H172	6:BP:38:LEU:HD22	1.78	0.65
6:B6:29:PHE:O	6:B6:33:VAL:HG23	1.97	0.65
5:B7:37:MET:N	14:B7:102:CRT:H2M1	2.10	0.65
1:AC:41:GLU:OE1	2:AL:153:HIS:CD2	2.50	0.65
3:AM:300:LYS:O	4:AH:8:TYR:HB2	1.97	0.65
14:AA:102:CRT:H35	5:AD:31:LEU:HD11	1.76	0.65
6:AB:20:ILE:CD1	14:AB:102:CRT:H10	2.21	0.65
6:AJ:34:ILE:HD13	6:AJ:35:ALA:N	2.12	0.65
6:AR:42:TYR:CD2	6:AR:43:ARG:HG3	2.32	0.65
5:AS:34:LEU:HB2	15:AS:101:PEF:C44	2.26	0.65
5:AW:4:MET:HG3	6:AZ:24:SER:HA	1.77	0.65
5:AW:27:PHE:CE2	5:AY:29:ILE:HG12	2.32	0.65
6:AX:43:ARG:HH12	5:AY:55:TYR:HB3	1.61	0.65
6:A2:21:PHE:HD1	14:A2:102:CRT:C14	2.09	0.65
6:A6:20:ILE:HG21	14:A7:102:CRT:C6	2.27	0.65
5:A7:18:ARG:O	5:A7:22:VAL:HG12	1.97	0.65
5:A7:31:LEU:O	5:A7:35:ILE:HG13	1.96	0.65
9:A8:101:BCL:C1C	9:A9:102:BCL:CBB	2.74	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:BL:237:ALA:HA	2:BL:240:ARG:HD2	1.79	0.65
9:BM:401:BCL:HBC1	9:BM:402:BCL:CBD	2.26	0.65
6:BE:18:HIS:CE1	6:BE:22:MET:HE2	2.31	0.65
5:BO:38:ILE:HG13	5:BO:39:VAL:H	1.62	0.65
6:BP:34:ILE:HD12	6:BP:34:ILE:C	2.22	0.65
5:BS:50:ASN:CG	5:BS:51:ILE:H	2.05	0.65
6:BZ:33:VAL:HG22	6:BZ:37:LEU:HD12	1.78	0.65
6:B4:40:TRP:HZ3	6:B4:45:TRP:N	1.93	0.65
9:B8:101:BCL:NC	9:B9:102:BCL:HBB3	2.11	0.65
6:B0:38:LEU:HD23	6:B0:38:LEU:O	1.96	0.65
1:AC:153:TYR:CE1	1:AC:157:ARG:HA	2.32	0.65
1:AC:210:ILE:HB	7:AC:503:HEM:O2D	1.96	0.65
2:AL:182:HIS:CE1	2:AL:186:ILE:HD11	2.32	0.65
2:AL:194:LEU:O	2:AL:198:MET:HG3	1.97	0.65
5:AA:37:MET:O	5:AA:41:SER:HB2	1.96	0.65
5:AO:10:LYS:O	5:AO:13:LEU:HD22	1.95	0.65
5:A7:36:HIS:CE1	9:A8:101:BCL:HMD3	2.31	0.65
5:A9:16:ASP:HB3	5:A9:19:ARG:HB2	1.77	0.65
2:BL:22:LEU:HB2	5:B7:19:ARG:HB2	1.77	0.65
2:BL:160:LEU:O	2:BL:163:LEU:HB2	1.96	0.65
2:BL:192:ASN:HD22	2:BL:193:CYS:N	1.94	0.65
2:BL:211:LYS:HD3	2:BL:212:GLY:H	1.62	0.65
3:BM:314:VAL:HG12	3:BM:315:ASN:H	1.62	0.65
4:BH:171:TRP:HE1	4:BH:183:GLU:HG3	1.61	0.65
14:B1:103:CRT:H392	9:B5:102:BCL:HMB2	1.77	0.65
14:B2:102:CRT:C34	9:B3:102:BCL:HBA1	2.23	0.65
5:B7:36:HIS:CB	14:B7:102:CRT:C39	2.74	0.65
1:AC:66:ASP:OD1	1:AC:67:SER:N	2.29	0.65
1:AC:91:THR:O	1:AC:95:VAL:HG23	1.96	0.65
2:AL:270:GLU:O	2:AL:273:ASN:N	2.29	0.65
5:AA:27:PHE:HE1	5:AD:29:ILE:CD1	2.10	0.65
5:AI:44:LEU:O	5:AI:44:LEU:HD12	1.96	0.65
9:AO:102:BCL:HAC2	9:AP:101:BCL:CBC	2.26	0.65
5:AW:36:HIS:CE1	9:AX:101:BCL:CMD	2.79	0.65
6:AZ:22:MET:O	6:AZ:25:MET:HB3	1.97	0.65
9:A1:102:BCL:CHD	9:A2:101:BCL:HMD1	2.27	0.65
6:A0:40:TRP:HZ3	6:A0:45:TRP:H	1.44	0.65
1:BC:269:ALA:O	1:BC:273:ILE:HG13	1.95	0.65
5:BD:22:VAL:HA	5:BD:25:VAL:HG23	1.76	0.65
6:BV:10:THR:HG22	6:BV:11:ASP:N	2.11	0.65
5:BW:24:ILE:CD1	9:BY:102:BCL:H18	2.23	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AC:142:LYS:O	1:AC:146:ALA:HA	1.96	0.65
2:AL:137:TYR:HD1	2:AL:138:LEU:HD12	1.60	0.65
3:AM:159:VAL:HG21	3:AM:281:GLY:HA3	1.76	0.65
6:AE:22:MET:HG3	6:AE:26:TYR:HE1	1.62	0.65
5:AK:5:ASN:HA	5:AK:8:LEU:HD12	1.79	0.65
5:AU:42:THR:HB	5:AW:48:ASP:CG	2.22	0.65
1:BC:169:ASP:OD1	1:BC:170:PRO:HD2	1.97	0.65
2:BL:140:LEU:O	2:BL:141:VAL:HB	1.96	0.65
3:BM:159:VAL:HG21	3:BM:281:GLY:HA3	1.78	0.65
3:BM:164:ARG:HD2	3:BM:284:ILE:O	1.97	0.65
5:BF:45:ASN:CB	5:BF:49:ASP:HB3	2.20	0.65
5:BI:33:LEU:HD12	5:BI:34:LEU:N	2.11	0.65
5:BO:50:ASN:ND2	6:BP:43:ARG:NH2	2.44	0.65
5:BO:50:ASN:OD1	6:BP:43:ARG:NH2	2.29	0.65
5:BQ:43:ASP:HA	5:BS:47:LEU:C	2.21	0.65
5:BW:21:LEU:O	5:BW:25:VAL:HG23	1.97	0.65
5:B7:35:ILE:HD12	9:B8:101:BCL:O1D	1.96	0.65
2:AL:202:LEU:HD13	2:AL:224:PHE:CD2	2.31	0.65
3:AM:215:LEU:C	3:AM:217:ALA:N	2.53	0.65
3:AM:297:TRP:HZ3	3:AM:303:MET:SD	2.20	0.65
6:AN:45:TRP:CD1	6:AN:46:LEU:HG	2.32	0.65
6:AR:45:TRP:CZ3	9:AR:101:BCL:HAC2	2.32	0.65
9:AY:102:BCL:CMD	6:AZ:36:HIS:CD2	2.80	0.65
1:BC:53:ILE:HA	1:BC:319:TYR:CE1	2.31	0.65
9:BQ:103:BCL:HMB1	9:BQ:103:BCL:HBB3	1.78	0.65
5:BU:12:TRP:CZ2	6:BV:21:PHE:CD2	2.85	0.65
5:BU:19:ARG:NE	5:BW:18:ARG:HH22	1.95	0.65
6:BX:40:TRP:O	6:BX:44:PRO:HG3	1.96	0.65
6:B0:17:PHE:CB	14:B0:101:CRT:H6	2.18	0.65
1:AC:167:VAL:HG21	1:AC:298:PRO:HD2	1.78	0.65
2:AL:204:LEU:HD11	3:AM:267:ARG:CD	2.27	0.65
3:AM:75:MET:O	3:AM:78:SER:HB3	1.96	0.65
3:AM:264:SER:O	3:AM:267:ARG:CB	2.45	0.65
4:AH:47:GLU:HG3	5:AA:19:ARG:CA	2.27	0.65
4:AH:135:PRO:HB3	4:AH:171:TRP:NE1	2.12	0.65
6:AB:23:GLN:HG3	5:A9:4:MET:CE	2.18	0.65
6:AG:30:GLY:O	6:AG:34:ILE:HG22	1.97	0.65
5:AY:38:ILE:HD12	5:AY:39:VAL:HG23	1.79	0.65
9:A3:103:BCL:C1D	9:A3:104:BCL:CMD	2.75	0.65
2:BL:43:THR:O	2:BL:47:VAL:HG23	1.96	0.65
3:BM:83:VAL:O	3:BM:87:LEU:HD23	1.97	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:BI:102:BCL:CHD	9:BJ:101:BCL:HMD1	2.27	0.65
6:BP:20:ILE:HG21	14:BP:102:CRT:H6	1.79	0.65
14:BP:102:CRT:O2	5:BQ:33:LEU:HD12	1.96	0.65
5:BU:12:TRP:CG	6:BV:17:PHE:CD2	2.85	0.65
5:BU:28:GLN:HB3	9:BU:102:BCL:C2	2.27	0.65
1:AC:173:LYS:HE3	5:AU:42:THR:HG22	1.78	0.65
1:AC:255:ALA:HB1	1:AC:258:ASP:HB3	1.79	0.65
1:AC:283:TYR:O	1:AC:286:PRO:HD2	1.97	0.65
4:AH:153:GLY:H	4:AH:167:VAL:CG2	2.10	0.65
6:AE:33:VAL:O	6:AE:37:LEU:HD23	1.96	0.65
5:AO:26:ALA:O	5:AO:29:ILE:HG22	1.97	0.65
5:AS:10:LYS:O	5:AS:13:LEU:HB2	1.96	0.65
5:A1:7:ASN:CB	5:A1:10:LYS:HE3	2.24	0.65
5:A1:12:TRP:CE2	6:A2:17:PHE:CE2	2.85	0.65
9:A1:102:BCL:H71	6:A2:28:TRP:CE3	2.32	0.65
3:BM:290:VAL:HG12	3:BM:291:VAL:N	2.12	0.65
14:BG:102:CRT:H2M3	5:BI:36:HIS:CB	2.27	0.65
6:BT:30:GLY:O	6:BT:34:ILE:HG23	1.97	0.65
5:B1:10:LYS:HD2	6:B4:20:ILE:CG1	2.27	0.65
5:B9:26:ALA:O	5:B9:29:ILE:HG22	1.95	0.65
6:B0:9:LEU:HB3	6:B0:13:GLU:HG3	1.77	0.65
2:AL:133:ALA:HB2	10:AL:302:BPH:HAC2	1.79	0.64
6:AG:45:TRP:HD1	6:AG:46:LEU:N	1.95	0.64
9:AJ:101:BCL:C1B	9:AK:102:BCL:CMB	2.75	0.64
9:AN:101:BCL:HMB3	9:AO:102:BCL:CHB	2.27	0.64
5:AY:51:ILE:HB	5:AY:52:PRO:CA	2.26	0.64
5:A1:10:LYS:HD2	6:A4:20:ILE:HG13	1.78	0.64
2:BL:29:PRO:HG2	3:BM:257:GLY:HA2	1.79	0.64
4:BH:106:PRO:HA	4:BH:109:SER:CB	2.28	0.64
5:BF:35:ILE:O	5:BF:38:ILE:HG22	1.96	0.64
5:BU:43:ASP:HB2	5:BW:47:LEU:HB3	1.80	0.64
14:BV:102:CRT:H2M3	5:BW:37:MET:CB	2.27	0.64
5:BW:4:MET:C	5:BW:6:ALA:H	2.05	0.64
5:BY:46:TRP:CZ3	9:BY:102:BCL:HBC3	2.32	0.64
5:BY:51:ILE:HB	5:BY:52:PRO:HA	1.79	0.64
5:B5:16:ASP:HB2	5:B5:19:ARG:HH21	1.63	0.64
2:AL:70:LEU:HB3	2:AL:159:ILE:HG12	1.79	0.64
2:AL:143:VAL:HG12	2:AL:258:LEU:HD11	1.77	0.64
3:AM:103:GLY:O	3:AM:104:LEU:HD13	1.96	0.64
6:AE:10:THR:HG22	6:AE:11:ASP:N	2.11	0.64
6:AG:28:TRP:HE1	6:AG:32:VAL:HG21	1.59	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:AI:102:BCL:HBC2	9:AJ:101:BCL:HHD	1.79	0.64
5:AO:9:TYR:CE1	6:AP:15:LYS:HD2	2.32	0.64
6:AR:29:PHE:O	6:AR:33:VAL:HB	1.97	0.64
1:BC:192:TYR:O	2:BL:269:PRO:HB3	1.97	0.64
2:BL:241:LEU:O	2:BL:244:PHE:HB3	1.97	0.64
3:BM:229:PHE:N	3:BM:229:PHE:CD1	2.64	0.64
4:BH:182:LEU:HD13	4:BH:195:LEU:HD23	1.79	0.64
5:BF:12:TRP:HE1	6:BG:17:PHE:HD1	1.43	0.64
6:BN:10:THR:HG22	6:BN:11:ASP:N	2.07	0.64
5:BO:9:TYR:CE1	6:BP:15:LYS:HB2	2.32	0.64
6:BP:17:PHE:HA	6:BP:20:ILE:HG22	1.78	0.64
5:BY:11:ILE:HD13	9:B1:102:BCL:H151	1.79	0.64
6:B0:32:VAL:CG1	6:B0:33:VAL:N	2.59	0.64
6:B0:40:TRP:HB2	9:B0:102:BCL:H191	1.79	0.64
1:AC:98:THR:O	1:AC:103:PRO:HD3	1.98	0.64
1:AC:100:TRP:HB3	1:AC:152:CYS:HB2	1.80	0.64
1:AC:112:VAL:HG12	1:AC:113:PRO:HD2	1.79	0.64
3:AM:104:LEU:HD21	3:AM:169:GLY:HA2	1.78	0.64
9:AI:102:BCL:C1D	9:AJ:101:BCL:CMD	2.76	0.64
5:AY:13:LEU:CD2	6:AZ:14:ALA:HB2	2.27	0.64
9:AY:102:BCL:ND	9:AZ:101:BCL:HMD3	2.13	0.64
6:A8:33:VAL:CG2	9:A8:101:BCL:H143	2.27	0.64
6:A0:36:HIS:CE1	9:A0:102:BCL:C1B	2.80	0.64
2:BL:182:HIS:CE1	2:BL:186:ILE:HD11	2.32	0.64
3:BM:200:PRO:HA	3:BM:203:MET:HG2	1.79	0.64
3:BM:240:HIS:NE2	4:BH:69:LEU:HD21	2.12	0.64
3:BM:244:ALA:C	3:BM:246:GLU:H	2.03	0.64
5:BD:46:TRP:CH2	9:BD:102:BCL:CBC	2.79	0.64
6:B6:20:ILE:O	6:B6:20:ILE:HD13	1.96	0.64
1:AC:130:MET:SD	7:AC:502:HEM:NC	2.70	0.64
1:AC:259:TRP:CD1	1:AC:259:TRP:H	2.15	0.64
2:AL:186:ILE:HD13	9:AM:401:BCL:HMD2	1.80	0.64
2:AL:196:LEU:HD13	2:AL:196:LEU:C	2.22	0.64
4:AH:48:ARG:HD3	15:AH:301:PEF:H42	1.78	0.64
4:AH:88:ASN:ND2	4:AH:109:SER:HB2	2.12	0.64
9:AT:101:BCL:HMA1	9:AU:102:BCL:HHB	1.78	0.64
5:AU:38:ILE:HD11	5:AW:40:LEU:CD2	2.28	0.64
5:A1:44:LEU:HG	5:A1:44:LEU:O	1.97	0.64
6:A4:13:GLU:H	6:A4:13:GLU:CD	2.06	0.64
1:BC:182:GLY:O	1:BC:196:PRO:HA	1.96	0.64
2:BL:252:TRP:O	2:BL:253:SER:C	2.41	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:BB:20:ILE:HG21	14:BB:102:CRT:H83	1.78	0.64
9:BF:102:BCL:HBA2	9:BG:101:BCL:OBD	1.98	0.64
5:B1:12:TRP:CH2	5:B1:20:VAL:HG11	2.32	0.64
5:B5:31:LEU:HA	5:B5:34:LEU:HB3	1.79	0.64
6:B0:17:PHE:HD1	14:B0:101:CRT:H9	1.53	0.64
1:AC:157:ARG:HH12	1:AC:318:LEU:CG	2.11	0.64
3:AM:34:PRO:HA	3:AM:48:ILE:O	1.98	0.64
9:AA:101:BCL:CHD	9:AB:101:BCL:HMD1	2.27	0.64
5:AD:5:ASN:HD22	6:AE:22:MET:HG2	1.62	0.64
6:AG:17:PHE:CD1	6:AG:17:PHE:C	2.75	0.64
5:AS:10:LYS:HG2	14:AS:104:CRT:H1M1	1.79	0.64
5:AS:30:VAL:HG22	15:AS:101:PEF:C40	2.25	0.64
6:A8:43:ARG:HH21	5:A9:55:TYR:HB2	1.62	0.64
6:A0:18:HIS:O	6:A0:22:MET:HB2	1.98	0.64
1:BC:211:ARG:C	1:BC:212:ILE:HD13	2.22	0.64
3:BM:34:PRO:HG2	3:BM:50:PRO:HD3	1.80	0.64
3:BM:199:ASN:HA	3:BM:294:TRP:CE3	2.32	0.64
5:BD:31:LEU:O	5:BD:35:ILE:HG12	1.98	0.64
5:BW:45:ASN:O	5:BW:49:ASP:HB3	1.98	0.64
5:BW:51:ILE:HB	5:BW:52:PRO:C	2.23	0.64
5:B5:20:VAL:HA	5:B5:23:SER:HB3	1.79	0.64
5:B5:31:LEU:HD12	5:B5:34:LEU:HD23	1.78	0.64
6:B8:34:ILE:HG12	6:B8:37:LEU:HD23	1.77	0.64
6:B0:11:ASP:O	6:B0:15:LYS:HG3	1.98	0.64
6:B0:34:ILE:HD13	6:B0:35:ALA:N	2.13	0.64
1:AC:264:PRO:HG2	1:AC:265:LYS:HD2	1.77	0.64
2:AL:160:LEU:C	2:AL:160:LEU:HD12	2.23	0.64
2:AL:211:LYS:HD3	2:AL:212:GLY:H	1.62	0.64
3:AM:103:GLY:C	3:AM:104:LEU:HD22	2.22	0.64
5:AD:35:ILE:HG13	9:AE:101:BCL:O1D	1.98	0.64
6:AG:12:ASP:O	6:AG:16:GLU:HG3	1.97	0.64
5:AW:5:ASN:CA	5:AW:8:LEU:HD12	2.26	0.64
5:AW:50:ASN:CG	5:AW:51:ILE:H	2.05	0.64
9:AZ:101:BCL:CHB	9:A1:102:BCL:HMB2	2.28	0.64
6:A8:34:ILE:HG12	6:A8:37:LEU:HD23	1.77	0.64
1:BC:20:LEU:HG	2:BL:271:TRP:HE1	1.62	0.64
3:BM:40:LEU:CD1	3:BM:48:ILE:HD11	2.27	0.64
4:BH:47:GLU:HG3	5:BA:19:ARG:HG3	1.78	0.64
6:BN:30:GLY:O	6:BN:34:ILE:HG22	1.98	0.64
5:BQ:35:ILE:HA	5:BQ:38:ILE:CG2	2.28	0.64
5:BU:43:ASP:HB2	5:BW:47:LEU:HD22	1.78	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:BY:9:TYR:CD1	6:BZ:15:LYS:HG2	2.33	0.64
9:B6:101:BCL:C1C	9:B7:103:BCL:HBB3	2.27	0.64
5:B7:10:LYS:HD3	6:B0:20:ILE:HD12	1.79	0.64
6:B0:21:PHE:CD1	6:B0:21:PHE:C	2.75	0.64
1:AC:137:ALA:HA	1:AC:141:TRP:HD1	1.62	0.64
1:AC:301:ASP:HB2	1:AC:302:PRO:CD	2.28	0.64
2:AL:82:TYR:HB3	2:AL:85:ARG:HE	1.63	0.64
2:AL:112:ARG:NH2	3:AM:255:THR:HA	2.12	0.64
2:AL:178:TYR:HB3	2:AL:272:TRP:CD1	2.33	0.64
2:AL:196:LEU:HD12	3:AM:273:ALA:HB2	1.78	0.64
3:AM:268:TRP:CD1	3:AM:268:TRP:H	2.14	0.64
5:AA:8:LEU:HB3	6:AE:20:ILE:CG2	2.27	0.64
5:AD:9:TYR:HB2	6:AE:15:LYS:HA	1.79	0.64
5:AK:12:TRP:HB2	6:AN:14:ALA:HB1	1.80	0.64
9:AK:102:BCL:HBD	9:AN:101:BCL:OBD	1.98	0.64
6:AR:38:LEU:HD12	6:AR:38:LEU:O	1.98	0.64
5:AS:20:VAL:HB	9:AU:102:BCL:H203	1.78	0.64
5:AW:18:ARG:HG2	5:AW:18:ARG:HH11	1.62	0.64
5:AW:19:ARG:HH12	5:AY:22:VAL:CG2	2.10	0.64
5:A5:5:ASN:HA	5:A5:8:LEU:HD12	1.80	0.64
5:A5:31:LEU:HD21	14:A7:102:CRT:H32	1.79	0.64
5:A5:43:ASP:CA	5:A7:47:LEU:HB3	2.27	0.64
6:A0:11:ASP:O	6:A0:15:LYS:HG3	1.98	0.64
2:BL:188:PHE:CB	2:BL:249:ALA:HB2	2.22	0.64
3:BM:7:ILE:HG22	3:BM:8:PHE:N	2.12	0.64
3:BM:208:PHE:HB3	3:BM:276:THR:OG1	1.97	0.64
4:BH:18:ALA:O	4:BH:21:LEU:HB3	1.97	0.64
14:BA:102:CRT:H392	9:BF:102:BCL:HMB2	1.78	0.64
6:BR:10:THR:HG22	6:BR:11:ASP:N	2.12	0.64
6:BX:45:TRP:CZ3	9:BX:101:BCL:HAC2	2.33	0.64
5:B1:16:ASP:HB2	5:B1:19:ARG:HD3	1.79	0.64
2:AL:148:MET:HE2	2:AL:262:PRO:HD3	1.80	0.64
2:AL:252:TRP:CE3	2:AL:252:TRP:HA	2.33	0.64
3:AM:35:ILE:CD1	15:AM:409:PEF:H321	2.28	0.64
3:AM:97:PRO:HB2	3:AM:171:TRP:O	1.97	0.64
3:AM:133:THR:HG22	3:AM:134:TYR:N	2.12	0.64
4:AH:5:ILE:HG23	4:AH:6:THR:N	2.10	0.64
5:AK:44:LEU:HD22	5:AK:44:LEU:O	1.97	0.64
5:AO:44:LEU:HD12	5:AO:46:TRP:H	1.63	0.64
6:AP:44:PRO:HG2	5:AQ:52:PRO:CB	2.28	0.64
5:AU:14:ILE:HD12	14:AX:102:CRT:H82	1.79	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:AW:21:LEU:HD22	14:AX:102:CRT:H132	1.78	0.64
9:AY:102:BCL:HMD1	6:AZ:36:HIS:CD2	2.33	0.64
6:A2:16:GLU:CB	14:A2:102:CRT:H1M1	2.28	0.64
3:BM:265:ILE:CG2	3:BM:266:HIS:N	2.60	0.64
6:BJ:10:THR:HB	6:BJ:13:GLU:CD	2.23	0.64
5:BK:47:LEU:H	5:BK:47:LEU:HD22	1.61	0.64
5:BO:21:LEU:O	5:BO:25:VAL:HG23	1.98	0.64
5:BY:50:ASN:HA	5:B1:60:LYS:O	1.98	0.64
5:B1:11:ILE:HG23	5:B1:15:LEU:HD12	1.80	0.64
14:B5:103:CRT:C32	5:B7:31:LEU:HD21	2.28	0.64
1:AC:110:CYS:HA	1:AC:123:THR:OG1	1.97	0.64
1:AC:187:SER:O	1:AC:189:THR:N	2.30	0.64
2:AL:12:VAL:CG2	2:AL:13:ARG:H	2.10	0.64
3:AM:229:PHE:HB3	3:AM:243:THR:HG23	1.80	0.64
5:AD:49:ASP:HB2	5:AF:56:GLN:CB	2.28	0.64
5:AO:11:ILE:HG12	14:AR:102:CRT:H10	1.78	0.64
5:AS:31:LEU:HD21	14:AT:102:CRT:H32	1.79	0.64
14:AS:104:CRT:C2M	5:AW:37:MET:N	2.61	0.64
5:AU:22:VAL:HG13	5:AU:23:SER:N	2.12	0.64
5:AW:2:PHE:CA	5:AW:5:ASN:HD22	2.01	0.64
6:A8:29:PHE:CE1	9:A8:101:BCL:C2	2.81	0.64
2:BL:35:PHE:CZ	2:BL:111:LEU:HD12	2.33	0.64
2:BL:38:VAL:HG23	2:BL:39:GLY:N	2.12	0.64
3:BM:159:VAL:HA	3:BM:163:ILE:HG22	1.80	0.64
3:BM:247:ARG:NH2	4:BH:244:ALA:HB1	2.13	0.64
4:BH:259:LEU:HD21	5:B5:19:ARG:C	2.23	0.64
5:BF:44:LEU:HD22	6:BG:43:ARG:HD2	1.79	0.64
6:BP:21:PHE:HB2	14:BP:102:CRT:H11	1.80	0.64
5:BS:9:TYR:HA	6:BT:18:HIS:CG	2.33	0.64
6:BX:17:PHE:HA	6:BX:20:ILE:HG22	1.79	0.64
6:B0:32:VAL:HG12	6:B0:33:VAL:H	1.63	0.64
1:AC:272:ALA:O	1:AC:276:VAL:HG12	1.98	0.64
4:AH:35:LYS:O	4:AH:36:ARG:C	2.41	0.64
14:AB:102:CRT:C8	5:A9:10:LYS:HB2	2.28	0.64
5:AO:7:ASN:H	5:AO:7:ASN:HD22	1.43	0.64
6:AP:36:HIS:CE1	9:AP:101:BCL:NA	2.65	0.64
5:AU:44:LEU:HD22	6:AV:43:ARG:CD	2.28	0.64
6:AZ:12:ASP:HA	6:AZ:15:LYS:HD2	1.80	0.64
5:A1:10:LYS:HD2	6:A4:20:ILE:CB	2.27	0.64
5:A3:29:ILE:O	5:A3:33:LEU:HG	1.96	0.64
5:A5:46:TRP:CZ2	9:A5:102:BCL:HHC	2.33	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:BL:112:ARG:NH2	3:BM:255:THR:HA	2.13	0.64
2:BL:230:GLY:CA	3:BM:51:ILE:HB	2.26	0.64
4:BH:77:VAL:O	4:BH:80:ARG:HD3	1.98	0.64
4:BH:159:LEU:C	4:BH:159:LEU:HD12	2.23	0.64
5:BF:50:ASN:ND2	6:BG:43:ARG:HH22	1.95	0.64
6:BG:25:MET:CE	9:BI:102:BCL:H203	2.28	0.64
9:BI:102:BCL:HMD3	6:BJ:36:HIS:CD2	2.33	0.64
9:BS:102:BCL:HBC1	9:BT:101:BCL:HBC3	1.79	0.64
14:BW:103:CRT:H9	6:BZ:17:PHE:HE1	1.61	0.64
5:B7:29:ILE:HB	9:B7:103:BCL:H43	1.80	0.64
2:AL:120:LEU:CD2	3:AM:250:LEU:HD23	2.27	0.63
2:AL:177:HIS:CB	3:AM:183:LEU:HD22	2.26	0.63
3:AM:137:ALA:O	3:AM:142:MET:HB2	1.97	0.63
3:AM:160:LEU:CD2	3:AM:284:ILE:HG21	2.28	0.63
3:AM:208:PHE:CE1	3:AM:275:LEU:HB3	2.33	0.63
3:AM:254:TRP:CD1	3:AM:254:TRP:N	2.64	0.63
4:AH:57:GLY:HA2	15:AH:301:PEF:O1P	1.98	0.63
5:AF:11:ILE:O	5:AF:14:ILE:HG12	1.98	0.63
6:A8:34:ILE:O	6:A8:34:ILE:HD13	1.98	0.63
6:A0:34:ILE:HD13	6:A0:35:ALA:N	2.13	0.63
1:BC:164:TYR:CB	1:BC:309:THR:HA	2.28	0.63
10:BL:302:BPH:H162	9:BL:303:BCL:HMB1	1.80	0.63
3:BM:2:PRO:HB3	4:BH:201:ARG:HH12	1.63	0.63
3:BM:34:PRO:HG3	3:BM:50:PRO:CD	2.28	0.63
3:BM:56:THR:HG21	3:BM:131:VAL:CG1	2.28	0.63
3:BM:109:LEU:CD2	5:BS:45:ASN:HD21	2.10	0.63
9:BQ:103:BCL:HMB1	9:BQ:103:BCL:CBB	2.28	0.63
5:BS:29:ILE:HG23	5:BS:30:VAL:N	2.12	0.63
5:BU:12:TRP:CD2	6:BV:17:PHE:CD2	2.85	0.63
9:BV:101:BCL:HMA1	9:BW:102:BCL:HMA1	1.80	0.63
14:BV:102:CRT:C34	9:BW:102:BCL:HBA1	2.23	0.63
5:B1:12:TRP:HH2	5:B1:20:VAL:HG11	1.62	0.63
5:B1:12:TRP:HH2	9:B3:102:BCL:H202	1.62	0.63
6:B2:32:VAL:CG1	9:B2:101:BCL:HBA2	2.27	0.63
5:B9:5:ASN:HA	5:B9:8:LEU:HD12	1.79	0.63
5:B9:12:TRP:HE1	6:B0:18:HIS:CA	2.11	0.63
9:B9:102:BCL:HMD2	6:B0:36:HIS:HD2	1.63	0.63
1:AC:123:THR:O	1:AC:126:VAL:HG22	1.97	0.63
1:AC:243:LEU:HD12	1:AC:243:LEU:N	2.12	0.63
1:AC:259:TRP:C	1:AC:261:GLN:H	2.05	0.63
2:AL:270:GLU:O	2:AL:271:TRP:C	2.41	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:AM:98:PRO:HG3	3:AM:107:PRO:HG3	1.80	0.63
5:AA:14:ILE:HG13	5:AA:15:LEU:CD2	2.29	0.63
5:AK:9:TYR:HB2	6:AN:15:LYS:HA	1.80	0.63
5:AK:48:ASP:HB3	5:AK:56:GLN:NE2	2.12	0.63
14:AS:104:CRT:H182	9:AU:102:BCL:C9	2.27	0.63
5:AW:7:ASN:H	5:AW:7:ASN:ND2	1.95	0.63
5:A3:32:GLY:N	9:A3:104:BCL:HED2	2.13	0.63
1:BC:270:TRP:O	1:BC:274:ARG:HD2	1.98	0.63
3:BM:131:VAL:C	3:BM:133:THR:H	2.05	0.63
4:BH:48:ARG:O	4:BH:53:VAL:HG23	1.98	0.63
9:BA:101:BCL:H202	5:B9:24:ILE:HD13	1.79	0.63
5:BW:16:ASP:O	5:BW:19:ARG:CG	2.46	0.63
6:BZ:11:ASP:O	6:BZ:15:LYS:HG3	1.98	0.63
9:B1:102:BCL:HBC2	9:B2:101:BCL:HBC3	1.79	0.63
6:B2:46:LEU:HB2	5:B3:52:PRO:CD	2.27	0.63
5:B3:9:TYR:CE1	5:B3:10:LYS:HD2	2.33	0.63
1:AC:17:SER:HB3	3:AM:91:PHE:CZ	2.33	0.63
1:AC:291:LEU:CD2	1:AC:292:PRO:HD2	2.23	0.63
2:AL:53:GLY:HA3	2:AL:75:ILE:HD11	1.79	0.63
2:AL:112:ARG:HH21	3:AM:255:THR:HA	1.64	0.63
4:AH:164:ALA:HB2	4:AH:216:ALA:HB1	1.80	0.63
5:AQ:51:ILE:HG23	5:AQ:52:PRO:CA	2.28	0.63
5:AU:19:ARG:NE	5:AW:18:ARG:HH22	1.97	0.63
9:AY:102:BCL:HBC1	9:AZ:101:BCL:HBC3	1.80	0.63
6:A2:21:PHE:CE1	14:A2:102:CRT:H16	2.32	0.63
5:A3:31:LEU:HD21	9:A3:104:BCL:HMA3	1.78	0.63
6:A4:45:TRP:O	6:A4:46:LEU:HG	1.99	0.63
5:A5:4:MET:O	5:A5:8:LEU:HG	1.99	0.63
5:A9:16:ASP:O	5:A9:20:VAL:HG22	1.98	0.63
3:BM:59:LEU:HG	3:BM:128:LEU:HD21	1.81	0.63
3:BM:138:GLU:C	3:BM:140:LEU:H	2.05	0.63
3:BM:260:VAL:HG23	3:BM:261:THR:H	1.61	0.63
3:BM:279:THR:HA	3:BM:282:ILE:CD1	2.27	0.63
4:BH:54:LYS:HE2	5:BD:23:SER:CA	2.28	0.63
5:BD:11:ILE:HG23	5:BD:12:TRP:CE3	2.33	0.63
9:BE:101:BCL:C1B	9:BF:102:BCL:CMB	2.76	0.63
5:BS:42:THR:HG22	5:BS:43:ASP:N	2.13	0.63
9:BV:101:BCL:CMA	9:BW:102:BCL:HMA1	2.29	0.63
14:BV:102:CRT:H2M1	5:BW:33:LEU:O	1.98	0.63
5:B1:16:ASP:CB	5:B1:18:ARG:HE	2.11	0.63
5:B9:40:LEU:CD1	5:B9:47:LEU:HD23	2.29	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AC:24:GLU:OE1	1:AC:45:ASN:HB2	1.98	0.63
2:AL:240:ARG:HH21	3:AM:6:ASN:C	2.05	0.63
2:AL:241:LEU:O	2:AL:244:PHE:HB3	1.99	0.63
4:AH:251:THR:HG22	4:AH:253:GLU:H	1.62	0.63
5:AD:48:ASP:HB2	5:AD:56:GLN:OE1	1.97	0.63
9:AO:102:BCL:CHD	9:AP:101:BCL:HMD1	2.29	0.63
5:AS:10:LYS:HG2	14:AS:104:CRT:C1M	2.29	0.63
5:AS:34:LEU:HA	15:AS:101:PEF:C44	2.29	0.63
5:AS:46:TRP:CD1	5:AS:47:LEU:HD22	2.34	0.63
14:AS:104:CRT:H2M1	5:AW:33:LEU:O	1.99	0.63
6:AV:10:THR:HG22	6:AV:11:ASP:H	1.62	0.63
1:BC:40:MET:HA	1:BC:248:THR:HG22	1.80	0.63
6:BB:42:TYR:HH	6:B0:46:LEU:HB3	1.63	0.63
14:BF:103:CRT:H41	6:BJ:17:PHE:CD2	2.34	0.63
9:BG:101:BCL:CBB	9:BI:102:BCL:CHC	2.76	0.63
5:BO:4:MET:CE	6:BR:23:GLN:HB2	2.27	0.63
6:BV:21:PHE:HB2	14:BV:102:CRT:H11	1.80	0.63
5:B7:47:LEU:HD22	5:B7:47:LEU:N	2.14	0.63
5:AA:11:ILE:HD13	14:AA:102:CRT:C10	2.28	0.63
5:AD:14:ILE:N	5:AD:14:ILE:HD12	2.13	0.63
9:AW:101:BCL:HBC1	9:AX:101:BCL:HBC3	1.80	0.63
5:A1:5:ASN:HA	5:A1:8:LEU:CB	2.20	0.63
5:A5:44:LEU:HD12	5:A5:44:LEU:O	1.98	0.63
6:A0:7:THR:HG23	6:A0:8:GLY:N	2.13	0.63
1:BC:94:MET:SD	7:BC:501:HEM:FE	1.89	0.63
2:BL:185:ALA:N	2:BL:252:TRP:HD1	1.97	0.63
2:BL:201:SER:HG	3:BM:270:TRP:CD1	2.17	0.63
4:BH:251:THR:HG22	4:BH:253:GLU:H	1.61	0.63
9:BA:101:BCL:C15	5:B9:24:ILE:HD11	2.29	0.63
5:BF:14:ILE:HD12	5:BI:21:LEU:HD22	1.79	0.63
9:BS:102:BCL:HHC	9:BS:102:BCL:OBB	1.98	0.63
5:BU:46:TRP:CZ2	9:BU:102:BCL:H2C	2.34	0.63
9:B1:102:BCL:C1D	9:B2:101:BCL:HMD1	2.28	0.63
9:B8:101:BCL:HMA3	9:B9:102:BCL:CMA	2.29	0.63
1:AC:170:PRO:HG2	1:AC:171:GLY:H	1.62	0.63
2:AL:22:LEU:O	5:A9:18:ARG:NH1	2.31	0.63
9:AL:301:BCL:H71	9:AL:301:BCL:H41	1.81	0.63
3:AM:148:TRP:HA	3:AM:148:TRP:HE3	1.62	0.63
14:AB:102:CRT:H5	5:A9:10:LYS:CB	2.29	0.63
6:AJ:14:ALA:O	6:AJ:18:HIS:HB2	1.98	0.63
5:AO:34:LEU:HA	5:AO:37:MET:HB2	1.80	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:AS:10:LYS:HA	5:AS:13:LEU:HD12	1.80	0.63
6:AV:10:THR:HG22	6:AV:11:ASP:N	2.13	0.63
5:AW:36:HIS:NE2	9:AX:101:BCL:HMD3	2.13	0.63
5:A1:14:ILE:HD12	5:A1:15:LEU:N	2.13	0.63
2:BL:244:PHE:O	2:BL:245:LEU:C	2.41	0.63
14:BB:102:CRT:H2M2	5:BD:37:MET:HE3	1.79	0.63
5:BI:18:ARG:HB3	5:BI:18:ARG:CZ	2.28	0.63
6:BP:10:THR:HB	6:BP:13:GLU:OE1	1.97	0.63
5:BQ:5:ASN:ND2	6:BR:22:MET:HE3	2.14	0.63
5:BU:42:THR:HB	5:BW:48:ASP:HB3	1.80	0.63
14:BU:103:CRT:H2M3	5:BY:36:HIS:C	2.23	0.63
9:BY:102:BCL:HMD2	6:BZ:36:HIS:HD2	1.63	0.63
5:B1:16:ASP:HB3	5:B1:18:ARG:NE	2.13	0.63
5:B1:18:ARG:HD2	5:B1:19:ARG:N	2.13	0.63
4:AH:41:LEU:HD13	4:AH:48:ARG:NH1	2.14	0.63
5:AA:35:ILE:HG13	9:AB:101:BCL:O1D	1.97	0.63
5:AI:32:GLY:N	9:AJ:101:BCL:HED2	2.14	0.63
9:AK:102:BCL:C4D	9:AN:101:BCL:CMD	2.77	0.63
5:AO:3:THR:HB	5:AO:4:MET:SD	2.39	0.63
14:AP:102:CRT:H2M1	5:AQ:37:MET:HG2	1.80	0.63
6:AT:42:TYR:CD2	6:AT:43:ARG:HG2	2.34	0.63
6:AX:42:TYR:CE2	6:AX:43:ARG:HD2	2.33	0.63
5:AY:8:LEU:HD22	5:AY:11:ILE:CD1	2.29	0.63
3:BM:186:THR:HG23	3:BM:187:ALA:H	1.63	0.63
5:BD:9:TYR:CZ	6:BE:11:ASP:HB3	2.33	0.63
5:BS:42:THR:HG21	5:BU:47:LEU:HB3	1.80	0.63
6:BT:45:TRP:CE3	9:BT:101:BCL:H2C	2.33	0.63
5:BY:43:ASP:N	5:B1:48:ASP:HB3	2.13	0.63
5:B1:10:LYS:HB2	14:B1:103:CRT:H83	1.81	0.63
6:B8:34:ILE:O	6:B8:34:ILE:HD13	1.99	0.63
6:B0:18:HIS:O	6:B0:22:MET:HB2	1.98	0.63
1:AC:135:ARG:HG2	1:AC:330:LEU:CA	2.28	0.63
1:AC:190:VAL:C	1:AC:192:TYR:N	2.56	0.63
1:AC:266:ARG:O	1:AC:269:ALA:N	2.24	0.63
2:AL:38:VAL:HG23	2:AL:39:GLY:N	2.13	0.63
2:AL:110:ALA:O	2:AL:113:GLU:HB2	1.98	0.63
3:AM:47:GLN:NE2	3:AM:49:GLY:O	2.31	0.63
5:AD:12:TRP:HA	5:AD:12:TRP:CE3	2.34	0.63
5:AD:49:ASP:HB2	5:AF:56:GLN:CG	2.29	0.63
5:AK:5:ASN:HD21	6:AN:22:MET:CE	2.11	0.63
6:AN:30:GLY:O	6:AN:33:VAL:HG12	1.97	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:AQ:15:LEU:CD2	5:AS:18:ARG:HD3	2.27	0.63
5:A3:46:TRP:CZ3	9:A3:103:BCL:H2C	2.33	0.63
6:A8:44:PRO:O	5:A9:52:PRO:CD	2.46	0.63
1:BC:175:PRO:CD	1:BC:179:LYS:HB2	2.26	0.63
2:BL:138:LEU:C	2:BL:140:LEU:H	2.04	0.63
5:BA:21:LEU:HD23	5:B9:14:ILE:HG21	1.81	0.63
5:BU:13:LEU:HB2	14:BU:103:CRT:H1M2	1.80	0.63
6:BV:20:ILE:O	6:BV:20:ILE:HD13	1.99	0.63
9:BZ:101:BCL:HBB3	9:B1:102:BCL:CHC	2.28	0.63
5:B5:12:TRP:HZ3	5:B5:17:PRO:HA	1.64	0.63
6:B0:36:HIS:HE1	9:B0:102:BCL:CHB	2.11	0.63
1:AC:269:ALA:O	1:AC:273:ILE:HG13	1.98	0.63
3:AM:84:PHE:HA	5:AW:37:MET:CE	2.29	0.63
3:AM:201:PHE:CZ	4:AH:16:ILE:HA	2.34	0.63
3:AM:237:GLN:OE1	3:AM:244:ALA:HB3	1.99	0.63
3:AM:314:VAL:HG12	3:AM:315:ASN:H	1.64	0.63
5:AA:50:ASN:HA	5:AD:60:LYS:N	2.14	0.63
6:AJ:21:PHE:CD1	6:AJ:21:PHE:C	2.73	0.63
5:AQ:31:LEU:CG	9:AR:101:BCL:HED3	2.29	0.63
6:A0:36:HIS:HE1	9:A0:102:BCL:CHB	2.12	0.63
1:BC:24:GLU:CD	2:BL:266:ARG:HH22	2.06	0.63
1:BC:94:MET:SD	7:BC:501:HEM:NA	2.72	0.63
2:BL:168:ASN:O	2:BL:171:TYR:N	2.32	0.63
4:BH:55:VAL:CG1	4:BH:56:VAL:H	2.09	0.63
4:BH:69:LEU:HD23	4:BH:70:PRO:HD2	1.81	0.63
6:BN:31:LEU:HA	6:BN:34:ILE:HG22	1.79	0.63
5:BY:42:THR:HB	5:B1:48:ASP:CG	2.24	0.63
5:B1:50:ASN:HB3	5:B3:60:LYS:CA	2.27	0.63
6:B4:13:GLU:CD	6:B4:13:GLU:H	2.06	0.63
2:AL:138:LEU:C	2:AL:140:LEU:H	2.07	0.62
3:AM:81:TRP:O	5:AU:41:SER:HB3	1.99	0.62
3:AM:290:VAL:HG11	4:AH:12:ALA:HB2	1.79	0.62
4:AH:258:LEU:O	5:A5:19:ARG:HD3	2.00	0.62
5:AF:14:ILE:HD12	5:AI:21:LEU:CD2	2.28	0.62
6:AR:44:PRO:HG2	5:AS:52:PRO:HG3	1.81	0.62
5:AW:21:LEU:O	5:AW:21:LEU:HD12	1.99	0.62
1:BC:52:SER:O	1:BC:56:ASN:HB2	1.99	0.62
5:BA:27:PHE:CE1	5:BD:29:ILE:CD1	2.80	0.62
5:BS:29:ILE:O	5:BS:33:LEU:HD13	1.99	0.62
5:BW:40:LEU:HD12	5:BW:45:ASN:HA	1.80	0.62
6:BX:38:LEU:C	6:BX:38:LEU:HD23	2.24	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:B4:45:TRP:O	6:B4:46:LEU:HG	1.99	0.62
1:AC:20:LEU:HG	2:AL:271:TRP:NE1	2.14	0.62
1:AC:20:LEU:HG	2:AL:271:TRP:HE1	1.64	0.62
4:AH:213:ALA:O	4:AH:246:GLY:HA3	1.99	0.62
5:AD:31:LEU:HG	9:AE:101:BCL:HED3	1.80	0.62
5:AK:43:ASP:OD1	5:AO:47:LEU:HB3	1.99	0.62
6:AN:13:GLU:CD	6:AN:13:GLU:H	2.07	0.62
6:AN:31:LEU:HA	6:AN:34:ILE:CG2	2.29	0.62
5:AO:13:LEU:HA	6:AP:9:LEU:HB2	1.80	0.62
5:AO:43:ASP:CA	5:AQ:48:ASP:HB3	2.29	0.62
5:AU:9:TYR:HB2	6:AV:15:LYS:HD3	1.81	0.62
6:AV:20:ILE:HD13	6:AV:20:ILE:C	2.24	0.62
6:A6:40:TRP:HZ3	6:A6:44:PRO:HA	1.64	0.62
5:A9:12:TRP:HA	5:A9:12:TRP:HE3	1.64	0.62
9:A9:102:BCL:CHD	9:A0:102:BCL:HMD1	2.29	0.62
6:A0:32:VAL:HG12	6:A0:33:VAL:H	1.63	0.62
1:BC:97:VAL:O	1:BC:97:VAL:HG12	1.98	0.62
6:BB:22:MET:C	6:BB:26:TYR:HE1	2.06	0.62
6:BX:13:GLU:O	6:BX:16:GLU:HB3	1.98	0.62
5:B5:18:ARG:HH11	5:B5:18:ARG:HG3	1.64	0.62
5:B7:46:TRP:CD1	5:B7:47:LEU:CD2	2.82	0.62
6:B0:17:PHE:CD1	14:B0:101:CRT:C9	2.74	0.62
1:AC:128:ARG:O	1:AC:131:PHE:HB2	1.99	0.62
1:AC:148:THR:HG23	1:AC:322:GLN:CA	2.29	0.62
3:AM:275:LEU:HD21	4:AH:19:PHE:HE2	1.63	0.62
9:AM:401:BCL:HBC1	9:AM:402:BCL:CBD	2.29	0.62
4:AH:52:ARG:NH1	4:AH:52:ARG:HB3	2.14	0.62
9:AE:101:BCL:CHC	9:AF:102:BCL:HBB3	2.29	0.62
14:A1:103:CRT:H9	6:A4:17:PHE:HE1	1.64	0.62
1:BC:284:ILE:HG21	1:BC:304:ARG:HA	1.80	0.62
3:BM:250:LEU:O	3:BM:254:TRP:CD1	2.52	0.62
4:BH:88:ASN:ND2	4:BH:109:SER:HB2	2.14	0.62
4:BH:166:THR:O	4:BH:184:VAL:HG13	2.00	0.62
5:BA:36:HIS:CD2	9:BB:101:BCL:CMD	2.83	0.62
9:BA:101:BCL:HBB3	9:B0:102:BCL:C4B	2.28	0.62
14:BA:102:CRT:H9	6:BE:17:PHE:CD1	2.34	0.62
6:BB:29:PHE:HE1	9:BB:101:BCL:H11	1.64	0.62
9:BO:102:BCL:HAC2	9:BP:101:BCL:CAC	2.29	0.62
5:BU:13:LEU:HD22	6:BV:9:LEU:HB2	1.79	0.62
5:BW:14:ILE:HG21	5:BY:21:LEU:HD12	1.80	0.62
14:BW:103:CRT:H35	5:BY:31:LEU:HD11	1.81	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:B0:10:THR:HB	6:B0:13:GLU:OE2	2.00	0.62
1:AC:36:ARG:HG3	1:AC:36:ARG:O	1.99	0.62
1:AC:166:TRP:O	1:AC:166:TRP:CE3	2.53	0.62
1:AC:236:MET:SD	7:AC:503:HEM:NC	2.72	0.62
4:AH:125:LEU:HB2	4:AH:129:GLY:O	1.99	0.62
5:AI:16:ASP:O	5:AI:20:VAL:HG22	1.99	0.62
5:AI:36:HIS:NE2	9:AJ:101:BCL:HMD3	2.15	0.62
9:AI:102:BCL:ND	9:AJ:101:BCL:HMD3	2.14	0.62
5:AO:8:LEU:CA	6:AR:20:ILE:HD11	2.20	0.62
9:AQ:102:BCL:C1D	9:AR:101:BCL:HMD1	2.30	0.62
6:AV:44:PRO:HG2	5:AW:52:PRO:HG2	1.80	0.62
5:AY:43:ASP:N	5:A1:48:ASP:HB3	2.13	0.62
3:BM:75:MET:HE1	3:BM:90:PHE:HE1	1.64	0.62
5:BA:36:HIS:HB3	14:B0:101:CRT:C39	2.27	0.62
9:BO:102:BCL:CHD	9:BP:101:BCL:HMD1	2.30	0.62
5:BS:55:TYR:HD1	5:BS:56:GLN:N	1.97	0.62
1:AC:259:TRP:O	1:AC:261:GLN:N	2.32	0.62
2:AL:231:TYR:OH	2:AL:233:ILE:HA	1.99	0.62
4:AH:164:ALA:HB2	4:AH:216:ALA:CB	2.30	0.62
5:AQ:44:LEU:HD22	6:AR:43:ARG:HD3	1.82	0.62
5:AS:10:LYS:HB3	14:AS:104:CRT:H1M2	1.79	0.62
5:AS:35:ILE:O	5:AS:36:HIS:C	2.42	0.62
5:AY:15:LEU:HG	5:A1:21:LEU:HD21	1.79	0.62
5:A1:43:ASP:HB2	5:A3:47:LEU:CD1	2.29	0.62
5:A5:33:LEU:HD12	5:A5:34:LEU:N	2.13	0.62
6:A0:40:TRP:HH2	6:A0:46:LEU:CG	2.09	0.62
2:BL:264:TRP:CH2	2:BL:271:TRP:HA	2.35	0.62
5:BF:29:ILE:HA	9:BF:102:BCL:C1	2.25	0.62
5:BF:43:ASP:OD2	5:BI:47:LEU:O	2.18	0.62
5:BO:36:HIS:CE1	9:BP:101:BCL:CMD	2.83	0.62
5:BU:11:ILE:HG12	14:BU:103:CRT:C7	2.29	0.62
5:BY:25:VAL:O	5:BY:29:ILE:HB	1.99	0.62
5:BY:36:HIS:NE2	9:BZ:101:BCL:HMD3	2.14	0.62
1:AC:154:THR:O	1:AC:157:ARG:HG3	1.98	0.62
2:AL:26:TRP:HE3	4:AH:97:GLY:O	1.81	0.62
2:AL:71:TRP:N	2:AL:71:TRP:CE3	2.68	0.62
3:AM:35:ILE:HG22	3:AM:36:PHE:H	1.64	0.62
3:AM:102:TYR:HD1	3:AM:102:TYR:H	1.47	0.62
3:AM:261:THR:HG23	4:AH:34:ASP:O	1.99	0.62
9:AI:102:BCL:C1D	9:AJ:101:BCL:HMD1	2.29	0.62
9:AO:102:BCL:H111	9:AO:102:BCL:H192	1.82	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:AP:10:THR:HG22	6:AP:11:ASP:N	2.14	0.62
5:AS:13:LEU:HB2	14:AS:104:CRT:H31A	1.81	0.62
5:AU:12:TRP:HA	5:AU:12:TRP:HE3	1.64	0.62
5:AU:52:PRO:HB3	5:AU:55:TYR:HE1	1.64	0.62
5:AY:50:ASN:HD22	5:A1:58:LEU:CB	2.12	0.62
5:A1:11:ILE:N	14:A1:103:CRT:H82	2.15	0.62
6:A4:41:LEU:HD23	6:A4:41:LEU:O	1.99	0.62
6:A0:10:THR:HB	6:A0:13:GLU:OE2	2.00	0.62
1:BC:80:GLN:HG3	1:BC:128:ARG:NH2	2.15	0.62
1:BC:196:PRO:HG3	1:BC:231:TRP:CD1	2.34	0.62
1:BC:249:PHE:CZ	1:BC:265:LYS:HG2	2.34	0.62
9:BL:303:BCL:H2A	9:BL:303:BCL:O1D	2.00	0.62
3:BM:35:ILE:HG22	3:BM:36:PHE:H	1.64	0.62
3:BM:98:PRO:HB2	3:BM:171:TRP:HB3	1.81	0.62
3:BM:215:LEU:O	3:BM:217:ALA:N	2.32	0.62
9:BD:102:BCL:HAC2	9:BE:101:BCL:HAC1	1.81	0.62
9:BQ:104:BCL:HMA1	9:BS:102:BCL:HMA1	1.81	0.62
6:BT:32:VAL:HG21	9:BT:101:BCL:HBA2	1.81	0.62
5:BU:9:TYR:HA	6:BV:18:HIS:CG	2.34	0.62
5:BU:56:GLN:NE2	5:BU:57:ALA:N	2.48	0.62
14:BV:102:CRT:H2M3	5:BW:37:MET:N	2.15	0.62
6:B2:45:TRP:O	6:B2:46:LEU:CG	2.47	0.62
4:AH:31:ARG:O	4:AH:34:ASP:HB2	2.00	0.62
5:AO:46:TRP:HA	5:AO:49:ASP:OD2	2.00	0.62
5:AO:51:ILE:CG1	5:AO:52:PRO:HD2	2.29	0.62
14:AS:104:CRT:C14	6:AV:24:SER:OG	2.48	0.62
6:AT:10:THR:HG22	6:AT:11:ASP:N	2.13	0.62
6:A8:45:TRP:O	6:A8:46:LEU:CG	2.48	0.62
5:A9:8:LEU:HD22	5:A9:11:ILE:HD11	1.80	0.62
1:BC:134:VAL:HG13	1:BC:150:VAL:HG22	1.80	0.62
2:BL:257:ILE:HG22	9:BL:301:BCL:HED2	1.80	0.62
6:BB:24:SER:O	6:BB:27:ALA:HB3	1.99	0.62
5:BI:27:PHE:HE2	5:BK:29:ILE:HD11	1.65	0.62
6:BR:38:LEU:C	6:BR:38:LEU:HD12	2.24	0.62
5:BW:22:VAL:O	5:BW:25:VAL:HB	2.00	0.62
5:B5:18:ARG:HB2	5:B5:19:ARG:NH2	2.15	0.62
6:B6:45:TRP:HD1	6:B6:46:LEU:H	1.47	0.62
5:B7:44:LEU:HD22	5:B7:44:LEU:O	1.99	0.62
6:B8:45:TRP:O	6:B8:46:LEU:CG	2.48	0.62
3:AM:207:ALA:O	3:AM:210:TYR:HB2	1.99	0.62
9:AA:101:BCL:H71	6:AB:28:TRP:CZ3	2.34	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:AO:9:TYR:HA	6:AP:18:HIS:ND1	2.14	0.62
9:AO:102:BCL:HBD	9:AP:101:BCL:OBD	1.99	0.62
14:AS:104:CRT:C39	5:AW:36:HIS:HB2	2.28	0.62
5:AW:12:TRP:HA	5:AW:12:TRP:HE3	1.63	0.62
5:A9:12:TRP:HA	5:A9:12:TRP:CE3	2.33	0.62
2:BL:87:ALA:N	2:BL:96:GLN:HE22	1.98	0.62
2:BL:160:LEU:HA	2:BL:163:LEU:HD13	1.81	0.62
2:BL:189:PHE:HE2	2:BL:253:SER:OG	1.82	0.62
3:BM:98:PRO:HG3	3:BM:107:PRO:HG3	1.80	0.62
4:BH:14:ILE:O	4:BH:17:TRP:HB2	2.00	0.62
5:BD:8:LEU:C	5:BD:10:LYS:H	2.08	0.62
6:BE:45:TRP:HA	5:BF:52:PRO:CD	2.29	0.62
5:BF:35:ILE:HA	5:BF:38:ILE:HG22	1.80	0.62
5:BQ:20:VAL:O	5:BQ:24:ILE:HD13	1.99	0.62
5:BU:35:ILE:O	5:BU:38:ILE:HG22	1.99	0.62
14:BU:103:CRT:C34	9:BY:102:BCL:CBA	2.76	0.62
9:BX:101:BCL:HMC3	5:BY:47:LEU:HD21	1.80	0.62
14:B2:102:CRT:H2M2	5:B3:36:HIS:HB3	1.81	0.62
6:B4:41:LEU:HD23	6:B4:41:LEU:O	1.99	0.62
5:B7:27:PHE:CZ	5:B9:29:ILE:HD11	2.35	0.62
3:AM:89:HIS:O	3:AM:93:LEU:HG	1.99	0.62
3:AM:234:GLU:O	3:AM:235:ILE:C	2.43	0.62
4:AH:171:TRP:CE2	4:AH:194:LEU:HD21	2.35	0.62
9:AA:101:BCL:HED1	6:AB:31:LEU:HB3	1.82	0.62
9:AE:101:BCL:CHB	9:AF:102:BCL:HMB2	2.29	0.62
5:AF:19:ARG:HH12	5:AI:18:ARG:HH22	1.47	0.62
5:AU:13:LEU:HD21	6:AV:14:ALA:HB2	1.82	0.62
5:AW:26:ALA:O	5:AW:30:VAL:HG12	1.99	0.62
5:AW:31:LEU:HD13	14:AX:102:CRT:H35	1.80	0.62
5:A1:57:ALA:C	5:A1:59:GLY:H	2.06	0.62
1:BC:298:PRO:C	1:BC:300:GLY:H	2.07	0.62
2:BL:117:CYS:SG	2:BL:124:PHE:HA	2.39	0.62
2:BL:119:LYS:C	2:BL:121:GLY:H	2.07	0.62
2:BL:151:TRP:C	2:BL:153:HIS:H	2.08	0.62
3:BM:287:SER:CB	3:BM:294:TRP:HE1	2.12	0.62
5:BA:2:PHE:HB2	5:BA:5:ASN:OD1	2.00	0.62
14:BU:103:CRT:C2M	5:BY:36:HIS:C	2.72	0.62
14:BU:103:CRT:H342	9:BY:102:BCL:HBA1	1.80	0.62
9:B1:102:BCL:C1D	9:B2:101:BCL:CMD	2.77	0.62
6:B8:31:LEU:O	6:B8:34:ILE:HG22	2.00	0.62
2:AL:131:SER:O	2:AL:134:ILE:HB	2.00	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:AM:17:ALA:O	3:AM:19:PRO:HD3	2.00	0.62
3:AM:152:ALA:HB2	3:AM:274:VAL:HG13	1.82	0.62
3:AM:274:VAL:HG12	3:AM:278:ILE:HD11	1.82	0.62
4:AH:32:ARG:HG3	4:AH:59:PRO:HB2	1.82	0.62
4:AH:63:ASP:O	4:AH:79:PRO:HD2	1.99	0.62
5:AK:16:ASP:O	5:AK:20:VAL:HG22	1.99	0.62
6:AN:34:ILE:HD13	6:AN:34:ILE:C	2.25	0.62
6:AN:38:LEU:O	6:AN:38:LEU:HD23	2.00	0.62
6:AP:12:ASP:O	6:AP:16:GLU:HG3	2.00	0.62
5:AQ:40:LEU:HD12	5:AQ:45:ASN:HA	1.82	0.62
6:AV:28:TRP:HA	6:AV:31:LEU:HD12	1.81	0.62
5:AW:26:ALA:HA	5:AW:29:ILE:HG22	1.82	0.62
5:A1:12:TRP:CD1	6:A2:17:PHE:CD2	2.88	0.62
3:BM:276:THR:HG22	3:BM:277:VAL:N	2.15	0.62
4:BH:48:ARG:HG2	4:BH:57:GLY:H	1.65	0.62
6:BE:40:TRP:HZ3	6:BE:45:TRP:H	1.47	0.62
5:BF:11:ILE:HB	14:BF:103:CRT:C8	2.24	0.62
14:BG:102:CRT:H2M2	5:BI:37:MET:CE	2.18	0.62
5:BK:44:LEU:O	5:BK:44:LEU:HD22	1.99	0.62
5:BQ:50:ASN:HB3	5:BS:56:GLN:HA	1.82	0.62
9:BS:102:BCL:CHD	9:BS:102:BCL:HBC2	2.28	0.62
14:BV:102:CRT:H2M3	5:BW:37:MET:CA	2.30	0.62
9:B7:103:BCL:HMD3	6:B8:36:HIS:CD2	2.35	0.62
5:B9:4:MET:O	5:B9:8:LEU:HG	2.00	0.62
2:AL:94:LEU:HA	2:AL:97:ILE:CD1	2.29	0.61
2:AL:188:PHE:CD2	2:AL:248:SER:HB3	2.35	0.61
2:AL:252:TRP:O	2:AL:253:SER:C	2.43	0.61
10:AM:403:BPH:H9C3	15:AM:409:PEF:C22	2.27	0.61
5:AA:45:ASN:OD1	5:AA:47:LEU:HB2	1.99	0.61
6:AB:20:ILE:HD12	14:AB:102:CRT:H81	1.82	0.61
5:AD:7:ASN:HD22	5:AD:7:ASN:H	1.46	0.61
6:AJ:33:VAL:O	6:AJ:37:LEU:HD23	2.00	0.61
5:AU:9:TYR:HA	6:AV:18:HIS:ND1	2.15	0.61
9:A8:101:BCL:H152	9:A8:101:BCL:C20	2.25	0.61
9:A0:102:BCL:C2B	9:A0:102:BCL:H141	2.30	0.61
2:BL:206:VAL:HG12	3:BM:142:MET:SD	2.39	0.61
5:BK:14:ILE:HG23	5:BO:18:ARG:HG2	1.82	0.61
6:BP:38:LEU:HD23	6:BP:38:LEU:C	2.25	0.61
5:BU:13:LEU:C	6:BV:7:THR:HA	2.25	0.61
5:BU:36:HIS:CE1	9:BU:102:BCL:NA	2.68	0.61
6:BV:46:LEU:HD13	6:BX:42:TYR:CE1	2.35	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:B3:43:ASP:HB2	5:B5:47:LEU:CD1	2.13	0.61
1:AC:167:VAL:HG23	1:AC:301:ASP:CG	2.25	0.61
2:AL:196:LEU:HB2	3:AM:216:PHE:CG	2.35	0.61
3:AM:42:LYS:HE3	15:AM:408:PEF:H42	1.81	0.61
5:AD:15:LEU:HB3	5:AD:20:VAL:CG2	2.30	0.61
5:AK:47:LEU:HD22	5:AK:47:LEU:H	1.65	0.61
9:AN:101:BCL:CMB	9:AO:102:BCL:C1B	2.79	0.61
5:AO:9:TYR:HA	6:AP:18:HIS:CE1	2.34	0.61
5:AO:26:ALA:HA	5:AO:29:ILE:HG22	1.81	0.61
6:AP:44:PRO:HG2	5:AQ:52:PRO:HB3	1.81	0.61
5:A7:44:LEU:HD23	6:A8:43:ARG:NH1	2.15	0.61
5:A7:44:LEU:HD22	5:A7:46:TRP:HE3	1.62	0.61
2:BL:13:ARG:HG3	2:BL:13:ARG:O	1.99	0.61
3:BM:77:ALA:O	3:BM:78:SER:C	2.41	0.61
6:BB:45:TRP:O	6:BB:46:LEU:HB2	2.00	0.61
9:BF:102:BCL:H62	6:BG:28:TRP:CH2	2.35	0.61
5:BY:49:ASP:HB2	5:B1:56:GLN:NE2	2.16	0.61
3:AM:157:TYR:CD1	3:AM:157:TYR:C	2.79	0.61
3:AM:208:PHE:C	3:AM:210:TYR:N	2.57	0.61
5:AA:50:ASN:CG	5:AA:51:ILE:N	2.58	0.61
5:AD:31:LEU:O	5:AD:35:ILE:HG12	2.00	0.61
5:AO:9:TYR:CE1	5:AO:10:LYS:HD3	2.34	0.61
9:AO:102:BCL:H2A	9:AO:102:BCL:O1D	2.01	0.61
6:AR:45:TRP:HD1	6:AR:46:LEU:H	1.48	0.61
2:BL:181:ALA:O	2:BL:183:MET:N	2.34	0.61
5:BO:44:LEU:HD11	5:BO:46:TRP:HE3	1.64	0.61
6:BZ:45:TRP:CG	9:BZ:101:BCL:H2C	2.35	0.61
9:B7:103:BCL:CMD	6:B8:36:HIS:CD2	2.83	0.61
1:AC:126:VAL:HG23	1:AC:127:SER:N	2.15	0.61
1:AC:307:CYS:O	1:AC:311:HIS:HB2	2.00	0.61
2:AL:10:TYR:HA	4:AH:112:GLY:HA2	1.81	0.61
2:AL:16:THR:OG1	4:AH:257:PRO:HB3	2.00	0.61
2:AL:188:PHE:C	2:AL:190:PHE:N	2.57	0.61
2:AL:214:PRO:HA	4:AH:68:VAL:O	2.01	0.61
3:AM:134:TYR:HA	3:AM:144:GLN:HE22	1.64	0.61
4:AH:123:CYS:N	4:AH:232:THR:HG22	2.16	0.61
5:AD:19:ARG:O	5:AD:23:SER:HB3	2.01	0.61
5:AI:4:MET:SD	6:AN:23:GLN:HB3	2.40	0.61
6:AJ:30:GLY:O	6:AJ:33:VAL:HG12	2.00	0.61
5:AO:4:MET:O	5:AO:7:ASN:ND2	2.32	0.61
5:AO:14:ILE:CG2	5:AO:15:LEU:HG	2.30	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:AS:13:LEU:HB2	14:AS:104:CRT:H32A	1.81	0.61
14:AW:102:CRT:H6	6:AZ:17:PHE:HD1	1.65	0.61
6:A2:29:PHE:N	6:A2:29:PHE:HD1	1.99	0.61
5:A3:46:TRP:CZ3	9:A3:103:BCL:HBC3	2.36	0.61
2:BL:181:ALA:HB3	2:BL:256:CYS:HA	1.82	0.61
2:BL:195:ALA:HB3	3:BM:216:PHE:HE2	1.64	0.61
3:BM:35:ILE:HG22	3:BM:36:PHE:N	2.15	0.61
3:BM:137:ALA:O	3:BM:142:MET:HB2	2.00	0.61
4:BH:235:GLU:HA	4:BH:238:LYS:HG2	1.82	0.61
14:BA:102:CRT:H342	9:BF:102:BCL:CBA	2.25	0.61
6:BB:27:ALA:O	6:BB:31:LEU:HG	2.01	0.61
9:BD:102:BCL:ND	9:BE:101:BCL:CMD	2.63	0.61
6:BE:23:GLN:HG3	6:BE:24:SER:N	2.16	0.61
5:BK:16:ASP:HB3	5:BK:18:ARG:HE	1.66	0.61
6:BN:20:ILE:HD12	6:BN:20:ILE:N	2.16	0.61
5:BO:12:TRP:O	6:BP:9:LEU:HD12	1.99	0.61
6:BP:30:GLY:O	6:BP:33:VAL:HG12	2.01	0.61
6:BP:31:LEU:O	6:BP:34:ILE:HG13	2.00	0.61
6:B0:32:VAL:HG12	6:B0:33:VAL:N	2.14	0.61
2:AL:164:ASP:O	2:AL:167:SER:N	2.33	0.61
4:AH:5:ILE:HG12	4:AH:6:THR:N	2.15	0.61
4:AH:123:CYS:CA	4:AH:232:THR:HA	2.29	0.61
5:AA:22:VAL:HA	5:AA:25:VAL:HG23	1.82	0.61
6:AB:24:SER:O	6:AB:27:ALA:HB3	2.00	0.61
5:AD:40:LEU:CD1	5:AD:47:LEU:HD23	2.28	0.61
9:AJ:101:BCL:C4B	9:AK:102:BCL:HBB3	2.30	0.61
5:AS:30:VAL:CG2	15:AS:101:PEF:H392	2.25	0.61
5:AU:36:HIS:CE1	9:AV:102:BCL:HMD3	2.35	0.61
5:AW:2:PHE:HB2	5:AW:5:ASN:HB2	1.81	0.61
5:AY:19:ARG:O	5:AY:23:SER:HB3	2.00	0.61
5:A1:5:ASN:O	5:A1:8:LEU:HD22	2.01	0.61
14:A1:103:CRT:H9	6:A4:17:PHE:CE1	2.36	0.61
2:BL:171:TYR:C	2:BL:173:PHE:H	2.08	0.61
3:BM:7:ILE:HB	15:BM:407:PEF:HN1	1.64	0.61
4:BH:197:ILE:HD13	4:BH:197:ILE:O	2.01	0.61
6:BB:29:PHE:HZ	9:BB:101:BCL:H42	1.66	0.61
5:BU:53:VAL:HA	5:BU:55:TYR:CZ	2.36	0.61
5:B7:43:ASP:CB	5:B9:47:LEU:HD12	2.21	0.61
1:AC:285:TRP:HB3	1:AC:286:PRO:HD3	1.81	0.61
2:AL:231:TYR:CD2	3:AM:48:ILE:HD13	2.35	0.61
4:AH:54:LYS:HE3	5:AD:23:SER:CB	2.30	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:AO:7:ASN:ND2	6:AR:23:GLN:OE1	2.34	0.61
5:AO:50:ASN:OD1	6:AP:43:ARG:NH2	2.34	0.61
6:AP:34:ILE:O	6:AP:38:LEU:HB3	2.00	0.61
14:AR:102:CRT:H342	9:AS:103:BCL:CBA	2.28	0.61
5:AW:9:TYR:CE1	6:AX:15:LYS:HB2	2.34	0.61
5:AW:14:ILE:HD13	5:AY:17:PRO:HB2	1.83	0.61
5:AY:11:ILE:HD13	9:A1:102:BCL:C15	2.28	0.61
14:A1:103:CRT:C34	9:A5:102:BCL:HBA1	2.31	0.61
6:A2:16:GLU:HB3	14:A2:102:CRT:H1M1	1.82	0.61
6:A8:22:MET:SD	6:A8:26:TYR:HE2	2.23	0.61
2:BL:40:PHE:O	2:BL:43:THR:HB	2.00	0.61
2:BL:139:VAL:HG23	2:BL:143:VAL:HB	1.81	0.61
9:BM:401:BCL:CMB	9:BM:402:BCL:H152	2.30	0.61
4:BH:164:ALA:HB2	4:BH:216:ALA:HB1	1.81	0.61
5:BU:44:LEU:HB3	5:BW:55:TYR:CD1	2.35	0.61
5:B3:20:VAL:O	5:B3:24:ILE:HG12	2.00	0.61
1:AC:264:PRO:HG2	1:AC:265:LYS:H	1.65	0.61
1:AC:292:PRO:O	1:AC:296:LYS:HG3	2.01	0.61
3:AM:59:LEU:CG	3:AM:128:LEU:HD21	2.31	0.61
3:AM:261:THR:C	3:AM:263:GLU:N	2.57	0.61
4:AH:19:PHE:CD1	4:AH:20:TRP:N	2.68	0.61
5:AD:49:ASP:HB2	5:AF:56:GLN:HB3	1.83	0.61
5:AS:10:LYS:HD3	14:AS:104:CRT:C1M	2.29	0.61
5:AY:31:LEU:HD23	9:AZ:101:BCL:HED3	1.81	0.61
9:A1:102:BCL:H143	14:A2:102:CRT:H132	1.80	0.61
5:A7:36:HIS:NE2	9:A8:101:BCL:HMD3	2.15	0.61
1:BC:33:ILE:HD12	1:BC:33:ILE:N	2.15	0.61
1:BC:276:VAL:HG13	1:BC:277:ARG:H	1.65	0.61
2:BL:15:GLY:O	2:BL:118:ARG:HD3	1.99	0.61
2:BL:159:ILE:H	2:BL:159:ILE:HD12	1.65	0.61
3:BM:12:GLN:HB2	4:BH:145:ALA:CB	2.29	0.61
4:BH:215:LYS:N	4:BH:218:HIS:HD2	1.99	0.61
5:BA:29:ILE:HG12	5:B9:27:PHE:CE2	2.36	0.61
9:BA:101:BCL:CBA	14:B0:101:CRT:H342	2.25	0.61
6:BB:37:LEU:HD22	9:BB:101:BCL:H193	1.83	0.61
14:BW:103:CRT:H83	6:BZ:20:ILE:HD13	1.81	0.61
6:B6:22:MET:SD	6:B6:25:MET:HE3	2.41	0.61
6:B8:22:MET:SD	6:B8:26:TYR:HE2	2.23	0.61
9:B9:102:BCL:ND	9:B0:102:BCL:HMD3	2.15	0.61
2:AL:97:ILE:HA	2:AL:100:ILE:CD1	2.28	0.61
2:AL:170:GLY:HA3	9:AL:301:BCL:CBC	2.30	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:AL:184:LEU:CB	2:AL:252:TRP:HE1	2.13	0.61
2:AL:186:ILE:HD13	9:AM:401:BCL:CMD	2.31	0.61
3:AM:156:PHE:HA	3:AM:159:VAL:HG23	1.82	0.61
3:AM:180:PHE:O	3:AM:183:LEU:N	2.34	0.61
4:AH:69:LEU:CD2	4:AH:70:PRO:HD2	2.30	0.61
14:AB:102:CRT:O2	5:AD:33:LEU:HD12	2.00	0.61
5:AD:12:TRP:HA	5:AD:12:TRP:HE3	1.64	0.61
9:AD:102:BCL:CHD	9:AE:101:BCL:HMD1	2.31	0.61
5:AF:10:LYS:HB2	14:AJ:102:CRT:H5	1.81	0.61
5:AF:49:ASP:CG	5:AF:50:ASN:N	2.58	0.61
6:AN:37:LEU:HD22	9:AN:101:BCL:H193	1.82	0.61
5:AS:9:TYR:HA	6:AT:18:HIS:HB2	1.82	0.61
9:A7:103:BCL:C1	9:A7:103:BCL:O1A	2.49	0.61
2:BL:188:PHE:CE2	2:BL:248:SER:HB3	2.35	0.61
2:BL:270:GLU:O	2:BL:271:TRP:C	2.42	0.61
4:BH:151:PRO:O	4:BH:154:MET:HG3	2.00	0.61
5:BK:44:LEU:HD22	5:BK:46:TRP:H	1.66	0.61
9:BN:101:BCL:H172	6:BP:38:LEU:CD2	2.30	0.61
6:B2:40:TRP:CE3	6:B2:44:PRO:HA	2.36	0.61
5:B3:31:LEU:O	5:B3:35:ILE:HG12	2.01	0.61
1:AC:28:PRO:CD	2:AL:262:PRO:HA	2.30	0.61
1:AC:130:MET:O	1:AC:133:LEU:HB3	2.00	0.61
1:AC:270:TRP:O	1:AC:274:ARG:HD2	1.99	0.61
2:AL:52:TRP:HA	2:AL:52:TRP:CE3	2.36	0.61
5:AA:29:ILE:HG12	5:A9:27:PHE:CE2	2.36	0.61
6:AN:45:TRP:O	6:AN:46:LEU:HB2	2.00	0.61
5:AS:16:ASP:HB2	5:AS:19:ARG:HD3	1.82	0.61
5:AS:17:PRO:O	5:AS:20:VAL:HG22	2.00	0.61
5:AW:21:LEU:HD22	14:AX:102:CRT:C13	2.30	0.61
5:AY:7:ASN:O	6:A2:20:ILE:CG1	2.49	0.61
5:AY:9:TYR:CZ	5:AY:10:LYS:HE3	2.36	0.61
5:AY:16:ASP:HB3	5:AY:18:ARG:HD2	1.82	0.61
6:A8:17:PHE:CD1	6:A8:20:ILE:HG21	2.36	0.61
4:BH:48:ARG:HE	4:BH:57:GLY:HA2	1.64	0.61
4:BH:96:PRO:O	5:B9:19:ARG:HD3	2.01	0.61
9:BI:102:BCL:H13	9:BI:102:BCL:H192	1.83	0.61
6:BJ:45:TRP:O	6:BJ:46:LEU:HB2	2.01	0.61
5:BU:12:TRP:NE1	6:BV:17:PHE:CD2	2.69	0.61
5:BU:35:ILE:HG22	5:BU:36:HIS:N	2.14	0.61
5:BU:46:TRP:CD1	5:BU:47:LEU:HD13	2.36	0.61
5:B3:56:GLN:NE2	5:B3:56:GLN:N	2.49	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AC:190:VAL:HG12	1:AC:237:MET:HB2	1.82	0.61
2:AL:22:LEU:HB2	5:A7:19:ARG:HB2	1.80	0.61
3:AM:2:PRO:HB3	4:AH:201:ARG:HH12	1.65	0.61
3:AM:179:ILE:HD13	3:AM:179:ILE:H	1.65	0.61
4:AH:241:ALA:O	4:AH:244:ALA:HB3	2.01	0.61
5:AA:40:LEU:HB2	5:AA:46:TRP:CH2	2.36	0.61
5:AA:47:LEU:HB3	5:A9:43:ASP:HB2	1.83	0.61
9:AA:101:BCL:CHB	14:A0:101:CRT:H372	2.31	0.61
5:AD:46:TRP:CH2	9:AD:102:BCL:HBC3	2.36	0.61
6:AE:38:LEU:O	6:AE:38:LEU:HD23	1.99	0.61
5:AF:9:TYR:CD1	6:AG:15:LYS:HG3	2.36	0.61
5:AS:9:TYR:CB	6:AT:15:LYS:HA	2.26	0.61
9:A7:103:BCL:HAC2	9:A8:101:BCL:HAC1	1.82	0.61
6:A8:46:LEU:HB3	6:A0:42:TYR:OH	1.99	0.61
1:BC:42:ASN:HA	2:BL:172:GLN:OE1	2.01	0.61
1:BC:80:GLN:HG3	1:BC:128:ARG:HH22	1.65	0.61
1:BC:270:TRP:CE3	1:BC:271:TYR:CD1	2.88	0.61
2:BL:280:LEU:HD21	5:BY:37:MET:CE	2.31	0.61
5:BA:26:ALA:O	5:BA:29:ILE:HG22	2.01	0.61
5:BD:8:LEU:O	5:BD:10:LYS:N	2.33	0.61
5:BF:10:LYS:HB3	14:BF:103:CRT:O1	2.01	0.61
5:BF:11:ILE:O	5:BF:14:ILE:HG12	2.01	0.61
9:BG:101:BCL:HMB3	9:BI:102:BCL:C4A	2.31	0.61
6:BN:17:PHE:HD1	14:BN:102:CRT:H6	1.62	0.61
14:BS:103:CRT:H6	6:BT:17:PHE:CD2	2.36	0.61
5:BW:26:ALA:HA	5:BW:29:ILE:CG2	2.31	0.61
14:B5:103:CRT:C40	5:B7:38:ILE:HG21	2.30	0.61
1:AC:97:VAL:HG21	1:AC:131:PHE:CZ	2.37	0.60
2:AL:174:LEU:N	2:AL:174:LEU:HD12	2.15	0.60
2:AL:253:SER:HA	2:AL:256:CYS:HB2	1.82	0.60
2:AL:257:ILE:HG22	9:AL:301:BCL:HED2	1.82	0.60
3:AM:4:TYR:HD1	3:AM:4:TYR:O	1.83	0.60
3:AM:128:LEU:O	3:AM:131:VAL:HB	2.01	0.60
3:AM:163:ILE:O	3:AM:167:MET:HB2	2.01	0.60
3:AM:204:LEU:HD21	4:AH:19:PHE:CE1	2.36	0.60
4:AH:45:ARG:HD3	4:AH:97:GLY:H	1.66	0.60
9:AA:101:BCL:HMB3	9:A0:102:BCL:C1B	2.31	0.60
9:AF:102:BCL:ND	9:AG:101:BCL:CMD	2.64	0.60
6:AG:45:TRP:CH2	9:AG:101:BCL:H2C	2.36	0.60
5:AI:31:LEU:HD21	14:AJ:102:CRT:H32	1.82	0.60
5:AQ:44:LEU:HD22	6:AR:43:ARG:CD	2.31	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:AS:32:GLY:HA3	9:AS:103:BCL:O1A	2.00	0.60
5:AW:24:ILE:HG21	14:AX:102:CRT:H22	1.83	0.60
9:A3:103:BCL:ND	9:A3:104:BCL:HMD3	2.15	0.60
2:BL:10:TYR:HD1	4:BH:112:GLY:HA2	1.65	0.60
2:BL:235:ALA:HA	11:BL:304:UQ8:H3MB	1.83	0.60
3:BM:63:PHE:CE2	3:BM:124:LEU:HD12	2.36	0.60
3:BM:97:PRO:HB2	3:BM:171:TRP:O	2.01	0.60
6:BJ:14:ALA:O	6:BJ:18:HIS:HB2	2.01	0.60
9:BK:102:BCL:O1D	9:BK:102:BCL:H2A	2.01	0.60
9:BU:102:BCL:H2A	9:BU:102:BCL:O1D	2.01	0.60
5:BW:46:TRP:CH2	9:BW:102:BCL:HBC3	2.36	0.60
14:BW:103:CRT:H9	6:BZ:17:PHE:CE1	2.36	0.60
6:B8:17:PHE:CD1	6:B8:20:ILE:HG21	2.36	0.60
2:AL:52:TRP:NE1	5:A9:41:SER:OG	2.34	0.60
2:AL:82:TYR:HB3	2:AL:85:ARG:HG3	1.83	0.60
2:AL:268:TRP:O	2:AL:269:PRO:C	2.44	0.60
4:AH:182:LEU:N	4:AH:182:LEU:HD12	2.16	0.60
5:AA:14:ILE:HG13	5:AA:15:LEU:HD22	1.83	0.60
5:AA:36:HIS:ND1	14:A0:101:CRT:H392	2.16	0.60
5:AA:36:HIS:O	5:AA:40:LEU:HB3	2.01	0.60
5:AA:46:TRP:HA	6:AB:43:ARG:NH1	2.15	0.60
5:AD:52:PRO:C	5:AD:54:SER:H	2.09	0.60
6:AE:29:PHE:HE1	9:AE:101:BCL:H11	1.64	0.60
5:AF:43:ASP:OD1	5:AF:44:LEU:CD2	2.50	0.60
9:AF:102:BCL:ND	9:AG:101:BCL:HMD3	2.17	0.60
9:AK:102:BCL:C2D	9:AN:101:BCL:C2D	2.80	0.60
5:AQ:51:ILE:CG1	5:AQ:52:PRO:HA	2.25	0.60
5:AS:36:HIS:CE1	9:AT:101:BCL:CMD	2.84	0.60
1:BC:126:VAL:HG12	1:BC:287:LEU:HD22	1.84	0.60
1:BC:316:LYS:O	1:BC:317:PRO:C	2.43	0.60
2:BL:48:LEU:HA	2:BL:51:VAL:HG23	1.82	0.60
3:BM:240:HIS:CE1	4:BH:69:LEU:HD21	2.36	0.60
4:BH:106:PRO:HA	4:BH:109:SER:HB3	1.84	0.60
9:BA:101:BCL:C1D	9:BB:101:BCL:HMD3	2.31	0.60
5:BO:49:ASP:CG	5:BO:50:ASN:H	2.09	0.60
5:BS:30:VAL:HG13	5:BS:31:LEU:N	2.16	0.60
9:BY:102:BCL:C1D	9:BZ:101:BCL:CMD	2.79	0.60
6:BZ:18:HIS:NE2	6:BZ:22:MET:HE2	2.16	0.60
5:B3:18:ARG:HA	5:B3:21:LEU:HB3	1.83	0.60
6:B4:21:PHE:HZ	9:B5:102:BCL:H203	1.66	0.60
5:B5:28:GLN:NE2	9:B6:101:BCL:HED1	2.16	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:B9:5:ASN:HA	5:B9:8:LEU:CD1	2.31	0.60
2:AL:12:VAL:CG2	2:AL:13:ARG:N	2.64	0.60
2:AL:40:PHE:O	2:AL:43:THR:HB	2.00	0.60
2:AL:196:LEU:HD23	3:AM:216:PHE:CB	2.27	0.60
4:AH:182:LEU:HD13	4:AH:195:LEU:CG	2.31	0.60
6:AE:13:GLU:H	6:AE:13:GLU:CD	2.09	0.60
5:AS:30:VAL:CG2	15:AS:101:PEF:C40	2.79	0.60
5:AS:36:HIS:CE1	9:AT:101:BCL:HMD3	2.36	0.60
6:AV:27:ALA:C	6:AV:31:LEU:HG	2.24	0.60
6:A2:20:ILE:HD13	6:A2:20:ILE:O	2.00	0.60
6:A2:32:VAL:HG11	9:A2:101:BCL:HBA2	1.84	0.60
5:A7:43:ASP:HB2	5:A9:47:LEU:CD1	2.25	0.60
5:A7:44:LEU:CD2	5:A7:46:TRP:CE3	2.78	0.60
6:A0:10:THR:H	6:A0:13:GLU:CG	2.15	0.60
1:BC:148:THR:HG23	1:BC:322:GLN:HA	1.82	0.60
2:BL:82:TYR:HB3	2:BL:85:ARG:HG3	1.83	0.60
3:BM:59:LEU:CG	3:BM:128:LEU:HD21	2.31	0.60
4:BH:35:LYS:O	4:BH:36:ARG:C	2.43	0.60
9:BE:101:BCL:HMC1	9:BF:102:BCL:HBB1	1.82	0.60
5:BK:26:ALA:O	5:BK:29:ILE:HG22	2.01	0.60
5:BO:24:ILE:O	5:BO:27:PHE:HB3	2.00	0.60
5:BU:51:ILE:HB	5:BU:52:PRO:CA	2.29	0.60
5:BW:12:TRP:HZ2	6:BX:21:PHE:CG	2.18	0.60
5:BY:38:ILE:HD12	5:BY:39:VAL:N	2.16	0.60
5:B1:14:ILE:CD1	5:B1:15:LEU:HG	2.31	0.60
14:B1:103:CRT:C2M	5:B5:36:HIS:HB3	2.32	0.60
2:AL:221:GLU:C	2:AL:223:THR:H	2.09	0.60
3:AM:56:THR:HG21	3:AM:131:VAL:HG11	1.84	0.60
4:AH:159:LEU:HB3	4:AH:212:ASP:HA	1.83	0.60
6:AE:42:TYR:CD2	6:AE:43:ARG:HG3	2.36	0.60
5:AI:44:LEU:HA	5:AK:56:GLN:CB	2.31	0.60
6:AV:20:ILE:HD13	6:AV:20:ILE:O	2.01	0.60
9:AX:101:BCL:CMC	5:AY:47:LEU:HD21	2.32	0.60
5:AY:8:LEU:HD12	6:AZ:22:MET:CE	2.31	0.60
5:AY:21:LEU:O	5:AY:25:VAL:HG23	2.02	0.60
14:A5:103:CRT:H11	5:A7:21:LEU:HD13	1.83	0.60
2:BL:231:TYR:OH	2:BL:233:ILE:HA	2.01	0.60
3:BM:148:TRP:HA	3:BM:148:TRP:HE3	1.65	0.60
5:BA:33:LEU:H	5:BA:33:LEU:HD12	1.65	0.60
5:BQ:44:LEU:HD12	5:BQ:44:LEU:O	2.01	0.60
6:BT:10:THR:HG22	6:BT:11:ASP:N	2.12	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:BT:101:BCL:C4A	9:BU:102:BCL:HMB2	2.31	0.60
5:B3:5:ASN:HA	5:B3:8:LEU:CG	2.31	0.60
1:AC:185:TYR:O	3:AM:89:HIS:ND1	2.21	0.60
1:AC:212:ILE:O	1:AC:222:ASN:ND2	2.34	0.60
1:AC:276:VAL:HG13	1:AC:277:ARG:N	2.17	0.60
3:AM:200:PRO:O	3:AM:203:MET:HG2	2.01	0.60
3:AM:229:PHE:O	3:AM:244:ALA:HB2	2.00	0.60
5:AA:47:LEU:HD12	5:A9:43:ASP:CB	2.31	0.60
14:AA:102:CRT:H11	6:AE:17:PHE:CE1	2.36	0.60
6:AB:20:ILE:O	6:AB:20:ILE:HD13	2.02	0.60
5:AF:40:LEU:CD2	5:AF:45:ASN:HA	2.31	0.60
5:AI:18:ARG:HH11	5:AI:18:ARG:CG	2.14	0.60
14:AS:104:CRT:H393	5:AW:36:HIS:CB	2.31	0.60
5:AU:5:ASN:HA	5:AU:8:LEU:HG	1.82	0.60
6:AV:46:LEU:HB3	6:AX:42:TYR:HH	1.66	0.60
5:A1:40:LEU:HB2	5:A1:46:TRP:CH2	2.35	0.60
5:A7:7:ASN:HD22	5:A7:7:ASN:H	1.49	0.60
5:A7:10:LYS:O	14:A0:101:CRT:H82	2.02	0.60
1:BC:266:ARG:O	1:BC:269:ALA:N	2.29	0.60
2:BL:82:TYR:HA	2:BL:85:ARG:HE	1.64	0.60
3:BM:208:PHE:C	3:BM:210:TYR:H	2.08	0.60
5:BD:8:LEU:C	5:BD:10:LYS:N	2.56	0.60
5:BI:55:TYR:CD1	5:BI:56:GLN:HG3	2.37	0.60
9:BN:101:BCL:C4B	9:BO:102:BCL:HBB3	2.32	0.60
5:BU:27:PHE:CE2	5:BW:29:ILE:HD11	2.36	0.60
14:BV:102:CRT:H2M3	5:BW:37:MET:HB2	1.84	0.60
5:B1:10:LYS:HB3	14:B1:103:CRT:C5	2.30	0.60
5:B7:32:GLY:HA3	9:B7:103:BCL:O1A	2.02	0.60
9:B9:102:BCL:HBA2	9:B0:102:BCL:OBD	2.01	0.60
2:AL:196:LEU:C	2:AL:198:MET:H	2.10	0.60
2:AL:257:ILE:HG13	2:AL:257:ILE:O	2.00	0.60
3:AM:201:PHE:CZ	4:AH:15:THR:HG22	2.36	0.60
5:AF:35:ILE:HA	5:AF:38:ILE:HG22	1.84	0.60
5:AS:11:ILE:CA	14:AS:104:CRT:H82	2.27	0.60
5:AU:13:LEU:HD22	6:AV:9:LEU:HB2	1.83	0.60
5:AY:44:LEU:HD22	6:AZ:43:ARG:CD	2.30	0.60
5:A1:27:PHE:CE2	5:A3:29:ILE:HD11	2.34	0.60
6:A8:29:PHE:CZ	9:A8:101:BCL:H101	2.36	0.60
6:A8:31:LEU:O	6:A8:34:ILE:HG22	2.00	0.60
5:A9:43:ASP:OD1	5:A9:44:LEU:HD12	2.02	0.60
6:A0:17:PHE:CE1	6:A0:21:PHE:CD2	2.90	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BC:225:SER:H	1:BC:228:GLN:NE2	2.00	0.60
2:BL:172:GLN:HE21	2:BL:172:GLN:HA	1.67	0.60
2:BL:196:LEU:CD1	3:BM:216:PHE:HB2	2.30	0.60
2:BL:211:LYS:HD3	2:BL:212:GLY:N	2.15	0.60
3:BM:4:TYR:HE2	3:BM:10:ALA:HB2	1.66	0.60
3:BM:242:GLY:O	3:BM:246:GLU:HB2	2.01	0.60
4:BH:5:ILE:HG13	5:BF:40:LEU:HD21	1.81	0.60
4:BH:215:LYS:H	4:BH:218:HIS:HD2	1.47	0.60
5:BA:36:HIS:HB3	14:B0:101:CRT:H391	1.78	0.60
14:BA:102:CRT:C2	6:BE:16:GLU:HG3	2.32	0.60
9:BO:102:BCL:CBC	9:BP:101:BCL:HAC1	2.31	0.60
6:B2:17:PHE:HD1	14:B2:102:CRT:H6	1.62	0.60
1:AC:265:LYS:HD2	1:AC:265:LYS:H	1.66	0.60
2:AL:148:MET:HE3	2:AL:153:HIS:CE1	2.37	0.60
2:AL:230:GLY:CA	3:AM:51:ILE:HB	2.30	0.60
3:AM:179:ILE:O	3:AM:183:LEU:HB2	2.01	0.60
3:AM:193:TYR:O	3:AM:194:GLY:C	2.44	0.60
4:AH:45:ARG:O	4:AH:96:PRO:HB3	2.01	0.60
6:AB:33:VAL:O	6:AB:37:LEU:HB2	2.02	0.60
6:AE:21:PHE:CD1	6:AE:21:PHE:C	2.80	0.60
6:AR:30:GLY:O	6:AR:34:ILE:HG22	2.02	0.60
5:AS:10:LYS:C	14:AS:104:CRT:H33	2.26	0.60
14:AW:102:CRT:H291	9:A1:102:BCL:O2A	2.02	0.60
6:A0:38:LEU:HD23	6:A0:38:LEU:C	2.27	0.60
1:BC:148:THR:OG1	1:BC:322:GLN:HG2	2.02	0.60
2:BL:194:LEU:HD23	2:BL:194:LEU:C	2.26	0.60
3:BM:59:LEU:HD23	3:BM:128:LEU:HD21	1.84	0.60
3:BM:222:THR:O	3:BM:225:SER:OG	2.20	0.60
6:BB:22:MET:CG	6:BB:26:TYR:OH	2.50	0.60
14:BG:102:CRT:H2M1	5:BI:33:LEU:O	2.02	0.60
5:BI:55:TYR:HD1	5:BI:56:GLN:H	1.48	0.60
9:BK:102:BCL:C1D	9:BN:101:BCL:CMD	2.79	0.60
5:BO:44:LEU:HD12	5:BO:46:TRP:N	2.16	0.60
6:BT:42:TYR:CE2	6:BT:43:ARG:HG2	2.36	0.60
9:BW:102:BCL:C2D	9:BX:101:BCL:C2D	2.80	0.60
14:BW:103:CRT:H14	6:BZ:21:PHE:CD2	2.36	0.60
5:BY:44:LEU:HD22	6:BZ:43:ARG:HD2	1.84	0.60
1:AC:52:SER:O	1:AC:56:ASN:HB2	2.01	0.60
1:AC:80:GLN:HG3	1:AC:128:ARG:HH22	1.66	0.60
2:AL:117:CYS:HA	2:AL:122:ILE:HD11	1.84	0.60
2:AL:248:SER:O	2:AL:249:ALA:C	2.45	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:AM:12:GLN:HB2	4:AH:145:ALA:CB	2.32	0.60
3:AM:32:GLY:O	3:AM:34:PRO:HD3	2.01	0.60
3:AM:104:LEU:HD22	3:AM:104:LEU:N	2.17	0.60
3:AM:275:LEU:HD21	4:AH:19:PHE:CE2	2.37	0.60
4:AH:5:ILE:HD11	5:AF:47:LEU:HD12	1.84	0.60
4:AH:113:PRO:HB2	4:AH:249:TYR:CE2	2.36	0.60
5:AF:14:ILE:HD12	5:AI:21:LEU:HD22	1.84	0.60
5:AI:9:TYR:CD1	5:AI:9:TYR:C	2.80	0.60
6:AP:22:MET:HG3	6:AP:26:TYR:HE2	1.67	0.60
9:AQ:102:BCL:C3D	6:AR:35:ALA:HB1	2.32	0.60
5:AW:26:ALA:HA	5:AW:29:ILE:CG2	2.32	0.60
5:AW:27:PHE:HE1	14:AX:102:CRT:H30	1.65	0.60
14:AW:102:CRT:H2M1	5:A1:36:HIS:HB3	1.84	0.60
9:AX:101:BCL:CHC	9:AY:102:BCL:HBB3	2.31	0.60
5:AY:4:MET:O	5:AY:8:LEU:N	2.35	0.60
6:AZ:34:ILE:HD13	6:AZ:34:ILE:C	2.26	0.60
5:A1:30:VAL:HA	5:A1:33:LEU:HG	1.83	0.60
14:A1:103:CRT:H2M3	5:A5:36:HIS:HB3	1.84	0.60
5:A7:50:ASN:CG	5:A7:51:ILE:N	2.60	0.60
6:A0:30:GLY:O	6:A0:34:ILE:HG22	2.02	0.60
1:BC:71:LYS:HD2	1:BC:71:LYS:N	2.16	0.60
1:BC:170:PRO:HG2	1:BC:171:GLY:H	1.67	0.60
2:BL:185:ALA:HB2	2:BL:252:TRP:HB3	1.82	0.60
3:BM:73:PHE:HA	14:BM:406:CRT:H1M1	1.84	0.60
3:BM:84:PHE:HA	5:BW:37:MET:CE	2.32	0.60
3:BM:98:PRO:HD2	3:BM:171:TRP:HB3	1.83	0.60
3:BM:163:ILE:O	3:BM:167:MET:HB2	2.02	0.60
3:BM:229:PHE:HD1	3:BM:229:PHE:H	1.49	0.60
4:BH:22:PHE:C	4:BH:22:PHE:CD1	2.79	0.60
4:BH:172:VAL:HG23	4:BH:173:ASP:N	2.16	0.60
5:BD:51:ILE:CG2	5:BD:52:PRO:HA	2.32	0.60
5:BK:26:ALA:HA	5:BK:29:ILE:HG22	1.83	0.60
5:BS:31:LEU:HD21	14:BS:103:CRT:H32	1.82	0.60
6:B6:40:TRP:HZ3	6:B6:45:TRP:H	1.50	0.60
1:AC:304:ARG:HH11	1:AC:304:ARG:HG3	1.65	0.60
2:AL:75:ILE:HD12	2:AL:94:LEU:HD22	1.81	0.60
2:AL:218:SER:C	2:AL:220:HIS:H	2.09	0.60
2:AL:276:LEU:H	2:AL:276:LEU:CD2	2.13	0.60
3:AM:34:PRO:HG2	3:AM:50:PRO:HD3	1.82	0.60
4:AH:69:LEU:HD11	4:AH:76:VAL:HG23	1.83	0.60
4:AH:121:LYS:NZ	4:BH:73:GLY:HA2	2.16	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:AF:40:LEU:HD22	5:AF:45:ASN:HA	1.82	0.60
14:AP:102:CRT:O2	5:AQ:33:LEU:HD12	2.01	0.60
5:AQ:36:HIS:CE1	9:AR:101:BCL:HMD3	2.37	0.60
6:AT:33:VAL:O	6:AT:37:LEU:HG	2.01	0.60
9:AW:101:BCL:O1A	6:AX:28:TRP:CH2	2.54	0.60
5:A1:4:MET:HE2	6:A4:27:ALA:CB	2.31	0.60
5:A3:14:ILE:O	5:A5:18:ARG:HD3	2.02	0.60
6:A0:10:THR:HG22	6:A0:11:ASP:N	2.14	0.60
1:BC:237:MET:SD	2:BL:174:LEU:HD23	2.42	0.60
3:BM:200:PRO:HA	3:BM:203:MET:CG	2.32	0.60
5:BD:43:ASP:HB2	5:BF:47:LEU:HD22	1.84	0.60
5:BD:46:TRP:CZ3	9:BD:102:BCL:CBC	2.84	0.60
5:BI:7:ASN:O	5:BI:10:LYS:HD3	2.02	0.60
5:BO:45:ASN:HB3	5:BO:48:ASP:OD1	2.01	0.60
5:BQ:50:ASN:ND2	5:BS:56:GLN:HA	2.10	0.60
5:BU:12:TRP:HA	5:BU:12:TRP:HE3	1.63	0.60
9:B1:102:BCL:HAC2	9:B2:101:BCL:CBC	2.32	0.60
9:B3:102:BCL:OBD	6:B4:32:VAL:HG13	2.02	0.60
5:B5:36:HIS:NE2	9:B6:101:BCL:HMD3	2.17	0.60
6:B0:10:THR:H	6:B0:13:GLU:CG	2.14	0.60
6:B0:40:TRP:HZ3	6:B0:45:TRP:N	1.99	0.60
2:AL:44:LEU:HB2	5:A9:30:VAL:HG11	1.83	0.60
3:AM:34:PRO:CG	3:AM:50:PRO:HD3	2.31	0.60
3:AM:261:THR:C	3:AM:263:GLU:H	2.10	0.60
4:AH:113:PRO:HG2	4:AH:248:LEU:HD22	1.84	0.60
5:AI:36:HIS:O	5:AI:40:LEU:HD13	2.02	0.60
5:AI:46:TRP:HA	5:AI:49:ASP:OD2	2.02	0.60
9:AN:101:BCL:H2	9:AN:101:BCL:H72	1.84	0.60
5:AO:31:LEU:HD21	14:AP:102:CRT:H32	1.83	0.60
6:AR:45:TRP:CD1	6:AR:46:LEU:N	2.69	0.60
5:AW:50:ASN:ND2	5:AW:51:ILE:HG12	2.17	0.60
5:AY:8:LEU:HD23	6:A2:20:ILE:HD11	1.84	0.60
6:A0:45:TRP:HD1	6:A0:46:LEU:H	1.49	0.60
2:BL:230:GLY:HA2	3:BM:51:ILE:CB	2.26	0.60
3:BM:299:VAL:CB	3:BM:304:ALA:HB3	2.29	0.60
5:BF:32:GLY:N	9:BG:101:BCL:HED2	2.17	0.60
6:BX:45:TRP:CD2	9:BX:101:BCL:H2C	2.37	0.60
5:BY:50:ASN:HB3	5:B1:60:LYS:CA	2.32	0.60
5:BY:51:ILE:HA	5:BY:52:PRO:C	2.27	0.60
5:B5:28:GLN:HB3	9:B5:102:BCL:C2	2.32	0.60
6:B0:38:LEU:HD23	6:B0:38:LEU:C	2.26	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AC:33:ILE:N	1:AC:33:ILE:HD12	2.17	0.59
1:AC:170:PRO:HG2	1:AC:171:GLY:N	2.17	0.59
3:AM:102:TYR:CD1	3:AM:102:TYR:N	2.69	0.59
4:AH:13:GLN:O	4:AH:16:ILE:HG22	2.02	0.59
6:AB:23:GLN:O	5:A9:4:MET:SD	2.61	0.59
6:AJ:17:PHE:CD2	14:AJ:102:CRT:H41	2.37	0.59
9:AJ:101:BCL:C1B	9:AK:102:BCL:HMB3	2.32	0.59
6:AP:20:ILE:CG2	6:AP:21:PHE:N	2.65	0.59
6:AR:34:ILE:HD13	6:AR:34:ILE:C	2.27	0.59
14:AS:104:CRT:C6	6:AV:20:ILE:HG21	2.32	0.59
6:AX:30:GLY:HA2	6:AX:33:VAL:HG12	1.84	0.59
5:A7:7:ASN:CB	5:A7:10:LYS:NZ	2.65	0.59
5:A7:44:LEU:HD22	5:A7:46:TRP:CE3	2.37	0.59
2:BL:146:LEU:C	2:BL:148:MET:H	2.10	0.59
3:BM:131:VAL:C	3:BM:133:THR:N	2.58	0.59
3:BM:131:VAL:O	3:BM:133:THR:N	2.35	0.59
5:BA:36:HIS:NE2	9:BB:101:BCL:CMD	2.65	0.59
5:BF:49:ASP:HB2	5:BI:56:GLN:OE1	2.02	0.59
5:BI:17:PRO:O	5:BI:21:LEU:CB	2.50	0.59
5:BO:36:HIS:CE1	9:BP:101:BCL:HMD2	2.37	0.59
5:BU:27:PHE:CD2	5:BW:29:ILE:HD11	2.37	0.59
5:B3:12:TRP:HE1	6:B4:18:HIS:HB2	1.66	0.59
5:B3:36:HIS:O	5:B3:40:LEU:HG	2.02	0.59
1:AC:32:GLN:CB	2:AL:80:LEU:HD12	2.28	0.59
1:AC:226:LEU:HD12	3:AM:192:ARG:HB2	1.84	0.59
2:AL:216:LYS:HD2	2:AL:220:HIS:NE2	2.18	0.59
2:AL:242:GLY:HA2	3:AM:216:PHE:CE2	2.37	0.59
2:AL:279:PRO:O	2:AL:280:LEU:HD23	2.03	0.59
3:AM:14:ARG:HH11	3:AM:14:ARG:HG3	1.66	0.59
4:AH:249:TYR:O	4:AH:251:THR:N	2.35	0.59
14:AB:102:CRT:H33	5:A9:10:LYS:HB3	0.71	0.59
9:AJ:101:BCL:CHB	9:AK:102:BCL:HMB3	2.32	0.59
9:AW:101:BCL:HBC2	9:AX:101:BCL:HHD	1.84	0.59
5:AY:36:HIS:CE1	9:AZ:101:BCL:CMD	2.86	0.59
5:A7:25:VAL:HA	9:A7:103:BCL:H52	1.83	0.59
14:A0:101:CRT:H35	9:A0:102:BCL:HMA3	1.84	0.59
1:BC:259:TRP:C	1:BC:261:GLN:H	2.10	0.59
4:BH:29:TYR:CD1	4:BH:29:TYR:C	2.79	0.59
5:BA:37:MET:O	5:BA:41:SER:HB2	2.01	0.59
5:BD:40:LEU:CD1	5:BD:47:LEU:HD23	2.32	0.59
9:BD:102:BCL:HBA2	9:BE:101:BCL:OBD	2.01	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:BQ:31:LEU:CD2	9:BQ:104:BCL:HED3	2.32	0.59
5:BY:50:ASN:ND2	6:BZ:43:ARG:HH12	1.99	0.59
5:B5:16:ASP:HB3	5:B5:17:PRO:HD2	1.83	0.59
5:B7:37:MET:HB2	14:B7:102:CRT:H2M1	1.83	0.59
1:AC:71:LYS:HD3	1:AC:74:GLU:OE1	2.02	0.59
1:AC:243:LEU:H	1:AC:243:LEU:CD1	2.16	0.59
2:AL:46:GLY:O	2:AL:50:ILE:HG22	2.02	0.59
3:AM:196:LEU:C	3:AM:198:TYR:H	2.10	0.59
15:AM:409:PEF:O1P	15:AM:409:PEF:H52	2.03	0.59
6:AB:29:PHE:HE1	9:AB:101:BCL:C1	2.07	0.59
6:BJ:34:ILE:HD13	6:BJ:35:ALA:N	2.16	0.59
9:BJ:101:BCL:HMB1	9:BJ:101:BCL:HBB3	1.84	0.59
5:BQ:5:ASN:HD22	6:BR:22:MET:HE3	1.66	0.59
6:BZ:29:PHE:N	6:BZ:29:PHE:CD1	2.70	0.59
5:B3:46:TRP:CZ3	9:B3:102:BCL:H2C	2.36	0.59
2:AL:170:GLY:HA3	9:AL:301:BCL:HBC2	1.84	0.59
3:AM:236:ASP:OD1	3:AM:237:GLN:N	2.35	0.59
4:AH:108:LEU:C	4:AH:110:GLY:H	2.08	0.59
5:AF:40:LEU:HD11	5:AF:47:LEU:HD12	1.85	0.59
5:AU:45:ASN:C	5:AU:49:ASP:HB3	2.26	0.59
6:A2:21:PHE:CD1	14:A2:102:CRT:C14	2.85	0.59
1:BC:176:SER:HA	5:BU:48:ASP:CG	2.28	0.59
3:BM:84:PHE:CZ	5:BU:38:ILE:HD12	2.38	0.59
3:BM:226:VAL:O	3:BM:226:VAL:HG13	2.02	0.59
5:BA:11:ILE:N	14:BA:102:CRT:H82	2.17	0.59
5:BF:21:LEU:O	5:BF:25:VAL:HG23	2.03	0.59
6:BG:30:GLY:O	6:BG:34:ILE:HG23	2.02	0.59
9:BI:102:BCL:HMB1	9:BI:102:BCL:HBB3	1.84	0.59
5:BU:12:TRP:CZ2	6:BV:17:PHE:CE2	2.91	0.59
14:BV:102:CRT:H392	5:BW:36:HIS:HB3	1.75	0.59
6:B2:46:LEU:HD22	6:B4:42:TYR:CE2	2.32	0.59
6:B4:20:ILE:HD13	6:B4:20:ILE:O	2.02	0.59
1:AC:236:MET:HA	1:AC:239:ILE:HD12	1.84	0.59
2:AL:137:TYR:O	2:AL:141:VAL:HG12	2.02	0.59
3:AM:172:ALA:C	3:AM:174:ALA:H	2.09	0.59
4:AH:135:PRO:HB3	4:AH:171:TRP:CE2	2.37	0.59
5:AF:36:HIS:CE1	9:AG:101:BCL:HMD3	2.37	0.59
6:AN:20:ILE:H	6:AN:20:ILE:CD1	2.14	0.59
6:AR:46:LEU:HD22	6:AT:42:TYR:CE2	2.37	0.59
5:AU:18:ARG:H	5:AU:18:ARG:CD	2.07	0.59
5:AU:49:ASP:CG	5:AU:50:ASN:N	2.59	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:AV:42:TYR:CE2	6:AV:43:ARG:HG3	2.37	0.59
5:AY:40:LEU:HD13	5:AY:46:TRP:CE2	2.37	0.59
9:AY:102:BCL:HBC2	9:AY:102:BCL:CHD	2.32	0.59
6:AZ:36:HIS:CE1	9:AZ:101:BCL:NA	2.70	0.59
6:A6:45:TRP:HD1	6:A6:46:LEU:H	1.45	0.59
2:BL:154:GLY:HA2	17:BL:403:HOH:O	2.00	0.59
2:BL:223:THR:HG21	3:BM:20:GLY:HA2	1.85	0.59
3:BM:98:PRO:HB3	3:BM:107:PRO:HB3	1.83	0.59
3:BM:156:PHE:CZ	9:BM:402:BCL:HBD	2.27	0.59
3:BM:208:PHE:C	3:BM:210:TYR:N	2.59	0.59
4:BH:69:LEU:CD1	4:BH:76:VAL:HG23	2.32	0.59
4:BH:170:VAL:HA	4:BH:182:LEU:HA	1.84	0.59
9:BF:102:BCL:HBC2	9:BG:101:BCL:HHD	1.85	0.59
9:BG:101:BCL:HHB	9:BI:102:BCL:HMA1	1.84	0.59
5:BI:27:PHE:CE2	5:BK:29:ILE:HD11	2.37	0.59
6:BJ:46:LEU:HB3	6:BN:42:TYR:CZ	2.38	0.59
6:BR:31:LEU:O	6:BR:34:ILE:HG13	2.03	0.59
5:BY:42:THR:O	5:BY:43:ASP:C	2.46	0.59
6:B2:21:PHE:HA	14:B2:102:CRT:C12	2.32	0.59
5:B3:33:LEU:O	5:B3:37:MET:HG2	2.03	0.59
5:B7:33:LEU:H	5:B7:33:LEU:HD12	1.67	0.59
9:B7:103:BCL:C3D	9:B8:101:BCL:C3D	2.80	0.59
2:AL:71:TRP:CD1	3:AM:303:MET:HG2	2.36	0.59
3:AM:35:ILE:HG22	3:AM:36:PHE:N	2.18	0.59
4:AH:55:VAL:CG1	5:AD:19:ARG:HD3	2.32	0.59
6:A2:25:MET:CE	9:A3:103:BCL:H171	2.27	0.59
14:A2:102:CRT:C31	9:A3:103:BCL:H3A	2.32	0.59
2:BL:206:VAL:CG1	3:BM:142:MET:HE1	2.29	0.59
2:BL:248:SER:O	2:BL:249:ALA:C	2.45	0.59
3:BM:260:VAL:HG23	3:BM:261:THR:N	2.16	0.59
4:BH:185:GLU:HA	4:BH:191:LYS:O	2.01	0.59
5:BA:47:LEU:HB3	5:B9:43:ASP:CA	2.32	0.59
5:BI:8:LEU:HB3	6:BJ:18:HIS:CE1	2.38	0.59
5:BI:19:ARG:O	5:BI:23:SER:HB3	2.03	0.59
9:BI:102:BCL:HBA2	9:BJ:101:BCL:OBD	2.03	0.59
6:BN:30:GLY:O	6:BN:33:VAL:HG12	2.02	0.59
5:BS:36:HIS:O	5:BS:40:LEU:N	2.31	0.59
6:BV:17:PHE:CD1	14:BV:102:CRT:C6	2.68	0.59
6:BX:34:ILE:HD13	6:BX:34:ILE:O	2.02	0.59
6:BX:46:LEU:HD22	6:BZ:42:TYR:OH	2.02	0.59
6:B2:42:TYR:CD1	6:B2:43:ARG:HG3	2.38	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:B4:18:HIS:C	6:B4:18:HIS:CD2	2.81	0.59
6:B4:29:PHE:CZ	9:B4:101:BCL:H101	2.36	0.59
5:B7:36:HIS:HB3	14:B7:102:CRT:C39	2.31	0.59
9:B7:103:BCL:C1D	9:B8:101:BCL:CMD	2.80	0.59
6:B0:33:VAL:CG2	9:B0:102:BCL:H143	2.31	0.59
2:AL:171:TYR:C	2:AL:173:PHE:N	2.60	0.59
3:AM:286:LEU:HD23	3:AM:290:VAL:HG21	1.84	0.59
14:AA:102:CRT:H32	5:AD:31:LEU:HD21	1.84	0.59
5:AF:9:TYR:CZ	5:AF:10:LYS:HD3	2.38	0.59
9:AJ:101:BCL:CHC	9:AK:102:BCL:CBB	2.80	0.59
5:AK:36:HIS:O	5:AK:40:LEU:HB2	2.02	0.59
5:AO:38:ILE:HG13	5:AO:39:VAL:N	2.16	0.59
5:AS:4:MET:HB2	5:AS:8:LEU:HD11	1.84	0.59
5:AY:44:LEU:HD13	6:AZ:43:ARG:CD	2.32	0.59
5:A1:44:LEU:HD23	5:A1:44:LEU:H	1.65	0.59
6:A4:10:THR:HG22	6:A4:11:ASP:N	2.17	0.59
3:BM:234:GLU:O	3:BM:235:ILE:C	2.45	0.59
3:BM:287:SER:OG	3:BM:294:TRP:NE1	2.36	0.59
4:BH:95:ALA:HB2	5:B9:16:ASP:OD2	2.02	0.59
5:BA:29:ILE:HD12	9:BA:101:BCL:H11	1.83	0.59
9:BD:102:BCL:C2D	9:BE:101:BCL:C2D	2.81	0.59
5:BF:4:MET:O	5:BF:8:LEU:HG	2.02	0.59
6:BN:17:PHE:O	6:BN:21:PHE:HB3	2.03	0.59
5:BQ:31:LEU:HD23	9:BQ:104:BCL:HED3	1.85	0.59
5:BQ:53:VAL:O	5:BQ:54:SER:C	2.45	0.59
5:BS:34:LEU:O	5:BS:38:ILE:HG22	2.02	0.59
9:BV:101:BCL:HMA1	9:BW:102:BCL:CHB	2.28	0.59
6:B2:46:LEU:HB3	6:B4:42:TYR:OH	2.03	0.59
5:B9:16:ASP:CG	5:B9:17:PRO:HD2	2.27	0.59
1:AC:40:MET:HE1	1:AC:252:ASN:ND2	2.18	0.59
1:AC:111:HIS:HE1	1:AC:124:LYS:HE2	1.68	0.59
2:AL:145:PRO:HB3	2:AL:150:ALA:O	2.02	0.59
4:AH:151:PRO:O	4:AH:154:MET:HG3	2.01	0.59
5:AI:52:PRO:HG2	5:AI:55:TYR:HE2	1.68	0.59
5:AO:27:PHE:HE2	5:AQ:29:ILE:CD1	2.16	0.59
5:AU:20:VAL:HG11	9:AW:101:BCL:H202	1.84	0.59
5:AW:8:LEU:O	5:AW:11:ILE:HG13	2.02	0.59
5:A5:24:ILE:CG1	9:A7:103:BCL:H201	2.32	0.59
6:A8:34:ILE:HD13	6:A8:34:ILE:C	2.27	0.59
1:BC:96:ALA:O	1:BC:98:THR:N	2.31	0.59
1:BC:107:CYS:O	1:BC:109:TYR:N	2.36	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:BM:80:HIS:O	5:BU:41:SER:HB2	2.02	0.59
5:BU:30:VAL:HG13	5:BU:31:LEU:N	2.18	0.59
6:B6:29:PHE:CD1	9:B6:101:BCL:H11	2.38	0.59
5:B7:35:ILE:O	5:B7:38:ILE:HG22	2.02	0.59
6:B0:10:THR:HG22	6:B0:11:ASP:N	2.15	0.59
2:AL:70:LEU:O	2:AL:159:ILE:HB	2.03	0.59
2:AL:199:HIS:O	2:AL:201:SER:N	2.35	0.59
3:AM:25:LYS:HG2	5:AO:16:ASP:OD1	2.02	0.59
3:AM:176:PRO:HD3	3:AM:185:TRP:HD1	1.66	0.59
3:AM:211:GLY:O	3:AM:214:LEU:HB3	2.02	0.59
4:AH:108:LEU:C	4:AH:110:GLY:N	2.59	0.59
4:AH:235:GLU:HA	4:AH:238:LYS:CB	2.31	0.59
5:AA:22:VAL:HA	5:AA:25:VAL:CG2	2.32	0.59
9:AA:101:BCL:HMD1	6:AB:35:ALA:HB1	1.85	0.59
6:AN:13:GLU:HA	6:AN:16:GLU:OE1	2.03	0.59
5:AO:43:ASP:HB2	5:AQ:47:LEU:HB3	1.85	0.59
6:AP:32:VAL:HG12	6:AP:36:HIS:HD1	1.66	0.59
6:AT:17:PHE:CE1	14:AT:102:CRT:H9	2.38	0.59
5:AY:35:ILE:HA	5:AY:38:ILE:HG13	1.85	0.59
5:A5:35:ILE:HA	5:A5:38:ILE:HG22	1.85	0.59
6:A8:45:TRP:HA	5:A9:52:PRO:HD3	1.85	0.59
5:A9:46:TRP:NE1	5:A9:47:LEU:HD22	2.18	0.59
9:A9:102:BCL:CMD	6:A0:36:HIS:HD2	2.15	0.59
3:BM:83:VAL:HG23	3:BM:84:PHE:HD1	1.68	0.59
3:BM:241:ARG:O	4:BH:119:ARG:HD3	2.03	0.59
5:BA:11:ILE:HD11	5:BD:21:LEU:HD21	1.85	0.59
5:BA:42:THR:HG22	5:BD:48:ASP:OD2	2.02	0.59
14:BA:102:CRT:H32	5:BD:31:LEU:HD21	1.84	0.59
6:BB:25:MET:HG2	6:BB:29:PHE:CE2	2.37	0.59
6:BB:40:TRP:HZ3	6:BB:45:TRP:N	2.00	0.59
5:BI:9:TYR:HA	6:BJ:18:HIS:ND1	2.17	0.59
5:BK:19:ARG:HG3	5:BK:20:VAL:H	1.68	0.59
5:BK:36:HIS:CE1	9:BN:101:BCL:CMD	2.83	0.59
6:BT:29:PHE:CE1	9:BT:101:BCL:H11	2.38	0.59
5:BU:11:ILE:HG12	14:BU:103:CRT:C8	2.32	0.59
6:BX:46:LEU:HB2	5:BY:52:PRO:CD	2.26	0.59
5:BY:45:ASN:O	5:BY:48:ASP:N	2.36	0.59
9:BY:102:BCL:CMD	6:BZ:36:HIS:HD2	2.15	0.59
5:B9:31:LEU:HD23	9:B0:102:BCL:HED3	1.84	0.59
1:AC:275:HIS:O	1:AC:279:ILE:HG13	2.03	0.59
3:AM:156:PHE:CE2	3:AM:280:ALA:HB1	2.38	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:AA:17:PRO:HG2	5:AA:18:ARG:HD2	1.84	0.59
6:AB:44:PRO:HD2	5:AD:55:TYR:HE2	1.67	0.59
5:AO:46:TRP:CD1	5:AO:47:LEU:HD13	2.38	0.59
5:AQ:42:THR:O	5:AS:48:ASP:HB3	2.02	0.59
6:AR:25:MET:HE2	14:AR:102:CRT:H242	1.84	0.59
6:AV:29:PHE:CD1	9:AV:102:BCL:H11	2.38	0.59
5:A1:5:ASN:HB3	5:A1:8:LEU:HD13	1.83	0.59
5:A3:14:ILE:CD1	6:A6:17:PHE:HE2	2.15	0.59
6:A6:40:TRP:HZ3	6:A6:45:TRP:N	2.00	0.59
5:A7:33:LEU:O	14:A7:102:CRT:H2M1	2.02	0.59
6:A0:17:PHE:CE1	6:A0:21:PHE:HD2	2.20	0.59
6:A0:32:VAL:HG12	6:A0:33:VAL:N	2.18	0.59
1:BC:270:TRP:C	1:BC:273:ILE:HD12	2.27	0.59
3:BM:7:ILE:HG22	3:BM:8:PHE:CG	2.38	0.59
3:BM:53:LEU:HG	3:BM:58:THR:HG23	1.85	0.59
3:BM:204:LEU:C	3:BM:206:ILE:N	2.60	0.59
5:BD:14:ILE:HD12	5:BD:14:ILE:N	2.18	0.59
6:BE:21:PHE:HZ	9:BF:102:BCL:H203	1.67	0.59
5:BI:55:TYR:HD1	5:BI:56:GLN:N	2.01	0.59
5:BO:29:ILE:HB	9:BO:102:BCL:C4	2.32	0.59
6:BX:18:HIS:CE1	6:BX:22:MET:HE2	2.38	0.59
5:B1:19:ARG:O	5:B1:23:SER:HB3	2.02	0.59
5:B3:12:TRP:HE1	6:B4:18:HIS:CB	2.16	0.59
9:B7:103:BCL:CAD	9:B8:101:BCL:CAD	2.80	0.59
9:B8:101:BCL:CMC	9:B9:102:BCL:CBB	2.80	0.59
5:B9:46:TRP:CH2	9:B9:102:BCL:HBC3	2.38	0.59
1:AC:53:ILE:HA	1:AC:319:TYR:CE1	2.37	0.58
2:AL:89:LEU:HG	2:AL:97:ILE:CD1	2.30	0.58
2:AL:184:LEU:HD22	2:AL:252:TRP:HE1	1.66	0.58
5:AA:50:ASN:CG	5:AA:51:ILE:HG12	2.28	0.58
5:AI:7:ASN:ND2	6:AN:20:ILE:HG13	2.16	0.58
5:AI:22:VAL:HA	5:AI:25:VAL:HG23	1.85	0.58
9:AJ:101:BCL:C1C	9:AK:102:BCL:HBB3	2.32	0.58
14:AS:104:CRT:H242	6:AV:25:MET:HE2	1.85	0.58
14:AW:102:CRT:H183	9:AY:102:BCL:H92	1.73	0.58
6:A2:17:PHE:O	6:A2:20:ILE:HG22	2.03	0.58
5:A7:35:ILE:CD1	9:A8:101:BCL:O1D	2.50	0.58
1:BC:36:ARG:HB3	2:BL:79:ASP:CG	2.28	0.58
2:BL:4:LEU:HD12	3:BM:250:LEU:CD1	2.31	0.58
3:BM:274:VAL:HG12	3:BM:278:ILE:HD11	1.84	0.58
5:BI:11:ILE:N	14:BN:102:CRT:H82	2.18	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:BN:31:LEU:HA	6:BN:34:ILE:CG2	2.33	0.58
14:BN:102:CRT:H2M3	5:BO:36:HIS:HB3	1.85	0.58
5:BO:14:ILE:HG23	5:BO:15:LEU:HG	1.83	0.58
5:BO:50:ASN:HD21	6:BP:43:ARG:NH2	2.02	0.58
14:BV:102:CRT:H31	9:BW:102:BCL:HBA1	1.85	0.58
5:B7:26:ALA:O	5:B7:29:ILE:HG22	2.03	0.58
5:B7:40:LEU:HD11	5:B7:47:LEU:HD23	1.85	0.58
6:B0:30:GLY:O	6:B0:34:ILE:HG22	2.02	0.58
1:AC:31:GLU:HB2	1:AC:42:ASN:HB3	1.84	0.58
2:AL:148:MET:HB3	2:AL:153:HIS:CE1	2.38	0.58
2:AL:192:ASN:C	2:AL:192:ASN:ND2	2.61	0.58
4:AH:67:PHE:N	4:AH:76:VAL:O	2.31	0.58
6:AB:20:ILE:HG21	14:AB:102:CRT:H83	1.85	0.58
5:AD:51:ILE:HA	5:AD:53:VAL:H	1.68	0.58
9:AG:101:BCL:CHC	9:AI:102:BCL:HBB3	2.33	0.58
6:AJ:46:LEU:HB3	6:AN:42:TYR:CZ	2.38	0.58
6:AN:10:THR:C	6:AN:13:GLU:OE2	2.46	0.58
6:AN:41:LEU:HD23	6:AN:42:TYR:CA	2.33	0.58
9:AS:103:BCL:C4D	9:AT:101:BCL:HMD3	2.33	0.58
5:AU:45:ASN:O	5:AU:49:ASP:N	2.35	0.58
6:AX:21:PHE:HA	14:AX:102:CRT:H14	1.85	0.58
9:AZ:101:BCL:H203	6:A2:38:LEU:HD11	1.83	0.58
6:A4:20:ILE:O	6:A4:20:ILE:HD13	2.03	0.58
5:A5:37:MET:HG2	5:A5:38:ILE:N	2.17	0.58
1:BC:251:HIS:CE1	1:BC:256:PHE:O	2.55	0.58
1:BC:266:ARG:HG3	7:BC:503:HEM:CMD	2.33	0.58
2:BL:116:ILE:HG22	2:BL:117:CYS:N	2.16	0.58
2:BL:196:LEU:HD11	3:BM:269:ALA:CB	2.29	0.58
3:BM:5:GLN:O	3:BM:7:ILE:N	2.36	0.58
3:BM:14:ARG:HH11	3:BM:14:ARG:HG3	1.67	0.58
3:BM:259:ASN:HD22	3:BM:259:ASN:N	2.01	0.58
5:BI:44:LEU:HD12	5:BI:46:TRP:HE3	1.68	0.58
5:BI:49:ASP:OD1	5:BI:50:ASN:N	2.30	0.58
5:BU:14:ILE:CD1	14:BU:103:CRT:H32A	2.33	0.58
6:B4:10:THR:HG22	6:B4:11:ASP:N	2.17	0.58
1:AC:47:ARG:HD3	5:A1:42:THR:HG22	1.85	0.58
1:AC:62:LEU:HD11	1:AC:95:VAL:HB	1.85	0.58
1:AC:127:SER:OG	7:AC:502:HEM:HMA1	2.04	0.58
1:AC:142:LYS:HA	1:AC:145:VAL:CG2	2.31	0.58
1:AC:176:SER:OG	5:AS:42:THR:HA	2.02	0.58
1:AC:316:LYS:HD2	7:AC:504:HEM:O2D	2.02	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:AM:208:PHE:C	3:AM:210:TYR:H	2.09	0.58
9:AM:401:BCL:HBC1	9:AM:402:BCL:HBD	1.84	0.58
9:AA:101:BCL:CMB	9:A0:102:BCL:C1B	2.81	0.58
9:AD:102:BCL:C1D	9:AE:101:BCL:HMD1	2.33	0.58
5:AQ:15:LEU:HA	5:AS:18:ARG:NH1	2.19	0.58
14:AS:104:CRT:H2M1	5:AW:37:MET:N	2.18	0.58
5:AU:38:ILE:HD11	5:AW:40:LEU:HD21	1.85	0.58
5:AU:43:ASP:OD2	5:AW:47:LEU:HA	2.03	0.58
6:AZ:46:LEU:OXT	5:A1:51:ILE:HG13	2.03	0.58
9:A7:103:BCL:HBC2	9:A8:101:BCL:HMD1	1.85	0.58
6:A0:32:VAL:HG21	9:A0:102:BCL:CGA	2.32	0.58
1:BC:135:ARG:HA	1:BC:330:LEU:O	2.03	0.58
1:BC:301:ASP:HB2	1:BC:302:PRO:HD2	1.85	0.58
2:BL:10:TYR:CD1	4:BH:112:GLY:HA2	2.38	0.58
2:BL:72:ARG:HA	3:BM:305:PRO:HB3	1.86	0.58
2:BL:150:ALA:O	2:BL:153:HIS:HB3	2.03	0.58
3:BM:66:VAL:HG11	3:BM:121:PHE:HD2	1.68	0.58
3:BM:151:ALA:O	3:BM:155:PHE:N	2.36	0.58
3:BM:170:SER:C	3:BM:172:ALA:N	2.58	0.58
3:BM:201:PHE:CZ	4:BH:16:ILE:HA	2.39	0.58
5:BF:42:THR:O	5:BI:48:ASP:HB3	2.03	0.58
5:BI:15:LEU:HB3	5:BI:20:VAL:HG21	1.84	0.58
5:BI:18:ARG:NH1	5:BI:18:ARG:CB	2.64	0.58
5:BO:55:TYR:N	5:BO:55:TYR:HD1	2.02	0.58
5:BU:38:ILE:HD11	5:BW:40:LEU:CD2	2.33	0.58
9:BY:102:BCL:HMD2	6:BZ:36:HIS:CD2	2.38	0.58
6:B8:34:ILE:HD13	6:B8:34:ILE:C	2.28	0.58
6:B0:21:PHE:CB	14:B0:101:CRT:C11	2.79	0.58
1:AC:134:VAL:O	1:AC:137:ALA:HB3	2.02	0.58
1:AC:267:THR:HG21	3:AM:314:VAL:CB	2.29	0.58
2:AL:168:ASN:O	2:AL:170:GLY:N	2.36	0.58
6:AN:22:MET:HG3	6:AN:26:TYR:CE2	2.35	0.58
5:AQ:43:ASP:CA	5:AS:47:LEU:HB3	2.33	0.58
5:AU:8:LEU:O	5:AU:11:ILE:HG13	2.02	0.58
5:AY:28:GLN:HG3	9:AY:102:BCL:H62	1.85	0.58
6:A6:19:ALA:O	6:A6:23:GLN:HG2	2.03	0.58
1:BC:53:ILE:HG12	1:BC:319:TYR:CE1	2.38	0.58
1:BC:70:PRO:HG2	1:BC:71:LYS:H	1.68	0.58
1:BC:325:LYS:HA	1:BC:331:TYR:OH	2.04	0.58
3:BM:70:ILE:CG2	3:BM:118:ALA:HB2	2.32	0.58
3:BM:246:GLU:O	3:BM:250:LEU:HB2	2.02	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:BH:32:ARG:HG3	4:BH:59:PRO:HB2	1.85	0.58
9:BA:101:BCL:CMB	9:B0:102:BCL:C1B	2.81	0.58
5:BO:26:ALA:HA	5:BO:29:ILE:HG22	1.85	0.58
14:BU:103:CRT:C2M	5:BY:37:MET:HG2	2.34	0.58
6:BV:42:TYR:CD2	6:BV:43:ARG:HG3	2.39	0.58
14:B5:103:CRT:C9	6:B8:17:PHE:HZ	2.16	0.58
9:B8:101:BCL:HMA3	9:B9:102:BCL:HMA2	1.85	0.58
2:AL:59:THR:HB	2:AL:63:SER:CB	2.31	0.58
2:AL:83:GLY:O	2:AL:150:ALA:HA	2.04	0.58
2:AL:223:THR:HA	2:AL:226:ARG:CB	2.31	0.58
3:AM:13:VAL:O	4:AH:177:PRO:HB2	2.03	0.58
3:AM:235:ILE:H	3:AM:235:ILE:CD1	2.12	0.58
4:AH:16:ILE:O	4:AH:16:ILE:HD13	2.03	0.58
4:AH:24:PHE:C	4:AH:27:ILE:HG22	2.29	0.58
6:AN:41:LEU:CD2	6:AN:42:TYR:N	2.60	0.58
5:AQ:54:SER:CB	5:AQ:57:ALA:HB3	2.33	0.58
9:AU:102:BCL:O1D	9:AU:102:BCL:C2A	2.49	0.58
6:AX:38:LEU:HD23	6:AX:38:LEU:C	2.28	0.58
6:A6:28:TRP:O	6:A6:31:LEU:N	2.36	0.58
1:BC:85:LEU:HD11	1:BC:329:GLY:CA	2.34	0.58
2:BL:10:TYR:O	2:BL:12:VAL:N	2.36	0.58
2:BL:129:ALA:HA	2:BL:247:LEU:HD11	1.83	0.58
3:BM:28:LEU:H	3:BM:28:LEU:HD12	1.67	0.58
3:BM:61:ILE:HG12	3:BM:129:TRP:HZ3	1.67	0.58
3:BM:207:ALA:O	3:BM:210:TYR:HB2	2.03	0.58
3:BM:242:GLY:C	4:BH:117:PRO:HG3	2.27	0.58
3:BM:267:ARG:HB3	4:BH:30:LEU:HD21	1.86	0.58
4:BH:137:ARG:HG2	4:BH:137:ARG:HH11	1.68	0.58
6:BB:45:TRP:O	6:BB:46:LEU:CB	2.50	0.58
14:BF:103:CRT:H25	5:BI:28:GLN:NE2	2.18	0.58
6:BP:34:ILE:HD12	6:BP:35:ALA:N	2.19	0.58
5:BQ:50:ASN:HA	5:BS:60:LYS:CB	2.33	0.58
6:BV:13:GLU:CD	6:BV:13:GLU:H	2.10	0.58
9:B3:102:BCL:HHC	9:B3:102:BCL:OBB	2.03	0.58
5:B7:37:MET:N	14:B7:102:CRT:H2M3	2.18	0.58
14:B7:102:CRT:H343	9:B7:103:BCL:HBA1	1.80	0.58
1:AC:33:ILE:HD12	1:AC:33:ILE:H	1.69	0.58
2:AL:35:PHE:CZ	2:AL:111:LEU:HD12	2.39	0.58
2:AL:172:GLN:HA	2:AL:172:GLN:HE21	1.68	0.58
5:AK:5:ASN:HA	5:AK:8:LEU:CD1	2.34	0.58
5:AO:50:ASN:HB3	5:AQ:55:TYR:HB2	1.85	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:A2:29:PHE:N	6:A2:29:PHE:CD1	2.69	0.58
6:A2:33:VAL:O	6:A2:37:LEU:HD23	2.03	0.58
6:A2:46:LEU:HB3	6:A4:42:TYR:OH	2.04	0.58
5:A7:16:ASP:O	5:A7:20:VAL:HG22	2.04	0.58
1:BC:166:TRP:CE3	1:BC:166:TRP:O	2.56	0.58
2:BL:36:GLY:HA2	2:BL:112:ARG:HD3	1.86	0.58
3:BM:32:GLY:O	3:BM:34:PRO:HD3	2.04	0.58
4:BH:37:GLU:C	4:BH:39:TYR:H	2.12	0.58
5:BA:27:PHE:CA	5:BA:30:VAL:HG12	2.34	0.58
6:BG:10:THR:HG22	6:BG:11:ASP:N	2.19	0.58
6:BV:17:PHE:HA	14:BV:102:CRT:H6	1.84	0.58
5:BW:20:VAL:O	5:BW:24:ILE:HG12	2.04	0.58
9:B1:102:BCL:HED2	6:B2:31:LEU:O	2.03	0.58
5:B3:56:GLN:NE2	5:B3:56:GLN:H	2.00	0.58
1:AC:90:PHE:O	1:AC:93:THR:HB	2.02	0.58
1:AC:122:TYR:CA	1:AC:125:VAL:HG23	2.33	0.58
2:AL:82:TYR:CA	2:AL:85:ARG:HE	2.16	0.58
2:AL:148:MET:CB	2:AL:153:HIS:ND1	2.65	0.58
3:AM:271:TRP:CD2	4:AH:26:LEU:HD21	2.38	0.58
4:AH:18:ALA:O	4:AH:19:PHE:C	2.45	0.58
6:AB:40:TRP:CE3	6:AB:40:TRP:HA	2.37	0.58
14:AB:102:CRT:H342	9:AD:102:BCL:CBA	2.34	0.58
5:AI:22:VAL:HA	5:AI:25:VAL:CG2	2.34	0.58
5:AS:33:LEU:O	5:AS:37:MET:HB2	2.04	0.58
5:AS:44:LEU:HD23	5:AS:44:LEU:H	1.68	0.58
14:AS:104:CRT:H6	6:AV:20:ILE:HG21	1.85	0.58
14:AT:102:CRT:H291	9:AU:102:BCL:O2A	2.03	0.58
5:AY:43:ASP:HB2	5:A1:47:LEU:HG	1.86	0.58
5:A1:10:LYS:NZ	6:A4:20:ILE:HB	2.19	0.58
5:A3:12:TRP:CD1	6:A4:18:HIS:HB2	2.37	0.58
5:A5:10:LYS:HB3	14:A5:103:CRT:O1	2.03	0.58
5:A5:44:LEU:C	5:A5:46:TRP:H	2.11	0.58
5:A7:11:ILE:CD1	5:A7:15:LEU:HD11	2.34	0.58
14:A7:102:CRT:H342	9:A7:103:BCL:HBA1	1.84	0.58
9:A8:101:BCL:HMA3	9:A9:102:BCL:CMA	2.34	0.58
1:BC:255:ALA:HB1	1:BC:258:ASP:HB3	1.86	0.58
3:BM:215:LEU:HA	3:BM:218:MET:HG3	1.84	0.58
5:BD:36:HIS:NE2	9:BE:101:BCL:HMD3	2.18	0.58
5:BI:55:TYR:HD1	5:BI:56:GLN:HG3	1.68	0.58
5:BO:44:LEU:HD12	5:BO:46:TRP:H	1.69	0.58
5:BO:55:TYR:N	5:BO:55:TYR:CD1	2.71	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:BO:102:BCL:CBD	9:BP:101:BCL:OBD	2.52	0.58
9:BO:102:BCL:CAC	9:BP:101:BCL:HAC1	2.34	0.58
9:BZ:101:BCL:CBB	9:B1:102:BCL:HMC1	2.31	0.58
9:BZ:101:BCL:HHC	9:BZ:101:BCL:OBB	2.03	0.58
6:B4:42:TYR:C	6:B4:42:TYR:HD1	2.12	0.58
6:B6:28:TRP:O	6:B6:31:LEU:N	2.36	0.58
6:B6:40:TRP:HA	6:B6:40:TRP:HE3	1.69	0.58
6:B8:22:MET:O	6:B8:26:TYR:HD2	1.87	0.58
5:B9:12:TRP:HZ2	6:B0:21:PHE:HE2	1.52	0.58
1:AC:316:LYS:O	1:AC:317:PRO:C	2.47	0.58
3:AM:40:LEU:CD1	3:AM:48:ILE:HD11	2.31	0.58
3:AM:120:LEU:HB2	14:AM:406:CRT:H35	1.84	0.58
3:AM:256:MET:HE1	13:AM:405:MQ8:H171	1.85	0.58
6:AE:23:GLN:HG3	6:AE:24:SER:N	2.14	0.58
5:AI:55:TYR:HD1	5:AI:56:GLN:H	1.51	0.58
5:AQ:43:ASP:HA	5:AS:47:LEU:O	2.03	0.58
6:AT:13:GLU:CD	6:AT:13:GLU:H	2.12	0.58
5:AU:20:VAL:CG1	9:AW:101:BCL:C20	2.81	0.58
6:AV:13:GLU:H	6:AV:13:GLU:CD	2.11	0.58
9:AY:102:BCL:HBC2	9:AZ:101:BCL:HMD1	1.84	0.58
2:BL:77:PRO:HB2	2:BL:152:GLY:HA2	1.85	0.58
2:BL:155:PHE:HB2	2:BL:156:PRO:HD2	1.86	0.58
2:BL:279:PRO:O	2:BL:280:LEU:HD23	2.03	0.58
3:BM:27:ASN:ND2	5:BO:19:ARG:HH11	2.00	0.58
3:BM:167:MET:HG2	3:BM:289:THR:HG21	1.85	0.58
3:BM:178:GLY:HA3	3:BM:181:PRO:HG2	1.85	0.58
3:BM:291:VAL:HG11	3:BM:297:TRP:HB2	1.85	0.58
4:BH:100:LEU:HB2	4:BH:111:PHE:CZ	2.39	0.58
4:BH:189:ASN:HB3	4:BH:191:LYS:HG3	1.84	0.58
5:BK:35:ILE:O	5:BK:38:ILE:HG22	2.04	0.58
14:B1:103:CRT:H2M3	5:B5:36:HIS:HB3	1.84	0.58
5:B3:46:TRP:CE3	9:B3:102:BCL:H2C	2.39	0.58
6:B6:19:ALA:O	6:B6:23:GLN:HG2	2.03	0.58
5:B9:32:GLY:CA	9:B0:102:BCL:HED2	2.33	0.58
1:AC:151:THR:HG21	1:AC:323:MET:HB2	1.86	0.58
1:AC:166:TRP:NE1	1:AC:305:VAL:C	2.59	0.58
1:AC:275:HIS:O	1:AC:278:ASP:HB3	2.04	0.58
3:AM:117:MET:SD	5:AQ:37:MET:HB3	2.43	0.58
3:AM:166:VAL:HG22	3:AM:171:TRP:CH2	2.38	0.58
14:AJ:102:CRT:H342	9:AK:102:BCL:CBA	2.18	0.58
9:AK:102:BCL:CAD	9:AN:101:BCL:C3D	2.82	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:AV:102:BCL:H192	9:AV:102:BCL:HBB1	1.84	0.58
5:AY:7:ASN:O	6:A2:20:ILE:HG13	2.03	0.58
5:A9:29:ILE:O	5:A9:33:LEU:HD13	2.04	0.58
1:BC:134:VAL:O	1:BC:137:ALA:HB3	2.03	0.58
1:BC:280:ASN:OD1	1:BC:304:ARG:CB	2.46	0.58
3:BM:215:LEU:C	3:BM:217:ALA:N	2.60	0.58
3:BM:300:LYS:O	4:BH:8:TYR:HB2	2.04	0.58
9:BA:101:BCL:HMB3	9:B0:102:BCL:CHB	2.33	0.58
6:BG:38:LEU:HA	6:BG:41:LEU:HD12	1.86	0.58
5:BK:28:GLN:HB2	9:BK:102:BCL:H43	1.85	0.58
5:BY:40:LEU:HD13	5:BY:46:TRP:CE2	2.39	0.58
5:B5:46:TRP:CZ3	9:B5:102:BCL:HBC3	2.39	0.58
6:B8:23:GLN:HG3	6:B8:24:SER:H	1.69	0.58
2:AL:178:TYR:CD2	2:AL:269:PRO:HG3	2.34	0.58
3:AM:274:VAL:HG12	3:AM:278:ILE:CD1	2.34	0.58
4:AH:69:LEU:CD1	4:AH:76:VAL:HG23	2.34	0.58
4:AH:202:PHE:HB3	4:AH:204:LYS:NZ	2.19	0.58
5:AA:44:LEU:HD12	5:AA:46:TRP:H	1.69	0.58
5:AF:42:THR:HG22	5:AI:47:LEU:HB3	1.85	0.58
14:AN:102:CRT:H2M3	5:AO:36:HIS:HB2	1.86	0.58
6:AP:46:LEU:CB	5:AQ:51:ILE:HG21	2.22	0.58
5:AQ:22:VAL:O	5:AQ:25:VAL:HG12	2.04	0.58
9:AZ:101:BCL:H203	6:A2:38:LEU:CD2	2.32	0.58
6:A4:18:HIS:C	6:A4:18:HIS:CD2	2.81	0.58
2:BL:184:LEU:CB	2:BL:252:TRP:HE1	2.17	0.58
2:BL:196:LEU:HD22	3:BM:216:PHE:HB2	1.84	0.58
3:BM:14:ARG:HG3	3:BM:14:ARG:NH1	2.19	0.58
3:BM:114:TRP:HZ3	3:BM:117:MET:HE2	1.69	0.58
3:BM:261:THR:O	3:BM:263:GLU:N	2.36	0.58
5:BA:18:ARG:CG	5:B9:14:ILE:HG23	2.33	0.58
5:BA:47:LEU:HB3	5:B9:43:ASP:HB2	1.86	0.58
6:BB:17:PHE:O	6:BB:20:ILE:HG22	2.03	0.58
5:BF:49:ASP:O	5:BI:56:GLN:HB3	2.03	0.58
6:BJ:23:GLN:CD	6:BJ:23:GLN:C	2.72	0.58
5:BO:7:ASN:O	6:BR:20:ILE:HD11	2.04	0.58
5:BO:12:TRP:NE1	6:BP:18:HIS:HA	2.18	0.58
14:BP:102:CRT:H2M2	5:BQ:37:MET:HG2	1.86	0.58
5:BY:46:TRP:CH2	9:BY:102:BCL:HBC3	2.38	0.58
5:B1:11:ILE:HD13	9:B3:102:BCL:C15	2.34	0.58
5:B3:12:TRP:HE1	6:B4:18:HIS:HA	1.68	0.58
5:B3:13:LEU:HB2	14:B7:102:CRT:H21A	1.85	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:B7:56:GLN:H	5:B7:56:GLN:CD	2.11	0.58
6:B8:46:LEU:O	5:B9:51:ILE:O	2.21	0.58
2:AL:237:ALA:O	2:AL:238:ILE:C	2.47	0.57
3:AM:157:TYR:CE1	3:AM:158:LEU:HD23	2.39	0.57
3:AM:200:PRO:HA	3:AM:203:MET:SD	2.44	0.57
3:AM:280:ALA:O	3:AM:282:ILE:N	2.37	0.57
5:AA:16:ASP:HB2	5:AA:18:ARG:HH11	1.69	0.57
9:AO:102:BCL:C3D	9:AP:101:BCL:C3D	2.82	0.57
14:AS:104:CRT:H182	9:AU:102:BCL:C8	2.34	0.57
5:AU:45:ASN:H	5:AW:56:GLN:HE21	1.52	0.57
6:AV:30:GLY:O	6:AV:33:VAL:HG12	2.04	0.57
6:AZ:45:TRP:O	6:AZ:46:LEU:HG	2.04	0.57
5:A1:11:ILE:CA	14:A1:103:CRT:H82	2.34	0.57
5:A3:12:TRP:HE1	6:A4:18:HIS:CA	2.15	0.57
5:A5:39:VAL:C	5:A5:41:SER:H	2.12	0.57
1:BC:110:CYS:HA	1:BC:123:THR:OG1	2.04	0.57
2:BL:164:ASP:O	2:BL:167:SER:N	2.37	0.57
2:BL:221:GLU:C	2:BL:223:THR:H	2.12	0.57
3:BM:34:PRO:HG3	3:BM:50:PRO:HD3	1.86	0.57
3:BM:83:VAL:HA	3:BM:86:PHE:HB3	1.86	0.57
4:BH:130:LEU:HG	4:BH:131:PRO:CD	2.29	0.57
4:BH:164:ALA:HB2	4:BH:216:ALA:CB	2.33	0.57
5:BQ:50:ASN:CB	5:BS:56:GLN:HA	2.34	0.57
5:BW:26:ALA:HA	5:BW:29:ILE:HG22	1.86	0.57
5:B1:51:ILE:HA	5:B1:52:PRO:O	2.03	0.57
9:B3:102:BCL:H2	6:B4:28:TRP:CH2	2.39	0.57
2:AL:48:LEU:HA	2:AL:51:VAL:CG2	2.34	0.57
2:AL:88:PRO:O	2:AL:91:GLU:N	2.36	0.57
3:AM:233:ARG:O	3:AM:234:GLU:C	2.47	0.57
3:AM:287:SER:OG	3:AM:294:TRP:NE1	2.37	0.57
5:AD:11:ILE:HG23	5:AD:12:TRP:CE3	2.38	0.57
14:A1:103:CRT:H183	9:A3:103:BCL:H8	1.86	0.57
2:BL:174:LEU:HD12	2:BL:174:LEU:N	2.19	0.57
5:BA:11:ILE:HD11	5:BD:21:LEU:CD2	2.33	0.57
9:BF:102:BCL:C1D	9:BG:101:BCL:CMD	2.81	0.57
9:BY:102:BCL:CMD	6:BZ:36:HIS:CD2	2.88	0.57
6:B2:17:PHE:CA	14:B2:102:CRT:H41	2.32	0.57
5:B7:28:GLN:O	9:B8:101:BCL:HED1	2.04	0.57
6:B8:31:LEU:O	6:B8:34:ILE:CG2	2.52	0.57
5:B9:29:ILE:HB	9:B9:102:BCL:H43	1.86	0.57
2:AL:75:ILE:HG22	2:AL:95:TRP:HD1	1.69	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:AL:89:LEU:CA	2:AL:94:LEU:H	2.00	0.57
3:AM:158:LEU:O	3:AM:163:ILE:HG22	2.03	0.57
3:AM:248:ALA:O	3:AM:251:PHE:N	2.35	0.57
4:AH:36:ARG:HE	4:AH:65:LYS:HD2	1.68	0.57
4:AH:178:GLN:HG3	4:AH:178:GLN:O	2.04	0.57
5:AD:36:HIS:O	5:AD:40:LEU:CB	2.52	0.57
5:AF:43:ASP:HB3	5:AI:47:LEU:HG	1.86	0.57
5:AI:14:ILE:HG23	5:AK:18:ARG:HB3	1.87	0.57
5:A7:4:MET:O	5:A7:8:LEU:HB2	2.04	0.57
6:A0:29:PHE:O	6:A0:32:VAL:HG12	2.03	0.57
2:BL:163:LEU:HD23	3:BM:197:TYR:O	2.04	0.57
4:BH:36:ARG:HE	4:BH:65:LYS:HD2	1.68	0.57
4:BH:215:LYS:HB2	4:BH:218:HIS:CD2	2.40	0.57
5:BD:46:TRP:CD1	5:BD:47:LEU:HD22	2.39	0.57
5:BD:50:ASN:ND2	5:BD:51:ILE:H	2.03	0.57
6:BZ:29:PHE:N	6:BZ:29:PHE:HD1	2.01	0.57
5:B9:12:TRP:HD1	6:B0:14:ALA:O	1.88	0.57
1:AC:271:TYR:O	1:AC:274:ARG:N	2.36	0.57
5:AD:8:LEU:O	5:AD:11:ILE:HG22	2.04	0.57
9:AK:102:BCL:CAC	9:AN:101:BCL:HBC3	2.33	0.57
9:AO:102:BCL:CBD	9:AP:101:BCL:CAD	2.82	0.57
5:AU:42:THR:C	5:AW:48:ASP:HB3	2.29	0.57
9:AW:101:BCL:O1A	6:AX:28:TRP:HH2	1.87	0.57
9:A2:101:BCL:HMA1	9:A3:103:BCL:HHB	1.85	0.57
5:A3:26:ALA:O	5:A3:29:ILE:HG22	2.05	0.57
5:A5:10:LYS:HB2	14:A5:103:CRT:H5	1.86	0.57
6:A6:40:TRP:HE3	6:A6:40:TRP:HA	1.69	0.57
2:BL:50:ILE:HA	2:BL:98:ILE:HD11	1.85	0.57
3:BM:260:VAL:HB	3:BM:264:SER:OG	2.03	0.57
9:BM:401:BCL:OBB	14:BM:406:CRT:H243	2.05	0.57
4:BH:204:LYS:H	4:BH:204:LYS:CD	2.14	0.57
14:BA:102:CRT:C14	6:BE:21:PHE:HA	2.35	0.57
9:BE:101:BCL:NB	9:BF:102:BCL:CMB	2.67	0.57
6:BG:12:ASP:O	6:BG:16:GLU:HG3	2.04	0.57
6:BG:42:TYR:OH	6:BG:43:ARG:NH2	2.38	0.57
5:BI:36:HIS:CE1	9:BI:102:BCL:NA	2.73	0.57
5:BK:20:VAL:O	5:BK:24:ILE:HG13	2.03	0.57
5:BO:9:TYR:HA	6:BP:18:HIS:ND1	2.18	0.57
6:BP:17:PHE:HA	6:BP:20:ILE:CG2	2.35	0.57
5:BQ:46:TRP:CZ2	9:BQ:103:BCL:H2C	2.39	0.57
14:BS:103:CRT:H14	6:BT:21:PHE:CD2	2.38	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:BZ:45:TRP:CE3	9:BZ:101:BCL:HAC2	2.40	0.57
1:AC:28:PRO:HD3	2:AL:262:PRO:HA	1.85	0.57
1:AC:325:LYS:HA	1:AC:331:TYR:HH	1.67	0.57
2:AL:219:GLU:CG	4:AH:127:PHE:HB2	2.34	0.57
2:AL:235:ALA:HA	11:AL:304:UQ8:C3M	2.34	0.57
4:AH:54:LYS:HE3	5:AD:23:SER:HB2	1.84	0.57
4:AH:182:LEU:HD13	4:AH:195:LEU:HG	1.87	0.57
4:AH:189:ASN:OD1	4:AH:220:ALA:HA	2.04	0.57
6:AB:40:TRP:HA	6:AB:40:TRP:HE3	1.69	0.57
14:AB:102:CRT:C3	5:A9:10:LYS:CA	2.82	0.57
5:AO:49:ASP:OD2	6:AP:43:ARG:NH2	2.37	0.57
6:AR:13:GLU:H	6:AR:13:GLU:CD	2.13	0.57
5:AY:5:ASN:CG	6:AZ:18:HIS:CD2	2.83	0.57
5:A3:40:LEU:HD21	5:A3:46:TRP:CH2	2.40	0.57
5:A7:26:ALA:O	5:A7:29:ILE:HG22	2.05	0.57
1:BC:285:TRP:CE3	1:BC:302:PRO:HG3	2.40	0.57
2:BL:186:ILE:HD13	9:BM:401:BCL:HMD2	1.85	0.57
3:BM:154:ILE:O	3:BM:154:ILE:HG22	2.03	0.57
3:BM:159:VAL:HG21	3:BM:281:GLY:CA	2.33	0.57
4:BH:105:ASP:OD1	4:BH:107:MET:HB3	2.05	0.57
5:BA:37:MET:SD	14:B0:101:CRT:H2M2	2.45	0.57
5:BA:44:LEU:HD11	5:BA:46:TRP:HE3	1.69	0.57
6:BB:18:HIS:O	6:BB:18:HIS:ND1	2.35	0.57
6:BV:43:ARG:HB3	5:BW:55:TYR:CE2	2.40	0.57
9:BW:102:BCL:HED1	6:BX:31:LEU:O	2.05	0.57
5:B1:51:ILE:HA	5:B1:52:PRO:C	2.29	0.57
5:B3:4:MET:SD	6:B6:23:GLN:NE2	2.77	0.57
6:B4:13:GLU:O	6:B4:16:GLU:HG2	2.04	0.57
2:AL:52:TRP:O	2:AL:55:THR:HB	2.04	0.57
2:AL:146:LEU:C	2:AL:148:MET:H	2.12	0.57
2:AL:154:GLY:O	2:AL:165:TRP:NE1	2.38	0.57
2:AL:204:LEU:CD1	3:AM:267:ARG:HD2	2.35	0.57
3:AM:271:TRP:O	3:AM:272:CYS:C	2.47	0.57
5:AQ:51:ILE:HG23	5:AQ:52:PRO:N	2.18	0.57
6:AV:30:GLY:O	6:AV:34:ILE:HG13	2.04	0.57
9:AY:102:BCL:C1D	9:AZ:101:BCL:CMD	2.82	0.57
6:A2:20:ILE:HD13	6:A2:20:ILE:C	2.30	0.57
5:A3:13:LEU:HD12	14:A7:102:CRT:C1M	2.33	0.57
5:A7:29:ILE:CG2	5:A7:30:VAL:H	2.18	0.57
5:A7:49:ASP:O	5:A9:59:GLY:HA3	2.03	0.57
6:A8:29:PHE:CZ	9:A8:101:BCL:C6	2.88	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BC:283:TYR:O	1:BC:286:PRO:HD2	2.05	0.57
1:BC:325:LYS:HD3	1:BC:325:LYS:O	2.05	0.57
3:BM:103:GLY:C	3:BM:104:LEU:HD22	2.28	0.57
5:BA:43:ASP:HA	5:BD:48:ASP:CB	2.27	0.57
9:BA:101:BCL:HBC1	9:BB:101:BCL:CBC	2.33	0.57
5:BQ:40:LEU:HD12	5:BQ:45:ASN:HA	1.86	0.57
6:BV:17:PHE:HA	14:BV:102:CRT:C6	2.34	0.57
5:B3:12:TRP:HE1	6:B4:18:HIS:CA	2.17	0.57
14:B5:103:CRT:H32	5:B7:31:LEU:HD21	1.85	0.57
6:B6:40:TRP:HA	6:B6:40:TRP:CE3	2.38	0.57
6:B8:26:TYR:HA	6:B8:29:PHE:HB3	1.87	0.57
1:AC:126:VAL:O	1:AC:129:ARG:N	2.38	0.57
2:AL:10:TYR:CD1	4:AH:112:GLY:HA2	2.40	0.57
3:AM:204:LEU:C	3:AM:206:ILE:H	2.12	0.57
4:AH:186:VAL:HG12	4:AH:187:ALA:N	2.19	0.57
9:AA:101:BCL:HBB1	9:A0:102:BCL:HMC1	1.86	0.57
6:AE:18:HIS:HE1	6:AE:22:MET:HE2	1.68	0.57
6:AR:10:THR:HG22	6:AR:11:ASP:N	2.16	0.57
5:AS:34:LEU:CA	15:AS:101:PEF:H442	2.34	0.57
9:AU:102:BCL:HBC2	9:AU:102:BCL:HHD	1.87	0.57
6:A4:42:TYR:C	6:A4:42:TYR:HD1	2.12	0.57
6:A8:22:MET:O	6:A8:26:TYR:HD2	1.87	0.57
5:A9:17:PRO:O	5:A9:21:LEU:HB2	2.05	0.57
1:BC:268:THR:CG2	7:BC:504:HEM:HAA1	2.34	0.57
2:BL:124:PHE:O	2:BL:127:PRO:HD2	2.03	0.57
3:BM:204:LEU:C	3:BM:206:ILE:H	2.12	0.57
3:BM:208:PHE:CE1	3:BM:275:LEU:HB3	2.40	0.57
5:BA:36:HIS:CE1	9:BB:101:BCL:HMD3	2.39	0.57
9:BK:102:BCL:ND	9:BN:101:BCL:HMD3	2.20	0.57
5:BS:39:VAL:O	5:BS:42:THR:HB	2.03	0.57
6:BT:36:HIS:CE1	9:BT:101:BCL:NA	2.72	0.57
5:B3:27:PHE:HE2	5:B5:29:ILE:HD12	1.70	0.57
6:B0:45:TRP:HD1	6:B0:46:LEU:H	1.52	0.57
1:AC:233:PHE:O	1:AC:236:MET:HB2	2.04	0.57
1:AC:276:VAL:O	1:AC:277:ARG:C	2.46	0.57
2:AL:22:LEU:HD22	5:A7:19:ARG:HB2	1.86	0.57
2:AL:52:TRP:HE1	5:A9:38:ILE:HA	1.69	0.57
2:AL:177:HIS:CD2	9:AL:301:BCL:CMC	2.88	0.57
2:AL:252:TRP:O	2:AL:255:VAL:N	2.38	0.57
9:AL:301:BCL:CHC	9:AM:402:BCL:CHC	2.83	0.57
3:AM:214:LEU:C	3:AM:214:LEU:CD2	2.78	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:AM:215:LEU:O	3:AM:217:ALA:N	2.38	0.57
14:AA:102:CRT:C14	6:AE:21:PHE:HA	2.34	0.57
9:AE:101:BCL:C1B	9:AF:102:BCL:CMB	2.77	0.57
5:AF:9:TYR:CE1	6:AG:15:LYS:HG3	2.40	0.57
6:AG:38:LEU:HD23	6:AG:39:ALA:N	2.20	0.57
6:AJ:10:THR:HB	6:AJ:13:GLU:CD	2.29	0.57
5:AO:9:TYR:CD1	6:AP:15:LYS:HD2	2.40	0.57
5:AU:43:ASP:HB2	5:AW:47:LEU:HD22	1.85	0.57
5:AW:10:LYS:NZ	14:AW:102:CRT:H1M2	2.19	0.57
14:AX:102:CRT:C2M	5:AY:36:HIS:HB2	2.35	0.57
5:A1:12:TRP:CD1	6:A2:18:HIS:HA	2.40	0.57
1:BC:53:ILE:O	1:BC:55:ALA:N	2.37	0.57
1:BC:155:CYS:O	1:BC:162:PRO:HB3	2.05	0.57
1:BC:164:TYR:CD2	1:BC:312:GLN:HG2	2.40	0.57
9:BL:301:BCL:H71	9:BL:301:BCL:H41	1.87	0.57
10:BL:302:BPH:H162	9:BL:303:BCL:CMB	2.34	0.57
14:BB:102:CRT:H2M2	5:BD:37:MET:HE2	1.85	0.57
5:BD:9:TYR:HB2	6:BE:15:LYS:CA	2.35	0.57
5:BD:31:LEU:HB3	9:BE:101:BCL:HED3	1.86	0.57
5:BF:40:LEU:HD11	5:BF:47:LEU:HD12	1.86	0.57
5:BI:18:ARG:CB	5:BI:18:ARG:HH11	2.17	0.57
5:BU:19:ARG:CZ	5:BW:18:ARG:NH2	2.67	0.57
6:BX:45:TRP:CE2	9:BX:101:BCL:H2C	2.39	0.57
6:B8:10:THR:HG22	6:B8:11:ASP:N	2.20	0.57
2:AL:129:ALA:HA	2:AL:247:LEU:CD1	2.33	0.57
2:AL:206:VAL:HG13	3:AM:142:MET:HE1	1.87	0.57
3:AM:5:GLN:O	3:AM:7:ILE:N	2.38	0.57
3:AM:240:HIS:CE1	4:AH:69:LEU:HD21	2.40	0.57
5:AI:22:VAL:O	5:AI:25:VAL:HB	2.05	0.57
6:AZ:46:LEU:HD22	6:A2:42:TYR:CZ	2.39	0.57
5:A1:8:LEU:CD2	5:A1:9:TYR:N	2.60	0.57
14:A1:103:CRT:H401	5:A3:38:ILE:HD12	1.87	0.57
5:A3:32:GLY:CA	9:A3:104:BCL:HED2	2.34	0.57
9:A0:102:BCL:H18	9:A0:102:BCL:HBB1	1.87	0.57
1:BC:29:GLY:O	1:BC:30:THR:HG23	2.05	0.57
1:BC:135:ARG:O	1:BC:136:ALA:C	2.47	0.57
2:BL:12:VAL:HG22	2:BL:13:ARG:H	1.67	0.57
3:BM:14:ARG:HD2	4:BH:146:GLU:OE1	2.04	0.57
4:BH:47:GLU:HG3	5:BA:19:ARG:CG	2.35	0.57
4:BH:134:VAL:HG21	4:BH:174:ARG:HH21	1.70	0.57
4:BH:194:LEU:HG	4:BH:225:LEU:HD23	1.87	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:BA:52:PRO:HD3	6:B0:45:TRP:O	2.05	0.57
6:BE:45:TRP:HA	5:BF:52:PRO:CG	2.34	0.57
6:BP:20:ILE:HG21	14:BP:102:CRT:C6	2.34	0.57
9:BP:101:BCL:HHB	9:BQ:103:BCL:HMA1	1.87	0.57
5:BS:29:ILE:HG23	5:BS:30:VAL:H	1.70	0.57
5:BW:4:MET:O	5:BW:6:ALA:N	2.38	0.57
5:B1:18:ARG:HG2	5:B1:18:ARG:NH1	2.19	0.57
14:B1:103:CRT:H372	5:B3:35:ILE:HD11	1.87	0.57
6:B2:38:LEU:O	6:B2:41:LEU:HG	2.05	0.57
5:B3:20:VAL:HA	5:B3:23:SER:HB3	1.86	0.57
5:B7:43:ASP:HA	5:B9:48:ASP:CB	2.29	0.57
5:B7:43:ASP:N	5:B9:48:ASP:HB3	2.19	0.57
9:B8:101:BCL:C1C	9:B9:102:BCL:CBB	2.81	0.57
1:AC:196:PRO:HG3	1:AC:231:TRP:CD1	2.40	0.57
1:AC:284:ILE:HD12	7:AC:502:HEM:HMD3	1.87	0.57
2:AL:109:TRP:HZ2	10:AL:302:BPH:HED3	1.70	0.57
3:AM:244:ALA:C	3:AM:246:GLU:H	2.13	0.57
3:AM:275:LEU:CD2	4:AH:19:PHE:HE2	2.17	0.57
5:AA:21:LEU:O	5:AA:25:VAL:HG23	2.05	0.57
6:AG:10:THR:HG22	6:AG:11:ASP:N	2.20	0.57
5:AS:18:ARG:O	5:AS:22:VAL:HG23	2.05	0.57
5:AS:46:TRP:CZ2	9:AS:103:BCL:CHC	2.87	0.57
5:AU:19:ARG:CZ	5:AW:18:ARG:NH2	2.68	0.57
5:AW:9:TYR:C	5:AW:11:ILE:H	2.13	0.57
6:AX:29:PHE:CZ	14:AX:102:CRT:H242	2.40	0.57
6:A4:42:TYR:C	6:A4:42:TYR:CD1	2.83	0.57
5:A9:36:HIS:CE1	9:A0:102:BCL:CMD	2.88	0.57
1:BC:142:LYS:O	1:BC:146:ALA:HA	2.05	0.57
2:BL:218:SER:C	2:BL:220:HIS:H	2.12	0.57
3:BM:59:LEU:CD2	3:BM:128:LEU:HD21	2.34	0.57
3:BM:132:ARG:O	3:BM:132:ARG:HD3	2.05	0.57
3:BM:160:LEU:CD2	3:BM:284:ILE:HG21	2.34	0.57
4:BH:113:PRO:HB2	4:BH:249:TYR:CE2	2.40	0.57
9:BB:101:BCL:C4B	9:BD:102:BCL:HBB3	2.34	0.57
9:BE:101:BCL:CHB	9:BF:102:BCL:HMB3	2.35	0.57
5:BK:44:LEU:O	5:BK:44:LEU:HD13	2.04	0.57
5:BO:4:MET:CB	6:BR:23:GLN:HG3	2.32	0.57
6:BT:24:SER:O	6:BT:27:ALA:HB3	2.05	0.57
5:BU:31:LEU:HD12	5:BU:34:LEU:HD23	1.86	0.57
5:BU:38:ILE:HD11	5:BW:40:LEU:HD23	1.87	0.57
5:BU:42:THR:C	5:BW:48:ASP:HB3	2.30	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:BZ:30:GLY:O	6:BZ:34:ILE:HG12	2.04	0.57
6:BZ:42:TYR:CD1	6:BZ:43:ARG:HG3	2.40	0.57
6:B2:20:ILE:HG21	14:B2:102:CRT:C6	2.32	0.57
6:B2:38:LEU:HD23	6:B2:38:LEU:C	2.29	0.57
1:AC:102:SER:C	1:AC:104:LYS:H	2.12	0.56
1:AC:173:LYS:HB3	3:AM:80:HIS:HB2	1.86	0.56
1:AC:265:LYS:HD2	1:AC:265:LYS:N	2.19	0.56
2:AL:52:TRP:O	2:AL:56:ILE:HG12	2.05	0.56
3:AM:31:ILE:O	3:AM:50:PRO:HB2	2.05	0.56
3:AM:241:ARG:HG2	3:AM:242:GLY:N	2.16	0.56
6:AB:29:PHE:O	6:AB:32:VAL:HG12	2.03	0.56
5:AO:5:ASN:HD21	6:AP:22:MET:HE3	1.71	0.56
6:AT:36:HIS:CE1	9:AT:101:BCL:NA	2.73	0.56
5:AU:14:ILE:HB	14:AX:102:CRT:C8	2.35	0.56
5:AW:4:MET:O	5:AW:8:LEU:HG	2.05	0.56
14:A5:103:CRT:C9	6:A8:17:PHE:HZ	2.17	0.56
6:A8:23:GLN:HG3	6:A8:24:SER:H	1.69	0.56
5:A9:2:PHE:HE1	6:A0:26:TYR:OH	1.88	0.56
1:BC:224:ALA:HB1	1:BC:228:GLN:OE1	2.05	0.56
2:BL:12:VAL:CG2	2:BL:13:ARG:N	2.68	0.56
2:BL:195:ALA:HA	2:BL:198:MET:CE	2.34	0.56
3:BM:84:PHE:CZ	5:BU:38:ILE:CD1	2.88	0.56
3:BM:186:THR:HA	9:BM:402:BCL:HMD1	1.87	0.56
3:BM:216:PHE:O	3:BM:216:PHE:CD1	2.58	0.56
3:BM:236:ASP:OD1	3:BM:237:GLN:N	2.38	0.56
3:BM:307:TYR:N	3:BM:307:TYR:CD1	2.72	0.56
5:BA:27:PHE:HA	5:BA:30:VAL:CG1	2.35	0.56
5:BD:8:LEU:O	5:BD:11:ILE:HG22	2.04	0.56
6:BE:21:PHE:C	6:BE:21:PHE:CD1	2.83	0.56
9:BF:102:BCL:ND	9:BG:101:BCL:HMD3	2.20	0.56
5:BK:16:ASP:O	5:BK:20:VAL:HG22	2.05	0.56
6:BP:21:PHE:CD1	14:BP:102:CRT:H14	2.40	0.56
6:BR:45:TRP:O	6:BR:46:LEU:HB2	2.05	0.56
5:BU:14:ILE:H	14:BU:103:CRT:C2	2.17	0.56
6:BV:25:MET:HE1	9:BW:102:BCL:H18	1.87	0.56
9:B7:103:BCL:HBA2	9:B8:101:BCL:OBD	2.04	0.56
6:B8:17:PHE:O	6:B8:20:ILE:HG22	2.05	0.56
1:AC:276:VAL:HG13	1:AC:277:ARG:H	1.70	0.56
2:AL:177:HIS:CD2	9:AL:301:BCL:HMC3	2.40	0.56
3:AM:76:LEU:HD21	5:AU:37:MET:HE3	1.87	0.56
3:AM:156:PHE:HD2	9:AM:402:BCL:H52	1.68	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:AM:200:PRO:CA	3:AM:203:MET:HG2	2.31	0.56
3:AM:254:TRP:CD1	3:AM:254:TRP:H	2.22	0.56
4:AH:69:LEU:CB	4:AH:70:PRO:HD2	2.36	0.56
6:AE:9:LEU:HB3	6:AE:13:GLU:OE2	2.05	0.56
6:AE:46:LEU:HD22	6:AG:42:TYR:CZ	2.40	0.56
5:AI:9:TYR:HA	6:AJ:18:HIS:CG	2.40	0.56
5:AK:18:ARG:O	5:AK:22:VAL:HG12	2.05	0.56
9:AK:102:BCL:HBD	9:AN:101:BCL:CAD	2.34	0.56
5:AY:45:ASN:HB3	5:AY:49:ASP:HB3	1.86	0.56
5:A3:20:VAL:HG13	9:A5:102:BCL:H203	1.86	0.56
6:A8:31:LEU:O	6:A8:34:ILE:CG2	2.52	0.56
5:A9:51:ILE:CB	5:A9:52:PRO:HA	2.34	0.56
1:BC:205:ASP:HB2	1:BC:304:ARG:HE	1.69	0.56
1:BC:271:TYR:O	1:BC:274:ARG:N	2.38	0.56
2:BL:71:TRP:N	2:BL:71:TRP:HE3	2.03	0.56
2:BL:182:HIS:HB2	2:BL:256:CYS:SG	2.45	0.56
3:BM:265:ILE:HG23	3:BM:266:HIS:N	2.18	0.56
4:BH:125:LEU:HB2	4:BH:129:GLY:O	2.04	0.56
4:BH:203:ASP:O	4:BH:205:LYS:N	2.39	0.56
6:BB:33:VAL:O	6:BB:37:LEU:HB2	2.06	0.56
6:BE:29:PHE:CD1	9:BE:101:BCL:H2	2.40	0.56
5:BF:12:TRP:NE1	6:BG:17:PHE:CD1	2.73	0.56
6:BG:17:PHE:CE1	6:BG:21:PHE:HD2	2.23	0.56
5:BK:9:TYR:CD1	5:BK:9:TYR:C	2.83	0.56
5:BO:34:LEU:HA	5:BO:37:MET:HB2	1.87	0.56
5:BS:8:LEU:HD22	5:BS:11:ILE:HD11	1.87	0.56
6:BZ:21:PHE:CE1	6:BZ:25:MET:HE3	2.40	0.56
5:B1:44:LEU:CD1	6:B2:43:ARG:HD2	2.31	0.56
6:B4:20:ILE:HD13	6:B4:20:ILE:C	2.30	0.56
1:AC:17:SER:HB3	3:AM:91:PHE:HZ	1.70	0.56
1:AC:68:THR:O	1:AC:86:SER:HB2	2.06	0.56
1:AC:242:SER:O	1:AC:313:ALA:CA	2.52	0.56
2:AL:52:TRP:HA	2:AL:52:TRP:HE3	1.69	0.56
2:AL:131:SER:O	2:AL:132:PHE:C	2.47	0.56
2:AL:139:VAL:HG11	2:AL:254:ALA:HB1	1.88	0.56
2:AL:144:ARG:CB	2:AL:145:PRO:HD3	2.35	0.56
3:AM:168:MET:HG2	3:AM:289:THR:HG22	1.86	0.56
3:AM:316:PRO:HG2	3:AM:317:TYR:CD1	2.40	0.56
4:AH:35:LYS:HG2	4:AH:39:TYR:CE2	2.41	0.56
4:AH:63:ASP:C	4:AH:79:PRO:HD2	2.30	0.56
4:AH:166:THR:O	4:AH:184:VAL:HG13	2.06	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:AH:173:ASP:OD1	4:AH:174:ARG:N	2.38	0.56
5:AA:47:LEU:HB3	5:A9:43:ASP:CA	2.35	0.56
5:AD:9:TYR:CE1	6:AE:11:ASP:HB3	2.40	0.56
5:AI:44:LEU:HD12	5:AI:46:TRP:HE3	1.70	0.56
6:AJ:40:TRP:HZ3	6:AJ:46:LEU:HG	1.70	0.56
5:AK:12:TRP:CD1	6:AN:17:PHE:HB3	2.40	0.56
5:AQ:43:ASP:HB2	5:AS:47:LEU:HB3	1.88	0.56
14:AS:104:CRT:H2M1	5:AW:37:MET:CA	2.35	0.56
6:AT:38:LEU:C	6:AT:38:LEU:HD23	2.30	0.56
5:AU:17:PRO:O	5:AU:21:LEU:HG	2.04	0.56
9:AW:101:BCL:CMD	6:AX:36:HIS:HA	2.36	0.56
6:AX:17:PHE:CZ	14:AX:102:CRT:H42	2.40	0.56
9:AX:101:BCL:HMC1	9:AY:102:BCL:HBB1	1.87	0.56
5:AY:30:VAL:O	5:AY:31:LEU:C	2.47	0.56
5:AY:52:PRO:HD2	5:AY:55:TYR:OH	2.05	0.56
5:A1:26:ALA:O	5:A1:29:ILE:HG22	2.05	0.56
5:A1:44:LEU:HD23	5:A1:44:LEU:N	2.21	0.56
9:A1:102:BCL:H71	6:A2:28:TRP:CD2	2.40	0.56
9:A2:101:BCL:CHB	9:A3:103:BCL:HMB3	2.35	0.56
5:A3:31:LEU:HD21	9:A3:104:BCL:CMA	2.35	0.56
9:A3:103:BCL:CHD	9:A3:104:BCL:HMD1	2.34	0.56
6:A4:20:ILE:HD13	6:A4:20:ILE:C	2.30	0.56
1:BC:233:PHE:O	1:BC:234:GLY:C	2.48	0.56
1:BC:264:PRO:HG2	1:BC:265:LYS:H	1.68	0.56
3:BM:13:VAL:HG12	4:BH:144:ILE:HA	1.86	0.56
3:BM:91:PHE:O	3:BM:180:PHE:HB2	2.04	0.56
4:BH:52:ARG:HB2	4:BH:54:LYS:HZ2	1.70	0.56
5:BA:9:TYR:CB	6:BB:18:HIS:HD2	2.19	0.56
5:BA:10:LYS:HD2	14:BA:102:CRT:H1M1	1.88	0.56
6:BB:17:PHE:CD1	14:BB:102:CRT:H6	2.40	0.56
6:BB:20:ILE:HD13	6:BB:20:ILE:O	2.04	0.56
6:BB:22:MET:C	6:BB:26:TYR:CE1	2.82	0.56
6:BB:36:HIS:CB	9:BB:101:BCL:H192	2.36	0.56
5:BD:28:GLN:HB3	9:BD:102:BCL:H12	1.88	0.56
5:BD:31:LEU:HB3	9:BE:101:BCL:CED	2.35	0.56
5:BI:36:HIS:NE2	9:BJ:101:BCL:HMD3	2.20	0.56
5:BK:54:SER:CB	5:BK:56:GLN:HE22	2.18	0.56
6:BN:34:ILE:HD13	6:BN:34:ILE:C	2.29	0.56
5:BO:43:ASP:HB2	5:BQ:47:LEU:CB	2.34	0.56
5:BS:51:ILE:HA	5:BS:52:PRO:C	2.30	0.56
14:BW:103:CRT:H35	5:BY:31:LEU:CD1	2.36	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:B3:8:LEU:HD22	5:B3:11:ILE:HD11	1.86	0.56
5:B9:46:TRP:CZ2	9:B9:102:BCL:H2C	2.40	0.56
1:AC:135:ARG:O	1:AC:136:ALA:C	2.49	0.56
1:AC:263:THR:HG22	3:AM:311:VAL:HB	1.86	0.56
1:AC:304:ARG:HG3	1:AC:304:ARG:NH1	2.21	0.56
2:AL:240:ARG:NH2	3:AM:6:ASN:C	2.63	0.56
3:AM:2:PRO:HB3	4:AH:201:ARG:NH1	2.19	0.56
6:AB:21:PHE:HD1	14:AB:102:CRT:H14	1.70	0.56
5:AD:18:ARG:O	5:AD:22:VAL:HG12	2.05	0.56
9:AN:101:BCL:HMB1	9:AO:102:BCL:C1B	2.36	0.56
5:AQ:29:ILE:HG23	5:AQ:30:VAL:N	2.20	0.56
5:AU:45:ASN:O	5:AU:49:ASP:HB3	2.06	0.56
5:A7:29:ILE:HD12	9:A7:103:BCL:H42	1.88	0.56
6:A8:26:TYR:HA	6:A8:29:PHE:HB3	1.87	0.56
9:A9:102:BCL:HMD2	6:A0:36:HIS:HD2	1.70	0.56
2:BL:88:PRO:O	2:BL:91:GLU:N	2.38	0.56
4:BH:102:PRO:CG	4:BH:106:PRO:HB3	2.35	0.56
5:BA:47:LEU:HB3	5:B9:43:ASP:CB	2.36	0.56
9:BA:101:BCL:CHD	9:BB:101:BCL:HMD1	2.35	0.56
5:BI:8:LEU:HB3	6:BJ:18:HIS:NE2	2.20	0.56
9:BJ:101:BCL:HMB1	9:BJ:101:BCL:CBB	2.36	0.56
5:BK:53:VAL:O	5:BK:55:TYR:N	2.39	0.56
6:BN:29:PHE:CZ	9:BN:101:BCL:H2	2.40	0.56
6:BT:45:TRP:O	5:BU:52:PRO:HD2	2.05	0.56
5:BU:44:LEU:HD22	6:BV:43:ARG:HD3	1.88	0.56
9:BV:101:BCL:OBB	9:BV:101:BCL:HHC	2.06	0.56
5:BY:55:TYR:HD1	5:BY:56:GLN:N	1.96	0.56
5:B5:12:TRP:CZ3	5:B5:17:PRO:HA	2.40	0.56
6:B6:28:TRP:C	6:B6:30:GLY:N	2.63	0.56
9:B7:103:BCL:C2D	9:B8:101:BCL:C2D	2.83	0.56
5:B9:12:TRP:HA	5:B9:12:TRP:HE3	1.70	0.56
1:AC:326:ASP:O	1:AC:327:TYR:CD1	2.59	0.56
2:AL:28:GLY:H	4:AH:46:THR:HG21	1.70	0.56
2:AL:264:TRP:CH2	2:AL:271:TRP:HA	2.41	0.56
2:AL:278:LEU:O	2:AL:280:LEU:N	2.38	0.56
3:AM:83:VAL:O	3:AM:87:LEU:HD23	2.06	0.56
3:AM:126:ILE:HD12	3:AM:157:TYR:CE2	2.41	0.56
5:AA:47:LEU:HB3	5:A9:43:ASP:CB	2.34	0.56
5:AF:9:TYR:HA	6:AG:18:HIS:ND1	2.20	0.56
6:AN:45:TRP:HD1	6:AN:46:LEU:HG	1.69	0.56
6:AP:45:TRP:O	6:AP:46:LEU:CG	2.54	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:AS:46:TRP:CD1	5:AS:47:LEU:HD13	2.40	0.56
6:AX:13:GLU:O	6:AX:16:GLU:HB3	2.05	0.56
6:AX:27:ALA:O	6:AX:31:LEU:HG	2.05	0.56
6:A2:45:TRP:NE1	9:A2:101:BCL:HHC	2.21	0.56
6:A6:40:TRP:HA	6:A6:40:TRP:CE3	2.39	0.56
1:BC:176:SER:OG	5:BU:48:ASP:HB3	2.05	0.56
1:BC:272:ALA:C	1:BC:274:ARG:H	2.13	0.56
2:BL:88:PRO:HB2	2:BL:91:GLU:HB2	1.86	0.56
2:BL:203:ILE:C	2:BL:205:SER:H	2.13	0.56
9:BK:102:BCL:C4D	9:BN:101:BCL:HMD3	2.35	0.56
14:BO:103:CRT:C14	6:BR:21:PHE:HB2	2.35	0.56
5:BQ:38:ILE:O	5:BQ:42:THR:HG22	2.05	0.56
6:BR:42:TYR:CD2	6:BR:43:ARG:HG3	2.40	0.56
6:BR:46:LEU:HB3	6:BT:42:TYR:OH	2.06	0.56
5:BS:44:LEU:HD23	5:BS:44:LEU:H	1.69	0.56
9:BT:101:BCL:OBB	9:BT:101:BCL:HHC	2.05	0.56
5:BU:30:VAL:HG13	5:BU:31:LEU:H	1.70	0.56
6:BV:21:PHE:CE1	14:BV:102:CRT:H14	2.36	0.56
5:BY:39:VAL:HG22	5:B1:47:LEU:HD11	1.87	0.56
5:B1:44:LEU:HD23	5:B1:44:LEU:H	1.71	0.56
5:B9:32:GLY:CA	9:B9:102:BCL:O1A	2.51	0.56
7:AC:503:HEM:HBB1	2:AL:174:LEU:HG	1.87	0.56
2:AL:125:HIS:NE2	3:AM:5:GLN:HG3	2.20	0.56
2:AL:233:ILE:HG12	2:AL:237:ALA:HB1	1.88	0.56
3:AM:280:ALA:HA	9:AM:402:BCL:O1D	2.05	0.56
5:AA:44:LEU:HD11	5:AA:46:TRP:HE3	1.70	0.56
5:AI:44:LEU:HA	5:AK:56:GLN:HB2	1.87	0.56
6:AJ:21:PHE:C	6:AJ:21:PHE:HD1	2.13	0.56
5:AQ:17:PRO:O	5:AQ:21:LEU:HG	2.05	0.56
5:AS:13:LEU:HD12	14:AS:104:CRT:H1M2	1.86	0.56
5:AU:26:ALA:HA	5:AU:29:ILE:HG22	1.88	0.56
5:AY:42:THR:O	5:AY:43:ASP:C	2.48	0.56
6:AZ:44:PRO:O	5:A1:55:TYR:CZ	2.59	0.56
6:A4:34:ILE:O	6:A4:34:ILE:HD13	2.05	0.56
6:A6:10:THR:HG22	6:A6:11:ASP:N	2.21	0.56
5:A7:36:HIS:CG	14:A7:102:CRT:H393	2.40	0.56
5:A7:46:TRP:CZ3	9:A7:103:BCL:HBC3	2.40	0.56
9:A0:102:BCL:C14	9:A0:102:BCL:CMB	2.84	0.56
1:BC:291:LEU:CD2	1:BC:292:PRO:HD2	2.36	0.56
2:BL:184:LEU:CD2	2:BL:252:TRP:HE1	2.18	0.56
3:BM:63:PHE:HB3	3:BM:125:SER:CB	2.26	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:BM:83:VAL:O	3:BM:86:PHE:HB3	2.06	0.56
3:BM:276:THR:C	3:BM:278:ILE:H	2.13	0.56
4:BH:19:PHE:C	4:BH:21:LEU:H	2.14	0.56
5:BA:45:ASN:OD1	5:BA:47:LEU:HB2	2.06	0.56
6:BB:33:VAL:O	6:BB:37:LEU:HD23	2.05	0.56
5:BQ:50:ASN:CG	5:BQ:51:ILE:H	2.12	0.56
5:BS:28:GLN:HB2	9:BS:102:BCL:H43	1.87	0.56
5:BU:35:ILE:HA	5:BU:38:ILE:HG22	1.86	0.56
5:BW:5:ASN:CB	6:BX:22:MET:HE3	2.35	0.56
9:BY:102:BCL:HBC2	9:BZ:101:BCL:HMD1	1.88	0.56
6:B4:34:ILE:O	6:B4:34:ILE:HD13	2.05	0.56
5:B7:46:TRP:NE1	5:B7:47:LEU:HD21	2.21	0.56
1:AC:94:MET:SD	7:AC:501:HEM:NA	2.78	0.56
1:AC:190:VAL:O	1:AC:192:TYR:N	2.38	0.56
1:AC:200:LEU:HG	1:AC:204:LEU:HD12	1.86	0.56
3:AM:27:ASN:HD21	5:AO:19:ARG:HH11	1.52	0.56
3:AM:187:ALA:HA	9:AM:402:BCL:HBC1	1.88	0.56
3:AM:233:ARG:O	3:AM:235:ILE:N	2.39	0.56
4:AH:105:ASP:OD1	4:AH:107:MET:HB3	2.05	0.56
9:AA:101:BCL:C1D	9:AB:101:BCL:HMD3	2.34	0.56
6:AB:38:LEU:HD23	6:AB:38:LEU:C	2.31	0.56
5:AQ:20:VAL:O	5:AQ:24:ILE:HD13	2.06	0.56
5:AS:13:LEU:CB	14:AS:104:CRT:H32A	2.35	0.56
9:AX:101:BCL:C4B	9:AY:102:BCL:HBB3	2.35	0.56
5:A3:12:TRP:HE1	6:A4:18:HIS:HA	1.71	0.56
5:A9:13:LEU:O	6:A0:7:THR:HB	2.06	0.56
9:A0:102:BCL:C14	9:A0:102:BCL:HMB1	2.36	0.56
1:BC:251:HIS:ND1	1:BC:256:PHE:HA	2.21	0.56
4:BH:5:ILE:CG1	5:BF:40:LEU:HD21	2.36	0.56
4:BH:232:THR:O	4:BH:235:GLU:HG2	2.06	0.56
5:BD:9:TYR:CB	6:BE:15:LYS:HA	2.34	0.56
6:BG:29:PHE:O	6:BG:33:VAL:HB	2.06	0.56
5:BK:12:TRP:CD1	6:BN:14:ALA:O	2.59	0.56
5:BO:46:TRP:CD1	5:BO:47:LEU:HD13	2.39	0.56
9:BO:102:BCL:OBD	6:BP:32:VAL:HG13	2.05	0.56
5:BQ:43:ASP:N	5:BS:47:LEU:HB3	2.21	0.56
6:BV:30:GLY:O	6:BV:33:VAL:HG12	2.06	0.56
5:BW:10:LYS:HD2	6:BZ:20:ILE:HD12	1.87	0.56
6:BX:10:THR:H	6:BX:13:GLU:CD	2.13	0.56
6:B2:24:SER:O	6:B2:27:ALA:HB3	2.06	0.56
6:B0:36:HIS:HD1	9:B0:102:BCL:H141	1.70	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AC:164:TYR:HB2	1:AC:309:THR:HA	1.88	0.56
1:AC:199:PRO:O	1:AC:203:PHE:HB2	2.06	0.56
1:AC:212:ILE:HD11	7:AC:503:HEM:HAA1	1.88	0.56
1:AC:283:TYR:N	1:AC:283:TYR:CD1	2.74	0.56
2:AL:131:SER:HA	2:AL:134:ILE:HD12	1.86	0.56
3:AM:165:PRO:HB3	3:AM:174:ALA:HB2	1.87	0.56
5:AI:39:VAL:HG21	9:AI:102:BCL:HBC1	1.88	0.56
5:AK:33:LEU:C	5:AK:33:LEU:HD12	2.30	0.56
5:AU:20:VAL:CG1	9:AW:101:BCL:H202	2.35	0.56
5:A1:4:MET:SD	6:A4:24:SER:HA	2.45	0.56
6:A0:24:SER:CB	14:A0:101:CRT:H183	2.35	0.56
1:BC:94:MET:SD	7:BC:501:HEM:NC	2.79	0.56
1:BC:226:LEU:HD12	3:BM:192:ARG:HB2	1.87	0.56
1:BC:274:ARG:HA	1:BC:277:ARG:HG2	1.87	0.56
2:BL:71:TRP:N	2:BL:71:TRP:CE3	2.74	0.56
2:BL:106:PHE:HD1	10:BL:302:BPH:H1C1	1.70	0.56
2:BL:170:GLY:HA3	9:BL:301:BCL:HBC3	1.86	0.56
2:BL:195:ALA:O	2:BL:198:MET:HB2	2.06	0.56
2:BL:197:SER:HB3	3:BM:273:ALA:CB	2.32	0.56
9:BL:301:BCL:HBC3	9:BL:301:BCL:HHD	1.87	0.56
3:BM:81:TRP:O	5:BU:41:SER:HB3	2.05	0.56
3:BM:197:TYR:CE1	9:BM:402:BCL:CMC	2.88	0.56
4:BH:135:PRO:HB3	4:BH:171:TRP:CE2	2.40	0.56
14:BA:102:CRT:H11	6:BE:17:PHE:HE1	1.71	0.56
5:BD:50:ASN:CG	5:BD:51:ILE:H	2.11	0.56
6:BE:29:PHE:CE1	9:BE:101:BCL:C2	2.89	0.56
9:BE:101:BCL:NB	9:BF:102:BCL:HMB1	2.21	0.56
5:BF:30:VAL:HG13	5:BF:31:LEU:N	2.20	0.56
6:BJ:40:TRP:HZ3	6:BJ:46:LEU:HG	1.71	0.56
6:BN:45:TRP:O	6:BN:46:LEU:HB2	2.05	0.56
5:BQ:50:ASN:HD22	5:BS:56:GLN:CA	2.12	0.56
9:BQ:103:BCL:CHD	9:BQ:104:BCL:CMD	2.79	0.56
5:BY:9:TYR:CD1	5:BY:9:TYR:C	2.84	0.56
5:BY:16:ASP:HB2	5:BY:19:ARG:NH2	2.21	0.56
5:B3:55:TYR:O	5:B3:59:GLY:HA3	2.06	0.56
5:B7:47:LEU:H	5:B7:47:LEU:CD2	2.18	0.56
1:AC:43:TYR:HE1	2:AL:153:HIS:HE2	1.54	0.56
1:AC:148:THR:HA	1:AC:322:GLN:CB	2.36	0.56
2:AL:10:TYR:O	2:AL:12:VAL:N	2.37	0.56
3:AM:196:LEU:C	3:AM:198:TYR:N	2.62	0.56
3:AM:261:THR:O	3:AM:263:GLU:N	2.39	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:AM:265:ILE:HG22	3:AM:266:HIS:N	2.10	0.56
3:AM:265:ILE:O	3:AM:267:ARG:N	2.38	0.56
4:AH:235:GLU:HA	4:AH:238:LYS:HG2	1.87	0.56
5:AQ:30:VAL:HG13	5:AQ:31:LEU:N	2.21	0.56
5:AS:34:LEU:HA	15:AS:101:PEF:H442	1.88	0.56
5:A1:12:TRP:CD1	6:A2:18:HIS:CA	2.89	0.56
5:A3:51:ILE:HG22	5:A3:54:SER:CB	2.36	0.56
5:A7:19:ARG:O	5:A7:23:SER:HB2	2.06	0.56
6:A8:20:ILE:HG23	6:A8:21:PHE:N	2.20	0.56
1:BC:112:VAL:HG12	1:BC:113:PRO:HD2	1.88	0.56
1:BC:153:TYR:CE1	1:BC:157:ARG:HA	2.40	0.56
2:BL:177:HIS:CD2	9:BL:301:BCL:HMC3	2.41	0.56
3:BM:12:GLN:C	4:BH:145:ALA:HB2	2.30	0.56
4:BH:55:VAL:HG11	5:BD:19:ARG:HD3	1.87	0.56
5:BD:15:LEU:HB3	5:BD:20:VAL:CG2	2.31	0.56
5:BK:8:LEU:O	5:BK:11:ILE:HG13	2.06	0.56
5:BK:18:ARG:HG2	5:BK:18:ARG:HH11	1.70	0.56
9:BQ:104:BCL:H2C	6:BR:45:TRP:CE3	2.41	0.56
6:BR:34:ILE:HD12	6:BR:35:ALA:N	2.20	0.56
5:BS:49:ASP:CG	5:BS:50:ASN:N	2.64	0.56
5:BU:2:PHE:HA	5:BU:5:ASN:HB2	1.88	0.56
5:BU:42:THR:HB	5:BW:48:ASP:CB	2.35	0.56
1:AC:97:VAL:HG21	1:AC:131:PHE:HZ	1.70	0.56
1:AC:253:THR:HG22	2:AL:171:TYR:CD2	2.41	0.56
2:AL:94:LEU:HD23	2:AL:94:LEU:C	2.30	0.56
2:AL:188:PHE:HE2	2:AL:248:SER:HB3	1.69	0.56
3:AM:66:VAL:HG11	3:AM:121:PHE:CD2	2.31	0.56
3:AM:132:ARG:HD3	3:AM:132:ARG:O	2.05	0.56
3:AM:168:MET:HG2	3:AM:289:THR:CG2	2.36	0.56
5:AA:14:ILE:C	5:AA:15:LEU:HD22	2.31	0.56
6:AB:22:MET:HG3	6:AB:26:TYR:HE1	1.71	0.56
5:AD:7:ASN:H	5:AD:7:ASN:ND2	2.04	0.56
6:AG:21:PHE:HB2	14:AG:102:CRT:H11	1.88	0.56
9:AJ:101:BCL:CMC	9:AK:102:BCL:HBB1	2.35	0.56
5:AO:50:ASN:CG	6:AP:43:ARG:NH2	2.62	0.56
6:AT:24:SER:O	6:AT:27:ALA:HB3	2.05	0.56
5:AW:8:LEU:HD21	6:AZ:24:SER:OG	2.06	0.56
14:AW:102:CRT:H83	6:AZ:20:ILE:HD13	1.88	0.56
5:A7:49:ASP:OD2	6:A8:43:ARG:NH2	2.38	0.56
6:A0:21:PHE:HB2	14:A0:101:CRT:C15	2.35	0.56
1:BC:20:LEU:HG	2:BL:271:TRP:CE2	2.40	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BC:24:GLU:O	2:BL:263:PHE:HA	2.06	0.56
2:BL:178:TYR:CD1	3:BM:183:LEU:HD13	2.41	0.56
5:BD:27:PHE:O	5:BD:30:VAL:HG12	2.06	0.56
5:BI:35:ILE:HA	5:BI:38:ILE:CG2	2.32	0.56
6:BJ:23:GLN:CD	6:BJ:24:SER:N	2.64	0.56
9:BK:102:BCL:CMD	6:BN:36:HIS:HD2	2.18	0.56
5:BO:8:LEU:O	5:BO:11:ILE:HG13	2.06	0.56
5:BO:46:TRP:CD1	5:BO:47:LEU:HD22	2.41	0.56
14:BP:102:CRT:H2M3	5:BQ:36:HIS:HB3	1.86	0.56
6:BT:29:PHE:CD1	6:BT:29:PHE:N	2.72	0.56
5:BW:18:ARG:O	5:BW:22:VAL:HG22	2.05	0.56
5:BW:18:ARG:O	5:BW:22:VAL:HG23	2.04	0.56
5:BW:49:ASP:OD1	5:BW:50:ASN:N	2.24	0.56
6:B6:10:THR:HG22	6:B6:11:ASP:N	2.21	0.56
5:B9:36:HIS:CE1	9:B0:102:BCL:CMD	2.88	0.56
1:AC:71:LYS:N	1:AC:71:LYS:HD2	2.21	0.55
1:AC:283:TYR:HD1	1:AC:283:TYR:N	2.03	0.55
2:AL:20:GLY:C	2:AL:22:LEU:H	2.15	0.55
2:AL:255:VAL:HG12	2:AL:256:CYS:N	2.21	0.55
3:AM:200:PRO:CA	3:AM:203:MET:HE3	2.36	0.55
3:AM:206:ILE:HG12	9:AM:402:BCL:HMA2	1.87	0.55
3:AM:234:GLU:HA	3:AM:262:MET:HE3	1.87	0.55
4:AH:69:LEU:HB2	4:AH:74:GLY:O	2.05	0.55
5:AA:11:ILE:CA	14:AA:102:CRT:H82	2.36	0.55
5:AF:27:PHE:HA	5:AF:30:VAL:CG1	2.34	0.55
5:AU:14:ILE:HG13	14:AX:102:CRT:H5	1.87	0.55
14:AW:102:CRT:H35	5:AY:31:LEU:CD1	2.35	0.55
6:A8:17:PHE:O	6:A8:20:ILE:HG22	2.05	0.55
6:A0:7:THR:HG23	6:A0:8:GLY:H	1.71	0.55
6:A0:20:ILE:HD13	6:A0:20:ILE:C	2.31	0.55
1:BC:283:TYR:N	1:BC:283:TYR:CD1	2.73	0.55
3:BM:132:ARG:HG2	3:BM:132:ARG:HH11	1.71	0.55
5:BA:2:PHE:HA	5:BA:5:ASN:ND2	2.20	0.55
9:BI:102:BCL:ND	9:BJ:101:BCL:CMD	2.69	0.55
14:BU:103:CRT:H2M3	5:BY:36:HIS:CB	2.36	0.55
5:B9:16:ASP:O	5:B9:20:VAL:HG22	2.06	0.55
1:AC:212:ILE:HD13	1:AC:212:ILE:N	2.21	0.55
2:AL:139:VAL:HA	2:AL:143:VAL:HB	1.87	0.55
3:AM:133:THR:O	3:AM:137:ALA:N	2.39	0.55
5:AI:50:ASN:CG	5:AI:51:ILE:H	2.15	0.55
9:AO:102:BCL:OBB	9:AO:102:BCL:HHC	2.07	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:AR:45:TRP:CE3	9:AR:101:BCL:HBC2	2.41	0.55
5:AS:34:LEU:CA	15:AS:101:PEF:H453	2.31	0.55
6:AT:22:MET:O	6:AT:26:TYR:HD1	1.89	0.55
6:AX:34:ILE:C	6:AX:34:ILE:HD13	2.31	0.55
5:A3:28:GLN:CG	9:A3:103:BCL:H12	2.36	0.55
1:BC:199:PRO:C	1:BC:202:PRO:HD2	2.32	0.55
2:BL:78:PRO:HB3	2:BL:92:GLY:HA3	1.89	0.55
2:BL:210:GLN:HB2	2:BL:213:GLU:HG3	1.86	0.55
3:BM:243:THR:O	3:BM:247:ARG:HG3	2.06	0.55
3:BM:284:ILE:HG13	9:BM:402:BCL:OBD	2.06	0.55
5:BI:22:VAL:HA	5:BI:25:VAL:CG2	2.36	0.55
5:BK:51:ILE:HA	5:BK:52:PRO:C	2.31	0.55
6:BN:10:THR:C	6:BN:13:GLU:OE2	2.49	0.55
5:BO:37:MET:O	5:BO:41:SER:HB2	2.07	0.55
6:B4:42:TYR:C	6:B4:42:TYR:CD1	2.83	0.55
6:B6:40:TRP:HZ3	6:B6:44:PRO:CA	2.17	0.55
5:B9:16:ASP:OD2	5:B9:17:PRO:HD2	2.05	0.55
5:B9:36:HIS:NE2	9:B0:102:BCL:CMD	2.70	0.55
1:AC:35:TYR:CD2	3:AM:308:PRO:HD2	2.41	0.55
1:AC:293:ALA:C	1:AC:295:ARG:H	2.15	0.55
2:AL:160:LEU:HA	2:AL:163:LEU:HD13	1.88	0.55
3:AM:268:TRP:H	3:AM:268:TRP:HD1	1.53	0.55
4:AH:55:VAL:CG1	4:AH:56:VAL:H	2.09	0.55
4:AH:55:VAL:HG11	5:AD:19:ARG:HD3	1.87	0.55
4:AH:66:THR:HA	4:AH:77:VAL:HG12	1.89	0.55
4:AH:159:LEU:CB	4:AH:212:ASP:HA	2.36	0.55
5:AA:45:ASN:O	5:AA:49:ASP:HB3	2.06	0.55
6:AB:17:PHE:O	6:AB:20:ILE:HG22	2.06	0.55
6:AB:45:TRP:O	6:AB:46:LEU:CB	2.53	0.55
6:AB:46:LEU:HD13	6:AE:42:TYR:CZ	2.41	0.55
14:AB:102:CRT:H2M2	5:AD:37:MET:CE	2.35	0.55
5:AF:20:VAL:HB	9:AI:102:BCL:C20	2.37	0.55
5:AF:49:ASP:HB2	5:AI:56:GLN:CD	2.32	0.55
5:AK:36:HIS:CE1	9:AK:102:BCL:NA	2.74	0.55
6:AP:21:PHE:O	6:AP:22:MET:C	2.49	0.55
5:AW:21:LEU:HD13	14:AX:102:CRT:H131	1.87	0.55
6:AZ:10:THR:CG2	6:AZ:11:ASP:H	2.13	0.55
6:A6:44:PRO:O	5:A7:52:PRO:HD2	2.05	0.55
5:A7:29:ILE:CG2	5:A7:30:VAL:N	2.69	0.55
5:A7:29:ILE:CA	9:A7:103:BCL:H11	2.30	0.55
1:BC:141:TRP:CE3	1:BC:275:HIS:HB2	2.40	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BC:183:GLN:O	1:BC:195:LEU:O	2.25	0.55
1:BC:200:LEU:HD11	1:BC:238:ASN:ND2	2.21	0.55
1:BC:242:SER:O	1:BC:313:ALA:CA	2.53	0.55
3:BM:176:PRO:HD2	3:BM:185:TRP:HB2	1.87	0.55
3:BM:271:TRP:CD2	4:BH:26:LEU:HD21	2.41	0.55
4:BH:225:LEU:HD12	4:BH:225:LEU:O	2.05	0.55
5:BI:26:ALA:O	5:BI:29:ILE:N	2.39	0.55
5:BO:9:TYR:CD2	6:BP:15:LYS:HD2	2.41	0.55
9:BQ:104:BCL:HAC2	6:BR:45:TRP:CZ3	2.41	0.55
9:B1:102:BCL:HMD3	6:B2:36:HIS:CE1	2.42	0.55
5:B5:10:LYS:CB	14:B5:103:CRT:H5	2.30	0.55
5:B5:45:ASN:O	5:B5:49:ASP:HB3	2.06	0.55
9:B9:102:BCL:OBB	9:B9:102:BCL:HHC	2.05	0.55
6:B0:33:VAL:CG1	6:B0:37:LEU:CD1	2.77	0.55
2:AL:119:LYS:C	2:AL:121:GLY:H	2.14	0.55
4:AH:130:LEU:HD13	4:AH:174:ARG:HH12	1.72	0.55
4:AH:135:PRO:C	4:AH:137:ARG:H	2.14	0.55
9:AA:101:BCL:C1D	9:AB:101:BCL:HMD1	2.36	0.55
9:AE:101:BCL:C4B	9:AF:102:BCL:HBB3	2.37	0.55
5:AF:45:ASN:HB3	5:AF:49:ASP:HB3	1.88	0.55
9:AU:102:BCL:CHD	9:AU:102:BCL:CBC	2.84	0.55
5:A3:32:GLY:HA2	9:A3:104:BCL:O1D	2.05	0.55
6:A6:17:PHE:O	6:A6:20:ILE:HG22	2.06	0.55
1:BC:327:TYR:HB2	1:BC:330:LEU:HD12	1.87	0.55
3:BM:155:PHE:O	3:BM:159:VAL:HG23	2.06	0.55
3:BM:229:PHE:O	3:BM:244:ALA:HB2	2.06	0.55
4:BH:206:ALA:C	4:BH:208:LYS:H	2.14	0.55
14:BA:102:CRT:H9	6:BE:17:PHE:HD1	1.70	0.55
14:BU:103:CRT:H2M2	5:BY:37:MET:HG2	1.88	0.55
5:BW:36:HIS:CE1	9:BX:101:BCL:HMD3	2.41	0.55
6:BX:36:HIS:HE1	9:BX:101:BCL:C1B	2.19	0.55
5:BY:49:ASP:HB2	5:B1:56:GLN:CD	2.31	0.55
6:B2:16:GLU:CD	14:B2:102:CRT:H1M1	2.30	0.55
5:B3:28:GLN:HA	5:B3:28:GLN:NE2	2.21	0.55
6:B6:17:PHE:O	6:B6:20:ILE:HG22	2.07	0.55
5:B9:43:ASP:OD1	5:B9:44:LEU:HD12	2.06	0.55
3:AM:197:TYR:CE1	9:AM:402:BCL:CMC	2.88	0.55
3:AM:259:ASN:C	3:AM:259:ASN:HD22	2.14	0.55
4:AH:35:LYS:HZ2	4:AH:59:PRO:HG2	1.72	0.55
5:AA:9:TYR:CZ	5:AA:10:LYS:HD3	2.42	0.55
5:AA:47:LEU:CD1	5:A9:43:ASP:HB2	2.31	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:AA:50:ASN:CG	6:AB:43:ARG:HH21	2.14	0.55
5:AQ:43:ASP:OD1	5:AQ:44:LEU:HD23	2.06	0.55
14:AS:104:CRT:H14	6:AV:21:PHE:CD1	2.41	0.55
5:AW:12:TRP:CE3	5:AW:12:TRP:CA	2.89	0.55
9:AY:102:BCL:CMD	6:AZ:36:HIS:HD2	2.20	0.55
5:A1:12:TRP:CD2	6:A2:17:PHE:CE2	2.95	0.55
5:A1:51:ILE:HA	5:A1:52:PRO:C	2.32	0.55
9:A7:103:BCL:CMD	6:A8:36:HIS:CD2	2.90	0.55
9:A9:102:BCL:OBB	9:A9:102:BCL:HHC	2.06	0.55
3:BM:104:LEU:HD22	3:BM:104:LEU:N	2.22	0.55
5:BF:10:LYS:HB3	14:BF:103:CRT:C1M	2.36	0.55
5:BK:45:ASN:CA	5:BK:49:ASP:HB3	2.36	0.55
6:BN:20:ILE:H	6:BN:20:ILE:CD1	2.18	0.55
5:BO:4:MET:HE3	6:BR:20:ILE:HA	1.87	0.55
6:BT:29:PHE:N	6:BT:29:PHE:HD1	2.05	0.55
5:BU:12:TRP:NE1	6:BV:18:HIS:CA	2.53	0.55
9:BU:102:BCL:HBC2	9:BU:102:BCL:HHD	1.88	0.55
5:BW:32:GLY:N	9:BX:101:BCL:HED2	2.21	0.55
6:BX:10:THR:HG22	6:BX:11:ASP:N	2.22	0.55
6:BX:30:GLY:HA2	6:BX:33:VAL:HG12	1.88	0.55
6:BZ:18:HIS:CE1	6:BZ:22:MET:HE2	2.41	0.55
5:B7:37:MET:H	14:B7:102:CRT:C2M	2.14	0.55
6:B8:20:ILE:HG23	6:B8:21:PHE:N	2.21	0.55
6:B8:46:LEU:HB3	6:B0:42:TYR:OH	2.06	0.55
1:AC:97:VAL:CG1	7:AC:501:HEM:HBC2	2.37	0.55
3:AM:132:ARG:HG2	3:AM:132:ARG:HH11	1.72	0.55
5:AA:26:ALA:O	5:AA:29:ILE:HG22	2.06	0.55
6:AB:20:ILE:HD13	6:AB:20:ILE:C	2.31	0.55
5:AO:10:LYS:HB2	14:AR:102:CRT:H5	1.87	0.55
5:AQ:45:ASN:O	5:AQ:47:LEU:N	2.39	0.55
6:AV:9:LEU:HB3	6:AV:13:GLU:OE1	2.06	0.55
6:A2:40:TRP:CE3	6:A2:44:PRO:HA	2.42	0.55
5:A3:36:HIS:CE1	9:A3:104:BCL:HMD3	2.42	0.55
9:A3:103:BCL:HED1	6:A4:32:VAL:HA	1.87	0.55
5:A5:51:ILE:HB	5:A5:52:PRO:CA	2.31	0.55
6:A8:10:THR:HG22	6:A8:11:ASP:N	2.20	0.55
1:BC:207:ALA:CB	1:BC:277:ARG:HE	2.19	0.55
2:BL:87:ALA:O	2:BL:93:GLY:HA3	2.07	0.55
2:BL:233:ILE:HG12	2:BL:237:ALA:CB	2.36	0.55
3:BM:102:TYR:CE2	3:BM:108:PRO:HD3	2.41	0.55
3:BM:281:GLY:O	3:BM:285:LEU:HB2	2.06	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:BH:35:LYS:HG2	4:BH:39:TYR:CE2	2.42	0.55
5:BA:7:ASN:C	5:BA:8:LEU:HD23	2.31	0.55
9:BA:101:BCL:HBC2	9:BB:101:BCL:HHD	1.88	0.55
6:BB:22:MET:HG3	6:BB:26:TYR:HE1	1.68	0.55
5:BW:35:ILE:HA	5:BW:38:ILE:CG2	2.37	0.55
5:B7:33:LEU:O	14:B7:102:CRT:C2M	2.48	0.55
5:B7:36:HIS:CB	14:B7:102:CRT:H391	2.32	0.55
1:AC:80:GLN:H	1:AC:128:ARG:NH2	2.04	0.55
1:AC:254:ARG:HD3	1:AC:255:ALA:N	2.21	0.55
2:AL:237:ALA:O	2:AL:240:ARG:HG3	2.06	0.55
2:AL:273:ASN:C	2:AL:275:TRP:N	2.63	0.55
3:AM:242:GLY:O	3:AM:246:GLU:HB2	2.07	0.55
9:AA:101:BCL:HMB3	9:A0:102:BCL:CHB	2.37	0.55
5:AF:33:LEU:N	5:AF:33:LEU:HD12	2.22	0.55
6:AG:40:TRP:HH2	6:AG:46:LEU:CD1	2.19	0.55
6:AN:45:TRP:CZ3	9:AN:101:BCL:HAC2	2.42	0.55
5:AS:34:LEU:HB2	15:AS:101:PEF:H431	1.87	0.55
5:AU:27:PHE:CD2	5:AW:29:ILE:HD11	2.42	0.55
5:AY:2:PHE:O	5:AY:5:ASN:HB2	2.06	0.55
5:A1:10:LYS:CB	14:A1:103:CRT:H5	2.36	0.55
6:A2:38:LEU:C	6:A2:38:LEU:HD23	2.31	0.55
5:A5:26:ALA:O	5:A5:29:ILE:HG22	2.07	0.55
6:A6:28:TRP:C	6:A6:30:GLY:N	2.62	0.55
5:A7:35:ILE:O	5:A7:36:HIS:C	2.49	0.55
5:A7:42:THR:O	5:A7:43:ASP:C	2.50	0.55
5:A9:8:LEU:O	6:A0:18:HIS:CE1	2.60	0.55
1:BC:121:ILE:HG22	1:BC:123:THR:HG23	1.87	0.55
2:BL:87:ALA:H	2:BL:96:GLN:NE2	2.00	0.55
3:BM:248:ALA:O	3:BM:251:PHE:N	2.40	0.55
9:BF:102:BCL:H143	14:BG:102:CRT:H132	1.88	0.55
9:BI:102:BCL:OBB	9:BI:102:BCL:HHC	2.06	0.55
5:BO:43:ASP:CB	5:BQ:47:LEU:HB3	2.34	0.55
5:BU:17:PRO:HG2	5:BU:18:ARG:HD2	1.89	0.55
6:BZ:10:THR:HG22	6:BZ:11:ASP:N	2.15	0.55
9:B7:103:BCL:ND	9:B8:101:BCL:HMD3	2.22	0.55
1:AC:135:ARG:HG2	1:AC:330:LEU:C	2.32	0.55
2:AL:22:LEU:HB2	5:A7:19:ARG:CG	2.36	0.55
3:AM:104:LEU:CD1	3:AM:169:GLY:HA2	2.29	0.55
3:AM:107:PRO:CG	3:AM:113:GLY:HA2	2.36	0.55
3:AM:194:GLY:N	3:AM:293:ASN:HA	2.21	0.55
3:AM:204:LEU:C	3:AM:206:ILE:N	2.63	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:AM:248:ALA:O	3:AM:250:LEU:N	2.39	0.55
3:AM:260:VAL:CB	3:AM:264:SER:OG	2.54	0.55
5:AA:50:ASN:OD1	5:AA:51:ILE:N	2.40	0.55
5:AF:10:LYS:HB3	14:AJ:102:CRT:O1	2.07	0.55
5:AK:47:LEU:HD22	5:AK:47:LEU:N	2.22	0.55
6:AN:41:LEU:C	6:AN:41:LEU:CD2	2.69	0.55
14:AS:104:CRT:C10	6:AV:20:ILE:HD11	2.33	0.55
5:AU:25:VAL:HG13	9:AU:102:BCL:C5	2.34	0.55
6:AV:34:ILE:HG22	6:AV:38:LEU:HD21	1.88	0.55
9:AW:101:BCL:HED1	6:AX:31:LEU:O	2.07	0.55
5:AY:35:ILE:HA	5:AY:38:ILE:CG1	2.37	0.55
6:A2:24:SER:O	6:A2:27:ALA:HB3	2.06	0.55
5:A9:51:ILE:HB	5:A9:52:PRO:HA	1.89	0.55
1:BC:151:THR:HG21	1:BC:323:MET:HB2	1.87	0.55
1:BC:281:GLN:OE1	1:BC:285:TRP:HD1	1.88	0.55
2:BL:28:GLY:HA2	4:BH:46:THR:HB	1.89	0.55
2:BL:57:GLY:HA3	2:BL:66:GLN:CG	2.36	0.55
2:BL:132:PHE:HD2	2:BL:247:LEU:HD22	1.71	0.55
2:BL:192:ASN:HA	2:BL:245:LEU:CD1	2.37	0.55
6:BT:9:LEU:HB3	6:BT:13:GLU:CG	2.37	0.55
6:BT:32:VAL:O	6:BT:35:ALA:HB3	2.07	0.55
6:BX:45:TRP:O	6:BX:46:LEU:CG	2.53	0.55
6:B2:40:TRP:CZ3	6:B2:44:PRO:HA	2.42	0.55
9:B2:101:BCL:C4A	9:B3:102:BCL:HMB3	2.37	0.55
9:B5:102:BCL:H71	6:B6:28:TRP:CZ3	2.41	0.55
14:B5:103:CRT:H2M3	5:B9:36:HIS:HB2	1.89	0.55
6:B0:40:TRP:HZ3	6:B0:45:TRP:H	1.55	0.55
1:AC:233:PHE:O	1:AC:234:GLY:C	2.50	0.55
2:AL:168:ASN:C	2:AL:170:GLY:N	2.65	0.55
2:AL:206:VAL:HG12	3:AM:142:MET:SD	2.46	0.55
2:AL:228:ILE:HG21	10:AM:403:BPH:HED1	1.87	0.55
3:AM:102:TYR:CE2	3:AM:108:PRO:HD3	2.41	0.55
3:AM:175:VAL:HG22	3:AM:185:TRP:CD2	2.41	0.55
3:AM:242:GLY:C	4:AH:117:PRO:HG3	2.32	0.55
4:AH:31:ARG:HA	4:AH:34:ASP:OD2	2.06	0.55
5:AA:38:ILE:HD12	5:AA:38:ILE:C	2.32	0.55
5:AI:35:ILE:HA	5:AI:38:ILE:HG22	1.88	0.55
9:AO:102:BCL:HBD	9:AP:101:BCL:CAD	2.37	0.55
5:AY:50:ASN:CG	5:AY:51:ILE:N	2.65	0.55
9:A3:103:BCL:O1D	9:A3:103:BCL:C2A	2.52	0.55
9:A9:102:BCL:C1D	9:A0:102:BCL:HMD3	2.36	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BC:40:MET:HA	1:BC:248:THR:CG2	2.37	0.55
1:BC:148:THR:HA	1:BC:322:GLN:CG	2.37	0.55
1:BC:275:HIS:O	1:BC:278:ASP:HB3	2.06	0.55
2:BL:119:LYS:O	2:BL:121:GLY:N	2.40	0.55
2:BL:177:HIS:CD2	9:BL:301:BCL:CMC	2.90	0.55
3:BM:205:SER:CB	9:BM:402:BCL:HMA1	2.36	0.55
3:BM:248:ALA:O	3:BM:250:LEU:N	2.40	0.55
4:BH:123:CYS:N	4:BH:232:THR:HG22	2.21	0.55
5:BD:22:VAL:HA	5:BD:25:VAL:CG2	2.37	0.55
5:BF:13:LEU:HD11	6:BG:11:ASP:OD2	2.06	0.55
9:BQ:103:BCL:HBC2	9:BQ:104:BCL:HHD	1.88	0.55
5:BU:40:LEU:HD11	5:BU:47:LEU:HD23	1.89	0.55
9:BU:102:BCL:HED1	6:BV:32:VAL:HA	1.89	0.55
6:BX:17:PHE:CA	6:BX:20:ILE:HG22	2.37	0.55
1:AC:20:LEU:HD13	1:AC:20:LEU:C	2.32	0.55
1:AC:100:TRP:HE3	1:AC:152:CYS:SG	2.30	0.55
1:AC:135:ARG:NH1	1:AC:332:LYS:HA	2.15	0.55
1:AC:311:HIS:ND1	1:AC:317:PRO:HD3	2.21	0.55
3:AM:98:PRO:HB2	3:AM:171:TRP:HB3	1.88	0.55
5:AA:34:LEU:O	5:AA:38:ILE:HG23	2.07	0.55
6:AN:10:THR:HB	6:AN:13:GLU:OE2	2.07	0.55
6:AT:9:LEU:HB3	6:AT:13:GLU:CG	2.37	0.55
5:AU:40:LEU:HD11	5:AU:47:LEU:HD23	1.89	0.55
5:AW:45:ASN:O	5:AW:49:ASP:HB3	2.07	0.55
6:AX:34:ILE:HD13	6:AX:34:ILE:O	2.07	0.55
6:A6:38:LEU:HD23	6:A6:38:LEU:C	2.32	0.55
2:BL:18:ILE:HG23	4:BH:259:LEU:HB2	1.89	0.55
2:BL:142:PHE:CD1	2:BL:143:VAL:N	2.74	0.55
3:BM:228:ARG:HD2	3:BM:228:ARG:N	2.22	0.55
3:BM:237:GLN:OE1	3:BM:244:ALA:HB3	2.06	0.55
14:BA:102:CRT:H401	5:BD:38:ILE:HD13	1.89	0.55
6:BJ:10:THR:HG22	6:BJ:11:ASP:N	2.16	0.55
5:BO:44:LEU:HD11	5:BO:46:TRP:CE3	2.42	0.55
6:BP:21:PHE:O	6:BP:22:MET:C	2.50	0.55
5:BQ:19:ARG:NH1	15:BQ:101:PEF:C5	2.70	0.55
6:BR:30:GLY:O	6:BR:34:ILE:HG23	2.07	0.55
5:BS:26:ALA:C	5:BS:29:ILE:HG22	2.32	0.55
5:BW:10:LYS:HB3	14:BW:103:CRT:H5	1.87	0.55
5:BW:12:TRP:HZ2	6:BX:21:PHE:CD2	2.25	0.55
6:BX:28:TRP:HA	6:BX:31:LEU:HD12	1.89	0.55
5:B5:45:ASN:O	5:B5:47:LEU:N	2.39	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:B6:30:GLY:O	6:B6:34:ILE:HG22	2.07	0.55
1:AC:48:GLN:C	1:AC:50:ALA:N	2.64	0.54
2:AL:49:LEU:HD21	5:A9:37:MET:HG2	1.89	0.54
2:AL:204:LEU:CD2	3:AM:267:ARG:HD2	2.36	0.54
3:AM:35:ILE:O	3:AM:48:ILE:N	2.39	0.54
3:AM:83:VAL:HA	3:AM:86:PHE:HB3	1.90	0.54
3:AM:151:ALA:O	3:AM:155:PHE:N	2.37	0.54
3:AM:179:ILE:N	3:AM:179:ILE:CD1	2.70	0.54
3:AM:240:HIS:O	3:AM:241:ARG:O	2.25	0.54
3:AM:301:HIS:CE1	4:AH:8:TYR:HD2	2.25	0.54
10:AM:403:BPH:H112	15:AS:101:PEF:H162	1.90	0.54
4:AH:19:PHE:CD1	4:AH:19:PHE:C	2.85	0.54
4:AH:28:ILE:O	4:AH:29:TYR:C	2.49	0.54
4:AH:52:ARG:HB2	4:AH:54:LYS:HZ3	1.70	0.54
5:AA:11:ILE:HD12	5:AA:14:ILE:HD11	1.88	0.54
9:AO:102:BCL:CBD	9:AP:101:BCL:OBD	2.54	0.54
9:AO:102:BCL:H62	6:AP:28:TRP:CZ3	2.42	0.54
5:AU:8:LEU:HB2	6:AV:18:HIS:CE1	2.43	0.54
5:A1:12:TRP:CZ2	6:A2:21:PHE:CE2	2.95	0.54
6:A0:34:ILE:HD13	6:A0:34:ILE:C	2.32	0.54
1:BC:47:ARG:HG2	1:BC:47:ARG:O	2.07	0.54
1:BC:71:LYS:HD3	1:BC:74:GLU:HG2	1.87	0.54
3:BM:74:ASN:CG	3:BM:95:LEU:HD13	2.32	0.54
5:BA:38:ILE:C	5:BA:38:ILE:HD12	2.31	0.54
6:BB:20:ILE:HG13	5:B9:7:ASN:O	2.06	0.54
6:BB:44:PRO:HG2	5:BD:52:PRO:HB2	1.89	0.54
6:BG:30:GLY:HA2	6:BG:33:VAL:HG12	1.89	0.54
6:B0:28:TRP:O	6:B0:31:LEU:HB2	2.07	0.54
1:AC:245:VAL:HG21	1:AC:249:PHE:CG	2.43	0.54
3:AM:78:SER:O	3:AM:80:HIS:N	2.36	0.54
4:AH:181:TYR:CD1	4:AH:196:PRO:HA	2.42	0.54
5:AO:36:HIS:NE2	9:AP:101:BCL:CMD	2.70	0.54
5:AO:44:LEU:HD12	5:AO:46:TRP:N	2.21	0.54
6:AT:40:TRP:CZ3	6:AT:44:PRO:HA	2.42	0.54
5:AU:12:TRP:NE1	6:AV:18:HIS:CA	2.62	0.54
5:AU:12:TRP:CZ2	6:AV:21:PHE:HD2	2.25	0.54
5:AU:46:TRP:CZ3	9:AU:102:BCL:HAC1	2.42	0.54
9:A1:102:BCL:C4D	9:A2:101:BCL:HMD3	2.37	0.54
5:A5:32:GLY:HA3	9:A5:102:BCL:O1A	2.08	0.54
1:BC:259:TRP:C	1:BC:261:GLN:N	2.59	0.54
2:BL:44:LEU:C	2:BL:46:GLY:N	2.66	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:BL:228:ILE:O	3:BM:51:ILE:HD11	2.05	0.54
3:BM:177:PHE:CZ	14:BM:406:CRT:H25	2.42	0.54
4:BH:94:PRO:CG	6:B0:8:GLY:HA3	2.31	0.54
4:BH:154:MET:O	4:BH:167:VAL:HG13	2.08	0.54
5:BK:34:LEU:O	5:BK:37:MET:HB2	2.07	0.54
5:BK:38:ILE:O	5:BK:42:THR:OG1	2.20	0.54
14:BP:102:CRT:C2M	5:BQ:36:HIS:HB2	2.34	0.54
5:BQ:46:TRP:CZ2	9:BQ:103:BCL:HHC	2.42	0.54
9:BZ:101:BCL:CHB	9:B1:102:BCL:HMB2	2.38	0.54
6:B4:46:LEU:HB2	5:B5:52:PRO:CD	2.36	0.54
5:B5:44:LEU:C	5:B5:46:TRP:H	2.16	0.54
5:B9:26:ALA:O	5:B9:29:ILE:CG2	2.55	0.54
9:B9:102:BCL:C1D	9:B0:102:BCL:CMD	2.85	0.54
1:AC:156:HIS:O	1:AC:157:ARG:C	2.50	0.54
1:AC:236:MET:SD	7:AC:503:HEM:NA	2.80	0.54
3:AM:71:ILE:HD13	3:AM:177:PHE:CD1	2.42	0.54
3:AM:202:HIS:C	3:AM:204:LEU:H	2.15	0.54
3:AM:316:PRO:HG2	3:AM:317:TYR:HD1	1.71	0.54
5:AD:46:TRP:CD1	5:AD:47:LEU:HD22	2.42	0.54
6:AJ:17:PHE:CE1	14:AJ:102:CRT:H9	2.43	0.54
5:AS:21:LEU:O	5:AS:25:VAL:HG23	2.06	0.54
5:AS:27:PHE:CZ	5:AU:29:ILE:HG13	2.41	0.54
5:AS:40:LEU:CD1	5:AS:47:LEU:HD23	2.35	0.54
5:A1:14:ILE:O	5:A3:18:ARG:NH1	2.40	0.54
1:BC:54:GLN:HA	1:BC:54:GLN:HE21	1.73	0.54
1:BC:205:ASP:OD1	1:BC:304:ARG:NE	2.41	0.54
1:BC:251:HIS:HE1	1:BC:256:PHE:O	1.90	0.54
2:BL:87:ALA:HB3	2:BL:96:GLN:NE2	2.23	0.54
2:BL:104:GLY:HA2	2:BL:107:ILE:HB	1.88	0.54
2:BL:117:CYS:HB3	2:BL:124:PHE:CE2	2.42	0.54
5:BA:36:HIS:O	5:BA:40:LEU:HB3	2.07	0.54
5:BA:40:LEU:HB2	5:BA:46:TRP:CH2	2.42	0.54
5:BD:33:LEU:O	5:BD:37:MET:HG2	2.08	0.54
5:BQ:42:THR:OG1	5:BS:48:ASP:OD1	2.25	0.54
9:BS:102:BCL:OBD	6:BT:32:VAL:HG23	2.07	0.54
5:B1:14:ILE:HD12	5:B1:15:LEU:HG	1.89	0.54
6:B2:46:LEU:OXT	6:B4:43:ARG:NH2	2.39	0.54
9:B2:101:BCL:C1B	9:B3:102:BCL:CMB	2.84	0.54
5:B3:42:THR:O	5:B3:43:ASP:C	2.51	0.54
9:B4:101:BCL:CHC	9:B5:102:BCL:HBB3	2.38	0.54
6:B8:36:HIS:HD1	9:B8:101:BCL:H141	1.71	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AC:128:ARG:HE	7:AC:501:HEM:HAD1	1.73	0.54
1:AC:253:THR:HG22	2:AL:171:TYR:HD2	1.71	0.54
1:AC:326:ASP:C	1:AC:327:TYR:CD1	2.86	0.54
2:AL:48:LEU:HD13	5:AA:33:LEU:HD23	1.89	0.54
3:AM:84:PHE:CE1	5:AW:37:MET:HG2	2.42	0.54
3:AM:170:SER:C	3:AM:172:ALA:N	2.66	0.54
3:AM:182:HIS:HD1	3:AM:183:LEU:N	2.05	0.54
10:AM:403:BPH:C9	15:AM:409:PEF:C22	2.85	0.54
5:AA:11:ILE:HD11	5:AA:14:ILE:HD11	1.89	0.54
5:AA:18:ARG:O	5:AA:22:VAL:HG12	2.07	0.54
5:AA:36:HIS:CE1	9:AB:101:BCL:HMD3	2.42	0.54
5:AA:45:ASN:HB3	5:AA:49:ASP:HB3	1.90	0.54
5:AF:4:MET:SD	6:AJ:23:GLN:HG3	2.48	0.54
9:AF:102:BCL:C4D	9:AG:101:BCL:HMD3	2.38	0.54
5:AI:29:ILE:HA	9:AI:102:BCL:H11	1.89	0.54
5:AO:31:LEU:CD2	14:AP:102:CRT:H32	2.38	0.54
5:AO:46:TRP:CZ3	9:AO:102:BCL:HBC3	2.41	0.54
14:AS:104:CRT:H342	9:AW:101:BCL:CBA	2.32	0.54
5:AU:45:ASN:OD1	5:AU:47:LEU:HB2	2.07	0.54
9:AU:102:BCL:HMB3	9:AU:102:BCL:HBB2	1.89	0.54
5:A1:57:ALA:C	5:A1:59:GLY:N	2.64	0.54
6:A4:43:ARG:O	6:A4:45:TRP:N	2.40	0.54
5:A5:18:ARG:HB2	5:A5:19:ARG:NH1	2.22	0.54
6:A6:30:GLY:O	6:A6:34:ILE:HG22	2.06	0.54
2:BL:246:ALA:CB	3:BM:217:ALA:HB2	2.37	0.54
3:BM:64:GLY:C	3:BM:66:VAL:H	2.15	0.54
3:BM:223:ILE:HG22	3:BM:224:LEU:N	2.22	0.54
9:BA:101:BCL:HMD1	6:BB:36:HIS:ND1	2.22	0.54
5:BD:9:TYR:CE1	6:BE:11:ASP:CB	2.88	0.54
6:BE:21:PHE:CZ	9:BF:102:BCL:H203	2.42	0.54
5:BF:16:ASP:O	5:BF:20:VAL:HG22	2.06	0.54
6:BJ:17:PHE:CE1	6:BJ:21:PHE:HB2	2.38	0.54
5:BK:16:ASP:CB	5:BK:18:ARG:HE	2.21	0.54
6:BT:30:GLY:O	6:BT:34:ILE:HG12	2.08	0.54
5:BW:35:ILE:HG22	5:BW:36:HIS:N	2.23	0.54
5:B1:20:VAL:O	5:B1:24:ILE:HG12	2.07	0.54
9:B1:102:BCL:HMB3	9:B1:102:BCL:HBB2	1.88	0.54
6:B4:18:HIS:CD2	6:B4:22:MET:HB2	2.43	0.54
6:B6:38:LEU:HD23	6:B6:38:LEU:C	2.32	0.54
6:B0:34:ILE:HD13	6:B0:34:ILE:C	2.32	0.54
1:AC:91:THR:O	1:AC:92:ARG:C	2.51	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AC:249:PHE:CD1	1:AC:250:CYS:SG	2.98	0.54
1:AC:265:LYS:HE2	7:AC:504:HEM:O2A	2.06	0.54
3:AM:83:VAL:HG23	3:AM:84:PHE:N	2.22	0.54
3:AM:90:PHE:HA	3:AM:93:LEU:HD12	1.90	0.54
4:AH:30:LEU:O	4:AH:31:ARG:C	2.49	0.54
6:AJ:45:TRP:CZ3	9:AJ:101:BCL:HAC2	2.43	0.54
5:AK:16:ASP:O	5:AK:19:ARG:HG2	2.08	0.54
5:AW:25:VAL:O	5:AW:29:ILE:HB	2.08	0.54
5:AW:31:LEU:CD1	14:AX:102:CRT:H35	2.38	0.54
14:AX:102:CRT:H2M1	5:AY:36:HIS:HB2	1.89	0.54
5:A1:10:LYS:HD2	6:A4:20:ILE:CG1	2.38	0.54
5:A1:27:PHE:C	5:A1:27:PHE:CD1	2.85	0.54
9:A7:103:BCL:O1A	9:A7:103:BCL:H12	2.08	0.54
1:BC:90:PHE:O	1:BC:93:THR:HB	2.08	0.54
1:BC:102:SER:C	1:BC:104:LYS:H	2.15	0.54
3:BM:235:ILE:HD12	3:BM:235:ILE:N	2.14	0.54
4:BH:133:ILE:HD11	4:BH:171:TRP:HB3	1.89	0.54
5:BA:44:LEU:C	5:BA:44:LEU:HD12	2.31	0.54
9:BA:101:BCL:CBB	9:B0:102:BCL:CHC	2.85	0.54
5:BF:36:HIS:O	5:BF:40:LEU:N	2.40	0.54
5:BI:9:TYR:CB	6:BJ:15:LYS:HA	2.32	0.54
5:BQ:43:ASP:H	5:BS:47:LEU:HB3	1.73	0.54
9:BQ:104:BCL:HAC2	6:BR:45:TRP:CE3	2.43	0.54
5:BY:16:ASP:CB	5:BY:18:ARG:HE	2.18	0.54
9:B1:102:BCL:C4D	9:B2:101:BCL:HMD3	2.37	0.54
5:B7:44:LEU:O	5:B7:44:LEU:HD13	2.07	0.54
6:B8:40:TRP:HH2	6:B8:46:LEU:HD12	1.73	0.54
5:B9:12:TRP:HA	5:B9:12:TRP:CE3	2.41	0.54
6:B0:21:PHE:CD1	14:B0:101:CRT:C16	2.91	0.54
1:AC:203:PHE:CD1	1:AC:235:LEU:HD22	2.42	0.54
1:AC:265:LYS:O	1:AC:266:ARG:C	2.50	0.54
2:AL:75:ILE:HG22	2:AL:95:TRP:CD1	2.42	0.54
3:AM:14:ARG:HG3	3:AM:14:ARG:NH1	2.22	0.54
3:AM:260:VAL:HG13	13:AM:405:MQ8:H142	1.89	0.54
13:AM:405:MQ8:H2M3	13:AM:405:MQ8:C12	2.38	0.54
4:AH:16:ILE:HD13	4:AH:16:ILE:C	2.32	0.54
4:AH:27:ILE:HG23	4:AH:28:ILE:N	2.23	0.54
6:AB:34:ILE:O	6:AB:34:ILE:HD13	2.07	0.54
5:AI:49:ASP:OD2	6:AJ:43:ARG:NH2	2.41	0.54
5:AK:26:ALA:O	5:AK:29:ILE:HG22	2.07	0.54
5:AO:12:TRP:CD2	6:AP:17:PHE:HD2	2.25	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
14:AS:104:CRT:H16	6:AV:21:PHE:CE1	2.43	0.54
6:AT:20:ILE:HD13	6:AT:20:ILE:C	2.32	0.54
6:AT:32:VAL:O	6:AT:35:ALA:HB3	2.07	0.54
5:AW:35:ILE:HA	5:AW:38:ILE:HG22	1.90	0.54
6:A4:10:THR:HB	6:A4:13:GLU:OE2	2.08	0.54
6:A8:40:TRP:HH2	6:A8:46:LEU:HD12	1.73	0.54
6:A0:28:TRP:O	6:A0:31:LEU:HB2	2.07	0.54
2:BL:29:PRO:O	3:BM:254:TRP:HA	2.08	0.54
2:BL:228:ILE:HG23	3:BM:132:ARG:HD2	1.90	0.54
5:BA:40:LEU:HD11	5:BA:47:LEU:HD23	1.90	0.54
5:BF:9:TYR:HA	6:BG:18:HIS:CE1	2.43	0.54
9:BI:102:BCL:HMB1	9:BI:102:BCL:CBB	2.37	0.54
5:BS:28:GLN:C	9:BS:102:BCL:H11	2.30	0.54
6:B2:17:PHE:HA	14:B2:102:CRT:C6	2.38	0.54
2:AL:44:LEU:HD21	13:AM:405:MQ8:H391	1.90	0.54
2:AL:196:LEU:CG	3:AM:216:PHE:HB2	2.37	0.54
3:AM:202:HIS:C	3:AM:204:LEU:N	2.65	0.54
3:AM:250:LEU:O	3:AM:253:ARG:HB3	2.07	0.54
5:AD:12:TRP:HE1	6:AE:18:HIS:HA	1.71	0.54
5:AF:7:ASN:CB	6:AJ:20:ILE:HD13	2.38	0.54
14:AS:104:CRT:H14	6:AV:21:PHE:HD1	1.73	0.54
6:AT:45:TRP:O	6:AT:46:LEU:HB2	2.08	0.54
6:AX:22:MET:HG3	6:AX:26:TYR:CE2	2.39	0.54
6:A8:44:PRO:HG2	5:A9:52:PRO:HB2	1.90	0.54
1:BC:53:ILE:C	1:BC:55:ALA:N	2.63	0.54
3:BM:8:PHE:HB3	3:BM:42:LYS:HG2	1.89	0.54
3:BM:98:PRO:CG	3:BM:107:PRO:HG3	2.37	0.54
3:BM:170:SER:O	3:BM:172:ALA:N	2.39	0.54
9:BA:101:BCL:HMB3	9:B0:102:BCL:C1B	2.38	0.54
5:BK:45:ASN:O	5:BK:49:ASP:HB3	2.08	0.54
14:BU:103:CRT:H291	9:BY:102:BCL:HAA2	1.90	0.54
5:BW:16:ASP:HB2	5:BW:19:ARG:HG2	1.89	0.54
5:BY:27:PHE:CD1	5:BY:27:PHE:C	2.86	0.54
6:B2:21:PHE:HE1	14:B2:102:CRT:H19	1.70	0.54
5:B3:12:TRP:CD1	6:B4:18:HIS:HB2	2.43	0.54
5:B7:32:GLY:N	9:B8:101:BCL:HED2	2.23	0.54
1:AC:249:PHE:CE1	1:AC:265:LYS:HG2	2.42	0.54
2:AL:15:GLY:O	2:AL:118:ARG:HD3	2.08	0.54
2:AL:196:LEU:HD21	3:AM:269:ALA:HB1	1.88	0.54
5:AF:11:ILE:CA	14:AJ:102:CRT:H82	2.37	0.54
5:AI:49:ASP:OD1	5:AI:49:ASP:N	2.40	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:AO:4:MET:SD	5:AO:4:MET:N	2.81	0.54
6:AP:23:GLN:O	6:AP:24:SER:C	2.51	0.54
9:AP:101:BCL:HHC	9:AP:101:BCL:OBB	2.08	0.54
9:AR:101:BCL:CMA	9:AS:103:BCL:HMA1	2.35	0.54
14:AS:104:CRT:H2M3	5:AW:37:MET:N	2.23	0.54
5:AU:9:TYR:HB2	6:AV:15:LYS:HD2	1.89	0.54
5:A3:27:PHE:CD1	5:A3:27:PHE:C	2.85	0.54
1:BC:195:LEU:HB3	1:BC:196:PRO:HD2	1.88	0.54
2:BL:37:VAL:HG23	2:BL:38:VAL:N	2.23	0.54
2:BL:180:PRO:O	2:BL:183:MET:HB2	2.08	0.54
2:BL:192:ASN:N	2:BL:245:LEU:HD13	2.23	0.54
3:BM:215:LEU:HA	3:BM:218:MET:CG	2.37	0.54
3:BM:244:ALA:C	3:BM:246:GLU:N	2.66	0.54
5:BD:17:PRO:O	5:BD:21:LEU:HB2	2.08	0.54
5:BO:10:LYS:CB	14:BO:103:CRT:H5	2.37	0.54
5:BW:9:TYR:CD1	6:BX:15:LYS:HB2	2.43	0.54
6:BX:46:LEU:HD13	6:BZ:42:TYR:OH	2.07	0.54
14:B2:102:CRT:C2M	5:B3:36:HIS:CB	2.71	0.54
5:B5:10:LYS:HB3	14:B5:103:CRT:C5	2.31	0.54
1:AC:66:ASP:C	1:AC:68:THR:H	2.16	0.54
1:AC:70:PRO:HG2	1:AC:71:LYS:H	1.72	0.54
1:AC:137:ALA:O	1:AC:139:SER:N	2.40	0.54
1:AC:231:TRP:O	1:AC:232:THR:C	2.50	0.54
2:AL:129:ALA:HB1	2:AL:247:LEU:HD21	1.90	0.54
2:AL:183:MET:HA	9:AM:401:BCL:OBD	2.07	0.54
2:AL:203:ILE:O	2:AL:205:SER:N	2.40	0.54
3:AM:4:TYR:HE2	3:AM:10:ALA:HB2	1.73	0.54
3:AM:53:LEU:HG	3:AM:58:THR:HG23	1.90	0.54
3:AM:114:TRP:HZ3	3:AM:117:MET:HE2	1.73	0.54
3:AM:138:GLU:HA	3:AM:142:MET:O	2.08	0.54
4:AH:18:ALA:O	4:AH:21:LEU:HB3	2.08	0.54
4:AH:119:ARG:NE	4:AH:237:ASP:OD2	2.41	0.54
5:AA:50:ASN:HD21	5:AA:51:ILE:HG12	1.71	0.54
5:AD:46:TRP:NE1	9:AD:102:BCL:HHC	2.23	0.54
5:AK:9:TYR:C	5:AK:9:TYR:CD1	2.86	0.54
5:AK:9:TYR:CD2	6:AN:15:LYS:HG3	2.43	0.54
14:AS:104:CRT:H2M3	5:AW:36:HIS:C	2.33	0.54
5:AU:35:ILE:HA	5:AU:38:ILE:HG22	1.90	0.54
5:AW:10:LYS:HZ3	14:AW:102:CRT:H1M2	1.71	0.54
9:AW:101:BCL:C1D	9:AX:101:BCL:HMD3	2.37	0.54
6:AX:30:GLY:O	6:AX:34:ILE:HG22	2.08	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
14:A5:103:CRT:C40	5:A7:38:ILE:HG21	2.37	0.54
9:A6:101:BCL:HMC1	9:A7:103:BCL:CBB	2.33	0.54
5:A7:33:LEU:H	5:A7:33:LEU:HD12	1.73	0.54
1:BC:48:GLN:C	1:BC:50:ALA:N	2.66	0.54
3:BM:103:GLY:O	3:BM:104:LEU:HD13	2.07	0.54
3:BM:190:SER:HA	3:BM:196:LEU:HG	1.90	0.54
3:BM:204:LEU:O	3:BM:206:ILE:N	2.41	0.54
4:BH:80:ARG:HH11	4:BH:80:ARG:HG3	1.72	0.54
4:BH:123:CYS:H	4:BH:232:THR:HG22	1.73	0.54
6:BB:46:LEU:O	5:BD:51:ILE:O	2.26	0.54
5:BD:50:ASN:CG	5:BD:51:ILE:N	2.66	0.54
6:BE:45:TRP:CZ3	9:BE:101:BCL:HAC2	2.42	0.54
6:BJ:20:ILE:HG23	6:BJ:21:PHE:N	2.23	0.54
5:BK:36:HIS:CE1	9:BK:102:BCL:NA	2.75	0.54
6:BN:46:LEU:HB3	6:BP:42:TYR:OH	2.08	0.54
14:BN:102:CRT:H2M3	5:BO:36:HIS:CB	2.38	0.54
5:BS:26:ALA:O	5:BS:30:VAL:HG12	2.08	0.54
5:B5:36:HIS:CE1	9:B6:101:BCL:CMD	2.89	0.54
2:AL:6:PHE:CD2	3:AM:246:GLU:HG3	2.43	0.54
2:AL:280:LEU:HD21	5:AY:37:MET:SD	2.48	0.54
3:AM:199:ASN:HB2	3:AM:294:TRP:CG	2.43	0.54
3:AM:264:SER:O	3:AM:265:ILE:C	2.50	0.54
4:AH:219:PHE:HA	4:AH:222:VAL:HG23	1.89	0.54
5:AF:43:ASP:O	5:AF:44:LEU:HG	2.08	0.54
6:AN:31:LEU:HA	6:AN:34:ILE:HG22	1.89	0.54
9:AU:102:BCL:C1D	9:AV:102:BCL:CMD	2.85	0.54
5:AY:43:ASP:HA	5:A1:48:ASP:HB3	1.90	0.54
5:A1:12:TRP:HD1	6:A2:18:HIS:HB2	1.73	0.54
9:A1:102:BCL:H91	14:A2:102:CRT:H183	1.89	0.54
5:A7:46:TRP:HH2	9:A7:103:BCL:HBC3	1.69	0.54
2:BL:167:SER:HA	9:BL:301:BCL:HBC1	1.90	0.54
2:BL:177:HIS:CE1	9:BM:402:BCL:HAC2	2.43	0.54
3:BM:126:ILE:HD13	9:BM:402:BCL:H91	1.90	0.54
3:BM:200:PRO:HD2	3:BM:294:TRP:CZ3	2.43	0.54
4:BH:5:ILE:HD11	5:BF:40:LEU:HD11	1.88	0.54
4:BH:182:LEU:HD12	4:BH:195:LEU:O	2.08	0.54
5:BA:22:VAL:HA	5:BA:25:VAL:HG23	1.90	0.54
5:BA:33:LEU:H	5:BA:33:LEU:CD1	2.21	0.54
5:BU:12:TRP:CZ2	6:BV:21:PHE:HD2	2.16	0.54
5:B7:10:LYS:H	5:B7:10:LYS:HD2	1.73	0.54
9:B9:102:BCL:HMD1	6:B0:36:HIS:HA	1.90	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AC:70:PRO:C	1:AC:71:LYS:HD2	2.32	0.53
1:AC:225:SER:HB3	1:AC:228:GLN:HE21	1.73	0.53
1:AC:226:LEU:H	3:AM:173:LYS:HE3	1.74	0.53
1:AC:248:THR:O	1:AC:251:HIS:O	2.26	0.53
2:AL:87:ALA:HB3	2:AL:96:GLN:NE2	2.23	0.53
2:AL:193:CYS:O	10:AM:403:BPH:H3C	2.07	0.53
2:AL:244:PHE:O	2:AL:245:LEU:C	2.50	0.53
3:AM:35:ILE:HD11	15:AM:409:PEF:C32	2.37	0.53
3:AM:55:LEU:HD23	5:AQ:22:VAL:HG23	1.89	0.53
3:AM:70:ILE:C	3:AM:72:GLY:N	2.65	0.53
3:AM:84:PHE:HA	5:AW:37:MET:HE2	1.90	0.53
3:AM:221:ALA:O	3:AM:222:THR:C	2.51	0.53
4:AH:215:LYS:H	4:AH:218:HIS:HD2	1.56	0.53
5:AI:15:LEU:HD22	5:AI:15:LEU:N	2.23	0.53
5:AS:9:TYR:HA	6:AT:18:HIS:CG	2.43	0.53
5:AU:14:ILE:HB	14:AX:102:CRT:H82	1.89	0.53
6:AX:10:THR:H	6:AX:13:GLU:CD	2.16	0.53
5:AY:10:LYS:NZ	14:A2:102:CRT:H1M2	2.23	0.53
5:AY:12:TRP:HE1	6:AZ:18:HIS:CA	2.15	0.53
5:A3:19:ARG:O	5:A3:23:SER:CB	2.42	0.53
5:A7:36:HIS:HB3	14:A7:102:CRT:C39	2.38	0.53
6:A8:43:ARG:NH2	5:A9:55:TYR:CB	2.69	0.53
1:BC:122:TYR:HB2	1:BC:290:VAL:HB	1.89	0.53
1:BC:190:VAL:C	1:BC:192:TYR:H	2.16	0.53
1:BC:243:LEU:H	1:BC:243:LEU:CD1	2.20	0.53
2:BL:166:VAL:HG13	9:BL:301:BCL:CHD	2.38	0.53
3:BM:41:GLY:HA2	3:BM:44:GLY:O	2.08	0.53
3:BM:59:LEU:CD1	5:BQ:29:ILE:HG21	2.34	0.53
3:BM:244:ALA:O	3:BM:246:GLU:N	2.41	0.53
3:BM:261:THR:N	3:BM:264:SER:OG	2.37	0.53
3:BM:268:TRP:CD1	4:BH:30:LEU:HD22	2.42	0.53
6:BG:20:ILE:O	6:BG:24:SER:OG	2.26	0.53
9:BG:101:BCL:C2B	9:BI:102:BCL:C2B	2.86	0.53
9:BI:102:BCL:HBC2	9:BI:102:BCL:CHD	2.38	0.53
5:BK:16:ASP:HB3	5:BK:18:ARG:NE	2.23	0.53
5:BK:43:ASP:OD1	5:BO:47:LEU:HB3	2.08	0.53
6:BN:38:LEU:HA	6:BN:41:LEU:HD12	1.89	0.53
6:BP:45:TRP:CE2	9:BP:101:BCL:H2C	2.43	0.53
5:BY:27:PHE:HE1	5:BY:31:LEU:HD22	1.73	0.53
5:B7:42:THR:O	5:B7:43:ASP:C	2.50	0.53
5:B9:44:LEU:HD22	5:B9:46:TRP:HB3	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AC:157:ARG:NE	1:AC:312:GLN:HE22	2.06	0.53
2:AL:273:ASN:O	2:AL:274:TRP:C	2.51	0.53
3:AM:91:PHE:O	3:AM:180:PHE:HB2	2.09	0.53
3:AM:156:PHE:CD1	3:AM:281:GLY:N	2.76	0.53
3:AM:191:ILE:O	3:AM:193:TYR:N	2.41	0.53
6:AJ:17:PHE:HE1	6:AJ:21:PHE:HB2	1.73	0.53
6:AJ:21:PHE:CE1	14:AJ:102:CRT:H16	2.43	0.53
5:AQ:31:LEU:O	5:AQ:34:LEU:HB3	2.08	0.53
5:AS:39:VAL:O	5:AS:42:THR:HB	2.08	0.53
6:AT:45:TRP:CD1	6:AT:46:LEU:H	2.26	0.53
6:AV:43:ARG:NH1	5:AW:55:TYR:HB3	2.23	0.53
6:AX:28:TRP:HE3	6:AX:31:LEU:HD12	1.73	0.53
5:AY:8:LEU:HD13	6:AZ:22:MET:HE1	1.90	0.53
5:AY:51:ILE:HA	5:AY:52:PRO:C	2.33	0.53
5:A1:43:ASP:HB2	5:A3:47:LEU:HD12	1.90	0.53
5:A1:46:TRP:O	5:A1:49:ASP:OD1	2.26	0.53
5:A5:29:ILE:O	5:A5:29:ILE:HD13	2.07	0.53
1:BC:176:SER:OG	5:BS:42:THR:HA	2.08	0.53
2:BL:10:TYR:CE1	4:BH:115:ALA:HB3	2.43	0.53
2:BL:177:HIS:NE2	9:BL:301:BCL:HMC2	2.23	0.53
2:BL:255:VAL:O	2:BL:257:ILE:N	2.42	0.53
2:BL:276:LEU:H	2:BL:276:LEU:CD2	2.18	0.53
3:BM:271:TRP:NE1	4:BH:26:LEU:HD11	2.23	0.53
4:BH:242:TYR:O	4:BH:243:TYR:C	2.50	0.53
5:BI:50:ASN:CG	5:BI:51:ILE:N	2.57	0.53
6:BJ:42:TYR:CE2	6:BJ:43:ARG:HG3	2.43	0.53
5:BQ:43:ASP:HB2	5:BS:47:LEU:CB	2.38	0.53
5:BU:29:ILE:N	9:BU:102:BCL:H43	2.23	0.53
6:BV:28:TRP:HA	6:BV:31:LEU:HB2	1.90	0.53
5:B7:36:HIS:CB	14:B7:102:CRT:H393	2.37	0.53
1:AC:285:TRP:CB	1:AC:286:PRO:HD3	2.39	0.53
2:AL:21:ASP:HB3	5:A7:19:ARG:NE	2.23	0.53
2:AL:142:PHE:CD1	2:AL:143:VAL:N	2.76	0.53
2:AL:148:MET:CE	2:AL:262:PRO:HD3	2.38	0.53
3:AM:56:THR:HG21	3:AM:131:VAL:CG1	2.39	0.53
3:AM:63:PHE:HZ	5:AQ:33:LEU:CD2	2.12	0.53
3:AM:73:PHE:HA	14:AM:406:CRT:H1M1	1.90	0.53
3:AM:153:ALA:HA	3:AM:277:VAL:HG11	1.91	0.53
3:AM:240:HIS:CD2	4:AH:120:PRO:HG2	2.43	0.53
3:AM:248:ALA:O	3:AM:249:ALA:C	2.51	0.53
4:AH:171:TRP:CZ3	4:AH:231:VAL:HG12	2.43	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:AA:38:ILE:O	5:AA:41:SER:HB3	2.08	0.53
5:AK:44:LEU:O	5:AK:44:LEU:HD13	2.08	0.53
9:AN:101:BCL:CGD	9:AN:101:BCL:H2A	2.37	0.53
5:AO:12:TRP:CH2	6:AP:17:PHE:HE2	2.26	0.53
5:AO:36:HIS:O	5:AO:40:LEU:CB	2.54	0.53
5:AO:46:TRP:CD1	5:AO:47:LEU:HD22	2.43	0.53
5:AW:19:ARG:HH12	5:AY:22:VAL:HG23	1.71	0.53
9:A3:104:BCL:C6	6:A4:29:PHE:HE1	2.20	0.53
6:A6:17:PHE:CD1	6:A6:17:PHE:C	2.87	0.53
1:BC:204:LEU:HD22	7:BC:504:HEM:HBB1	1.91	0.53
2:BL:49:LEU:HD12	2:BL:98:ILE:HG13	1.91	0.53
3:BM:316:PRO:HG2	3:BM:317:TYR:CD1	2.44	0.53
4:BH:53:VAL:O	4:BH:53:VAL:HG13	2.08	0.53
4:BH:54:LYS:HA	4:BH:54:LYS:CE	2.34	0.53
6:BG:34:ILE:C	6:BG:34:ILE:HD12	2.34	0.53
5:BI:35:ILE:CA	5:BI:38:ILE:HG22	2.35	0.53
5:BO:7:ASN:ND2	6:BR:20:ILE:HD12	2.18	0.53
5:BS:45:ASN:O	5:BS:49:ASP:CG	2.51	0.53
14:BS:103:CRT:H2M1	5:BU:37:MET:HG2	1.90	0.53
5:BW:8:LEU:HD21	6:BZ:24:SER:OG	2.08	0.53
9:B1:102:BCL:HMB3	9:B1:102:BCL:CBB	2.38	0.53
6:B2:17:PHE:O	6:B2:20:ILE:HG22	2.08	0.53
6:B0:9:LEU:HB3	6:B0:13:GLU:CG	2.38	0.53
2:AL:48:LEU:O	2:AL:51:VAL:HB	2.09	0.53
2:AL:203:ILE:C	2:AL:205:SER:N	2.66	0.53
3:AM:214:LEU:HD22	3:AM:215:LEU:CD1	2.38	0.53
4:AH:24:PHE:O	4:AH:27:ILE:HG22	2.08	0.53
4:AH:126:THR:HG23	4:AH:130:LEU:O	2.08	0.53
5:AA:40:LEU:HD11	5:AA:47:LEU:HD23	1.90	0.53
6:AE:10:THR:N	6:AE:13:GLU:OE2	2.41	0.53
6:AE:27:ALA:O	6:AE:31:LEU:HG	2.09	0.53
5:AI:40:LEU:HD11	5:AI:46:TRP:CH2	2.42	0.53
5:AO:45:ASN:O	5:AO:47:LEU:N	2.40	0.53
6:AP:46:LEU:HD22	6:AR:42:TYR:OH	2.08	0.53
6:AR:20:ILE:C	6:AR:20:ILE:HD13	2.33	0.53
5:AS:47:LEU:HD22	5:AS:47:LEU:N	2.21	0.53
5:AW:27:PHE:HE2	5:AY:29:ILE:HG12	1.72	0.53
14:AW:102:CRT:H14	6:AZ:21:PHE:CD2	2.43	0.53
5:AY:43:ASP:HA	5:A1:48:ASP:HA	1.91	0.53
6:AZ:44:PRO:O	5:A1:55:TYR:OH	2.26	0.53
5:A1:46:TRP:CD1	5:A1:47:LEU:N	2.77	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:A5:17:PRO:O	5:A5:21:LEU:HB2	2.08	0.53
5:A9:12:TRP:NE1	6:A0:18:HIS:ND1	2.56	0.53
1:BC:18:VAL:HG22	1:BC:19:MET:N	2.24	0.53
1:BC:170:PRO:HG2	1:BC:171:GLY:N	2.22	0.53
3:BM:4:TYR:O	3:BM:4:TYR:HD1	1.92	0.53
3:BM:31:ILE:HD11	15:BQ:101:PEF:O2P	2.09	0.53
3:BM:164:ARG:HA	3:BM:167:MET:HE2	1.91	0.53
3:BM:199:ASN:HD21	3:BM:283:GLY:CA	2.22	0.53
5:BD:20:VAL:HA	5:BD:23:SER:OG	2.08	0.53
9:BQ:104:BCL:OBB	9:BQ:104:BCL:HHC	2.08	0.53
5:BS:24:ILE:CD1	9:BU:102:BCL:H202	2.38	0.53
6:B2:20:ILE:CG1	14:B2:102:CRT:C8	2.85	0.53
5:B3:9:TYR:HA	6:B4:18:HIS:ND1	2.24	0.53
9:B3:102:BCL:HBA2	9:B4:101:BCL:OBD	2.09	0.53
6:B6:17:PHE:C	6:B6:17:PHE:CD1	2.87	0.53
6:B6:45:TRP:O	6:B6:46:LEU:C	2.51	0.53
5:B9:46:TRP:CE2	9:B9:102:BCL:H2C	2.43	0.53
3:AM:35:ILE:HD11	15:AM:409:PEF:H321	1.90	0.53
3:AM:76:LEU:HA	3:AM:86:PHE:CD1	2.43	0.53
3:AM:98:PRO:HD2	3:AM:171:TRP:O	2.08	0.53
3:AM:131:VAL:C	3:AM:133:THR:N	2.65	0.53
4:AH:168:SER:HB3	4:AH:183:GLU:HB2	1.91	0.53
5:AF:12:TRP:HE1	6:AG:17:PHE:HD1	1.56	0.53
5:AO:44:LEU:HD12	5:AO:44:LEU:C	2.34	0.53
6:AP:31:LEU:C	6:AP:34:ILE:HG23	2.33	0.53
9:AQ:102:BCL:HBC2	9:AQ:102:BCL:CHD	2.37	0.53
9:AY:102:BCL:C9	6:AZ:28:TRP:HB2	2.39	0.53
6:AZ:29:PHE:N	6:AZ:29:PHE:CD1	2.77	0.53
9:A2:101:BCL:HMB1	9:A2:101:BCL:HBB3	1.90	0.53
5:A3:12:TRP:HA	5:A3:12:TRP:HE3	1.74	0.53
1:BC:66:ASP:C	1:BC:68:THR:H	2.17	0.53
1:BC:124:LYS:O	1:BC:125:VAL:C	2.51	0.53
1:BC:276:VAL:CG1	1:BC:277:ARG:N	2.71	0.53
2:BL:139:VAL:HA	2:BL:143:VAL:HB	1.91	0.53
2:BL:177:HIS:CB	3:BM:183:LEU:HD22	2.27	0.53
2:BL:207:THR:HA	2:BL:215:VAL:HG13	1.90	0.53
3:BM:13:VAL:HG12	4:BH:144:ILE:HD13	1.91	0.53
3:BM:17:ALA:O	3:BM:19:PRO:HD3	2.09	0.53
5:BF:11:ILE:CD1	5:BF:14:ILE:HD11	2.39	0.53
5:BF:14:ILE:HD13	6:BJ:17:PHE:CE2	2.44	0.53
6:BG:45:TRP:O	6:BG:46:LEU:HB2	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:BO:4:MET:CB	6:BR:23:GLN:CB	2.77	0.53
5:BQ:28:GLN:O	9:BQ:103:BCL:H11	2.08	0.53
5:BQ:36:HIS:NE2	9:BQ:104:BCL:HMD3	2.23	0.53
5:BU:5:ASN:HB3	6:BV:22:MET:HE3	1.89	0.53
6:BV:17:PHE:HA	14:BV:102:CRT:C4	2.38	0.53
5:BW:4:MET:C	5:BW:6:ALA:N	2.67	0.53
5:B1:53:VAL:O	5:B1:55:TYR:N	2.42	0.53
1:AC:29:GLY:O	1:AC:30:THR:HG23	2.08	0.53
1:AC:124:LYS:O	1:AC:125:VAL:C	2.51	0.53
1:AC:302:PRO:O	1:AC:302:PRO:HG2	2.08	0.53
2:AL:99:THR:HG23	2:AL:157:TYR:OH	2.08	0.53
3:AM:59:LEU:HG	3:AM:128:LEU:CD2	2.38	0.53
4:AH:234:TYR:CZ	4:AH:238:LYS:HE3	2.44	0.53
9:AE:101:BCL:HHC	9:AE:101:BCL:OBB	2.09	0.53
6:AG:24:SER:O	6:AG:27:ALA:HB3	2.09	0.53
5:AK:13:LEU:HD21	6:AN:10:THR:O	2.08	0.53
5:AK:16:ASP:HB2	5:AK:19:ARG:CG	2.35	0.53
5:AU:26:ALA:C	5:AU:29:ILE:HG22	2.32	0.53
9:AU:102:BCL:C1D	9:AV:102:BCL:HMD1	2.39	0.53
9:AW:101:BCL:CHD	9:AX:101:BCL:HMD1	2.37	0.53
5:A1:9:TYR:HA	6:A2:18:HIS:CE1	2.38	0.53
5:A3:32:GLY:HA3	9:A3:103:BCL:O1A	2.09	0.53
5:A7:21:LEU:O	5:A7:25:VAL:HG23	2.07	0.53
1:BC:128:ARG:O	1:BC:131:PHE:HB2	2.08	0.53
1:BC:245:VAL:HG23	1:BC:245:VAL:O	2.07	0.53
1:BC:272:ALA:O	1:BC:276:VAL:HG12	2.09	0.53
2:BL:97:ILE:O	2:BL:100:ILE:HB	2.09	0.53
2:BL:112:ARG:HH21	3:BM:255:THR:HA	1.72	0.53
2:BL:229:VAL:C	3:BM:51:ILE:HD12	2.34	0.53
2:BL:237:ALA:O	2:BL:240:ARG:N	2.42	0.53
3:BM:291:VAL:HG21	3:BM:297:TRP:CD1	2.44	0.53
4:BH:69:LEU:HB2	4:BH:74:GLY:O	2.09	0.53
5:BA:10:LYS:HB2	14:BA:102:CRT:H5	1.91	0.53
5:BD:40:LEU:HD13	5:BD:47:LEU:HD23	1.90	0.53
5:BD:45:ASN:HB3	5:BD:49:ASP:HB3	1.89	0.53
9:BD:102:BCL:HMD3	6:BE:36:HIS:CD2	2.44	0.53
6:BE:29:PHE:CD1	9:BE:101:BCL:C2	2.91	0.53
6:BE:45:TRP:O	6:BE:46:LEU:CG	2.48	0.53
6:BN:19:ALA:O	6:BN:23:GLN:HG2	2.09	0.53
6:BP:25:MET:HE2	14:BP:102:CRT:H242	1.89	0.53
5:BS:13:LEU:HD21	6:BT:10:THR:O	2.09	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:B3:44:LEU:HD13	5:B3:46:TRP:HE3	1.73	0.53
1:AC:270:TRP:HA	1:AC:273:ILE:HD12	1.90	0.53
2:AL:5:SER:HB3	4:AH:38:GLY:O	2.09	0.53
2:AL:29:PRO:HB2	3:AM:253:ARG:CD	2.37	0.53
3:AM:177:PHE:CD1	14:AM:406:CRT:H16	2.43	0.53
3:AM:290:VAL:HG12	3:AM:291:VAL:N	2.23	0.53
6:AB:17:PHE:CE1	14:AB:102:CRT:C9	2.91	0.53
6:AJ:34:ILE:HD13	6:AJ:34:ILE:C	2.34	0.53
5:AK:5:ASN:CG	6:AN:22:MET:HE3	2.32	0.53
5:AQ:35:ILE:HA	5:AQ:38:ILE:CG2	2.32	0.53
6:AX:45:TRP:O	6:AX:46:LEU:CG	2.56	0.53
5:AY:4:MET:C	5:AY:8:LEU:HB2	2.33	0.53
5:A3:35:ILE:O	5:A3:36:HIS:C	2.52	0.53
6:A4:18:HIS:CD2	6:A4:22:MET:HB2	2.43	0.53
6:A8:23:GLN:HG3	6:A8:24:SER:N	2.24	0.53
6:A8:29:PHE:HZ	9:A8:101:BCL:C6	2.21	0.53
1:BC:53:ILE:C	1:BC:55:ALA:H	2.17	0.53
1:BC:107:CYS:C	1:BC:109:TYR:H	2.17	0.53
1:BC:232:THR:O	1:BC:233:PHE:C	2.52	0.53
2:BL:212:GLY:O	2:BL:213:GLU:HG2	2.09	0.53
3:BM:71:ILE:HD13	3:BM:177:PHE:CD1	2.43	0.53
3:BM:77:ALA:O	3:BM:80:HIS:N	2.42	0.53
3:BM:79:VAL:O	3:BM:79:VAL:HG22	2.08	0.53
3:BM:102:TYR:H	3:BM:102:TYR:HD1	1.56	0.53
4:BH:108:LEU:C	4:BH:110:GLY:N	2.65	0.53
4:BH:133:ILE:CD1	4:BH:171:TRP:HB3	2.39	0.53
4:BH:235:GLU:HA	4:BH:238:LYS:HB2	1.91	0.53
9:BA:101:BCL:H151	5:B9:24:ILE:HD11	1.90	0.53
9:BN:101:BCL:C1B	9:BO:102:BCL:HMB3	2.39	0.53
5:BO:7:ASN:C	6:BR:20:ILE:HD11	2.34	0.53
9:BQ:103:BCL:HBA2	9:BQ:104:BCL:OBD	2.09	0.53
5:BS:8:LEU:HB2	6:BT:18:HIS:HE2	1.74	0.53
5:BW:7:ASN:HD22	5:BW:7:ASN:N	2.07	0.53
5:BW:14:ILE:HG21	5:BY:21:LEU:CD1	2.38	0.53
5:BY:33:LEU:HD12	5:BY:34:LEU:N	2.23	0.53
5:B1:11:ILE:HD13	9:B3:102:BCL:H151	1.90	0.53
6:B2:13:GLU:C	14:B2:102:CRT:C3	2.80	0.53
6:B2:30:GLY:O	6:B2:33:VAL:HG12	2.09	0.53
6:B4:10:THR:HB	6:B4:13:GLU:OE2	2.08	0.53
5:B7:10:LYS:HD3	6:B0:20:ILE:CD1	2.38	0.53
5:B7:36:HIS:HB3	14:B7:102:CRT:H2M3	1.87	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:AL:51:VAL:CG1	5:AA:37:MET:HG2	2.39	0.53
2:AL:278:LEU:HD12	2:AL:281:TRP:CZ2	2.44	0.53
3:AM:34:PRO:HG3	3:AM:50:PRO:CD	2.38	0.53
3:AM:136:ARG:NH1	3:AM:136:ARG:CA	2.68	0.53
4:AH:185:GLU:HA	4:AH:191:LYS:O	2.09	0.53
4:AH:197:ILE:O	4:AH:197:ILE:HD13	2.08	0.53
5:AA:27:PHE:HA	5:AA:30:VAL:CG1	2.38	0.53
6:AB:28:TRP:HA	6:AB:31:LEU:HG	1.90	0.53
6:AG:45:TRP:O	6:AG:46:LEU:CB	2.53	0.53
9:AI:102:BCL:C3D	6:AJ:35:ALA:HB1	2.38	0.53
14:AS:104:CRT:H16	6:AV:21:PHE:HE1	1.73	0.53
9:AV:102:BCL:H172	6:AX:39:ALA:HA	1.90	0.53
5:AY:40:LEU:HD13	5:AY:46:TRP:CZ2	2.43	0.53
9:A1:102:BCL:H92	6:A2:28:TRP:HB2	1.89	0.53
5:A5:46:TRP:CZ2	9:A5:102:BCL:CHC	2.92	0.53
9:A7:103:BCL:C1D	9:A8:101:BCL:CMD	2.87	0.53
6:A8:7:THR:HG23	6:A8:8:GLY:N	2.24	0.53
1:BC:120:ASP:CG	1:BC:120:ASP:O	2.52	0.53
3:BM:25:LYS:HG2	5:BO:16:ASP:OD1	2.09	0.53
3:BM:316:PRO:HG2	3:BM:317:TYR:HD1	1.73	0.53
4:BH:27:ILE:O	4:BH:27:ILE:HD13	2.07	0.53
5:BD:16:ASP:OD2	5:BD:18:ARG:HG2	2.07	0.53
5:BQ:35:ILE:CA	5:BQ:38:ILE:HG22	2.30	0.53
5:BQ:43:ASP:CB	5:BS:47:LEU:HB3	2.39	0.53
5:BU:55:TYR:O	5:BU:59:GLY:HA3	2.09	0.53
5:BW:35:ILE:HA	5:BW:38:ILE:HG22	1.90	0.53
5:B1:19:ARG:O	5:B1:23:SER:CB	2.57	0.53
5:B1:27:PHE:O	5:B1:30:VAL:HG12	2.08	0.53
5:B1:38:ILE:HG23	5:B1:39:VAL:H	1.73	0.53
5:B3:14:ILE:CG2	5:B5:17:PRO:HB2	2.39	0.53
5:B9:46:TRP:NE1	5:B9:47:LEU:HD22	2.23	0.53
1:AC:120:ASP:CG	1:AC:120:ASP:O	2.52	0.53
1:AC:130:MET:SD	7:AC:502:HEM:NA	2.82	0.53
1:AC:157:ARG:HH12	1:AC:318:LEU:HG	1.73	0.53
1:AC:194:SER:C	1:AC:195:LEU:HG	2.33	0.53
1:AC:272:ALA:C	1:AC:274:ARG:H	2.16	0.53
2:AL:228:ILE:HG23	3:AM:132:ARG:HD2	1.90	0.53
2:AL:242:GLY:HA2	3:AM:216:PHE:HE2	1.74	0.53
2:AL:242:GLY:O	2:AL:243:LEU:C	2.51	0.53
3:AM:214:LEU:O	3:AM:218:MET:HG3	2.09	0.53
4:AH:66:THR:HG23	4:AH:66:THR:O	2.07	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:AH:196:PRO:HG2	4:AH:199:PHE:HB2	1.91	0.53
6:AB:20:ILE:HG13	5:A9:7:ASN:HB2	1.88	0.53
6:AE:20:ILE:O	6:AE:23:GLN:HG3	2.08	0.53
9:AG:101:BCL:C1B	9:AI:102:BCL:HMB3	2.39	0.53
5:AI:39:VAL:O	5:AI:43:ASP:HB3	2.08	0.53
9:AO:102:BCL:C2D	9:AP:101:BCL:C2D	2.86	0.53
5:AQ:14:ILE:O	5:AS:18:ARG:NH2	2.41	0.53
6:AT:42:TYR:CE2	6:AT:43:ARG:HG2	2.44	0.53
5:AY:9:TYR:HA	6:AZ:18:HIS:CG	2.44	0.53
5:AY:36:HIS:CE1	9:AY:102:BCL:NA	2.77	0.53
5:A9:2:PHE:HE1	6:A0:26:TYR:HH	1.56	0.53
2:BL:182:HIS:CE1	9:BL:301:BCL:NB	2.76	0.53
3:BM:157:TYR:CD1	3:BM:158:LEU:HD23	2.44	0.53
4:BH:5:ILE:HD11	5:BF:47:LEU:HD12	1.91	0.53
4:BH:36:ARG:HD2	4:BH:78:ALA:CB	2.39	0.53
4:BH:54:LYS:O	4:BH:55:VAL:C	2.52	0.53
6:BB:17:PHE:HD1	14:BB:102:CRT:C6	2.21	0.53
5:BD:12:TRP:HA	5:BD:12:TRP:HE3	1.74	0.53
5:BD:39:VAL:HG12	5:BD:46:TRP:HZ3	1.73	0.53
6:BT:21:PHE:CD1	6:BT:21:PHE:C	2.87	0.53
6:BV:7:THR:OG1	6:BV:8:GLY:N	2.40	0.53
5:B1:44:LEU:O	5:B1:44:LEU:HG	2.08	0.53
6:B2:26:TYR:HA	6:B2:29:PHE:HD2	1.74	0.53
5:B9:35:ILE:O	5:B9:39:VAL:HG23	2.09	0.53
5:B9:50:ASN:ND2	5:B9:51:ILE:HG12	2.21	0.53
6:B0:17:PHE:HD1	14:B0:101:CRT:C9	2.19	0.53
1:AC:284:ILE:O	1:AC:284:ILE:HG22	2.08	0.53
3:AM:85:GLN:HG3	3:AM:89:HIS:HD2	1.73	0.53
3:AM:132:ARG:O	3:AM:132:ARG:CD	2.57	0.53
4:AH:219:PHE:HA	4:AH:222:VAL:CG2	2.39	0.53
5:AF:7:ASN:HB3	6:AJ:20:ILE:HD13	1.90	0.53
5:AI:55:TYR:CD1	5:AI:56:GLN:N	2.75	0.53
6:AP:13:GLU:HA	6:AP:16:GLU:CG	2.38	0.53
5:AQ:42:THR:HG23	5:AQ:43:ASP:H	1.74	0.53
6:AV:33:VAL:HG13	6:AV:34:ILE:N	2.23	0.53
5:AY:44:LEU:HD12	5:AY:44:LEU:O	2.09	0.53
6:AZ:30:GLY:O	6:AZ:34:ILE:HG22	2.08	0.53
9:A3:103:BCL:OBD	6:A4:32:VAL:HG13	2.09	0.53
6:A6:10:THR:HG22	6:A6:11:ASP:H	1.74	0.53
5:A7:40:LEU:HD13	5:A7:46:TRP:CE2	2.44	0.53
1:BC:276:VAL:HG22	1:BC:280:ASN:ND2	2.23	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:BL:111:LEU:HA	2:BL:114:VAL:HG23	1.91	0.53
2:BL:246:ALA:C	2:BL:248:SER:H	2.16	0.53
6:BB:20:ILE:HD12	14:BB:102:CRT:H81	1.89	0.53
5:BD:31:LEU:HD12	5:BD:34:LEU:HD23	1.91	0.53
6:BN:22:MET:HG3	6:BN:26:TYR:HE2	1.74	0.53
9:BO:102:BCL:O1D	9:BO:102:BCL:C2A	2.57	0.53
5:BS:55:TYR:CD1	5:BS:56:GLN:N	2.76	0.53
5:BW:33:LEU:C	5:BW:33:LEU:HD12	2.34	0.53
9:BW:102:BCL:C4D	9:BX:101:BCL:HMD3	2.38	0.53
5:B9:5:ASN:HA	5:B9:8:LEU:CG	2.39	0.53
1:AC:104:LYS:HB3	1:AC:105:GLU:OE2	2.09	0.52
1:AC:183:GLN:O	1:AC:195:LEU:O	2.27	0.52
3:AM:4:TYR:CD1	3:AM:4:TYR:C	2.87	0.52
3:AM:102:TYR:O	3:AM:104:LEU:N	2.42	0.52
3:AM:146:LEU:O	3:AM:147:SER:C	2.51	0.52
3:AM:204:LEU:O	3:AM:206:ILE:N	2.42	0.52
3:AM:317:TYR:HD1	3:AM:317:TYR:H	1.55	0.52
9:AK:102:BCL:H2	6:AN:28:TRP:CH2	2.44	0.52
9:AR:101:BCL:OBB	9:AR:101:BCL:HHC	2.09	0.52
5:AU:2:PHE:CD1	5:AU:2:PHE:O	2.62	0.52
5:AU:25:VAL:HG21	9:AU:102:BCL:H142	1.90	0.52
14:AW:102:CRT:H1M3	6:AZ:16:GLU:HB3	1.91	0.52
5:A3:56:GLN:N	5:A3:56:GLN:HE21	2.05	0.52
6:A0:20:ILE:HG23	6:A0:21:PHE:N	2.25	0.52
1:BC:263:THR:O	1:BC:264:PRO:C	2.52	0.52
2:BL:93:GLY:HA2	2:BL:96:GLN:NE2	2.24	0.52
2:BL:202:LEU:HD21	2:BL:221:GLU:CB	2.35	0.52
3:BM:76:LEU:HD23	5:BU:37:MET:CE	2.35	0.52
3:BM:85:GLN:HG3	3:BM:89:HIS:HD2	1.72	0.52
3:BM:222:THR:HG21	3:BM:252:TRP:NE1	2.22	0.52
5:BK:38:ILE:HD13	14:BN:102:CRT:H401	1.91	0.52
5:BO:11:ILE:N	14:BO:103:CRT:H82	2.24	0.52
5:BO:50:ASN:CG	5:BO:51:ILE:N	2.67	0.52
5:BQ:46:TRP:NE1	5:BQ:47:LEU:HG	2.24	0.52
5:BW:26:ALA:CA	5:BW:29:ILE:HG22	2.39	0.52
5:BY:20:VAL:O	5:BY:24:ILE:HG12	2.09	0.52
5:B1:18:ARG:HD2	5:B1:19:ARG:HG3	1.91	0.52
5:B5:29:ILE:HB	9:B5:102:BCL:H43	1.91	0.52
5:B7:19:ARG:O	5:B7:23:SER:HB2	2.09	0.52
1:AC:81:VAL:HG11	1:AC:131:PHE:CB	2.34	0.52
2:AL:38:VAL:HA	2:AL:41:CYS:SG	2.49	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:AL:78:PRO:CB	2:AL:92:GLY:HA3	2.31	0.52
2:AL:156:PRO:O	2:AL:157:TYR:CD1	2.62	0.52
3:AM:56:THR:HA	3:AM:59:LEU:HB3	1.91	0.52
3:AM:63:PHE:HE1	5:AQ:30:VAL:HA	1.75	0.52
3:AM:64:GLY:HA3	10:AM:403:BPH:H5C1	1.91	0.52
3:AM:179:ILE:H	3:AM:179:ILE:CD1	2.22	0.52
4:AH:132:LYS:NZ	4:AH:173:ASP:OD2	2.32	0.52
5:AA:18:ARG:HD2	5:AA:18:ARG:N	2.18	0.52
6:AJ:17:PHE:CA	6:AJ:20:ILE:HG22	2.39	0.52
9:AJ:101:BCL:HMB3	9:AK:102:BCL:CHB	2.39	0.52
5:AK:31:LEU:O	5:AK:35:ILE:HG12	2.09	0.52
5:AQ:44:LEU:HD12	5:AQ:46:TRP:CE3	2.37	0.52
6:AR:20:ILE:HD13	6:AR:20:ILE:O	2.09	0.52
9:AR:101:BCL:HMB3	9:AS:103:BCL:C4A	2.39	0.52
5:AW:54:SER:CB	5:AW:57:ALA:HB3	2.39	0.52
14:AW:102:CRT:C7	6:AZ:20:ILE:HD13	2.40	0.52
6:AX:40:TRP:O	6:AX:44:PRO:HG3	2.09	0.52
5:AY:30:VAL:O	5:AY:33:LEU:HG	2.08	0.52
1:BC:70:PRO:C	1:BC:71:LYS:HD2	2.34	0.52
2:BL:276:LEU:HD22	2:BL:276:LEU:N	2.21	0.52
3:BM:190:SER:O	3:BM:194:GLY:O	2.27	0.52
4:BH:5:ILE:HG21	5:BD:42:THR:HG21	1.88	0.52
4:BH:37:GLU:O	4:BH:39:TYR:N	2.42	0.52
4:BH:234:TYR:CE1	4:BH:238:LYS:HE3	2.45	0.52
6:BP:10:THR:CB	6:BP:13:GLU:OE1	2.58	0.52
5:BQ:51:ILE:HA	5:BQ:52:PRO:C	2.33	0.52
5:BU:15:LEU:HD11	9:BW:102:BCL:H141	1.91	0.52
6:B2:21:PHE:CA	14:B2:102:CRT:C11	2.85	0.52
5:B3:14:ILE:HG21	5:B5:17:PRO:HB2	1.91	0.52
1:AC:270:TRP:HE3	1:AC:271:TYR:CD1	2.27	0.52
2:AL:7:GLU:CD	2:AL:11:ARG:HH21	2.17	0.52
2:AL:120:LEU:O	2:AL:121:GLY:C	2.51	0.52
2:AL:151:TRP:C	2:AL:153:HIS:H	2.18	0.52
2:AL:177:HIS:CG	3:AM:183:LEU:HD22	2.44	0.52
2:AL:181:ALA:O	2:AL:183:MET:N	2.42	0.52
2:AL:230:GLY:N	3:AM:51:ILE:HD12	2.25	0.52
3:AM:156:PHE:HA	3:AM:159:VAL:CG2	2.39	0.52
3:AM:161:GLY:O	3:AM:162:PHE:C	2.52	0.52
3:AM:221:ALA:HA	3:AM:224:LEU:HD12	1.91	0.52
3:AM:290:VAL:HG12	3:AM:291:VAL:H	1.75	0.52
4:AH:45:ARG:HH22	4:AH:100:LEU:HD21	1.74	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:AA:9:TYR:HB2	6:AB:18:HIS:CD2	2.45	0.52
5:AO:46:TRP:CE3	9:AO:102:BCL:H2C	2.44	0.52
6:AP:21:PHE:CD1	6:AP:21:PHE:O	2.62	0.52
5:AW:12:TRP:HZ2	6:AX:21:PHE:CB	2.21	0.52
5:AW:18:ARG:HG2	5:AW:18:ARG:NH1	2.24	0.52
5:AW:36:HIS:O	5:AW:40:LEU:HB3	2.09	0.52
6:AX:18:HIS:NE2	6:AX:22:MET:HE2	2.25	0.52
6:A2:40:TRP:CZ3	6:A2:44:PRO:HA	2.44	0.52
5:A5:8:LEU:CD2	14:A5:103:CRT:H133	2.40	0.52
14:A5:103:CRT:H402	5:A7:38:ILE:HG21	1.92	0.52
6:A0:9:LEU:HB3	6:A0:13:GLU:CG	2.38	0.52
1:BC:148:THR:CB	1:BC:322:GLN:HG2	2.39	0.52
2:BL:125:HIS:CE1	3:BM:5:GLN:HG3	2.44	0.52
3:BM:102:TYR:CD1	3:BM:102:TYR:N	2.77	0.52
3:BM:249:ALA:HB2	13:BM:405:MQ8:H61	1.92	0.52
3:BM:261:THR:C	3:BM:263:GLU:N	2.65	0.52
6:BB:17:PHE:CE1	14:BB:102:CRT:H9	2.37	0.52
5:BK:45:ASN:HB3	5:BK:49:ASP:HB3	1.92	0.52
5:BK:47:LEU:HD22	5:BK:47:LEU:N	2.24	0.52
5:BO:44:LEU:HD12	5:BO:45:ASN:N	2.24	0.52
6:BP:23:GLN:O	6:BP:24:SER:C	2.51	0.52
9:B4:101:BCL:HMC1	9:B5:102:BCL:HBB1	1.91	0.52
9:B8:101:BCL:HHB	9:B9:102:BCL:HMA1	1.91	0.52
1:AC:191:ALA:HB3	1:AC:237:MET:HE3	1.90	0.52
2:AL:238:ILE:HG22	2:AL:239:HIS:N	2.25	0.52
3:AM:134:TYR:O	3:AM:144:GLN:NE2	2.40	0.52
3:AM:151:ALA:O	3:AM:152:ALA:C	2.52	0.52
4:AH:106:PRO:HA	4:AH:109:SER:OG	2.09	0.52
6:AG:43:ARG:HD3	5:AI:55:TYR:OH	2.09	0.52
5:AK:11:ILE:HG12	14:AP:102:CRT:H81	1.91	0.52
5:AQ:52:PRO:HG2	5:AQ:53:VAL:H	1.74	0.52
5:AS:10:LYS:CB	14:AS:104:CRT:H1M2	2.39	0.52
5:A1:12:TRP:CD1	6:A2:18:HIS:HB2	2.44	0.52
14:A1:103:CRT:H342	9:A5:102:BCL:HBA1	1.91	0.52
14:A5:103:CRT:H22A	6:A8:17:PHE:CE1	2.44	0.52
6:A6:28:TRP:C	6:A6:30:GLY:H	2.17	0.52
2:BL:26:TRP:HE3	4:BH:97:GLY:O	1.92	0.52
2:BL:192:ASN:C	2:BL:194:LEU:H	2.17	0.52
3:BM:177:PHE:CD1	14:BM:406:CRT:H19	2.44	0.52
3:BM:276:THR:C	3:BM:278:ILE:N	2.68	0.52
4:BH:114:ALA:HB2	4:BH:245:GLY:CA	2.37	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:BE:45:TRP:C	5:BF:52:PRO:HD3	2.34	0.52
5:BF:36:HIS:CE1	9:BG:101:BCL:HMD3	2.45	0.52
9:BG:101:BCL:CMB	9:BI:102:BCL:CHB	2.86	0.52
6:BN:13:GLU:CD	6:BN:13:GLU:H	2.17	0.52
5:BO:36:HIS:CE1	9:BP:101:BCL:HMD3	2.43	0.52
6:BX:21:PHE:HD1	6:BX:22:MET:N	2.07	0.52
9:BZ:101:BCL:HBB3	9:B1:102:BCL:C1C	2.39	0.52
5:B1:11:ILE:CG2	5:B1:15:LEU:HD12	2.39	0.52
5:B5:21:LEU:O	5:B5:25:VAL:HG23	2.10	0.52
6:B6:10:THR:HG22	6:B6:11:ASP:H	1.74	0.52
6:B6:28:TRP:C	6:B6:30:GLY:H	2.18	0.52
5:B7:36:HIS:CE1	9:B7:103:BCL:NA	2.76	0.52
5:B7:42:THR:C	5:B9:48:ASP:HB3	2.34	0.52
6:B8:34:ILE:O	6:B8:37:LEU:HB3	2.09	0.52
2:AL:69:ASN:O	2:AL:70:LEU:C	2.52	0.52
2:AL:88:PRO:O	2:AL:89:LEU:C	2.53	0.52
2:AL:124:PHE:O	2:AL:127:PRO:HD2	2.10	0.52
3:AM:124:LEU:O	3:AM:127:LEU:N	2.43	0.52
3:AM:214:LEU:C	3:AM:214:LEU:HD23	2.34	0.52
3:AM:214:LEU:HD22	3:AM:215:LEU:HD12	1.91	0.52
4:AH:77:VAL:HG23	4:AH:80:ARG:HB3	1.90	0.52
6:AB:34:ILE:HD13	6:AB:34:ILE:C	2.35	0.52
6:AG:46:LEU:CB	6:AJ:42:TYR:OH	2.56	0.52
6:AP:36:HIS:HE1	9:AP:101:BCL:C1A	2.23	0.52
5:AQ:48:ASP:O	5:AQ:49:ASP:CB	2.55	0.52
5:AS:42:THR:CG2	5:AU:47:LEU:HB3	2.35	0.52
5:AU:36:HIS:O	5:AU:40:LEU:CB	2.57	0.52
9:AU:102:BCL:C4D	9:AV:102:BCL:HMD3	2.39	0.52
5:AW:14:ILE:HG21	5:AY:21:LEU:CD1	2.39	0.52
5:A3:8:LEU:HD21	6:A6:24:SER:OG	2.10	0.52
5:A5:50:ASN:CG	5:A5:51:ILE:H	2.17	0.52
9:A8:101:BCL:OBB	9:A8:101:BCL:HHC	2.10	0.52
6:A0:40:TRP:HA	6:A0:40:TRP:CE3	2.45	0.52
9:A0:102:BCL:H141	9:A0:102:BCL:CMB	2.40	0.52
1:BC:153:TYR:CD2	1:BC:323:MET:HE1	2.43	0.52
1:BC:245:VAL:HG21	1:BC:249:PHE:CG	2.45	0.52
2:BL:240:ARG:CZ	3:BM:7:ILE:O	2.57	0.52
9:BL:301:BCL:CHC	9:BM:402:BCL:CHC	2.88	0.52
3:BM:120:LEU:HB2	14:BM:406:CRT:H35	1.91	0.52
3:BM:241:ARG:HG2	3:BM:242:GLY:N	2.25	0.52
9:BM:401:BCL:HMB2	9:BM:402:BCL:H152	1.90	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:BA:36:HIS:CG	9:BB:101:BCL:HMD2	2.44	0.52
5:BD:12:TRP:CE3	5:BD:12:TRP:HA	2.45	0.52
5:BF:33:LEU:O	5:BF:37:MET:HG2	2.10	0.52
5:BF:43:ASP:O	5:BF:44:LEU:HG	2.10	0.52
5:BF:43:ASP:OD1	5:BF:44:LEU:HD23	2.09	0.52
9:BF:102:BCL:ND	9:BG:101:BCL:CMD	2.73	0.52
5:BS:9:TYR:HB2	6:BT:15:LYS:HA	1.91	0.52
6:BT:38:LEU:HD23	6:BT:38:LEU:C	2.35	0.52
9:BU:102:BCL:C1D	9:BV:101:BCL:CMD	2.87	0.52
9:BW:102:BCL:CAD	9:BX:101:BCL:CAD	2.87	0.52
5:BY:18:ARG:HG2	5:BY:18:ARG:HH11	1.75	0.52
5:BY:53:VAL:C	5:BY:55:TYR:H	2.17	0.52
1:AC:94:MET:SD	7:AC:501:HEM:NB	2.82	0.52
1:AC:164:TYR:O	1:AC:309:THR:HG23	2.10	0.52
1:AC:270:TRP:O	1:AC:274:ARG:CD	2.57	0.52
3:AM:61:ILE:HD12	15:AM:409:PEF:H191	1.92	0.52
4:AH:106:PRO:HA	4:AH:109:SER:CB	2.39	0.52
4:AH:136:MET:HG2	4:AH:172:VAL:HG13	1.90	0.52
6:AB:32:VAL:HG21	9:AB:101:BCL:CBA	2.27	0.52
5:AF:14:ILE:HD13	6:AJ:17:PHE:CE2	2.45	0.52
6:AG:28:TRP:C	6:AG:30:GLY:N	2.66	0.52
5:AO:11:ILE:CD1	14:AR:102:CRT:H132	2.39	0.52
6:AR:34:ILE:HD13	6:AR:34:ILE:O	2.10	0.52
5:AS:27:PHE:O	5:AS:31:LEU:HB3	2.10	0.52
5:AS:53:VAL:O	5:AS:55:TYR:N	2.41	0.52
9:AU:102:BCL:HMB3	9:AU:102:BCL:CBB	2.40	0.52
5:A5:45:ASN:HD21	5:A5:48:ASP:CG	2.17	0.52
5:A7:4:MET:HA	6:A0:23:GLN:OE1	2.10	0.52
6:A8:28:TRP:HA	6:A8:31:LEU:HB2	1.92	0.52
6:A8:38:LEU:O	6:A8:38:LEU:HD23	2.10	0.52
5:A9:44:LEU:O	5:A9:46:TRP:N	2.39	0.52
1:BC:110:CYS:O	1:BC:111:HIS:ND1	2.43	0.52
1:BC:157:ARG:NH1	1:BC:318:LEU:CD2	2.73	0.52
2:BL:12:VAL:CG2	2:BL:13:ARG:H	2.22	0.52
2:BL:221:GLU:O	2:BL:223:THR:N	2.42	0.52
2:BL:237:ALA:HA	2:BL:240:ARG:CD	2.40	0.52
2:BL:267:GLY:O	2:BL:270:GLU:HB2	2.09	0.52
3:BM:121:PHE:CD1	3:BM:121:PHE:N	2.77	0.52
3:BM:165:PRO:CB	3:BM:174:ALA:HB2	2.39	0.52
3:BM:228:ARG:HB2	3:BM:229:PHE:CD1	2.45	0.52
4:BH:5:ILE:CD1	5:BF:47:LEU:HD12	2.40	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:BA:46:TRP:O	6:B0:46:LEU:OXT	2.26	0.52
5:BD:45:ASN:O	5:BD:49:ASP:CG	2.53	0.52
6:BG:21:PHE:CZ	9:BI:102:BCL:H202	2.45	0.52
6:BP:21:PHE:CD1	6:BP:21:PHE:C	2.88	0.52
5:BW:16:ASP:HB2	5:BW:19:ARG:CD	2.39	0.52
5:BY:25:VAL:HG11	9:BY:102:BCL:H202	1.91	0.52
6:BZ:42:TYR:CE1	6:BZ:43:ARG:HG3	2.45	0.52
5:B3:11:ILE:CG1	14:B7:102:CRT:C8	2.81	0.52
5:B5:30:VAL:CG1	5:B5:31:LEU:H	2.16	0.52
5:B5:44:LEU:HD23	5:B5:44:LEU:H	1.75	0.52
2:AL:10:TYR:CE2	3:AM:246:GLU:HG2	2.45	0.52
2:AL:171:TYR:O	2:AL:173:PHE:N	2.41	0.52
2:AL:243:LEU:HD13	3:AM:221:ALA:HB2	1.92	0.52
3:AM:268:TRP:CD1	4:AH:30:LEU:HD22	2.45	0.52
3:AM:301:HIS:NE2	4:AH:10:ASP:OD2	2.43	0.52
5:AF:27:PHE:CA	5:AF:30:VAL:HG12	2.37	0.52
9:AF:102:BCL:C1D	9:AG:101:BCL:CMD	2.88	0.52
9:AJ:101:BCL:NB	9:AK:102:BCL:HMB1	2.24	0.52
9:AN:101:BCL:C2B	9:AO:102:BCL:C2B	2.88	0.52
9:AN:101:BCL:C1B	9:AO:102:BCL:CMB	2.88	0.52
5:AO:12:TRP:CD1	6:AP:18:HIS:HB2	2.45	0.52
5:AO:22:VAL:O	5:AO:25:VAL:HB	2.10	0.52
5:AS:43:ASP:C	5:AS:45:ASN:H	2.17	0.52
6:AX:29:PHE:HZ	14:AX:102:CRT:H242	1.73	0.52
1:BC:24:GLU:OE2	2:BL:266:ARG:NH2	2.43	0.52
1:BC:97:VAL:HG13	7:BC:502:HEM:HMB2	1.91	0.52
1:BC:167:VAL:HG23	1:BC:301:ASP:CG	2.34	0.52
2:BL:52:TRP:CE3	2:BL:52:TRP:HA	2.45	0.52
2:BL:188:PHE:HD2	2:BL:249:ALA:N	2.08	0.52
2:BL:190:PHE:O	2:BL:191:THR:C	2.53	0.52
2:BL:246:ALA:HB1	10:BL:302:BPH:HBC3	1.91	0.52
3:BM:34:PRO:HA	3:BM:48:ILE:O	2.09	0.52
3:BM:59:LEU:HD11	5:BQ:29:ILE:CG2	2.36	0.52
3:BM:161:GLY:O	3:BM:163:ILE:N	2.43	0.52
3:BM:254:TRP:CD1	3:BM:254:TRP:N	2.76	0.52
4:BH:35:LYS:HZ1	4:BH:57:GLY:HA3	1.74	0.52
4:BH:123:CYS:SG	4:BH:231:VAL:O	2.67	0.52
5:BA:22:VAL:C	5:BA:24:ILE:H	2.17	0.52
6:BG:24:SER:O	6:BG:27:ALA:HB3	2.10	0.52
5:BO:34:LEU:O	5:BO:38:ILE:HG23	2.08	0.52
5:BU:44:LEU:HD22	6:BV:43:ARG:CD	2.40	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:BY:45:ASN:O	5:BY:48:ASP:O	2.26	0.52
9:B2:101:BCL:CHC	9:B3:102:BCL:HBB3	2.40	0.52
5:B5:36:HIS:CE1	9:B5:102:BCL:NA	2.77	0.52
6:B8:7:THR:HG23	6:B8:8:GLY:N	2.24	0.52
2:AL:166:VAL:HG13	9:AL:301:BCL:HMD1	1.92	0.52
2:AL:279:PRO:HG3	5:AY:41:SER:HB2	1.91	0.52
3:AM:250:LEU:O	3:AM:254:TRP:CD1	2.62	0.52
4:AH:102:PRO:CG	4:AH:106:PRO:HB3	2.40	0.52
5:AI:17:PRO:O	5:AI:21:LEU:HB3	2.10	0.52
5:AK:36:HIS:CE1	9:AN:101:BCL:CMD	2.93	0.52
6:AP:28:TRP:HA	6:AP:31:LEU:HD12	1.91	0.52
5:AQ:32:GLY:HA2	9:AR:101:BCL:O1D	2.10	0.52
5:AS:9:TYR:HA	6:AT:18:HIS:CB	2.38	0.52
5:AW:42:THR:HB	5:AY:48:ASP:CB	2.40	0.52
5:A3:13:LEU:HD21	6:A4:10:THR:C	2.34	0.52
5:A9:44:LEU:HD22	5:A9:46:TRP:HB3	1.91	0.52
5:A9:44:LEU:CD1	5:A9:44:LEU:H	2.22	0.52
3:BM:90:PHE:C	3:BM:92:TRP:H	2.17	0.52
3:BM:138:GLU:C	3:BM:140:LEU:N	2.67	0.52
3:BM:214:LEU:HD22	3:BM:215:LEU:HD12	1.92	0.52
4:BH:123:CYS:SG	4:BH:230:GLN:HB2	2.50	0.52
5:BA:10:LYS:O	5:BA:13:LEU:HD13	2.10	0.52
5:BA:55:TYR:HE1	5:B9:44:LEU:CB	2.12	0.52
6:BE:17:PHE:CE1	6:BE:21:PHE:HB2	2.44	0.52
6:BG:43:ARG:HD3	5:BI:55:TYR:CZ	2.44	0.52
9:BP:101:BCL:HHC	9:BP:101:BCL:OBB	2.10	0.52
5:BW:35:ILE:O	5:BW:36:HIS:C	2.52	0.52
6:B2:38:LEU:O	6:B2:38:LEU:HD23	2.10	0.52
6:B8:23:GLN:HG3	6:B8:24:SER:N	2.24	0.52
2:AL:202:LEU:O	2:AL:205:SER:HB2	2.10	0.52
3:AM:79:VAL:O	3:AM:79:VAL:HG13	2.08	0.52
3:AM:101:GLN:C	3:AM:103:GLY:H	2.17	0.52
5:AA:27:PHE:HE1	5:AD:29:ILE:HD12	1.72	0.52
6:AE:9:LEU:HD13	6:AE:13:GLU:HG2	1.91	0.52
5:AF:11:ILE:CD1	5:AF:14:ILE:HD11	2.37	0.52
5:AF:36:HIS:CE1	9:AG:101:BCL:CMD	2.93	0.52
6:AJ:17:PHE:O	6:AJ:18:HIS:C	2.53	0.52
9:AJ:101:BCL:C2B	9:AK:102:BCL:C2B	2.88	0.52
6:AT:45:TRP:CD1	6:AT:46:LEU:N	2.78	0.52
5:AU:13:LEU:HD13	6:AV:9:LEU:O	2.10	0.52
6:AX:17:PHE:CD1	14:AX:102:CRT:H6	2.44	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:AX:101:BCL:OBB	9:AX:101:BCL:HHC	2.09	0.52
5:AY:43:ASP:HA	5:A1:48:ASP:CA	2.40	0.52
5:A1:10:LYS:HD2	6:A4:20:ILE:HB	1.91	0.52
5:A7:44:LEU:O	5:A7:44:LEU:HD13	2.09	0.52
5:A9:36:HIS:NE2	9:A0:102:BCL:CMD	2.70	0.52
6:A0:24:SER:HB2	14:A0:101:CRT:H183	1.91	0.52
1:BC:176:SER:HA	5:BU:48:ASP:OD2	2.09	0.52
1:BC:265:LYS:O	1:BC:268:THR:HB	2.10	0.52
1:BC:325:LYS:HD3	1:BC:325:LYS:C	2.34	0.52
2:BL:48:LEU:O	2:BL:51:VAL:HB	2.10	0.52
2:BL:148:MET:CE	2:BL:262:PRO:HD3	2.38	0.52
14:BA:102:CRT:H372	5:BD:35:ILE:HD11	1.92	0.52
14:BF:103:CRT:H6	6:BJ:17:PHE:CZ	2.44	0.52
6:BN:45:TRP:CD1	6:BN:46:LEU:HG	2.45	0.52
9:BN:101:BCL:HBB3	9:BO:102:BCL:C4B	2.39	0.52
5:BS:26:ALA:O	5:BS:29:ILE:HG22	2.10	0.52
14:BS:103:CRT:H6	6:BT:17:PHE:HD2	1.75	0.52
5:B1:32:GLY:N	9:B2:101:BCL:HED2	2.25	0.52
6:B2:13:GLU:HA	14:B2:102:CRT:H1M3	1.91	0.52
6:B2:17:PHE:CG	14:B2:102:CRT:H6	2.45	0.52
6:B2:21:PHE:HD1	14:B2:102:CRT:C15	2.21	0.52
5:B5:18:ARG:HB2	5:B5:19:ARG:HH22	1.74	0.52
6:B0:40:TRP:CE3	6:B0:40:TRP:HA	2.44	0.52
2:AL:47:VAL:HA	9:AL:303:BCL:H191	1.90	0.52
2:AL:273:ASN:O	2:AL:275:TRP:N	2.43	0.52
3:AM:62:PHE:C	3:AM:64:GLY:H	2.18	0.52
3:AM:130:TRP:HA	3:AM:150:PHE:CD2	2.45	0.52
3:AM:177:PHE:HD1	14:AM:406:CRT:H16	1.74	0.52
3:AM:234:GLU:O	3:AM:238:ILE:HG12	2.10	0.52
4:AH:189:ASN:HB3	4:AH:191:LYS:CG	2.40	0.52
5:AA:44:LEU:HD12	5:AA:44:LEU:C	2.35	0.52
9:AA:101:BCL:C4A	9:A0:102:BCL:HMB2	2.40	0.52
5:AD:16:ASP:OD2	5:AD:18:ARG:CG	2.55	0.52
5:AF:44:LEU:O	5:AF:46:TRP:N	2.37	0.52
5:AF:50:ASN:HB3	5:AI:56:GLN:HA	1.91	0.52
5:AI:17:PRO:O	5:AI:21:LEU:CB	2.58	0.52
5:AI:31:LEU:HD12	5:AI:34:LEU:HD23	1.91	0.52
5:AS:33:LEU:O	5:AS:37:MET:CB	2.58	0.52
5:AS:34:LEU:HD13	15:AS:101:PEF:C44	2.36	0.52
5:AW:19:ARG:NH1	5:AY:22:VAL:HG21	2.24	0.52
5:AW:36:HIS:CE1	9:AW:101:BCL:NA	2.77	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
14:AX:102:CRT:H2M1	5:AY:37:MET:N	2.25	0.52
6:A2:28:TRP:O	6:A2:32:VAL:HG23	2.10	0.52
5:A7:35:ILE:O	5:A7:38:ILE:HG22	2.10	0.52
5:A9:36:HIS:HE1	9:A0:102:BCL:OBD	1.92	0.52
6:A0:29:PHE:N	6:A0:29:PHE:HD1	2.08	0.52
1:BC:243:LEU:HD12	1:BC:243:LEU:N	2.23	0.52
1:BC:326:ASP:O	1:BC:327:TYR:CD1	2.62	0.52
2:BL:196:LEU:C	2:BL:198:MET:H	2.16	0.52
3:BM:215:LEU:HD21	13:BM:405:MQ8:H193	1.91	0.52
3:BM:267:ARG:NH1	4:BH:33:GLU:OE1	2.43	0.52
4:BH:119:ARG:NE	4:BH:237:ASP:OD2	2.43	0.52
6:BE:18:HIS:HE1	6:BE:22:MET:HE2	1.72	0.52
5:BF:50:ASN:CG	6:BG:43:ARG:NH2	2.55	0.52
9:BF:102:BCL:OBB	9:BF:102:BCL:HHC	2.08	0.52
9:BG:101:BCL:CHC	9:BI:102:BCL:HBB3	2.40	0.52
6:BJ:33:VAL:O	6:BJ:37:LEU:HD23	2.10	0.52
6:BP:7:THR:HG23	6:BP:8:GLY:N	2.24	0.52
5:B5:49:ASP:CG	5:B5:50:ASN:N	2.68	0.52
2:AL:107:ILE:O	2:AL:111:LEU:HG	2.09	0.51
2:AL:233:ILE:HG12	2:AL:237:ALA:CB	2.40	0.51
3:AM:59:LEU:HD11	5:AQ:29:ILE:HG21	1.91	0.51
3:AM:84:PHE:CE2	5:AW:37:MET:HG2	2.44	0.51
5:AA:11:ILE:HD11	5:AD:21:LEU:CD2	2.40	0.51
5:AD:43:ASP:CG	5:AD:44:LEU:N	2.68	0.51
6:AV:34:ILE:O	6:AV:37:LEU:HB2	2.09	0.51
5:AW:33:LEU:O	5:AW:37:MET:HB2	2.11	0.51
9:AZ:101:BCL:HMB3	9:A1:102:BCL:CHB	2.39	0.51
5:A7:35:ILE:HA	5:A7:38:ILE:HG22	1.91	0.51
1:BC:253:THR:HA	1:BC:256:PHE:CE1	2.45	0.51
2:BL:89:LEU:CD1	2:BL:89:LEU:H	2.23	0.51
2:BL:160:LEU:C	2:BL:160:LEU:HD12	2.36	0.51
2:BL:199:HIS:CE1	2:BL:239:HIS:CE1	2.98	0.51
3:BM:25:LYS:O	5:BO:18:ARG:NH2	2.43	0.51
3:BM:114:TRP:CZ2	5:BS:37:MET:SD	3.03	0.51
3:BM:178:GLY:CA	3:BM:181:PRO:HG2	2.41	0.51
4:BH:120:PRO:O	4:BH:234:TYR:N	2.40	0.51
5:BA:12:TRP:O	6:BB:9:LEU:HD22	2.09	0.51
6:BG:30:GLY:C	6:BG:33:VAL:HG12	2.34	0.51
5:BQ:32:GLY:HA2	9:BQ:104:BCL:O1D	2.10	0.51
6:BT:22:MET:O	6:BT:26:TYR:HD1	1.93	0.51
6:BT:44:PRO:O	5:BU:55:TYR:OH	2.28	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:BU:102:BCL:CHD	9:BU:102:BCL:CBC	2.88	0.51
6:B8:38:LEU:HD23	6:B8:38:LEU:O	2.10	0.51
1:AC:35:TYR:CZ	3:AM:308:PRO:HG2	2.46	0.51
1:AC:85:LEU:HD22	1:AC:89:GLU:CG	2.40	0.51
1:AC:100:TRP:CB	1:AC:152:CYS:HB2	2.39	0.51
1:AC:102:SER:O	1:AC:104:LYS:N	2.44	0.51
1:AC:138:ASN:HB3	1:AC:331:TYR:CE1	2.45	0.51
2:AL:86:MET:HG3	5:A7:37:MET:HG3	1.92	0.51
2:AL:166:VAL:HG13	9:AL:301:BCL:CHD	2.41	0.51
2:AL:259:ILE:HA	2:AL:263:PHE:HB2	1.92	0.51
3:AM:115:TRP:CD1	3:AM:177:PHE:HD2	2.27	0.51
4:AH:100:LEU:HB2	4:AH:111:PHE:CZ	2.45	0.51
4:AH:259:LEU:HD21	5:A5:19:ARG:HB3	1.93	0.51
5:AA:27:PHE:CE1	5:AD:29:ILE:HD12	2.44	0.51
5:AA:29:ILE:O	5:AA:33:LEU:HD13	2.10	0.51
5:AI:20:VAL:O	5:AI:24:ILE:HG12	2.09	0.51
5:AQ:8:LEU:HD23	6:AR:22:MET:CE	2.38	0.51
5:AS:34:LEU:CA	15:AS:101:PEF:C44	2.89	0.51
5:AU:42:THR:HB	5:AW:48:ASP:HB3	1.92	0.51
5:AY:13:LEU:CD2	6:AZ:14:ALA:HB1	2.37	0.51
5:A5:10:LYS:HB2	14:A5:103:CRT:H83	1.90	0.51
5:A5:14:ILE:CD1	14:A5:103:CRT:H41	2.41	0.51
5:A7:10:LYS:HB3	14:A0:101:CRT:C8	2.40	0.51
6:A8:34:ILE:O	6:A8:37:LEU:HB3	2.09	0.51
1:BC:235:LEU:C	1:BC:237:MET:N	2.67	0.51
1:BC:265:LYS:H	1:BC:265:LYS:HD2	1.75	0.51
2:BL:126:VAL:HB	2:BL:127:PRO:CD	2.40	0.51
3:BM:102:TYR:O	3:BM:104:LEU:N	2.44	0.51
3:BM:121:PHE:N	3:BM:121:PHE:HD1	2.07	0.51
3:BM:228:ARG:H	3:BM:228:ARG:CD	2.23	0.51
6:BV:45:TRP:O	6:BV:46:LEU:HB2	2.10	0.51
5:BY:36:HIS:O	5:BY:40:LEU:HB3	2.11	0.51
5:B3:37:MET:HA	5:B3:40:LEU:HD12	1.91	0.51
5:B5:17:PRO:O	5:B5:21:LEU:HB2	2.09	0.51
5:B5:18:ARG:HG3	5:B5:18:ARG:NH1	2.23	0.51
14:B5:103:CRT:H22A	6:B8:17:PHE:CE1	2.46	0.51
5:B7:13:LEU:O	6:B8:7:THR:HB	2.10	0.51
1:AC:199:PRO:HG2	1:AC:200:LEU:HD12	1.92	0.51
2:AL:150:ALA:O	2:AL:153:HIS:HB3	2.09	0.51
2:AL:218:SER:C	2:AL:220:HIS:N	2.68	0.51
2:AL:237:ALA:CA	2:AL:240:ARG:HG3	2.38	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:AM:80:HIS:O	3:AM:82:ASP:N	2.44	0.51
3:AM:166:VAL:HG22	3:AM:171:TRP:HZ3	1.75	0.51
3:AM:208:PHE:HZ	3:AM:275:LEU:HD13	1.72	0.51
4:AH:106:PRO:HB2	4:AH:249:TYR:CE1	2.44	0.51
4:AH:171:TRP:HE1	4:AH:183:GLU:HG3	1.76	0.51
5:AA:13:LEU:HA	6:AB:9:LEU:HD22	1.92	0.51
5:AA:36:HIS:NE2	9:AB:101:BCL:HMD3	2.25	0.51
5:AA:50:ASN:OD1	5:AA:51:ILE:HG12	2.10	0.51
14:AA:102:CRT:H401	5:AD:38:ILE:HD13	1.91	0.51
5:AD:50:ASN:HB3	5:AF:56:GLN:HA	1.92	0.51
9:AD:102:BCL:HAC2	9:AE:101:BCL:HAC1	1.93	0.51
5:AS:12:TRP:HE1	6:AT:18:HIS:CB	2.23	0.51
9:AV:102:BCL:H191	9:AW:101:BCL:CMC	2.39	0.51
5:AW:26:ALA:C	5:AW:29:ILE:HG22	2.34	0.51
6:AX:21:PHE:CE1	6:AX:25:MET:HE3	2.46	0.51
9:AY:102:BCL:HMD2	6:AZ:36:HIS:CD2	2.45	0.51
9:A1:102:BCL:C8	14:A2:102:CRT:H182	2.39	0.51
5:A3:12:TRP:HA	5:A3:12:TRP:CE3	2.45	0.51
6:A8:20:ILE:HD13	6:A8:20:ILE:C	2.36	0.51
3:BM:83:VAL:HA	3:BM:86:PHE:CB	2.40	0.51
3:BM:84:PHE:CG	5:BW:37:MET:SD	3.04	0.51
3:BM:170:SER:HB3	3:BM:173:LYS:HB2	1.92	0.51
14:BN:102:CRT:H342	9:BO:102:BCL:CBA	2.29	0.51
5:BQ:17:PRO:O	5:BQ:21:LEU:HG	2.10	0.51
5:BW:44:LEU:CD1	5:BY:56:GLN:HB3	2.39	0.51
5:BY:18:ARG:HD2	5:BY:19:ARG:N	2.25	0.51
5:B1:53:VAL:O	5:B1:54:SER:C	2.54	0.51
5:B5:43:ASP:OD1	5:B7:47:LEU:O	2.28	0.51
5:B7:9:TYR:CD1	5:B7:9:TYR:C	2.88	0.51
1:AC:173:LYS:CB	3:AM:80:HIS:HB2	2.40	0.51
2:AL:134:ILE:O	2:AL:137:TYR:HB3	2.10	0.51
2:AL:140:LEU:HD11	9:AL:301:BCL:OBD	2.10	0.51
2:AL:178:TYR:CE1	3:AM:180:PHE:CD2	2.98	0.51
9:AL:303:BCL:O2D	3:AM:203:MET:HB2	2.09	0.51
3:AM:199:ASN:HB2	3:AM:294:TRP:CD2	2.46	0.51
3:AM:274:VAL:C	3:AM:276:THR:H	2.17	0.51
5:AA:13:LEU:O	6:AB:9:LEU:CD1	2.58	0.51
5:AI:14:ILE:CG2	5:AK:18:ARG:HB3	2.40	0.51
5:AU:22:VAL:HG13	5:AU:23:SER:H	1.74	0.51
9:AX:101:BCL:CHC	9:AY:102:BCL:CBB	2.88	0.51
5:AY:7:ASN:O	6:A2:20:ILE:HG12	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:AY:9:TYR:OH	5:AY:10:LYS:HE3	2.10	0.51
5:AY:9:TYR:CD1	6:AZ:15:LYS:HG3	2.45	0.51
14:A7:102:CRT:H342	9:A7:103:BCL:CBA	2.40	0.51
5:A9:44:LEU:CD2	5:A9:46:TRP:HB3	2.41	0.51
1:BC:52:SER:C	1:BC:319:TYR:OH	2.53	0.51
1:BC:167:VAL:CG2	1:BC:297:GLY:HA3	2.40	0.51
1:BC:184:ASN:HD21	3:BM:96:GLU:HG2	1.75	0.51
2:BL:196:LEU:CD1	3:BM:269:ALA:HB1	2.31	0.51
3:BM:104:LEU:HD21	3:BM:169:GLY:CA	2.41	0.51
3:BM:260:VAL:HG13	13:BM:405:MQ8:H142	1.91	0.51
5:BI:36:HIS:CE1	9:BJ:101:BCL:CMD	2.94	0.51
5:BO:4:MET:HE2	6:BR:23:GLN:CG	2.40	0.51
5:BS:53:VAL:O	5:BS:55:TYR:N	2.38	0.51
5:BU:2:PHE:CA	5:BU:5:ASN:HD22	2.23	0.51
5:BU:11:ILE:HG12	14:BU:103:CRT:H83	1.89	0.51
5:BU:20:VAL:O	5:BU:24:ILE:HG12	2.10	0.51
5:BU:36:HIS:CE1	9:BV:101:BCL:CMD	2.92	0.51
5:BW:16:ASP:CA	5:BW:19:ARG:HE	2.24	0.51
5:B1:10:LYS:HD2	6:B4:20:ILE:HB	1.93	0.51
5:B1:18:ARG:CD	5:B1:19:ARG:HG3	2.41	0.51
5:B3:36:HIS:CD2	9:B4:101:BCL:CMD	2.93	0.51
5:B9:44:LEU:CD2	5:B9:46:TRP:HB3	2.40	0.51
1:AC:232:THR:O	1:AC:233:PHE:C	2.54	0.51
3:AM:159:VAL:CG1	3:AM:285:LEU:HD13	2.38	0.51
5:AA:27:PHE:CA	5:AA:30:VAL:HG12	2.39	0.51
9:AP:101:BCL:H2A	9:AP:101:BCL:O1D	2.11	0.51
5:AU:31:LEU:O	5:AU:35:ILE:HG12	2.11	0.51
6:AV:21:PHE:CD1	6:AV:21:PHE:C	2.89	0.51
5:AY:36:HIS:O	5:AY:40:LEU:HB3	2.11	0.51
5:AY:38:ILE:CD1	5:AY:39:VAL:HG23	2.40	0.51
5:AY:45:ASN:O	5:AY:47:LEU:N	2.44	0.51
6:A2:44:PRO:HG2	5:A3:52:PRO:HB2	1.93	0.51
5:A7:7:ASN:HB2	5:A7:10:LYS:HZ2	1.73	0.51
9:A7:103:BCL:HHC	9:A7:103:BCL:OBB	2.10	0.51
5:A9:12:TRP:NE1	6:A0:17:PHE:CE1	2.78	0.51
9:A9:102:BCL:HMD1	6:A0:36:HIS:HA	1.93	0.51
1:BC:46:LYS:C	1:BC:48:GLN:H	2.19	0.51
1:BC:96:ALA:C	1:BC:98:THR:N	2.65	0.51
2:BL:13:ARG:HA	4:BH:99:PRO:HB2	1.91	0.51
2:BL:22:LEU:HB2	5:B7:19:ARG:CG	2.41	0.51
2:BL:184:LEU:HB2	2:BL:252:TRP:HE1	1.76	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:BM:193:TYR:O	3:BM:194:GLY:C	2.52	0.51
3:BM:202:HIS:C	3:BM:204:LEU:H	2.18	0.51
3:BM:202:HIS:C	3:BM:204:LEU:N	2.66	0.51
6:BJ:37:LEU:HD13	9:BJ:101:BCL:H193	1.91	0.51
5:BO:50:ASN:HD21	6:BP:43:ARG:HH22	1.51	0.51
9:BQ:103:BCL:C3D	6:BR:35:ALA:HB1	2.41	0.51
1:AC:157:ARG:NE	1:AC:312:GLN:NE2	2.58	0.51
2:AL:6:PHE:CE2	3:AM:246:GLU:HA	2.45	0.51
2:AL:116:ILE:O	2:AL:118:ARG:N	2.44	0.51
5:AD:35:ILE:HA	5:AD:38:ILE:HG22	1.93	0.51
9:AF:102:BCL:HAC2	9:AG:101:BCL:HBC1	1.92	0.51
9:AG:101:BCL:HMC1	9:AI:102:BCL:HBB1	1.93	0.51
5:AI:43:ASP:O	5:AI:44:LEU:HB3	2.11	0.51
5:AK:9:TYR:HA	6:AN:18:HIS:CG	2.45	0.51
5:AO:51:ILE:O	5:AO:52:PRO:C	2.53	0.51
6:AP:20:ILE:HG23	6:AP:21:PHE:N	2.24	0.51
6:AP:24:SER:O	6:AP:27:ALA:HB3	2.11	0.51
6:AP:36:HIS:CE1	9:AP:101:BCL:C1A	2.93	0.51
6:AT:12:ASP:O	6:AT:15:LYS:HD2	2.10	0.51
5:AW:32:GLY:O	5:AW:35:ILE:HG22	2.10	0.51
5:A5:36:HIS:NE2	9:A6:101:BCL:HMD3	2.25	0.51
5:A7:19:ARG:O	5:A7:23:SER:CB	2.58	0.51
3:BM:84:PHE:CD1	5:BW:37:MET:HE2	2.46	0.51
3:BM:115:TRP:CD1	3:BM:177:PHE:HD2	2.28	0.51
4:BH:58:PHE:N	4:BH:59:PRO:HD2	2.25	0.51
4:BH:258:LEU:HG	5:B5:19:ARG:HE	1.76	0.51
6:BJ:17:PHE:O	6:BJ:18:HIS:C	2.53	0.51
6:BV:34:ILE:O	6:BV:37:LEU:HB2	2.11	0.51
6:BX:21:PHE:CD1	6:BX:21:PHE:C	2.88	0.51
6:BX:36:HIS:CE1	9:BX:101:BCL:C1B	2.93	0.51
6:BX:46:LEU:HD22	6:BZ:42:TYR:CE2	2.46	0.51
5:BY:52:PRO:HD2	5:BY:55:TYR:HE2	1.75	0.51
5:B5:43:ASP:CB	5:B7:47:LEU:HB3	2.40	0.51
1:AC:54:GLN:HA	1:AC:54:GLN:HE21	1.76	0.51
1:AC:225:SER:OG	1:AC:227:LYS:HB3	2.10	0.51
1:AC:254:ARG:HD3	1:AC:254:ARG:C	2.36	0.51
2:AL:71:TRP:CD1	3:AM:303:MET:HE3	2.45	0.51
2:AL:86:MET:HE2	2:AL:96:GLN:OE1	2.11	0.51
2:AL:111:LEU:O	2:AL:114:VAL:N	2.43	0.51
2:AL:115:GLU:O	2:AL:118:ARG:HB2	2.11	0.51
2:AL:116:ILE:HD11	3:AM:254:TRP:O	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:AL:192:ASN:HD22	2:AL:192:ASN:C	2.19	0.51
2:AL:225:PHE:O	2:AL:229:VAL:HG22	2.10	0.51
9:AL:301:BCL:H13	9:AL:301:BCL:H203	1.92	0.51
3:AM:34:PRO:HD3	3:AM:50:PRO:HB3	1.92	0.51
3:AM:138:GLU:C	3:AM:140:LEU:N	2.64	0.51
4:AH:137:ARG:HG2	4:AH:137:ARG:HH11	1.75	0.51
4:AH:195:LEU:HD12	4:AH:196:PRO:CD	2.31	0.51
5:AA:49:ASP:HB2	5:AD:56:GLN:O	2.10	0.51
9:AA:101:BCL:HBC2	9:AB:101:BCL:HHD	1.93	0.51
5:AF:7:ASN:C	6:AJ:20:ILE:HD13	2.35	0.51
5:AI:40:LEU:HD12	5:AI:40:LEU:N	2.26	0.51
9:AK:102:BCL:CED	6:AN:31:LEU:HB3	2.40	0.51
5:AO:12:TRP:CZ3	6:AP:17:PHE:HE2	2.29	0.51
6:AP:44:PRO:HG2	5:AQ:52:PRO:HG3	1.92	0.51
5:AQ:27:PHE:HD1	5:AQ:28:GLN:HE21	1.59	0.51
6:AR:46:LEU:HB3	6:AT:42:TYR:HH	1.72	0.51
5:AU:50:ASN:CB	5:AW:59:GLY:HA3	2.41	0.51
5:A3:35:ILE:HA	5:A3:38:ILE:HG22	1.91	0.51
6:A6:46:LEU:O	5:A7:46:TRP:HB2	2.11	0.51
1:BC:65:ALA:HB1	1:BC:89:GLU:OE1	2.11	0.51
1:BC:190:VAL:C	1:BC:192:TYR:N	2.65	0.51
2:BL:26:TRP:HZ2	3:BM:254:TRP:CZ3	2.28	0.51
2:BL:184:LEU:HB2	2:BL:252:TRP:NE1	2.25	0.51
2:BL:238:ILE:CG2	2:BL:239:HIS:N	2.73	0.51
3:BM:106:ILE:HG12	5:BO:42:THR:HG21	1.92	0.51
3:BM:176:PRO:HD3	3:BM:185:TRP:CD1	2.45	0.51
3:BM:176:PRO:CD	3:BM:185:TRP:HB2	2.41	0.51
3:BM:260:VAL:HG12	4:BH:34:ASP:CB	2.37	0.51
3:BM:277:VAL:HG22	10:BM:403:BPH:HBC1	1.92	0.51
6:BB:34:ILE:HD13	6:BB:34:ILE:C	2.36	0.51
5:BF:44:LEU:CB	6:BG:43:ARG:HH11	2.14	0.51
6:BG:8:GLY:O	6:BG:9:LEU:HD23	2.11	0.51
5:BK:9:TYR:CE1	6:BN:15:LYS:HB2	2.45	0.51
6:BR:43:ARG:NH1	5:BS:55:TYR:CE1	2.79	0.51
9:BS:102:BCL:H192	9:BS:102:BCL:H111	1.93	0.51
5:BU:19:ARG:HB2	5:BU:19:ARG:NH2	2.26	0.51
6:BZ:36:HIS:O	6:BZ:45:TRP:HH2	1.93	0.51
9:B5:102:BCL:OBD	6:B6:32:VAL:HG23	2.10	0.51
5:B7:44:LEU:HD23	6:B8:43:ARG:NH1	2.25	0.51
9:B8:101:BCL:OBB	9:B8:101:BCL:HHC	2.11	0.51
5:B9:5:ASN:HA	5:B9:8:LEU:HG	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:B9:9:TYR:HA	6:B0:18:HIS:CG	2.46	0.51
5:B9:33:LEU:HD12	5:B9:33:LEU:N	2.26	0.51
6:B0:17:PHE:CD1	6:B0:17:PHE:C	2.88	0.51
1:AC:126:VAL:CG2	1:AC:127:SER:N	2.74	0.51
1:AC:276:VAL:CG2	1:AC:280:ASN:ND2	2.74	0.51
2:AL:44:LEU:C	2:AL:46:GLY:N	2.68	0.51
2:AL:226:ARG:O	3:AM:50:PRO:O	2.29	0.51
3:AM:4:TYR:O	3:AM:4:TYR:CD1	2.64	0.51
3:AM:77:ALA:O	3:AM:78:SER:C	2.54	0.51
3:AM:267:ARG:NE	4:AH:29:TYR:OH	2.38	0.51
4:AH:253:GLU:O	4:AH:254:ARG:C	2.53	0.51
5:AS:10:LYS:HB3	14:AS:104:CRT:C3	2.41	0.51
5:AS:27:PHE:CG	5:AU:29:ILE:HD11	2.45	0.51
6:AT:29:PHE:CE1	9:AT:101:BCL:C1	2.92	0.51
5:AW:34:LEU:O	5:AW:38:ILE:HG22	2.11	0.51
5:AY:15:LEU:HG	5:A1:21:LEU:CD2	2.41	0.51
5:AY:44:LEU:CD2	6:AZ:43:ARG:HD2	2.39	0.51
5:A1:5:ASN:C	5:A1:8:LEU:HB3	2.36	0.51
9:A1:102:BCL:HMB3	9:A1:102:BCL:HBB2	1.93	0.51
1:BC:305:VAL:HG11	7:BC:502:HEM:HBC1	1.92	0.51
2:BL:164:ASP:C	2:BL:166:VAL:N	2.68	0.51
2:BL:206:VAL:HG23	2:BL:207:THR:N	2.25	0.51
2:BL:236:LEU:HD13	3:BM:232:ASP:HB3	1.92	0.51
3:BM:259:ASN:N	3:BM:259:ASN:ND2	2.58	0.51
4:BH:5:ILE:O	4:BH:6:THR:HG23	2.11	0.51
4:BH:47:GLU:HG3	5:BA:19:ARG:HA	1.91	0.51
5:BA:33:LEU:CA	14:B0:101:CRT:H2M3	2.37	0.51
9:BG:101:BCL:C1B	9:BI:102:BCL:CMB	2.89	0.51
6:BJ:15:LYS:O	6:BJ:18:HIS:HB3	2.10	0.51
5:BK:49:ASP:OD2	6:BN:43:ARG:NH1	2.40	0.51
5:BO:10:LYS:C	14:BO:103:CRT:H82	2.36	0.51
6:BP:24:SER:O	6:BP:27:ALA:HB3	2.11	0.51
5:BS:46:TRP:CZ3	9:BS:102:BCL:HAC1	2.46	0.51
5:BU:5:ASN:HB3	6:BV:22:MET:CE	2.40	0.51
5:BW:7:ASN:ND2	5:BW:7:ASN:H	2.09	0.51
6:B2:17:PHE:HD1	14:B2:102:CRT:C9	2.20	0.51
5:B3:17:PRO:O	5:B3:21:LEU:HB2	2.11	0.51
5:B5:29:ILE:HG23	5:B5:30:VAL:N	2.26	0.51
6:B8:21:PHE:CG	6:B8:22:MET:N	2.78	0.51
1:AC:85:LEU:HD22	1:AC:89:GLU:HG2	1.93	0.51
1:AC:97:VAL:O	1:AC:97:VAL:CG1	2.57	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AC:135:ARG:CB	1:AC:135:ARG:HH11	2.23	0.51
1:AC:235:LEU:HG	1:AC:239:ILE:CD1	2.40	0.51
1:AC:313:ALA:O	1:AC:314:VAL:HG22	2.11	0.51
3:AM:222:THR:HG21	3:AM:252:TRP:HE1	1.76	0.51
3:AM:224:LEU:HA	3:AM:227:SER:HB2	1.92	0.51
3:AM:301:HIS:HA	4:AH:8:TYR:HB2	1.93	0.51
4:AH:53:VAL:HG13	4:AH:53:VAL:O	2.10	0.51
4:AH:56:VAL:HG23	4:AH:56:VAL:O	2.10	0.51
5:AF:9:TYR:CD1	5:AF:9:TYR:C	2.89	0.51
5:AS:50:ASN:CB	5:AU:60:LYS:HA	2.41	0.51
9:AT:101:BCL:C4A	9:AU:102:BCL:HMB2	2.41	0.51
6:AX:21:PHE:O	6:AX:22:MET:C	2.54	0.51
6:AX:46:LEU:HD13	6:AZ:42:TYR:CZ	2.45	0.51
9:AY:102:BCL:CBC	9:AZ:101:BCL:HHD	2.41	0.51
5:A3:14:ILE:HD13	6:A6:17:PHE:CE2	2.38	0.51
5:A7:17:PRO:O	5:A7:21:LEU:HG	2.11	0.51
2:BL:47:VAL:O	2:BL:50:ILE:HG22	2.11	0.51
2:BL:48:LEU:HD13	5:BA:33:LEU:CD2	2.41	0.51
3:BM:199:ASN:HB2	3:BM:294:TRP:CD2	2.45	0.51
3:BM:279:THR:HA	3:BM:282:ILE:CG1	2.41	0.51
4:BH:19:PHE:CD1	4:BH:20:TRP:N	2.79	0.51
4:BH:36:ARG:HD2	4:BH:78:ALA:HB3	1.93	0.51
4:BH:126:THR:HG22	4:BH:132:LYS:HA	1.93	0.51
9:BG:101:BCL:C1B	9:BI:102:BCL:HMB3	2.40	0.51
9:BG:101:BCL:H203	6:BJ:38:LEU:HD21	1.93	0.51
5:BQ:7:ASN:HB3	6:BT:20:ILE:CG2	2.40	0.51
5:BQ:7:ASN:HB3	6:BT:20:ILE:HG22	1.93	0.51
5:BQ:31:LEU:O	5:BQ:34:LEU:HB3	2.10	0.51
6:BT:40:TRP:CZ3	6:BT:44:PRO:HA	2.45	0.51
5:BY:9:TYR:CG	6:BZ:15:LYS:HG2	2.45	0.51
5:BY:30:VAL:O	5:BY:31:LEU:C	2.53	0.51
5:B1:14:ILE:HD12	5:B1:15:LEU:H	1.74	0.51
5:B3:40:LEU:HD21	5:B3:46:TRP:CZ2	2.45	0.51
5:B3:44:LEU:CD1	5:B3:46:TRP:CE3	2.94	0.51
6:B8:20:ILE:O	6:B8:23:GLN:CG	2.59	0.51
5:B9:17:PRO:O	5:B9:21:LEU:CB	2.58	0.51
5:B9:31:LEU:CD1	5:B9:35:ILE:HD11	2.41	0.51
5:B9:44:LEU:O	5:B9:46:TRP:N	2.41	0.51
6:B0:21:PHE:CB	14:B0:101:CRT:C14	2.78	0.51
1:AC:130:MET:SD	7:AC:502:HEM:NB	2.84	0.51
2:AL:68:TYR:CA	2:AL:73:ILE:HD11	2.35	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:AL:101:CYS:O	2:AL:103:ALA:N	2.44	0.51
2:AL:199:HIS:C	2:AL:201:SER:N	2.69	0.51
3:AM:4:TYR:HE1	3:AM:6:ASN:HA	1.75	0.51
3:AM:58:THR:HB	3:AM:62:PHE:CE2	2.46	0.51
4:AH:54:LYS:HG3	4:AH:58:PHE:HD1	1.76	0.51
6:AJ:38:LEU:HD23	6:AJ:38:LEU:O	2.10	0.51
9:AO:102:BCL:H111	9:AO:102:BCL:C19	2.41	0.51
5:AS:42:THR:HG22	5:AS:43:ASP:N	2.26	0.51
5:AW:27:PHE:CE1	14:AX:102:CRT:H30	2.45	0.51
9:A1:102:BCL:H2A	9:A1:102:BCL:O1D	2.11	0.51
5:A3:8:LEU:HD23	6:A6:20:ILE:HD11	1.93	0.51
6:A8:45:TRP:CE2	9:A8:101:BCL:H2C	2.46	0.51
1:BC:275:HIS:O	1:BC:275:HIS:HD2	1.93	0.51
2:BL:224:PHE:HE1	3:BM:137:ALA:HA	1.76	0.51
3:BM:220:GLY:O	3:BM:224:LEU:HG	2.11	0.51
3:BM:248:ALA:O	3:BM:249:ALA:C	2.54	0.51
4:BH:69:LEU:CB	4:BH:70:PRO:CD	2.89	0.51
9:BA:101:BCL:HHC	9:BA:101:BCL:OBB	2.11	0.51
5:BF:12:TRP:HA	5:BF:12:TRP:CE3	2.46	0.51
9:BI:102:BCL:CBC	9:BJ:101:BCL:HAC1	2.41	0.51
6:BN:45:TRP:CZ3	9:BN:101:BCL:HAC2	2.46	0.51
9:BO:102:BCL:HAC2	9:BP:101:BCL:HBC1	1.92	0.51
5:BU:35:ILE:HG21	9:BV:101:BCL:C2D	2.40	0.51
6:BV:20:ILE:CG2	14:BV:102:CRT:C9	2.87	0.51
5:B5:28:GLN:HE22	9:B6:101:BCL:HED1	1.74	0.51
5:B7:46:TRP:CZ3	9:B7:103:BCL:CBC	2.94	0.51
6:B8:20:ILE:O	6:B8:20:ILE:HD13	2.11	0.51
1:AC:212:ILE:CD1	7:AC:503:HEM:HAA1	2.41	0.50
2:AL:156:PRO:CG	2:AL:162:HIS:HA	2.41	0.50
2:AL:203:ILE:O	2:AL:206:VAL:HG22	2.12	0.50
2:AL:217:THR:H	2:AL:220:HIS:HD1	1.58	0.50
3:AM:63:PHE:CD2	3:AM:124:LEU:HB2	2.46	0.50
3:AM:264:SER:O	3:AM:267:ARG:N	2.37	0.50
9:AM:401:BCL:HHC	9:AM:401:BCL:OBB	2.11	0.50
4:AH:27:ILE:C	4:AH:27:ILE:HD13	2.37	0.50
4:AH:153:GLY:H	4:AH:167:VAL:HG23	1.76	0.50
5:AF:52:PRO:HB2	5:AF:55:TYR:CE1	2.32	0.50
9:AK:102:BCL:CAD	9:AN:101:BCL:CAD	2.89	0.50
9:AK:102:BCL:CBD	9:AN:101:BCL:CAD	2.89	0.50
9:AO:102:BCL:HAC2	9:AP:101:BCL:HBC1	1.92	0.50
6:AX:32:VAL:O	6:AX:36:HIS:N	2.41	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:A3:104:BCL:C2	6:A4:29:PHE:HD1	2.24	0.50
9:BL:301:BCL:H191	9:BL:303:BCL:H8	1.93	0.50
3:BM:52:TYR:CE2	3:BM:136:ARG:NE	2.79	0.50
4:BH:52:ARG:HB2	4:BH:54:LYS:NZ	2.26	0.50
4:BH:66:THR:O	4:BH:66:THR:HG23	2.11	0.50
4:BH:138:VAL:C	4:BH:140:LYS:HD3	2.36	0.50
5:BA:31:LEU:HD12	5:BA:34:LEU:HD23	1.92	0.50
9:BE:101:BCL:CBB	9:BE:101:BCL:CMB	2.89	0.50
5:BF:23:SER:O	5:BF:26:ALA:HB3	2.11	0.50
5:BK:4:MET:SD	6:BP:27:ALA:HB2	2.51	0.50
6:BP:40:TRP:CZ3	6:BP:44:PRO:HA	2.45	0.50
6:BX:21:PHE:CD1	6:BX:22:MET:N	2.79	0.50
9:B2:101:BCL:C4B	9:B3:102:BCL:HBB3	2.41	0.50
5:B7:29:ILE:HG23	5:B7:30:VAL:N	2.26	0.50
1:AC:153:TYR:O	1:AC:157:ARG:N	2.44	0.50
1:AC:187:SER:C	1:AC:189:THR:N	2.68	0.50
1:AC:200:LEU:O	1:AC:204:LEU:N	2.44	0.50
1:AC:302:PRO:O	1:AC:304:ARG:N	2.42	0.50
2:AL:4:LEU:HD12	3:AM:250:LEU:CD1	2.39	0.50
2:AL:155:PHE:HA	2:AL:165:TRP:CD1	2.46	0.50
2:AL:185:ALA:CB	2:AL:252:TRP:HB3	2.41	0.50
2:AL:187:SER:O	2:AL:190:PHE:HB2	2.11	0.50
6:AB:20:ILE:HG21	14:AB:102:CRT:C8	2.41	0.50
6:AE:42:TYR:CE2	6:AE:43:ARG:HG3	2.45	0.50
6:AG:17:PHE:C	6:AG:17:PHE:HD1	2.18	0.50
6:AG:30:GLY:O	6:AG:34:ILE:CG2	2.60	0.50
6:AG:38:LEU:HA	6:AG:41:LEU:CD1	2.39	0.50
5:AO:29:ILE:HA	9:AO:102:BCL:C1	2.37	0.50
6:AR:10:THR:HB	6:AR:13:GLU:OE2	2.11	0.50
6:AR:29:PHE:N	6:AR:29:PHE:CD1	2.79	0.50
5:AS:52:PRO:O	5:AS:53:VAL:C	2.53	0.50
5:AY:11:ILE:CG2	5:AY:15:LEU:HD12	2.41	0.50
5:A1:19:ARG:HH21	5:A3:18:ARG:HH21	1.58	0.50
14:A1:103:CRT:H343	9:A5:102:BCL:HBA1	1.92	0.50
5:A3:51:ILE:HA	5:A3:54:SER:H	1.75	0.50
14:A7:102:CRT:C34	9:A7:103:BCL:HBA1	2.41	0.50
1:BC:92:ARG:O	1:BC:95:VAL:HB	2.11	0.50
2:BL:4:LEU:HB2	2:BL:7:GLU:HB2	1.93	0.50
2:BL:48:LEU:HD23	2:BL:51:VAL:HG21	1.94	0.50
2:BL:110:ALA:HB2	2:BL:134:ILE:HD11	1.93	0.50
3:BM:166:VAL:HG22	3:BM:171:TRP:HZ3	1.75	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:BM:268:TRP:CD1	3:BM:268:TRP:H	2.29	0.50
4:BH:94:PRO:HG2	6:B0:8:GLY:CA	2.35	0.50
14:BA:102:CRT:C32	5:BD:31:LEU:HD21	2.41	0.50
6:BB:34:ILE:HD13	6:BB:34:ILE:O	2.10	0.50
9:BG:101:BCL:CHB	9:BI:102:BCL:HMB3	2.41	0.50
5:BK:9:TYR:HA	6:BN:18:HIS:CG	2.46	0.50
5:BK:46:TRP:HA	5:BK:49:ASP:CG	2.35	0.50
5:BO:27:PHE:C	5:BO:27:PHE:CD1	2.89	0.50
9:BO:102:BCL:H71	6:BP:28:TRP:CE3	2.46	0.50
6:BP:45:TRP:O	6:BP:46:LEU:CG	2.59	0.50
5:BQ:46:TRP:CZ2	9:BQ:103:BCL:CHC	2.94	0.50
5:BS:4:MET:HG3	5:BS:5:ASN:N	2.26	0.50
5:BY:28:GLN:O	9:BY:102:BCL:H11	2.11	0.50
9:B3:102:BCL:HMD1	6:B4:36:HIS:ND1	2.26	0.50
5:B7:24:ILE:O	5:B7:28:GLN:HB2	2.11	0.50
1:AC:147:GLU:HB2	1:AC:322:GLN:HE22	1.77	0.50
1:AC:236:MET:CE	7:AC:503:HEM:ND	2.74	0.50
2:AL:26:TRP:HZ2	3:AM:254:TRP:CZ3	2.29	0.50
2:AL:140:LEU:O	2:AL:141:VAL:HB	2.10	0.50
3:AM:229:PHE:CD1	3:AM:229:PHE:N	2.78	0.50
4:AH:5:ILE:HG21	5:AD:42:THR:OG1	2.11	0.50
4:AH:13:GLN:HE21	16:AH:302:PO4:P	2.33	0.50
4:AH:123:CYS:SG	4:AH:230:GLN:HB2	2.51	0.50
4:AH:159:LEU:HD23	4:AH:214:ILE:N	2.27	0.50
4:AH:193:VAL:O	4:AH:193:VAL:HG23	2.11	0.50
5:AA:17:PRO:HB2	5:A9:14:ILE:HD13	1.93	0.50
5:AF:10:LYS:CB	14:AJ:102:CRT:H5	2.41	0.50
5:AF:44:LEU:HD22	6:AG:43:ARG:CD	2.36	0.50
5:AI:18:ARG:CG	5:AI:18:ARG:NH1	2.73	0.50
5:AK:18:ARG:HG2	5:AK:18:ARG:HH11	1.76	0.50
5:AO:52:PRO:C	5:AO:54:SER:H	2.19	0.50
6:AR:29:PHE:N	6:AR:29:PHE:HD1	2.08	0.50
6:AR:29:PHE:HD1	6:AR:29:PHE:H	1.58	0.50
6:AV:46:LEU:HD13	6:AX:42:TYR:CE1	2.46	0.50
6:AZ:38:LEU:HD23	6:AZ:38:LEU:C	2.36	0.50
9:A2:101:BCL:C4A	9:A3:103:BCL:HMB3	2.41	0.50
6:A8:20:ILE:O	6:A8:20:ILE:HD13	2.11	0.50
6:A0:29:PHE:N	6:A0:29:PHE:CD1	2.78	0.50
2:BL:113:GLU:CB	2:BL:127:PRO:HG3	2.36	0.50
3:BM:28:LEU:HB3	3:BM:29:PRO:CD	2.37	0.50
5:BF:12:TRP:HA	5:BF:12:TRP:HE3	1.76	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:BJ:45:TRP:CD1	6:BJ:46:LEU:N	2.80	0.50
5:BO:7:ASN:CB	6:BR:20:ILE:HD12	2.35	0.50
5:BO:33:LEU:O	5:BO:37:MET:HG2	2.11	0.50
9:BP:101:BCL:CMC	5:BQ:47:LEU:CD2	2.89	0.50
5:B1:48:ASP:O	5:B1:49:ASP:HB3	2.10	0.50
14:B2:102:CRT:H2M1	5:B3:36:HIS:CB	2.21	0.50
5:B7:35:ILE:CD1	9:B8:101:BCL:O1D	2.59	0.50
2:AL:213:GLU:OE2	2:AL:214:PRO:HD2	2.11	0.50
3:AM:98:PRO:CD	3:AM:171:TRP:HB3	2.41	0.50
4:AH:106:PRO:HA	4:AH:109:SER:HB3	1.92	0.50
4:AH:126:THR:HG22	4:AH:132:LYS:HA	1.94	0.50
5:AA:22:VAL:C	5:AA:24:ILE:H	2.19	0.50
5:AD:26:ALA:O	5:AD:29:ILE:HG22	2.11	0.50
5:AF:19:ARG:HH22	5:AI:18:ARG:CZ	2.21	0.50
5:AQ:18:ARG:HA	5:AQ:21:LEU:HD12	1.92	0.50
6:AR:45:TRP:HD1	6:AR:46:LEU:N	2.09	0.50
5:AU:15:LEU:HB3	5:AU:20:VAL:HG21	1.93	0.50
9:AV:102:BCL:HHC	9:AV:102:BCL:OBB	2.12	0.50
5:A1:14:ILE:CD1	5:A1:15:LEU:HG	2.41	0.50
6:A2:42:TYR:CD1	6:A2:43:ARG:HG3	2.45	0.50
9:A6:101:BCL:OBB	9:A6:101:BCL:HHC	2.11	0.50
5:A7:7:ASN:CB	5:A7:10:LYS:HZ2	2.24	0.50
5:A7:40:LEU:HD13	5:A7:46:TRP:CZ2	2.47	0.50
5:A9:4:MET:HA	5:A9:4:MET:HE3	1.93	0.50
5:A9:53:VAL:O	5:A9:54:SER:C	2.55	0.50
1:BC:91:THR:O	1:BC:92:ARG:C	2.51	0.50
1:BC:204:LEU:CD2	7:BC:504:HEM:HBB1	2.41	0.50
2:BL:237:ALA:O	2:BL:238:ILE:C	2.54	0.50
14:BA:102:CRT:H35	5:BD:31:LEU:HD21	1.93	0.50
9:BB:101:BCL:HBB3	9:BD:102:BCL:C4B	2.40	0.50
6:BJ:22:MET:O	6:BJ:26:TYR:HD1	1.94	0.50
5:BY:35:ILE:O	5:BY:38:ILE:HG13	2.11	0.50
5:BY:40:LEU:HB2	5:BY:46:TRP:CH2	2.47	0.50
5:BY:43:ASP:HB2	5:B1:47:LEU:CD1	2.41	0.50
5:B1:31:LEU:HD23	9:B2:101:BCL:HED3	1.94	0.50
5:B5:32:GLY:HA2	9:B6:101:BCL:HED2	1.93	0.50
5:B5:43:ASP:HB2	5:B7:47:LEU:HB3	1.94	0.50
5:B7:35:ILE:O	5:B7:36:HIS:C	2.54	0.50
6:B8:20:ILE:HD13	6:B8:20:ILE:C	2.36	0.50
5:B9:44:LEU:CD1	5:B9:44:LEU:H	2.25	0.50
1:AC:285:TRP:CZ3	1:AC:302:PRO:CD	2.89	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:AL:11:ARG:HB3	2:AL:26:TRP:CH2	2.45	0.50
2:AL:57:GLY:HA3	2:AL:66:GLN:HG2	1.93	0.50
2:AL:87:ALA:HB3	2:AL:96:GLN:HE21	1.76	0.50
3:AM:134:TYR:CA	3:AM:144:GLN:HE22	2.23	0.50
3:AM:222:THR:O	3:AM:223:ILE:C	2.55	0.50
9:AA:101:BCL:HED1	6:AB:31:LEU:CB	2.41	0.50
5:AD:36:HIS:NE2	9:AE:101:BCL:HMD3	2.26	0.50
6:AE:23:GLN:CG	6:AE:24:SER:N	2.75	0.50
9:AF:102:BCL:H172	9:AF:102:BCL:H111	1.94	0.50
9:AF:102:BCL:HBC2	9:AG:101:BCL:HHD	1.94	0.50
6:AG:28:TRP:CE2	6:AG:32:VAL:HG23	2.44	0.50
6:AR:20:ILE:HD12	14:AR:102:CRT:C10	2.41	0.50
5:AS:24:ILE:HD11	9:AU:102:BCL:H191	1.93	0.50
9:AS:103:BCL:HMB1	9:AS:103:BCL:HBB3	1.93	0.50
5:AW:4:MET:HE3	5:AW:7:ASN:HD21	1.76	0.50
5:AY:43:ASP:CA	5:A1:48:ASP:HB3	2.41	0.50
5:A3:42:THR:O	5:A3:43:ASP:C	2.54	0.50
6:A8:45:TRP:HA	5:A9:52:PRO:CD	2.41	0.50
6:A0:33:VAL:O	6:A0:37:LEU:HB2	2.11	0.50
1:BC:43:TYR:HE1	2:BL:153:HIS:HE2	1.57	0.50
1:BC:187:SER:C	1:BC:189:THR:N	2.68	0.50
1:BC:236:MET:HG3	7:BC:503:HEM:C4A	2.47	0.50
1:BC:270:TRP:CE3	1:BC:271:TYR:HD1	2.29	0.50
1:BC:270:TRP:CZ2	1:BC:274:ARG:NH1	2.78	0.50
2:BL:184:LEU:O	2:BL:185:ALA:C	2.54	0.50
2:BL:192:ASN:HA	2:BL:245:LEU:HD12	1.92	0.50
2:BL:199:HIS:HA	11:BL:304:UQ8:O2	2.12	0.50
2:BL:202:LEU:HD13	2:BL:224:PHE:CD2	2.47	0.50
2:BL:246:ALA:C	2:BL:248:SER:N	2.69	0.50
3:BM:71:ILE:O	3:BM:75:MET:HB2	2.11	0.50
3:BM:253:ARG:NH2	4:BH:41:LEU:HD11	2.27	0.50
3:BM:259:ASN:HD22	3:BM:259:ASN:H	1.58	0.50
5:BA:43:ASP:HB2	5:BD:47:LEU:HD12	1.93	0.50
14:BF:103:CRT:H14	6:BJ:21:PHE:CD2	2.47	0.50
6:BG:45:TRP:O	6:BG:46:LEU:CB	2.60	0.50
9:BG:101:BCL:C2B	9:BI:102:BCL:C1B	2.89	0.50
5:BI:45:ASN:C	5:BI:49:ASP:HB3	2.37	0.50
6:BJ:21:PHE:HE1	6:BJ:25:MET:HE3	1.76	0.50
5:BU:11:ILE:CG1	14:BU:103:CRT:H83	2.40	0.50
9:BU:102:BCL:C2D	9:BV:101:BCL:CMD	2.90	0.50
5:BY:29:ILE:HG13	9:BY:102:BCL:H43	1.91	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:BZ:36:HIS:O	6:BZ:45:TRP:CH2	2.65	0.50
9:BZ:101:BCL:C1B	9:B1:102:BCL:HMB2	2.42	0.50
6:B2:21:PHE:CB	14:B2:102:CRT:H11	2.27	0.50
5:B3:46:TRP:HZ3	9:B3:102:BCL:CBC	2.21	0.50
6:B8:28:TRP:HA	6:B8:31:LEU:HB2	1.92	0.50
3:AM:63:PHE:HD2	3:AM:124:LEU:HB2	1.76	0.50
3:AM:265:ILE:HG23	3:AM:266:HIS:N	2.25	0.50
5:AA:40:LEU:HD12	5:AA:40:LEU:O	2.12	0.50
6:AB:43:ARG:HB3	5:AD:55:TYR:CZ	2.46	0.50
5:AI:16:ASP:HB2	5:AI:19:ARG:HG2	1.94	0.50
5:AK:12:TRP:CD1	6:AN:14:ALA:O	2.65	0.50
5:AU:2:PHE:CD1	5:AU:2:PHE:C	2.90	0.50
5:AW:21:LEU:HD11	9:AW:101:BCL:C14	2.41	0.50
6:AX:45:TRP:CE3	9:AX:101:BCL:HAC2	2.46	0.50
5:AY:4:MET:HB3	5:AY:8:LEU:CG	2.39	0.50
6:A2:16:GLU:C	14:A2:102:CRT:H1M1	2.37	0.50
5:A3:9:TYR:HA	6:A4:18:HIS:ND1	2.27	0.50
1:BC:142:LYS:HA	1:BC:145:VAL:HG23	1.92	0.50
1:BC:265:LYS:HD2	1:BC:265:LYS:N	2.26	0.50
1:BC:285:TRP:CZ3	1:BC:302:PRO:CD	2.91	0.50
1:BC:286:PRO:C	1:BC:288:ASN:H	2.20	0.50
1:BC:301:ASP:HB2	1:BC:302:PRO:CD	2.42	0.50
3:BM:284:ILE:HD11	9:BM:402:BCL:OBD	2.11	0.50
5:BA:47:LEU:HB3	5:B9:43:ASP:HA	1.93	0.50
5:BF:51:ILE:CG2	5:BF:52:PRO:HA	2.38	0.50
9:BJ:101:BCL:HMB3	9:BK:102:BCL:CHB	2.41	0.50
5:BK:12:TRP:CD1	6:BN:17:PHE:HB3	2.47	0.50
5:BK:46:TRP:CA	5:BK:49:ASP:OD1	2.52	0.50
6:BT:10:THR:H	6:BT:13:GLU:CD	2.20	0.50
14:BW:103:CRT:H392	5:BY:35:ILE:CD1	2.42	0.50
6:BX:21:PHE:O	6:BX:22:MET:C	2.54	0.50
5:B3:32:GLY:N	9:B4:101:BCL:HED2	2.26	0.50
1:AC:114:GLY:O	1:AC:116:TRP:CD1	2.65	0.50
1:AC:156:HIS:ND1	1:AC:160:PRO:O	2.44	0.50
1:AC:178:LEU:HD21	3:AM:110:SER:HB2	1.93	0.50
1:AC:254:ARG:HH21	3:AM:295:TYR:HE1	1.59	0.50
1:AC:258:ASP:O	1:AC:261:GLN:HB2	2.12	0.50
1:AC:312:GLN:O	1:AC:313:ALA:HB3	2.11	0.50
1:AC:314:VAL:HG12	1:AC:315:ASN:N	2.27	0.50
2:AL:117:CYS:HB3	2:AL:124:PHE:CD2	2.47	0.50
2:AL:162:HIS:CD2	10:AL:302:BPH:H191	2.46	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:AM:130:TRP:HA	3:AM:150:PHE:CE2	2.46	0.50
3:AM:254:TRP:N	3:AM:254:TRP:HD1	2.08	0.50
3:AM:297:TRP:CZ3	3:AM:303:MET:SD	3.05	0.50
4:AH:29:TYR:CZ	4:AH:33:GLU:HG3	2.46	0.50
4:AH:71:HIS:HE1	4:AH:125:LEU:HD22	1.77	0.50
5:AA:9:TYR:CE1	5:AA:10:LYS:HD3	2.47	0.50
5:AA:9:TYR:C	5:AA:11:ILE:H	2.20	0.50
5:AF:49:ASP:OD2	5:AI:56:GLN:HB3	2.12	0.50
5:AK:44:LEU:HD23	6:AN:43:ARG:NH2	2.27	0.50
5:AO:9:TYR:CB	6:AP:18:HIS:CD2	2.94	0.50
5:AQ:50:ASN:HD21	6:AR:43:ARG:HH22	1.59	0.50
6:AR:36:HIS:O	6:AR:39:ALA:N	2.45	0.50
9:A1:102:BCL:ND	9:A2:101:BCL:CMD	2.70	0.50
6:A2:30:GLY:O	6:A2:33:VAL:HG12	2.11	0.50
5:A5:10:LYS:C	14:A5:103:CRT:H5	2.36	0.50
6:A8:33:VAL:HG22	9:A8:101:BCL:H143	1.93	0.50
1:BC:90:PHE:HD1	1:BC:91:THR:N	2.10	0.50
2:BL:69:ASN:O	2:BL:70:LEU:C	2.54	0.50
2:BL:181:ALA:O	2:BL:182:HIS:C	2.55	0.50
2:BL:187:SER:O	2:BL:190:PHE:HB2	2.12	0.50
9:BL:301:BCL:CBA	9:BL:303:BCL:HBC1	2.34	0.50
4:BH:28:ILE:O	4:BH:29:TYR:C	2.55	0.50
4:BH:39:TYR:HD1	4:BH:40:PRO:HA	1.76	0.50
4:BH:136:MET:CA	4:BH:139:ALA:HB3	2.30	0.50
4:BH:178:GLN:HG3	4:BH:178:GLN:O	2.09	0.50
5:BA:9:TYR:HB2	6:BB:18:HIS:CD2	2.42	0.50
6:BB:17:PHE:CD1	14:BB:102:CRT:C6	2.95	0.50
6:BB:32:VAL:HG21	9:BB:101:BCL:CBA	2.31	0.50
9:BB:101:BCL:H12	14:BB:102:CRT:C25	2.41	0.50
6:BE:8:GLY:O	6:BE:9:LEU:HG	2.12	0.50
5:BF:22:VAL:O	5:BF:25:VAL:HB	2.11	0.50
6:BN:32:VAL:HG21	9:BN:101:BCL:HBA2	1.94	0.50
5:BU:44:LEU:C	5:BU:44:LEU:HD12	2.37	0.50
14:B5:103:CRT:H14	5:B7:21:LEU:HD22	1.93	0.50
5:B9:5:ASN:O	5:B9:6:ALA:C	2.54	0.50
5:B9:33:LEU:O	5:B9:37:MET:HB2	2.11	0.50
1:AC:316:LYS:O	1:AC:320:GLY:N	2.40	0.50
2:AL:5:SER:HB3	4:AH:38:GLY:HA2	1.92	0.50
3:AM:53:LEU:CG	3:AM:58:THR:HG23	2.42	0.50
3:AM:62:PHE:C	3:AM:64:GLY:N	2.69	0.50
3:AM:121:PHE:N	3:AM:121:PHE:HD1	2.09	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:AM:201:PHE:HZ	4:AH:15:THR:HG22	1.77	0.50
4:AH:71:HIS:HE1	4:AH:125:LEU:CD2	2.25	0.50
9:AA:101:BCL:HMD1	6:AB:35:ALA:CB	2.41	0.50
6:AB:25:MET:HE2	14:AB:102:CRT:H242	1.93	0.50
14:AB:102:CRT:C34	9:AD:102:BCL:HBA1	2.41	0.50
5:AD:7:ASN:HD22	5:AD:7:ASN:N	2.05	0.50
5:AF:19:ARG:CZ	5:AI:18:ARG:NH2	2.69	0.50
6:AG:36:HIS:HE1	9:AG:101:BCL:C4A	2.25	0.50
6:AN:45:TRP:CE3	9:AN:101:BCL:H2C	2.46	0.50
14:AP:102:CRT:C2M	5:AQ:37:MET:HG2	2.40	0.50
14:AS:104:CRT:C2M	5:AW:33:LEU:O	2.58	0.50
5:AW:9:TYR:C	5:AW:11:ILE:N	2.68	0.50
6:AZ:46:LEU:O	5:A1:51:ILE:O	2.30	0.50
5:A1:28:GLN:NE2	6:A2:28:TRP:NE1	2.60	0.50
5:A1:43:ASP:HB3	5:A1:44:LEU:HD23	1.94	0.50
1:BC:226:LEU:H	3:BM:173:LYS:HE3	1.77	0.50
2:BL:52:TRP:O	2:BL:55:THR:HB	2.11	0.50
2:BL:168:ASN:O	2:BL:169:VAL:C	2.54	0.50
3:BM:4:TYR:O	3:BM:4:TYR:CD1	2.65	0.50
3:BM:84:PHE:CD2	5:BW:37:MET:SD	3.05	0.50
3:BM:90:PHE:O	3:BM:92:TRP:N	2.45	0.50
3:BM:129:TRP:O	3:BM:150:PHE:HE2	1.95	0.50
3:BM:133:THR:HG22	3:BM:147:SER:OG	2.12	0.50
3:BM:276:THR:O	3:BM:278:ILE:N	2.45	0.50
6:BB:42:TYR:HE2	6:BB:43:ARG:HH21	1.60	0.50
5:BF:43:ASP:OD1	5:BF:43:ASP:C	2.55	0.50
5:BW:9:TYR:CD1	5:BW:9:TYR:C	2.90	0.50
5:BW:30:VAL:O	5:BW:33:LEU:HG	2.11	0.50
5:BY:43:ASP:HB2	5:B1:47:LEU:CG	2.42	0.50
6:B4:40:TRP:CZ3	6:B4:45:TRP:N	2.77	0.50
1:AC:73:SER:HB3	1:AC:83:LYS:CB	2.38	0.50
1:AC:286:PRO:C	1:AC:288:ASN:H	2.19	0.50
2:AL:82:TYR:HA	2:AL:85:ARG:NE	2.26	0.50
2:AL:98:ILE:HG22	2:AL:99:THR:N	2.26	0.50
2:AL:117:CYS:SG	2:AL:124:PHE:HA	2.50	0.50
2:AL:139:VAL:O	2:AL:139:VAL:HG13	2.11	0.50
2:AL:239:HIS:CG	3:AM:223:ILE:HG21	2.46	0.50
2:AL:244:PHE:HA	11:AL:304:UQ8:H45A	1.94	0.50
3:AM:149:ALA:O	3:AM:150:PHE:C	2.54	0.50
4:AH:182:LEU:HD12	4:AH:195:LEU:O	2.12	0.50
4:AH:182:LEU:HD13	4:AH:195:LEU:HD23	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:AG:101:BCL:CHB	9:AI:102:BCL:HMB3	2.42	0.50
5:AI:43:ASP:OD2	9:AJ:101:BCL:CMC	2.59	0.50
6:AJ:33:VAL:HG22	6:AJ:37:LEU:HD23	1.94	0.50
5:AO:49:ASP:CG	5:AO:50:ASN:N	2.61	0.50
5:AS:30:VAL:HG13	5:AS:31:LEU:N	2.26	0.50
5:AU:27:PHE:CE2	5:AW:29:ILE:HD11	2.46	0.50
9:AW:101:BCL:HBC2	9:AW:101:BCL:CHD	2.41	0.50
14:AW:102:CRT:C2M	5:A1:36:HIS:HB3	2.42	0.50
14:AW:102:CRT:H342	9:A1:102:BCL:CBA	2.42	0.50
5:A5:14:ILE:CG2	5:A7:18:ARG:HG2	2.39	0.50
5:A5:24:ILE:HD11	9:A7:103:BCL:H191	1.94	0.50
6:A6:45:TRP:CD1	6:A6:46:LEU:N	2.66	0.50
5:A7:49:ASP:OD2	6:A8:43:ARG:NH1	2.45	0.50
6:A0:21:PHE:HE1	6:A0:25:MET:HB2	1.77	0.50
1:BC:250:CYS:HA	1:BC:263:THR:OG1	2.11	0.50
1:BC:259:TRP:O	1:BC:261:GLN:N	2.45	0.50
1:BC:274:ARG:O	1:BC:277:ARG:HB2	2.11	0.50
2:BL:111:LEU:O	2:BL:114:VAL:HB	2.11	0.50
2:BL:126:VAL:O	2:BL:127:PRO:C	2.54	0.50
2:BL:231:TYR:CG	2:BL:232:SER:N	2.79	0.50
2:BL:279:PRO:CG	5:BY:37:MET:SD	2.99	0.50
3:BM:98:PRO:HB2	3:BM:171:TRP:CB	2.40	0.50
3:BM:251:PHE:O	3:BM:255:THR:OG1	2.25	0.50
4:BH:31:ARG:O	4:BH:34:ASP:HB2	2.11	0.50
4:BH:47:GLU:CG	5:BA:19:ARG:HG3	2.42	0.50
4:BH:138:VAL:HG13	4:BH:140:LYS:HZ3	1.76	0.50
5:BF:46:TRP:NE1	9:BF:102:BCL:OBB	2.45	0.50
6:BG:8:GLY:C	6:BG:9:LEU:HG	2.36	0.50
6:BJ:28:TRP:NE1	6:BJ:32:VAL:HG21	2.27	0.50
5:BO:29:ILE:HG23	5:BO:30:VAL:N	2.27	0.50
5:BS:43:ASP:HA	5:BU:56:GLN:HG3	1.92	0.50
6:BT:38:LEU:HD23	6:BT:38:LEU:O	2.12	0.50
6:BV:45:TRP:O	6:BV:46:LEU:CB	2.60	0.50
5:BW:15:LEU:O	5:BW:17:PRO:HD3	2.11	0.50
6:BX:34:ILE:HG23	6:BX:35:ALA:N	2.26	0.50
6:B0:21:PHE:HE1	6:B0:25:MET:HB2	1.77	0.50
6:B0:29:PHE:HE1	9:B0:102:BCL:H72	1.77	0.50
1:AC:111:HIS:HE1	1:AC:124:LYS:CE	2.25	0.49
1:AC:225:SER:O	1:AC:228:GLN:HB2	2.12	0.49
2:AL:184:LEU:O	2:AL:185:ALA:C	2.54	0.49
3:AM:83:VAL:HG23	3:AM:84:PHE:CD1	2.42	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:AA:10:LYS:HB2	14:AA:102:CRT:H5	1.94	0.49
6:AB:36:HIS:CE1	9:AB:101:BCL:ND	2.80	0.49
5:AI:12:TRP:CH2	6:AJ:17:PHE:CZ	3.01	0.49
5:AI:36:HIS:CE1	9:AJ:101:BCL:CMD	2.94	0.49
5:AI:44:LEU:HA	5:AK:56:GLN:HB3	1.94	0.49
9:AI:102:BCL:HED1	6:AJ:31:LEU:HB3	1.92	0.49
5:AK:20:VAL:O	5:AK:24:ILE:HG13	2.12	0.49
6:AP:41:LEU:HG	6:AP:42:TYR:N	2.27	0.49
5:AS:27:PHE:CD1	5:AU:29:ILE:HD11	2.47	0.49
5:AU:13:LEU:HD22	6:AV:9:LEU:CB	2.41	0.49
5:AU:46:TRP:CD1	5:AU:47:LEU:HD13	2.46	0.49
6:A2:40:TRP:HH2	6:A2:46:LEU:HG	1.77	0.49
6:A6:8:GLY:O	6:A6:9:LEU:HD23	2.12	0.49
5:A7:49:ASP:CG	5:A7:50:ASN:N	2.70	0.49
6:A8:33:VAL:CG1	6:A8:34:ILE:N	2.75	0.49
1:BC:138:ASN:ND2	1:BC:149:GLY:HA3	2.27	0.49
1:BC:173:LYS:NZ	5:BU:42:THR:HG22	2.26	0.49
2:BL:7:GLU:HB2	3:BM:250:LEU:HD11	1.93	0.49
2:BL:71:TRP:NE1	3:BM:303:MET:HE3	2.27	0.49
2:BL:170:GLY:HA3	9:BL:301:BCL:CBC	2.41	0.49
2:BL:188:PHE:C	2:BL:190:PHE:N	2.60	0.49
2:BL:257:ILE:CG2	9:BL:301:BCL:HED2	2.42	0.49
3:BM:234:GLU:O	3:BM:238:ILE:HG12	2.11	0.49
3:BM:252:TRP:CD1	3:BM:252:TRP:H	2.28	0.49
3:BM:264:SER:HB2	4:BH:34:ASP:OD1	2.12	0.49
5:BK:35:ILE:HG13	9:BN:101:BCL:O1D	2.12	0.49
14:BP:102:CRT:H391	5:BQ:36:HIS:HB3	1.94	0.49
6:BR:20:ILE:HG23	6:BR:21:PHE:N	2.26	0.49
14:BW:103:CRT:C2M	5:B1:36:HIS:HB3	2.41	0.49
5:B7:9:TYR:HA	6:B8:18:HIS:ND1	2.27	0.49
2:AL:224:PHE:HE1	3:AM:137:ALA:HA	1.77	0.49
2:AL:233:ILE:H	11:AL:304:UQ8:H10A	1.76	0.49
4:AH:54:LYS:O	4:AH:55:VAL:C	2.55	0.49
5:AA:16:ASP:OD1	5:AA:19:ARG:HB2	2.13	0.49
5:AA:46:TRP:CD1	5:AA:47:LEU:HD22	2.47	0.49
6:AB:20:ILE:CD1	14:AB:102:CRT:H133	2.43	0.49
6:AB:38:LEU:HD23	6:AB:38:LEU:O	2.11	0.49
5:AF:11:ILE:HD12	5:AF:14:ILE:CD1	2.35	0.49
9:AF:102:BCL:H2	6:AG:28:TRP:HH2	1.71	0.49
5:AI:10:LYS:HB3	14:AN:102:CRT:H5	1.94	0.49
5:AK:44:LEU:HD22	5:AK:46:TRP:H	1.77	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:AO:10:LYS:O	5:AO:13:LEU:CD2	2.59	0.49
5:AQ:40:LEU:HD21	5:AQ:47:LEU:HD12	1.94	0.49
6:AR:33:VAL:HG23	9:AR:101:BCL:H143	1.94	0.49
5:AW:26:ALA:CA	5:AW:29:ILE:HG22	2.42	0.49
14:AX:102:CRT:H2M2	5:AY:37:MET:HG2	1.93	0.49
5:AY:5:ASN:OD1	6:AZ:18:HIS:HD2	1.94	0.49
6:A2:17:PHE:CD1	14:A2:102:CRT:H41	2.47	0.49
9:A3:104:BCL:OBB	9:A3:104:BCL:HHC	2.12	0.49
9:A3:104:BCL:HAC2	6:A4:45:TRP:CE3	2.48	0.49
5:A5:17:PRO:O	5:A5:21:LEU:CB	2.60	0.49
6:A0:17:PHE:CD1	6:A0:18:HIS:CA	2.92	0.49
9:A0:102:BCL:HMB1	9:A0:102:BCL:H142	1.94	0.49
1:BC:46:LYS:C	1:BC:48:GLN:N	2.70	0.49
1:BC:123:THR:O	1:BC:127:SER:OG	2.22	0.49
1:BC:153:TYR:O	1:BC:157:ARG:N	2.44	0.49
2:BL:116:ILE:CD1	3:BM:254:TRP:HB2	2.42	0.49
2:BL:250:ALA:HB2	10:BL:302:BPH:HBC2	1.93	0.49
9:BL:301:BCL:HBB3	9:BL:301:BCL:HMB1	1.93	0.49
9:BL:301:BCL:HAC1	3:BM:197:TYR:OH	2.12	0.49
3:BM:12:GLN:CB	4:BH:145:ALA:HB2	2.40	0.49
3:BM:27:ASN:N	3:BM:27:ASN:ND2	2.56	0.49
3:BM:75:MET:HG3	3:BM:94:GLY:N	2.28	0.49
3:BM:195:ASN:HD21	3:BM:197:TYR:HB2	1.77	0.49
4:BH:173:ASP:OD2	4:BH:175:SER:HB2	2.12	0.49
5:BA:22:VAL:HA	5:BA:25:VAL:CG2	2.41	0.49
5:BA:38:ILE:O	5:BA:41:SER:HB3	2.12	0.49
9:BD:102:BCL:C4D	9:BE:101:BCL:HMD3	2.42	0.49
6:BE:9:LEU:HB3	6:BE:13:GLU:HG2	1.94	0.49
6:BE:43:ARG:HH11	5:BF:55:TYR:HD2	1.59	0.49
6:BJ:30:GLY:O	6:BJ:33:VAL:HG12	2.13	0.49
5:BK:18:ARG:HG2	5:BK:18:ARG:NH1	2.27	0.49
9:BU:102:BCL:HMD1	9:BV:101:BCL:HAC1	1.94	0.49
5:BY:31:LEU:HD23	9:BZ:101:BCL:HED3	1.94	0.49
6:B6:16:GLU:OE1	14:B7:102:CRT:H1M1	2.11	0.49
3:AM:209:LEU:HD22	9:AM:402:BCL:H3A	1.94	0.49
3:AM:239:THR:OG1	3:AM:240:HIS:N	2.42	0.49
14:AM:406:CRT:H402	5:AO:38:ILE:HG22	1.93	0.49
4:AH:141:GLU:H	4:AH:141:GLU:CD	2.19	0.49
4:AH:242:TYR:O	4:AH:243:TYR:C	2.55	0.49
5:AA:16:ASP:OD2	5:AA:19:ARG:HB2	2.12	0.49
9:AN:101:BCL:HBB3	9:AO:102:BCL:CHC	2.43	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:AS:36:HIS:CE1	9:AT:101:BCL:HMD2	2.47	0.49
14:A2:102:CRT:C24	9:A3:103:BCL:H18	2.43	0.49
9:A3:103:BCL:OBB	9:A3:103:BCL:HHC	2.12	0.49
5:A5:43:ASP:HB2	5:A7:47:LEU:HD12	1.95	0.49
9:A6:101:BCL:C1B	9:A7:103:BCL:HMB3	2.42	0.49
6:A8:21:PHE:CG	6:A8:22:MET:N	2.78	0.49
1:BC:111:HIS:HE1	1:BC:124:LYS:HE2	1.78	0.49
1:BC:263:THR:HG22	3:BM:311:VAL:CG2	2.42	0.49
1:BC:276:VAL:O	1:BC:277:ARG:C	2.56	0.49
2:BL:86:MET:HE2	2:BL:96:GLN:CD	2.37	0.49
2:BL:92:GLY:O	2:BL:93:GLY:C	2.54	0.49
2:BL:144:ARG:CB	2:BL:145:PRO:HD3	2.41	0.49
3:BM:199:ASN:O	3:BM:203:MET:HE3	2.12	0.49
3:BM:256:MET:HE1	13:BM:405:MQ8:H171	1.93	0.49
4:BH:52:ARG:O	4:BH:54:LYS:N	2.45	0.49
4:BH:168:SER:HB3	4:BH:183:GLU:HB2	1.93	0.49
5:BA:46:TRP:CD1	5:BA:47:LEU:HD22	2.46	0.49
9:BA:101:BCL:HMD2	6:BB:36:HIS:CE1	2.47	0.49
5:BF:51:ILE:HA	5:BF:52:PRO:C	2.37	0.49
6:BJ:17:PHE:HA	6:BJ:20:ILE:HG22	1.93	0.49
5:BS:30:VAL:HG13	5:BS:31:LEU:H	1.76	0.49
5:BW:33:LEU:HD12	5:BW:34:LEU:N	2.27	0.49
14:BW:103:CRT:H36	5:B1:33:LEU:HA	1.94	0.49
5:B5:16:ASP:CB	5:B5:19:ARG:HH21	2.25	0.49
5:B9:31:LEU:HD21	9:B0:102:BCL:HMA3	1.93	0.49
1:AC:236:MET:CA	1:AC:239:ILE:HD12	2.42	0.49
1:AC:293:ALA:O	1:AC:295:ARG:N	2.45	0.49
2:AL:137:TYR:CE2	9:AL:303:BCL:HBB1	2.47	0.49
3:AM:176:PRO:CD	3:AM:185:TRP:HB2	2.42	0.49
9:AM:401:BCL:HMB1	9:AM:401:BCL:CBB	2.42	0.49
4:AH:157:VAL:HG23	4:AH:210:LYS:HA	1.93	0.49
9:AG:101:BCL:C1B	9:AI:102:BCL:CMB	2.90	0.49
5:AI:35:ILE:C	5:AI:37:MET:N	2.70	0.49
6:AJ:20:ILE:HG12	14:AJ:102:CRT:H83	1.93	0.49
5:AO:34:LEU:HG	5:AO:34:LEU:O	2.13	0.49
9:AO:102:BCL:CAD	9:AP:101:BCL:CAD	2.90	0.49
6:AP:17:PHE:HD1	14:AP:102:CRT:H6	1.78	0.49
5:AQ:10:LYS:HB2	14:AT:102:CRT:C8	2.39	0.49
5:AQ:31:LEU:CD2	9:AR:101:BCL:HED3	2.43	0.49
14:AT:102:CRT:H31	9:AU:102:BCL:HBA1	1.92	0.49
5:AU:13:LEU:HD22	6:AV:9:LEU:C	2.37	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:AU:44:LEU:HD13	6:AV:43:ARG:CD	2.42	0.49
5:AW:42:THR:HB	5:AY:48:ASP:HB2	1.95	0.49
5:A1:20:VAL:O	5:A1:24:ILE:HG12	2.12	0.49
1:BC:21:LEU:HD23	1:BC:21:LEU:C	2.36	0.49
1:BC:126:VAL:HG23	1:BC:127:SER:N	2.27	0.49
2:BL:131:SER:O	2:BL:134:ILE:HB	2.11	0.49
2:BL:280:LEU:HD21	5:BY:37:MET:HE1	1.95	0.49
3:BM:60:SER:N	3:BM:128:LEU:HD23	2.27	0.49
3:BM:84:PHE:CD1	3:BM:84:PHE:N	2.80	0.49
9:BM:401:BCL:CBB	9:BM:401:BCL:HMB1	2.42	0.49
4:BH:27:ILE:HD13	4:BH:27:ILE:C	2.37	0.49
4:BH:113:PRO:HD2	4:BH:249:TYR:OH	2.13	0.49
4:BH:154:MET:HB3	4:BH:207:ARG:O	2.12	0.49
4:BH:189:ASN:O	4:BH:191:LYS:HG3	2.11	0.49
6:BN:21:PHE:CD1	6:BN:21:PHE:C	2.91	0.49
5:BY:12:TRP:HE1	6:BZ:18:HIS:CA	2.12	0.49
5:BY:36:HIS:CE1	9:BZ:101:BCL:CMD	2.94	0.49
6:B6:8:GLY:O	6:B6:9:LEU:HD23	2.12	0.49
6:B8:33:VAL:CG1	6:B8:34:ILE:N	2.75	0.49
6:B0:20:ILE:HG23	6:B0:21:PHE:N	2.27	0.49
1:AC:270:TRP:CZ2	1:AC:274:ARG:NH1	2.80	0.49
2:AL:13:ARG:HA	4:AH:99:PRO:HB2	1.94	0.49
2:AL:38:VAL:O	2:AL:39:GLY:C	2.56	0.49
2:AL:116:ILE:O	2:AL:117:CYS:C	2.55	0.49
2:AL:168:ASN:O	2:AL:171:TYR:N	2.44	0.49
2:AL:204:LEU:HD21	3:AM:267:ARG:NH1	2.28	0.49
2:AL:237:ALA:HA	2:AL:240:ARG:CG	2.42	0.49
2:AL:238:ILE:C	2:AL:240:ARG:H	2.19	0.49
9:AM:401:BCL:OBB	14:AM:406:CRT:H243	2.12	0.49
4:AH:27:ILE:HD13	4:AH:27:ILE:O	2.12	0.49
5:AO:13:LEU:O	6:AP:7:THR:N	2.45	0.49
5:AO:36:HIS:NE2	9:AP:101:BCL:HMD3	2.28	0.49
6:AV:46:LEU:HB3	6:AX:42:TYR:CZ	2.47	0.49
5:AY:36:HIS:NE2	9:AZ:101:BCL:CMD	2.72	0.49
5:A7:37:MET:N	14:A7:102:CRT:H2M3	2.25	0.49
1:BC:121:ILE:HG22	1:BC:123:THR:H	1.78	0.49
1:BC:242:SER:O	1:BC:313:ALA:N	2.46	0.49
1:BC:276:VAL:HG22	1:BC:280:ASN:HD22	1.77	0.49
2:BL:93:GLY:O	2:BL:94:LEU:C	2.55	0.49
2:BL:106:PHE:CE1	9:BL:301:BCL:H121	2.48	0.49
3:BM:168:MET:HG2	3:BM:289:THR:CG2	2.41	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:BM:226:VAL:HG22	3:BM:229:PHE:HB2	1.94	0.49
3:BM:261:THR:C	3:BM:263:GLU:H	2.19	0.49
9:BM:401:BCL:HMB1	9:BM:401:BCL:HBB3	1.94	0.49
4:BH:186:VAL:HG12	4:BH:187:ALA:N	2.24	0.49
5:BA:11:ILE:CD1	5:BA:14:ILE:HD11	2.42	0.49
5:BD:36:HIS:NE2	9:BE:101:BCL:CMD	2.76	0.49
9:BE:101:BCL:HMB3	9:BF:102:BCL:CHB	2.42	0.49
5:BK:56:GLN:NE2	5:BK:57:ALA:H	2.09	0.49
9:BK:102:BCL:ND	9:BN:101:BCL:CMD	2.76	0.49
6:BP:17:PHE:CA	6:BP:20:ILE:HG22	2.42	0.49
6:BP:29:PHE:CZ	9:BP:101:BCL:H42	2.48	0.49
6:BX:36:HIS:HE1	9:BX:101:BCL:CHB	2.25	0.49
9:BY:102:BCL:HBC1	9:BZ:101:BCL:HBC3	1.95	0.49
5:B1:29:ILE:O	5:B1:33:LEU:HG	2.12	0.49
5:B3:8:LEU:HD21	6:B6:24:SER:OG	2.13	0.49
9:B4:101:BCL:HHC	9:B4:101:BCL:OBB	2.12	0.49
5:B5:26:ALA:O	5:B5:29:ILE:HG22	2.12	0.49
6:B0:21:PHE:CE1	6:B0:25:MET:HB2	2.48	0.49
6:B0:33:VAL:O	6:B0:37:LEU:CG	2.58	0.49
1:AC:186:GLY:O	3:AM:89:HIS:HE1	1.95	0.49
2:AL:214:PRO:O	2:AL:216:LYS:N	2.46	0.49
2:AL:268:TRP:O	2:AL:270:GLU:N	2.46	0.49
4:AH:232:THR:O	4:AH:235:GLU:HG2	2.12	0.49
9:AG:101:BCL:HMB2	9:AI:102:BCL:C1B	2.42	0.49
6:AR:20:ILE:HG23	6:AR:21:PHE:N	2.26	0.49
5:AS:33:LEU:C	15:AS:101:PEF:H453	2.37	0.49
5:AW:31:LEU:HD22	14:AX:102:CRT:H32	1.95	0.49
9:AY:102:BCL:ND	9:AZ:101:BCL:CMD	2.76	0.49
5:A1:14:ILE:HD12	5:A1:15:LEU:HG	1.94	0.49
5:A1:46:TRP:CZ3	9:A1:102:BCL:H2C	2.48	0.49
5:A3:39:VAL:HA	5:A5:47:LEU:HD11	1.93	0.49
1:BC:138:ASN:ND2	1:BC:150:VAL:H	2.11	0.49
1:BC:138:ASN:HD21	1:BC:150:VAL:H	1.60	0.49
2:BL:268:TRP:O	2:BL:270:GLU:N	2.45	0.49
3:BM:76:LEU:HD21	5:BU:37:MET:HE3	1.95	0.49
3:BM:90:PHE:HA	3:BM:93:LEU:HD12	1.93	0.49
3:BM:163:ILE:HG23	3:BM:285:LEU:HD11	1.94	0.49
3:BM:185:TRP:CH2	3:BM:189:PHE:CD1	3.00	0.49
3:BM:200:PRO:C	3:BM:203:MET:HG2	2.37	0.49
3:BM:253:ARG:HB3	3:BM:254:TRP:HD1	1.77	0.49
3:BM:265:ILE:CD1	13:BM:405:MQ8:H143	2.42	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:BD:102:BCL:CMD	6:BE:36:HIS:CD2	2.96	0.49
6:BE:29:PHE:CD1	9:BE:101:BCL:C1	2.95	0.49
6:BR:45:TRP:CD1	6:BR:46:LEU:N	2.81	0.49
5:BS:9:TYR:CE1	6:BT:15:LYS:HG2	2.48	0.49
5:BY:30:VAL:O	5:BY:33:LEU:HG	2.13	0.49
6:BZ:29:PHE:O	6:BZ:33:VAL:HG12	2.13	0.49
5:B1:38:ILE:HG23	5:B1:39:VAL:N	2.26	0.49
5:B3:51:ILE:HB	5:B3:52:PRO:HA	1.95	0.49
5:B7:33:LEU:HD12	5:B7:33:LEU:N	2.28	0.49
6:B0:21:PHE:O	6:B0:24:SER:N	2.45	0.49
1:AC:236:MET:HE1	7:AC:503:HEM:ND	2.27	0.49
2:AL:120:LEU:HD21	3:AM:254:TRP:CZ2	2.47	0.49
2:AL:155:PHE:HA	2:AL:165:TRP:NE1	2.27	0.49
2:AL:188:PHE:HD2	2:AL:249:ALA:N	2.11	0.49
2:AL:206:VAL:C	2:AL:209:PRO:HD3	2.37	0.49
2:AL:276:LEU:O	3:AM:88:LYS:HE3	2.13	0.49
9:AM:402:BCL:HHC	9:AM:402:BCL:OBB	2.13	0.49
4:AH:119:ARG:O	4:AH:234:TYR:HB2	2.11	0.49
5:AA:39:VAL:C	5:AA:41:SER:N	2.70	0.49
5:AI:30:VAL:HA	5:AI:33:LEU:CD2	2.43	0.49
15:AS:101:PEF:O1P	5:AU:22:VAL:HG21	2.13	0.49
5:AU:2:PHE:C	5:AU:2:PHE:HD1	2.20	0.49
9:AV:102:BCL:H191	9:AW:101:BCL:HMC3	1.94	0.49
5:AY:27:PHE:CD1	5:AY:27:PHE:C	2.91	0.49
5:A1:10:LYS:HB3	14:A1:103:CRT:O1	2.13	0.49
6:A2:43:ARG:HD3	5:A3:55:TYR:CD2	2.48	0.49
9:A2:101:BCL:HMB1	9:A2:101:BCL:CBB	2.42	0.49
5:A7:40:LEU:HD11	5:A7:47:LEU:HD23	1.93	0.49
5:A7:51:ILE:HB	5:A7:52:PRO:HA	1.94	0.49
6:A8:20:ILE:O	6:A8:23:GLN:CG	2.59	0.49
9:A8:101:BCL:HHB	9:A9:102:BCL:HMA1	1.95	0.49
9:A9:102:BCL:CHA	9:A0:102:BCL:OBD	2.61	0.49
6:A0:21:PHE:CE1	6:A0:25:MET:HB2	2.48	0.49
1:BC:205:ASP:OD1	1:BC:304:ARG:CZ	2.60	0.49
1:BC:270:TRP:O	1:BC:274:ARG:CD	2.61	0.49
2:BL:159:ILE:H	2:BL:159:ILE:CD1	2.24	0.49
3:BM:11:VAL:HG11	4:BH:151:PRO:HD3	1.94	0.49
4:BH:108:LEU:C	4:BH:110:GLY:H	2.20	0.49
5:BA:39:VAL:C	5:BA:41:SER:H	2.20	0.49
6:BB:29:PHE:HE1	9:BB:101:BCL:C1	2.25	0.49
6:BB:38:LEU:O	6:BB:38:LEU:HD23	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:BE:42:TYR:CE2	6:BE:43:ARG:HG3	2.48	0.49
6:BG:11:ASP:O	6:BG:15:LYS:HG3	2.12	0.49
6:BG:28:TRP:C	6:BG:30:GLY:N	2.66	0.49
5:BI:35:ILE:C	5:BI:37:MET:N	2.71	0.49
6:BJ:45:TRP:CZ3	9:BJ:101:BCL:HAC2	2.47	0.49
5:BK:18:ARG:O	5:BK:22:VAL:HG12	2.13	0.49
5:BS:49:ASP:CG	5:BS:50:ASN:OD1	2.55	0.49
6:BT:13:GLU:CD	6:BT:13:GLU:H	2.20	0.49
5:BY:48:ASP:O	5:BY:49:ASP:CB	2.57	0.49
5:BY:49:ASP:CA	5:B1:56:GLN:HE22	2.26	0.49
5:B1:10:LYS:HB3	14:B1:103:CRT:H5	1.94	0.49
5:B9:48:ASP:CG	5:B9:48:ASP:O	2.56	0.49
1:AC:66:ASP:O	1:AC:67:SER:HB3	2.12	0.49
1:AC:184:ASN:HD21	3:AM:96:GLU:HG2	1.77	0.49
1:AC:236:MET:HG3	7:AC:503:HEM:C4A	2.47	0.49
1:AC:259:TRP:O	1:AC:262:SER:N	2.45	0.49
2:AL:13:ARG:HD2	4:AH:101:VAL:HG22	1.93	0.49
2:AL:204:LEU:HD21	3:AM:267:ARG:CZ	2.43	0.49
2:AL:230:GLY:HA2	3:AM:51:ILE:CB	2.37	0.49
2:AL:238:ILE:CG2	2:AL:239:HIS:N	2.74	0.49
3:AM:131:VAL:O	3:AM:135:LYS:HB2	2.13	0.49
3:AM:200:PRO:HA	3:AM:203:MET:HE3	1.95	0.49
3:AM:244:ALA:C	3:AM:246:GLU:N	2.69	0.49
4:AH:146:GLU:H	4:AH:146:GLU:CD	2.19	0.49
4:AH:185:GLU:HB2	4:AH:192:LYS:NZ	2.27	0.49
5:AA:16:ASP:CG	5:AA:19:ARG:HB2	2.37	0.49
6:AG:46:LEU:HD13	6:AJ:42:TYR:CZ	2.48	0.49
9:AG:101:BCL:OBB	9:AG:101:BCL:HHC	2.13	0.49
6:AJ:46:LEU:HD13	6:AN:42:TYR:CE1	2.48	0.49
5:AK:49:ASP:OD2	6:AN:43:ARG:NH1	2.44	0.49
9:AN:101:BCL:HHC	9:AN:101:BCL:OBB	2.12	0.49
6:AP:23:GLN:O	6:AP:26:TYR:N	2.46	0.49
6:AP:45:TRP:O	6:AP:46:LEU:HD23	2.13	0.49
6:AR:13:GLU:OE2	6:AR:13:GLU:N	2.46	0.49
9:AS:103:BCL:C1D	9:AT:101:BCL:HMD1	2.43	0.49
5:AY:11:ILE:CD1	9:A1:102:BCL:H151	2.30	0.49
6:A2:16:GLU:HB3	14:A2:102:CRT:C1M	2.41	0.49
9:A3:103:BCL:C1D	9:A3:104:BCL:HMD1	2.41	0.49
6:A4:40:TRP:CZ3	6:A4:45:TRP:N	2.77	0.49
5:A5:33:LEU:HD12	5:A5:33:LEU:C	2.37	0.49
5:A7:14:ILE:HG22	5:A7:14:ILE:O	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:A7:103:BCL:HMD3	6:A8:36:HIS:CD2	2.47	0.49
5:A9:24:ILE:HG21	14:A0:101:CRT:H243	1.95	0.49
1:BC:41:GLU:OE1	1:BC:43:TYR:OH	2.21	0.49
1:BC:122:TYR:CA	1:BC:125:VAL:HG23	2.38	0.49
1:BC:138:ASN:HB3	1:BC:331:TYR:CE1	2.48	0.49
2:BL:78:PRO:O	2:BL:152:GLY:HA3	2.13	0.49
2:BL:129:ALA:CB	2:BL:247:LEU:HD11	2.42	0.49
2:BL:148:MET:HE3	2:BL:153:HIS:HE1	1.76	0.49
2:BL:171:TYR:C	2:BL:173:PHE:N	2.66	0.49
3:BM:41:GLY:HA3	3:BM:46:ALA:CB	2.38	0.49
3:BM:218:MET:O	3:BM:221:ALA:N	2.46	0.49
4:BH:108:LEU:HD23	4:BH:108:LEU:N	2.28	0.49
5:BA:33:LEU:HD12	5:BA:33:LEU:N	2.27	0.49
5:BF:38:ILE:HG23	5:BF:39:VAL:HG23	1.95	0.49
6:BG:20:ILE:HG23	6:BG:21:PHE:N	2.28	0.49
6:BJ:41:LEU:HD23	6:BJ:41:LEU:C	2.38	0.49
6:BP:23:GLN:O	6:BP:26:TYR:N	2.46	0.49
14:BW:103:CRT:H6	6:BZ:17:PHE:HD1	1.77	0.49
6:BX:38:LEU:HD23	6:BX:38:LEU:O	2.12	0.49
5:BY:51:ILE:HB	5:BY:52:PRO:CA	2.42	0.49
1:AC:166:TRP:O	1:AC:166:TRP:CD2	2.66	0.49
2:AL:3:MET:HG2	2:AL:11:ARG:CZ	2.42	0.49
2:AL:13:ARG:HA	4:AH:99:PRO:CB	2.42	0.49
2:AL:71:TRP:O	2:AL:160:LEU:HG	2.13	0.49
3:AM:275:LEU:HA	3:AM:278:ILE:HD12	1.93	0.49
5:AA:32:GLY:HA2	9:AB:101:BCL:HED2	1.95	0.49
5:AD:32:GLY:HA3	9:AD:102:BCL:O1A	2.13	0.49
6:AE:45:TRP:O	6:AE:46:LEU:CG	2.58	0.49
6:AE:46:LEU:O	5:AF:46:TRP:O	2.30	0.49
5:AI:34:LEU:O	5:AI:37:MET:HB2	2.12	0.49
5:AI:50:ASN:HA	5:AK:60:LYS:HA	1.94	0.49
9:AO:102:BCL:C1D	9:AP:101:BCL:C2D	2.90	0.49
5:AS:12:TRP:HE1	6:AT:18:HIS:HD1	1.59	0.49
9:AW:101:BCL:OBB	9:AW:101:BCL:HHC	2.12	0.49
5:A3:56:GLN:H	5:A3:56:GLN:CD	2.20	0.49
5:A9:17:PRO:O	5:A9:21:LEU:CB	2.61	0.49
6:A0:40:TRP:HA	6:A0:40:TRP:HE3	1.78	0.49
1:BC:196:PRO:HG3	1:BC:231:TRP:NE1	2.28	0.49
2:BL:218:SER:C	2:BL:220:HIS:N	2.69	0.49
2:BL:268:TRP:O	2:BL:269:PRO:C	2.56	0.49
3:BM:74:ASN:ND2	3:BM:95:LEU:HD13	2.28	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:BM:199:ASN:HB2	3:BM:294:TRP:CG	2.48	0.49
5:BA:37:MET:HE2	14:B0:101:CRT:H2M1	1.94	0.49
5:BI:15:LEU:HD22	5:BI:15:LEU:N	2.28	0.49
5:BI:39:VAL:HG12	5:BI:46:TRP:HZ3	1.78	0.49
9:BQ:103:BCL:CBB	9:BQ:103:BCL:CMB	2.90	0.49
5:BW:10:LYS:HD2	6:BZ:20:ILE:CD1	2.42	0.49
6:BX:10:THR:N	6:BX:13:GLU:OE1	2.37	0.49
5:BY:49:ASP:CB	5:B1:56:GLN:HE22	2.24	0.49
9:BY:102:BCL:C1D	9:BZ:101:BCL:HMD1	2.43	0.49
5:B3:2:PHE:CE1	5:B3:5:ASN:CG	2.84	0.49
6:B4:41:LEU:HD23	6:B4:41:LEU:C	2.38	0.49
5:B7:43:ASP:OD2	5:B9:47:LEU:HD13	2.12	0.49
5:B9:33:LEU:HD12	5:B9:33:LEU:H	1.78	0.49
5:B9:35:ILE:CG1	9:B0:102:BCL:O1D	2.58	0.49
1:AC:236:MET:HG3	7:AC:503:HEM:CHB	2.43	0.49
2:AL:89:LEU:HD12	2:AL:89:LEU:N	2.26	0.49
2:AL:89:LEU:HB3	2:AL:94:LEU:HB2	1.95	0.49
2:AL:175:HIS:CD2	2:AL:178:TYR:CE2	3.01	0.49
2:AL:253:SER:HB2	9:AL:301:BCL:C2A	2.37	0.49
3:AM:32:GLY:C	3:AM:34:PRO:HD3	2.38	0.49
3:AM:157:TYR:CD1	3:AM:158:LEU:HD23	2.47	0.49
3:AM:286:LEU:O	3:AM:290:VAL:HB	2.12	0.49
9:AM:401:BCL:HMB1	9:AM:401:BCL:HBB3	1.94	0.49
4:AH:46:THR:HG22	4:AH:47:GLU:N	2.28	0.49
5:AA:47:LEU:HD22	5:AA:47:LEU:N	2.27	0.49
6:AB:44:PRO:HG2	5:AD:52:PRO:HB3	1.93	0.49
5:AF:38:ILE:HD13	14:AG:102:CRT:H401	1.93	0.49
5:AO:51:ILE:HG12	5:AO:52:PRO:CD	2.40	0.49
5:AS:47:LEU:H	5:AS:47:LEU:CD2	2.24	0.49
6:AV:45:TRP:O	6:AV:46:LEU:CB	2.60	0.49
5:AY:26:ALA:O	5:AY:30:VAL:HG23	2.13	0.49
5:A1:24:ILE:C	5:A1:26:ALA:N	2.69	0.49
5:A5:5:ASN:HA	5:A5:8:LEU:CD1	2.42	0.49
5:A5:40:LEU:HD11	5:A5:47:LEU:HB2	1.95	0.49
1:BC:165:ALA:HB1	1:BC:303:LEU:CB	2.30	0.49
1:BC:212:ILE:H	7:BC:503:HEM:CGA	2.26	0.49
2:BL:16:THR:HG21	2:BL:20:GLY:C	2.38	0.49
2:BL:30:PHE:CD2	3:BM:255:THR:O	2.66	0.49
2:BL:138:LEU:C	2:BL:140:LEU:N	2.66	0.49
2:BL:151:TRP:C	2:BL:153:HIS:N	2.67	0.49
3:BM:150:PHE:CE1	3:BM:154:ILE:HD11	2.48	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:BA:101:BCL:HBB1	9:B0:102:BCL:CMC	2.42	0.49
6:BB:44:PRO:HB2	5:BD:52:PRO:HG3	1.95	0.49
5:BD:30:VAL:O	5:BD:34:LEU:N	2.42	0.49
9:BK:102:BCL:HMB1	9:BK:102:BCL:HBB3	1.95	0.49
6:BT:45:TRP:CD1	6:BT:46:LEU:N	2.81	0.49
14:BV:102:CRT:H2M2	5:BW:37:MET:HG2	1.94	0.49
5:BW:44:LEU:HD13	5:BY:56:GLN:HB3	1.95	0.49
6:B2:29:PHE:HE1	9:B2:101:BCL:C2	2.25	0.49
9:B2:101:BCL:OBB	9:B2:101:BCL:HHC	2.12	0.49
5:B3:39:VAL:CG1	5:B3:44:LEU:HG	2.43	0.49
5:B3:56:GLN:N	5:B3:56:GLN:HE21	2.10	0.49
5:B5:46:TRP:HA	5:B5:49:ASP:OD1	2.12	0.49
9:B5:102:BCL:OBB	9:B5:102:BCL:HHC	2.13	0.49
5:B7:29:ILE:O	5:B7:33:LEU:CD1	2.61	0.49
6:B8:28:TRP:HA	6:B8:31:LEU:HD12	1.94	0.49
1:AC:151:THR:HG21	1:AC:323:MET:CB	2.43	0.48
2:AL:10:TYR:CZ	3:AM:246:GLU:HG2	2.48	0.48
2:AL:16:THR:HG21	2:AL:20:GLY:C	2.38	0.48
2:AL:204:LEU:HD21	3:AM:267:ARG:CD	2.41	0.48
2:AL:220:HIS:O	2:AL:223:THR:HG23	2.13	0.48
3:AM:194:GLY:H	3:AM:293:ASN:HA	1.78	0.48
3:AM:196:LEU:O	3:AM:198:TYR:N	2.46	0.48
9:AA:101:BCL:CED	6:AB:31:LEU:HB3	2.43	0.48
14:AA:102:CRT:H81	6:AE:20:ILE:HG21	1.94	0.48
5:AI:28:GLN:O	9:AJ:101:BCL:HED1	2.12	0.48
5:AK:35:ILE:HA	5:AK:38:ILE:HG22	1.95	0.48
6:AN:32:VAL:HG11	9:AN:101:BCL:HBA2	1.94	0.48
6:AN:43:ARG:HD2	5:AO:55:TYR:CE2	2.48	0.48
5:AO:11:ILE:HG22	5:AO:15:LEU:CD1	2.42	0.48
6:AP:46:LEU:HD13	6:AR:42:TYR:OH	2.12	0.48
5:AY:39:VAL:HG22	5:A1:47:LEU:HD11	1.95	0.48
6:A4:41:LEU:HD23	6:A4:41:LEU:C	2.38	0.48
5:A5:5:ASN:HA	5:A5:8:LEU:CG	2.42	0.48
5:A7:43:ASP:CB	5:A9:47:LEU:HD12	2.31	0.48
5:A9:31:LEU:HD11	5:A9:35:ILE:HD11	1.95	0.48
9:A9:102:BCL:C1D	9:A0:102:BCL:HMD1	2.42	0.48
1:BC:94:MET:SD	7:BC:501:HEM:NB	2.86	0.48
1:BC:137:ALA:C	1:BC:139:SER:N	2.70	0.48
1:BC:167:VAL:CG2	1:BC:298:PRO:HD2	2.38	0.48
2:BL:41:CYS:HA	5:B9:30:VAL:CG2	2.42	0.48
2:BL:137:TYR:HD1	2:BL:138:LEU:HD12	1.78	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:BL:243:LEU:HD12	3:BM:217:ALA:O	2.13	0.48
2:BL:253:SER:O	2:BL:256:CYS:HB3	2.13	0.48
3:BM:173:LYS:O	3:BM:185:TRP:HZ2	1.96	0.48
5:BD:31:LEU:O	5:BD:35:ILE:N	2.39	0.48
5:BF:33:LEU:H	5:BF:33:LEU:CD1	2.24	0.48
5:BQ:35:ILE:HD11	9:BS:102:BCL:HMB2	1.95	0.48
6:BT:27:ALA:O	6:BT:31:LEU:HG	2.13	0.48
9:BW:102:BCL:CHA	9:BX:101:BCL:OBD	2.60	0.48
6:BZ:40:TRP:HB2	6:BZ:45:TRP:CH2	2.47	0.48
6:BZ:44:PRO:O	5:B1:55:TYR:CZ	2.66	0.48
6:B2:21:PHE:CB	14:B2:102:CRT:C11	2.90	0.48
5:B3:19:ARG:O	5:B3:23:SER:N	2.44	0.48
5:B5:14:ILE:O	5:B5:14:ILE:HG22	2.11	0.48
9:B7:103:BCL:HHC	9:B7:103:BCL:OBB	2.13	0.48
1:AC:225:SER:O	1:AC:226:LEU:C	2.55	0.48
1:AC:274:ARG:HH11	1:AC:274:ARG:CG	2.24	0.48
1:AC:282:ASN:C	1:AC:283:TYR:CD1	2.88	0.48
2:AL:110:ALA:HB2	2:AL:134:ILE:HD11	1.95	0.48
2:AL:182:HIS:HA	2:AL:256:CYS:SG	2.54	0.48
3:AM:98:PRO:CG	3:AM:107:PRO:HG3	2.44	0.48
3:AM:210:TYR:O	3:AM:214:LEU:N	2.46	0.48
9:AM:401:BCL:HBC1	9:AM:402:BCL:CAD	2.43	0.48
5:AA:8:LEU:CD2	6:AE:20:ILE:HG23	2.41	0.48
5:AD:30:VAL:HG13	5:AD:31:LEU:N	2.28	0.48
5:AF:28:GLN:CB	9:AF:102:BCL:H12	2.27	0.48
6:AJ:17:PHE:CZ	14:AJ:102:CRT:H6	2.48	0.48
9:AQ:102:BCL:HMD2	9:AR:101:BCL:C1D	2.43	0.48
6:AT:45:TRP:CD2	9:AT:101:BCL:H2C	2.48	0.48
5:AU:8:LEU:O	5:AU:10:LYS:N	2.46	0.48
5:AU:28:GLN:OE1	6:AV:28:TRP:CZ2	2.66	0.48
5:AU:44:LEU:C	5:AU:44:LEU:HD12	2.38	0.48
5:A1:40:LEU:O	5:A1:40:LEU:HG	2.12	0.48
5:A5:25:VAL:HG13	9:A5:102:BCL:H51	1.96	0.48
14:A5:103:CRT:C7	6:A8:17:PHE:HZ	2.26	0.48
14:A5:103:CRT:C32	5:A7:31:LEU:HD21	2.43	0.48
5:A7:36:HIS:CB	14:A7:102:CRT:C39	2.90	0.48
6:A8:28:TRP:HA	6:A8:31:LEU:HD12	1.94	0.48
1:BC:36:ARG:NH1	2:BL:91:GLU:C	2.71	0.48
1:BC:137:ALA:O	1:BC:139:SER:N	2.46	0.48
1:BC:183:GLN:NE2	1:BC:230:GLU:HG2	2.28	0.48
1:BC:200:LEU:O	1:BC:204:LEU:N	2.45	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BC:203:PHE:CG	1:BC:235:LEU:HD22	2.48	0.48
1:BC:213:THR:HG22	1:BC:214:GLY:O	2.13	0.48
2:BL:52:TRP:HA	2:BL:52:TRP:HE3	1.76	0.48
2:BL:53:GLY:C	2:BL:55:THR:N	2.68	0.48
2:BL:115:GLU:O	2:BL:118:ARG:HB2	2.13	0.48
2:BL:195:ALA:O	2:BL:198:MET:N	2.47	0.48
4:BH:27:ILE:HG23	4:BH:28:ILE:N	2.28	0.48
4:BH:76:VAL:HG12	4:BH:77:VAL:N	2.27	0.48
5:BI:17:PRO:HG3	6:BJ:9:LEU:HD11	1.94	0.48
6:BJ:34:ILE:HD13	6:BJ:34:ILE:C	2.37	0.48
6:BR:28:TRP:HA	6:BR:28:TRP:CE3	2.46	0.48
5:BS:35:ILE:O	5:BS:39:VAL:HG23	2.13	0.48
5:BU:38:ILE:HD12	14:BV:102:CRT:H401	1.94	0.48
5:BW:54:SER:CB	5:BW:57:ALA:HB3	2.43	0.48
5:BY:27:PHE:C	5:BY:27:PHE:HD1	2.21	0.48
5:B1:56:GLN:O	5:B1:60:LYS:N	2.39	0.48
5:B3:12:TRP:HA	5:B3:12:TRP:HE3	1.78	0.48
6:B0:40:TRP:CH2	6:B0:46:LEU:CG	2.83	0.48
1:AC:148:THR:HA	1:AC:322:GLN:CG	2.43	0.48
1:AC:324:ALA:C	1:AC:331:TYR:HH	2.19	0.48
2:AL:6:PHE:O	2:AL:9:LYS:HG2	2.12	0.48
2:AL:142:PHE:HD1	2:AL:143:VAL:N	2.11	0.48
2:AL:196:LEU:HD23	3:AM:216:PHE:CD1	2.48	0.48
9:AL:301:BCL:H202	10:AL:302:BPH:HMA3	1.94	0.48
3:AM:27:ASN:HD22	3:AM:27:ASN:N	2.11	0.48
3:AM:103:GLY:O	3:AM:171:TRP:N	2.47	0.48
3:AM:133:THR:O	3:AM:134:TYR:C	2.56	0.48
3:AM:221:ALA:O	3:AM:224:LEU:N	2.46	0.48
3:AM:264:SER:HB3	4:AH:34:ASP:HA	1.94	0.48
4:AH:156:VAL:HG12	4:AH:157:VAL:N	2.27	0.48
5:AA:36:HIS:CE1	9:AB:101:BCL:CMD	2.96	0.48
9:AD:102:BCL:OBD	6:AE:32:VAL:HG23	2.14	0.48
5:AF:43:ASP:CB	5:AI:47:LEU:HG	2.42	0.48
5:AO:50:ASN:O	5:AO:51:ILE:C	2.56	0.48
5:AS:33:LEU:O	15:AS:101:PEF:H453	2.12	0.48
9:AS:103:BCL:OBD	6:AT:32:VAL:HG23	2.13	0.48
9:AT:101:BCL:HBB2	9:AT:101:BCL:H151	1.95	0.48
6:AX:36:HIS:HE1	9:AX:101:BCL:NB	2.06	0.48
9:A5:102:BCL:HHC	9:A5:102:BCL:OBB	2.12	0.48
5:A7:10:LYS:HB3	14:A0:101:CRT:H83	1.94	0.48
5:A9:2:PHE:CE1	6:A0:26:TYR:OH	2.66	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BC:53:ILE:HG12	1:BC:319:TYR:CZ	2.48	0.48
1:BC:253:THR:CG2	2:BL:171:TYR:HD2	2.26	0.48
9:BL:301:BCL:HMB1	9:BL:301:BCL:CBB	2.43	0.48
3:BM:297:TRP:CZ2	4:BH:13:GLN:HB2	2.49	0.48
4:BH:52:ARG:HB3	4:BH:52:ARG:CZ	2.43	0.48
4:BH:151:PRO:O	4:BH:167:VAL:HG21	2.13	0.48
6:BB:24:SER:HA	5:B9:4:MET:CE	2.42	0.48
6:BE:38:LEU:HD23	6:BE:38:LEU:C	2.38	0.48
5:BI:9:TYR:HA	6:BJ:18:HIS:CG	2.48	0.48
5:BI:9:TYR:CD1	5:BI:9:TYR:C	2.91	0.48
5:BK:26:ALA:HA	5:BK:29:ILE:CG2	2.44	0.48
5:BQ:36:HIS:CD2	9:BQ:103:BCL:NB	2.81	0.48
6:BR:29:PHE:HD1	6:BR:29:PHE:N	2.06	0.48
5:BS:47:LEU:HD22	5:BS:47:LEU:H	1.78	0.48
6:BT:28:TRP:CE3	6:BT:28:TRP:HA	2.47	0.48
5:BW:12:TRP:HA	5:BW:12:TRP:CE3	2.48	0.48
9:BW:102:BCL:HMD2	6:BX:36:HIS:HA	1.94	0.48
5:B3:9:TYR:OH	5:B3:10:LYS:HE3	2.13	0.48
5:B5:51:ILE:HA	5:B5:52:PRO:C	2.36	0.48
5:B7:19:ARG:O	5:B7:23:SER:CB	2.60	0.48
5:B7:42:THR:CB	5:B9:48:ASP:CG	2.84	0.48
9:B7:103:BCL:HAC2	9:B8:101:BCL:HAC1	1.95	0.48
6:B8:44:PRO:HG2	5:B9:52:PRO:HB2	1.96	0.48
5:B9:29:ILE:HA	9:B9:102:BCL:H11	1.96	0.48
1:AC:53:ILE:C	1:AC:55:ALA:H	2.20	0.48
1:AC:66:ASP:OD2	1:AC:88:GLY:HA3	2.13	0.48
1:AC:138:ASN:HB3	1:AC:331:TYR:CD1	2.48	0.48
1:AC:306:SER:OG	1:AC:307:CYS:N	2.46	0.48
2:AL:22:LEU:CB	5:A7:19:ARG:HB3	2.34	0.48
2:AL:116:ILE:C	2:AL:118:ARG:N	2.71	0.48
2:AL:207:THR:HA	2:AL:215:VAL:HG13	1.94	0.48
3:AM:98:PRO:HA	3:AM:112:GLY:HA3	1.95	0.48
3:AM:264:SER:HB2	4:AH:34:ASP:OD1	2.13	0.48
4:AH:5:ILE:CD1	5:AF:47:LEU:CD1	2.89	0.48
4:AH:178:GLN:NE2	4:AH:180:ARG:CZ	2.77	0.48
14:AB:102:CRT:H2M2	5:AD:37:MET:HE3	1.94	0.48
9:AI:102:BCL:CBC	9:AJ:101:BCL:HHD	2.43	0.48
6:AN:7:THR:OG1	6:AN:8:GLY:N	2.45	0.48
6:AN:22:MET:O	6:AN:25:MET:HB3	2.13	0.48
6:AN:36:HIS:HD1	9:AN:101:BCL:H162	1.77	0.48
6:AT:45:TRP:CZ3	9:AT:101:BCL:HAC2	2.49	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:AY:28:GLN:HB3	9:AY:102:BCL:C1	2.43	0.48
5:A1:54:SER:CB	5:A1:57:ALA:HB2	2.44	0.48
1:BC:98:THR:O	1:BC:103:PRO:CD	2.58	0.48
2:BL:164:ASP:O	2:BL:166:VAL:N	2.46	0.48
3:BM:79:VAL:O	3:BM:80:HIS:C	2.55	0.48
3:BM:84:PHE:CE1	5:BW:37:MET:HG2	2.47	0.48
3:BM:133:THR:O	3:BM:137:ALA:N	2.40	0.48
3:BM:179:ILE:O	3:BM:183:LEU:N	2.43	0.48
9:BM:402:BCL:HMB1	9:BM:402:BCL:HBB3	1.96	0.48
4:BH:150:ASP:OD1	4:BH:152:ARG:HB2	2.13	0.48
6:BE:46:LEU:N	5:BF:52:PRO:HD3	2.28	0.48
9:BG:101:BCL:CBB	9:BI:102:BCL:C4B	2.88	0.48
5:BK:16:ASP:HB2	5:BK:19:ARG:CD	2.44	0.48
5:BS:42:THR:CG2	5:BS:43:ASP:N	2.76	0.48
6:BV:45:TRP:CZ3	9:BV:101:BCL:HAC2	2.48	0.48
5:B7:36:HIS:CE1	9:B8:101:BCL:CMD	2.90	0.48
6:B8:28:TRP:C	6:B8:30:GLY:N	2.70	0.48
5:B9:40:LEU:HD12	5:B9:45:ASN:HA	1.94	0.48
1:AC:47:ARG:O	1:AC:47:ARG:HG2	2.12	0.48
1:AC:167:VAL:O	1:AC:168:THR:C	2.56	0.48
1:AC:314:VAL:HG12	1:AC:315:ASN:H	1.79	0.48
2:AL:40:PHE:O	2:AL:41:CYS:C	2.56	0.48
2:AL:199:HIS:CE1	2:AL:239:HIS:CE1	3.01	0.48
2:AL:229:VAL:C	3:AM:51:ILE:HD12	2.38	0.48
2:AL:231:TYR:CZ	2:AL:233:ILE:HA	2.49	0.48
2:AL:243:LEU:O	2:AL:247:LEU:HB2	2.14	0.48
3:AM:83:VAL:O	3:AM:86:PHE:HB3	2.14	0.48
3:AM:182:HIS:ND1	3:AM:183:LEU:N	2.61	0.48
3:AM:204:LEU:HD23	3:AM:279:THR:CG2	2.43	0.48
4:AH:6:THR:HB	5:AF:41:SER:OG	2.13	0.48
4:AH:9:ILE:HG23	4:AH:9:ILE:O	2.14	0.48
4:AH:48:ARG:HD3	15:AH:301:PEF:C4	2.41	0.48
4:AH:154:MET:HB3	4:AH:207:ARG:O	2.13	0.48
5:AF:16:ASP:HB2	5:AF:19:ARG:CB	2.44	0.48
6:AP:7:THR:OG1	6:AP:8:GLY:N	2.45	0.48
5:AU:35:ILE:HA	5:AU:38:ILE:CG2	2.44	0.48
5:AW:10:LYS:HB3	14:AW:102:CRT:H5	1.95	0.48
9:AW:101:BCL:C1D	9:AX:101:BCL:HMD1	2.44	0.48
5:AY:18:ARG:HG2	5:AY:18:ARG:NH1	2.27	0.48
5:A1:10:LYS:HB2	14:A1:103:CRT:C8	2.37	0.48
5:A1:49:ASP:O	5:A1:50:ASN:HB3	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:A3:50:ASN:O	5:A3:54:SER:N	2.45	0.48
9:A7:103:BCL:OBD	6:A8:32:VAL:HG22	2.12	0.48
5:A9:12:TRP:CD1	6:A0:18:HIS:HB2	2.47	0.48
5:A9:13:LEU:O	6:A0:7:THR:CB	2.61	0.48
9:A9:102:BCL:HBC2	9:A0:102:BCL:CMD	2.38	0.48
6:A0:21:PHE:O	6:A0:24:SER:N	2.45	0.48
1:BC:247:CYS:O	1:BC:248:THR:C	2.54	0.48
2:BL:101:CYS:O	2:BL:102:ALA:C	2.56	0.48
9:BM:401:BCL:HBB2	9:BM:402:BCL:H111	1.95	0.48
5:BA:36:HIS:CE1	9:BB:101:BCL:CMD	2.97	0.48
9:BA:101:BCL:HBB3	9:B0:102:BCL:C1C	2.43	0.48
6:BB:28:TRP:O	6:BB:31:LEU:N	2.46	0.48
5:BD:12:TRP:HE1	6:BE:18:HIS:HA	1.78	0.48
6:BE:21:PHE:HD1	6:BE:22:MET:N	2.12	0.48
5:BF:7:ASN:CB	6:BJ:20:ILE:HD13	2.42	0.48
5:BF:8:LEU:HD21	6:BJ:24:SER:HG	1.77	0.48
14:BF:103:CRT:H401	5:BI:38:ILE:CD1	2.44	0.48
6:BG:21:PHE:HZ	9:BI:102:BCL:H202	1.79	0.48
9:BG:101:BCL:C3B	9:BI:102:BCL:C3B	2.90	0.48
9:BG:101:BCL:HHC	9:BG:101:BCL:OBB	2.14	0.48
5:BO:40:LEU:HD23	5:BO:40:LEU:O	2.13	0.48
5:BS:21:LEU:O	5:BS:25:VAL:HG23	2.13	0.48
5:BS:46:TRP:NE1	9:BS:102:BCL:OBB	2.47	0.48
5:BU:12:TRP:CE2	6:BV:17:PHE:HE2	2.14	0.48
5:B1:11:ILE:N	14:B1:103:CRT:H82	2.16	0.48
5:B1:13:LEU:CD1	14:B1:103:CRT:H1M3	2.43	0.48
6:B2:28:TRP:O	6:B2:32:VAL:HG23	2.13	0.48
5:B7:49:ASP:OD2	6:B8:43:ARG:NH1	2.45	0.48
9:B9:102:BCL:H12	9:B0:102:BCL:HED1	1.94	0.48
1:AC:37:GLY:HA3	2:AL:77:PRO:HG2	1.95	0.48
2:AL:80:LEU:HD21	2:AL:153:HIS:CD2	2.47	0.48
2:AL:129:ALA:CA	2:AL:247:LEU:HD11	2.38	0.48
3:AM:7:ILE:HG22	3:AM:8:PHE:CG	2.48	0.48
3:AM:14:ARG:HH12	4:AH:145:ALA:CB	2.27	0.48
3:AM:156:PHE:HZ	9:AM:402:BCL:HBD	1.78	0.48
3:AM:161:GLY:HA3	14:AM:406:CRT:C29	2.43	0.48
9:AB:101:BCL:CHB	9:AD:102:BCL:HMB3	2.43	0.48
5:AD:50:ASN:CG	5:AD:51:ILE:N	2.68	0.48
5:AF:17:PRO:HG2	5:AF:18:ARG:H	1.79	0.48
5:AK:51:ILE:O	5:AK:51:ILE:HG13	2.13	0.48
6:AR:42:TYR:CE2	6:AR:43:ARG:HG3	2.48	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:AU:102:BCL:HBC1	9:AV:102:BCL:HBC3	1.94	0.48
6:AV:19:ALA:O	6:AV:23:GLN:HG3	2.13	0.48
6:AV:21:PHE:CD1	6:AV:21:PHE:O	2.67	0.48
9:AV:102:BCL:H203	6:AX:39:ALA:HB1	1.96	0.48
9:A1:102:BCL:C2D	9:A2:101:BCL:C2D	2.91	0.48
9:A8:101:BCL:CMC	9:A9:102:BCL:CBB	2.92	0.48
1:BC:238:ASN:O	1:BC:239:ILE:C	2.57	0.48
2:BL:102:ALA:HB2	10:BL:302:BPH:H112	1.95	0.48
2:BL:116:ILE:O	2:BL:117:CYS:C	2.56	0.48
2:BL:139:VAL:HG23	2:BL:143:VAL:CB	2.43	0.48
2:BL:231:TYR:CZ	2:BL:233:ILE:HA	2.49	0.48
2:BL:253:SER:CB	9:BL:301:BCL:HED3	2.42	0.48
3:BM:61:ILE:HG12	3:BM:129:TRP:CZ3	2.47	0.48
3:BM:101:GLN:C	3:BM:103:GLY:H	2.22	0.48
3:BM:122:LEU:HD13	9:BM:402:BCL:H203	1.96	0.48
3:BM:206:ILE:O	3:BM:207:ALA:C	2.57	0.48
6:BB:44:PRO:HG2	5:BD:52:PRO:CB	2.44	0.48
5:BF:4:MET:CG	6:BJ:23:GLN:CG	2.76	0.48
9:BI:102:BCL:CMD	6:BJ:36:HIS:CD2	2.96	0.48
9:BN:101:BCL:OBB	9:BN:101:BCL:HHC	2.13	0.48
6:BT:29:PHE:HD1	6:BT:29:PHE:H	1.58	0.48
14:BV:102:CRT:H2M2	5:BW:37:MET:CG	2.44	0.48
9:BX:101:BCL:HHC	9:BX:101:BCL:OBB	2.13	0.48
5:BY:32:GLY:HA2	9:BZ:101:BCL:O1D	2.13	0.48
6:BZ:38:LEU:HD23	6:BZ:38:LEU:C	2.38	0.48
6:B2:33:VAL:O	6:B2:37:LEU:HD23	2.13	0.48
5:B3:10:LYS:C	14:B7:102:CRT:H82	2.37	0.48
9:B9:102:BCL:CHD	9:B0:102:BCL:HMD1	2.44	0.48
6:B0:45:TRP:CD1	6:B0:46:LEU:N	2.80	0.48
1:AC:152:CYS:O	1:AC:156:HIS:HB2	2.13	0.48
2:AL:20:GLY:C	2:AL:22:LEU:N	2.71	0.48
2:AL:207:THR:CB	3:AM:238:ILE:HG21	2.44	0.48
2:AL:221:GLU:O	2:AL:223:THR:N	2.46	0.48
3:AM:4:TYR:CE1	3:AM:6:ASN:HA	2.47	0.48
3:AM:10:ALA:HA	15:AM:408:PEF:C4	2.44	0.48
3:AM:35:ILE:HG13	15:AM:409:PEF:H321	1.96	0.48
3:AM:121:PHE:N	3:AM:121:PHE:CD1	2.79	0.48
3:AM:137:ALA:O	3:AM:140:LEU:HB3	2.12	0.48
3:AM:284:ILE:HG13	9:AM:402:BCL:OBD	2.14	0.48
4:AH:5:ILE:HG12	4:AH:6:THR:HG23	1.94	0.48
4:AH:259:LEU:CD2	5:A5:19:ARG:HB3	2.42	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:AG:34:ILE:O	6:AG:38:LEU:HB3	2.13	0.48
14:AG:102:CRT:H2M1	5:AI:36:HIS:HB3	1.95	0.48
5:AI:28:GLN:NE2	14:AJ:102:CRT:H25	2.29	0.48
9:AI:102:BCL:OBD	6:AJ:32:VAL:HG13	2.14	0.48
5:AK:12:TRP:NE1	6:AN:17:PHE:CD2	2.69	0.48
5:AK:46:TRP:CH2	9:AK:102:BCL:HBC3	2.49	0.48
5:AW:46:TRP:HA	5:AW:49:ASP:OD1	2.14	0.48
9:AX:101:BCL:HBA1	9:AX:101:BCL:H3A	1.53	0.48
5:AY:27:PHE:HE2	5:A1:29:ILE:CD1	2.27	0.48
5:A1:10:LYS:C	14:A1:103:CRT:H82	2.38	0.48
5:A1:21:LEU:CD1	9:A1:102:BCL:C14	2.88	0.48
5:A3:19:ARG:HH21	5:A3:20:VAL:HG23	1.79	0.48
1:BC:33:ILE:H	1:BC:33:ILE:CD1	2.24	0.48
1:BC:66:ASP:O	1:BC:67:SER:HB3	2.14	0.48
2:BL:38:VAL:O	2:BL:39:GLY:C	2.57	0.48
2:BL:105:ALA:HB1	10:BL:302:BPH:H2	1.94	0.48
2:BL:132:PHE:CD2	2:BL:247:LEU:HD13	2.49	0.48
2:BL:154:GLY:O	2:BL:165:TRP:NE1	2.46	0.48
2:BL:207:THR:CB	3:BM:238:ILE:HG21	2.43	0.48
3:BM:170:SER:OG	3:BM:173:LYS:HD3	2.14	0.48
4:BH:2:SER:HA	16:BH:301:PO4:O3	2.14	0.48
6:BE:24:SER:O	6:BE:27:ALA:HB3	2.14	0.48
5:BF:9:TYR:CD1	5:BF:9:TYR:C	2.91	0.48
5:BK:16:ASP:HB2	5:BK:19:ARG:CG	2.44	0.48
5:BS:9:TYR:CD1	5:BS:9:TYR:C	2.91	0.48
5:BU:45:ASN:O	5:BU:49:ASP:HB3	2.13	0.48
14:BU:103:CRT:C2M	5:BY:36:HIS:HB2	2.43	0.48
9:BW:102:BCL:HMB1	9:BW:102:BCL:HBB3	1.96	0.48
5:B1:36:HIS:CE1	9:B2:101:BCL:CMD	2.97	0.48
5:B3:12:TRP:HA	5:B3:12:TRP:CE3	2.49	0.48
6:B4:18:HIS:CD2	6:B4:18:HIS:O	2.67	0.48
6:B4:21:PHE:CZ	9:B5:102:BCL:H203	2.48	0.48
2:AL:50:ILE:HA	2:AL:98:ILE:HD11	1.96	0.48
2:AL:72:ARG:O	2:AL:73:ILE:C	2.57	0.48
2:AL:261:GLY:O	2:AL:263:PHE:N	2.47	0.48
2:AL:271:TRP:C	2:AL:273:ASN:H	2.21	0.48
3:AM:55:LEU:O	3:AM:59:LEU:N	2.43	0.48
3:AM:75:MET:HG3	3:AM:94:GLY:H	1.79	0.48
3:AM:229:PHE:CE2	3:AM:247:ARG:NE	2.81	0.48
3:AM:265:ILE:C	3:AM:267:ARG:N	2.71	0.48
3:AM:276:THR:C	3:AM:278:ILE:N	2.72	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:AH:77:VAL:O	4:AH:80:ARG:CD	2.56	0.48
4:AH:227:ASN:ND2	4:AH:228:PRO:CD	2.64	0.48
6:AE:24:SER:O	6:AE:27:ALA:HB3	2.14	0.48
5:AI:20:VAL:HA	5:AI:23:SER:OG	2.14	0.48
9:AN:101:BCL:H3A	9:AN:101:BCL:HBA1	1.46	0.48
5:AQ:35:ILE:CA	5:AQ:38:ILE:HG22	2.36	0.48
5:AU:36:HIS:O	5:AU:40:LEU:HB2	2.13	0.48
5:AY:35:ILE:O	5:AY:36:HIS:C	2.54	0.48
5:A1:39:VAL:HG11	9:A1:102:BCL:HBC1	1.95	0.48
9:A1:102:BCL:HBC2	9:A2:101:BCL:HMD1	1.95	0.48
1:BC:274:ARG:HG2	1:BC:274:ARG:HH11	1.77	0.48
2:BL:239:HIS:O	3:BM:224:LEU:HD11	2.13	0.48
3:BM:14:ARG:HH12	4:BH:145:ALA:CB	2.26	0.48
3:BM:79:VAL:O	3:BM:79:VAL:HG13	2.14	0.48
3:BM:180:PHE:O	3:BM:184:ASP:OD1	2.31	0.48
5:BI:39:VAL:C	5:BI:41:SER:H	2.22	0.48
5:BK:26:ALA:CA	5:BK:29:ILE:HG22	2.43	0.48
5:BK:46:TRP:CZ3	9:BK:102:BCL:HBC3	2.48	0.48
5:BO:36:HIS:NE2	9:BP:101:BCL:CMD	2.77	0.48
6:BP:33:VAL:O	6:BP:37:LEU:HG	2.13	0.48
5:BQ:30:VAL:HG13	5:BQ:31:LEU:N	2.28	0.48
5:BQ:46:TRP:CH2	9:BQ:103:BCL:H2C	2.49	0.48
6:BV:10:THR:HB	6:BV:13:GLU:OE2	2.14	0.48
5:B1:29:ILE:HG23	5:B1:30:VAL:N	2.28	0.48
9:B3:102:BCL:HMD1	6:B4:36:HIS:CE1	2.48	0.48
9:B7:103:BCL:H12	9:B8:101:BCL:HED1	1.96	0.48
5:B9:49:ASP:CG	5:B9:50:ASN:OD1	2.56	0.48
1:AC:210:ILE:O	1:AC:210:ILE:HG22	2.12	0.48
1:AC:327:TYR:HB2	1:AC:330:LEU:HD12	1.95	0.48
2:AL:35:PHE:CE1	2:AL:111:LEU:HD12	2.49	0.48
2:AL:93:GLY:O	2:AL:94:LEU:C	2.57	0.48
2:AL:167:SER:CA	9:AL:301:BCL:HBC1	2.44	0.48
2:AL:172:GLN:HA	2:AL:172:GLN:NE2	2.28	0.48
3:AM:23:LEU:HD22	3:AM:139:ALA:O	2.13	0.48
3:AM:152:ALA:O	3:AM:155:PHE:HB3	2.13	0.48
3:AM:180:PHE:O	3:AM:181:PRO:C	2.57	0.48
3:AM:247:ARG:HH11	3:AM:247:ARG:CG	2.27	0.48
3:AM:260:VAL:HG11	4:AH:34:ASP:OD1	2.14	0.48
3:AM:269:ALA:O	3:AM:271:TRP:N	2.47	0.48
4:AH:63:ASP:HB3	4:AH:64:PRO:CD	2.44	0.48
6:AB:17:PHE:HE1	14:AB:102:CRT:C9	2.20	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:AB:43:ARG:HB3	5:AD:55:TYR:OH	2.14	0.48
5:AF:12:TRP:HB2	6:AG:14:ALA:HB1	1.94	0.48
9:AN:101:BCL:C1B	9:AO:102:BCL:HMB3	2.44	0.48
14:AP:102:CRT:H342	9:AQ:102:BCL:CBA	2.15	0.48
9:A1:102:BCL:HMB3	9:A1:102:BCL:CBB	2.44	0.48
6:A8:20:ILE:CG2	6:A8:21:PHE:N	2.77	0.48
6:A8:28:TRP:C	6:A8:30:GLY:N	2.70	0.48
6:A8:38:LEU:HA	6:A8:41:LEU:HD12	1.96	0.48
9:A0:102:BCL:C1	9:A0:102:BCL:CGA	2.87	0.48
1:BC:267:THR:C	1:BC:269:ALA:N	2.72	0.48
2:BL:221:GLU:C	2:BL:223:THR:N	2.72	0.48
4:BH:156:VAL:HG12	4:BH:157:VAL:N	2.29	0.48
6:BR:29:PHE:N	6:BR:29:PHE:CD1	2.76	0.48
5:BS:43:ASP:CA	5:BU:56:GLN:HG3	2.43	0.48
6:BT:45:TRP:CD1	6:BT:46:LEU:H	2.31	0.48
5:BU:29:ILE:CA	9:BU:102:BCL:H43	2.44	0.48
5:BY:54:SER:O	5:BY:55:TYR:C	2.56	0.48
9:B2:101:BCL:C1B	9:B3:102:BCL:C2B	2.92	0.48
14:B0:101:CRT:H32	9:B0:102:BCL:HMA2	1.96	0.48
1:AC:315:ASN:OD1	1:AC:316:LYS:HE2	2.14	0.48
2:AL:29:PRO:HG2	3:AM:257:GLY:HA2	1.94	0.48
3:AM:156:PHE:CZ	3:AM:280:ALA:HB1	2.49	0.48
3:AM:172:ALA:C	3:AM:174:ALA:N	2.72	0.48
4:AH:215:LYS:N	4:AH:218:HIS:HD2	2.10	0.48
6:AB:20:ILE:CD1	14:AB:102:CRT:H81	2.44	0.48
9:AI:102:BCL:ND	9:AJ:101:BCL:CMD	2.76	0.48
9:AO:102:BCL:C3D	9:AP:101:BCL:C2D	2.92	0.48
5:AW:10:LYS:HD2	6:AZ:20:ILE:CD1	2.34	0.48
5:A1:5:ASN:CB	5:A1:8:LEU:HD13	2.44	0.48
6:A6:29:PHE:HA	6:A6:32:VAL:HG12	1.96	0.48
9:A6:101:BCL:H43	14:A7:102:CRT:C26	2.44	0.48
6:A0:17:PHE:CD1	6:A0:18:HIS:HA	2.48	0.48
9:A0:102:BCL:OBB	9:A0:102:BCL:HHC	2.13	0.48
1:BC:137:ALA:HA	1:BC:141:TRP:HD1	1.78	0.48
1:BC:213:THR:OG1	1:BC:257:ASN:HB2	2.13	0.48
2:BL:72:ARG:CG	3:BM:305:PRO:HA	2.44	0.48
2:BL:226:ARG:O	3:BM:51:ILE:HA	2.14	0.48
3:BM:122:LEU:O	3:BM:157:TYR:OH	2.27	0.48
4:BH:30:LEU:O	4:BH:31:ARG:C	2.57	0.48
4:BH:47:GLU:HG3	5:BA:19:ARG:CA	2.44	0.48
5:BA:33:LEU:O	5:BA:37:MET:HB2	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:BF:9:TYR:CE1	6:BG:15:LYS:CG	2.97	0.48
14:BG:102:CRT:C34	9:BI:102:BCL:HBA1	2.20	0.48
5:BI:35:ILE:C	5:BI:37:MET:H	2.21	0.48
5:BK:12:TRP:HD1	6:BN:14:ALA:O	1.97	0.48
5:BO:40:LEU:HD23	5:BO:40:LEU:C	2.39	0.48
5:BO:44:LEU:HD12	5:BO:44:LEU:C	2.39	0.48
9:BQ:104:BCL:H3A	9:BQ:104:BCL:HBA1	1.36	0.48
6:BZ:21:PHE:CD1	6:BZ:21:PHE:C	2.91	0.48
6:BZ:21:PHE:HE1	6:BZ:25:MET:HE3	1.78	0.48
6:B8:46:LEU:HD22	6:B0:42:TYR:HE2	1.79	0.48
1:AC:225:SER:HB3	1:AC:228:GLN:HG3	1.95	0.47
1:AC:266:ARG:O	1:AC:267:THR:C	2.57	0.47
2:AL:276:LEU:C	2:AL:278:LEU:H	2.22	0.47
3:AM:52:TYR:CE2	3:AM:136:ARG:NE	2.82	0.47
3:AM:79:VAL:O	3:AM:80:HIS:C	2.58	0.47
3:AM:218:MET:HE2	3:AM:252:TRP:CZ3	2.49	0.47
5:AA:10:LYS:C	5:AA:13:LEU:HD13	2.39	0.47
6:AB:46:LEU:O	5:AD:52:PRO:HD2	2.14	0.47
5:AD:12:TRP:CE3	5:AD:12:TRP:CA	2.97	0.47
5:AD:45:ASN:HB3	5:AD:49:ASP:HB3	1.97	0.47
5:AF:49:ASP:HB2	5:AI:56:GLN:CB	2.43	0.47
9:AF:102:BCL:HHC	9:AF:102:BCL:OBB	2.13	0.47
5:AI:26:ALA:O	5:AI:29:ILE:CG2	2.60	0.47
9:AK:102:BCL:C2D	9:AN:101:BCL:CMD	2.92	0.47
6:AP:20:ILE:HG21	14:AP:102:CRT:C6	2.44	0.47
14:AS:104:CRT:H36	5:AW:33:LEU:CA	2.33	0.47
5:AU:9:TYR:CE1	6:AV:14:ALA:HB3	2.49	0.47
5:AU:46:TRP:CE2	5:AU:47:LEU:HD22	2.49	0.47
5:AW:40:LEU:HD12	5:AW:45:ASN:HA	1.95	0.47
5:A1:34:LEU:O	5:A1:38:ILE:HG22	2.14	0.47
5:A7:42:THR:HB	5:A9:48:ASP:CG	2.39	0.47
1:BC:126:VAL:O	1:BC:127:SER:C	2.55	0.47
1:BC:137:ALA:C	1:BC:139:SER:H	2.22	0.47
1:BC:205:ASP:HB2	1:BC:304:ARG:NE	2.29	0.47
1:BC:268:THR:HG22	7:BC:504:HEM:HAA1	1.95	0.47
2:BL:48:LEU:HD13	5:BA:33:LEU:HD23	1.96	0.47
3:BM:77:ALA:HA	3:BM:81:TRP:CD1	2.49	0.47
3:BM:317:TYR:HD1	3:BM:317:TYR:H	1.59	0.47
4:BH:94:PRO:CG	6:B0:8:GLY:CA	2.92	0.47
6:BB:18:HIS:ND1	6:BB:22:MET:HB2	2.27	0.47
6:BB:18:HIS:HE1	6:BB:22:MET:SD	2.37	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:BG:36:HIS:HE1	9:BG:101:BCL:C4A	2.26	0.47
5:BI:14:ILE:CG2	5:BK:18:ARG:HB3	2.44	0.47
6:BJ:17:PHE:CD1	6:BJ:17:PHE:C	2.92	0.47
5:BK:13:LEU:HD21	6:BN:10:THR:O	2.14	0.47
5:B5:10:LYS:HG2	14:B5:103:CRT:C1M	2.44	0.47
5:B7:46:TRP:NE1	5:B7:47:LEU:CD2	2.77	0.47
5:B9:9:TYR:HA	6:B0:18:HIS:CD2	2.49	0.47
1:AC:53:ILE:C	1:AC:55:ALA:N	2.71	0.47
1:AC:172:PRO:O	1:AC:173:LYS:O	2.33	0.47
1:AC:273:ILE:O	1:AC:273:ILE:HG22	2.14	0.47
2:AL:82:TYR:HB3	2:AL:85:ARG:NE	2.28	0.47
2:AL:87:ALA:O	2:AL:93:GLY:HA3	2.14	0.47
2:AL:144:ARG:HB3	2:AL:145:PRO:CD	2.42	0.47
3:AM:115:TRP:NE1	3:AM:177:PHE:HD2	2.12	0.47
3:AM:218:MET:O	3:AM:221:ALA:N	2.47	0.47
4:AH:124:ASP:CB	4:AH:233:LEU:HD21	2.37	0.47
4:AH:167:VAL:HA	4:AH:183:GLU:O	2.14	0.47
5:AF:29:ILE:O	5:AF:33:LEU:HD13	2.13	0.47
5:AF:35:ILE:O	5:AF:36:HIS:C	2.58	0.47
9:AF:102:BCL:HBC2	9:AF:102:BCL:CHD	2.44	0.47
9:AI:102:BCL:H2	6:AJ:28:TRP:CZ2	2.49	0.47
5:AK:8:LEU:O	5:AK:11:ILE:HG13	2.14	0.47
5:AK:49:ASP:HB2	5:AO:56:GLN:HG2	1.96	0.47
9:AN:101:BCL:CMB	9:AO:102:BCL:CHB	2.93	0.47
9:AP:101:BCL:HMA1	9:AQ:102:BCL:HMA1	1.96	0.47
5:AS:46:TRP:CZ2	9:AS:103:BCL:HHC	2.49	0.47
9:AS:103:BCL:HHC	9:AS:103:BCL:OBB	2.13	0.47
6:AT:10:THR:HB	6:AT:13:GLU:OE2	2.14	0.47
6:AT:29:PHE:HA	6:AT:32:VAL:HG12	1.96	0.47
9:AV:102:BCL:H203	6:AX:39:ALA:CB	2.44	0.47
9:A5:102:BCL:HED1	6:A6:31:LEU:HB3	1.95	0.47
1:BC:105:GLU:OE2	1:BC:105:GLU:N	2.47	0.47
1:BC:212:ILE:HG21	1:BC:229:ALA:HB2	1.96	0.47
1:BC:269:ALA:HB2	7:BC:504:HEM:HMA2	1.96	0.47
2:BL:88:PRO:O	2:BL:89:LEU:C	2.56	0.47
3:BM:90:PHE:C	3:BM:92:TRP:N	2.72	0.47
3:BM:120:LEU:HB2	14:BM:406:CRT:H372	1.95	0.47
3:BM:166:VAL:HG22	3:BM:171:TRP:CH2	2.49	0.47
3:BM:176:PRO:HD3	3:BM:185:TRP:HD1	1.78	0.47
3:BM:203:MET:HA	3:BM:206:ILE:HD12	1.96	0.47
3:BM:286:LEU:O	3:BM:290:VAL:HB	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:BH:65:LYS:O	4:BH:77:VAL:CA	2.63	0.47
5:BF:14:ILE:HD13	6:BJ:17:PHE:HE2	1.79	0.47
5:BI:20:VAL:HA	5:BI:23:SER:OG	2.15	0.47
6:BN:21:PHE:HA	14:BN:102:CRT:C11	2.44	0.47
6:BP:45:TRP:O	6:BP:45:TRP:CD1	2.67	0.47
9:BP:101:BCL:HMB3	9:BQ:103:BCL:CHB	2.44	0.47
5:BU:46:TRP:CE2	5:BU:47:LEU:HD22	2.49	0.47
6:B2:43:ARG:HD3	5:B3:55:TYR:CG	2.49	0.47
1:AC:41:GLU:OE1	2:AL:153:HIS:NE2	2.48	0.47
1:AC:133:LEU:HD13	1:AC:283:TYR:CD2	2.49	0.47
1:AC:157:ARG:NH1	1:AC:318:LEU:CD2	2.75	0.47
1:AC:245:VAL:O	1:AC:245:VAL:HG23	2.14	0.47
2:AL:127:PRO:O	2:AL:128:PHE:C	2.56	0.47
2:AL:184:LEU:O	2:AL:187:SER:N	2.47	0.47
2:AL:192:ASN:C	2:AL:194:LEU:H	2.23	0.47
2:AL:207:THR:HB	3:AM:238:ILE:HG21	1.95	0.47
3:AM:156:PHE:CA	3:AM:159:VAL:HG23	2.44	0.47
3:AM:301:HIS:CE1	4:AH:8:TYR:HB3	2.48	0.47
4:AH:87:VAL:O	4:AH:89:ALA:N	2.46	0.47
4:AH:108:LEU:O	4:AH:110:GLY:N	2.46	0.47
4:AH:121:LYS:HZ1	4:BH:73:GLY:HA2	1.78	0.47
4:AH:215:LYS:HB2	4:AH:218:HIS:CD2	2.49	0.47
5:AF:12:TRP:CE3	5:AF:12:TRP:HA	2.50	0.47
5:AI:23:SER:OG	5:AI:24:ILE:N	2.47	0.47
5:AK:48:ASP:CB	5:AK:56:GLN:NE2	2.77	0.47
9:AK:102:BCL:C3D	9:AN:101:BCL:C3D	2.92	0.47
5:AO:9:TYR:HB3	6:AP:18:HIS:CD2	2.48	0.47
9:AP:101:BCL:H2A	9:AP:101:BCL:CGD	2.44	0.47
14:AP:102:CRT:H342	9:AQ:102:BCL:H3A	1.97	0.47
5:AQ:43:ASP:HB2	5:AS:47:LEU:CD1	2.38	0.47
9:AS:103:BCL:HBD	9:AT:101:BCL:OBD	2.13	0.47
5:A1:10:LYS:HD2	6:A4:20:ILE:HG21	1.97	0.47
5:A7:26:ALA:O	5:A7:30:VAL:HG12	2.14	0.47
6:A8:18:HIS:C	6:A8:18:HIS:CD2	2.92	0.47
6:A8:45:TRP:O	6:A8:46:LEU:CB	2.62	0.47
1:BC:254:ARG:HD3	1:BC:255:ALA:N	2.29	0.47
1:BC:270:TRP:HD1	7:BC:503:HEM:HBD2	1.79	0.47
2:BL:13:ARG:HA	4:BH:99:PRO:CB	2.44	0.47
2:BL:148:MET:CB	2:BL:153:HIS:ND1	2.74	0.47
5:BA:22:VAL:C	5:BA:24:ILE:N	2.72	0.47
6:BB:36:HIS:CE1	9:BB:101:BCL:ND	2.81	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:BB:101:BCL:CBB	9:BB:101:BCL:HMB1	2.45	0.47
5:BF:24:ILE:C	5:BF:26:ALA:N	2.72	0.47
6:BG:30:GLY:CA	6:BG:33:VAL:HG12	2.44	0.47
9:BI:102:BCL:CBC	9:BJ:101:BCL:HHD	2.39	0.47
6:BJ:18:HIS:NE2	6:BJ:22:MET:CE	2.77	0.47
6:BT:9:LEU:CD2	6:BT:13:GLU:HG3	2.30	0.47
6:BT:10:THR:C	6:BT:13:GLU:OE2	2.56	0.47
5:BW:17:PRO:HA	5:BW:20:VAL:HG22	1.96	0.47
14:BW:103:CRT:H6	6:BZ:17:PHE:CD1	2.48	0.47
6:BZ:22:MET:HG3	6:BZ:26:TYR:HE1	1.78	0.47
6:BZ:27:ALA:O	6:BZ:31:LEU:HG	2.15	0.47
6:B2:20:ILE:CG1	14:B2:102:CRT:H83	2.45	0.47
6:B2:42:TYR:HD1	6:B2:43:ARG:HG3	1.79	0.47
5:B3:2:PHE:O	5:B3:3:THR:O	2.32	0.47
9:B8:101:BCL:HMC1	9:B9:102:BCL:CAB	2.42	0.47
1:AC:213:THR:HB	1:AC:257:ASN:OD1	2.15	0.47
2:AL:164:ASP:O	2:AL:165:TRP:C	2.56	0.47
3:AM:98:PRO:HB2	3:AM:171:TRP:CB	2.44	0.47
3:AM:221:ALA:O	3:AM:224:LEU:HB2	2.15	0.47
4:AH:19:PHE:C	4:AH:21:LEU:H	2.21	0.47
4:AH:35:LYS:NZ	4:AH:57:GLY:CA	2.72	0.47
4:AH:58:PHE:N	4:AH:59:PRO:HD2	2.29	0.47
4:AH:123:CYS:SG	4:AH:231:VAL:O	2.66	0.47
4:AH:136:MET:HG3	4:AH:170:VAL:O	2.14	0.47
5:AA:33:LEU:O	5:AA:37:MET:HB2	2.14	0.47
5:AA:42:THR:HG22	5:AD:48:ASP:OD2	2.14	0.47
6:AB:44:PRO:HD2	5:AD:55:TYR:CE2	2.49	0.47
9:AQ:102:BCL:C2D	9:AR:101:BCL:C2D	2.93	0.47
9:AT:101:BCL:OBB	9:AT:101:BCL:HHC	2.13	0.47
5:AU:9:TYR:CB	6:AV:15:LYS:HD3	2.45	0.47
5:AW:21:LEU:HD11	9:AW:101:BCL:H141	1.96	0.47
5:AW:30:VAL:HA	5:AW:33:LEU:HG	1.96	0.47
6:AX:34:ILE:HG23	6:AX:35:ALA:N	2.29	0.47
5:A9:43:ASP:CG	5:A9:44:LEU:HD12	2.38	0.47
6:A0:29:PHE:HD1	6:A0:29:PHE:H	1.61	0.47
1:BC:139:SER:O	1:BC:142:LYS:HG3	2.14	0.47
1:BC:273:ILE:HG22	1:BC:273:ILE:O	2.14	0.47
1:BC:285:TRP:HB3	1:BC:286:PRO:HD3	1.95	0.47
2:BL:9:LYS:HA	4:BH:111:PHE:HE1	1.79	0.47
2:BL:18:ILE:HG23	4:BH:259:LEU:CB	2.45	0.47
3:BM:51:ILE:HG12	3:BM:52:TYR:N	2.29	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:BM:84:PHE:HD1	3:BM:84:PHE:N	2.12	0.47
3:BM:154:ILE:O	3:BM:154:ILE:CG2	2.62	0.47
3:BM:279:THR:HG22	3:BM:282:ILE:HD12	1.96	0.47
10:BM:403:BPH:HBA2	15:BQ:101:PEF:H431	1.97	0.47
4:BH:136:MET:SD	4:BH:170:VAL:HG23	2.54	0.47
4:BH:145:ALA:O	4:BH:147:GLY:N	2.48	0.47
5:BA:11:ILE:C	5:BA:13:LEU:H	2.20	0.47
5:BD:27:PHE:CE2	5:BF:29:ILE:HD11	2.48	0.47
9:BF:102:BCL:C4D	9:BG:101:BCL:HMD3	2.44	0.47
5:BI:27:PHE:HA	5:BI:30:VAL:HG12	1.96	0.47
6:BJ:34:ILE:O	6:BJ:38:LEU:HB2	2.13	0.47
14:BP:102:CRT:C2M	5:BQ:37:MET:HG2	2.43	0.47
5:BU:44:LEU:HD12	5:BU:44:LEU:O	2.13	0.47
5:BY:8:LEU:CA	6:B2:20:ILE:HD11	2.28	0.47
6:B2:42:TYR:CE1	6:B2:43:ARG:HG3	2.50	0.47
5:B5:10:LYS:HG2	14:B5:103:CRT:H1M1	1.96	0.47
5:B5:18:ARG:HB3	5:B5:19:ARG:NH1	2.30	0.47
5:B9:33:LEU:H	5:B9:33:LEU:CD1	2.28	0.47
1:AC:110:CYS:HA	1:AC:123:THR:HG1	1.79	0.47
1:AC:195:LEU:HB3	1:AC:196:PRO:HD2	1.95	0.47
1:AC:196:PRO:O	1:AC:197:PHE:CG	2.68	0.47
1:AC:317:PRO:HD2	7:AC:504:HEM:C2D	2.48	0.47
2:AL:230:GLY:O	3:AM:49:GLY:HA2	2.14	0.47
3:AM:26:GLY:HA2	5:AO:16:ASP:OD2	2.14	0.47
3:AM:64:GLY:C	3:AM:66:VAL:H	2.22	0.47
3:AM:84:PHE:CZ	5:AU:38:ILE:HD12	2.50	0.47
3:AM:184:ASP:O	3:AM:187:ALA:HB3	2.14	0.47
3:AM:265:ILE:C	3:AM:267:ARG:H	2.21	0.47
4:AH:65:LYS:O	4:AH:77:VAL:CA	2.61	0.47
4:AH:170:VAL:HA	4:AH:182:LEU:HA	1.96	0.47
5:AA:17:PRO:HD2	5:AA:18:ARG:NH1	2.28	0.47
5:AD:43:ASP:OD2	5:AD:44:LEU:HG	2.14	0.47
5:AF:12:TRP:NE1	6:AG:17:PHE:CD1	2.82	0.47
6:AG:19:ALA:O	6:AG:23:GLN:HG3	2.13	0.47
9:AG:101:BCL:C1C	9:AI:102:BCL:HBB3	2.45	0.47
9:AK:102:BCL:CHD	9:AN:101:BCL:HMD1	2.43	0.47
5:AO:17:PRO:O	5:AO:21:LEU:N	2.44	0.47
9:AQ:102:BCL:HMB1	9:AQ:102:BCL:HBB3	1.96	0.47
6:AT:40:TRP:CE3	6:AT:44:PRO:HA	2.50	0.47
6:AV:44:PRO:CG	5:AW:52:PRO:HG2	2.45	0.47
14:AW:102:CRT:H342	9:A1:102:BCL:HBA2	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:A3:46:TRP:CD1	5:A3:47:LEU:N	2.83	0.47
9:A5:102:BCL:H162	9:A5:102:BCL:H111	1.95	0.47
1:BC:275:HIS:O	1:BC:275:HIS:CD2	2.68	0.47
1:BC:313:ALA:O	1:BC:314:VAL:HG22	2.14	0.47
2:BL:116:ILE:HD13	3:BM:254:TRP:HB2	1.97	0.47
2:BL:186:ILE:HD13	9:BM:401:BCL:HMD1	1.96	0.47
3:BM:196:LEU:C	3:BM:198:TYR:H	2.22	0.47
3:BM:204:LEU:CD1	3:BM:279:THR:HG21	2.38	0.47
3:BM:262:MET:HG3	3:BM:262:MET:O	2.14	0.47
5:BF:29:ILE:HG23	5:BF:30:VAL:N	2.28	0.47
5:BK:2:PHE:O	5:BK:5:ASN:HB3	2.14	0.47
6:BT:29:PHE:HA	6:BT:32:VAL:HG12	1.97	0.47
9:BW:102:BCL:H2A	9:BW:102:BCL:O1D	2.14	0.47
5:B1:10:LYS:CB	14:B1:103:CRT:H83	2.43	0.47
6:B2:17:PHE:CD1	14:B2:102:CRT:C6	2.87	0.47
1:AC:144:HIS:CE1	7:AC:504:HEM:NC	2.83	0.47
2:AL:6:PHE:HE2	3:AM:246:GLU:HA	1.79	0.47
2:AL:38:VAL:HG23	2:AL:108:SER:OG	2.15	0.47
2:AL:67:THR:OG1	2:AL:68:TYR:N	2.48	0.47
2:AL:180:PRO:HA	2:AL:183:MET:SD	2.54	0.47
3:AM:92:TRP:O	3:AM:93:LEU:O	2.33	0.47
3:AM:134:TYR:C	3:AM:144:GLN:HE22	2.22	0.47
3:AM:208:PHE:O	3:AM:210:TYR:N	2.47	0.47
3:AM:226:VAL:HG21	3:AM:244:ALA:HA	1.97	0.47
3:AM:276:THR:HG22	3:AM:277:VAL:H	1.74	0.47
4:AH:171:TRP:N	4:AH:181:TYR:O	2.47	0.47
5:AA:33:LEU:N	5:AA:33:LEU:HD12	2.30	0.47
9:AB:101:BCL:HHC	9:AB:101:BCL:OBB	2.14	0.47
5:AF:36:HIS:NE2	9:AG:101:BCL:CMD	2.78	0.47
5:AO:33:LEU:O	5:AO:37:MET:HG2	2.15	0.47
6:AP:13:GLU:HA	6:AP:16:GLU:HG3	1.96	0.47
6:AP:13:GLU:HA	6:AP:16:GLU:OE1	2.14	0.47
6:AP:45:TRP:O	6:AP:46:LEU:CD2	2.63	0.47
9:AS:103:BCL:HMB1	9:AS:103:BCL:CBB	2.45	0.47
5:AU:17:PRO:HG2	5:AU:18:ARG:HD2	1.96	0.47
5:AW:19:ARG:NH1	5:AY:22:VAL:CG2	2.77	0.47
5:A1:18:ARG:O	5:A1:22:VAL:HG12	2.15	0.47
5:A1:19:ARG:O	5:A1:23:SER:CB	2.62	0.47
5:A1:28:GLN:HE22	6:A2:28:TRP:NE1	2.13	0.47
5:A3:18:ARG:O	5:A3:22:VAL:HG12	2.14	0.47
5:A5:39:VAL:C	5:A5:41:SER:N	2.72	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:A8:17:PHE:CE1	6:A8:20:ILE:HG21	2.50	0.47
9:A9:102:BCL:H92	6:A0:28:TRP:HB2	1.97	0.47
1:BC:36:ARG:NH2	2:BL:90:THR:O	2.47	0.47
1:BC:53:ILE:CG1	1:BC:319:TYR:CZ	2.97	0.47
2:BL:13:ARG:HG3	2:BL:13:ARG:HH11	1.79	0.47
2:BL:168:ASN:O	2:BL:170:GLY:N	2.48	0.47
2:BL:184:LEU:HB2	2:BL:252:TRP:CD1	2.49	0.47
2:BL:268:TRP:C	2:BL:270:GLU:N	2.73	0.47
4:BH:23:PHE:O	4:BH:25:GLY:N	2.48	0.47
4:BH:54:LYS:HG3	4:BH:58:PHE:HA	1.96	0.47
4:BH:112:GLY:N	4:BH:115:ALA:HB2	2.29	0.47
4:BH:135:PRO:C	4:BH:137:ARG:H	2.21	0.47
6:BN:7:THR:OG1	6:BN:8:GLY:N	2.46	0.47
6:BN:44:PRO:HD2	5:BO:55:TYR:HH	1.77	0.47
9:BO:102:BCL:CBC	9:BP:101:BCL:HHD	2.39	0.47
5:BU:11:ILE:C	14:BU:103:CRT:H21A	2.40	0.47
14:BU:103:CRT:O2	5:BY:33:LEU:O	2.32	0.47
6:BV:20:ILE:HD13	6:BV:20:ILE:C	2.40	0.47
5:BY:26:ALA:O	5:BY:30:VAL:HG23	2.14	0.47
5:B5:20:VAL:HA	5:B5:23:SER:CB	2.43	0.47
6:B6:40:TRP:HZ3	6:B6:45:TRP:N	2.11	0.47
5:B7:30:VAL:HG13	5:B7:31:LEU:N	2.30	0.47
1:AC:263:THR:O	1:AC:264:PRO:C	2.58	0.47
1:AC:280:ASN:OD1	1:AC:305:VAL:N	2.45	0.47
1:AC:291:LEU:HD22	1:AC:295:ARG:HB2	1.96	0.47
2:AL:139:VAL:HG22	2:AL:258:LEU:HB2	1.97	0.47
2:AL:164:ASP:OD2	3:AM:307:TYR:OH	2.33	0.47
2:AL:196:LEU:C	2:AL:198:MET:N	2.73	0.47
2:AL:226:ARG:NH2	3:AM:47:GLN:HB3	2.30	0.47
3:AM:12:GLN:C	4:AH:145:ALA:HB2	2.40	0.47
3:AM:51:ILE:HG12	3:AM:52:TYR:O	2.15	0.47
3:AM:76:LEU:CD2	5:AU:37:MET:HE3	2.44	0.47
3:AM:195:ASN:O	3:AM:197:TYR:N	2.48	0.47
3:AM:240:HIS:CE1	4:AH:69:LEU:HD11	2.38	0.47
3:AM:246:GLU:O	3:AM:250:LEU:HB2	2.14	0.47
3:AM:284:ILE:HD11	9:AM:402:BCL:OBD	2.15	0.47
4:AH:5:ILE:CG2	4:AH:6:THR:H	2.10	0.47
4:AH:151:PRO:O	4:AH:167:VAL:HG21	2.15	0.47
4:AH:218:HIS:O	4:AH:222:VAL:HG23	2.15	0.47
5:AA:36:HIS:NE2	9:AB:101:BCL:CMD	2.78	0.47
9:AA:101:BCL:HAC2	9:AB:101:BCL:CBC	2.45	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
14:AA:102:CRT:H372	5:AD:35:ILE:HD11	1.97	0.47
5:AD:9:TYR:CD1	5:AD:9:TYR:C	2.93	0.47
5:AF:8:LEU:N	6:AJ:20:ILE:HD11	2.29	0.47
5:AF:26:ALA:O	5:AF:27:PHE:C	2.58	0.47
5:AI:15:LEU:HD12	5:AI:20:VAL:HG11	1.95	0.47
14:AN:102:CRT:H342	9:AO:102:BCL:CBA	2.36	0.47
6:AP:28:TRP:O	6:AP:31:LEU:N	2.48	0.47
6:AR:16:GLU:CB	14:AR:102:CRT:H23	2.45	0.47
5:AS:26:ALA:O	5:AS:30:VAL:HG12	2.14	0.47
5:AS:49:ASP:CG	5:AS:50:ASN:N	2.73	0.47
6:AT:38:LEU:HD23	6:AT:38:LEU:O	2.14	0.47
5:AU:19:ARG:HE	5:AW:18:ARG:HH22	1.63	0.47
5:AU:50:ASN:HB2	5:AW:59:GLY:HA3	1.97	0.47
9:AU:102:BCL:OBD	6:AV:32:VAL:HG22	2.14	0.47
6:AV:14:ALA:O	6:AV:18:HIS:HB2	2.14	0.47
6:AX:33:VAL:CG2	6:AX:37:LEU:HD23	2.42	0.47
6:AX:45:TRP:CZ3	9:AX:101:BCL:HAC2	2.49	0.47
5:AY:44:LEU:CD1	6:AZ:43:ARG:HD2	2.43	0.47
6:AZ:36:HIS:HE1	9:AZ:101:BCL:C4A	2.27	0.47
5:A5:9:TYR:CD1	5:A5:9:TYR:C	2.92	0.47
5:A5:12:TRP:CH2	5:A5:20:VAL:HG23	2.49	0.47
5:A5:44:LEU:C	5:A5:46:TRP:N	2.72	0.47
6:A6:38:LEU:HD23	6:A6:38:LEU:O	2.15	0.47
1:BC:29:GLY:N	1:BC:44:TYR:O	2.47	0.47
1:BC:40:MET:HB3	1:BC:248:THR:HB	1.97	0.47
1:BC:196:PRO:O	1:BC:197:PHE:CG	2.67	0.47
1:BC:265:LYS:O	1:BC:266:ARG:C	2.56	0.47
1:BC:282:ASN:HB3	1:BC:283:TYR:CD1	2.50	0.47
1:BC:283:TYR:N	1:BC:283:TYR:HD1	2.13	0.47
1:BC:285:TRP:HZ3	1:BC:302:PRO:HD3	1.74	0.47
2:BL:7:GLU:CD	2:BL:11:ARG:HH21	2.22	0.47
2:BL:129:ALA:CB	2:BL:247:LEU:HD21	2.42	0.47
2:BL:178:TYR:CE1	3:BM:180:PHE:CD2	3.03	0.47
2:BL:181:ALA:C	2:BL:183:MET:N	2.73	0.47
11:BL:304:UQ8:H42	11:BL:304:UQ8:H45	1.68	0.47
3:BM:222:THR:O	3:BM:226:VAL:HG12	2.13	0.47
3:BM:271:TRP:CE2	4:BH:26:LEU:HD11	2.50	0.47
4:BH:196:PRO:O	4:BH:197:ILE:C	2.57	0.47
9:BA:101:BCL:HBA1	14:B0:101:CRT:C34	2.31	0.47
6:BB:18:HIS:HD1	6:BB:18:HIS:C	2.21	0.47
5:BD:9:TYR:C	5:BD:9:TYR:CD1	2.92	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:BD:44:LEU:HD22	5:BF:55:TYR:CZ	2.50	0.47
5:BF:26:ALA:HA	5:BF:29:ILE:HG22	1.95	0.47
9:BK:102:BCL:HMB1	9:BK:102:BCL:CBB	2.44	0.47
6:BN:29:PHE:CE2	9:BN:101:BCL:H2	2.50	0.47
6:BP:15:LYS:O	6:BP:16:GLU:C	2.58	0.47
5:BQ:12:TRP:HA	5:BQ:12:TRP:HE3	1.80	0.47
6:BV:33:VAL:CG1	6:BV:34:ILE:N	2.77	0.47
5:BW:16:ASP:HB3	5:BW:19:ARG:HG2	1.96	0.47
6:BX:30:GLY:O	6:BX:34:ILE:HG22	2.14	0.47
6:BX:38:LEU:C	6:BX:38:LEU:CD2	2.88	0.47
9:BY:102:BCL:HMB1	9:BY:102:BCL:HBB3	1.96	0.47
5:B1:44:LEU:HD23	5:B1:44:LEU:N	2.29	0.47
5:B3:36:HIS:CE1	9:B4:101:BCL:CMD	2.96	0.47
9:B3:102:BCL:H2	6:B4:28:TRP:CZ2	2.50	0.47
5:B5:33:LEU:C	5:B5:33:LEU:HD12	2.39	0.47
9:B5:102:BCL:H13	9:B5:102:BCL:H192	1.97	0.47
9:B5:102:BCL:HBC2	9:B6:101:BCL:HMD1	1.97	0.47
6:B6:21:PHE:CD2	14:B7:102:CRT:H14	2.49	0.47
6:B8:17:PHE:CE1	6:B8:20:ILE:HG21	2.50	0.47
9:B9:102:BCL:HAC2	9:B0:102:BCL:CBC	2.45	0.47
6:B0:29:PHE:O	6:B0:32:VAL:HG12	2.14	0.47
6:B0:40:TRP:CE3	6:B0:44:PRO:HA	2.50	0.47
6:B0:40:TRP:HA	6:B0:40:TRP:HE3	1.78	0.47
1:AC:157:ARG:NE	1:AC:312:GLN:OE1	2.47	0.47
2:AL:195:ALA:O	2:AL:198:MET:HB2	2.14	0.47
3:AM:76:LEU:HA	3:AM:86:PHE:CE1	2.50	0.47
3:AM:164:ARG:HA	3:AM:167:MET:CB	2.45	0.47
4:AH:180:ARG:HG2	4:AH:180:ARG:HH11	1.80	0.47
5:AF:11:ILE:HG23	5:AF:12:TRP:CE3	2.50	0.47
5:AI:35:ILE:O	5:AI:36:HIS:C	2.51	0.47
5:AI:45:ASN:O	5:AI:49:ASP:CG	2.58	0.47
6:AJ:33:VAL:HG22	6:AJ:37:LEU:CD2	2.45	0.47
6:AJ:45:TRP:CD1	6:AJ:46:LEU:N	2.83	0.47
5:AK:48:ASP:CB	5:AK:56:GLN:HE22	2.28	0.47
9:AR:101:BCL:HHB	9:AS:103:BCL:HMA1	1.97	0.47
9:AV:102:BCL:H8	9:AV:102:BCL:HMB1	1.96	0.47
6:AZ:29:PHE:N	6:AZ:29:PHE:HD1	2.12	0.47
5:A3:56:GLN:HG2	5:A3:57:ALA:N	2.30	0.47
5:A5:27:PHE:CZ	5:A7:29:ILE:HD11	2.50	0.47
5:A7:9:TYR:C	5:A7:9:TYR:CD1	2.92	0.47
9:A7:103:BCL:CHD	9:A8:101:BCL:HMD1	2.45	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:A0:40:TRP:CE3	6:A0:44:PRO:HA	2.50	0.47
1:BC:35:TYR:CD2	3:BM:308:PRO:HD2	2.50	0.47
1:BC:141:TRP:O	1:BC:145:VAL:HG22	2.14	0.47
1:BC:184:ASN:O	1:BC:185:TYR:HB2	2.15	0.47
1:BC:261:GLN:O	1:BC:262:SER:O	2.33	0.47
2:BL:216:LYS:HD2	2:BL:220:HIS:NE2	2.30	0.47
6:BB:18:HIS:ND1	6:BB:18:HIS:C	2.72	0.47
9:BE:101:BCL:HMC1	9:BF:102:BCL:CBB	2.45	0.47
5:BF:10:LYS:O	5:BF:13:LEU:HG	2.15	0.47
5:BF:29:ILE:CB	9:BF:102:BCL:H43	2.39	0.47
5:BF:35:ILE:CA	5:BF:38:ILE:HG22	2.45	0.47
5:BI:11:ILE:HG23	5:BI:12:TRP:CD1	2.50	0.47
5:BO:46:TRP:CE3	9:BO:102:BCL:H2C	2.50	0.47
5:BQ:12:TRP:HA	5:BQ:12:TRP:CE3	2.50	0.47
5:BW:10:LYS:HB3	14:BW:103:CRT:H23	1.96	0.47
5:BY:9:TYR:CZ	5:BY:10:LYS:HE3	2.49	0.47
2:AL:118:ARG:O	2:AL:119:LYS:C	2.58	0.47
2:AL:173:PHE:C	2:AL:174:LEU:HD12	2.40	0.47
2:AL:181:ALA:O	2:AL:182:HIS:C	2.58	0.47
13:AM:405:MQ8:H493	5:AA:34:LEU:HD13	1.97	0.47
4:AH:27:ILE:CG2	4:AH:28:ILE:N	2.78	0.47
4:AH:60:ASP:O	4:AH:61:LEU:C	2.57	0.47
5:AA:20:VAL:O	5:AA:20:VAL:HG12	2.15	0.47
5:AD:40:LEU:HD13	5:AD:46:TRP:CZ2	2.50	0.47
9:AD:102:BCL:HMB1	9:AD:102:BCL:HBB3	1.95	0.47
5:AO:7:ASN:ND2	5:AO:7:ASN:N	2.60	0.47
9:AO:102:BCL:C4D	9:AP:101:BCL:CMD	2.89	0.47
9:AO:102:BCL:HBB3	9:AO:102:BCL:HMB1	1.97	0.47
6:AP:44:PRO:HD2	5:AQ:55:TYR:OH	2.14	0.47
5:AS:29:ILE:HG23	5:AS:30:VAL:N	2.30	0.47
14:AS:104:CRT:C2M	5:AW:37:MET:CA	2.93	0.47
6:AT:10:THR:C	6:AT:13:GLU:OE2	2.58	0.47
14:AW:102:CRT:C6	6:AZ:20:ILE:HG21	2.45	0.47
9:AX:101:BCL:CGA	9:AX:101:BCL:H12	2.41	0.47
5:A1:28:GLN:HE22	6:A2:28:TRP:CD1	2.32	0.47
6:A4:13:GLU:O	6:A4:16:GLU:CG	2.60	0.47
6:A4:18:HIS:CD2	6:A4:18:HIS:O	2.67	0.47
6:A4:40:TRP:CZ3	6:A4:44:PRO:CA	2.96	0.47
6:A6:31:LEU:HA	6:A6:34:ILE:HG22	1.97	0.47
5:A7:40:LEU:HD12	5:A7:45:ASN:HA	1.97	0.47
5:A7:46:TRP:CZ3	9:A7:103:BCL:H2C	2.49	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:A9:33:LEU:HD12	5:A9:33:LEU:N	2.30	0.47
1:BC:286:PRO:O	1:BC:288:ASN:N	2.46	0.47
2:BL:21:ASP:HB3	5:B7:19:ARG:NE	2.30	0.47
2:BL:188:PHE:CZ	11:BL:304:UQ8:H26A	2.49	0.47
3:BM:114:TRP:CE3	3:BM:117:MET:HG3	2.50	0.47
3:BM:151:ALA:O	3:BM:154:ILE:N	2.48	0.47
3:BM:156:PHE:HA	3:BM:159:VAL:HG23	1.96	0.47
3:BM:241:ARG:O	4:BH:119:ARG:CD	2.63	0.47
4:BH:36:ARG:HE	4:BH:65:LYS:HB2	1.80	0.47
4:BH:36:ARG:NE	4:BH:65:LYS:HB2	2.29	0.47
4:BH:71:HIS:HE1	4:BH:125:LEU:HD22	1.80	0.47
4:BH:138:VAL:O	4:BH:140:LYS:HD3	2.15	0.47
4:BH:146:GLU:H	4:BH:146:GLU:CD	2.22	0.47
5:BA:51:ILE:HD11	6:BB:42:TYR:OH	2.14	0.47
9:BA:101:BCL:HBA2	9:BB:101:BCL:OBD	2.14	0.47
6:BE:29:PHE:CD1	9:BE:101:BCL:H11	2.50	0.47
9:BG:101:BCL:HBB1	9:BI:102:BCL:HMC1	1.97	0.47
6:BP:45:TRP:CE3	9:BP:101:BCL:HBC2	2.50	0.47
9:BW:102:BCL:C3D	9:BX:101:BCL:C3D	2.93	0.47
5:BY:21:LEU:O	5:BY:25:VAL:HG23	2.14	0.47
5:B1:21:LEU:HD11	9:B1:102:BCL:H142	1.96	0.47
9:B1:102:BCL:CAC	9:B2:101:BCL:CBC	2.92	0.47
5:B3:20:VAL:HG23	5:B3:21:LEU:N	2.30	0.47
6:B8:45:TRP:O	6:B8:46:LEU:CB	2.62	0.47
6:B8:45:TRP:HA	5:B9:52:PRO:HD2	1.97	0.47
1:AC:47:ARG:HD3	5:A1:42:THR:CG2	2.46	0.47
3:AM:109:LEU:HD22	5:AQ:42:THR:HG21	1.96	0.47
3:AM:206:ILE:HA	9:AM:402:BCL:CMA	2.43	0.47
3:AM:244:ALA:O	3:AM:246:GLU:N	2.48	0.47
6:AB:8:GLY:C	6:AB:9:LEU:HG	2.40	0.47
6:AB:33:VAL:O	6:AB:37:LEU:HD23	2.15	0.47
5:AF:31:LEU:HD11	14:AG:102:CRT:H372	1.97	0.47
6:AJ:21:PHE:O	6:AJ:22:MET:C	2.58	0.47
5:AK:12:TRP:CD1	6:AN:17:PHE:HD2	2.32	0.47
5:AK:33:LEU:HD12	5:AK:34:LEU:N	2.30	0.47
5:AO:30:VAL:HG13	5:AO:31:LEU:N	2.30	0.47
6:AR:46:LEU:HD22	6:AT:42:TYR:CD2	2.50	0.47
5:AS:36:HIS:HD2	5:AS:46:TRP:HH2	1.63	0.47
6:AT:9:LEU:HB3	6:AT:13:GLU:HG3	1.96	0.47
6:AV:33:VAL:CG1	6:AV:34:ILE:N	2.78	0.47
6:AV:45:TRP:O	6:AV:46:LEU:HB2	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:A3:19:ARG:CG	5:A3:20:VAL:N	2.78	0.47
5:A5:29:ILE:HD13	5:A5:29:ILE:C	2.39	0.47
5:A7:10:LYS:CB	14:A0:101:CRT:C8	2.93	0.47
1:BC:211:ARG:HB3	7:BC:503:HEM:O2A	2.15	0.47
1:BC:254:ARG:HD3	1:BC:254:ARG:C	2.40	0.47
2:BL:207:THR:HB	3:BM:238:ILE:HG21	1.96	0.47
4:BH:189:ASN:HB3	4:BH:191:LYS:CG	2.45	0.47
4:BH:229:ASP:O	4:BH:230:GLN:HB3	2.15	0.47
6:BB:32:VAL:CG2	9:BB:101:BCL:HBA2	2.32	0.47
5:BF:44:LEU:HB3	5:BI:55:TYR:HH	1.79	0.47
14:BF:103:CRT:C34	9:BK:102:BCL:HBA1	2.32	0.47
5:BI:26:ALA:O	5:BI:27:PHE:C	2.56	0.47
9:BI:102:BCL:C2D	9:BJ:101:BCL:C2D	2.93	0.47
5:BQ:43:ASP:HB2	5:BS:47:LEU:HB3	1.97	0.47
9:BS:102:BCL:C1D	9:BT:101:BCL:CMD	2.92	0.47
5:B3:27:PHE:CE2	5:B5:29:ILE:HD12	2.50	0.47
6:B4:10:THR:O	6:B4:14:ALA:HB2	2.14	0.47
5:B7:11:ILE:CD1	5:B7:15:LEU:HD11	2.45	0.47
6:B8:18:HIS:C	6:B8:18:HIS:CD2	2.92	0.47
5:B9:12:TRP:HE1	6:B0:18:HIS:HA	1.80	0.47
1:AC:102:SER:HB2	1:AC:105:GLU:OE2	2.15	0.46
1:AC:264:PRO:O	1:AC:265:LYS:C	2.58	0.46
1:AC:313:ALA:C	1:AC:314:VAL:CG2	2.88	0.46
2:AL:142:PHE:CD1	2:AL:142:PHE:C	2.92	0.46
2:AL:159:ILE:CD1	2:AL:159:ILE:H	2.28	0.46
9:AL:301:BCL:H191	9:AL:303:BCL:H8	1.97	0.46
10:AL:302:BPH:HED1	3:AM:255:THR:CG2	2.45	0.46
3:AM:8:PHE:HB3	3:AM:42:LYS:HG2	1.96	0.46
3:AM:98:PRO:HG3	3:AM:112:GLY:O	2.14	0.46
3:AM:115:TRP:CZ3	3:AM:116:LEU:HD12	2.49	0.46
3:AM:204:LEU:HD23	3:AM:279:THR:HG22	1.97	0.46
3:AM:286:LEU:HA	3:AM:290:VAL:HG21	1.96	0.46
4:AH:35:LYS:O	4:AH:36:ARG:O	2.33	0.46
4:AH:45:ARG:O	4:AH:96:PRO:CB	2.62	0.46
5:AA:17:PRO:HG2	5:AA:18:ARG:CD	2.45	0.46
9:AF:102:BCL:HAC2	9:AG:101:BCL:HBC3	1.94	0.46
14:AJ:102:CRT:C2M	5:AK:36:HIS:HB3	2.45	0.46
5:AQ:28:GLN:O	9:AQ:102:BCL:H12	2.15	0.46
6:AR:44:PRO:O	5:AS:52:PRO:HG3	2.15	0.46
9:AV:102:BCL:H18	9:AW:101:BCL:H3C	1.97	0.46
9:AY:102:BCL:H92	6:AZ:28:TRP:HB2	1.96	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:A1:31:LEU:HD23	9:A2:101:BCL:HED3	1.96	0.46
14:A1:103:CRT:H32	5:A3:31:LEU:HD21	1.96	0.46
6:A2:41:LEU:HD23	6:A2:41:LEU:C	2.39	0.46
5:A3:44:LEU:HD21	9:A3:104:BCL:CBC	2.45	0.46
5:A3:46:TRP:HZ3	9:A3:103:BCL:HBC3	1.79	0.46
6:A4:38:LEU:HD23	6:A4:38:LEU:O	2.15	0.46
1:BC:273:ILE:HG22	1:BC:277:ARG:NH1	2.30	0.46
3:BM:64:GLY:O	3:BM:66:VAL:N	2.46	0.46
3:BM:206:ILE:HG22	3:BM:210:TYR:CE2	2.50	0.46
3:BM:268:TRP:HZ2	4:BH:34:ASP:OD2	1.98	0.46
3:BM:310:VAL:HG12	3:BM:310:VAL:O	2.15	0.46
4:BH:54:LYS:HG3	4:BH:58:PHE:HD1	1.80	0.46
4:BH:141:GLU:CD	4:BH:141:GLU:H	2.23	0.46
4:BH:235:GLU:O	4:BH:239:VAL:HG23	2.15	0.46
5:BA:27:PHE:C	5:BA:30:VAL:HG12	2.40	0.46
5:BA:39:VAL:C	5:BA:41:SER:N	2.69	0.46
6:BB:9:LEU:HD12	6:BB:9:LEU:N	2.30	0.46
5:BF:35:ILE:O	5:BF:36:HIS:C	2.58	0.46
6:BP:17:PHE:CD1	14:BP:102:CRT:H6	2.50	0.46
6:BP:42:TYR:CD2	6:BP:43:ARG:HG3	2.50	0.46
5:B1:12:TRP:CH2	9:B3:102:BCL:H202	2.46	0.46
6:B6:31:LEU:HA	6:B6:34:ILE:HG22	1.97	0.46
5:B7:51:ILE:HD12	5:B7:51:ILE:C	2.40	0.46
5:B9:43:ASP:CG	5:B9:44:LEU:HD12	2.40	0.46
1:AC:298:PRO:C	1:AC:300:GLY:N	2.66	0.46
2:AL:168:ASN:C	2:AL:170:GLY:H	2.23	0.46
3:AM:84:PHE:N	3:AM:84:PHE:CD1	2.83	0.46
3:AM:164:ARG:HB2	3:AM:285:LEU:HD12	1.96	0.46
3:AM:260:VAL:HG12	4:AH:31:ARG:NH2	2.30	0.46
4:AH:184:VAL:HB	4:AH:193:VAL:HG23	1.96	0.46
9:AE:101:BCL:HBB3	9:AE:101:BCL:HMB1	1.98	0.46
5:AF:10:LYS:HD2	14:AJ:102:CRT:H1M1	1.97	0.46
5:AF:35:ILE:CD1	14:AG:102:CRT:H403	2.44	0.46
5:AI:36:HIS:NE2	9:AJ:101:BCL:CMD	2.78	0.46
9:AJ:101:BCL:OBB	9:AJ:101:BCL:HHC	2.14	0.46
9:AK:102:BCL:HBB3	9:AK:102:BCL:HMB1	1.97	0.46
5:AO:52:PRO:C	5:AO:54:SER:N	2.72	0.46
6:AP:17:PHE:CD1	14:AP:102:CRT:H6	2.50	0.46
6:AR:40:TRP:HH2	6:AR:46:LEU:CD1	2.29	0.46
6:AR:46:LEU:CB	6:AT:42:TYR:CZ	2.77	0.46
5:AS:11:ILE:CG1	14:AS:104:CRT:C8	2.92	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:A0:45:TRP:CD1	6:A0:46:LEU:N	2.77	0.46
1:BC:148:THR:HG23	1:BC:322:GLN:HG2	1.96	0.46
1:BC:199:PRO:HG2	1:BC:200:LEU:HD12	1.97	0.46
2:BL:35:PHE:HA	2:BL:38:VAL:HG22	1.97	0.46
2:BL:160:LEU:O	2:BL:160:LEU:HD12	2.14	0.46
2:BL:182:HIS:O	2:BL:186:ILE:HG13	2.15	0.46
2:BL:204:LEU:CD1	3:BM:267:ARG:HG3	2.45	0.46
3:BM:71:ILE:HD13	3:BM:177:PHE:CE1	2.49	0.46
3:BM:265:ILE:HD12	3:BM:265:ILE:HA	1.78	0.46
3:BM:273:ALA:O	3:BM:276:THR:HB	2.15	0.46
4:BH:56:VAL:O	4:BH:56:VAL:HG23	2.15	0.46
4:BH:130:LEU:HD23	4:BH:131:PRO:O	2.16	0.46
5:BA:2:PHE:CA	5:BA:5:ASN:HD21	2.26	0.46
5:BD:16:ASP:OD1	5:BD:17:PRO:HD2	2.14	0.46
9:BD:102:BCL:HBC1	9:BE:101:BCL:HBC3	1.97	0.46
5:BK:5:ASN:ND2	6:BN:22:MET:HE3	2.31	0.46
9:BK:102:BCL:H143	14:BN:102:CRT:H132	1.97	0.46
6:BN:44:PRO:CD	5:BO:55:TYR:OH	2.62	0.46
5:BQ:44:LEU:HD12	5:BQ:46:TRP:CE3	2.37	0.46
5:BS:42:THR:HG22	5:BS:43:ASP:H	1.80	0.46
14:BU:103:CRT:H9	6:BX:20:ILE:HG23	1.97	0.46
9:BW:102:BCL:OBB	9:BW:102:BCL:HHC	2.14	0.46
9:BY:102:BCL:HED1	6:BZ:31:LEU:HB3	1.96	0.46
6:B6:29:PHE:HA	6:B6:32:VAL:HG12	1.96	0.46
6:B0:32:VAL:HG21	9:B0:102:BCL:CGA	2.44	0.46
1:AC:29:GLY:N	1:AC:44:TYR:O	2.48	0.46
2:AL:48:LEU:HD13	5:AA:33:LEU:CD2	2.45	0.46
2:AL:184:LEU:HB2	2:AL:252:TRP:HE1	1.79	0.46
3:AM:83:VAL:O	3:AM:86:PHE:N	2.47	0.46
3:AM:91:PHE:N	3:AM:91:PHE:CD1	2.82	0.46
3:AM:191:ILE:O	3:AM:192:ARG:C	2.58	0.46
3:AM:215:LEU:HA	3:AM:218:MET:HG3	1.97	0.46
9:AM:401:BCL:CBC	9:AM:402:BCL:CAD	2.93	0.46
4:AH:168:SER:N	4:AH:183:GLU:O	2.48	0.46
4:AH:234:TYR:CE1	4:AH:238:LYS:HE3	2.50	0.46
5:AA:39:VAL:C	5:AA:41:SER:H	2.24	0.46
9:AA:101:BCL:CED	9:A0:102:BCL:H92	2.44	0.46
5:AD:36:HIS:NE2	9:AE:101:BCL:CMD	2.79	0.46
5:AF:30:VAL:HG13	5:AF:31:LEU:N	2.30	0.46
5:AF:31:LEU:HB3	9:AG:101:BCL:CED	2.45	0.46
6:AG:45:TRP:CZ2	9:AG:101:BCL:H2C	2.50	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:AI:27:PHE:CD1	5:AI:27:PHE:C	2.93	0.46
5:AO:27:PHE:HE2	5:AQ:29:ILE:HD11	1.79	0.46
5:AW:36:HIS:NE2	9:AX:101:BCL:CMD	2.79	0.46
6:AX:24:SER:O	6:AX:27:ALA:HB3	2.15	0.46
6:AX:33:VAL:O	6:AX:37:LEU:HB2	2.15	0.46
9:AX:101:BCL:H43	14:AX:102:CRT:H26	1.96	0.46
14:AX:102:CRT:H131	14:AX:102:CRT:H15	1.71	0.46
9:AY:102:BCL:C1D	9:AZ:101:BCL:HMD1	2.44	0.46
9:A3:104:BCL:C2	6:A4:29:PHE:CD1	2.98	0.46
5:A5:4:MET:HE1	9:A7:103:BCL:H91	1.96	0.46
6:A6:31:LEU:HA	6:A6:34:ILE:CG2	2.44	0.46
6:A8:29:PHE:HZ	9:A8:101:BCL:H72	1.80	0.46
5:A9:32:GLY:CA	9:A0:102:BCL:HED2	2.46	0.46
6:A0:24:SER:HB3	14:A0:101:CRT:H183	1.97	0.46
2:BL:142:PHE:HD1	2:BL:143:VAL:N	2.13	0.46
2:BL:175:HIS:NE2	3:BM:184:ASP:OD2	2.47	0.46
2:BL:207:THR:C	2:BL:209:PRO:HD3	2.41	0.46
3:BM:138:GLU:HA	3:BM:142:MET:O	2.16	0.46
3:BM:191:ILE:O	3:BM:192:ARG:C	2.57	0.46
3:BM:221:ALA:O	3:BM:222:THR:C	2.57	0.46
4:BH:169:ASP:H	4:BH:183:GLU:HB2	1.81	0.46
5:BA:44:LEU:HD12	5:BA:46:TRP:H	1.80	0.46
5:BF:33:LEU:HD12	5:BF:33:LEU:N	2.28	0.46
5:BK:25:VAL:O	5:BK:29:ILE:HG22	2.15	0.46
6:BN:31:LEU:CA	6:BN:34:ILE:HG22	2.44	0.46
9:B1:102:BCL:ND	9:B2:101:BCL:CMD	2.77	0.46
6:B2:36:HIS:CE1	9:B2:101:BCL:C4D	2.98	0.46
5:B3:38:ILE:HG23	5:B3:39:VAL:N	2.31	0.46
9:B6:101:BCL:CHC	9:B7:103:BCL:CBB	2.92	0.46
9:B7:103:BCL:H62	6:B8:28:TRP:CZ3	2.51	0.46
6:B8:20:ILE:CG2	6:B8:21:PHE:N	2.77	0.46
2:AL:35:PHE:HA	2:AL:38:VAL:HG22	1.98	0.46
2:AL:218:SER:O	2:AL:220:HIS:N	2.49	0.46
9:AL:301:BCL:OBB	9:AL:301:BCL:HHC	2.14	0.46
9:AL:301:BCL:HMB3	9:AL:301:BCL:HBB2	1.98	0.46
3:AM:13:VAL:HG12	4:AH:144:ILE:HA	1.98	0.46
3:AM:73:PHE:O	3:AM:74:ASN:C	2.59	0.46
3:AM:176:PRO:HD2	3:AM:185:TRP:HB2	1.97	0.46
5:AA:13:LEU:O	6:AB:9:LEU:HD13	2.15	0.46
6:AB:20:ILE:HD11	5:A9:8:LEU:HD23	1.97	0.46
9:AD:102:BCL:C1D	9:AE:101:BCL:HMD3	2.45	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:AE:44:PRO:HG2	5:AF:52:PRO:HG2	1.98	0.46
9:AI:102:BCL:OBB	9:AI:102:BCL:HHC	2.14	0.46
6:AJ:17:PHE:C	6:AJ:17:PHE:CD1	2.93	0.46
5:AU:19:ARG:HB2	5:AU:19:ARG:HH21	1.79	0.46
5:AW:10:LYS:HA	5:AW:13:LEU:HD12	1.98	0.46
9:AW:101:BCL:HMB1	9:AW:101:BCL:HBB3	1.96	0.46
5:A1:29:ILE:O	5:A1:33:LEU:HG	2.15	0.46
5:A3:47:LEU:HB3	5:A3:48:ASP:H	1.52	0.46
5:A7:9:TYR:CE2	5:A7:10:LYS:HE3	2.50	0.46
6:A0:26:TYR:O	6:A0:27:ALA:C	2.59	0.46
1:BC:148:THR:HA	1:BC:322:GLN:HG2	1.96	0.46
2:BL:17:LEU:HD21	2:BL:114:VAL:HG12	1.97	0.46
2:BL:75:ILE:HG22	2:BL:95:TRP:HB2	1.96	0.46
2:BL:98:ILE:HG22	2:BL:99:THR:N	2.29	0.46
2:BL:159:ILE:HD12	2:BL:159:ILE:N	2.29	0.46
2:BL:191:THR:CG2	11:BL:304:UQ8:H16	2.45	0.46
2:BL:218:SER:O	2:BL:220:HIS:N	2.48	0.46
2:BL:253:SER:O	2:BL:254:ALA:C	2.57	0.46
3:BM:178:GLY:O	3:BM:182:HIS:CB	2.61	0.46
3:BM:196:LEU:HD12	9:BM:402:BCL:CHD	2.45	0.46
4:BH:67:PHE:N	4:BH:76:VAL:O	2.39	0.46
4:BH:215:LYS:HE3	4:BH:250:ALA:O	2.15	0.46
6:BE:33:VAL:HG22	6:BE:37:LEU:HD23	1.95	0.46
6:BJ:21:PHE:O	6:BJ:22:MET:C	2.58	0.46
5:BK:5:ASN:HD22	6:BN:22:MET:HE3	1.80	0.46
5:BO:50:ASN:OD1	5:BO:51:ILE:N	2.49	0.46
9:BO:102:BCL:HBB3	9:BO:102:BCL:HMB1	1.98	0.46
5:BQ:19:ARG:O	5:BQ:22:VAL:HG12	2.16	0.46
5:BS:7:ASN:HB3	5:BS:10:LYS:CE	2.42	0.46
5:B3:55:TYR:HB2	5:B3:56:GLN:NE2	2.30	0.46
5:B5:18:ARG:CB	5:B5:19:ARG:NH1	2.78	0.46
6:B6:38:LEU:HD23	6:B6:38:LEU:O	2.15	0.46
1:AC:170:PRO:CG	1:AC:171:GLY:N	2.78	0.46
2:AL:151:TRP:C	2:AL:153:HIS:N	2.72	0.46
2:AL:231:TYR:CG	2:AL:232:SER:N	2.83	0.46
3:AM:85:GLN:O	3:AM:89:HIS:N	2.47	0.46
3:AM:181:PRO:HA	3:AM:184:ASP:OD1	2.16	0.46
4:AH:15:THR:O	4:AH:18:ALA:HB3	2.16	0.46
4:AH:197:ILE:CG2	4:AH:198:GLY:N	2.79	0.46
6:AJ:22:MET:O	6:AJ:26:TYR:HD1	1.99	0.46
5:AK:5:ASN:ND2	6:AN:22:MET:CE	2.60	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:AN:19:ALA:HB3	6:AN:20:ILE:HD12	1.98	0.46
5:AO:43:ASP:OD2	5:AQ:47:LEU:HB3	2.15	0.46
6:AP:44:PRO:HG2	5:AQ:52:PRO:CG	2.45	0.46
6:AT:16:GLU:HG2	6:AT:17:PHE:N	2.30	0.46
5:AW:30:VAL:HG13	5:AW:31:LEU:N	2.30	0.46
14:AW:102:CRT:C8	6:AZ:20:ILE:HD13	2.46	0.46
5:A3:15:LEU:CD1	9:A5:102:BCL:H151	2.46	0.46
5:A7:25:VAL:HG13	9:A7:103:BCL:H51	1.97	0.46
1:BC:153:TYR:HD2	1:BC:323:MET:HE1	1.80	0.46
2:BL:257:ILE:O	2:BL:257:ILE:HG13	2.16	0.46
2:BL:271:TRP:CZ2	2:BL:274:TRP:NE1	2.80	0.46
3:BM:250:LEU:O	3:BM:253:ARG:HB3	2.16	0.46
9:BM:402:BCL:HHC	9:BM:402:BCL:OBB	2.16	0.46
4:BH:23:PHE:C	4:BH:25:GLY:N	2.68	0.46
6:BB:18:HIS:CE1	6:BB:22:MET:HE2	2.50	0.46
5:BD:35:ILE:O	5:BD:39:VAL:HG23	2.16	0.46
6:BE:44:PRO:CG	5:BF:55:TYR:OH	2.58	0.46
6:BJ:38:LEU:HD23	6:BJ:38:LEU:C	2.40	0.46
9:BJ:101:BCL:HHC	9:BJ:101:BCL:OBB	2.14	0.46
9:BN:101:BCL:C1B	9:BO:102:BCL:CMB	2.94	0.46
6:BP:39:ALA:O	6:BP:42:TYR:N	2.49	0.46
9:BS:102:BCL:C4D	9:BT:101:BCL:HMD3	2.45	0.46
5:BU:53:VAL:HA	5:BU:55:TYR:CE1	2.50	0.46
6:BX:33:VAL:O	6:BX:37:LEU:HB2	2.15	0.46
6:BX:42:TYR:CE2	6:BX:43:ARG:HD2	2.50	0.46
5:BY:35:ILE:O	5:BY:36:HIS:C	2.59	0.46
5:B1:40:LEU:HB2	5:B1:46:TRP:CH2	2.51	0.46
9:B2:101:BCL:HBB3	9:B3:102:BCL:CHC	2.45	0.46
5:B3:51:ILE:HA	5:B3:53:VAL:N	2.22	0.46
6:B4:38:LEU:O	6:B4:38:LEU:HD23	2.16	0.46
5:B5:46:TRP:O	5:B5:49:ASP:OD1	2.33	0.46
5:B7:48:ASP:HA	5:B7:53:VAL:CB	2.46	0.46
6:B0:37:LEU:HA	9:B0:102:BCL:H193	1.96	0.46
1:AC:20:LEU:HD13	1:AC:21:LEU:N	2.30	0.46
1:AC:82:LEU:HD13	1:AC:93:THR:HG21	1.97	0.46
2:AL:3:MET:HE1	2:AL:8:LYS:HG3	1.97	0.46
2:AL:160:LEU:HD12	2:AL:161:SER:N	2.30	0.46
3:AM:74:ASN:CG	3:AM:95:LEU:HD13	2.40	0.46
3:AM:137:ALA:HB3	3:AM:144:GLN:HE22	1.81	0.46
3:AM:189:PHE:O	3:AM:190:SER:C	2.59	0.46
5:AA:36:HIS:CD2	9:AB:101:BCL:HMD2	2.51	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:AD:43:ASP:CG	5:AD:44:LEU:H	2.24	0.46
6:AJ:17:PHE:CE1	6:AJ:21:PHE:HB2	2.51	0.46
9:AJ:101:BCL:C3B	9:AK:102:BCL:C3B	2.93	0.46
5:AQ:12:TRP:HA	5:AQ:12:TRP:CE3	2.51	0.46
5:AS:34:LEU:CB	15:AS:101:PEF:C44	2.86	0.46
5:AU:26:ALA:CA	5:AU:29:ILE:HG22	2.45	0.46
9:AU:102:BCL:O2D	6:AV:32:VAL:HG23	2.16	0.46
5:A1:36:HIS:CE1	9:A2:101:BCL:CMD	2.98	0.46
5:A5:10:LYS:HB3	14:A5:103:CRT:H5	1.97	0.46
9:A8:101:BCL:HMA3	9:A9:102:BCL:HMA2	1.97	0.46
1:BC:175:PRO:HG3	3:BM:80:HIS:HA	1.98	0.46
2:BL:22:LEU:HB2	5:B7:19:ARG:HG3	1.97	0.46
2:BL:147:LEU:HB3	2:BL:262:PRO:HB3	1.97	0.46
3:BM:153:ALA:HA	3:BM:277:VAL:HG11	1.98	0.46
3:BM:175:VAL:HG22	3:BM:185:TRP:CE3	2.51	0.46
3:BM:238:ILE:HD13	3:BM:262:MET:HG3	1.98	0.46
4:BH:168:SER:N	4:BH:183:GLU:O	2.49	0.46
4:BH:171:TRP:N	4:BH:181:TYR:O	2.49	0.46
4:BH:184:VAL:HB	4:BH:193:VAL:HG23	1.97	0.46
4:BH:197:ILE:HD13	4:BH:197:ILE:C	2.41	0.46
9:BB:101:BCL:HBB3	9:BB:101:BCL:HMB1	1.98	0.46
5:BI:20:VAL:O	5:BI:24:ILE:HG12	2.15	0.46
5:BO:44:LEU:CD1	5:BO:46:TRP:N	2.79	0.46
5:BS:17:PRO:HA	5:BS:20:VAL:HG22	1.98	0.46
5:BS:43:ASP:CB	5:BU:56:GLN:HG3	2.45	0.46
6:BT:7:THR:OG1	6:BT:8:GLY:N	2.47	0.46
5:BU:43:ASP:HB2	5:BW:47:LEU:CB	2.45	0.46
6:BX:24:SER:O	6:BX:27:ALA:HB3	2.16	0.46
6:BX:29:PHE:N	6:BX:29:PHE:CD1	2.83	0.46
5:BY:33:LEU:HD12	5:BY:33:LEU:C	2.41	0.46
9:BZ:101:BCL:H41	9:B1:102:BCL:HMA1	1.97	0.46
6:B2:25:MET:HG2	6:B2:29:PHE:HE2	1.80	0.46
9:B3:102:BCL:O2D	6:B4:32:VAL:HG22	2.15	0.46
6:B4:29:PHE:CE1	9:B4:101:BCL:H72	2.42	0.46
5:B9:2:PHE:HA	5:B9:5:ASN:HD22	1.80	0.46
1:AC:247:CYS:O	1:AC:251:HIS:HB2	2.16	0.46
1:AC:264:PRO:HG2	1:AC:265:LYS:CD	2.43	0.46
2:AL:46:GLY:HA2	2:AL:49:LEU:HB3	1.98	0.46
3:AM:102:TYR:O	3:AM:103:GLY:C	2.59	0.46
3:AM:115:TRP:HZ3	3:AM:116:LEU:HD12	1.81	0.46
3:AM:277:VAL:HG22	10:AM:403:BPH:HBC1	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:AM:307:TYR:N	3:AM:307:TYR:CD1	2.84	0.46
10:AM:403:BPH:H112	15:AS:101:PEF:C16	2.46	0.46
4:AH:30:LEU:O	4:AH:34:ASP:N	2.48	0.46
9:AA:101:BCL:C2D	6:AB:35:ALA:HB1	2.46	0.46
5:AF:28:GLN:HB2	9:AF:102:BCL:H43	1.97	0.46
9:AF:102:BCL:HMD3	6:AG:36:HIS:CD2	2.51	0.46
9:AI:102:BCL:H143	14:AJ:102:CRT:H132	1.98	0.46
6:AP:15:LYS:O	6:AP:16:GLU:C	2.58	0.46
5:AS:13:LEU:CB	14:AS:104:CRT:H31A	2.42	0.46
6:AT:33:VAL:HG13	6:AT:34:ILE:N	2.30	0.46
14:AT:102:CRT:H2M1	5:AU:37:MET:HG2	1.96	0.46
5:AU:42:THR:HB	5:AW:48:ASP:CB	2.45	0.46
6:AV:45:TRP:CE3	9:AV:102:BCL:HAC2	2.51	0.46
5:AY:30:VAL:CA	5:AY:33:LEU:HG	2.45	0.46
6:A4:10:THR:O	6:A4:14:ALA:HB2	2.14	0.46
5:A5:29:ILE:HG23	5:A5:30:VAL:N	2.31	0.46
5:A5:35:ILE:HA	5:A5:38:ILE:CG2	2.45	0.46
5:A5:43:ASP:HB2	5:A7:47:LEU:CG	2.45	0.46
5:A7:32:GLY:N	9:A8:101:BCL:HED2	2.31	0.46
9:A8:101:BCL:CHC	9:A9:102:BCL:CAB	2.93	0.46
1:BC:52:SER:CB	1:BC:319:TYR:OH	2.64	0.46
1:BC:133:LEU:HA	1:BC:283:TYR:CE2	2.51	0.46
1:BC:164:TYR:O	1:BC:309:THR:HG23	2.15	0.46
1:BC:201:THR:N	1:BC:202:PRO:HD2	2.31	0.46
1:BC:210:ILE:O	1:BC:210:ILE:CG2	2.60	0.46
1:BC:316:LYS:O	1:BC:320:GLY:N	2.44	0.46
2:BL:30:PHE:HD2	3:BM:255:THR:O	1.99	0.46
2:BL:94:LEU:HD23	2:BL:95:TRP:N	2.31	0.46
2:BL:192:ASN:ND2	2:BL:193:CYS:N	2.63	0.46
2:BL:193:CYS:O	10:BM:403:BPH:H3C	2.15	0.46
2:BL:214:PRO:HA	4:BH:68:VAL:O	2.16	0.46
3:BM:196:LEU:C	3:BM:198:TYR:N	2.73	0.46
4:BH:80:ARG:HG3	4:BH:80:ARG:NH1	2.31	0.46
4:BH:182:LEU:HD12	4:BH:182:LEU:N	2.30	0.46
5:BA:31:LEU:HD21	14:BB:102:CRT:H32	1.98	0.46
6:BB:29:PHE:HE1	9:BB:101:BCL:C2	2.28	0.46
5:BD:46:TRP:NE1	9:BD:102:BCL:HHC	2.31	0.46
6:BE:23:GLN:CG	6:BE:24:SER:N	2.78	0.46
5:BF:9:TYR:HA	6:BG:18:HIS:ND1	2.31	0.46
5:BF:12:TRP:CZ2	6:BG:21:PHE:CE2	3.04	0.46
6:BJ:17:PHE:C	6:BJ:17:PHE:HD1	2.24	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:BN:19:ALA:HB3	6:BN:20:ILE:HD12	1.98	0.46
5:BO:43:ASP:OD2	5:BQ:47:LEU:O	2.33	0.46
9:BQ:103:BCL:C2D	9:BQ:104:BCL:C2D	2.93	0.46
5:BS:12:TRP:NE1	6:BT:18:HIS:ND1	2.64	0.46
9:BS:102:BCL:C1D	9:BT:101:BCL:HMD1	2.45	0.46
9:BU:102:BCL:C2D	9:BV:101:BCL:HMD1	2.45	0.46
6:BV:10:THR:CG2	6:BV:11:ASP:N	2.79	0.46
6:BV:17:PHE:HB2	14:BV:102:CRT:H42	1.97	0.46
5:BY:32:GLY:HA3	9:BY:102:BCL:O1A	2.15	0.46
5:B1:46:TRP:CZ3	9:B1:102:BCL:HBC3	2.51	0.46
6:B2:21:PHE:HE1	14:B2:102:CRT:C19	2.28	0.46
6:B4:13:GLU:O	6:B4:16:GLU:CG	2.63	0.46
6:B6:31:LEU:HA	6:B6:34:ILE:CG2	2.45	0.46
6:B8:22:MET:HG3	6:B8:26:TYR:HE2	1.80	0.46
9:B0:102:BCL:OBB	9:B0:102:BCL:HHC	2.15	0.46
1:AC:129:ARG:HG2	1:AC:287:LEU:HD11	1.98	0.46
1:AC:167:VAL:CG2	1:AC:297:GLY:HA3	2.41	0.46
2:AL:46:GLY:O	2:AL:48:LEU:N	2.49	0.46
2:AL:46:GLY:C	2:AL:48:LEU:N	2.70	0.46
2:AL:101:CYS:O	2:AL:102:ALA:C	2.59	0.46
2:AL:199:HIS:C	2:AL:201:SER:H	2.24	0.46
2:AL:238:ILE:HD12	2:AL:238:ILE:HA	1.84	0.46
2:AL:268:TRP:O	2:AL:271:TRP:N	2.49	0.46
11:AL:304:UQ8:H10	11:AL:304:UQ8:H7	1.76	0.46
4:AH:152:ARG:HG2	4:AH:168:SER:O	2.16	0.46
4:AH:180:ARG:HG2	4:AH:180:ARG:NH1	2.31	0.46
5:AA:47:LEU:HB3	5:A9:43:ASP:HA	1.98	0.46
6:AE:38:LEU:HD23	6:AE:38:LEU:C	2.41	0.46
6:AJ:23:GLN:C	6:AJ:23:GLN:CD	2.84	0.46
5:AO:43:ASP:CB	5:AQ:47:LEU:HB3	2.46	0.46
5:AY:4:MET:HB2	5:AY:8:LEU:HD12	1.98	0.46
5:A1:21:LEU:HD11	9:A1:102:BCL:H142	1.97	0.46
5:A1:36:HIS:CE1	9:A1:102:BCL:NA	2.83	0.46
5:A5:4:MET:SD	6:A8:24:SER:O	2.74	0.46
5:A5:44:LEU:CD1	5:A5:46:TRP:HB3	2.45	0.46
5:A7:10:LYS:CB	14:A0:101:CRT:H83	2.45	0.46
5:A7:33:LEU:O	5:A7:37:MET:HB2	2.15	0.46
5:A9:4:MET:HE3	5:A9:4:MET:CA	2.46	0.46
2:BL:213:GLU:OE2	2:BL:213:GLU:HA	2.16	0.46
2:BL:226:ARG:O	3:BM:50:PRO:O	2.33	0.46
3:BM:98:PRO:HA	3:BM:99:PRO:HD3	1.83	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:BM:200:PRO:HG2	3:BM:201:PHE:H	1.81	0.46
9:BM:402:BCL:HMB1	9:BM:402:BCL:CBB	2.46	0.46
5:BA:15:LEU:O	5:BA:16:ASP:C	2.58	0.46
6:BB:33:VAL:HG13	6:BB:34:ILE:N	2.31	0.46
5:BD:40:LEU:HD11	5:BD:47:LEU:HD23	1.98	0.46
5:BF:12:TRP:HB2	6:BG:14:ALA:HB1	1.96	0.46
9:BF:102:BCL:HMD3	6:BG:36:HIS:CD2	2.51	0.46
5:BO:26:ALA:HA	5:BO:29:ILE:CG2	2.45	0.46
5:BQ:35:ILE:CD1	9:BS:102:BCL:HMB2	2.46	0.46
6:BT:45:TRP:O	6:BT:46:LEU:HB2	2.16	0.46
5:BW:54:SER:O	5:BW:58:LEU:N	2.43	0.46
9:BY:102:BCL:CHD	9:BZ:101:BCL:HMD1	2.45	0.46
5:B1:13:LEU:O	6:B2:7:THR:HA	2.15	0.46
9:B2:101:BCL:HBB3	9:B2:101:BCL:HMB1	1.98	0.46
5:B5:44:LEU:HD23	5:B5:44:LEU:N	2.31	0.46
5:B7:40:LEU:HD12	5:B7:45:ASN:HA	1.97	0.46
1:AC:20:LEU:HG	2:AL:271:TRP:CE2	2.50	0.46
1:AC:26:PRO:HB2	1:AC:27:PRO:HA	1.98	0.46
2:AL:50:ILE:HG23	2:AL:51:VAL:N	2.30	0.46
10:AL:302:BPH:HED1	3:AM:255:THR:HG21	1.98	0.46
3:AM:164:ARG:C	3:AM:164:ARG:HD3	2.41	0.46
3:AM:199:ASN:HA	3:AM:294:TRP:CE3	2.50	0.46
3:AM:205:SER:C	9:AM:402:BCL:CMA	2.89	0.46
15:AM:407:PEF:H32	4:AH:29:TYR:CZ	2.51	0.46
4:AH:35:LYS:HZ1	4:AH:59:PRO:HD2	1.80	0.46
4:AH:182:LEU:HD13	4:AH:195:LEU:CD2	2.46	0.46
5:AA:27:PHE:C	5:AA:30:VAL:HG12	2.41	0.46
9:AA:101:BCL:OBB	9:AA:101:BCL:HHC	2.15	0.46
5:AO:11:ILE:CG1	14:AR:102:CRT:H81	2.45	0.46
9:AQ:102:BCL:HMB1	9:AQ:102:BCL:CBB	2.46	0.46
5:AU:30:VAL:CG1	5:AU:31:LEU:H	2.28	0.46
5:AW:35:ILE:HG23	5:AW:36:HIS:N	2.30	0.46
5:A3:8:LEU:O	5:A3:11:ILE:HG13	2.16	0.46
5:A5:8:LEU:O	5:A5:11:ILE:HG13	2.16	0.46
5:A5:19:ARG:HG2	5:A5:19:ARG:HH21	1.81	0.46
5:A7:7:ASN:O	5:A7:10:LYS:HD3	2.16	0.46
9:A8:101:BCL:C20	9:A8:101:BCL:C15	2.94	0.46
6:A0:21:PHE:CG	6:A0:22:MET:N	2.84	0.46
6:A0:45:TRP:O	6:A0:46:LEU:CB	2.60	0.46
1:BC:166:TRP:O	1:BC:166:TRP:CD2	2.69	0.46
2:BL:50:ILE:CA	2:BL:98:ILE:HD11	2.46	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:BL:179:ASN:HB3	2:BL:182:HIS:HB3	1.96	0.46
2:BL:207:THR:CG2	3:BM:238:ILE:HG13	2.41	0.46
2:BL:253:SER:C	9:BL:301:BCL:CED	2.86	0.46
3:BM:189:PHE:O	3:BM:190:SER:C	2.59	0.46
3:BM:211:GLY:O	3:BM:214:LEU:N	2.42	0.46
4:BH:141:GLU:OE1	4:BH:141:GLU:N	2.46	0.46
5:BA:18:ARG:H	5:BA:18:ARG:CD	2.19	0.46
5:BA:47:LEU:HD22	5:BA:47:LEU:N	2.31	0.46
6:BE:20:ILE:O	6:BE:23:GLN:HG3	2.15	0.46
6:BG:33:VAL:O	6:BG:37:LEU:HB2	2.15	0.46
9:BN:101:BCL:HMB3	9:BO:102:BCL:CHB	2.46	0.46
5:BO:3:THR:C	5:BO:5:ASN:H	2.23	0.46
5:BO:9:TYR:HA	6:BP:18:HIS:CE1	2.51	0.46
5:BQ:18:ARG:HA	5:BQ:21:LEU:HD12	1.98	0.46
5:BQ:22:VAL:HA	5:BQ:25:VAL:HG12	1.98	0.46
9:BQ:104:BCL:HHB	9:BS:102:BCL:HMA1	1.98	0.46
14:B1:103:CRT:H81	14:B1:103:CRT:H10	1.79	0.46
9:B6:101:BCL:HHC	9:B6:101:BCL:OBB	2.14	0.46
6:B0:25:MET:HG3	14:B0:101:CRT:C20	2.46	0.46
1:AC:235:LEU:C	1:AC:239:ILE:HD12	2.41	0.46
2:AL:51:VAL:HG11	5:AA:37:MET:HG2	1.98	0.46
2:AL:78:PRO:O	2:AL:152:GLY:HA3	2.16	0.46
2:AL:86:MET:HE2	2:AL:96:GLN:CD	2.41	0.46
2:AL:106:PHE:CD2	9:AL:301:BCL:H91	2.50	0.46
2:AL:138:LEU:O	2:AL:142:PHE:N	2.49	0.46
3:AM:39:TRP:O	3:AM:40:LEU:C	2.60	0.46
3:AM:66:VAL:CG1	3:AM:121:PHE:HD2	2.20	0.46
3:AM:156:PHE:CD1	3:AM:281:GLY:CA	2.99	0.46
4:AH:48:ARG:HH21	15:AH:301:PEF:P	2.38	0.46
4:AH:49:SER:O	15:AH:301:PEF:H32	2.15	0.46
4:AH:95:ALA:HB3	4:AH:98:SER:OG	2.15	0.46
4:AH:130:LEU:HD11	4:AH:174:ARG:HH22	1.81	0.46
5:AD:29:ILE:HB	9:AD:102:BCL:C4	2.46	0.46
5:AF:24:ILE:C	5:AF:26:ALA:N	2.73	0.46
5:AI:51:ILE:HA	5:AI:52:PRO:HA	1.69	0.46
9:AI:102:BCL:HBB3	9:AI:102:BCL:HMB1	1.98	0.46
6:AJ:16:GLU:OE2	14:AJ:102:CRT:H23	2.16	0.46
9:AK:102:BCL:HAC2	9:AN:101:BCL:HBC1	1.96	0.46
5:AS:9:TYR:C	5:AS:9:TYR:CD1	2.94	0.46
9:AS:103:BCL:C1D	9:AT:101:BCL:CMD	2.93	0.46
9:AW:101:BCL:HMB1	9:AW:101:BCL:CBB	2.46	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:A2:21:PHE:HE1	14:A2:102:CRT:C16	2.24	0.46
5:A3:19:ARG:HG2	5:A3:20:VAL:N	2.31	0.46
5:A5:32:GLY:HA2	9:A6:101:BCL:HED2	1.98	0.46
5:A5:53:VAL:O	5:A5:54:SER:C	2.59	0.46
5:A7:7:ASN:CA	5:A7:10:LYS:HZ2	2.29	0.46
9:A7:103:BCL:HBB3	9:A7:103:BCL:HMB1	1.98	0.46
1:BC:232:THR:O	1:BC:235:LEU:HB3	2.15	0.46
2:BL:17:LEU:HD11	2:BL:114:VAL:CB	2.39	0.46
2:BL:137:TYR:CD1	2:BL:137:TYR:C	2.93	0.46
2:BL:192:ASN:HD21	3:BM:212:SER:C	2.24	0.46
3:BM:152:ALA:CB	3:BM:274:VAL:HG13	2.46	0.46
4:BH:170:VAL:HG12	4:BH:182:LEU:HB3	1.98	0.46
5:BA:29:ILE:O	5:BA:33:LEU:HD13	2.16	0.46
5:BA:33:LEU:CA	14:B0:101:CRT:C2M	2.94	0.46
5:BA:47:LEU:CB	5:B9:43:ASP:HB2	2.46	0.46
5:BK:12:TRP:CD1	6:BN:17:PHE:HD2	2.34	0.46
5:BO:8:LEU:HG	6:BP:18:HIS:NE2	2.31	0.46
6:BP:30:GLY:HA2	6:BP:33:VAL:HG12	1.98	0.46
5:BU:13:LEU:HD21	6:BV:14:ALA:HB2	1.98	0.46
9:BU:102:BCL:HHC	9:BU:102:BCL:OBB	2.15	0.46
9:BU:102:BCL:OBD	6:BV:32:VAL:HG23	2.16	0.46
6:BV:33:VAL:HG13	6:BV:34:ILE:N	2.31	0.46
6:BX:9:LEU:HD22	6:BX:13:GLU:HG3	1.98	0.46
6:B2:10:THR:HG23	6:B2:12:ASP:H	1.81	0.46
6:B2:21:PHE:CG	6:B2:22:MET:N	2.84	0.46
5:B3:56:GLN:H	5:B3:56:GLN:CD	2.24	0.46
6:B4:40:TRP:CZ3	6:B4:44:PRO:CA	2.96	0.46
9:B7:103:BCL:CHD	9:B8:101:BCL:HMD1	2.46	0.46
1:AC:121:ILE:O	1:AC:124:LYS:N	2.48	0.45
2:AL:171:TYR:OH	3:AM:191:ILE:HD11	2.16	0.45
2:AL:242:GLY:CA	3:AM:216:PHE:CE2	2.99	0.45
10:AL:302:BPH:H162	9:AL:303:BCL:CMB	2.46	0.45
10:AL:302:BPH:H6C1	9:AL:303:BCL:H202	1.98	0.45
3:AM:40:LEU:C	3:AM:40:LEU:HD23	2.42	0.45
3:AM:74:ASN:ND2	3:AM:95:LEU:HD13	2.32	0.45
3:AM:79:VAL:O	3:AM:79:VAL:HG22	2.16	0.45
15:AM:408:PEF:C5	4:AH:204:LYS:HE2	2.41	0.45
4:AH:246:GLY:O	4:AH:254:ARG:NH1	2.49	0.45
9:AA:101:BCL:CMD	6:AB:35:ALA:HB1	2.44	0.45
6:AE:31:LEU:HA	6:AE:34:ILE:HG22	1.97	0.45
5:AF:8:LEU:HD22	14:AJ:102:CRT:H133	1.96	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:AF:50:ASN:CG	5:AF:51:ILE:N	2.73	0.45
5:AK:22:VAL:O	5:AK:25:VAL:HB	2.15	0.45
6:AR:17:PHE:O	6:AR:20:ILE:HG22	2.16	0.45
6:AT:15:LYS:HG2	6:AT:16:GLU:N	2.30	0.45
6:AT:22:MET:HB3	6:AT:26:TYR:HE1	1.80	0.45
5:AU:19:ARG:HB2	5:AU:19:ARG:NH2	2.31	0.45
9:AU:102:BCL:O2D	6:AV:32:VAL:CG2	2.64	0.45
5:AW:4:MET:HE1	6:AZ:23:GLN:HE21	1.81	0.45
6:AZ:40:TRP:O	6:AZ:40:TRP:CD1	2.69	0.45
5:A1:21:LEU:O	5:A1:25:VAL:HG23	2.15	0.45
5:A3:9:TYR:CD1	5:A3:9:TYR:C	2.95	0.45
9:A3:103:BCL:C2D	9:A3:104:BCL:C2D	2.94	0.45
5:A5:25:VAL:HG13	9:A5:102:BCL:C5	2.46	0.45
5:A7:7:ASN:HD22	5:A7:7:ASN:N	2.09	0.45
9:A8:101:BCL:H203	9:A8:101:BCL:C15	2.32	0.45
6:A0:26:TYR:O	6:A0:29:PHE:HB2	2.16	0.45
1:BC:107:CYS:C	1:BC:109:TYR:N	2.73	0.45
1:BC:135:ARG:HG2	1:BC:330:LEU:HA	1.98	0.45
1:BC:212:ILE:O	1:BC:222:ASN:ND2	2.49	0.45
1:BC:225:SER:O	1:BC:226:LEU:C	2.58	0.45
2:BL:184:LEU:CB	2:BL:252:TRP:NE1	2.79	0.45
3:BM:157:TYR:CD1	3:BM:157:TYR:C	2.93	0.45
3:BM:191:ILE:O	3:BM:193:TYR:N	2.49	0.45
4:BH:173:ASP:OD1	4:BH:174:ARG:N	2.49	0.45
4:BH:235:GLU:HA	4:BH:238:LYS:CB	2.45	0.45
4:BH:246:GLY:O	4:BH:254:ARG:NH1	2.47	0.45
9:BF:102:BCL:HAC2	9:BG:101:BCL:CBC	2.46	0.45
5:BK:5:ASN:ND2	6:BN:22:MET:HG2	2.31	0.45
5:BO:26:ALA:C	5:BO:29:ILE:HG22	2.39	0.45
5:BQ:17:PRO:O	5:BQ:20:VAL:HG22	2.15	0.45
5:BS:36:HIS:ND1	9:BT:101:BCL:HMD2	2.31	0.45
5:BW:8:LEU:HD22	5:BW:11:ILE:HD11	1.98	0.45
5:B1:13:LEU:CB	14:B1:103:CRT:C1M	2.78	0.45
6:B2:17:PHE:HD1	14:B2:102:CRT:C6	2.27	0.45
6:B8:22:MET:HG3	6:B8:26:TYR:CE2	2.51	0.45
2:AL:30:PHE:CZ	3:AM:257:GLY:HA3	2.51	0.45
2:AL:32:VAL:HG12	2:AL:37:VAL:HG13	1.98	0.45
2:AL:275:TRP:O	2:AL:276:LEU:C	2.59	0.45
2:AL:279:PRO:C	2:AL:280:LEU:HG	2.39	0.45
3:AM:34:PRO:HG3	3:AM:50:PRO:HD3	1.98	0.45
3:AM:70:ILE:C	3:AM:72:GLY:H	2.23	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:AM:291:VAL:HG21	3:AM:297:TRP:CD1	2.52	0.45
4:AH:122:HIS:N	4:AH:232:THR:HB	2.32	0.45
9:AA:101:BCL:HBC1	9:AB:101:BCL:HBC3	1.99	0.45
14:AA:102:CRT:H32	5:AD:31:LEU:CD2	2.46	0.45
5:AD:29:ILE:HG23	5:AD:30:VAL:N	2.31	0.45
6:AE:10:THR:CG2	6:AE:11:ASP:H	2.23	0.45
5:AF:26:ALA:O	5:AF:29:ILE:N	2.49	0.45
6:AN:17:PHE:HD1	14:AN:102:CRT:H6	1.80	0.45
6:AP:33:VAL:HG22	6:AP:37:LEU:HD23	1.97	0.45
5:AQ:43:ASP:HB2	5:AS:47:LEU:CB	2.46	0.45
5:AS:34:LEU:HG	5:AS:34:LEU:O	2.17	0.45
9:AS:103:BCL:ND	9:AT:101:BCL:HMD3	2.31	0.45
6:AT:45:TRP:O	5:AU:52:PRO:HD2	2.15	0.45
9:AT:101:BCL:HBB2	9:AT:101:BCL:HMB3	1.98	0.45
9:AU:102:BCL:ND	9:AV:102:BCL:HMD3	2.31	0.45
5:AW:22:VAL:O	5:AW:25:VAL:HB	2.15	0.45
5:AW:33:LEU:HD12	5:AW:34:LEU:N	2.31	0.45
9:AW:101:BCL:H71	6:AX:28:TRP:CE3	2.51	0.45
9:AW:101:BCL:CBC	9:AX:101:BCL:HHD	2.44	0.45
5:AY:9:TYR:CG	6:AZ:15:LYS:HG2	2.52	0.45
6:AZ:45:TRP:HE3	9:AZ:101:BCL:HAC2	1.72	0.45
5:A3:46:TRP:NE1	5:A3:47:LEU:HD22	2.32	0.45
5:A5:4:MET:HG3	6:A8:27:ALA:HB1	1.89	0.45
6:A6:44:PRO:O	5:A7:52:PRO:CD	2.63	0.45
1:BC:48:GLN:C	1:BC:50:ALA:H	2.24	0.45
1:BC:68:THR:O	1:BC:86:SER:HB2	2.17	0.45
1:BC:76:TYR:HB3	7:BC:501:HEM:O2A	2.16	0.45
1:BC:268:THR:HG21	7:BC:504:HEM:HAA1	1.97	0.45
2:BL:16:THR:HG1	4:BH:257:PRO:HB3	1.79	0.45
2:BL:22:LEU:HB2	5:B7:19:ARG:HB3	1.95	0.45
2:BL:89:LEU:HD11	5:B7:38:ILE:HD11	1.98	0.45
2:BL:106:PHE:CD1	10:BL:302:BPH:H1C1	2.51	0.45
2:BL:240:ARG:HH21	3:BM:6:ASN:C	2.24	0.45
3:BM:237:GLN:CD	3:BM:244:ALA:HB3	2.41	0.45
3:BM:284:ILE:CG1	9:BM:402:BCL:OBD	2.64	0.45
3:BM:284:ILE:CD1	9:BM:402:BCL:OBD	2.64	0.45
4:BH:90:THR:HG23	4:BH:103:ASN:OD1	2.16	0.45
5:BA:35:ILE:C	5:BA:37:MET:N	2.73	0.45
5:BO:44:LEU:CD1	5:BO:46:TRP:H	2.30	0.45
5:BU:38:ILE:HD12	14:BV:102:CRT:C40	2.46	0.45
6:BV:18:HIS:NE2	6:BV:22:MET:CE	2.79	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:B1:8:LEU:C	5:B1:10:LYS:N	2.73	0.45
6:B2:31:LEU:C	6:B2:34:ILE:HG22	2.41	0.45
9:B6:101:BCL:HMC1	9:B7:103:BCL:CBB	2.42	0.45
6:B0:26:TYR:O	6:B0:27:ALA:C	2.58	0.45
1:AC:153:TYR:HB3	1:AC:323:MET:CE	2.23	0.45
1:AC:161:VAL:HG22	7:AC:502:HEM:O1D	2.17	0.45
2:AL:30:PHE:HA	3:AM:254:TRP:HA	1.98	0.45
2:AL:30:PHE:HD2	3:AM:255:THR:O	2.00	0.45
2:AL:31:TYR:CD1	2:AL:32:VAL:N	2.84	0.45
2:AL:108:SER:O	2:AL:111:LEU:N	2.50	0.45
2:AL:144:ARG:CB	2:AL:145:PRO:CD	2.95	0.45
3:AM:81:TRP:C	5:AU:41:SER:HB3	2.40	0.45
3:AM:84:PHE:HA	5:AW:37:MET:HE1	1.96	0.45
3:AM:134:TYR:CA	3:AM:144:GLN:NE2	2.78	0.45
3:AM:170:SER:O	3:AM:172:ALA:N	2.48	0.45
3:AM:199:ASN:C	3:AM:203:MET:HE3	2.42	0.45
3:AM:267:ARG:NH2	15:AM:407:PEF:O2P	2.50	0.45
3:AM:286:LEU:HD22	3:AM:290:VAL:HG11	1.99	0.45
4:AH:169:ASP:N	4:AH:183:GLU:HB2	2.31	0.45
5:AD:40:LEU:O	5:AD:41:SER:O	2.34	0.45
5:AF:49:ASP:CB	5:AI:56:GLN:HB2	2.46	0.45
6:AG:38:LEU:HD23	6:AG:38:LEU:C	2.41	0.45
6:AJ:38:LEU:HD23	6:AJ:38:LEU:C	2.41	0.45
5:AK:5:ASN:O	5:AK:8:LEU:HB2	2.16	0.45
5:AO:12:TRP:CH2	6:AP:17:PHE:CE2	3.04	0.45
6:AR:10:THR:H	6:AR:13:GLU:CD	2.25	0.45
5:AS:49:ASP:CG	5:AS:50:ASN:H	2.24	0.45
5:AW:2:PHE:CB	5:AW:5:ASN:HB2	2.46	0.45
5:AW:45:ASN:O	5:AW:47:LEU:N	2.50	0.45
9:AY:102:BCL:CHD	9:AZ:101:BCL:HMD1	2.46	0.45
9:AZ:101:BCL:H2A	9:AZ:101:BCL:CGD	2.46	0.45
5:A1:32:GLY:N	9:A2:101:BCL:HED2	2.32	0.45
1:BC:82:LEU:CD1	1:BC:93:THR:HG21	2.47	0.45
1:BC:84:ASP:OD2	1:BC:333:THR:HG21	2.16	0.45
1:BC:226:LEU:HD12	3:BM:192:ARG:CB	2.46	0.45
1:BC:233:PHE:O	1:BC:236:MET:N	2.50	0.45
1:BC:251:HIS:CG	1:BC:256:PHE:HD1	2.34	0.45
2:BL:38:VAL:O	2:BL:41:CYS:N	2.50	0.45
2:BL:40:PHE:O	2:BL:41:CYS:C	2.60	0.45
3:BM:116:LEU:HD11	3:BM:171:TRP:CZ2	2.52	0.45
3:BM:234:GLU:O	3:BM:237:GLN:N	2.50	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:BM:236:ASP:O	3:BM:240:HIS:N	2.49	0.45
3:BM:243:THR:HG22	4:BH:237:ASP:OD1	2.16	0.45
4:BH:193:VAL:HG23	4:BH:193:VAL:O	2.16	0.45
4:BH:249:TYR:C	4:BH:251:THR:H	2.24	0.45
4:BH:253:GLU:O	4:BH:254:ARG:C	2.59	0.45
5:BA:35:ILE:O	5:BA:36:HIS:C	2.59	0.45
9:BA:101:BCL:HMB1	9:B0:102:BCL:NB	2.31	0.45
5:BF:11:ILE:HG23	5:BF:12:TRP:CE3	2.51	0.45
9:BG:101:BCL:HBB3	9:BI:102:BCL:C1C	2.46	0.45
5:BI:16:ASP:HB2	5:BI:19:ARG:CB	2.46	0.45
5:BI:49:ASP:OD2	6:BJ:43:ARG:NH2	2.49	0.45
6:BJ:22:MET:O	6:BJ:25:MET:HB3	2.17	0.45
5:BQ:50:ASN:OD1	5:BQ:51:ILE:N	2.48	0.45
5:BW:11:ILE:C	5:BW:13:LEU:H	2.24	0.45
5:BY:45:ASN:O	5:BY:46:TRP:C	2.58	0.45
9:B2:101:BCL:CHB	9:B3:102:BCL:CMB	2.94	0.45
5:B5:28:GLN:O	5:B5:32:GLY:N	2.49	0.45
14:B5:103:CRT:H11	5:B7:21:LEU:HD13	1.98	0.45
1:AC:41:GLU:O	2:AL:172:GLN:NE2	2.50	0.45
1:AC:135:ARG:CG	1:AC:330:LEU:HA	2.38	0.45
2:AL:96:GLN:O	2:AL:100:ILE:N	2.45	0.45
2:AL:135:GLY:O	2:AL:138:LEU:N	2.44	0.45
2:AL:221:GLU:C	2:AL:223:THR:N	2.73	0.45
9:AL:301:BCL:HAC1	3:AM:197:TYR:OH	2.16	0.45
3:AM:27:ASN:ND2	5:AO:19:ARG:NH1	2.61	0.45
3:AM:99:PRO:O	3:AM:101:GLN:N	2.50	0.45
3:AM:268:TRP:CE2	4:AH:30:LEU:HD13	2.51	0.45
4:AH:196:PRO:O	4:AH:197:ILE:C	2.59	0.45
5:AA:35:ILE:HG21	9:AB:101:BCL:C4D	2.46	0.45
14:AB:102:CRT:H391	5:AD:36:HIS:CG	2.51	0.45
6:AJ:17:PHE:C	6:AJ:17:PHE:HD1	2.24	0.45
9:AK:102:BCL:CBB	9:AK:102:BCL:HMB1	2.47	0.45
6:AR:45:TRP:O	6:AR:46:LEU:HB2	2.17	0.45
5:AS:49:ASP:OD2	5:AS:50:ASN:ND2	2.48	0.45
9:AT:101:BCL:HMB3	9:AT:101:BCL:CBB	2.47	0.45
5:AW:34:LEU:HD21	14:AX:102:CRT:C40	2.41	0.45
5:AW:43:ASP:OD1	5:AW:44:LEU:N	2.50	0.45
5:AY:5:ASN:HA	6:AZ:18:HIS:NE2	2.31	0.45
6:A6:32:VAL:O	6:A6:35:ALA:HB3	2.16	0.45
5:A9:36:HIS:CD2	9:A0:102:BCL:HMD2	2.52	0.45
1:BC:50:ALA:O	1:BC:51:LEU:C	2.59	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:BL:82:TYR:CB	2:BL:85:ARG:HE	2.28	0.45
2:BL:120:LEU:O	2:BL:121:GLY:C	2.59	0.45
2:BL:129:ALA:HA	2:BL:247:LEU:CD1	2.46	0.45
2:BL:182:HIS:CB	2:BL:256:CYS:SG	3.04	0.45
2:BL:224:PHE:O	2:BL:228:ILE:HG13	2.17	0.45
2:BL:238:ILE:C	2:BL:240:ARG:H	2.23	0.45
3:BM:102:TYR:O	3:BM:103:GLY:C	2.58	0.45
3:BM:249:ALA:O	3:BM:259:ASN:OD1	2.34	0.45
4:BH:54:LYS:HE2	5:BD:23:SER:CB	2.45	0.45
4:BH:173:ASP:CG	4:BH:175:SER:H	2.23	0.45
4:BH:182:LEU:HD13	4:BH:195:LEU:CG	2.46	0.45
4:BH:197:ILE:HG23	4:BH:198:GLY:N	2.31	0.45
5:BK:32:GLY:N	9:BN:101:BCL:HED2	2.31	0.45
9:BO:102:BCL:HMB1	9:BO:102:BCL:CBB	2.47	0.45
5:BS:27:PHE:CG	5:BU:29:ILE:HD11	2.52	0.45
5:BU:6:ALA:HA	6:BV:15:LYS:NZ	2.30	0.45
5:BW:30:VAL:HG13	5:BW:31:LEU:N	2.30	0.45
5:BW:36:HIS:CE1	9:BX:101:BCL:CMD	2.99	0.45
5:BY:52:PRO:HD2	5:BY:55:TYR:CE2	2.51	0.45
5:B5:46:TRP:CD1	5:B5:47:LEU:N	2.85	0.45
6:B6:32:VAL:O	6:B6:35:ALA:HB3	2.15	0.45
5:B7:8:LEU:O	5:B7:11:ILE:HG22	2.16	0.45
5:B7:11:ILE:HG12	5:B7:15:LEU:HG	1.99	0.45
2:AL:10:TYR:HA	4:AH:112:GLY:CA	2.46	0.45
2:AL:10:TYR:HD1	4:AH:112:GLY:HA2	1.81	0.45
2:AL:192:ASN:HA	2:AL:245:LEU:HD12	1.99	0.45
2:AL:231:TYR:CE1	2:AL:233:ILE:HA	2.52	0.45
2:AL:261:GLY:C	2:AL:263:PHE:H	2.25	0.45
3:AM:61:ILE:HG23	3:AM:62:PHE:N	2.31	0.45
3:AM:155:PHE:O	3:AM:159:VAL:HG23	2.15	0.45
3:AM:226:VAL:CG2	3:AM:244:ALA:HA	2.46	0.45
3:AM:286:LEU:CD2	3:AM:290:VAL:HG21	2.46	0.45
4:AH:138:VAL:C	4:AH:140:LYS:H	2.23	0.45
4:AH:139:ALA:HA	4:AH:141:GLU:OE1	2.15	0.45
5:AA:21:LEU:HD23	5:A9:14:ILE:HG21	1.98	0.45
5:AD:46:TRP:HE1	9:AD:102:BCL:HHC	1.81	0.45
5:AI:43:ASP:O	5:AI:44:LEU:CB	2.65	0.45
5:AI:52:PRO:HB2	5:AI:55:TYR:CE2	2.52	0.45
6:AJ:20:ILE:HG23	6:AJ:21:PHE:N	2.32	0.45
9:AK:102:BCL:HHC	9:AK:102:BCL:OBB	2.16	0.45
6:AR:40:TRP:HH2	6:AR:46:LEU:HD11	1.81	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:AR:46:LEU:CB	6:AT:42:TYR:CE2	2.88	0.45
5:AU:43:ASP:OD1	5:AU:44:LEU:HD23	2.17	0.45
5:AU:43:ASP:CB	5:AW:47:LEU:HB3	2.46	0.45
5:AW:32:GLY:N	9:AX:101:BCL:HED2	2.32	0.45
6:AX:28:TRP:O	6:AX:31:LEU:N	2.50	0.45
6:AX:29:PHE:N	6:AX:29:PHE:CD1	2.84	0.45
5:AY:20:VAL:HA	5:AY:23:SER:OG	2.16	0.45
5:AY:43:ASP:HB2	5:A1:47:LEU:CD1	2.47	0.45
9:AZ:101:BCL:OBB	9:AZ:101:BCL:HHC	2.16	0.45
5:A5:43:ASP:OD2	5:A7:47:LEU:CA	2.58	0.45
14:A5:103:CRT:H31	9:A9:102:BCL:O2A	2.16	0.45
9:A6:101:BCL:CHC	9:A7:103:BCL:CBB	2.91	0.45
5:A7:28:GLN:O	5:A7:31:LEU:HB3	2.16	0.45
1:BC:21:LEU:HD23	1:BC:21:LEU:O	2.16	0.45
1:BC:276:VAL:CG1	1:BC:277:ARG:H	2.27	0.45
1:BC:314:VAL:HG12	1:BC:315:ASN:N	2.31	0.45
2:BL:129:ALA:CA	2:BL:247:LEU:HD11	2.46	0.45
3:BM:8:PHE:O	3:BM:10:ALA:N	2.49	0.45
4:BH:16:ILE:HD13	4:BH:16:ILE:C	2.41	0.45
4:BH:39:TYR:CD1	4:BH:40:PRO:HA	2.50	0.45
4:BH:132:LYS:HG3	4:BH:173:ASP:OD1	2.16	0.45
5:BA:46:TRP:CA	6:BB:43:ARG:HH12	2.27	0.45
9:BA:101:BCL:HED1	6:BB:31:LEU:HB3	1.99	0.45
6:BN:43:ARG:HB3	5:BO:55:TYR:HE2	1.78	0.45
5:BO:26:ALA:CA	5:BO:29:ILE:HG22	2.46	0.45
6:BP:20:ILE:HG23	6:BP:21:PHE:N	2.32	0.45
6:BR:21:PHE:CD1	6:BR:22:MET:N	2.85	0.45
9:BS:102:BCL:ND	9:BT:101:BCL:HMD3	2.31	0.45
14:BU:103:CRT:H342	9:BY:102:BCL:H3A	1.99	0.45
9:B2:101:BCL:HMB1	9:B2:101:BCL:CBB	2.47	0.45
9:B2:101:BCL:C2B	9:B3:102:BCL:C2B	2.94	0.45
9:AL:301:BCL:HMB3	9:AL:301:BCL:CBB	2.47	0.45
3:AM:8:PHE:HB3	3:AM:42:LYS:O	2.16	0.45
3:AM:70:ILE:HG22	3:AM:71:ILE:N	2.31	0.45
4:AH:47:GLU:HG3	5:AA:19:ARG:CB	2.46	0.45
5:AO:21:LEU:HD11	14:AP:102:CRT:H14	1.99	0.45
6:AR:45:TRP:CD2	9:AR:101:BCL:H2C	2.51	0.45
5:AW:50:ASN:CG	5:AW:51:ILE:N	2.74	0.45
5:AW:50:ASN:HA	5:AY:60:LYS:HA	1.99	0.45
6:AX:25:MET:HE2	14:AX:102:CRT:H21	1.98	0.45
5:A1:12:TRP:CD2	6:A2:17:PHE:HE2	2.34	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:A2:20:ILE:HD12	14:A2:102:CRT:H83	1.98	0.45
5:A5:11:ILE:CA	14:A5:103:CRT:H82	2.45	0.45
6:A8:43:ARG:HH21	5:A9:55:TYR:CB	2.26	0.45
6:A8:44:PRO:C	5:A9:52:PRO:HG2	2.41	0.45
6:A0:45:TRP:NE1	9:A0:102:BCL:H193	2.28	0.45
14:A0:101:CRT:H32	9:A0:102:BCL:CMA	2.46	0.45
1:BC:157:ARG:HH12	1:BC:318:LEU:CD2	2.30	0.45
1:BC:252:ASN:O	1:BC:254:ARG:N	2.49	0.45
1:BC:273:ILE:HA	7:BC:504:HEM:HBB2	1.99	0.45
1:BC:291:LEU:HD22	1:BC:295:ARG:HB2	1.98	0.45
2:BL:86:MET:HE2	2:BL:96:GLN:OE1	2.16	0.45
2:BL:216:LYS:HB3	2:BL:220:HIS:CD2	2.51	0.45
9:BL:303:BCL:HMB1	9:BL:303:BCL:HBB3	1.98	0.45
3:BM:85:GLN:HG3	3:BM:89:HIS:CD2	2.51	0.45
3:BM:107:PRO:HG2	3:BM:113:GLY:HA2	1.99	0.45
3:BM:172:ALA:C	3:BM:174:ALA:H	2.24	0.45
3:BM:184:ASP:O	3:BM:185:TRP:C	2.59	0.45
3:BM:265:ILE:HD13	13:BM:405:MQ8:H143	1.99	0.45
3:BM:267:ARG:HB3	4:BH:30:LEU:CD2	2.45	0.45
6:BB:38:LEU:HD23	6:BB:38:LEU:C	2.42	0.45
5:BF:49:ASP:OD1	5:BF:50:ASN:N	2.50	0.45
5:BI:46:TRP:CD1	5:BI:47:LEU:HD12	2.52	0.45
6:BN:38:LEU:HD23	6:BN:38:LEU:C	2.41	0.45
5:BQ:16:ASP:O	5:BQ:19:ARG:HB3	2.17	0.45
5:BQ:35:ILE:O	5:BQ:38:ILE:HG22	2.16	0.45
5:BQ:50:ASN:CG	5:BQ:51:ILE:N	2.75	0.45
6:BT:33:VAL:HG13	6:BT:34:ILE:N	2.32	0.45
5:BY:35:ILE:C	5:BY:37:MET:N	2.71	0.45
5:BY:45:ASN:O	5:BY:47:LEU:N	2.49	0.45
2:AL:109:TRP:O	2:AL:113:GLU:HG3	2.17	0.45
2:AL:138:LEU:C	2:AL:140:LEU:N	2.73	0.45
2:AL:156:PRO:HD3	2:AL:165:TRP:CD1	2.52	0.45
2:AL:196:LEU:CD2	3:AM:269:ALA:HB1	2.47	0.45
9:AL:303:BCL:O1A	3:AM:207:ALA:HB1	2.17	0.45
3:AM:12:GLN:CB	4:AH:145:ALA:HB2	2.45	0.45
3:AM:58:THR:HA	3:AM:61:ILE:HG22	1.98	0.45
3:AM:122:LEU:O	3:AM:157:TYR:OH	2.29	0.45
3:AM:274:VAL:O	3:AM:278:ILE:HG13	2.17	0.45
3:AM:296:LEU:O	3:AM:297:TRP:C	2.59	0.45
14:AM:406:CRT:H36	14:AM:406:CRT:H341	1.86	0.45
4:AH:206:ALA:C	4:AH:208:LYS:H	2.25	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:AA:101:BCL:HBB3	9:A0:102:BCL:C4B	2.47	0.45
5:AD:27:PHE:CZ	5:AF:29:ILE:HD12	2.51	0.45
6:AE:9:LEU:HD22	6:AE:13:GLU:CG	2.45	0.45
5:AF:36:HIS:NE2	9:AG:101:BCL:HMD3	2.32	0.45
6:AG:21:PHE:CD1	6:AG:22:MET:HA	2.52	0.45
6:AJ:17:PHE:C	6:AJ:20:ILE:HG22	2.41	0.45
5:AK:39:VAL:O	5:AK:39:VAL:HG12	2.17	0.45
9:AO:102:BCL:HMB1	9:AO:102:BCL:CBB	2.46	0.45
5:AQ:25:VAL:HG13	5:AQ:26:ALA:N	2.31	0.45
5:AS:20:VAL:HG12	9:AU:102:BCL:H202	1.97	0.45
9:AS:103:BCL:HED1	6:AT:31:LEU:O	2.17	0.45
5:AU:12:TRP:CZ2	6:AV:21:PHE:CD2	3.05	0.45
9:A1:102:BCL:C9	6:A2:28:TRP:HB2	2.47	0.45
9:A2:101:BCL:H3A	9:A2:101:BCL:HBA1	1.53	0.45
5:A9:46:TRP:CE2	9:A9:102:BCL:H2C	2.52	0.45
1:BC:51:LEU:O	1:BC:55:ALA:N	2.50	0.45
1:BC:130:MET:HE2	1:BC:133:LEU:CD2	2.47	0.45
1:BC:266:ARG:O	1:BC:267:THR:C	2.60	0.45
1:BC:313:ALA:O	1:BC:314:VAL:CG2	2.64	0.45
2:BL:18:ILE:O	2:BL:18:ILE:HG22	2.16	0.45
2:BL:240:ARG:NH2	3:BM:6:ASN:O	2.50	0.45
9:BL:303:BCL:HMB1	9:BL:303:BCL:CBB	2.47	0.45
3:BM:138:GLU:O	3:BM:141:GLY:N	2.46	0.45
9:BM:401:BCL:OBB	9:BM:401:BCL:HHC	2.15	0.45
4:BH:55:VAL:CG1	5:BD:19:ARG:HD3	2.46	0.45
4:BH:59:PRO:O	4:BH:60:ASP:C	2.57	0.45
5:BA:10:LYS:C	14:BA:102:CRT:H82	2.42	0.45
5:BA:18:ARG:O	5:BA:22:VAL:HG12	2.16	0.45
5:BD:9:TYR:CG	6:BE:15:LYS:HB2	2.52	0.45
5:BD:21:LEU:O	5:BD:25:VAL:HG23	2.16	0.45
5:BF:27:PHE:CE1	5:BI:29:ILE:CD1	2.91	0.45
5:BI:39:VAL:C	5:BI:41:SER:N	2.74	0.45
5:BK:39:VAL:O	5:BK:39:VAL:HG12	2.15	0.45
5:BO:8:LEU:CD2	6:BP:18:HIS:HE2	2.30	0.45
6:BP:18:HIS:O	6:BP:22:MET:HB2	2.17	0.45
14:BP:102:CRT:O2	5:BQ:33:LEU:HA	2.16	0.45
5:BS:12:TRP:HZ3	5:BS:20:VAL:HG11	1.82	0.45
5:BY:30:VAL:CA	5:BY:33:LEU:HG	2.44	0.45
5:BY:38:ILE:CD1	5:BY:39:VAL:HG23	2.47	0.45
6:B0:40:TRP:CZ3	6:B0:44:PRO:HA	2.51	0.45
1:AC:122:TYR:HA	1:AC:125:VAL:HG21	1.95	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AC:263:THR:HB	1:AC:264:PRO:CD	2.43	0.45
1:AC:313:ALA:O	1:AC:314:VAL:CG2	2.65	0.45
2:AL:12:VAL:HA	4:AH:111:PHE:HE2	1.82	0.45
2:AL:119:LYS:O	2:AL:121:GLY:N	2.50	0.45
3:AM:146:LEU:O	3:AM:148:TRP:N	2.50	0.45
4:AH:69:LEU:HB3	4:AH:70:PRO:CD	2.43	0.45
4:AH:125:LEU:CA	4:AH:131:PRO:HA	2.28	0.45
4:AH:169:ASP:H	4:AH:183:GLU:HB2	1.81	0.45
4:AH:173:ASP:OD1	4:AH:173:ASP:C	2.59	0.45
14:AG:102:CRT:C2M	5:AI:36:HIS:HB3	2.47	0.45
9:AI:102:BCL:HAC2	9:AJ:101:BCL:CBC	2.47	0.45
9:AO:102:BCL:CHA	9:AP:101:BCL:OBD	2.65	0.45
9:AX:101:BCL:CMC	9:AY:102:BCL:HBB1	2.47	0.45
5:AY:30:VAL:HA	5:AY:33:LEU:CG	2.46	0.45
6:A2:29:PHE:CE1	9:A2:101:BCL:H11	2.52	0.45
5:A3:19:ARG:HG2	5:A3:20:VAL:H	1.81	0.45
5:A3:28:GLN:OE1	9:A3:104:BCL:O1A	2.33	0.45
5:A5:21:LEU:CD1	9:A5:102:BCL:C14	2.88	0.45
5:A7:44:LEU:HD21	5:A7:46:TRP:CE3	2.35	0.45
6:A8:22:MET:HG3	6:A8:26:TYR:CE2	2.51	0.45
6:A8:30:GLY:O	6:A8:33:VAL:N	2.50	0.45
9:A9:102:BCL:HBB3	9:A9:102:BCL:HMB1	1.99	0.45
6:A0:21:PHE:HB2	14:A0:101:CRT:C16	2.47	0.45
1:BC:201:THR:N	1:BC:202:PRO:CD	2.80	0.45
1:BC:231:TRP:O	1:BC:232:THR:C	2.58	0.45
2:BL:63:SER:C	2:BL:65:LEU:N	2.74	0.45
2:BL:125:HIS:CD2	3:BM:224:LEU:HB3	2.52	0.45
2:BL:134:ILE:O	2:BL:138:LEU:HD13	2.17	0.45
2:BL:196:LEU:HD22	3:BM:216:PHE:CB	2.47	0.45
11:BL:304:UQ8:H15	11:BL:304:UQ8:H12A	1.81	0.45
11:BL:304:UQ8:H40	11:BL:304:UQ8:H37A	1.78	0.45
3:BM:8:PHE:HB3	3:BM:42:LYS:O	2.17	0.45
3:BM:11:VAL:HA	4:BH:148:ASP:CG	2.41	0.45
3:BM:170:SER:CB	3:BM:173:LYS:HD3	2.46	0.45
4:BH:69:LEU:HB3	4:BH:70:PRO:CD	2.33	0.45
4:BH:91:PRO:HA	4:BH:100:LEU:HD23	1.99	0.45
4:BH:123:CYS:SG	4:BH:231:VAL:C	3.00	0.45
4:BH:135:PRO:HA	4:BH:171:TRP:HA	1.99	0.45
4:BH:169:ASP:N	4:BH:183:GLU:HB2	2.32	0.45
4:BH:197:ILE:CG2	4:BH:198:GLY:N	2.80	0.45
4:BH:203:ASP:C	4:BH:205:LYS:H	2.25	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:BE:31:LEU:C	6:BE:34:ILE:HG22	2.41	0.45
5:BF:9:TYR:CD1	5:BF:10:LYS:N	2.85	0.45
9:BG:101:BCL:H3A	9:BG:101:BCL:HBA1	1.48	0.45
5:BK:45:ASN:O	5:BK:47:LEU:N	2.50	0.45
5:BQ:52:PRO:O	5:BQ:53:VAL:C	2.59	0.45
6:BR:22:MET:SD	6:BR:26:TYR:HE1	2.39	0.45
5:BU:13:LEU:HD22	6:BV:9:LEU:C	2.42	0.45
14:BU:103:CRT:C2M	5:BY:37:MET:HA	2.47	0.45
6:BV:28:TRP:O	6:BV:32:VAL:HG12	2.16	0.45
5:BW:7:ASN:HD22	5:BW:8:LEU:N	2.15	0.45
5:BW:10:LYS:C	14:BW:103:CRT:H82	2.41	0.45
5:BW:18:ARG:HH11	5:BW:18:ARG:HG2	1.82	0.45
5:BW:55:TYR:O	5:BW:59:GLY:HA3	2.17	0.45
5:BY:47:LEU:HD23	5:BY:47:LEU:HA	1.87	0.45
5:B7:9:TYR:CE2	5:B7:10:LYS:HE3	2.52	0.45
14:B0:101:CRT:H36	14:B0:101:CRT:H393	1.68	0.45
1:AC:179:LYS:N	1:AC:180:PRO:HD3	2.32	0.45
1:AC:234:GLY:O	1:AC:235:LEU:C	2.60	0.45
2:AL:22:LEU:HB2	5:A7:19:ARG:HG3	1.99	0.45
2:AL:93:GLY:O	2:AL:96:GLN:HB2	2.17	0.45
2:AL:194:LEU:HD23	2:AL:194:LEU:C	2.41	0.45
2:AL:276:LEU:O	2:AL:278:LEU:N	2.48	0.45
3:AM:233:ARG:NH1	4:AH:236:GLU:HG2	2.32	0.45
3:AM:257:GLY:O	3:AM:258:PHE:O	2.34	0.45
4:AH:5:ILE:HD12	5:AF:47:LEU:CD1	2.46	0.45
4:AH:31:ARG:HH21	4:AH:34:ASP:CB	2.30	0.45
4:AH:46:THR:HG22	4:AH:47:GLU:H	1.82	0.45
4:AH:154:MET:O	4:AH:167:VAL:HG13	2.17	0.45
5:AA:11:ILE:HD11	5:AD:21:LEU:HD21	1.99	0.45
6:AB:25:MET:HG2	6:AB:29:PHE:CE2	2.52	0.45
6:AG:28:TRP:CD1	6:AG:32:VAL:CG2	3.00	0.45
6:AJ:22:MET:HG3	6:AJ:26:TYR:HE1	1.82	0.45
5:AK:46:TRP:HA	5:AK:49:ASP:CG	2.41	0.45
6:AN:16:GLU:HB3	14:AN:102:CRT:H23	1.99	0.45
6:AP:21:PHE:O	6:AP:21:PHE:HD1	1.99	0.45
9:AP:101:BCL:HBB3	9:AQ:102:BCL:CHC	2.46	0.45
5:AS:40:LEU:O	5:AS:45:ASN:ND2	2.50	0.45
9:AU:102:BCL:C2D	9:AV:102:BCL:CMD	2.94	0.45
5:AY:44:LEU:HD13	6:AZ:43:ARG:NE	2.32	0.45
9:AY:102:BCL:HMD1	6:AZ:36:HIS:HA	1.98	0.45
5:A3:14:ILE:CD1	6:A6:17:PHE:CE2	2.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:A0:40:TRP:CZ3	6:A0:44:PRO:HA	2.51	0.45
1:BC:153:TYR:HB3	1:BC:323:MET:CE	2.34	0.45
1:BC:235:LEU:O	1:BC:237:MET:N	2.50	0.45
1:BC:269:ALA:O	1:BC:273:ILE:CD1	2.65	0.45
1:BC:282:ASN:HB3	1:BC:283:TYR:CE1	2.52	0.45
3:BM:40:LEU:HD23	3:BM:40:LEU:C	2.42	0.45
3:BM:130:TRP:HA	3:BM:150:PHE:CE2	2.52	0.45
4:BH:80:ARG:HG3	4:BH:80:ARG:O	2.17	0.45
4:BH:206:ALA:C	4:BH:208:LYS:N	2.73	0.45
5:BA:13:LEU:O	6:BB:9:LEU:HD13	2.17	0.45
6:BB:20:ILE:HG12	5:B9:7:ASN:CB	2.47	0.45
6:BB:36:HIS:CE1	9:BB:101:BCL:C4D	3.00	0.45
9:BD:102:BCL:CAC	9:BE:101:BCL:CBC	2.95	0.45
6:BE:22:MET:O	6:BE:26:TYR:HD1	2.00	0.45
5:BI:46:TRP:HE1	5:BI:47:LEU:CD1	2.29	0.45
5:BO:14:ILE:HA	6:BP:7:THR:N	2.31	0.45
9:BP:101:BCL:H3A	9:BP:101:BCL:HBA1	1.66	0.45
5:BS:4:MET:C	5:BS:6:ALA:H	2.24	0.45
9:BU:102:BCL:HBD	9:BV:101:BCL:OBD	2.15	0.45
6:BV:21:PHE:CB	14:BV:102:CRT:H11	2.47	0.45
6:B6:29:PHE:CE1	9:B6:101:BCL:C1	2.90	0.45
5:B9:12:TRP:CE3	5:B9:12:TRP:CA	3.00	0.45
1:AC:53:ILE:HA	1:AC:319:TYR:HE1	1.82	0.45
2:AL:184:LEU:CB	2:AL:252:TRP:NE1	2.78	0.45
2:AL:231:TYR:CE1	2:AL:233:ILE:N	2.85	0.45
4:AH:235:GLU:HG2	4:AH:235:GLU:H	1.52	0.45
9:AA:101:BCL:C3D	6:AB:35:ALA:HB1	2.47	0.45
9:AA:101:BCL:HHB	14:A0:101:CRT:H372	1.99	0.45
5:AD:52:PRO:C	5:AD:54:SER:N	2.73	0.45
5:AF:12:TRP:HA	5:AF:12:TRP:HE3	1.82	0.45
9:AO:102:BCL:HBC2	9:AO:102:BCL:CHD	2.46	0.45
6:AT:40:TRP:HZ3	6:AT:44:PRO:CA	2.30	0.45
5:AU:36:HIS:CE1	9:AU:102:BCL:NA	2.85	0.45
5:AW:16:ASP:OD1	5:AW:16:ASP:N	2.50	0.45
14:AW:102:CRT:H83	6:AZ:20:ILE:CD1	2.47	0.45
5:AY:27:PHE:HE1	5:AY:31:LEU:HD22	1.82	0.45
9:A3:103:BCL:H71	6:A4:28:TRP:CD2	2.51	0.45
9:A3:103:BCL:HMB1	9:A3:103:BCL:HBB3	1.99	0.45
9:A0:102:BCL:HBA1	9:A0:102:BCL:H3A	1.73	0.45
1:BC:156:HIS:O	1:BC:157:ARG:C	2.60	0.45
1:BC:280:ASN:CG	1:BC:304:ARG:HD2	2.41	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:BL:52:TRP:HE1	5:B9:38:ILE:HA	1.75	0.45
2:BL:125:HIS:CB	3:BM:221:ALA:HB1	2.47	0.45
3:BM:34:PRO:CG	3:BM:50:PRO:CD	2.90	0.45
3:BM:98:PRO:HA	3:BM:112:GLY:HA3	2.00	0.45
3:BM:146:LEU:C	3:BM:148:TRP:N	2.75	0.45
3:BM:214:LEU:O	3:BM:218:MET:HG2	2.17	0.45
4:BH:31:ARG:HA	4:BH:34:ASP:OD2	2.17	0.45
4:BH:145:ALA:HB3	4:BH:148:ASP:HB2	1.99	0.45
4:BH:259:LEU:HD11	5:B5:19:ARG:O	2.17	0.45
5:BD:28:GLN:C	9:BD:102:BCL:H11	2.42	0.45
9:BD:102:BCL:OBD	6:BE:32:VAL:HG23	2.17	0.45
5:BF:36:HIS:O	5:BF:40:LEU:HB2	2.17	0.45
9:BF:102:BCL:H62	6:BG:28:TRP:CZ2	2.52	0.45
5:BI:22:VAL:HA	5:BI:25:VAL:HG23	1.99	0.45
6:BJ:23:GLN:OE1	6:BJ:24:SER:N	2.50	0.45
5:BK:18:ARG:HD2	5:BK:19:ARG:N	2.31	0.45
5:BO:31:LEU:HD11	5:BO:35:ILE:HD11	1.98	0.45
5:BS:28:GLN:CB	9:BS:102:BCL:H43	2.47	0.45
6:BV:17:PHE:HB2	14:BV:102:CRT:C4	2.47	0.45
5:BW:32:GLY:CA	9:BX:101:BCL:HED2	2.47	0.45
5:B1:10:LYS:NZ	6:B4:20:ILE:HB	2.31	0.45
6:B0:37:LEU:HD23	9:B0:102:BCL:H202	1.99	0.45
1:AC:65:ALA:HB2	1:AC:89:GLU:OE1	2.17	0.44
1:AC:141:TRP:C	1:AC:143:ALA:N	2.75	0.44
1:AC:148:THR:OG1	1:AC:322:GLN:HG2	2.16	0.44
1:AC:236:MET:N	1:AC:239:ILE:HD12	2.32	0.44
1:AC:325:LYS:O	1:AC:325:LYS:HD3	2.18	0.44
2:AL:139:VAL:CG2	2:AL:258:LEU:HD13	2.47	0.44
3:AM:68:ILE:HG12	10:AM:403:BPH:H141	1.98	0.44
3:AM:222:THR:OG1	3:AM:252:TRP:HZ2	1.99	0.44
3:AM:242:GLY:CA	4:AH:117:PRO:HG3	2.48	0.44
14:AB:102:CRT:H33	14:AB:102:CRT:H5	1.63	0.44
5:AD:31:LEU:CG	9:AE:101:BCL:HED3	2.47	0.44
5:AF:8:LEU:HD23	6:AJ:20:ILE:CD1	2.44	0.44
6:AP:24:SER:O	6:AP:27:ALA:CB	2.65	0.44
5:AQ:17:PRO:O	5:AQ:20:VAL:HG22	2.16	0.44
6:AV:30:GLY:C	6:AV:33:VAL:HG12	2.42	0.44
6:A6:24:SER:O	6:A6:27:ALA:HB3	2.18	0.44
9:A6:101:BCL:HBA1	9:A6:101:BCL:H3A	1.79	0.44
6:A8:22:MET:HG3	6:A8:26:TYR:HE2	1.81	0.44
1:BC:35:TYR:CE1	1:BC:36:ARG:HG2	2.51	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BC:122:TYR:HA	1:BC:125:VAL:HG21	1.95	0.44
1:BC:253:THR:HG21	2:BL:168:ASN:HA	1.98	0.44
2:BL:243:LEU:O	2:BL:247:LEU:N	2.43	0.44
9:BL:301:BCL:HHC	9:BL:301:BCL:OBB	2.16	0.44
3:BM:33:ARG:O	3:BM:34:PRO:C	2.59	0.44
3:BM:163:ILE:HG23	3:BM:285:LEU:CD1	2.47	0.44
4:BH:37:GLU:C	4:BH:39:TYR:N	2.73	0.44
4:BH:184:VAL:CG2	4:BH:195:LEU:HB2	2.47	0.44
4:BH:222:VAL:O	4:BH:223:PRO:C	2.60	0.44
5:BA:46:TRP:HB3	6:BB:43:ARG:NH2	2.32	0.44
14:BA:102:CRT:C35	5:BD:31:LEU:HD21	2.47	0.44
6:BG:8:GLY:O	6:BG:9:LEU:HG	2.17	0.44
6:BG:21:PHE:O	6:BG:24:SER:OG	2.34	0.44
9:BG:101:BCL:HMB1	9:BI:102:BCL:NB	2.32	0.44
5:BK:12:TRP:CE3	5:BK:12:TRP:HA	2.52	0.44
6:BN:32:VAL:HG21	9:BN:101:BCL:CGA	2.47	0.44
6:BP:17:PHE:C	6:BP:20:ILE:HG22	2.41	0.44
6:BP:20:ILE:CG2	6:BP:21:PHE:N	2.80	0.44
6:BR:42:TYR:CE2	6:BR:43:ARG:HG3	2.52	0.44
5:BS:31:LEU:O	5:BS:35:ILE:HG12	2.17	0.44
5:BU:13:LEU:CD2	6:BV:9:LEU:HB2	2.46	0.44
14:BU:103:CRT:H2M1	5:BY:37:MET:CB	2.47	0.44
5:BW:4:MET:SD	6:BZ:27:ALA:HB2	2.57	0.44
5:B3:39:VAL:HA	5:B5:47:LEU:HD11	1.99	0.44
5:B7:17:PRO:O	5:B7:21:LEU:CG	2.62	0.44
1:AC:112:VAL:HG12	1:AC:113:PRO:CD	2.47	0.44
1:AC:130:MET:HE1	1:AC:284:ILE:CD1	2.47	0.44
1:AC:170:PRO:CG	1:AC:171:GLY:H	2.28	0.44
2:AL:97:ILE:O	2:AL:98:ILE:C	2.57	0.44
2:AL:225:PHE:C	2:AL:227:ASP:H	2.23	0.44
3:AM:90:PHE:C	3:AM:92:TRP:H	2.25	0.44
3:AM:176:PRO:CD	3:AM:185:TRP:CD1	3.00	0.44
3:AM:179:ILE:CG1	3:AM:180:PHE:H	2.22	0.44
3:AM:242:GLY:HA2	4:AH:117:PRO:HG3	2.00	0.44
3:AM:276:THR:O	3:AM:278:ILE:N	2.50	0.44
4:AH:172:VAL:HG23	4:AH:173:ASP:H	1.78	0.44
9:AA:101:BCL:HBB3	9:AA:101:BCL:HMB1	1.99	0.44
14:AB:102:CRT:H132	5:A9:11:ILE:HD11	1.98	0.44
5:AD:47:LEU:HD11	9:AD:102:BCL:HBB1	1.98	0.44
5:AF:28:GLN:CB	9:AF:102:BCL:H11	2.36	0.44
5:AI:39:VAL:CG1	5:AI:46:TRP:HZ3	2.29	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:AJ:101:BCL:CMB	9:AK:102:BCL:C1B	2.96	0.44
9:AK:102:BCL:CBD	9:AN:101:BCL:OBD	2.65	0.44
5:AU:43:ASP:HB2	5:AW:47:LEU:CB	2.45	0.44
5:AW:2:PHE:N	5:AW:2:PHE:CD1	2.85	0.44
5:AW:4:MET:O	5:AW:7:ASN:ND2	2.50	0.44
5:AW:49:ASP:O	5:AY:60:LYS:CB	2.65	0.44
6:AX:7:THR:OG1	6:AX:8:GLY:N	2.48	0.44
5:AY:48:ASP:OD1	5:AY:48:ASP:C	2.60	0.44
5:A1:46:TRP:CH2	9:A1:102:BCL:H2C	2.52	0.44
5:A3:36:HIS:O	5:A3:40:LEU:HD12	2.16	0.44
14:A5:103:CRT:H2M3	5:A9:36:HIS:CB	2.47	0.44
6:A6:37:LEU:HD23	6:A6:37:LEU:O	2.17	0.44
5:A7:30:VAL:HG13	5:A7:31:LEU:N	2.31	0.44
5:A9:31:LEU:HD23	9:A0:102:BCL:HED3	1.99	0.44
1:BC:70:PRO:HG2	1:BC:71:LYS:N	2.33	0.44
2:BL:23:PHE:HA	2:BL:25:PHE:CE2	2.52	0.44
2:BL:122:ILE:HG13	2:BL:123:GLY:O	2.17	0.44
2:BL:124:PHE:O	2:BL:125:HIS:C	2.60	0.44
2:BL:138:LEU:CD1	2:BL:138:LEU:N	2.81	0.44
3:BM:253:ARG:HH21	4:BH:41:LEU:HD11	1.82	0.44
9:BM:401:BCL:H3A	9:BM:401:BCL:HBA1	1.73	0.44
4:BH:182:LEU:HD13	4:BH:195:LEU:CD2	2.44	0.44
4:BH:185:GLU:O	4:BH:185:GLU:CD	2.60	0.44
6:BE:32:VAL:HG21	9:BE:101:BCL:CBA	2.42	0.44
5:BF:12:TRP:HZ2	6:BG:21:PHE:CD2	2.36	0.44
5:BF:30:VAL:HG13	5:BF:31:LEU:H	1.81	0.44
6:BJ:8:GLY:O	6:BJ:9:LEU:HD23	2.18	0.44
5:BO:18:ARG:HH11	5:BO:18:ARG:CB	2.29	0.44
5:BS:8:LEU:O	5:BS:11:ILE:HG13	2.17	0.44
5:B3:36:HIS:CD2	9:B4:101:BCL:HMD3	2.51	0.44
1:AC:36:ARG:NH1	2:AL:91:GLU:O	2.50	0.44
1:AC:199:PRO:C	1:AC:202:PRO:HD2	2.43	0.44
1:AC:243:LEU:N	1:AC:243:LEU:CD1	2.80	0.44
1:AC:286:PRO:O	1:AC:288:ASN:N	2.50	0.44
1:AC:293:ALA:C	1:AC:295:ARG:N	2.72	0.44
2:AL:47:VAL:HG23	10:AL:302:BPH:H7C2	1.98	0.44
2:AL:54:ALA:C	2:AL:68:TYR:HE1	2.25	0.44
2:AL:184:LEU:HB2	2:AL:252:TRP:NE1	2.33	0.44
2:AL:206:VAL:HG23	2:AL:207:THR:N	2.31	0.44
9:AL:303:BCL:CBB	9:AL:303:BCL:HMB1	2.47	0.44
3:AM:236:ASP:O	3:AM:239:THR:N	2.50	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:AA:8:LEU:HB3	6:AE:20:ILE:HG23	1.98	0.44
5:AA:27:PHE:O	5:AA:30:VAL:HG12	2.17	0.44
6:AB:15:LYS:O	6:AB:16:GLU:C	2.60	0.44
6:AN:40:TRP:CD1	6:AN:40:TRP:C	2.95	0.44
5:AQ:12:TRP:HA	5:AQ:12:TRP:HE3	1.83	0.44
5:AQ:36:HIS:CE1	9:AR:101:BCL:CMD	3.00	0.44
6:AR:40:TRP:C	6:AR:40:TRP:CD1	2.95	0.44
6:AT:29:PHE:CD1	6:AT:29:PHE:N	2.80	0.44
5:AU:8:LEU:O	5:AU:9:TYR:C	2.60	0.44
5:AU:45:ASN:H	5:AW:56:GLN:NE2	2.13	0.44
6:AV:46:LEU:HD13	6:AX:42:TYR:CZ	2.53	0.44
5:AW:38:ILE:HG23	5:AW:39:VAL:N	2.32	0.44
6:AX:10:THR:HG22	6:AX:11:ASP:N	2.32	0.44
5:AY:9:TYR:CD1	5:AY:9:TYR:C	2.94	0.44
6:A2:29:PHE:HD1	6:A2:29:PHE:H	1.58	0.44
14:A2:102:CRT:H243	9:A3:103:BCL:H18	2.00	0.44
5:A3:28:GLN:NE2	5:A3:28:GLN:HA	2.32	0.44
5:A5:44:LEU:HD12	5:A5:46:TRP:HE3	1.82	0.44
9:A7:103:BCL:C2D	9:A8:101:BCL:C2D	2.95	0.44
5:A9:12:TRP:CE3	5:A9:12:TRP:CA	3.00	0.44
1:BC:29:GLY:HA3	1:BC:44:TYR:CD2	2.53	0.44
1:BC:46:LYS:O	1:BC:48:GLN:N	2.50	0.44
1:BC:176:SER:OG	5:BS:42:THR:HG23	2.18	0.44
1:BC:276:VAL:CG2	1:BC:280:ASN:ND2	2.80	0.44
2:BL:97:ILE:O	2:BL:98:ILE:C	2.61	0.44
2:BL:240:ARG:NH2	3:BM:7:ILE:O	2.49	0.44
2:BL:261:GLY:O	2:BL:263:PHE:N	2.50	0.44
3:BM:264:SER:O	3:BM:265:ILE:C	2.61	0.44
4:BH:200:SER:HA	4:BH:211:VAL:HG22	1.98	0.44
4:BH:227:ASN:ND2	4:BH:228:PRO:CD	2.66	0.44
5:BA:17:PRO:HG2	5:BA:18:ARG:HD2	1.99	0.44
5:BA:39:VAL:HG11	9:BB:101:BCL:H3C	1.98	0.44
6:BB:30:GLY:O	6:BB:34:ILE:HG22	2.17	0.44
9:BB:101:BCL:CHC	9:BD:102:BCL:CBB	2.91	0.44
9:BE:101:BCL:C2B	9:BF:102:BCL:C2B	2.95	0.44
6:BG:46:LEU:O	5:BI:51:ILE:O	2.35	0.44
6:BJ:40:TRP:CD1	6:BJ:40:TRP:C	2.95	0.44
9:BJ:101:BCL:CBB	9:BJ:101:BCL:CMB	2.95	0.44
5:BO:34:LEU:C	5:BO:37:MET:HB2	2.42	0.44
9:BQ:104:BCL:HMA1	9:BS:102:BCL:HHB	1.99	0.44
5:BS:12:TRP:HE1	6:BT:18:HIS:CB	2.29	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:BU:18:ARG:O	5:BU:22:VAL:CG1	2.50	0.44
6:BV:44:PRO:O	5:BW:55:TYR:OH	2.30	0.44
5:BY:50:ASN:HD21	6:BZ:43:ARG:CZ	2.29	0.44
5:B1:12:TRP:CZ3	5:B1:20:VAL:HG21	2.42	0.44
5:B3:11:ILE:HA	14:B7:102:CRT:C8	2.41	0.44
6:B6:24:SER:O	6:B6:27:ALA:HB3	2.18	0.44
6:B6:37:LEU:O	6:B6:37:LEU:HD23	2.17	0.44
6:B0:21:PHE:CG	6:B0:22:MET:N	2.84	0.44
1:AC:187:SER:C	1:AC:189:THR:H	2.26	0.44
1:AC:252:ASN:O	1:AC:254:ARG:N	2.50	0.44
2:AL:114:VAL:O	2:AL:115:GLU:C	2.61	0.44
2:AL:131:SER:O	2:AL:134:ILE:N	2.51	0.44
2:AL:139:VAL:HG23	2:AL:143:VAL:CB	2.44	0.44
2:AL:240:ARG:O	2:AL:241:LEU:C	2.61	0.44
3:AM:281:GLY:O	3:AM:285:LEU:HB2	2.16	0.44
3:AM:284:ILE:HD11	9:AM:402:BCL:CAD	2.48	0.44
9:AM:402:BCL:HMB3	9:AM:402:BCL:HBB2	1.99	0.44
4:AH:55:VAL:HG12	5:AD:19:ARG:HD3	2.00	0.44
5:AA:11:ILE:HD13	14:AA:102:CRT:C9	2.47	0.44
6:AB:33:VAL:HG13	6:AB:34:ILE:N	2.31	0.44
14:AB:102:CRT:H2M1	5:AD:33:LEU:O	2.18	0.44
5:AI:35:ILE:HA	5:AI:38:ILE:CG2	2.47	0.44
6:AJ:17:PHE:CE1	14:AJ:102:CRT:H6	2.52	0.44
5:AU:49:ASP:OD2	6:AV:43:ARG:NH2	2.50	0.44
5:AU:51:ILE:HA	5:AU:53:VAL:N	2.31	0.44
14:AW:102:CRT:H392	5:AY:35:ILE:CD1	2.48	0.44
9:AX:101:BCL:H43	14:AX:102:CRT:H292	1.99	0.44
5:AY:9:TYR:CD1	5:AY:10:LYS:N	2.85	0.44
5:A1:10:LYS:HZ3	6:A4:20:ILE:HB	1.80	0.44
6:A2:42:TYR:CE1	6:A2:43:ARG:HG3	2.52	0.44
6:A4:34:ILE:CG2	6:A4:35:ALA:N	2.80	0.44
5:A5:16:ASP:HB2	5:A5:19:ARG:HG3	1.95	0.44
5:A5:27:PHE:HA	5:A5:30:VAL:HG12	1.99	0.44
5:A5:53:VAL:HA	5:A5:56:GLN:HE21	1.82	0.44
9:A6:101:BCL:NB	9:A7:103:BCL:HMB1	2.33	0.44
5:A7:4:MET:O	5:A7:4:MET:HE3	2.18	0.44
9:A7:103:BCL:O2D	9:A7:103:BCL:HAA1	2.17	0.44
1:BC:35:TYR:O	1:BC:36:ARG:C	2.58	0.44
1:BC:132:GLU:O	1:BC:136:ALA:CB	2.66	0.44
2:BL:63:SER:C	2:BL:65:LEU:H	2.25	0.44
2:BL:148:MET:O	2:BL:150:ALA:N	2.50	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:BL:160:LEU:CA	2:BL:163:LEU:HD13	2.46	0.44
2:BL:180:PRO:HA	2:BL:183:MET:SD	2.58	0.44
2:BL:195:ALA:HB3	3:BM:216:PHE:CE2	2.49	0.44
3:BM:2:PRO:HB3	4:BH:201:ARG:NH1	2.29	0.44
4:BH:107:MET:HG3	4:BH:242:TYR:HE1	1.82	0.44
4:BH:139:ALA:HA	4:BH:141:GLU:OE1	2.17	0.44
6:BB:11:ASP:HA	6:BB:14:ALA:HB3	1.98	0.44
14:BB:102:CRT:C8	5:B9:11:ILE:HG12	2.48	0.44
5:BD:29:ILE:HG23	5:BD:30:VAL:N	2.33	0.44
9:BD:102:BCL:CBB	9:BD:102:BCL:HMB1	2.47	0.44
5:BK:12:TRP:HA	5:BK:12:TRP:HE3	1.82	0.44
5:BO:29:ILE:CG2	5:BO:30:VAL:N	2.80	0.44
5:BQ:17:PRO:HB3	6:BR:17:PHE:CE2	2.53	0.44
5:BQ:43:ASP:OD1	5:BQ:44:LEU:N	2.50	0.44
5:BQ:49:ASP:OD1	5:BQ:50:ASN:N	2.43	0.44
6:BR:26:TYR:O	6:BR:27:ALA:C	2.61	0.44
6:BX:28:TRP:O	6:BX:31:LEU:N	2.50	0.44
5:B1:10:LYS:C	14:B1:103:CRT:H82	2.39	0.44
5:B3:36:HIS:NE2	9:B4:101:BCL:CMD	2.75	0.44
6:B4:21:PHE:CD1	6:B4:22:MET:N	2.86	0.44
6:B4:34:ILE:CG2	6:B4:35:ALA:N	2.80	0.44
6:B4:38:LEU:HD23	6:B4:38:LEU:C	2.42	0.44
6:B8:9:LEU:HD22	6:B8:13:GLU:CD	2.43	0.44
1:AC:70:PRO:HB2	1:AC:71:LYS:HD2	2.00	0.44
1:AC:211:ARG:HD3	3:AM:317:TYR:CZ	2.51	0.44
2:AL:4:LEU:CD1	3:AM:250:LEU:HD12	2.40	0.44
2:AL:63:SER:C	2:AL:65:LEU:N	2.73	0.44
2:AL:181:ALA:C	2:AL:183:MET:N	2.76	0.44
9:AL:303:BCL:HMB1	9:AL:303:BCL:HBB3	2.00	0.44
3:AM:149:ALA:O	3:AM:152:ALA:N	2.51	0.44
3:AM:151:ALA:O	3:AM:154:ILE:N	2.50	0.44
3:AM:210:TYR:O	3:AM:213:ALA:N	2.50	0.44
3:AM:256:MET:SD	13:AM:405:MQ8:H151	2.57	0.44
3:AM:314:VAL:HG12	3:AM:315:ASN:N	2.33	0.44
9:AA:101:BCL:HBB3	9:A0:102:BCL:C1C	2.47	0.44
14:AA:102:CRT:H83	6:AE:20:ILE:CD1	2.34	0.44
6:AB:28:TRP:O	6:AB:31:LEU:N	2.46	0.44
6:AB:43:ARG:HD3	5:AD:55:TYR:CE1	2.53	0.44
6:AG:40:TRP:CZ3	6:AG:46:LEU:HG	2.52	0.44
9:AI:102:BCL:HAC2	9:AJ:101:BCL:HAC1	1.99	0.44
6:AJ:15:LYS:O	6:AJ:18:HIS:HB3	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:AR:28:TRP:CE3	6:AR:28:TRP:HA	2.53	0.44
5:AS:35:ILE:C	5:AS:37:MET:N	2.70	0.44
6:AT:22:MET:O	6:AT:25:MET:N	2.49	0.44
6:AT:45:TRP:HD1	6:AT:46:LEU:N	2.15	0.44
5:AU:20:VAL:HG11	9:AW:101:BCL:C20	2.48	0.44
9:A3:103:BCL:H2	6:A4:28:TRP:CZ2	2.52	0.44
6:A4:21:PHE:CD1	6:A4:22:MET:N	2.86	0.44
6:A6:17:PHE:CD2	14:A7:102:CRT:H42	2.53	0.44
9:A6:101:BCL:C1B	9:A7:103:BCL:CMB	2.96	0.44
5:A7:13:LEU:O	6:A8:7:THR:HB	2.17	0.44
5:A7:46:TRP:CH2	9:A7:103:BCL:H2C	2.52	0.44
1:BC:126:VAL:HG12	1:BC:287:LEU:HB3	1.99	0.44
1:BC:135:ARG:HB3	1:BC:332:LYS:N	2.32	0.44
1:BC:271:TYR:O	1:BC:274:ARG:HB2	2.17	0.44
2:BL:165:TRP:HE3	2:BL:166:VAL:HG23	1.83	0.44
3:BM:149:ALA:O	3:BM:150:PHE:C	2.60	0.44
3:BM:176:PRO:HG2	3:BM:182:HIS:HA	1.99	0.44
3:BM:215:LEU:O	3:BM:218:MET:N	2.41	0.44
5:BD:17:PRO:O	5:BD:21:LEU:CB	2.66	0.44
5:BD:30:VAL:HG13	5:BD:31:LEU:N	2.32	0.44
9:BD:102:BCL:HBB3	9:BD:102:BCL:HMB1	1.98	0.44
5:BK:51:ILE:HA	5:BK:52:PRO:O	2.17	0.44
5:BO:38:ILE:HG13	5:BO:39:VAL:HG23	2.00	0.44
9:BP:101:BCL:O1D	9:BP:101:BCL:H2A	2.17	0.44
5:BQ:2:PHE:O	5:BQ:5:ASN:HB3	2.16	0.44
5:BQ:44:LEU:CD1	5:BQ:46:TRP:HE3	2.23	0.44
9:BQ:103:BCL:HMD2	9:BQ:104:BCL:C1D	2.48	0.44
5:BW:45:ASN:O	5:BW:49:ASP:CB	2.65	0.44
6:BX:29:PHE:N	6:BX:29:PHE:HD1	2.15	0.44
9:BY:102:BCL:HMB1	9:BY:102:BCL:CBB	2.47	0.44
6:B2:17:PHE:CD1	14:B2:102:CRT:C9	2.84	0.44
5:B3:24:ILE:O	5:B3:27:PHE:N	2.50	0.44
5:B5:49:ASP:OD1	5:B5:50:ASN:N	2.48	0.44
14:B0:101:CRT:H241	14:B0:101:CRT:H26	1.81	0.44
1:AC:46:LYS:C	1:AC:48:GLN:N	2.76	0.44
1:AC:297:GLY:H	1:AC:301:ASP:H	1.66	0.44
2:AL:192:ASN:N	2:AL:245:LEU:HD13	2.33	0.44
2:AL:252:TRP:O	2:AL:254:ALA:N	2.50	0.44
3:AM:63:PHE:CE1	5:AQ:30:VAL:HA	2.52	0.44
4:AH:11:ALA:O	4:AH:14:ILE:HG13	2.18	0.44
4:AH:59:PRO:O	4:AH:60:ASP:C	2.61	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:AD:102:BCL:HAC2	9:AE:101:BCL:CBC	2.48	0.44
5:AI:31:LEU:HB3	9:AJ:101:BCL:HED2	1.99	0.44
5:AO:11:ILE:HD13	14:AR:102:CRT:H132	1.98	0.44
5:AO:31:LEU:HD11	14:AP:102:CRT:H35	1.99	0.44
5:AO:43:ASP:HA	5:AQ:48:ASP:CB	2.35	0.44
5:AO:43:ASP:OD2	5:AQ:47:LEU:HD22	2.17	0.44
9:AO:102:BCL:CBC	9:AP:101:BCL:HHD	2.41	0.44
5:AQ:35:ILE:HD11	14:AR:102:CRT:H372	2.00	0.44
9:AR:101:BCL:HMB1	9:AR:101:BCL:HBB3	1.99	0.44
5:AU:14:ILE:HD11	14:AX:102:CRT:H23	2.00	0.44
6:AZ:20:ILE:HG23	6:AZ:21:PHE:N	2.32	0.44
9:AZ:101:BCL:CMB	9:A1:102:BCL:C1B	2.95	0.44
5:A1:12:TRP:CE3	5:A1:12:TRP:CA	3.01	0.44
5:A1:51:ILE:HB	5:A1:52:PRO:CA	2.46	0.44
6:A2:31:LEU:C	6:A2:34:ILE:HG22	2.43	0.44
9:A2:101:BCL:C1B	9:A3:103:BCL:CMB	2.96	0.44
5:A5:25:VAL:CG1	9:A5:102:BCL:H192	2.36	0.44
1:BC:21:LEU:HD21	2:BL:263:PHE:CE1	2.53	0.44
1:BC:141:TRP:CH2	1:BC:275:HIS:HA	2.52	0.44
1:BC:226:LEU:HD11	3:BM:189:PHE:HA	2.00	0.44
1:BC:227:LYS:O	1:BC:230:GLU:HB3	2.17	0.44
1:BC:258:ASP:O	1:BC:261:GLN:HB2	2.17	0.44
2:BL:6:PHE:CD1	3:BM:246:GLU:HG3	2.53	0.44
2:BL:48:LEU:HA	2:BL:51:VAL:CG2	2.46	0.44
2:BL:56:ILE:O	2:BL:66:GLN:HG3	2.17	0.44
2:BL:109:TRP:CD1	2:BL:109:TRP:C	2.96	0.44
3:BM:68:ILE:HG21	10:BM:403:BPH:H141	1.99	0.44
3:BM:200:PRO:CA	3:BM:203:MET:HG2	2.45	0.44
4:BH:69:LEU:HD13	4:BH:76:VAL:HG23	1.99	0.44
4:BH:80:ARG:O	4:BH:81:VAL:C	2.61	0.44
4:BH:249:TYR:C	4:BH:251:THR:N	2.75	0.44
5:BF:9:TYR:CD1	6:BG:15:LYS:HG2	2.52	0.44
6:BG:28:TRP:O	6:BG:29:PHE:C	2.60	0.44
5:BK:36:HIS:NE2	9:BN:101:BCL:HMD3	2.30	0.44
9:BK:102:BCL:H92	14:BN:102:CRT:C18	2.47	0.44
5:BO:45:ASN:O	5:BO:49:ASP:CG	2.61	0.44
14:BO:103:CRT:H27	5:BQ:28:GLN:NE2	2.32	0.44
6:BP:24:SER:O	6:BP:27:ALA:CB	2.66	0.44
5:BQ:19:ARG:NH1	15:BQ:101:PEF:N	2.66	0.44
6:BX:32:VAL:O	6:BX:36:HIS:HB2	2.18	0.44
5:BY:13:LEU:HD23	5:BY:13:LEU:HA	1.85	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:B1:11:ILE:HG12	14:B1:103:CRT:H81	1.98	0.44
6:B2:10:THR:HG23	6:B2:11:ASP:N	2.32	0.44
9:B2:101:BCL:HMB3	9:B3:102:BCL:CHB	2.48	0.44
5:B3:29:ILE:HB	9:B3:102:BCL:H42	2.00	0.44
14:B5:103:CRT:C7	6:B8:17:PHE:HZ	2.30	0.44
6:B8:30:GLY:O	6:B8:33:VAL:N	2.50	0.44
6:B0:29:PHE:CE1	9:B0:102:BCL:H72	2.53	0.44
1:AC:293:ALA:O	1:AC:296:LYS:N	2.32	0.44
2:AL:23:PHE:CE1	5:A9:22:VAL:HG21	2.53	0.44
2:AL:264:TRP:CZ3	2:AL:271:TRP:HD1	2.35	0.44
3:AM:8:PHE:O	3:AM:9:THR:C	2.61	0.44
4:AH:144:ILE:HG13	4:AH:150:ASP:OD2	2.16	0.44
9:AA:101:BCL:C2B	9:A0:102:BCL:C2B	2.95	0.44
9:AB:101:BCL:HMA1	9:AD:102:BCL:HHB	2.00	0.44
5:AI:44:LEU:CD1	5:AI:46:TRP:HE3	2.31	0.44
6:AJ:40:TRP:HA	6:AJ:44:PRO:HA	2.00	0.44
9:AK:102:BCL:H193	9:AK:102:BCL:H111	1.99	0.44
5:AU:32:GLY:HA3	9:AU:102:BCL:O1A	2.18	0.44
6:AV:10:THR:CG2	6:AV:11:ASP:N	2.80	0.44
6:AV:45:TRP:CZ3	9:AV:102:BCL:HAC2	2.52	0.44
5:AW:35:ILE:HA	5:AW:38:ILE:CG2	2.47	0.44
14:AX:102:CRT:H181	14:AX:102:CRT:H20	1.80	0.44
5:AY:15:LEU:CG	5:A1:21:LEU:HD21	2.46	0.44
9:AY:102:BCL:HMD2	6:AZ:36:HIS:HD2	1.82	0.44
6:AZ:45:TRP:CE3	9:AZ:101:BCL:CAC	2.88	0.44
5:A3:51:ILE:HA	5:A3:52:PRO:C	2.43	0.44
5:A5:30:VAL:HG13	5:A5:31:LEU:N	2.32	0.44
5:A7:7:ASN:HA	5:A7:10:LYS:HZ2	1.81	0.44
5:A9:49:ASP:OD2	5:A9:50:ASN:OD1	2.35	0.44
1:BC:112:VAL:CG1	1:BC:113:PRO:HD2	2.46	0.44
1:BC:226:LEU:H	3:BM:173:LYS:CE	2.31	0.44
1:BC:326:ASP:C	1:BC:327:TYR:CD1	2.96	0.44
2:BL:79:ASP:C	2:BL:81:SER:N	2.74	0.44
2:BL:196:LEU:HD13	3:BM:216:PHE:CB	2.36	0.44
2:BL:204:LEU:HD13	3:BM:267:ARG:HG3	1.99	0.44
3:BM:53:LEU:HD11	3:BM:58:THR:HA	2.00	0.44
3:BM:73:PHE:O	3:BM:76:LEU:N	2.47	0.44
3:BM:84:PHE:HD1	3:BM:84:PHE:H	1.66	0.44
3:BM:168:MET:HE3	3:BM:289:THR:HG22	2.00	0.44
3:BM:229:PHE:CE2	3:BM:247:ARG:NE	2.86	0.44
3:BM:279:THR:HA	3:BM:282:ILE:HG13	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:BH:259:LEU:CD2	5:B5:19:ARG:HB3	2.48	0.44
5:BA:12:TRP:CD2	6:BB:17:PHE:HE2	2.36	0.44
5:BA:21:LEU:O	5:BA:25:VAL:HG23	2.18	0.44
5:BA:31:LEU:HD11	14:BB:102:CRT:H35	2.00	0.44
9:BA:101:BCL:HBB3	9:BA:101:BCL:HMB1	2.00	0.44
5:BD:9:TYR:HE1	6:BE:11:ASP:CB	2.30	0.44
6:BJ:28:TRP:NE1	6:BJ:32:VAL:CG2	2.80	0.44
9:BJ:101:BCL:H3A	9:BJ:101:BCL:HBA1	1.60	0.44
5:BK:27:PHE:CD1	5:BK:27:PHE:C	2.94	0.44
6:BN:18:HIS:NE2	6:BN:22:MET:HE2	2.32	0.44
5:BO:9:TYR:HA	6:BP:18:HIS:CD2	2.51	0.44
6:BP:40:TRP:HZ3	6:BP:45:TRP:H	1.65	0.44
5:BQ:27:PHE:CD1	5:BQ:27:PHE:C	2.95	0.44
6:BT:31:LEU:HA	6:BT:34:ILE:CD1	2.48	0.44
6:BV:10:THR:CG2	6:BV:11:ASP:H	2.27	0.44
9:BW:102:BCL:HMB1	9:BW:102:BCL:CBB	2.48	0.44
9:BW:102:BCL:CBD	9:BX:101:BCL:OBD	2.66	0.44
5:BY:36:HIS:NE2	9:BZ:101:BCL:CMD	2.80	0.44
5:B5:18:ARG:HB2	5:B5:19:ARG:CZ	2.48	0.44
6:B0:45:TRP:O	6:B0:46:LEU:CB	2.62	0.44
14:B0:101:CRT:H30	9:B0:102:BCL:HMA2	1.99	0.44
1:AC:212:ILE:HG21	1:AC:229:ALA:HB2	2.00	0.44
2:AL:22:LEU:HD23	2:AL:23:PHE:CZ	2.53	0.44
2:AL:44:LEU:O	2:AL:48:LEU:HB2	2.18	0.44
2:AL:93:GLY:O	2:AL:96:GLN:N	2.50	0.44
2:AL:95:TRP:HE3	2:AL:96:GLN:N	2.16	0.44
2:AL:156:PRO:HG2	2:AL:162:HIS:HA	2.00	0.44
2:AL:186:ILE:HD12	17:AL:403:HOH:O	2.17	0.44
2:AL:281:TRP:CG	3:AM:88:LYS:HB2	2.53	0.44
3:AM:41:GLY:C	3:AM:43:ILE:H	2.24	0.44
3:AM:131:VAL:O	3:AM:133:THR:N	2.50	0.44
3:AM:232:ASP:OD2	4:AH:180:ARG:NH2	2.50	0.44
3:AM:243:THR:HA	3:AM:246:GLU:HB2	2.00	0.44
3:AM:271:TRP:NE1	4:AH:26:LEU:HD11	2.33	0.44
4:AH:189:ASN:O	4:AH:191:LYS:HG3	2.18	0.44
5:AD:14:ILE:N	5:AD:14:ILE:CD1	2.81	0.44
5:AF:31:LEU:HD21	14:AG:102:CRT:H32	1.99	0.44
9:AG:101:BCL:HMB2	9:AI:102:BCL:CHB	2.48	0.44
5:AI:4:MET:SD	6:AN:23:GLN:OE1	2.76	0.44
5:AI:27:PHE:CE1	5:AI:31:LEU:HD22	2.53	0.44
5:AO:4:MET:HG3	6:AR:23:GLN:HB3	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:AO:34:LEU:O	5:AO:38:ILE:HG23	2.18	0.44
6:AX:46:LEU:HB3	6:AZ:42:TYR:OH	2.17	0.44
5:A1:26:ALA:O	5:A1:29:ILE:CG2	2.66	0.44
5:A3:40:LEU:HD21	5:A3:46:TRP:CZ2	2.52	0.44
14:A7:102:CRT:C21	14:A7:102:CRT:C24	2.95	0.44
1:BC:40:MET:HG2	1:BC:251:HIS:O	2.18	0.44
1:BC:235:LEU:C	1:BC:237:MET:H	2.26	0.44
1:BC:316:LYS:HB3	1:BC:320:GLY:H	1.82	0.44
2:BL:3:MET:O	4:BH:41:LEU:HG	2.17	0.44
2:BL:117:CYS:SG	2:BL:122:ILE:HD11	2.58	0.44
2:BL:193:CYS:O	2:BL:193:CYS:SG	2.76	0.44
2:BL:281:TRP:CG	3:BM:88:LYS:HB2	2.53	0.44
3:BM:132:ARG:HH11	3:BM:132:ARG:CG	2.30	0.44
3:BM:241:ARG:HB2	4:BH:37:GLU:OE1	2.18	0.44
5:BF:44:LEU:HD12	5:BF:46:TRP:HE3	1.82	0.44
5:BI:45:ASN:OD1	5:BI:48:ASP:OD1	2.35	0.44
5:BO:25:VAL:HG12	9:BO:102:BCL:C4	2.45	0.44
6:BP:32:VAL:O	6:BP:35:ALA:HB3	2.18	0.44
9:BQ:104:BCL:H2C	6:BR:45:TRP:CD2	2.53	0.44
5:BS:5:ASN:HA	5:BS:8:LEU:CG	2.46	0.44
5:BU:19:ARG:HB2	5:BU:19:ARG:HH21	1.82	0.44
5:BU:45:ASN:CG	5:BU:48:ASP:OD1	2.61	0.44
14:BU:103:CRT:H181	14:BU:103:CRT:H20	1.80	0.44
5:B3:29:ILE:HG23	5:B3:30:VAL:N	2.33	0.44
5:B5:38:ILE:O	5:B5:41:SER:HB3	2.18	0.44
6:B6:37:LEU:C	6:B6:37:LEU:CD2	2.91	0.44
1:AC:53:ILE:O	1:AC:55:ALA:N	2.51	0.44
1:AC:86:SER:O	1:AC:87:VAL:C	2.58	0.44
1:AC:302:PRO:O	1:AC:302:PRO:CG	2.65	0.44
2:AL:119:LYS:HD3	3:AM:254:TRP:HZ3	1.83	0.44
2:AL:189:PHE:CE2	9:AL:301:BCL:HMA2	2.53	0.44
2:AL:196:LEU:HD13	2:AL:197:SER:N	2.33	0.44
2:AL:253:SER:HB2	9:AL:301:BCL:CAA	2.48	0.44
3:AM:132:ARG:HH11	3:AM:132:ARG:CG	2.31	0.44
3:AM:163:ILE:HG23	3:AM:285:LEU:CD1	2.48	0.44
9:AA:101:BCL:CBC	9:AB:101:BCL:HHD	2.48	0.44
5:AD:51:ILE:HA	5:AD:53:VAL:N	2.33	0.44
9:AD:102:BCL:HMB1	9:AD:102:BCL:CBB	2.47	0.44
6:AE:10:THR:CG2	6:AE:11:ASP:N	2.81	0.44
6:AG:22:MET:HG3	6:AG:26:TYR:CZ	2.53	0.44
9:AI:102:BCL:H62	6:AJ:28:TRP:CZ3	2.53	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:AI:102:BCL:H62	6:AJ:28:TRP:CH2	2.52	0.44
6:AJ:34:ILE:O	6:AJ:38:LEU:HB2	2.17	0.44
6:AP:39:ALA:O	6:AP:42:TYR:N	2.49	0.44
5:AS:34:LEU:O	5:AS:38:ILE:HG22	2.17	0.44
14:AS:104:CRT:H81	14:AS:104:CRT:H10	1.81	0.44
6:AT:10:THR:H	6:AT:13:GLU:CD	2.26	0.44
5:AU:59:GLY:O	5:AU:60:LYS:O	2.35	0.44
6:AV:30:GLY:O	6:AV:34:ILE:CG1	2.66	0.44
6:AX:22:MET:O	6:AX:26:TYR:CD2	2.71	0.44
5:A1:54:SER:O	5:A1:57:ALA:HB3	2.18	0.44
9:A3:104:BCL:H3A	9:A3:104:BCL:HBA1	1.64	0.44
6:A8:32:VAL:O	6:A8:33:VAL:C	2.60	0.44
6:A8:42:TYR:CG	6:A8:43:ARG:N	2.85	0.44
6:A0:27:ALA:O	6:A0:31:LEU:HG	2.18	0.44
1:BC:153:TYR:O	1:BC:157:ARG:HG2	2.18	0.44
1:BC:267:THR:C	1:BC:269:ALA:H	2.24	0.44
1:BC:316:LYS:CB	1:BC:320:GLY:H	2.31	0.44
2:BL:95:TRP:CE3	2:BL:96:GLN:HA	2.53	0.44
2:BL:194:LEU:O	2:BL:195:ALA:C	2.61	0.44
2:BL:194:LEU:C	2:BL:194:LEU:CD2	2.91	0.44
2:BL:203:ILE:C	2:BL:205:SER:N	2.75	0.44
2:BL:252:TRP:HA	2:BL:252:TRP:HE3	1.81	0.44
3:BM:204:LEU:HD11	4:BH:19:PHE:CZ	2.52	0.44
14:BA:102:CRT:H32	5:BD:31:LEU:CD2	2.48	0.44
6:BB:44:PRO:C	5:BD:52:PRO:HG3	2.43	0.44
5:BD:8:LEU:O	5:BD:9:TYR:C	2.60	0.44
5:BD:43:ASP:CG	5:BD:44:LEU:N	2.75	0.44
5:BI:49:ASP:CG	5:BI:50:ASN:N	2.76	0.44
5:BK:49:ASP:CG	5:BK:50:ASN:N	2.76	0.44
5:BS:51:ILE:CA	5:BS:52:PRO:C	2.91	0.44
14:BV:102:CRT:C39	5:BW:36:HIS:HB2	2.15	0.44
5:B1:11:ILE:HG12	14:B1:103:CRT:H10	2.00	0.44
14:B1:103:CRT:H32	5:B3:31:LEU:HD21	1.99	0.44
6:B2:17:PHE:HE1	14:B2:102:CRT:H9	1.75	0.44
6:B2:21:PHE:O	6:B2:22:MET:C	2.61	0.44
5:B3:43:ASP:OD1	5:B3:44:LEU:N	2.51	0.44
6:B8:32:VAL:O	6:B8:33:VAL:C	2.60	0.44
5:B9:12:TRP:CD1	6:B0:18:HIS:HB2	2.53	0.44
5:B9:15:LEU:HB3	5:B9:20:VAL:HG21	1.99	0.44
6:B0:34:ILE:C	6:B0:34:ILE:CD1	2.91	0.44
1:AC:82:LEU:CD1	1:AC:93:THR:HG21	2.48	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AC:121:ILE:CG2	1:AC:123:THR:CG2	2.96	0.43
1:AC:280:ASN:CG	1:AC:304:ARG:HB3	2.35	0.43
2:AL:137:TYR:HB2	9:AL:301:BCL:H52	2.00	0.43
2:AL:139:VAL:CG2	2:AL:258:LEU:HD22	2.48	0.43
2:AL:164:ASP:O	2:AL:166:VAL:N	2.51	0.43
3:AM:164:ARG:O	3:AM:167:MET:N	2.50	0.43
3:AM:180:PHE:O	3:AM:184:ASP:OD1	2.36	0.43
3:AM:280:ALA:C	3:AM:282:ILE:N	2.76	0.43
3:AM:286:LEU:HD22	4:AH:12:ALA:HB2	2.00	0.43
4:AH:107:MET:HG3	4:AH:242:TYR:CE1	2.53	0.43
4:AH:171:TRP:NE1	4:AH:194:LEU:HD21	2.33	0.43
4:AH:203:ASP:O	4:AH:205:LYS:N	2.51	0.43
6:AB:11:ASP:HA	6:AB:14:ALA:HB3	1.99	0.43
5:AD:46:TRP:NE1	5:AD:47:LEU:HD22	2.33	0.43
6:AG:46:LEU:CB	6:AJ:42:TYR:CZ	2.86	0.43
5:AQ:43:ASP:CB	5:AS:47:LEU:HB3	2.47	0.43
5:AU:9:TYR:HE1	6:AV:14:ALA:HB3	1.82	0.43
5:AW:9:TYR:O	5:AW:11:ILE:N	2.51	0.43
6:AZ:22:MET:HG3	6:AZ:26:TYR:CE1	2.41	0.43
6:A2:45:TRP:HE1	9:A2:101:BCL:HHC	1.82	0.43
6:A4:38:LEU:HD23	6:A4:38:LEU:C	2.42	0.43
5:A5:14:ILE:O	5:A5:14:ILE:HG22	2.18	0.43
5:A5:16:ASP:HB2	5:A5:19:ARG:HH21	1.82	0.43
9:A0:102:BCL:CBB	9:A0:102:BCL:H18	2.47	0.43
1:BC:123:THR:OG1	1:BC:124:LYS:N	2.50	0.43
2:BL:56:ILE:O	2:BL:60:GLY:C	2.60	0.43
2:BL:261:GLY:C	2:BL:263:PHE:H	2.26	0.43
3:BM:84:PHE:HZ	14:BV:102:CRT:H401	1.83	0.43
3:BM:248:ALA:C	3:BM:250:LEU:N	2.77	0.43
4:BH:184:VAL:O	4:BH:193:VAL:HG22	2.18	0.43
5:BA:27:PHE:O	5:BA:30:VAL:HG12	2.18	0.43
5:BA:55:TYR:CE1	5:B9:44:LEU:CB	2.94	0.43
5:BD:43:ASP:O	5:BD:45:ASN:N	2.50	0.43
6:BG:24:SER:O	6:BG:27:ALA:N	2.50	0.43
6:BJ:45:TRP:HD1	6:BJ:46:LEU:HD23	1.83	0.43
6:BN:18:HIS:O	6:BN:22:MET:CB	2.66	0.43
5:BO:11:ILE:O	5:BO:11:ILE:HG22	2.17	0.43
6:BP:46:LEU:O	5:BQ:51:ILE:O	2.35	0.43
5:BS:29:ILE:CG2	5:BS:30:VAL:N	2.79	0.43
9:BT:101:BCL:HBB2	9:BT:101:BCL:HMB3	1.99	0.43
9:BU:102:BCL:HED2	6:BV:35:ALA:HB2	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:BW:12:TRP:CE3	5:BW:12:TRP:CA	3.01	0.43
5:BY:40:LEU:HD13	5:BY:46:TRP:CD2	2.53	0.43
9:BZ:101:BCL:CBB	9:B1:102:BCL:CHC	2.95	0.43
5:B1:30:VAL:HG13	5:B1:31:LEU:N	2.33	0.43
6:B2:17:PHE:CD1	6:B2:17:PHE:C	2.94	0.43
6:B4:40:TRP:HZ3	6:B4:44:PRO:C	2.26	0.43
5:B5:14:ILE:CD1	14:B5:103:CRT:H41	2.48	0.43
5:B7:44:LEU:HB2	6:B8:43:ARG:NH1	2.32	0.43
6:B8:26:TYR:O	6:B8:30:GLY:N	2.50	0.43
1:AC:36:ARG:NH1	2:AL:91:GLU:C	2.77	0.43
3:AM:11:VAL:HG11	4:AH:151:PRO:HD3	1.99	0.43
3:AM:27:ASN:ND2	5:AO:19:ARG:HH11	2.13	0.43
3:AM:83:VAL:O	3:AM:84:PHE:C	2.61	0.43
14:AA:102:CRT:H23	6:AE:16:GLU:CG	2.45	0.43
9:AB:101:BCL:HBB3	9:AB:101:BCL:HMB1	2.00	0.43
5:AD:49:ASP:HB2	5:AF:56:GLN:HG2	1.98	0.43
6:AE:31:LEU:C	6:AE:34:ILE:HG22	2.43	0.43
5:AI:10:LYS:CB	14:AN:102:CRT:H5	2.49	0.43
6:AJ:17:PHE:O	6:AJ:17:PHE:HD1	2.02	0.43
5:AK:19:ARG:O	5:AK:23:SER:HB3	2.17	0.43
6:AN:40:TRP:CE2	6:AN:44:PRO:HB3	2.53	0.43
6:AN:40:TRP:NE1	6:AN:44:PRO:HB3	2.33	0.43
9:AR:101:BCL:HBB3	9:AS:103:BCL:NC	2.33	0.43
5:AS:20:VAL:O	5:AS:24:ILE:HG12	2.17	0.43
9:AU:102:BCL:HBD	9:AV:102:BCL:OBD	2.17	0.43
5:AW:52:PRO:HD2	5:AW:55:TYR:OH	2.18	0.43
6:AX:45:TRP:CE2	9:AX:101:BCL:H2C	2.53	0.43
6:A2:22:MET:O	6:A2:26:TYR:HD1	2.01	0.43
5:A7:39:VAL:HG11	9:A7:103:BCL:HBC1	2.00	0.43
6:A0:36:HIS:CE1	9:A0:102:BCL:CHB	2.98	0.43
1:BC:129:ARG:HG2	1:BC:287:LEU:HD11	2.00	0.43
2:BL:77:PRO:HB3	2:BL:95:TRP:CE2	2.53	0.43
3:BM:63:PHE:C	3:BM:125:SER:HB2	2.44	0.43
3:BM:73:PHE:O	3:BM:74:ASN:C	2.61	0.43
3:BM:75:MET:HG2	3:BM:93:LEU:HB3	1.99	0.43
4:BH:222:VAL:HA	4:BH:242:TYR:CD2	2.53	0.43
5:BD:51:ILE:HG22	5:BD:52:PRO:HA	2.00	0.43
6:BE:10:THR:CG2	6:BE:11:ASP:N	2.80	0.43
5:BF:9:TYR:HA	6:BG:18:HIS:CG	2.53	0.43
6:BG:17:PHE:CD2	14:BG:102:CRT:H6	2.53	0.43
5:BI:30:VAL:HG13	5:BI:31:LEU:N	2.31	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:BI:102:BCL:HBC2	9:BJ:101:BCL:HMD1	2.00	0.43
6:BN:10:THR:CG2	6:BN:11:ASP:H	2.08	0.43
9:BN:101:BCL:HMB3	9:BO:102:BCL:HMA1	2.00	0.43
9:BS:102:BCL:O1A	6:BT:28:TRP:HZ2	2.01	0.43
5:BU:14:ILE:CD1	14:BU:103:CRT:H31A	2.28	0.43
9:BU:102:BCL:HMB3	9:BU:102:BCL:HBB2	2.00	0.43
14:BU:103:CRT:H15	14:BU:103:CRT:H131	1.72	0.43
5:BY:46:TRP:CZ3	9:BY:102:BCL:CBC	3.00	0.43
5:B1:9:TYR:HA	6:B2:18:HIS:ND1	2.33	0.43
6:B2:46:LEU:HB2	5:B3:52:PRO:HD3	1.97	0.43
6:B8:42:TYR:CG	6:B8:43:ARG:N	2.85	0.43
5:B9:35:ILE:CG2	9:B0:102:BCL:C1D	2.96	0.43
1:AC:175:PRO:HG3	3:AM:77:ALA:O	2.17	0.43
1:AC:246:GLY:C	1:AC:248:THR:N	2.74	0.43
1:AC:275:HIS:O	1:AC:275:HIS:CD2	2.71	0.43
2:AL:46:GLY:O	2:AL:47:VAL:C	2.61	0.43
2:AL:47:VAL:C	2:AL:50:ILE:HG22	2.43	0.43
2:AL:136:ALA:HA	2:AL:139:VAL:HG12	2.00	0.43
2:AL:190:PHE:O	2:AL:191:THR:C	2.61	0.43
3:AM:67:ALA:O	3:AM:71:ILE:N	2.51	0.43
3:AM:78:SER:C	3:AM:80:HIS:N	2.77	0.43
3:AM:212:SER:C	3:AM:214:LEU:N	2.74	0.43
5:AA:10:LYS:C	14:AA:102:CRT:H82	2.43	0.43
5:AD:50:ASN:O	5:AD:53:VAL:CB	2.66	0.43
5:AF:33:LEU:HD12	5:AF:33:LEU:H	1.81	0.43
6:AG:24:SER:O	6:AG:27:ALA:N	2.50	0.43
9:AG:101:BCL:HBA1	9:AG:101:BCL:H3A	1.58	0.43
5:AI:10:LYS:HB2	14:AN:102:CRT:H83	1.99	0.43
5:AI:27:PHE:CE2	5:AK:29:ILE:CD1	2.98	0.43
9:AQ:102:BCL:HHC	9:AQ:102:BCL:OBB	2.18	0.43
5:AU:34:LEU:O	5:AU:38:ILE:HG22	2.18	0.43
6:AX:21:PHE:CD2	14:AX:102:CRT:C16	3.01	0.43
5:AY:16:ASP:O	5:AY:20:VAL:HG22	2.18	0.43
6:A4:40:TRP:CZ3	6:A4:44:PRO:C	2.96	0.43
6:A4:40:TRP:CE3	6:A4:44:PRO:HA	2.54	0.43
6:A6:45:TRP:CG	6:A6:46:LEU:H	2.27	0.43
1:BC:65:ALA:CB	1:BC:89:GLU:OE1	2.67	0.43
1:BC:90:PHE:HZ	7:BC:501:HEM:C2A	2.37	0.43
1:BC:129:ARG:HA	1:BC:132:GLU:HG3	1.99	0.43
1:BC:156:HIS:CE1	1:BC:160:PRO:O	2.71	0.43
1:BC:174:TYR:O	1:BC:174:TYR:HD1	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BC:203:PHE:CD1	1:BC:210:ILE:HG12	2.50	0.43
1:BC:272:ALA:C	1:BC:274:ARG:N	2.77	0.43
1:BC:303:LEU:O	7:BC:502:HEM:HMD2	2.18	0.43
2:BL:142:PHE:CD1	2:BL:142:PHE:C	2.95	0.43
2:BL:150:ALA:O	2:BL:153:HIS:CB	2.67	0.43
9:BL:301:BCL:H203	9:BL:301:BCL:H13	2.00	0.43
3:BM:39:TRP:O	3:BM:40:LEU:C	2.58	0.43
3:BM:70:ILE:C	3:BM:72:GLY:N	2.73	0.43
3:BM:297:TRP:HZ3	3:BM:303:MET:SD	2.41	0.43
6:BB:46:LEU:OXT	6:BE:43:ARG:NH2	2.48	0.43
6:BE:10:THR:CG2	6:BE:11:ASP:H	2.22	0.43
5:BI:44:LEU:CD1	5:BI:46:TRP:HE3	2.30	0.43
9:BO:102:BCL:HHC	9:BO:102:BCL:OBB	2.18	0.43
6:BP:28:TRP:O	6:BP:29:PHE:C	2.62	0.43
5:BQ:50:ASN:HB3	5:BS:56:GLN:CA	2.46	0.43
6:BV:21:PHE:CD1	14:BV:102:CRT:C14	2.68	0.43
9:BX:101:BCL:H3A	9:BX:101:BCL:HBA1	1.35	0.43
5:B1:12:TRP:HD1	6:B2:18:HIS:HB2	1.83	0.43
9:B1:102:BCL:OBB	9:B1:102:BCL:HHC	2.18	0.43
9:B2:101:BCL:CMB	9:B3:102:BCL:C1B	2.95	0.43
6:B4:40:TRP:CZ3	6:B4:44:PRO:C	2.96	0.43
9:B7:103:BCL:C1D	9:B8:101:BCL:HMD1	2.48	0.43
6:B8:31:LEU:C	6:B8:34:ILE:HG22	2.43	0.43
5:B9:36:HIS:CD2	9:B0:102:BCL:HMD2	2.54	0.43
1:AC:20:LEU:HG	2:AL:271:TRP:CZ2	2.53	0.43
1:AC:57:GLN:NE2	1:AC:58:PRO:HD2	2.33	0.43
1:AC:62:LEU:O	1:AC:92:ARG:NH2	2.51	0.43
1:AC:111:HIS:CE1	1:AC:124:LYS:CE	3.01	0.43
2:AL:198:MET:O	2:AL:201:SER:HB3	2.18	0.43
2:AL:230:GLY:CA	3:AM:51:ILE:HD12	2.49	0.43
2:AL:231:TYR:CD1	2:AL:231:TYR:C	2.96	0.43
2:AL:241:LEU:O	2:AL:245:LEU:HG	2.17	0.43
3:AM:83:VAL:HG23	3:AM:84:PHE:H	1.83	0.43
3:AM:84:PHE:HD1	3:AM:84:PHE:N	2.16	0.43
3:AM:193:TYR:HA	3:AM:292:ASP:O	2.19	0.43
3:AM:226:VAL:O	3:AM:226:VAL:HG22	2.19	0.43
4:AH:4:GLY:O	4:AH:5:ILE:HB	2.18	0.43
4:AH:80:ARG:O	4:AH:81:VAL:C	2.62	0.43
4:AH:134:VAL:HG21	4:AH:174:ARG:HH21	1.83	0.43
5:AA:33:LEU:HA	14:A0:101:CRT:H391	2.00	0.43
9:AF:102:BCL:CBB	9:AF:102:BCL:HMB3	2.49	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:AG:28:TRP:O	6:AG:29:PHE:C	2.60	0.43
6:AG:46:LEU:HD13	6:AJ:42:TYR:CE1	2.52	0.43
5:AI:8:LEU:O	5:AI:11:ILE:HG22	2.18	0.43
5:AI:33:LEU:HD12	5:AI:33:LEU:C	2.44	0.43
6:AN:36:HIS:CG	9:AN:101:BCL:H162	2.54	0.43
6:AP:18:HIS:O	6:AP:22:MET:HB2	2.18	0.43
5:AQ:10:LYS:C	14:AT:102:CRT:H82	2.43	0.43
5:AS:4:MET:HB2	5:AS:8:LEU:CD1	2.47	0.43
5:AS:24:ILE:HG21	14:AT:102:CRT:C20	2.48	0.43
5:AU:16:ASP:CG	5:AU:19:ARG:HD3	2.43	0.43
9:AY:102:BCL:HAC2	9:AZ:101:BCL:CBC	2.48	0.43
9:AZ:101:BCL:H201	6:A2:38:LEU:HD21	1.94	0.43
5:A5:18:ARG:HB2	5:A5:19:ARG:HH12	1.82	0.43
5:A5:40:LEU:HD13	5:A5:46:TRP:CE3	2.53	0.43
5:A5:50:ASN:OD1	5:A5:51:ILE:N	2.52	0.43
9:A8:101:BCL:HMA3	9:A9:102:BCL:HMA1	1.99	0.43
2:BL:3:MET:HG2	2:BL:11:ARG:NH2	2.32	0.43
2:BL:89:LEU:H	2:BL:89:LEU:HD12	1.83	0.43
2:BL:192:ASN:C	2:BL:192:ASN:ND2	2.75	0.43
2:BL:209:PRO:O	2:BL:210:GLN:O	2.36	0.43
2:BL:216:LYS:HB3	2:BL:220:HIS:CG	2.54	0.43
5:BD:35:ILE:HA	5:BD:38:ILE:HG22	1.99	0.43
6:BE:43:ARG:NH1	5:BF:55:TYR:CD2	2.78	0.43
5:BF:39:VAL:O	5:BF:39:VAL:HG12	2.18	0.43
9:BF:102:BCL:CHD	9:BG:101:BCL:HMD1	2.48	0.43
6:BG:21:PHE:HB2	14:BG:102:CRT:H11	1.99	0.43
5:BO:5:ASN:O	5:BO:8:LEU:HD22	2.19	0.43
14:BO:103:CRT:H10	14:BO:103:CRT:H81	1.80	0.43
6:BP:40:TRP:HZ3	6:BP:45:TRP:N	2.17	0.43
5:BS:42:THR:CG2	5:BU:47:LEU:HB3	2.47	0.43
5:BW:9:TYR:HA	6:BX:18:HIS:ND1	2.28	0.43
5:BY:9:TYR:CE1	6:BZ:15:LYS:HG2	2.53	0.43
14:B2:102:CRT:H291	9:B3:102:BCL:O2A	2.18	0.43
1:AC:50:ALA:O	1:AC:51:LEU:C	2.60	0.43
1:AC:200:LEU:HD11	1:AC:238:ASN:HD22	1.78	0.43
2:AL:278:LEU:HD12	2:AL:281:TRP:HZ2	1.83	0.43
3:AM:56:THR:O	3:AM:60:SER:N	2.38	0.43
3:AM:106:ILE:H	5:AO:42:THR:HG21	1.83	0.43
3:AM:156:PHE:CZ	9:AM:402:BCL:HBD	2.52	0.43
3:AM:211:GLY:O	3:AM:215:LEU:N	2.44	0.43
5:AD:5:ASN:HD22	6:AE:22:MET:CG	2.30	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:AD:11:ILE:C	5:AD:13:LEU:H	2.26	0.43
5:AD:35:ILE:C	5:AD:37:MET:N	2.74	0.43
9:AD:102:BCL:CAC	9:AE:101:BCL:CBC	2.96	0.43
5:AF:17:PRO:O	5:AF:21:LEU:HB2	2.19	0.43
6:AG:36:HIS:CE1	9:AG:101:BCL:C4A	3.01	0.43
5:AK:22:VAL:HG13	5:AK:23:SER:N	2.34	0.43
5:AO:18:ARG:H	5:AO:18:ARG:HG2	1.42	0.43
6:AP:21:PHE:CE1	14:AP:102:CRT:H16	2.53	0.43
9:AP:101:BCL:HBB1	9:AQ:102:BCL:HMC1	1.99	0.43
6:AT:11:ASP:O	6:AT:15:LYS:HD2	2.18	0.43
5:AU:26:ALA:HA	5:AU:29:ILE:CG2	2.48	0.43
5:AU:30:VAL:CG1	5:AU:31:LEU:N	2.80	0.43
6:AV:44:PRO:O	5:AW:52:PRO:HD2	2.19	0.43
9:AV:102:BCL:H3A	9:AV:102:BCL:HBA1	1.76	0.43
5:AW:9:TYR:CD1	6:AX:15:LYS:HB2	2.54	0.43
5:AW:48:ASP:CG	5:AW:48:ASP:O	2.60	0.43
9:AZ:101:BCL:H3A	9:AZ:101:BCL:HBA1	1.31	0.43
5:A3:46:TRP:CZ2	9:A3:103:BCL:CHC	3.02	0.43
6:A6:29:PHE:HE1	9:A6:101:BCL:H11	1.70	0.43
6:A8:9:LEU:HD22	6:A8:13:GLU:CD	2.43	0.43
1:BC:36:ARG:HD2	2:BL:77:PRO:O	2.18	0.43
2:BL:6:PHE:O	2:BL:9:LYS:HG2	2.18	0.43
2:BL:17:LEU:CD2	2:BL:118:ARG:HD2	2.48	0.43
2:BL:68:TYR:CA	2:BL:73:ILE:HD11	2.47	0.43
2:BL:240:ARG:O	2:BL:241:LEU:C	2.61	0.43
10:BL:302:BPH:H161	9:BL:303:BCL:H171	2.00	0.43
3:BM:130:TRP:HA	3:BM:150:PHE:CD2	2.53	0.43
3:BM:144:GLN:HA	3:BM:144:GLN:OE1	2.19	0.43
3:BM:180:PHE:O	3:BM:183:LEU:N	2.52	0.43
3:BM:243:THR:HA	3:BM:246:GLU:HB2	2.01	0.43
4:BH:176:GLU:HG3	4:BH:178:GLN:CG	2.33	0.43
5:BD:38:ILE:HD11	5:BF:40:LEU:CD2	2.48	0.43
5:BF:45:ASN:O	5:BF:47:LEU:N	2.52	0.43
6:BJ:37:LEU:HD13	6:BJ:37:LEU:HA	1.92	0.43
5:BQ:17:PRO:HB3	6:BR:17:PHE:CZ	2.54	0.43
9:BQ:104:BCL:CMA	9:BS:102:BCL:HMA1	2.47	0.43
5:BS:12:TRP:HE1	6:BT:18:HIS:CG	2.37	0.43
5:BS:38:ILE:HG23	5:BS:39:VAL:N	2.33	0.43
14:BV:102:CRT:C2M	5:BW:37:MET:HG2	2.46	0.43
5:BW:29:ILE:CG2	5:BW:30:VAL:N	2.82	0.43
5:BW:51:ILE:HB	5:BW:52:PRO:O	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:BX:46:LEU:CD2	6:BZ:42:TYR:CE2	3.02	0.43
5:BY:13:LEU:O	6:BZ:7:THR:HA	2.19	0.43
9:B1:102:BCL:CBB	9:B1:102:BCL:CMB	2.96	0.43
6:B2:16:GLU:HB2	14:B2:102:CRT:C1M	2.47	0.43
5:B3:46:TRP:CZ3	9:B3:102:BCL:HAC1	2.53	0.43
9:B3:102:BCL:H13	9:B3:102:BCL:H193	2.00	0.43
5:B7:46:TRP:HH2	9:B7:103:BCL:HBC3	1.66	0.43
5:B9:39:VAL:C	5:B9:41:SER:N	2.76	0.43
1:AC:211:ARG:C	1:AC:212:ILE:HD13	2.44	0.43
1:AC:282:ASN:HB3	1:AC:283:TYR:CD1	2.53	0.43
2:AL:30:PHE:CD2	3:AM:255:THR:O	2.72	0.43
2:AL:138:LEU:N	2:AL:138:LEU:CD1	2.81	0.43
2:AL:140:LEU:CD1	9:AL:301:BCL:O2D	2.67	0.43
10:AL:302:BPH:HMB3	10:AL:302:BPH:CBB	2.48	0.43
11:AL:304:UQ8:H32	11:AL:304:UQ8:H35	1.58	0.43
3:AM:165:PRO:CB	3:AM:174:ALA:HB2	2.48	0.43
3:AM:172:ALA:O	3:AM:174:ALA:N	2.51	0.43
3:AM:223:ILE:HD11	3:AM:234:GLU:OE2	2.19	0.43
15:AM:409:PEF:H21	15:AM:409:PEF:H121	2.00	0.43
4:AH:159:LEU:C	4:AH:159:LEU:CD1	2.92	0.43
5:AA:29:ILE:HG12	5:A9:27:PHE:HE2	1.82	0.43
6:AG:21:PHE:CD2	14:AG:102:CRT:H14	2.54	0.43
6:AG:28:TRP:CD1	6:AG:32:VAL:HG21	2.49	0.43
5:AI:39:VAL:CG1	9:AI:102:BCL:HBC1	2.40	0.43
5:AI:40:LEU:N	5:AI:40:LEU:CD1	2.81	0.43
6:AJ:17:PHE:HA	6:AJ:20:ILE:CG2	2.43	0.43
5:AO:9:TYR:HA	6:AP:18:HIS:CG	2.54	0.43
5:AO:14:ILE:O	6:AP:7:THR:N	2.51	0.43
5:AO:21:LEU:HD13	6:AP:17:PHE:HZ	1.83	0.43
5:AS:28:GLN:HB2	9:AS:103:BCL:H43	2.01	0.43
6:AV:10:THR:HB	6:AV:13:GLU:OE2	2.19	0.43
6:AX:21:PHE:CD1	6:AX:22:MET:N	2.87	0.43
6:AX:46:LEU:HD23	6:AX:46:LEU:HA	1.86	0.43
5:AY:5:ASN:O	5:AY:6:ALA:C	2.61	0.43
5:AY:12:TRP:HH2	9:A1:102:BCL:H203	1.84	0.43
5:A1:7:ASN:O	5:A1:10:LYS:HE3	2.19	0.43
5:A1:39:VAL:HG13	5:A1:40:LEU:N	2.33	0.43
9:A1:102:BCL:H92	14:A2:102:CRT:H181	1.95	0.43
9:A2:101:BCL:CHC	9:A3:103:BCL:HBB3	2.49	0.43
9:A5:102:BCL:O1D	9:A5:102:BCL:C2A	2.63	0.43
5:A7:40:LEU:HD11	5:A7:47:LEU:CD2	2.49	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:A8:31:LEU:C	6:A8:34:ILE:HG22	2.43	0.43
1:BC:161:VAL:HG13	7:BC:502:HEM:O1D	2.17	0.43
1:BC:183:GLN:HE22	1:BC:230:GLU:CD	2.26	0.43
2:BL:246:ALA:HB3	3:BM:217:ALA:HB2	2.00	0.43
3:BM:174:ALA:O	3:BM:175:VAL:O	2.36	0.43
3:BM:216:PHE:O	3:BM:216:PHE:CG	2.72	0.43
4:BH:240:CYS:O	4:BH:241:ALA:C	2.62	0.43
6:BB:21:PHE:HD1	14:BB:102:CRT:H14	1.83	0.43
6:BB:26:TYR:HA	6:BB:29:PHE:HD2	1.84	0.43
5:BF:38:ILE:O	5:BF:38:ILE:HG12	2.18	0.43
6:BG:24:SER:C	6:BG:26:TYR:N	2.75	0.43
5:BI:15:LEU:HB3	5:BI:20:VAL:CG2	2.47	0.43
5:BK:8:LEU:HD22	5:BK:11:ILE:HD11	2.00	0.43
6:BT:10:THR:CG2	6:BT:11:ASP:H	2.19	0.43
14:BV:102:CRT:H2M1	5:BW:37:MET:CB	2.39	0.43
14:BV:102:CRT:H2M1	5:BW:37:MET:CG	2.49	0.43
5:BY:46:TRP:CD1	5:BY:47:LEU:N	2.86	0.43
14:B5:103:CRT:C34	9:B9:102:BCL:HBA1	2.24	0.43
5:B9:2:PHE:CD1	5:B9:2:PHE:N	2.85	0.43
9:B9:102:BCL:C1D	9:B0:102:BCL:HMD1	2.49	0.43
1:AC:236:MET:HE1	7:AC:503:HEM:C1D	2.53	0.43
1:AC:259:TRP:CH2	1:AC:266:ARG:NH2	2.86	0.43
1:AC:267:THR:C	1:AC:269:ALA:N	2.77	0.43
2:AL:43:THR:O	2:AL:47:VAL:N	2.45	0.43
2:AL:95:TRP:CE3	2:AL:96:GLN:HA	2.53	0.43
9:AL:301:BCL:CGA	9:AL:303:BCL:HBC1	2.48	0.43
10:AL:302:BPH:H141	9:AL:303:BCL:HBB3	2.00	0.43
9:AL:303:BCL:OBB	9:AL:303:BCL:HHC	2.18	0.43
3:AM:91:PHE:O	3:AM:180:PHE:CG	2.71	0.43
3:AM:164:ARG:C	3:AM:167:MET:H	2.27	0.43
3:AM:235:ILE:O	3:AM:238:ILE:HB	2.18	0.43
4:AH:125:LEU:CB	4:AH:129:GLY:O	2.65	0.43
4:AH:176:GLU:HA	4:AH:177:PRO:HD2	1.86	0.43
4:AH:226:SER:OG	4:AH:227:ASN:N	2.52	0.43
5:AD:13:LEU:C	5:AD:14:ILE:HD12	2.44	0.43
9:AD:102:BCL:H13	9:AD:102:BCL:H193	2.01	0.43
5:AF:35:ILE:CA	5:AF:38:ILE:HG22	2.47	0.43
9:AI:102:BCL:HMB1	9:AI:102:BCL:CBB	2.49	0.43
5:AK:14:ILE:HG23	5:AO:18:ARG:HD3	2.00	0.43
5:AK:49:ASP:OD1	5:AK:50:ASN:N	2.51	0.43
9:AO:102:BCL:HBA1	9:AO:102:BCL:H3A	1.81	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
14:AP:102:CRT:H2M3	5:AQ:36:HIS:HB3	2.00	0.43
6:AR:26:TYR:O	6:AR:27:ALA:C	2.60	0.43
9:AV:102:BCL:H8	9:AV:102:BCL:CMB	2.49	0.43
6:AX:28:TRP:HA	6:AX:31:LEU:HG	2.01	0.43
9:AZ:101:BCL:HMA1	9:A1:102:BCL:HHB	2.01	0.43
6:A6:25:MET:SD	6:A6:29:PHE:CE2	3.12	0.43
5:A9:26:ALA:O	5:A9:30:VAL:HG23	2.19	0.43
5:A9:46:TRP:CZ3	9:A9:102:BCL:HBC3	2.53	0.43
6:A0:10:THR:H	6:A0:13:GLU:HG2	1.84	0.43
6:A0:42:TYR:CE2	6:A0:43:ARG:HD2	2.54	0.43
1:BC:84:ASP:CG	1:BC:333:THR:HG21	2.43	0.43
1:BC:203:PHE:CD1	1:BC:235:LEU:HD22	2.53	0.43
1:BC:218:LEU:HA	3:BM:291:VAL:HA	2.00	0.43
1:BC:313:ALA:C	1:BC:314:VAL:CG2	2.91	0.43
2:BL:239:HIS:CG	3:BM:223:ILE:HG21	2.54	0.43
2:BL:250:ALA:HB2	10:BL:302:BPH:CBC	2.49	0.43
3:BM:189:PHE:HB3	9:BM:402:BCL:CMD	2.48	0.43
3:BM:204:LEU:C	3:BM:204:LEU:HD13	2.43	0.43
9:BA:101:BCL:CBD	9:BB:101:BCL:OBD	2.67	0.43
6:BB:22:MET:O	6:BB:26:TYR:HD1	1.97	0.43
5:BD:13:LEU:O	6:BE:7:THR:HB	2.18	0.43
6:BE:42:TYR:CD2	6:BE:43:ARG:HG3	2.53	0.43
5:BF:18:ARG:CZ	5:BF:18:ARG:HB2	2.48	0.43
5:BI:9:TYR:CD2	6:BJ:15:LYS:HG2	2.53	0.43
9:BN:101:BCL:HBA1	9:BN:101:BCL:H3A	1.75	0.43
5:BS:46:TRP:CZ3	9:BS:102:BCL:HBC3	2.54	0.43
5:BS:46:TRP:CE3	9:BS:102:BCL:H2C	2.54	0.43
5:BU:9:TYR:CE2	5:BU:10:LYS:HG2	2.54	0.43
6:BV:30:GLY:O	6:BV:34:ILE:HG13	2.19	0.43
5:BY:32:GLY:CA	9:BZ:101:BCL:HED2	2.49	0.43
14:B5:103:CRT:H2M3	5:B9:36:HIS:CB	2.48	0.43
6:B6:25:MET:SD	6:B6:29:PHE:CE2	3.12	0.43
1:AC:224:ALA:O	3:AM:192:ARG:NE	2.29	0.43
1:AC:232:THR:O	1:AC:235:LEU:HB3	2.18	0.43
1:AC:316:LYS:O	1:AC:317:PRO:O	2.37	0.43
2:AL:194:LEU:C	2:AL:194:LEU:CD2	2.92	0.43
2:AL:280:LEU:HA	5:AW:38:ILE:HG13	1.99	0.43
3:AM:241:ARG:CG	3:AM:242:GLY:H	2.14	0.43
14:AB:102:CRT:H2M2	5:AD:37:MET:HE2	2.00	0.43
5:AD:35:ILE:O	5:AD:36:HIS:C	2.59	0.43
5:AF:43:ASP:OD1	5:AF:44:LEU:CG	2.66	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:AG:101:BCL:CMC	9:AI:102:BCL:HBB1	2.49	0.43
5:AI:26:ALA:C	5:AI:29:ILE:HG22	2.42	0.43
6:AP:27:ALA:O	6:AP:31:LEU:CG	2.56	0.43
6:AP:28:TRP:O	6:AP:29:PHE:C	2.62	0.43
5:AQ:29:ILE:CG2	5:AQ:30:VAL:N	2.82	0.43
5:AQ:42:THR:C	5:AS:48:ASP:HB3	2.43	0.43
6:AT:45:TRP:HD1	6:AT:46:LEU:H	1.66	0.43
5:AY:34:LEU:O	5:AY:37:MET:HB2	2.19	0.43
6:AZ:33:VAL:HG13	6:AZ:34:ILE:N	2.33	0.43
5:A1:11:ILE:HG22	14:A1:103:CRT:C8	2.42	0.43
6:A2:21:PHE:O	6:A2:22:MET:C	2.61	0.43
6:A4:40:TRP:HZ3	6:A4:44:PRO:C	2.26	0.43
5:A5:46:TRP:CZ3	9:A5:102:BCL:HBC3	2.54	0.43
5:A9:36:HIS:CE1	9:A0:102:BCL:OBD	2.71	0.43
2:BL:50:ILE:HG21	9:BL:303:BCL:H191	1.99	0.43
3:BM:83:VAL:O	3:BM:84:PHE:C	2.62	0.43
3:BM:314:VAL:HG12	3:BM:315:ASN:N	2.33	0.43
5:BA:47:LEU:CG	5:B9:43:ASP:HB2	2.48	0.43
9:BA:101:BCL:CMD	6:BB:36:HIS:ND1	2.82	0.43
5:BF:54:SER:O	5:BF:58:LEU:CB	2.66	0.43
5:BK:22:VAL:HA	5:BK:25:VAL:HG23	2.01	0.43
5:BK:46:TRP:CD2	9:BK:102:BCL:H2C	2.54	0.43
5:BO:10:LYS:HB2	14:BO:103:CRT:H83	2.00	0.43
5:BO:36:HIS:O	5:BO:40:LEU:CB	2.67	0.43
6:BT:29:PHE:CE1	9:BT:101:BCL:C1	3.01	0.43
5:BY:9:TYR:HB2	6:BZ:15:LYS:HA	2.01	0.43
5:BY:55:TYR:CD1	5:BY:56:GLN:N	2.76	0.43
6:B2:33:VAL:HG13	6:B2:34:ILE:N	2.33	0.43
6:B2:40:TRP:CZ3	6:B2:44:PRO:CA	3.01	0.43
9:B7:103:BCL:CMD	9:B8:101:BCL:C1D	2.97	0.43
6:B0:33:VAL:HG12	6:B0:37:LEU:HD12	1.94	0.43
1:AC:162:PRO:C	1:AC:164:TYR:H	2.27	0.43
1:AC:184:ASN:ND2	3:AM:96:GLU:HG2	2.34	0.43
2:AL:7:GLU:O	2:AL:9:LYS:N	2.52	0.43
2:AL:18:ILE:HG23	4:AH:259:LEU:HB2	2.01	0.43
2:AL:49:LEU:HD12	2:AL:98:ILE:CG1	2.49	0.43
2:AL:75:ILE:HG21	2:AL:95:TRP:HA	2.01	0.43
2:AL:155:PHE:HB2	2:AL:156:PRO:CD	2.44	0.43
2:AL:208:ASN:N	2:AL:209:PRO:CD	2.82	0.43
2:AL:257:ILE:HD13	9:AL:301:BCL:OBD	2.18	0.43
3:AM:197:TYR:HA	17:AM:503:HOH:O	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:AM:401:BCL:HBA1	9:AM:401:BCL:H3A	1.72	0.43
4:AH:189:ASN:N	4:AH:189:ASN:HD22	2.16	0.43
5:AA:47:LEU:CB	5:A9:43:ASP:HB2	2.47	0.43
5:AF:40:LEU:HD22	5:AF:45:ASN:CA	2.48	0.43
5:AI:46:TRP:NE1	5:AI:47:LEU:HD13	2.34	0.43
5:AK:53:VAL:O	5:AK:54:SER:C	2.62	0.43
6:AN:46:LEU:O	5:AO:51:ILE:HD12	2.19	0.43
9:AR:101:BCL:C2B	9:AS:103:BCL:C1B	2.97	0.43
6:AX:17:PHE:O	6:AX:18:HIS:C	2.62	0.43
5:A3:35:ILE:O	5:A3:38:ILE:HG22	2.18	0.43
5:A3:43:ASP:OD2	5:A5:47:LEU:O	2.36	0.43
5:A5:4:MET:HB3	6:A8:24:SER:OG	2.19	0.43
5:A9:32:GLY:N	9:A0:102:BCL:HED2	2.33	0.43
2:BL:3:MET:HE2	4:BH:45:ARG:HE	1.83	0.43
2:BL:71:TRP:O	2:BL:160:LEU:HG	2.19	0.43
2:BL:104:GLY:HA2	2:BL:107:ILE:HD12	2.00	0.43
2:BL:166:VAL:HG13	9:BL:301:BCL:HMD1	2.00	0.43
2:BL:259:ILE:O	2:BL:261:GLY:N	2.52	0.43
3:BM:83:VAL:HG23	3:BM:84:PHE:N	2.34	0.43
3:BM:168:MET:HG2	3:BM:289:THR:HG22	2.01	0.43
5:BF:45:ASN:C	5:BF:49:ASP:OD1	2.62	0.43
6:BG:38:LEU:O	6:BG:38:LEU:HD23	2.17	0.43
5:BK:51:ILE:CA	5:BK:52:PRO:C	2.92	0.43
5:BQ:51:ILE:CG1	5:BQ:52:PRO:CA	2.94	0.43
9:BT:101:BCL:HMB3	9:BT:101:BCL:CBB	2.48	0.43
6:BX:46:LEU:HD23	6:BX:46:LEU:HA	1.80	0.43
5:B1:52:PRO:O	5:B1:53:VAL:C	2.62	0.43
6:B2:41:LEU:HD12	6:B2:42:TYR:N	2.34	0.43
9:B2:101:BCL:HMB3	9:B3:102:BCL:C1B	2.48	0.43
5:B3:32:GLY:C	5:B3:34:LEU:N	2.77	0.43
5:B7:21:LEU:O	5:B7:25:VAL:HG23	2.19	0.43
5:B7:33:LEU:H	5:B7:33:LEU:CD1	2.30	0.43
5:B7:33:LEU:HG	14:B7:102:CRT:C36	2.48	0.43
1:AC:53:ILE:CG1	1:AC:319:TYR:CZ	3.00	0.43
1:AC:96:ALA:C	1:AC:98:THR:N	2.62	0.43
1:AC:148:THR:N	1:AC:322:GLN:NE2	2.67	0.43
1:AC:253:THR:C	1:AC:255:ALA:N	2.77	0.43
2:AL:79:ASP:C	2:AL:81:SER:N	2.76	0.43
3:AM:191:ILE:C	3:AM:193:TYR:N	2.75	0.43
3:AM:226:VAL:O	3:AM:226:VAL:HG13	2.18	0.43
15:AM:409:PEF:H52	5:AQ:19:ARG:NH2	2.33	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:AH:17:TRP:O	4:AH:18:ALA:O	2.37	0.43
4:AH:29:TYR:CD1	4:AH:30:LEU:N	2.87	0.43
4:AH:31:ARG:HD2	15:AH:301:PEF:O4	2.19	0.43
4:AH:48:ARG:HB3	15:AH:301:PEF:H42	2.00	0.43
4:AH:94:PRO:HG2	6:A0:8:GLY:HA3	1.99	0.43
5:AA:29:ILE:C	5:AA:31:LEU:H	2.26	0.43
6:AB:24:SER:HG	5:A9:4:MET:HE2	1.79	0.43
5:AD:9:TYR:CZ	6:AE:11:ASP:HB3	2.54	0.43
5:AD:9:TYR:OH	6:AE:11:ASP:HB3	2.18	0.43
6:AE:9:LEU:HB3	6:AE:13:GLU:HG2	2.01	0.43
6:AJ:45:TRP:O	6:AJ:46:LEU:CB	2.64	0.43
5:AK:52:PRO:HB2	5:AK:55:TYR:HD1	1.82	0.43
5:AO:5:ASN:ND2	5:AO:8:LEU:HD21	2.32	0.43
5:AO:43:ASP:O	5:AQ:48:ASP:HB3	2.19	0.43
5:AS:13:LEU:HD21	6:AT:10:THR:O	2.19	0.43
5:AU:9:TYR:N	6:AV:18:HIS:CE1	2.87	0.43
5:AU:22:VAL:CG1	5:AU:23:SER:N	2.79	0.43
5:AW:5:ASN:HA	5:AW:8:LEU:CD1	2.33	0.43
5:AW:12:TRP:CZ2	6:AX:21:PHE:CD1	3.06	0.43
5:AW:38:ILE:O	5:AW:41:SER:OG	2.33	0.43
9:A3:103:BCL:CED	6:A4:32:VAL:HA	2.48	0.43
5:A5:4:MET:HG3	6:A8:27:ALA:HB2	1.90	0.43
14:A5:103:CRT:H15	14:A5:103:CRT:H131	1.89	0.43
6:A6:37:LEU:C	6:A6:37:LEU:CD2	2.91	0.43
9:A7:103:BCL:ND	9:A8:101:BCL:HMD3	2.33	0.43
5:A9:44:LEU:CD1	5:A9:44:LEU:N	2.82	0.43
1:BC:129:ARG:NH1	1:BC:132:GLU:HB2	2.33	0.43
1:BC:167:VAL:O	1:BC:168:THR:C	2.61	0.43
1:BC:245:VAL:HG21	1:BC:249:PHE:CD1	2.54	0.43
1:BC:264:PRO:O	1:BC:265:LYS:C	2.62	0.43
2:BL:170:GLY:O	2:BL:176:PHE:HD2	2.02	0.43
2:BL:184:LEU:O	2:BL:187:SER:HB3	2.18	0.43
2:BL:239:HIS:CE1	3:BM:234:GLU:CD	2.96	0.43
3:BM:136:ARG:HA	3:BM:136:ARG:CZ	2.49	0.43
3:BM:174:ALA:O	3:BM:175:VAL:C	2.62	0.43
9:BI:102:BCL:CBB	9:BI:102:BCL:CMB	2.97	0.43
6:BJ:26:TYR:O	6:BJ:27:ALA:C	2.62	0.43
6:BJ:29:PHE:N	6:BJ:29:PHE:CD1	2.87	0.43
5:BK:19:ARG:HG3	5:BK:20:VAL:N	2.31	0.43
9:BN:101:BCL:HMC3	9:BO:102:BCL:CBB	2.47	0.43
9:BO:102:BCL:H13	9:BO:102:BCL:H193	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:BQ:31:LEU:HD21	9:BQ:104:BCL:HMA3	2.01	0.43
9:BQ:103:BCL:C1D	9:BQ:104:BCL:HMD1	2.34	0.43
5:BS:27:PHE:CD1	5:BU:29:ILE:HD11	2.54	0.43
5:B1:35:ILE:O	5:B1:39:VAL:HG12	2.18	0.43
6:B2:28:TRP:C	6:B2:30:GLY:N	2.77	0.43
5:B3:14:ILE:CD1	6:B6:17:PHE:CE2	2.96	0.43
5:B7:41:SER:OG	5:B7:42:THR:N	2.51	0.43
6:B0:17:PHE:HA	6:B0:20:ILE:HG22	2.01	0.43
1:AC:33:ILE:H	1:AC:33:ILE:CD1	2.30	0.42
1:AC:190:VAL:HG12	1:AC:237:MET:CB	2.46	0.42
9:AA:101:BCL:HBB1	9:A0:102:BCL:CMC	2.49	0.42
6:AG:24:SER:C	6:AG:26:TYR:N	2.75	0.42
5:AI:18:ARG:NH1	5:AI:18:ARG:CB	2.82	0.42
5:AK:29:ILE:O	5:AK:29:ILE:HG12	2.18	0.42
6:AN:23:GLN:HA	6:AN:26:TYR:HD2	1.84	0.42
6:AR:16:GLU:HB2	14:AR:102:CRT:H23	2.01	0.42
5:AW:10:LYS:NZ	14:AW:102:CRT:C1M	2.81	0.42
5:A1:27:PHE:CE2	5:A3:29:ILE:CD1	3.02	0.42
14:A1:103:CRT:C40	5:A3:38:ILE:HD12	2.48	0.42
6:A2:29:PHE:CD1	9:A2:101:BCL:H11	2.54	0.42
9:A3:103:BCL:H172	9:A3:103:BCL:H111	2.01	0.42
1:BC:97:VAL:HG13	7:BC:502:HEM:CMB	2.48	0.42
2:BL:147:LEU:O	2:BL:262:PRO:HG3	2.19	0.42
2:BL:164:ASP:OD2	3:BM:307:TYR:OH	2.36	0.42
2:BL:240:ARG:NH2	3:BM:6:ASN:C	2.76	0.42
3:BM:120:LEU:O	3:BM:123:THR:HB	2.19	0.42
3:BM:167:MET:HE3	3:BM:289:THR:HG23	2.00	0.42
3:BM:254:TRP:N	3:BM:254:TRP:HD1	2.17	0.42
3:BM:255:THR:HB	3:BM:256:MET:H	1.74	0.42
4:BH:122:HIS:C	4:BH:232:THR:HB	2.44	0.42
4:BH:138:VAL:C	4:BH:140:LYS:H	2.27	0.42
5:BD:33:LEU:HD12	5:BD:33:LEU:HA	1.94	0.42
5:BF:36:HIS:NE2	9:BG:101:BCL:HMD3	2.34	0.42
5:BK:22:VAL:O	5:BK:25:VAL:HB	2.18	0.42
9:BP:101:BCL:HMA1	9:BQ:103:BCL:HMA1	2.01	0.42
9:BS:102:BCL:H61	9:BS:102:BCL:H112	2.01	0.42
6:BT:40:TRP:CE3	6:BT:44:PRO:HA	2.54	0.42
5:BU:11:ILE:HA	14:BU:103:CRT:H21A	2.01	0.42
5:BU:35:ILE:CA	5:BU:38:ILE:HG22	2.48	0.42
9:BU:102:BCL:ND	9:BV:101:BCL:HMD3	2.33	0.42
14:BU:103:CRT:H26	14:BU:103:CRT:H241	1.90	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:BY:43:ASP:HA	5:B1:48:ASP:CB	2.40	0.42
6:BZ:45:TRP:CD2	9:BZ:101:BCL:H2C	2.54	0.42
6:B2:21:PHE:CE1	14:B2:102:CRT:H19	2.52	0.42
1:AC:246:GLY:O	1:AC:248:THR:N	2.52	0.42
2:AL:270:GLU:O	2:AL:272:TRP:N	2.52	0.42
3:AM:7:ILE:HG22	3:AM:8:PHE:CD2	2.54	0.42
3:AM:98:PRO:CB	3:AM:171:TRP:HB3	2.49	0.42
3:AM:250:LEU:HG	3:AM:254:TRP:NE1	2.26	0.42
3:AM:267:ARG:O	3:AM:268:TRP:C	2.60	0.42
3:AM:301:HIS:CE1	4:AH:8:TYR:CD2	3.06	0.42
9:AM:401:BCL:CMB	9:AM:402:BCL:H171	2.49	0.42
4:AH:69:LEU:HD23	4:AH:70:PRO:CD	2.46	0.42
5:AA:47:LEU:HD22	5:AA:47:LEU:H	1.84	0.42
9:AB:101:BCL:HMB3	9:AD:102:BCL:CHB	2.49	0.42
5:AF:44:LEU:HD12	5:AF:46:TRP:HE3	1.84	0.42
6:AG:21:PHE:HD1	6:AG:22:MET:HA	1.81	0.42
5:AK:18:ARG:HG2	5:AK:18:ARG:NH1	2.34	0.42
14:AP:102:CRT:H36	14:AP:102:CRT:H341	1.92	0.42
5:AQ:50:ASN:HA	5:AS:60:LYS:HA	2.01	0.42
5:AS:10:LYS:HB3	14:AS:104:CRT:C1	2.49	0.42
6:AT:40:TRP:CE3	6:AT:40:TRP:O	2.72	0.42
9:AV:102:BCL:HMA1	9:AW:101:BCL:HHB	2.02	0.42
9:AY:102:BCL:O1D	9:AY:102:BCL:C2A	2.66	0.42
14:A1:103:CRT:C14	5:A3:21:LEU:HD11	2.49	0.42
5:A5:28:GLN:NE2	14:A7:102:CRT:H25	2.34	0.42
6:A6:29:PHE:CZ	9:A6:101:BCL:H42	2.54	0.42
5:A9:4:MET:CE	5:A9:4:MET:HA	2.47	0.42
1:BC:114:GLY:O	1:BC:115:ASN:C	2.61	0.42
1:BC:187:SER:C	1:BC:189:THR:H	2.25	0.42
1:BC:195:LEU:CB	1:BC:196:PRO:CD	2.96	0.42
2:BL:148:MET:HB3	2:BL:153:HIS:CE1	2.53	0.42
2:BL:188:PHE:HD2	2:BL:249:ALA:CA	2.33	0.42
2:BL:246:ALA:O	2:BL:248:SER:N	2.52	0.42
3:BM:265:ILE:HG22	3:BM:266:HIS:H	1.84	0.42
5:BA:2:PHE:CB	5:BA:5:ASN:HD21	2.32	0.42
5:BF:13:LEU:CD1	14:BF:103:CRT:H1M1	2.25	0.42
9:BF:102:BCL:H111	9:BF:102:BCL:H192	2.00	0.42
6:BJ:18:HIS:O	6:BJ:22:MET:HB2	2.19	0.42
6:BJ:40:TRP:HA	6:BJ:44:PRO:HA	2.00	0.42
5:BK:2:PHE:O	5:BK:2:PHE:CD1	2.73	0.42
5:BO:9:TYR:CZ	5:BO:10:LYS:HD3	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:BO:20:VAL:C	5:BO:24:ILE:HG12	2.44	0.42
5:BO:54:SER:O	5:BO:58:LEU:N	2.40	0.42
9:BO:102:BCL:O1D	9:BO:102:BCL:C1A	2.64	0.42
6:BP:7:THR:OG1	6:BP:8:GLY:N	2.50	0.42
5:BU:8:LEU:O	5:BU:9:TYR:C	2.60	0.42
6:BV:29:PHE:O	6:BV:33:VAL:HB	2.19	0.42
6:BV:46:LEU:HD13	6:BX:42:TYR:CZ	2.54	0.42
5:BY:19:ARG:O	5:BY:20:VAL:C	2.62	0.42
5:BY:51:ILE:CA	5:BY:52:PRO:C	2.91	0.42
9:BZ:101:BCL:CGD	9:BZ:101:BCL:H2A	2.48	0.42
5:B1:12:TRP:CD1	6:B2:18:HIS:HA	2.54	0.42
6:B2:36:HIS:HE1	9:B2:101:BCL:C4D	2.31	0.42
14:B2:102:CRT:H2M2	5:B3:40:LEU:HD11	1.95	0.42
5:B3:13:LEU:HD21	6:B4:10:THR:C	2.44	0.42
5:B3:16:ASP:O	5:B3:20:VAL:HG22	2.19	0.42
1:AC:111:HIS:CE1	1:AC:124:LYS:HE2	2.52	0.42
1:AC:138:ASN:ND2	1:AC:149:GLY:HA3	2.34	0.42
1:AC:142:LYS:HE3	1:AC:147:GLU:OE2	2.19	0.42
1:AC:267:THR:C	1:AC:269:ALA:H	2.26	0.42
2:AL:135:GLY:C	2:AL:137:TYR:N	2.74	0.42
2:AL:181:ALA:HB3	2:AL:256:CYS:SG	2.58	0.42
2:AL:203:ILE:HA	2:AL:206:VAL:HG22	2.00	0.42
6:AB:23:GLN:CD	5:A9:4:MET:HE1	2.44	0.42
5:AD:43:ASP:O	5:AD:45:ASN:N	2.52	0.42
6:AG:40:TRP:HH2	6:AG:46:LEU:HD12	1.84	0.42
6:AJ:29:PHE:CE1	9:AJ:101:BCL:H11	2.53	0.42
5:AO:27:PHE:HE2	5:AQ:29:ILE:HD12	1.84	0.42
6:AP:31:LEU:HA	6:AP:34:ILE:CG2	2.49	0.42
5:AS:4:MET:HE3	5:AS:8:LEU:CD1	2.49	0.42
5:AS:8:LEU:CB	6:AT:18:HIS:CE1	3.02	0.42
5:AS:55:TYR:HD1	5:AS:56:GLN:N	2.17	0.42
5:AU:12:TRP:CD1	6:AV:18:HIS:HB2	2.54	0.42
5:AU:13:LEU:O	6:AV:7:THR:HA	2.19	0.42
5:AW:10:LYS:HD3	14:AW:102:CRT:H23	2.01	0.42
5:AW:27:PHE:CE1	5:AW:31:LEU:HD22	2.54	0.42
5:AY:18:ARG:HG2	5:AY:18:ARG:HH11	1.84	0.42
5:AY:50:ASN:ND2	5:AY:51:ILE:HG12	2.33	0.42
6:A2:20:ILE:HG23	6:A2:21:PHE:N	2.33	0.42
6:A2:40:TRP:CE3	6:A2:40:TRP:HA	2.54	0.42
5:A5:5:ASN:OD1	5:A5:8:LEU:HD12	2.19	0.42
9:A6:101:BCL:CGA	9:A6:101:BCL:H12	2.46	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:A8:19:ALA:O	6:A8:20:ILE:C	2.61	0.42
1:BC:126:VAL:O	1:BC:129:ARG:N	2.53	0.42
1:BC:135:ARG:HG2	1:BC:330:LEU:CA	2.49	0.42
1:BC:200:LEU:HG	1:BC:204:LEU:HD12	2.01	0.42
1:BC:271:TYR:O	1:BC:272:ALA:C	2.63	0.42
2:BL:55:THR:HA	2:BL:68:TYR:HE1	1.83	0.42
2:BL:79:ASP:C	2:BL:81:SER:H	2.28	0.42
2:BL:138:LEU:O	2:BL:142:PHE:N	2.52	0.42
2:BL:151:TRP:O	2:BL:153:HIS:N	2.52	0.42
2:BL:196:LEU:HB2	3:BM:216:PHE:CG	2.54	0.42
3:BM:33:ARG:O	3:BM:34:PRO:O	2.38	0.42
3:BM:124:LEU:O	3:BM:127:LEU:N	2.52	0.42
4:BH:63:ASP:O	4:BH:79:PRO:HD2	2.19	0.42
4:BH:171:TRP:CZ3	4:BH:231:VAL:HG12	2.54	0.42
4:BH:178:GLN:NE2	4:BH:180:ARG:CZ	2.82	0.42
4:BH:203:ASP:C	4:BH:205:LYS:N	2.75	0.42
4:BH:248:LEU:O	4:BH:248:LEU:HD23	2.19	0.42
5:BA:32:GLY:O	5:BA:33:LEU:C	2.62	0.42
6:BG:10:THR:CG2	6:BG:11:ASP:N	2.82	0.42
5:BK:16:ASP:HA	5:BK:17:PRO:HD3	1.85	0.42
5:BO:43:ASP:C	5:BQ:48:ASP:HB3	2.40	0.42
5:BQ:38:ILE:O	5:BQ:42:THR:CG2	2.67	0.42
9:BT:101:BCL:CHB	9:BU:102:BCL:HMB2	2.50	0.42
5:BU:38:ILE:CD1	14:BV:102:CRT:C40	2.98	0.42
5:BU:56:GLN:CD	5:BU:56:GLN:C	2.87	0.42
5:B1:9:TYR:HB2	6:B2:15:LYS:HA	2.02	0.42
9:B5:102:BCL:H92	6:B6:28:TRP:CE3	2.54	0.42
9:B7:103:BCL:CMD	6:B8:36:HIS:HD2	2.31	0.42
5:B9:39:VAL:C	5:B9:41:SER:H	2.26	0.42
6:B0:10:THR:H	6:B0:13:GLU:HG2	1.83	0.42
6:B0:21:PHE:HB2	14:B0:101:CRT:H14	1.93	0.42
6:B0:27:ALA:O	6:B0:31:LEU:HG	2.18	0.42
1:AC:90:PHE:HD1	1:AC:91:THR:N	2.17	0.42
1:AC:280:ASN:CG	1:AC:304:ARG:HD2	2.45	0.42
2:AL:18:ILE:HG23	4:AH:259:LEU:CB	2.50	0.42
2:AL:159:ILE:HD12	2:AL:159:ILE:N	2.34	0.42
2:AL:168:ASN:O	2:AL:169:VAL:C	2.61	0.42
2:AL:178:TYR:O	2:AL:272:TRP:NE1	2.52	0.42
3:AM:200:PRO:C	3:AM:203:MET:HG2	2.44	0.42
3:AM:212:SER:C	3:AM:214:LEU:H	2.27	0.42
4:AH:28:ILE:O	4:AH:31:ARG:HB2	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:AH:142:PHE:CD2	4:AH:172:VAL:HG21	2.55	0.42
4:AH:222:VAL:O	4:AH:223:PRO:C	2.60	0.42
5:AA:22:VAL:C	5:AA:24:ILE:N	2.75	0.42
5:AA:24:ILE:HA	5:AA:27:PHE:HB3	2.01	0.42
5:AF:53:VAL:O	5:AF:54:SER:C	2.62	0.42
9:AF:102:BCL:H121	9:AF:102:BCL:H8	1.92	0.42
6:AJ:17:PHE:O	6:AJ:20:ILE:CG2	2.61	0.42
9:AJ:101:BCL:H3A	9:AJ:101:BCL:HBA1	1.70	0.42
5:AK:12:TRP:CE2	6:AN:17:PHE:HD2	2.37	0.42
5:AO:11:ILE:HG12	14:AR:102:CRT:C8	2.46	0.42
6:AV:32:VAL:HG11	9:AV:102:BCL:CBA	2.48	0.42
6:AX:13:GLU:HG2	6:AX:14:ALA:N	2.34	0.42
6:A2:38:LEU:HD23	6:A2:38:LEU:O	2.20	0.42
5:A3:5:ASN:HA	5:A3:8:LEU:HD12	2.02	0.42
5:A3:19:ARG:HH21	5:A3:19:ARG:CG	2.32	0.42
5:A3:38:ILE:HG23	5:A3:39:VAL:HG23	2.01	0.42
9:A3:103:BCL:H2	6:A4:28:TRP:CH2	2.54	0.42
9:A3:103:BCL:C3D	9:A3:104:BCL:C3D	2.97	0.42
9:A5:102:BCL:HMB1	9:A5:102:BCL:HBB3	2.01	0.42
6:A6:31:LEU:O	6:A6:34:ILE:HG23	2.19	0.42
5:A7:43:ASP:OD2	5:A9:47:LEU:HD13	2.20	0.42
6:A8:26:TYR:O	6:A8:30:GLY:N	2.49	0.42
6:A8:37:LEU:O	6:A8:41:LEU:HG	2.20	0.42
1:BC:93:THR:O	1:BC:94:MET:C	2.62	0.42
1:BC:135:ARG:C	1:BC:137:ALA:N	2.75	0.42
1:BC:269:ALA:HB2	7:BC:504:HEM:CMA	2.49	0.42
1:BC:270:TRP:CG	3:BM:316:PRO:HG3	2.55	0.42
2:BL:99:THR:HG23	2:BL:137:TYR:OH	2.19	0.42
2:BL:109:TRP:HZ3	13:BM:405:MQ8:H342	1.84	0.42
2:BL:184:LEU:HD12	2:BL:184:LEU:N	2.34	0.42
4:BH:27:ILE:CG2	4:BH:28:ILE:N	2.83	0.42
4:BH:138:VAL:CG1	4:BH:140:LYS:HZ3	2.32	0.42
4:BH:159:LEU:HB3	4:BH:212:ASP:HA	2.01	0.42
4:BH:219:PHE:HA	4:BH:222:VAL:HG23	2.00	0.42
6:BB:15:LYS:O	6:BB:16:GLU:C	2.60	0.42
9:BD:102:BCL:HAC2	9:BE:101:BCL:CBC	2.49	0.42
6:BG:8:GLY:O	6:BG:9:LEU:CD2	2.68	0.42
5:BI:46:TRP:NE1	5:BI:47:LEU:CD1	2.81	0.42
9:BV:101:BCL:C4A	9:BW:102:BCL:HMB3	2.49	0.42
5:BW:33:LEU:O	5:BW:37:MET:HB2	2.20	0.42
5:BW:43:ASP:OD1	5:BW:44:LEU:N	2.53	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
14:BW:103:CRT:H16	6:BZ:21:PHE:CE2	2.54	0.42
6:BZ:18:HIS:O	6:BZ:22:MET:HB2	2.19	0.42
5:B3:44:LEU:HD13	5:B3:46:TRP:CE3	2.54	0.42
5:B5:32:GLY:HA2	9:B5:102:BCL:O1A	2.18	0.42
5:B5:45:ASN:O	5:B5:46:TRP:C	2.63	0.42
9:B6:101:BCL:HBA1	9:B6:101:BCL:H3A	1.66	0.42
5:B7:42:THR:OG1	5:B9:48:ASP:OD1	2.31	0.42
1:AC:102:SER:C	1:AC:104:LYS:N	2.77	0.42
2:AL:36:GLY:HA2	2:AL:112:ARG:HD3	2.01	0.42
2:AL:111:LEU:O	2:AL:112:ARG:C	2.60	0.42
2:AL:146:LEU:C	2:AL:148:MET:N	2.77	0.42
2:AL:280:LEU:HD11	5:AY:37:MET:HE1	2.01	0.42
3:AM:47:GLN:HG2	3:AM:48:ILE:N	2.35	0.42
3:AM:134:TYR:CE2	3:AM:144:GLN:HG3	2.55	0.42
3:AM:162:PHE:CE1	5:AO:37:MET:SD	3.13	0.42
3:AM:185:TRP:CH2	3:AM:189:PHE:CD1	3.07	0.42
3:AM:206:ILE:O	3:AM:207:ALA:C	2.63	0.42
3:AM:206:ILE:CG1	9:AM:402:BCL:HMA2	2.48	0.42
4:AH:36:ARG:HE	4:AH:65:LYS:CD	2.31	0.42
4:AH:39:TYR:CD1	4:AH:40:PRO:HA	2.54	0.42
4:AH:45:ARG:CA	4:AH:96:PRO:HB3	2.48	0.42
4:AH:142:PHE:HZ	4:AH:173:ASP:O	2.02	0.42
5:AA:8:LEU:O	5:AA:11:ILE:HG22	2.19	0.42
9:AB:101:BCL:H3A	9:AB:101:BCL:HBA1	1.76	0.42
5:AD:36:HIS:CE1	9:AE:101:BCL:HMD2	2.52	0.42
9:AD:102:BCL:CAC	9:AE:101:BCL:HBC3	2.49	0.42
5:AF:11:ILE:HD12	5:AF:14:ILE:CG1	2.50	0.42
9:AJ:101:BCL:HMB1	9:AJ:101:BCL:CBB	2.50	0.42
5:AQ:35:ILE:O	5:AQ:38:ILE:HG22	2.19	0.42
14:AR:102:CRT:H391	5:AS:36:HIS:CG	2.54	0.42
5:AU:14:ILE:CB	14:AX:102:CRT:H82	2.49	0.42
9:AV:102:BCL:HMA1	9:AW:101:BCL:HMA1	2.00	0.42
5:A5:49:ASP:CG	5:A5:50:ASN:N	2.77	0.42
6:A8:38:LEU:HD23	6:A8:38:LEU:C	2.45	0.42
5:A9:12:TRP:CE2	6:A0:17:PHE:CE1	3.07	0.42
5:A9:31:LEU:O	5:A9:35:ILE:HG12	2.20	0.42
1:BC:115:ASN:OD1	1:BC:115:ASN:N	2.51	0.42
1:BC:266:ARG:HG3	7:BC:503:HEM:C2D	2.54	0.42
1:BC:314:VAL:HG12	1:BC:315:ASN:H	1.85	0.42
2:BL:113:GLU:OE1	2:BL:127:PRO:HG3	2.19	0.42
2:BL:168:ASN:C	2:BL:170:GLY:N	2.77	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:BL:177:HIS:CD2	9:BL:301:BCL:HMC2	2.54	0.42
2:BL:252:TRP:CZ2	11:BL:304:UQ8:H30B	2.54	0.42
3:BM:14:ARG:NH1	4:BH:145:ALA:HA	2.35	0.42
3:BM:151:ALA:O	3:BM:152:ALA:C	2.62	0.42
3:BM:208:PHE:O	3:BM:210:TYR:N	2.51	0.42
3:BM:224:LEU:O	3:BM:225:SER:C	2.62	0.42
4:BH:167:VAL:HA	4:BH:183:GLU:O	2.19	0.42
5:BA:45:ASN:O	5:BA:49:ASP:HB3	2.19	0.42
9:BA:101:BCL:C1D	9:BB:101:BCL:HMD1	2.50	0.42
5:BI:12:TRP:CH2	6:BJ:17:PHE:CE1	3.08	0.42
5:BI:50:ASN:CG	6:BJ:43:ARG:HH22	2.27	0.42
5:BK:11:ILE:HG12	14:BP:102:CRT:C8	2.34	0.42
5:BQ:42:THR:HG23	5:BQ:43:ASP:N	2.19	0.42
5:BQ:42:THR:HG21	5:BS:47:LEU:HG	2.02	0.42
5:BY:44:LEU:HD12	5:BY:44:LEU:O	2.20	0.42
9:B6:101:BCL:C1B	9:B7:103:BCL:CMB	2.97	0.42
6:B8:19:ALA:O	6:B8:20:ILE:C	2.61	0.42
9:B8:101:BCL:HMA3	9:B9:102:BCL:HMA1	1.98	0.42
5:B9:16:ASP:OD1	5:B9:18:ARG:N	2.52	0.42
1:AC:21:LEU:HG	2:AL:259:ILE:HG21	2.00	0.42
1:AC:176:SER:CB	5:AS:42:THR:HA	2.50	0.42
2:AL:18:ILE:O	2:AL:18:ILE:HG22	2.18	0.42
2:AL:24:ASP:O	2:AL:24:ASP:CG	2.62	0.42
2:AL:38:VAL:CG2	2:AL:39:GLY:H	2.30	0.42
2:AL:56:ILE:O	2:AL:66:GLN:HG3	2.19	0.42
2:AL:63:SER:C	2:AL:65:LEU:H	2.28	0.42
2:AL:164:ASP:C	2:AL:166:VAL:N	2.73	0.42
3:AM:84:PHE:CE2	5:AW:37:MET:HA	2.54	0.42
3:AM:210:TYR:O	3:AM:211:GLY:C	2.62	0.42
3:AM:230:GLY:CA	17:AM:502:HOH:O	2.67	0.42
3:AM:310:VAL:O	3:AM:310:VAL:HG12	2.19	0.42
4:AH:47:GLU:HG3	5:AA:19:ARG:HG3	2.00	0.42
5:AA:29:ILE:HD11	14:A0:101:CRT:C34	2.45	0.42
5:AF:3:THR:O	5:AF:4:MET:HB2	2.18	0.42
5:AI:36:HIS:O	5:AI:40:LEU:CD1	2.68	0.42
5:AO:45:ASN:HB3	5:AO:48:ASP:OD1	2.20	0.42
6:AP:21:PHE:CE1	14:AP:102:CRT:H19	2.55	0.42
5:AS:31:LEU:HG	9:AT:101:BCL:HED3	2.01	0.42
5:A1:43:ASP:HB2	5:A3:47:LEU:HD13	2.01	0.42
14:A1:103:CRT:H20	14:A1:103:CRT:H181	1.89	0.42
5:A3:14:ILE:CG2	5:A5:17:PRO:HB2	2.49	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:A3:43:ASP:OD1	5:A3:44:LEU:N	2.52	0.42
5:A7:28:GLN:O	9:A8:101:BCL:HED1	2.19	0.42
1:BC:82:LEU:HD13	1:BC:93:THR:HG21	2.01	0.42
2:BL:108:SER:O	2:BL:111:LEU:N	2.53	0.42
2:BL:111:LEU:O	2:BL:114:VAL:N	2.53	0.42
2:BL:181:ALA:CB	2:BL:256:CYS:HA	2.48	0.42
2:BL:266:ARG:HG3	2:BL:266:ARG:NH1	2.35	0.42
3:BM:66:VAL:HG11	3:BM:121:PHE:CD2	2.53	0.42
3:BM:152:ALA:O	3:BM:277:VAL:HB	2.19	0.42
3:BM:221:ALA:O	3:BM:225:SER:N	2.36	0.42
3:BM:299:VAL:HB	3:BM:304:ALA:CB	2.36	0.42
9:BM:402:BCL:C14	15:BQ:101:PEF:H442	2.39	0.42
4:BH:32:ARG:C	4:BH:34:ASP:N	2.76	0.42
4:BH:215:LYS:H	4:BH:218:HIS:CD2	2.32	0.42
5:BA:9:TYR:CD1	5:BA:9:TYR:C	2.97	0.42
5:BA:35:ILE:O	5:BA:37:MET:N	2.53	0.42
6:BB:20:ILE:HG12	5:B9:7:ASN:HB3	2.01	0.42
14:BB:102:CRT:C2M	5:BD:37:MET:CE	2.86	0.42
5:BD:51:ILE:HG22	5:BD:52:PRO:CA	2.50	0.42
6:BG:32:VAL:CG1	9:BG:101:BCL:HBA2	2.29	0.42
9:BG:101:BCL:C20	6:BJ:38:LEU:HD21	2.49	0.42
5:BI:29:ILE:HG23	5:BI:30:VAL:N	2.33	0.42
5:BK:44:LEU:HD22	5:BK:44:LEU:C	2.45	0.42
6:BN:34:ILE:HD13	6:BN:34:ILE:O	2.19	0.42
5:BO:7:ASN:HB3	6:BR:20:ILE:CD1	2.39	0.42
5:BS:46:TRP:CZ3	9:BS:102:BCL:CBC	3.03	0.42
5:BS:49:ASP:CG	5:BS:50:ASN:H	2.27	0.42
9:BU:102:BCL:C3D	9:BV:101:BCL:CMD	2.98	0.42
6:BV:45:TRP:CD1	6:BV:46:LEU:N	2.88	0.42
5:BW:46:TRP:CZ2	9:BW:102:BCL:H2C	2.54	0.42
6:BX:29:PHE:CE2	9:BX:101:BCL:H43	2.54	0.42
5:BY:38:ILE:HD12	5:BY:38:ILE:C	2.44	0.42
5:B1:33:LEU:C	5:B1:33:LEU:HD12	2.45	0.42
5:B7:17:PRO:HG2	5:B7:18:ARG:H	1.85	0.42
1:AC:46:LYS:C	1:AC:48:GLN:H	2.28	0.42
1:AC:48:GLN:C	1:AC:50:ALA:H	2.28	0.42
2:AL:123:GLY:HA2	3:AM:228:ARG:NH2	2.34	0.42
2:AL:129:ALA:CB	2:AL:247:LEU:HD11	2.50	0.42
2:AL:170:GLY:HA3	9:AL:301:BCL:HBC3	2.02	0.42
9:AL:303:BCL:CED	3:AM:203:MET:HB2	2.50	0.42
3:AM:41:GLY:HA2	3:AM:44:GLY:O	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:AM:105:ARG:HA	5:AO:42:THR:CG2	2.28	0.42
3:AM:122:LEU:O	3:AM:126:ILE:HD12	2.19	0.42
3:AM:260:VAL:HG23	3:AM:261:THR:N	2.34	0.42
3:AM:268:TRP:CG	4:AH:30:LEU:HD13	2.54	0.42
3:AM:317:TYR:CD1	3:AM:317:TYR:N	2.88	0.42
5:AA:20:VAL:HA	5:AA:23:SER:HB3	2.00	0.42
5:AD:27:PHE:O	5:AD:30:VAL:HG12	2.20	0.42
5:AF:9:TYR:HA	6:AG:18:HIS:CG	2.55	0.42
6:AG:28:TRP:O	6:AG:30:GLY:N	2.52	0.42
9:AJ:101:BCL:HMB1	9:AJ:101:BCL:HBB3	2.00	0.42
5:AK:19:ARG:HG3	5:AK:20:VAL:N	2.35	0.42
9:AK:102:BCL:CMD	9:AN:101:BCL:CHD	2.86	0.42
5:AO:9:TYR:HB2	6:AP:18:HIS:CG	2.54	0.42
6:AR:45:TRP:CE3	9:AR:101:BCL:H2C	2.54	0.42
6:AX:28:TRP:CE3	6:AX:31:LEU:HD12	2.53	0.42
9:A3:104:BCL:CGD	9:A3:104:BCL:H2A	2.50	0.42
5:A5:53:VAL:HA	5:A5:56:GLN:HG2	2.01	0.42
6:A6:33:VAL:O	6:A6:37:LEU:HB2	2.20	0.42
1:BC:24:GLU:CD	2:BL:266:ARG:NH2	2.76	0.42
1:BC:59:VAL:O	1:BC:59:VAL:HG23	2.20	0.42
1:BC:112:VAL:O	1:BC:114:GLY:N	2.52	0.42
1:BC:154:THR:HG22	1:BC:155:CYS:N	2.35	0.42
1:BC:155:CYS:O	1:BC:162:PRO:CB	2.66	0.42
1:BC:196:PRO:CG	1:BC:231:TRP:CD1	3.02	0.42
1:BC:269:ALA:O	1:BC:273:ILE:CG1	2.65	0.42
2:BL:131:SER:O	2:BL:132:PHE:C	2.62	0.42
2:BL:266:ARG:HH11	2:BL:266:ARG:CG	2.33	0.42
3:BM:35:ILE:CG2	3:BM:36:PHE:N	2.82	0.42
3:BM:122:LEU:C	3:BM:157:TYR:HH	2.22	0.42
3:BM:126:ILE:HD12	3:BM:157:TYR:CE2	2.54	0.42
3:BM:179:ILE:N	3:BM:179:ILE:CD1	2.68	0.42
3:BM:180:PHE:O	3:BM:181:PRO:C	2.61	0.42
3:BM:282:ILE:O	3:BM:283:GLY:C	2.62	0.42
4:BH:35:LYS:NZ	4:BH:57:GLY:CA	2.73	0.42
4:BH:129:GLY:O	4:BH:130:LEU:C	2.62	0.42
5:BF:12:TRP:CE3	5:BF:12:TRP:CA	3.03	0.42
6:BG:28:TRP:O	6:BG:30:GLY:N	2.52	0.42
6:BJ:45:TRP:O	6:BJ:46:LEU:CB	2.66	0.42
5:BK:45:ASN:C	5:BK:49:ASP:CB	2.85	0.42
5:BO:36:HIS:O	5:BO:40:LEU:HB3	2.20	0.42
6:BP:12:ASP:O	6:BP:16:GLU:HG3	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:BQ:50:ASN:HB3	5:BS:56:GLN:CG	2.49	0.42
5:BU:28:GLN:O	9:BU:102:BCL:H11	2.18	0.42
5:BY:34:LEU:O	5:BY:37:MET:HB2	2.19	0.42
5:B1:39:VAL:HG13	5:B1:40:LEU:N	2.33	0.42
6:B2:31:LEU:O	6:B2:34:ILE:HG22	2.19	0.42
5:B3:46:TRP:CZ3	9:B3:102:BCL:CAC	3.02	0.42
6:B4:45:TRP:O	6:B4:46:LEU:CG	2.67	0.42
6:B8:22:MET:O	6:B8:26:TYR:CD2	2.71	0.42
5:B9:27:PHE:CD1	5:B9:27:PHE:C	2.98	0.42
1:AC:138:ASN:O	1:AC:142:LYS:HG2	2.20	0.42
1:AC:165:ALA:HB1	1:AC:303:LEU:CB	2.22	0.42
1:AC:259:TRP:O	1:AC:260:THR:C	2.60	0.42
1:AC:313:ALA:C	1:AC:314:VAL:HG23	2.44	0.42
2:AL:159:ILE:H	2:AL:159:ILE:HD12	1.84	0.42
2:AL:273:ASN:HA	2:AL:276:LEU:CD2	2.34	0.42
3:AM:78:SER:C	3:AM:80:HIS:H	2.24	0.42
3:AM:234:GLU:OE1	3:AM:266:HIS:HE1	1.93	0.42
9:AM:402:BCL:HMB3	9:AM:402:BCL:CBB	2.49	0.42
4:AH:35:LYS:HZ1	4:AH:57:GLY:HA3	1.79	0.42
4:AH:121:LYS:HB2	4:BH:122:HIS:CD2	2.55	0.42
4:AH:130:LEU:HG	4:AH:131:PRO:CD	2.40	0.42
9:AA:101:BCL:HMB1	9:AA:101:BCL:CBB	2.50	0.42
9:AA:101:BCL:HMB1	9:A0:102:BCL:NB	2.34	0.42
14:AB:102:CRT:H342	9:AD:102:BCL:CGA	2.50	0.42
6:AE:33:VAL:HG22	6:AE:37:LEU:HD23	2.02	0.42
5:AI:30:VAL:HG13	5:AI:31:LEU:N	2.35	0.42
9:AJ:101:BCL:HBB3	9:AK:102:BCL:C4B	2.50	0.42
6:AP:17:PHE:HA	6:AP:20:ILE:CG2	2.50	0.42
5:AQ:36:HIS:NE2	9:AR:101:BCL:HMD3	2.35	0.42
9:AQ:102:BCL:C4D	9:AR:101:BCL:HMD3	2.49	0.42
6:AT:45:TRP:CE3	9:AT:101:BCL:HAC2	2.54	0.42
5:AU:42:THR:O	5:AU:43:ASP:C	2.63	0.42
5:AW:27:PHE:HE1	5:AW:31:LEU:HD22	1.84	0.42
5:AY:45:ASN:O	5:AY:46:TRP:C	2.63	0.42
14:A5:103:CRT:C30	5:A7:31:LEU:HD21	2.49	0.42
6:A6:17:PHE:CD2	14:A7:102:CRT:H6	2.54	0.42
6:A0:16:GLU:CD	14:A0:101:CRT:H41	2.45	0.42
1:BC:141:TRP:CZ3	1:BC:275:HIS:HA	2.55	0.42
1:BC:316:LYS:O	1:BC:317:PRO:O	2.37	0.42
2:BL:82:TYR:HA	2:BL:85:ARG:NE	2.33	0.42
2:BL:155:PHE:HB3	2:BL:165:TRP:CE2	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
11:BL:304:UQ8:H22	11:BL:304:UQ8:H25	1.80	0.42
3:BM:52:TYR:CD2	3:BM:136:ARG:NE	2.82	0.42
3:BM:264:SER:O	3:BM:267:ARG:HB2	2.20	0.42
4:BH:132:LYS:NZ	4:BH:173:ASP:OD2	2.46	0.42
5:BA:40:LEU:HD12	5:BA:40:LEU:O	2.19	0.42
6:BE:13:GLU:CD	6:BE:13:GLU:H	2.28	0.42
5:BF:9:TYR:CZ	5:BF:10:LYS:HD3	2.48	0.42
5:BF:48:ASP:OD1	5:BF:48:ASP:N	2.52	0.42
9:BG:101:BCL:CBB	9:BG:101:BCL:HMB1	2.49	0.42
5:BK:4:MET:O	5:BK:8:LEU:HG	2.19	0.42
5:BK:12:TRP:CE3	5:BK:12:TRP:CA	3.02	0.42
9:BK:102:BCL:C3D	9:BN:101:BCL:C2D	2.98	0.42
6:BN:32:VAL:CG2	9:BN:101:BCL:HBA2	2.50	0.42
5:BQ:43:ASP:CA	5:BS:47:LEU:HB3	2.50	0.42
6:BR:10:THR:CG2	6:BR:11:ASP:N	2.82	0.42
5:BS:52:PRO:O	5:BS:53:VAL:C	2.62	0.42
6:BZ:40:TRP:HB2	6:BZ:45:TRP:CZ3	2.54	0.42
9:BZ:101:BCL:HBB3	9:BZ:101:BCL:HMB1	2.02	0.42
6:B2:17:PHE:O	6:B2:18:HIS:C	2.63	0.42
14:B5:103:CRT:H401	5:B7:38:ILE:HG21	1.99	0.42
6:B6:31:LEU:O	6:B6:34:ILE:HG23	2.20	0.42
6:B8:23:GLN:HE21	6:B8:23:GLN:HB2	1.63	0.42
6:B0:42:TYR:CE2	6:B0:43:ARG:HD2	2.54	0.42
1:AC:116:TRP:CD1	1:AC:116:TRP:N	2.87	0.42
1:AC:317:PRO:O	1:AC:318:LEU:HB2	2.20	0.42
2:AL:38:VAL:CG2	2:AL:39:GLY:N	2.82	0.42
2:AL:82:TYR:HB3	2:AL:85:ARG:CG	2.50	0.42
2:AL:116:ILE:HG22	2:AL:117:CYS:N	2.35	0.42
2:AL:147:LEU:N	2:AL:147:LEU:CD1	2.83	0.42
2:AL:281:TRP:HD1	2:AL:281:TRP:OXT	2.02	0.42
9:AL:301:BCL:HBB3	9:AM:401:BCL:HMD2	2.02	0.42
3:AM:35:ILE:HD11	15:AM:409:PEF:H312	2.00	0.42
3:AM:250:LEU:O	3:AM:251:PHE:C	2.63	0.42
9:AM:401:BCL:H141	9:AM:401:BCL:HMA1	2.01	0.42
14:AM:406:CRT:H2M1	5:AO:41:SER:OG	2.20	0.42
4:AH:149:PRO:HG3	4:AH:204:LYS:HB3	2.02	0.42
4:AH:173:ASP:OD1	4:AH:175:SER:N	2.31	0.42
4:AH:203:ASP:C	4:AH:205:LYS:H	2.28	0.42
5:AA:44:LEU:HD12	5:AA:46:TRP:N	2.34	0.42
5:AF:28:GLN:CB	9:AF:102:BCL:C1	2.68	0.42
5:AI:39:VAL:O	5:AI:43:ASP:CB	2.68	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:AO:13:LEU:C	6:AP:7:THR:HA	2.41	0.42
6:AV:20:ILE:HG23	6:AV:21:PHE:N	2.35	0.42
6:AX:18:HIS:CE1	6:AX:22:MET:HE2	2.55	0.42
5:AY:10:LYS:HB3	14:A2:102:CRT:H23	2.01	0.42
5:A1:12:TRP:CD1	6:A2:18:HIS:CB	3.03	0.42
5:A3:56:GLN:HG2	5:A3:57:ALA:H	1.85	0.42
6:A4:46:LEU:HB2	5:A5:52:PRO:HD2	2.02	0.42
6:A8:45:TRP:CZ3	9:A8:101:BCL:HAC2	2.55	0.42
5:A9:46:TRP:CZ2	9:A9:102:BCL:H2C	2.55	0.42
9:A0:102:BCL:H13	9:A0:102:BCL:HMB3	2.01	0.42
1:BC:20:LEU:CD2	1:BC:21:LEU:N	2.75	0.42
1:BC:53:ILE:HG22	1:BC:54:GLN:N	2.35	0.42
1:BC:132:GLU:O	1:BC:136:ALA:HB2	2.20	0.42
1:BC:236:MET:HG3	7:BC:503:HEM:CHB	2.49	0.42
1:BC:237:MET:O	1:BC:238:ASN:C	2.63	0.42
1:BC:306:SER:O	1:BC:309:THR:N	2.52	0.42
2:BL:44:LEU:HD22	2:BL:48:LEU:HD12	2.01	0.42
2:BL:75:ILE:HD12	2:BL:94:LEU:HD22	2.01	0.42
2:BL:225:PHE:O	2:BL:229:VAL:HG22	2.19	0.42
3:BM:84:PHE:CE2	5:BW:37:MET:HG2	2.54	0.42
3:BM:161:GLY:HA3	14:BM:406:CRT:H291	2.00	0.42
3:BM:214:LEU:O	3:BM:218:MET:CG	2.67	0.42
9:BM:402:BCL:H141	9:BM:402:BCL:H161	1.85	0.42
5:BA:33:LEU:O	14:B0:101:CRT:H2M3	2.20	0.42
9:BA:101:BCL:CMD	6:BB:36:HIS:CE1	3.03	0.42
5:BO:17:PRO:HG2	5:BO:18:ARG:H	1.85	0.42
5:BY:11:ILE:CG2	5:BY:15:LEU:HD12	2.50	0.42
5:B3:17:PRO:O	5:B3:21:LEU:CB	2.67	0.42
9:B4:101:BCL:HBA1	9:B4:101:BCL:H3A	1.56	0.42
6:B6:33:VAL:O	6:B6:37:LEU:HB2	2.20	0.42
9:B7:103:BCL:HBB3	9:B7:103:BCL:HMB1	2.01	0.42
6:B0:22:MET:O	6:B0:26:TYR:HD2	2.03	0.42
1:AC:295:ARG:HD2	7:AC:502:HEM:CGD	2.50	0.42
2:AL:3:MET:O	4:AH:41:LEU:HG	2.19	0.42
2:AL:13:ARG:CA	4:AH:99:PRO:HB2	2.50	0.42
2:AL:59:THR:OG1	2:AL:65:LEU:HD12	2.20	0.42
2:AL:174:LEU:N	2:AL:174:LEU:CD1	2.83	0.42
2:AL:226:ARG:NE	3:AM:47:GLN:NE2	2.68	0.42
3:AM:234:GLU:O	3:AM:237:GLN:N	2.53	0.42
3:AM:268:TRP:HZ2	4:AH:34:ASP:OD2	2.02	0.42
9:AA:101:BCL:CAC	9:AB:101:BCL:HBC3	2.50	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
14:AB:102:CRT:H10	14:AB:102:CRT:H81	1.88	0.42
5:AF:44:LEU:O	5:AF:44:LEU:CD1	2.62	0.42
6:AG:25:MET:SD	6:AG:29:PHE:CZ	3.13	0.42
9:AG:101:BCL:HBB2	9:AG:101:BCL:HMB3	2.02	0.42
6:AN:38:LEU:HD23	6:AN:38:LEU:C	2.44	0.42
9:AP:101:BCL:HBA1	9:AP:101:BCL:H3A	1.76	0.42
14:AP:102:CRT:H2M3	5:AQ:36:HIS:CB	2.50	0.42
5:AQ:19:ARG:O	5:AQ:22:VAL:HG12	2.20	0.42
6:AR:29:PHE:CD1	9:AR:101:BCL:H11	2.55	0.42
5:AY:38:ILE:HG13	5:AY:38:ILE:H	1.75	0.42
6:AZ:36:HIS:CE1	9:AZ:101:BCL:C4A	3.03	0.42
5:A1:12:TRP:HZ2	6:A2:21:PHE:CE2	2.25	0.42
5:A1:28:GLN:CB	9:A1:102:BCL:C2	2.94	0.42
5:A7:35:ILE:O	5:A7:39:VAL:HG23	2.20	0.42
5:A7:36:HIS:C	14:A7:102:CRT:H2M3	2.44	0.42
1:BC:159:ASN:HA	1:BC:160:PRO:HD3	1.77	0.42
2:BL:44:LEU:O	2:BL:48:LEU:HB2	2.19	0.42
2:BL:146:LEU:C	2:BL:148:MET:N	2.77	0.42
2:BL:164:ASP:C	2:BL:166:VAL:H	2.28	0.42
2:BL:175:HIS:CE1	3:BM:184:ASP:OD2	2.73	0.42
2:BL:225:PHE:C	2:BL:227:ASP:H	2.27	0.42
3:BM:84:PHE:O	3:BM:87:LEU:HB2	2.20	0.42
3:BM:219:HIS:CE1	3:BM:223:ILE:HD11	2.54	0.42
4:BH:93:SER:O	4:BH:98:SER:HB2	2.20	0.42
9:BB:101:BCL:H92	9:BD:102:BCL:HED3	2.02	0.42
5:BD:46:TRP:CE2	9:BD:102:BCL:H2C	2.54	0.42
6:BG:21:PHE:CE1	14:BG:102:CRT:H16	2.54	0.42
6:BG:46:LEU:HB3	6:BJ:42:TYR:OH	2.18	0.42
9:BG:101:BCL:HBB3	9:BG:101:BCL:HMB1	2.02	0.42
5:BI:45:ASN:O	5:BI:47:LEU:N	2.53	0.42
5:BI:56:GLN:HG3	5:BI:56:GLN:H	1.66	0.42
6:BJ:21:PHE:CE1	6:BJ:25:MET:HE3	2.55	0.42
6:BJ:29:PHE:N	6:BJ:29:PHE:HD1	2.18	0.42
5:BK:47:LEU:H	5:BK:47:LEU:CD2	2.29	0.42
5:BK:55:TYR:HD1	5:BK:56:GLN:N	2.18	0.42
5:BO:32:GLY:N	9:BP:101:BCL:HED2	2.35	0.42
9:BO:102:BCL:CAD	9:BP:101:BCL:CAD	2.98	0.42
6:BP:27:ALA:O	6:BP:31:LEU:CG	2.67	0.42
5:BQ:42:THR:CG2	5:BS:47:LEU:HG	2.49	0.42
6:BR:45:TRP:CD1	6:BR:46:LEU:H	2.37	0.42
6:BV:28:TRP:HA	6:BV:31:LEU:HD12	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:BV:101:BCL:HMB1	9:BV:101:BCL:HBB3	2.02	0.42
5:BY:28:GLN:HB3	9:BY:102:BCL:H12	2.02	0.42
5:BY:38:ILE:HD12	5:BY:39:VAL:HG23	2.02	0.42
5:B3:5:ASN:CA	5:B3:8:LEU:HD12	2.38	0.42
5:B9:8:LEU:HD22	5:B9:11:ILE:HD11	2.02	0.42
9:B9:102:BCL:CMD	6:B0:36:HIS:CD2	2.96	0.42
9:B0:102:BCL:HMB1	9:B0:102:BCL:HBB3	2.01	0.42
1:AC:105:GLU:N	1:AC:105:GLU:CD	2.78	0.41
1:AC:153:TYR:O	1:AC:157:ARG:HG2	2.20	0.41
1:AC:274:ARG:O	1:AC:277:ARG:HB2	2.20	0.41
2:AL:6:PHE:CE1	3:AM:241:ARG:NH1	2.88	0.41
2:AL:196:LEU:HB2	3:AM:216:PHE:CB	2.50	0.41
2:AL:225:PHE:C	2:AL:227:ASP:N	2.78	0.41
3:AM:59:LEU:CD2	3:AM:128:LEU:HD21	2.49	0.41
3:AM:179:ILE:O	3:AM:182:HIS:ND1	2.49	0.41
14:AM:406:CRT:H10	14:AM:406:CRT:H81	1.87	0.41
4:AH:249:TYR:C	4:AH:251:THR:H	2.28	0.41
5:AA:45:ASN:O	5:AA:47:LEU:N	2.53	0.41
5:AD:5:ASN:O	5:AD:6:ALA:C	2.62	0.41
6:AG:36:HIS:ND1	9:AG:101:BCL:H142	2.35	0.41
6:AN:32:VAL:CG1	9:AN:101:BCL:H141	2.46	0.41
5:AO:11:ILE:HG22	5:AO:15:LEU:HD12	2.02	0.41
5:AS:12:TRP:HZ3	5:AS:20:VAL:HG11	1.85	0.41
5:AS:46:TRP:CD1	5:AS:47:LEU:N	2.88	0.41
5:AU:12:TRP:CD1	6:AV:18:HIS:CA	3.03	0.41
9:AY:102:BCL:HHC	9:AY:102:BCL:OBB	2.19	0.41
5:A1:60:LYS:O	5:A1:61:LYS:OXT	2.38	0.41
6:A2:31:LEU:O	6:A2:34:ILE:HG22	2.20	0.41
5:A3:51:ILE:HB	5:A3:52:PRO:HA	2.02	0.41
5:A7:11:ILE:CG1	14:A0:101:CRT:H132	2.49	0.41
9:A0:102:BCL:HBB2	9:A0:102:BCL:HMB3	2.01	0.41
2:BL:38:VAL:CG2	2:BL:39:GLY:H	2.25	0.41
2:BL:139:VAL:HA	2:BL:143:VAL:CG2	2.50	0.41
9:BL:303:BCL:H11	3:BM:210:TYR:HB3	2.01	0.41
3:BM:8:PHE:O	3:BM:9:THR:C	2.63	0.41
3:BM:61:ILE:CG1	3:BM:129:TRP:HZ3	2.33	0.41
3:BM:133:THR:O	3:BM:134:TYR:C	2.63	0.41
4:BH:31:ARG:HH21	4:BH:34:ASP:CB	2.33	0.41
4:BH:186:VAL:HG22	4:BH:193:VAL:HG22	2.02	0.41
5:BD:16:ASP:OD2	5:BD:18:ARG:CG	2.68	0.41
5:BF:44:LEU:O	5:BF:46:TRP:N	2.49	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:BG:17:PHE:O	6:BG:21:PHE:HB3	2.19	0.41
6:BG:30:GLY:HA2	6:BG:33:VAL:CG1	2.50	0.41
5:BI:15:LEU:HD12	5:BI:20:VAL:HG11	2.01	0.41
6:BN:22:MET:HG3	6:BN:26:TYR:CE2	2.54	0.41
14:BO:103:CRT:H23	6:BR:16:GLU:CB	2.50	0.41
6:BP:38:LEU:C	6:BP:38:LEU:CD2	2.93	0.41
5:BS:30:VAL:CG1	5:BS:31:LEU:N	2.82	0.41
5:BU:35:ILE:HA	5:BU:35:ILE:HD13	1.84	0.41
5:BU:43:ASP:OD2	5:BW:47:LEU:HA	2.20	0.41
5:BW:27:PHE:CD1	5:BW:27:PHE:C	2.97	0.41
5:B9:36:HIS:CD2	9:B0:102:BCL:CMD	3.02	0.41
1:AC:65:ALA:CB	1:AC:89:GLU:OE1	2.68	0.41
1:AC:270:TRP:CE3	1:AC:271:TYR:CD1	3.08	0.41
1:AC:276:VAL:HG23	1:AC:280:ASN:ND2	2.34	0.41
2:AL:48:LEU:HD23	2:AL:51:VAL:HG21	2.01	0.41
2:AL:193:CYS:O	2:AL:193:CYS:SG	2.77	0.41
2:AL:202:LEU:O	2:AL:203:ILE:C	2.63	0.41
2:AL:203:ILE:CG2	3:AM:266:HIS:ND1	2.62	0.41
3:AM:28:LEU:H	3:AM:28:LEU:HD12	1.85	0.41
3:AM:90:PHE:C	3:AM:92:TRP:N	2.78	0.41
5:AF:18:ARG:HH11	5:AF:18:ARG:HG3	1.85	0.41
5:AF:45:ASN:O	5:AF:47:LEU:N	2.54	0.41
9:AF:102:BCL:HMB3	9:AF:102:BCL:HBB2	2.01	0.41
6:AJ:26:TYR:O	6:AJ:27:ALA:C	2.62	0.41
6:AN:18:HIS:O	6:AN:22:MET:CB	2.68	0.41
5:AS:30:VAL:HG22	15:AS:101:PEF:H412	0.57	0.41
6:AT:16:GLU:OE1	14:AT:102:CRT:H31A	2.19	0.41
6:AV:15:LYS:O	6:AV:18:HIS:HB3	2.20	0.41
5:AY:36:HIS:O	5:AY:40:LEU:CB	2.67	0.41
9:AZ:101:BCL:CBB	9:AZ:101:BCL:HMB1	2.50	0.41
5:A3:22:VAL:HA	5:A3:25:VAL:CG2	2.51	0.41
5:A5:44:LEU:HD12	5:A5:46:TRP:HB3	2.01	0.41
1:BC:153:TYR:CE1	1:BC:158:GLY:N	2.88	0.41
1:BC:233:PHE:O	1:BC:235:LEU:N	2.53	0.41
1:BC:259:TRP:CH2	7:BC:503:HEM:HAD1	2.56	0.41
2:BL:12:VAL:HB	4:BH:113:PRO:HD3	2.01	0.41
2:BL:72:ARG:HG2	3:BM:305:PRO:HA	2.01	0.41
2:BL:86:MET:HE1	2:BL:96:GLN:HB2	2.02	0.41
2:BL:267:GLY:O	2:BL:268:TRP:C	2.63	0.41
3:BM:98:PRO:CB	3:BM:171:TRP:HB3	2.48	0.41
3:BM:210:TYR:O	3:BM:213:ALA:HB3	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:BM:268:TRP:NE1	4:BH:30:LEU:HD22	2.35	0.41
3:BM:284:ILE:HD11	9:BM:402:BCL:CAD	2.51	0.41
5:BA:37:MET:CE	14:B0:101:CRT:H2M1	2.50	0.41
9:BG:101:BCL:HMA1	9:BI:102:BCL:HHB	2.02	0.41
5:BI:49:ASP:O	5:BK:60:LYS:CB	2.68	0.41
5:BI:52:PRO:O	5:BI:53:VAL:C	2.62	0.41
6:BJ:18:HIS:NE2	6:BJ:22:MET:HE2	2.35	0.41
5:BO:38:ILE:CG1	5:BO:39:VAL:N	2.73	0.41
6:BP:22:MET:O	6:BP:25:MET:HB3	2.20	0.41
9:BQ:104:BCL:H11	6:BR:29:PHE:HA	2.02	0.41
5:B7:56:GLN:HG2	5:B7:57:ALA:N	2.35	0.41
6:B8:30:GLY:C	6:B8:32:VAL:N	2.76	0.41
9:B9:102:BCL:HMD2	6:B0:36:HIS:CD2	2.49	0.41
1:AC:21:LEU:O	1:AC:22:GLY:C	2.63	0.41
1:AC:63:PRO:O	1:AC:92:ARG:CZ	2.67	0.41
2:AL:21:ASP:HB3	5:A7:19:ARG:HE	1.84	0.41
2:AL:51:VAL:HG12	5:AA:37:MET:HG2	2.01	0.41
2:AL:144:ARG:O	2:AL:146:LEU:N	2.53	0.41
2:AL:175:HIS:CD2	2:AL:178:TYR:CZ	3.08	0.41
2:AL:252:TRP:C	2:AL:254:ALA:N	2.77	0.41
3:AM:12:GLN:O	3:AM:13:VAL:HG13	2.19	0.41
3:AM:61:ILE:CG2	3:AM:62:PHE:N	2.83	0.41
3:AM:83:VAL:HA	3:AM:86:PHE:CB	2.50	0.41
3:AM:261:THR:H	3:AM:264:SER:HG	1.68	0.41
3:AM:299:VAL:CB	3:AM:304:ALA:HB3	2.41	0.41
4:AH:29:TYR:C	4:AH:29:TYR:CD1	2.97	0.41
5:AA:15:LEU:HD22	5:AA:15:LEU:N	2.34	0.41
6:AE:8:GLY:O	6:AE:9:LEU:HG	2.20	0.41
6:AG:10:THR:CG2	6:AG:11:ASP:N	2.82	0.41
5:AI:44:LEU:HD13	6:AJ:43:ARG:CD	2.41	0.41
5:AO:14:ILE:C	6:AP:7:THR:N	2.78	0.41
6:AP:17:PHE:HA	6:AP:20:ILE:HG22	2.01	0.41
5:AQ:24:ILE:HD11	9:AS:103:BCL:H191	2.01	0.41
5:AQ:39:VAL:HG13	5:AQ:43:ASP:HB3	2.01	0.41
6:AT:29:PHE:N	6:AT:29:PHE:HD1	2.17	0.41
5:AW:28:GLN:NE2	14:AX:102:CRT:H27	2.35	0.41
5:AY:13:LEU:HD22	6:AZ:14:ALA:HB2	2.01	0.41
6:AZ:21:PHE:CD1	6:AZ:21:PHE:C	2.97	0.41
5:A1:10:LYS:HD2	6:A4:20:ILE:CG2	2.49	0.41
5:A3:28:GLN:CA	5:A3:28:GLN:HE21	2.33	0.41
5:A5:20:VAL:HB	9:A7:103:BCL:C20	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:A5:102:BCL:OBD	6:A6:32:VAL:HG23	2.20	0.41
14:A0:101:CRT:H32	9:A0:102:BCL:HMA3	2.02	0.41
1:BC:130:MET:HB3	7:BC:502:HEM:C4B	2.55	0.41
1:BC:170:PRO:CG	1:BC:171:GLY:H	2.33	0.41
2:BL:30:PHE:N	2:BL:30:PHE:CD1	2.88	0.41
2:BL:32:VAL:HG12	2:BL:33:GLY:N	2.36	0.41
2:BL:106:PHE:O	2:BL:110:ALA:HB2	2.20	0.41
2:BL:111:LEU:HA	2:BL:114:VAL:CG2	2.50	0.41
2:BL:231:TYR:OH	3:BM:40:LEU:HD21	2.20	0.41
5:BA:39:VAL:HG11	9:BB:101:BCL:HBC3	2.02	0.41
9:BA:101:BCL:C20	5:B9:24:ILE:HD13	2.48	0.41
6:BB:40:TRP:CZ3	6:BB:44:PRO:HA	2.55	0.41
9:BF:102:BCL:H143	14:BG:102:CRT:C13	2.49	0.41
14:BF:103:CRT:C2M	5:BK:36:HIS:HB3	2.49	0.41
6:BG:16:GLU:O	6:BG:20:ILE:HG22	2.19	0.41
5:BI:27:PHE:CE2	5:BK:29:ILE:CD1	3.02	0.41
5:BI:29:ILE:CG2	5:BI:30:VAL:N	2.83	0.41
5:BI:46:TRP:NE1	5:BI:47:LEU:HD12	2.35	0.41
5:BQ:25:VAL:HG13	5:BQ:26:ALA:N	2.35	0.41
5:BQ:42:THR:CG2	5:BQ:43:ASP:H	2.21	0.41
6:BR:45:TRP:O	6:BR:46:LEU:CB	2.67	0.41
9:BU:102:BCL:HBC1	9:BV:101:BCL:HBC3	2.02	0.41
5:BW:29:ILE:CA	9:BW:102:BCL:H11	2.42	0.41
5:B7:20:VAL:HG23	5:B7:21:LEU:N	2.35	0.41
6:B8:38:LEU:HD23	6:B8:38:LEU:C	2.46	0.41
6:B0:21:PHE:CD1	14:B0:101:CRT:H16	2.56	0.41
6:B0:36:HIS:CE1	9:B0:102:BCL:CHB	2.99	0.41
1:AC:125:VAL:O	1:AC:128:ARG:HB2	2.21	0.41
1:AC:274:ARG:NH1	1:AC:274:ARG:CG	2.82	0.41
2:AL:114:VAL:O	2:AL:118:ARG:HG3	2.20	0.41
2:AL:170:GLY:O	2:AL:176:PHE:HD1	2.03	0.41
2:AL:205:SER:O	3:AM:142:MET:SD	2.78	0.41
9:AL:301:BCL:HBC3	9:AL:301:BCL:CHD	2.43	0.41
3:AM:98:PRO:CG	3:AM:171:TRP:HB3	2.50	0.41
3:AM:148:TRP:O	3:AM:149:ALA:C	2.62	0.41
15:AM:408:PEF:O2P	15:AM:408:PEF:H51	2.21	0.41
4:AH:32:ARG:NH2	4:AH:60:ASP:HB2	2.30	0.41
4:AH:45:ARG:HA	4:AH:96:PRO:HB3	2.02	0.41
4:AH:111:PHE:C	4:AH:115:ALA:HB2	2.45	0.41
5:AA:12:TRP:O	6:AB:9:LEU:HD22	2.20	0.41
6:AB:21:PHE:CE1	14:AB:102:CRT:H16	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:AB:101:BCL:HMB1	9:AB:101:BCL:CBB	2.51	0.41
9:AE:101:BCL:HMB3	9:AF:102:BCL:CHB	2.50	0.41
5:AI:52:PRO:HB2	5:AI:55:TYR:CD2	2.56	0.41
5:AI:52:PRO:CG	5:AI:55:TYR:HE2	2.31	0.41
9:AJ:101:BCL:NB	9:AK:102:BCL:CMB	2.83	0.41
5:AS:27:PHE:CZ	5:AU:29:ILE:CG1	3.04	0.41
5:AU:19:ARG:CZ	5:AU:19:ARG:HB3	2.50	0.41
5:AU:33:LEU:HD12	5:AU:33:LEU:N	2.35	0.41
5:AU:50:ASN:HB3	5:AW:59:GLY:HA3	2.02	0.41
6:AX:21:PHE:CD2	14:AX:102:CRT:C15	3.03	0.41
6:AZ:46:LEU:HD12	5:A1:52:PRO:CG	2.49	0.41
6:A2:17:PHE:HB2	14:A2:102:CRT:H41	2.03	0.41
5:A5:45:ASN:ND2	5:A5:48:ASP:CG	2.79	0.41
6:A8:24:SER:O	6:A8:27:ALA:N	2.54	0.41
6:A8:46:LEU:HD22	6:A0:42:TYR:HE2	1.76	0.41
6:A8:46:LEU:HB3	6:A0:42:TYR:HH	1.86	0.41
6:A0:20:ILE:HD13	6:A0:20:ILE:O	2.20	0.41
6:A0:29:PHE:O	6:A0:30:GLY:C	2.63	0.41
1:BC:86:SER:O	1:BC:87:VAL:C	2.63	0.41
1:BC:248:THR:OG1	1:BC:249:PHE:N	2.54	0.41
7:BC:504:HEM:HHA	7:BC:504:HEM:HBD1	2.02	0.41
2:BL:69:ASN:ND2	2:BL:71:TRP:HB2	2.35	0.41
2:BL:82:TYR:CA	2:BL:85:ARG:HE	2.30	0.41
2:BL:106:PHE:CD1	2:BL:106:PHE:N	2.87	0.41
2:BL:138:LEU:HD12	2:BL:138:LEU:N	2.36	0.41
2:BL:233:ILE:HG12	2:BL:237:ALA:HB1	2.02	0.41
3:BM:85:GLN:O	3:BM:86:PHE:C	2.63	0.41
3:BM:132:ARG:O	3:BM:132:ARG:CD	2.67	0.41
3:BM:194:GLY:N	3:BM:293:ASN:HA	2.35	0.41
3:BM:290:VAL:HG12	3:BM:291:VAL:HG23	2.02	0.41
4:BH:125:LEU:CB	4:BH:129:GLY:O	2.67	0.41
5:BD:44:LEU:O	5:BD:45:ASN:C	2.63	0.41
6:BG:36:HIS:CE1	9:BG:101:BCL:C4A	3.03	0.41
6:BJ:28:TRP:C	6:BJ:30:GLY:N	2.78	0.41
5:BK:16:ASP:HB2	5:BK:19:ARG:HD3	2.02	0.41
5:BK:46:TRP:CE2	9:BK:102:BCL:H2C	2.56	0.41
5:BQ:45:ASN:HB2	5:BQ:49:ASP:HB3	2.01	0.41
5:BS:10:LYS:O	5:BS:13:LEU:HB2	2.21	0.41
5:BU:13:LEU:HD12	14:BU:103:CRT:H1M2	2.02	0.41
9:BX:101:BCL:HMC2	5:BY:47:LEU:HD21	2.01	0.41
5:BY:17:PRO:HB3	6:BZ:17:PHE:CE2	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:BY:18:ARG:HG2	5:BY:18:ARG:NH1	2.35	0.41
5:B1:11:ILE:CG1	14:B1:103:CRT:H81	2.50	0.41
5:B1:13:LEU:HD13	14:B1:103:CRT:H23	1.79	0.41
5:B3:3:THR:HB	5:B3:4:MET:H	1.56	0.41
5:B3:4:MET:SD	5:B3:4:MET:N	2.89	0.41
5:B5:29:ILE:HG23	5:B5:30:VAL:H	1.84	0.41
6:B8:32:VAL:O	6:B8:35:ALA:HB3	2.20	0.41
6:B8:45:TRP:C	5:B9:52:PRO:CD	2.93	0.41
1:AC:128:ARG:NE	7:AC:501:HEM:O2D	2.53	0.41
2:AL:139:VAL:HG23	2:AL:258:LEU:HD13	2.02	0.41
2:AL:215:VAL:HG11	3:AM:239:THR:CG2	2.51	0.41
3:AM:175:VAL:HA	3:AM:185:TRP:CG	2.53	0.41
3:AM:269:ALA:O	3:AM:270:TRP:C	2.62	0.41
4:AH:54:LYS:HE2	4:AH:58:PHE:CE1	2.55	0.41
4:AH:80:ARG:HH11	4:AH:80:ARG:HG3	1.85	0.41
4:AH:232:THR:OG1	4:AH:235:GLU:HG2	2.20	0.41
4:AH:240:CYS:O	4:AH:242:TYR:N	2.54	0.41
4:AH:249:TYR:C	4:AH:251:THR:N	2.78	0.41
9:AE:101:BCL:HMB1	9:AE:101:BCL:CBB	2.49	0.41
5:AK:11:ILE:CG1	14:AP:102:CRT:H81	2.51	0.41
9:AK:102:BCL:H2A	9:AK:102:BCL:O1D	2.19	0.41
6:AN:20:ILE:HG21	14:AN:102:CRT:C8	2.51	0.41
9:AT:101:BCL:CHB	9:AU:102:BCL:HMB2	2.50	0.41
9:AU:102:BCL:C1D	9:AV:102:BCL:HMD3	2.51	0.41
6:AV:10:THR:CG2	6:AV:11:ASP:H	2.29	0.41
9:AV:102:BCL:CBB	9:AV:102:BCL:HMB3	2.51	0.41
5:AY:8:LEU:CD2	5:AY:11:ILE:HD11	2.47	0.41
5:AY:11:ILE:HA	14:A2:102:CRT:H82	2.03	0.41
5:A5:35:ILE:O	5:A5:36:HIS:C	2.61	0.41
14:A5:103:CRT:H20	14:A5:103:CRT:H181	1.92	0.41
5:A7:17:PRO:HG2	5:A7:18:ARG:H	1.85	0.41
1:BC:28:PRO:HD3	2:BL:262:PRO:HA	2.01	0.41
1:BC:264:PRO:HG2	1:BC:265:LYS:HD2	2.02	0.41
2:BL:156:PRO:O	2:BL:157:TYR:CD1	2.73	0.41
2:BL:190:PHE:CZ	9:BM:402:BCL:CGA	3.03	0.41
2:BL:192:ASN:CA	2:BL:245:LEU:HD13	2.51	0.41
3:BM:134:TYR:CG	3:BM:144:GLN:NE2	2.88	0.41
3:BM:271:TRP:O	3:BM:272:CYS:C	2.62	0.41
4:BH:5:ILE:CG2	4:BH:6:THR:N	2.52	0.41
4:BH:58:PHE:N	4:BH:59:PRO:CD	2.83	0.41
4:BH:100:LEU:O	4:BH:111:PHE:HE2	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:BH:142:PHE:HZ	4:BH:173:ASP:O	2.03	0.41
4:BH:176:GLU:O	4:BH:178:GLN:HG2	2.20	0.41
5:BA:47:LEU:HA	5:B9:43:ASP:OD2	2.20	0.41
14:BA:102:CRT:H403	5:BD:35:ILE:CD1	2.47	0.41
5:BF:55:TYR:N	5:BF:55:TYR:CD1	2.88	0.41
9:BF:102:BCL:C2D	9:BG:101:BCL:C2D	2.98	0.41
6:BG:28:TRP:O	6:BG:31:LEU:N	2.53	0.41
6:BJ:33:VAL:CG1	6:BJ:34:ILE:N	2.82	0.41
5:BK:33:LEU:O	5:BK:37:MET:HG2	2.20	0.41
9:BO:102:BCL:CBD	9:BP:101:BCL:CAD	2.99	0.41
5:BS:8:LEU:HB3	6:BT:18:HIS:CE1	2.56	0.41
5:BU:43:ASP:HB2	5:BW:47:LEU:CD2	2.48	0.41
9:BZ:101:BCL:C4A	9:B1:102:BCL:HMB2	2.50	0.41
14:B1:103:CRT:H35	5:B3:31:LEU:HD11	2.02	0.41
6:B0:29:PHE:O	6:B0:32:VAL:N	2.54	0.41
1:AC:141:TRP:C	1:AC:143:ALA:H	2.28	0.41
1:AC:276:VAL:HG22	1:AC:280:ASN:HD22	1.85	0.41
2:AL:137:TYR:CD1	2:AL:137:TYR:C	2.99	0.41
2:AL:147:LEU:O	2:AL:148:MET:HG2	2.21	0.41
10:AL:302:BPH:HMB3	10:AL:302:BPH:HBB3	2.02	0.41
3:AM:104:LEU:CD2	3:AM:169:GLY:HA2	2.48	0.41
3:AM:123:THR:HA	3:AM:157:TYR:OH	2.21	0.41
3:AM:161:GLY:HA3	14:AM:406:CRT:H291	2.02	0.41
3:AM:226:VAL:HG22	3:AM:229:PHE:HB2	2.03	0.41
4:AH:184:VAL:O	4:AH:193:VAL:HG22	2.20	0.41
6:AB:42:TYR:CZ	6:A0:46:LEU:HD22	2.56	0.41
6:AB:46:LEU:HD13	6:AE:42:TYR:OH	2.20	0.41
6:AN:31:LEU:CA	6:AN:34:ILE:HG22	2.48	0.41
5:AO:43:ASP:OD1	5:AO:44:LEU:HD23	2.21	0.41
5:AO:43:ASP:C	5:AQ:48:ASP:HB3	2.46	0.41
9:AO:102:BCL:C4D	9:AP:101:BCL:C2D	2.98	0.41
5:AS:36:HIS:HD2	5:AS:46:TRP:CH2	2.38	0.41
14:AS:104:CRT:C36	5:AW:33:LEU:HA	2.34	0.41
14:A1:103:CRT:H10	14:A1:103:CRT:H81	1.57	0.41
5:A3:46:TRP:NE1	5:A3:47:LEU:CD2	2.84	0.41
6:A8:32:VAL:O	6:A8:35:ALA:HB3	2.20	0.41
6:A0:10:THR:H	6:A0:13:GLU:CD	2.29	0.41
1:BC:178:LEU:HD12	5:BS:41:SER:HB2	2.03	0.41
1:BC:236:MET:HE1	7:BC:503:HEM:ND	2.34	0.41
2:BL:220:HIS:O	2:BL:221:GLU:C	2.64	0.41
11:BL:304:UQ8:H35	11:BL:304:UQ8:H32A	1.83	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:BM:164:ARG:HG3	3:BM:167:MET:HE2	2.02	0.41
3:BM:184:ASP:O	3:BM:187:ALA:N	2.53	0.41
3:BM:238:ILE:HG23	3:BM:263:GLU:HB2	2.01	0.41
3:BM:296:LEU:O	3:BM:300:LYS:HG2	2.20	0.41
4:BH:35:LYS:HZ2	4:BH:59:PRO:HG2	1.86	0.41
4:BH:128:GLU:CD	4:BH:128:GLU:N	2.66	0.41
4:BH:151:PRO:HA	4:BH:154:MET:CG	2.50	0.41
5:BD:31:LEU:HD12	5:BD:31:LEU:HA	1.80	0.41
5:BF:35:ILE:HG13	9:BG:101:BCL:O1D	2.21	0.41
5:BF:40:LEU:HD13	5:BF:46:TRP:CZ2	2.56	0.41
5:BO:10:LYS:H	5:BO:10:LYS:HG2	1.67	0.41
5:BO:30:VAL:HG13	5:BO:31:LEU:N	2.35	0.41
5:BO:35:ILE:C	5:BO:37:MET:N	2.75	0.41
6:BR:45:TRP:HD1	6:BR:46:LEU:N	2.18	0.41
6:BZ:21:PHE:CD1	6:BZ:22:MET:N	2.89	0.41
5:B1:9:TYR:CD1	5:B1:9:TYR:C	2.97	0.41
5:B1:54:SER:C	5:B1:56:GLN:N	2.78	0.41
14:B1:103:CRT:H183	9:B3:102:BCL:H92	2.01	0.41
6:B6:23:GLN:HG3	6:B6:24:SER:H	1.85	0.41
9:B6:101:BCL:C1B	9:B7:103:BCL:HMB3	2.51	0.41
5:B9:50:ASN:C	5:B9:51:ILE:O	2.62	0.41
1:AC:36:ARG:NH2	2:AL:90:THR:O	2.54	0.41
1:AC:123:THR:OG1	1:AC:124:LYS:N	2.53	0.41
1:AC:237:MET:SD	2:AL:174:LEU:HB3	2.60	0.41
1:AC:271:TYR:O	1:AC:272:ALA:C	2.64	0.41
2:AL:5:SER:HB3	4:AH:38:GLY:C	2.46	0.41
2:AL:5:SER:CB	4:AH:38:GLY:O	2.69	0.41
2:AL:93:GLY:C	2:AL:97:ILE:HG13	2.45	0.41
2:AL:147:LEU:HB3	2:AL:262:PRO:HB3	2.02	0.41
2:AL:182:HIS:O	2:AL:186:ILE:HG13	2.21	0.41
2:AL:192:ASN:HA	2:AL:245:LEU:CD1	2.51	0.41
2:AL:251:PHE:O	2:AL:254:ALA:HB3	2.21	0.41
3:AM:35:ILE:CG1	15:AM:409:PEF:H321	2.50	0.41
3:AM:84:PHE:CD1	5:AW:37:MET:HE2	2.56	0.41
3:AM:85:GLN:O	3:AM:88:LYS:N	2.54	0.41
3:AM:107:PRO:CD	3:AM:116:LEU:HD22	2.51	0.41
3:AM:219:HIS:O	3:AM:223:ILE:HG12	2.21	0.41
3:AM:284:ILE:CD1	9:AM:402:BCL:OBD	2.68	0.41
15:AM:408:PEF:H52	4:AH:204:LYS:CE	2.46	0.41
4:AH:52:ARG:HB3	4:AH:52:ARG:HH11	1.83	0.41
4:AH:64:PRO:HA	4:AH:79:PRO:HD2	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:AF:3:THR:HG22	5:AF:4:MET:CE	2.51	0.41
6:AJ:45:TRP:O	5:AK:52:PRO:HD3	2.21	0.41
6:AN:25:MET:HE2	14:AN:102:CRT:H242	2.01	0.41
5:AO:51:ILE:O	5:AO:53:VAL:N	2.54	0.41
5:AQ:52:PRO:CG	5:AQ:53:VAL:H	2.34	0.41
9:AQ:102:BCL:H62	6:AR:28:TRP:CH2	2.56	0.41
9:AU:102:BCL:O1D	9:AU:102:BCL:C1A	2.64	0.41
5:AY:35:ILE:CA	5:AY:38:ILE:HG13	2.51	0.41
5:A3:55:TYR:H	5:A3:56:GLN:NE2	2.18	0.41
6:A4:27:ALA:O	6:A4:31:LEU:HG	2.21	0.41
6:A4:45:TRP:O	6:A4:46:LEU:CG	2.68	0.41
5:A5:25:VAL:HG13	9:A5:102:BCL:H41	2.01	0.41
9:A5:102:BCL:H143	14:A7:102:CRT:C13	2.39	0.41
6:A6:20:ILE:HD13	6:A6:20:ILE:C	2.45	0.41
6:A6:26:TYR:C	6:A6:28:TRP:N	2.76	0.41
1:BC:170:PRO:CG	1:BC:171:GLY:N	2.83	0.41
1:BC:239:ILE:O	1:BC:243:LEU:HD13	2.19	0.41
1:BC:253:THR:HG22	2:BL:171:TYR:CD2	2.56	0.41
1:BC:267:THR:O	1:BC:270:TRP:N	2.54	0.41
2:BL:166:VAL:O	9:BL:301:BCL:CBC	2.69	0.41
2:BL:184:LEU:HD22	2:BL:252:TRP:HE1	1.86	0.41
2:BL:273:ASN:O	2:BL:274:TRP:C	2.62	0.41
3:BM:51:ILE:CG1	3:BM:52:TYR:N	2.84	0.41
3:BM:148:TRP:HB3	3:BM:270:TRP:HZ2	1.84	0.41
3:BM:205:SER:OG	3:BM:279:THR:C	2.63	0.41
3:BM:242:GLY:HA2	4:BH:117:PRO:HG3	2.03	0.41
3:BM:268:TRP:CZ2	4:BH:30:LEU:HB3	2.56	0.41
4:BH:69:LEU:HD11	4:BH:76:VAL:HG23	2.01	0.41
4:BH:69:LEU:CD2	4:BH:70:PRO:HD2	2.49	0.41
9:BB:101:BCL:HHC	9:BB:101:BCL:OBB	2.19	0.41
14:BB:102:CRT:O1	5:B9:10:LYS:HB3	2.21	0.41
5:BD:51:ILE:HG23	5:BD:52:PRO:HA	2.02	0.41
6:BJ:20:ILE:CG2	6:BJ:21:PHE:N	2.84	0.41
5:BK:44:LEU:HD21	5:BK:46:TRP:HB3	1.99	0.41
6:BR:46:LEU:HD13	6:BT:42:TYR:CZ	2.56	0.41
6:BT:9:LEU:HB3	6:BT:13:GLU:HG3	2.02	0.41
5:BW:36:HIS:O	5:BW:40:LEU:HB3	2.20	0.41
9:BY:102:BCL:HED1	6:BZ:31:LEU:C	2.45	0.41
5:B1:13:LEU:CD1	14:B1:103:CRT:C2	2.46	0.41
9:B2:101:BCL:HMC1	9:B3:102:BCL:HBB1	2.02	0.41
1:AC:32:GLN:OE1	2:AL:80:LEU:N	2.40	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AC:53:ILE:HG12	1:AC:319:TYR:CD1	2.54	0.41
1:AC:320:GLY:O	1:AC:321:ALA:C	2.63	0.41
2:AL:126:VAL:O	2:AL:129:ALA:HB3	2.21	0.41
2:AL:270:GLU:C	2:AL:272:TRP:N	2.78	0.41
4:AH:32:ARG:O	4:AH:35:LYS:N	2.51	0.41
4:AH:36:ARG:NE	4:AH:65:LYS:HB2	2.36	0.41
5:AF:31:LEU:O	5:AF:35:ILE:HG12	2.20	0.41
6:AG:23:GLN:HA	6:AG:26:TYR:CD2	2.56	0.41
9:AN:101:BCL:CHB	9:AO:102:BCL:HMB3	2.51	0.41
14:AN:102:CRT:H10	14:AN:102:CRT:H81	1.87	0.41
5:AO:12:TRP:CZ3	6:AP:17:PHE:CE2	3.09	0.41
5:AQ:15:LEU:HD11	5:AS:21:LEU:HD13	2.02	0.41
5:AQ:55:TYR:O	5:AQ:59:GLY:HA3	2.20	0.41
5:AU:19:ARG:NH2	5:AU:19:ARG:CB	2.84	0.41
6:AV:24:SER:O	6:AV:27:ALA:HB3	2.21	0.41
6:AX:13:GLU:HA	6:AX:16:GLU:HB3	2.03	0.41
5:AY:9:TYR:CG	5:AY:10:LYS:N	2.89	0.41
5:AY:43:ASP:HA	5:A1:48:ASP:CB	2.50	0.41
9:A7:103:BCL:CBB	9:A7:103:BCL:HMB1	2.51	0.41
1:BC:123:THR:O	1:BC:126:VAL:HG22	2.20	0.41
1:BC:138:ASN:O	1:BC:142:LYS:HG2	2.20	0.41
1:BC:316:LYS:O	1:BC:319:TYR:N	2.52	0.41
2:BL:67:THR:OG1	2:BL:68:TYR:N	2.54	0.41
2:BL:144:ARG:O	2:BL:147:LEU:N	2.54	0.41
9:BL:301:BCL:H201	9:BL:303:BCL:H72	2.03	0.41
3:BM:25:LYS:HG2	5:BO:16:ASP:CG	2.46	0.41
3:BM:28:LEU:CB	3:BM:29:PRO:HD2	2.42	0.41
3:BM:265:ILE:C	3:BM:267:ARG:H	2.27	0.41
4:BH:151:PRO:HA	4:BH:154:MET:HG3	2.03	0.41
4:BH:211:VAL:HG12	4:BH:213:ALA:HB3	2.03	0.41
6:BB:27:ALA:HB1	5:B9:4:MET:HG3	2.03	0.41
6:BE:31:LEU:HA	6:BE:34:ILE:HG22	2.02	0.41
5:BF:4:MET:O	5:BF:8:LEU:CG	2.67	0.41
5:BF:31:LEU:HD11	5:BF:35:ILE:HD11	2.02	0.41
5:BF:38:ILE:HD12	5:BI:37:MET:HE1	2.00	0.41
5:BO:28:GLN:HB3	9:BP:101:BCL:HED1	2.02	0.41
5:BS:4:MET:C	5:BS:6:ALA:N	2.78	0.41
5:BU:42:THR:O	5:BW:48:ASP:HB3	2.20	0.41
5:BW:26:ALA:O	5:BW:30:VAL:HG12	2.21	0.41
5:BW:51:ILE:CB	5:BW:52:PRO:CA	2.94	0.41
6:B2:26:TYR:HA	6:B2:29:PHE:CD2	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:B3:37:MET:O	5:B3:38:ILE:C	2.64	0.41
9:B3:102:BCL:H8	9:B3:102:BCL:H121	1.75	0.41
5:B9:44:LEU:CD1	5:B9:44:LEU:N	2.84	0.41
1:AC:114:GLY:O	1:AC:115:ASN:C	2.63	0.41
2:AL:27:VAL:O	2:AL:28:GLY:C	2.63	0.41
2:AL:50:ILE:CB	2:AL:98:ILE:HD11	2.51	0.41
2:AL:93:GLY:HA2	2:AL:96:GLN:NE2	2.36	0.41
2:AL:101:CYS:O	2:AL:104:GLY:N	2.54	0.41
2:AL:119:LYS:C	2:AL:121:GLY:N	2.79	0.41
2:AL:157:TYR:CD1	10:AL:302:BPH:H151	2.55	0.41
2:AL:177:HIS:CG	3:AM:183:LEU:CD2	3.04	0.41
2:AL:188:PHE:CB	2:AL:249:ALA:HB2	2.44	0.41
2:AL:190:PHE:CZ	9:AM:402:BCL:O2A	2.74	0.41
2:AL:204:LEU:HD11	3:AM:267:ARG:HD3	2.02	0.41
2:AL:244:PHE:CZ	11:AL:304:UQ8:H30A	2.55	0.41
2:AL:252:TRP:HA	2:AL:252:TRP:HE3	1.84	0.41
3:AM:161:GLY:N	3:AM:165:PRO:HD2	2.36	0.41
3:AM:175:VAL:HB	14:AM:406:CRT:H242	2.02	0.41
3:AM:249:ALA:HB2	13:AM:405:MQ8:C6	2.46	0.41
3:AM:259:ASN:C	3:AM:259:ASN:ND2	2.76	0.41
3:AM:269:ALA:C	3:AM:271:TRP:N	2.78	0.41
3:AM:274:VAL:C	3:AM:276:THR:N	2.79	0.41
4:AH:13:GLN:C	4:AH:16:ILE:HG22	2.45	0.41
4:AH:14:ILE:O	4:AH:17:TRP:N	2.50	0.41
4:AH:31:ARG:HH21	4:AH:34:ASP:CG	2.28	0.41
4:AH:116:SER:OG	4:AH:117:PRO:HD2	2.21	0.41
4:AH:135:PRO:C	4:AH:137:ARG:N	2.77	0.41
4:AH:135:PRO:HD3	4:AH:171:TRP:CH2	2.55	0.41
4:AH:168:SER:OG	4:AH:169:ASP:N	2.54	0.41
4:AH:184:VAL:CG2	4:AH:195:LEU:HB2	2.50	0.41
4:AH:240:CYS:O	4:AH:241:ALA:C	2.64	0.41
5:AA:49:ASP:O	5:AD:60:LYS:N	2.54	0.41
5:AD:40:LEU:HD13	5:AD:46:TRP:CH2	2.56	0.41
5:AF:43:ASP:OD1	5:AF:43:ASP:C	2.64	0.41
6:AG:28:TRP:O	6:AG:31:LEU:N	2.53	0.41
5:AI:50:ASN:CG	5:AI:51:ILE:N	2.79	0.41
5:AI:50:ASN:HB3	5:AK:59:GLY:C	2.46	0.41
6:AJ:46:LEU:HB3	6:AN:42:TYR:OH	2.21	0.41
5:AK:45:ASN:O	5:AK:47:LEU:N	2.54	0.41
9:AK:102:BCL:HMD2	9:AN:101:BCL:C4C	2.47	0.41
5:AO:7:ASN:CG	6:AR:23:GLN:OE1	2.63	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:AO:35:ILE:HD13	5:AO:35:ILE:HA	1.91	0.41
5:AO:45:ASN:C	5:AO:47:LEU:N	2.79	0.41
5:AO:53:VAL:C	5:AO:55:TYR:H	2.29	0.41
6:AP:24:SER:O	6:AP:27:ALA:N	2.51	0.41
5:AQ:27:PHE:HE2	5:AS:29:ILE:HD12	1.81	0.41
5:AQ:43:ASP:H	5:AS:47:LEU:HB3	1.83	0.41
5:AQ:51:ILE:HG23	5:AQ:52:PRO:HA	1.99	0.41
6:AR:33:VAL:HG23	9:AR:101:BCL:C14	2.51	0.41
5:AS:5:ASN:O	6:AT:18:HIS:CD2	2.74	0.41
5:AS:20:VAL:HG23	5:AS:21:LEU:N	2.35	0.41
5:AS:27:PHE:CE2	5:AU:29:ILE:HG13	2.56	0.41
14:AS:104:CRT:H2M3	5:AW:36:HIS:CB	2.50	0.41
6:AT:33:VAL:CG1	6:AT:34:ILE:N	2.83	0.41
5:AU:27:PHE:CD1	5:AU:27:PHE:C	2.98	0.41
5:AU:42:THR:O	5:AW:48:ASP:HB3	2.21	0.41
5:AU:46:TRP:HA	5:AU:49:ASP:CG	2.45	0.41
6:AV:45:TRP:O	6:AV:46:LEU:HG	2.20	0.41
9:AV:102:BCL:HMB3	9:AV:102:BCL:HBB2	2.03	0.41
5:AW:21:LEU:HD12	5:AW:21:LEU:C	2.46	0.41
5:AW:30:VAL:HA	5:AW:33:LEU:CG	2.51	0.41
14:AW:102:CRT:H392	5:AY:35:ILE:HD11	2.02	0.41
6:AZ:34:ILE:HD13	6:AZ:34:ILE:O	2.21	0.41
5:A1:33:LEU:C	5:A1:33:LEU:HD12	2.46	0.41
6:A4:28:TRP:HA	6:A4:31:LEU:HB2	2.03	0.41
5:A5:51:ILE:HA	5:A5:53:VAL:N	2.36	0.41
14:A7:102:CRT:H36	14:A7:102:CRT:H341	1.88	0.41
6:A8:24:SER:O	6:A8:25:MET:C	2.64	0.41
6:A0:29:PHE:O	6:A0:32:VAL:N	2.54	0.41
1:BC:24:GLU:CG	2:BL:266:ARG:HH22	2.33	0.41
1:BC:53:ILE:HG13	1:BC:319:TYR:CZ	2.56	0.41
1:BC:90:PHE:CD1	1:BC:91:THR:N	2.87	0.41
1:BC:190:VAL:O	1:BC:192:TYR:N	2.54	0.41
1:BC:211:ARG:HD3	3:BM:317:TYR:CZ	2.56	0.41
1:BC:236:MET:HA	1:BC:239:ILE:HD12	2.03	0.41
1:BC:237:MET:SD	2:BL:174:LEU:HB3	2.61	0.41
1:BC:267:THR:HG21	3:BM:314:VAL:CA	2.51	0.41
1:BC:269:ALA:O	1:BC:270:TRP:C	2.64	0.41
1:BC:281:GLN:HA	1:BC:285:TRP:HD1	1.86	0.41
1:BC:313:ALA:C	1:BC:314:VAL:HG23	2.44	0.41
1:BC:327:TYR:HA	1:BC:328:PRO:HD2	1.79	0.41
2:BL:51:VAL:HG12	5:BA:37:MET:HG2	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:BL:51:VAL:CG1	5:BA:37:MET:HG2	2.51	0.41
2:BL:72:ARG:O	2:BL:73:ILE:C	2.64	0.41
2:BL:261:GLY:HA3	2:BL:262:PRO:HD3	1.95	0.41
2:BL:270:GLU:O	2:BL:273:ASN:HB2	2.21	0.41
2:BL:273:ASN:C	2:BL:275:TRP:N	2.78	0.41
3:BM:9:THR:O	3:BM:10:ALA:C	2.64	0.41
3:BM:85:GLN:O	3:BM:88:LYS:N	2.53	0.41
3:BM:98:PRO:CD	3:BM:171:TRP:HB3	2.48	0.41
3:BM:155:PHE:HE2	3:BM:278:ILE:O	2.04	0.41
3:BM:167:MET:HE3	3:BM:289:THR:CG2	2.51	0.41
3:BM:215:LEU:HD11	13:BM:405:MQ8:H193	2.03	0.41
4:BH:17:TRP:O	4:BH:18:ALA:C	2.63	0.41
4:BH:176:GLU:HA	4:BH:177:PRO:HD2	1.86	0.41
5:BA:20:VAL:HA	5:BA:23:SER:HB3	2.03	0.41
5:BA:30:VAL:HG22	5:BA:30:VAL:O	2.21	0.41
5:BA:35:ILE:HG22	5:BA:36:HIS:N	2.35	0.41
14:BA:102:CRT:C40	5:BD:38:ILE:HG21	2.50	0.41
6:BJ:32:VAL:HG11	9:BJ:101:BCL:CBA	2.40	0.41
5:BK:26:ALA:O	5:BK:29:ILE:CG2	2.68	0.41
5:BO:12:TRP:HE1	6:BP:18:HIS:HA	1.82	0.41
5:BO:35:ILE:O	5:BO:38:ILE:HG12	2.20	0.41
6:BP:30:GLY:C	6:BP:32:VAL:N	2.79	0.41
5:BS:29:ILE:CG2	5:BS:30:VAL:H	2.33	0.41
6:BT:22:MET:O	6:BT:25:MET:N	2.50	0.41
5:BU:40:LEU:O	5:BU:45:ASN:ND2	2.54	0.41
9:BU:102:BCL:CED	6:BV:35:ALA:HB2	2.50	0.41
6:BV:17:PHE:CA	14:BV:102:CRT:H6	2.51	0.41
6:BV:17:PHE:N	14:BV:102:CRT:H41	2.36	0.41
6:BV:20:ILE:HG21	14:BV:102:CRT:C8	2.48	0.41
5:BW:7:ASN:HD22	5:BW:7:ASN:H	1.65	0.41
14:BW:103:CRT:H36	14:BW:103:CRT:H341	1.93	0.41
9:BY:102:BCL:HMD1	6:BZ:36:HIS:CD2	2.56	0.41
6:BZ:26:TYR:O	6:BZ:27:ALA:C	2.64	0.41
5:B1:13:LEU:HD13	14:B1:103:CRT:H1M3	2.02	0.41
6:B4:40:TRP:CE3	6:B4:44:PRO:HA	2.56	0.41
5:B5:14:ILE:CG2	5:B7:18:ARG:HG2	2.43	0.41
5:B5:44:LEU:C	5:B5:46:TRP:N	2.79	0.41
6:B6:20:ILE:HD13	6:B6:20:ILE:C	2.44	0.41
5:B9:9:TYR:CD1	5:B9:9:TYR:C	2.99	0.41
5:B9:46:TRP:CZ3	9:B9:102:BCL:HBC3	2.56	0.41
6:B0:10:THR:CG2	6:B0:11:ASP:H	2.23	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:B0:40:TRP:HB2	9:B0:102:BCL:C19	2.49	0.41
1:AC:153:TYR:O	1:AC:157:ARG:CG	2.69	0.41
1:AC:157:ARG:HH12	1:AC:318:LEU:HD11	1.86	0.41
1:AC:193:ALA:HB3	1:AC:195:LEU:CD1	2.46	0.41
1:AC:195:LEU:HB3	1:AC:196:PRO:HD3	1.99	0.41
7:AC:501:HEM:CBC	7:AC:502:HEM:HMB1	2.51	0.41
2:AL:109:TRP:CD1	2:AL:109:TRP:C	2.99	0.41
3:AM:206:ILE:HG22	3:AM:210:TYR:CE2	2.56	0.41
3:AM:253:ARG:HD3	3:AM:257:GLY:O	2.21	0.41
3:AM:257:GLY:O	3:AM:258:PHE:C	2.63	0.41
3:AM:298:ALA:C	3:AM:303:MET:HB2	2.46	0.41
4:AH:106:PRO:O	4:AH:107:MET:C	2.64	0.41
5:AD:43:ASP:HB2	5:AF:47:LEU:HD22	2.03	0.41
6:AG:45:TRP:CG	6:AG:46:LEU:N	2.79	0.41
5:AK:12:TRP:HD1	6:AN:17:PHE:HB3	1.86	0.41
9:AR:101:BCL:HMA1	9:AS:103:BCL:HHB	2.03	0.41
5:AS:36:HIS:CD2	5:AS:46:TRP:HH2	2.39	0.41
5:AU:44:LEU:HD13	6:AV:43:ARG:NE	2.36	0.41
6:AV:44:PRO:O	5:AW:55:TYR:OH	2.39	0.41
5:AW:4:MET:CE	6:AZ:23:GLN:HB3	2.51	0.41
5:AW:9:TYR:C	5:AW:9:TYR:CD1	2.99	0.41
5:AW:9:TYR:CA	6:AX:18:HIS:CG	2.99	0.41
5:AW:24:ILE:HG21	14:AX:102:CRT:H20	2.03	0.41
9:AW:101:BCL:O1A	6:AX:28:TRP:CZ2	2.74	0.41
5:AY:8:LEU:O	6:AZ:18:HIS:CE1	2.74	0.41
5:AY:46:TRP:HA	5:AY:49:ASP:OD1	2.21	0.41
9:AY:102:BCL:C3D	9:AZ:101:BCL:C3D	2.98	0.41
5:A5:55:TYR:O	5:A5:59:GLY:HA3	2.21	0.41
6:A0:22:MET:O	6:A0:26:TYR:HD2	2.03	0.41
6:A0:33:VAL:O	6:A0:37:LEU:N	2.47	0.41
1:BC:129:ARG:O	1:BC:132:GLU:N	2.53	0.41
1:BC:133:LEU:HD11	1:BC:279:ILE:HG12	2.02	0.41
1:BC:287:LEU:O	1:BC:290:VAL:HG22	2.21	0.41
2:BL:97:ILE:HG21	5:B9:37:MET:HE3	2.03	0.41
2:BL:139:VAL:CG2	2:BL:258:LEU:HD22	2.51	0.41
2:BL:228:ILE:HG23	3:BM:132:ARG:CD	2.51	0.41
3:BM:76:LEU:O	3:BM:77:ALA:C	2.63	0.41
3:BM:156:PHE:HD1	3:BM:281:GLY:HA2	1.86	0.41
3:BM:293:ASN:ND2	3:BM:296:LEU:HG	2.36	0.41
4:BH:106:PRO:C	4:BH:108:LEU:N	2.78	0.41
6:BB:20:ILE:HD13	6:BB:20:ILE:C	2.46	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
14:BB:102:CRT:H341	14:BB:102:CRT:H36	1.83	0.41
5:BD:36:HIS:CE1	9:BE:101:BCL:HMD2	2.54	0.41
5:BF:43:ASP:HB2	5:BI:47:LEU:C	2.46	0.41
6:BT:45:TRP:HD1	6:BT:46:LEU:N	2.18	0.41
9:BT:101:BCL:HBA1	9:BT:101:BCL:H3A	1.80	0.41
14:BU:103:CRT:H81	14:BU:103:CRT:H10	1.66	0.41
6:BV:24:SER:O	6:BV:27:ALA:HB3	2.20	0.41
6:BX:17:PHE:O	6:BX:18:HIS:C	2.62	0.41
9:BX:101:BCL:H172	6:BZ:38:LEU:HD21	2.03	0.41
5:B3:28:GLN:HG3	9:B3:102:BCL:C2	2.50	0.41
5:B5:10:LYS:C	14:B5:103:CRT:H5	2.46	0.41
5:B7:33:LEU:HG	14:B7:102:CRT:C35	2.51	0.41
6:B8:22:MET:CG	6:B8:26:TYR:HE2	2.33	0.41
9:B9:102:BCL:HMD1	6:B0:36:HIS:HD2	1.85	0.41
14:B0:101:CRT:H291	14:B0:101:CRT:H31	1.83	0.41
1:AC:26:PRO:HB3	2:AL:262:PRO:O	2.21	0.40
1:AC:33:ILE:N	1:AC:33:ILE:CD1	2.84	0.40
1:AC:188:LYS:O	1:AC:189:THR:C	2.64	0.40
1:AC:263:THR:HG22	3:AM:311:VAL:CB	2.49	0.40
2:AL:29:PRO:O	3:AM:253:ARG:O	2.39	0.40
2:AL:78:PRO:HG2	2:AL:152:GLY:CA	2.52	0.40
2:AL:147:LEU:O	2:AL:262:PRO:HG3	2.21	0.40
2:AL:150:ALA:O	2:AL:153:HIS:CB	2.69	0.40
9:AL:303:BCL:H11	3:AM:210:TYR:HB3	2.03	0.40
3:AM:124:LEU:O	3:AM:125:SER:C	2.63	0.40
3:AM:251:PHE:HD1	3:AM:252:TRP:N	2.19	0.40
4:AH:102:PRO:O	4:AH:103:ASN:C	2.64	0.40
5:AA:8:LEU:HD21	6:AE:23:GLN:OE1	2.22	0.40
6:AB:44:PRO:O	5:AD:52:PRO:HG3	2.20	0.40
9:AK:102:BCL:H202	9:AK:102:BCL:H161	1.90	0.40
6:AN:43:ARG:HA	6:AN:44:PRO:HD2	1.94	0.40
5:AU:16:ASP:HB2	5:AU:19:ARG:HD3	2.02	0.40
5:AU:44:LEU:HD12	5:AU:44:LEU:O	2.21	0.40
9:AU:102:BCL:CBB	9:AU:102:BCL:CMB	2.99	0.40
9:A1:102:BCL:H2A	9:A1:102:BCL:CGD	2.51	0.40
6:A4:10:THR:CG2	6:A4:11:ASP:H	2.18	0.40
6:A4:42:TYR:CD1	6:A4:43:ARG:HG3	2.56	0.40
6:A6:23:GLN:HG3	6:A6:24:SER:H	1.85	0.40
6:A6:28:TRP:O	6:A6:30:GLY:N	2.53	0.40
5:A7:7:ASN:CB	5:A7:10:LYS:HZ3	2.27	0.40
5:A7:11:ILE:HG12	5:A7:15:LEU:CD1	2.50	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:A8:22:MET:CG	6:A8:26:TYR:HE2	2.33	0.40
1:BC:31:GLU:O	1:BC:33:ILE:HD12	2.21	0.40
1:BC:70:PRO:HB2	1:BC:71:LYS:HD2	2.03	0.40
1:BC:135:ARG:O	1:BC:137:ALA:N	2.54	0.40
1:BC:237:MET:O	1:BC:240:SER:N	2.54	0.40
2:BL:30:PHE:CZ	3:BM:257:GLY:HA3	2.55	0.40
2:BL:177:HIS:CB	3:BM:183:LEU:CD2	2.92	0.40
2:BL:185:ALA:CB	2:BL:252:TRP:HB3	2.47	0.40
2:BL:192:ASN:HA	2:BL:245:LEU:HD13	2.03	0.40
2:BL:240:ARG:HG2	3:BM:224:LEU:HD21	2.04	0.40
2:BL:278:LEU:O	2:BL:280:LEU:N	2.51	0.40
3:BM:232:ASP:C	3:BM:234:GLU:H	2.29	0.40
4:BH:28:ILE:O	4:BH:31:ARG:N	2.54	0.40
5:BA:43:ASP:O	5:BD:48:ASP:HB3	2.21	0.40
9:BE:101:BCL:C1C	9:BF:102:BCL:HBB3	2.50	0.40
6:BJ:27:ALA:O	6:BJ:31:LEU:HG	2.21	0.40
5:BK:56:GLN:H	5:BK:56:GLN:CD	2.29	0.40
5:BQ:43:ASP:OD1	5:BQ:44:LEU:CD2	2.69	0.40
5:BS:43:ASP:HB3	5:BU:56:GLN:CB	2.50	0.40
5:BW:21:LEU:HD11	9:BW:102:BCL:H142	2.04	0.40
5:BW:39:VAL:HG22	5:BY:47:LEU:HD11	2.03	0.40
6:BX:17:PHE:C	6:BX:20:ILE:HG22	2.46	0.40
5:B1:46:TRP:CD1	5:B1:47:LEU:N	2.89	0.40
9:B6:101:BCL:CHB	9:B7:103:BCL:HMB3	2.51	0.40
6:B0:17:PHE:CE1	6:B0:21:PHE:HB3	2.57	0.40
1:AC:48:GLN:O	1:AC:50:ALA:N	2.54	0.40
1:AC:148:THR:CG2	1:AC:322:GLN:HA	2.38	0.40
2:AL:4:LEU:H	2:AL:7:GLU:HB3	1.87	0.40
2:AL:233:ILE:O	3:AM:44:GLY:HA3	2.21	0.40
2:AL:246:ALA:CB	3:AM:217:ALA:HB2	2.51	0.40
9:AL:301:BCL:H201	9:AL:303:BCL:H72	2.03	0.40
3:AM:262:MET:O	3:AM:262:MET:HG3	2.21	0.40
4:AH:189:ASN:C	4:AH:191:LYS:N	2.78	0.40
5:AA:49:ASP:O	5:AD:55:TYR:O	2.39	0.40
9:AA:101:BCL:H92	6:AB:28:TRP:CE3	2.56	0.40
9:AB:101:BCL:C1B	9:AD:102:BCL:CMB	3.00	0.40
5:AD:35:ILE:O	5:AD:38:ILE:HG22	2.21	0.40
5:AF:4:MET:C	5:AF:6:ALA:N	2.79	0.40
6:AG:32:VAL:HG11	9:AG:101:BCL:CBA	2.18	0.40
5:AI:18:ARG:H	5:AI:18:ARG:HG2	1.63	0.40
5:AK:16:ASP:HA	5:AK:17:PRO:HD2	1.94	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:AK:40:LEU:O	5:AK:45:ASN:HA	2.20	0.40
5:AK:47:LEU:H	5:AK:47:LEU:CD2	2.31	0.40
9:AK:102:BCL:C3D	9:AN:101:BCL:CMD	3.00	0.40
5:AQ:30:VAL:HG13	5:AQ:31:LEU:H	1.84	0.40
5:AQ:38:ILE:O	5:AQ:42:THR:HG22	2.21	0.40
6:AT:29:PHE:CA	6:AT:32:VAL:HG12	2.52	0.40
5:A1:33:LEU:O	5:A1:37:MET:HB2	2.21	0.40
5:A3:12:TRP:CE3	5:A3:12:TRP:CA	3.04	0.40
5:A5:21:LEU:HD11	9:A5:102:BCL:H142	1.97	0.40
5:A7:48:ASP:O	5:A7:48:ASP:CG	2.64	0.40
5:A7:50:ASN:ND2	5:A7:51:ILE:HG12	2.37	0.40
1:BC:24:GLU:HG3	2:BL:266:ARG:HH22	1.86	0.40
1:BC:98:THR:HG22	1:BC:99:THR:N	2.37	0.40
1:BC:252:ASN:HB3	1:BC:255:ALA:HB3	2.03	0.40
1:BC:304:ARG:HB3	1:BC:305:VAL:H	1.59	0.40
2:BL:78:PRO:HG2	2:BL:152:GLY:HA3	2.04	0.40
2:BL:173:PHE:C	2:BL:174:LEU:HD12	2.46	0.40
2:BL:195:ALA:O	2:BL:196:LEU:C	2.64	0.40
2:BL:235:ALA:HA	11:BL:304:UQ8:C3M	2.50	0.40
3:BM:62:PHE:O	3:BM:66:VAL:HG23	2.21	0.40
3:BM:186:THR:HG23	3:BM:187:ALA:N	2.32	0.40
3:BM:218:MET:HG3	3:BM:218:MET:H	1.64	0.40
14:BM:406:CRT:H20	14:BM:406:CRT:H181	1.87	0.40
4:BH:36:ARG:NE	4:BH:65:LYS:HD2	2.34	0.40
5:BA:10:LYS:C	5:BA:13:LEU:HD13	2.46	0.40
5:BA:11:ILE:HD13	14:BA:102:CRT:C10	2.51	0.40
9:BA:101:BCL:CAD	9:BB:101:BCL:CAD	3.00	0.40
14:BA:102:CRT:H11	6:BE:17:PHE:CE1	2.52	0.40
9:BB:101:BCL:HBA1	9:BB:101:BCL:H3A	1.64	0.40
5:BI:14:ILE:HG23	5:BK:18:ARG:HB3	2.02	0.40
5:BO:8:LEU:HD21	6:BP:18:HIS:HE2	1.85	0.40
9:BO:102:BCL:H202	9:BO:102:BCL:H161	1.95	0.40
6:BP:17:PHE:O	6:BP:18:HIS:C	2.65	0.40
6:BP:21:PHE:HE1	6:BP:25:MET:HB2	1.86	0.40
6:BP:22:MET:HG3	6:BP:26:TYR:HE2	1.86	0.40
5:BU:35:ILE:O	5:BU:36:HIS:C	2.63	0.40
14:BU:103:CRT:H2M3	5:BY:36:HIS:HB3	2.03	0.40
9:BW:102:BCL:CAC	9:BX:101:BCL:HBC3	2.51	0.40
5:B5:30:VAL:O	5:B5:34:LEU:N	2.53	0.40
6:B6:37:LEU:HD23	6:B6:37:LEU:C	2.46	0.40
6:B8:24:SER:O	6:B8:27:ALA:N	2.54	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AC:46:LYS:O	1:AC:48:GLN:N	2.55	0.40
1:AC:107:CYS:O	1:AC:109:TYR:N	2.55	0.40
1:AC:311:HIS:CE1	1:AC:317:PRO:HD3	2.56	0.40
2:AL:37:VAL:HG23	2:AL:38:VAL:N	2.32	0.40
2:AL:253:SER:C	9:AL:301:BCL:CED	2.93	0.40
3:AM:145:HIS:O	3:AM:270:TRP:NE1	2.42	0.40
3:AM:161:GLY:C	3:AM:165:PRO:HD2	2.46	0.40
3:AM:242:GLY:HA2	4:AH:117:PRO:CG	2.51	0.40
4:AH:184:VAL:HG21	4:AH:195:LEU:HB2	2.04	0.40
4:AH:206:ALA:C	4:AH:208:LYS:N	2.80	0.40
4:AH:259:LEU:HD21	5:A5:19:ARG:O	2.19	0.40
14:AB:102:CRT:H36	14:AB:102:CRT:H341	1.91	0.40
5:AD:51:ILE:CG2	5:AD:52:PRO:HA	2.52	0.40
9:AG:101:BCL:HMB3	9:AG:101:BCL:CBB	2.51	0.40
5:AI:34:LEU:O	5:AI:37:MET:N	2.51	0.40
6:AJ:22:MET:O	6:AJ:25:MET:HB3	2.22	0.40
6:AJ:33:VAL:CG1	6:AJ:34:ILE:N	2.83	0.40
6:AN:43:ARG:HD2	5:AO:55:TYR:HE2	1.85	0.40
5:AO:22:VAL:HA	5:AO:25:VAL:CG2	2.51	0.40
5:AO:29:ILE:HG23	5:AO:30:VAL:N	2.36	0.40
5:AS:42:THR:CG2	5:AS:43:ASP:N	2.84	0.40
6:AT:30:GLY:HA2	6:AT:33:VAL:HG12	2.04	0.40
5:AU:12:TRP:HD1	6:AV:18:HIS:HB2	1.87	0.40
5:AU:14:ILE:HB	14:AX:102:CRT:H83	2.03	0.40
6:AX:21:PHE:CE2	14:AX:102:CRT:C16	3.04	0.40
5:AY:27:PHE:CE2	5:A1:29:ILE:CD1	3.04	0.40
6:A2:46:LEU:OXT	6:A4:43:ARG:NH2	2.49	0.40
9:A3:103:BCL:HMB1	9:A3:103:BCL:CBB	2.50	0.40
6:A4:43:ARG:NH1	5:A5:55:TYR:HB2	2.37	0.40
5:A7:45:ASN:O	5:A7:46:TRP:C	2.64	0.40
5:A9:2:PHE:HB2	5:A9:3:THR:H	1.58	0.40
6:A0:17:PHE:HA	6:A0:20:ILE:HG22	2.03	0.40
14:A0:101:CRT:H81	14:A0:101:CRT:H10	1.64	0.40
9:A0:102:BCL:CBB	9:A0:102:BCL:HMB3	2.51	0.40
1:BC:80:GLN:CG	1:BC:128:ARG:HH22	2.31	0.40
1:BC:219:ALA:HB2	3:BM:290:VAL:O	2.22	0.40
2:BL:82:TYR:HB3	2:BL:85:ARG:CG	2.51	0.40
2:BL:160:LEU:O	2:BL:163:LEU:HD13	2.21	0.40
2:BL:229:VAL:HG23	2:BL:231:TYR:H	1.85	0.40
2:BL:268:TRP:HB2	2:BL:269:PRO:HD3	2.03	0.40
2:BL:276:LEU:O	3:BM:88:LYS:HE3	2.20	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:BM:40:LEU:HD23	3:BM:40:LEU:O	2.21	0.40
3:BM:84:PHE:CE1	5:BU:38:ILE:HD12	2.56	0.40
3:BM:164:ARG:C	3:BM:167:MET:H	2.29	0.40
3:BM:205:SER:OG	3:BM:280:ALA:N	2.55	0.40
3:BM:208:PHE:CZ	3:BM:275:LEU:HB3	2.56	0.40
3:BM:265:ILE:CG2	3:BM:266:HIS:H	2.30	0.40
3:BM:285:LEU:HD12	3:BM:285:LEU:HA	1.75	0.40
3:BM:297:TRP:HZ2	4:BH:13:GLN:HB2	1.84	0.40
6:BB:9:LEU:HA	6:BB:13:GLU:OE1	2.21	0.40
14:BB:102:CRT:H81	14:BB:102:CRT:H10	1.97	0.40
5:BF:9:TYR:CE1	6:BG:15:LYS:HG3	2.57	0.40
5:BF:33:LEU:O	5:BF:37:MET:CG	2.69	0.40
6:BN:22:MET:O	6:BN:25:MET:HB3	2.21	0.40
5:BO:18:ARG:HB2	5:BO:18:ARG:CZ	2.48	0.40
5:BS:36:HIS:CE1	9:BT:101:BCL:HMD2	2.53	0.40
5:BU:13:LEU:HD22	6:BV:9:LEU:CB	2.48	0.40
5:BU:35:ILE:HA	5:BU:38:ILE:CG2	2.51	0.40
6:BV:20:ILE:HG23	6:BV:21:PHE:N	2.36	0.40
9:BW:102:BCL:CED	6:BX:31:LEU:O	2.70	0.40
5:B1:24:ILE:C	5:B1:26:ALA:N	2.79	0.40
5:B5:16:ASP:H	5:B5:19:ARG:HG3	1.86	0.40
1:AC:135:ARG:HH11	1:AC:135:ARG:HB3	1.85	0.40
1:AC:241:ASP:O	1:AC:242:SER:C	2.64	0.40
2:AL:147:LEU:C	2:AL:148:MET:HG2	2.47	0.40
2:AL:228:ILE:O	3:AM:51:ILE:HD11	2.21	0.40
2:AL:237:ALA:C	2:AL:240:ARG:HG3	2.46	0.40
2:AL:268:TRP:N	2:AL:269:PRO:CD	2.85	0.40
3:AM:59:LEU:HD13	5:AQ:29:ILE:HG21	1.99	0.40
3:AM:135:LYS:C	3:AM:137:ALA:N	2.80	0.40
3:AM:178:GLY:O	3:AM:179:ILE:C	2.64	0.40
3:AM:230:GLY:O	3:AM:231:GLY:C	2.65	0.40
3:AM:251:PHE:C	3:AM:251:PHE:CD1	2.99	0.40
3:AM:263:GLU:O	3:AM:264:SER:C	2.65	0.40
4:AH:105:ASP:HB3	4:AH:108:LEU:HD21	2.04	0.40
6:AB:28:TRP:CE3	6:AB:31:LEU:HD12	2.56	0.40
6:AE:45:TRP:CD1	6:AE:46:LEU:N	2.90	0.40
5:AF:7:ASN:CG	6:AJ:20:ILE:HD13	2.46	0.40
5:AI:45:ASN:C	5:AI:49:ASP:CG	2.89	0.40
5:AK:9:TYR:CE1	6:AN:11:ASP:O	2.75	0.40
14:AN:102:CRT:H2M3	5:AO:36:HIS:CB	2.51	0.40
9:AQ:102:BCL:CAD	6:AR:35:ALA:CB	2.99	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:AR:101:BCL:HMC1	9:AS:103:BCL:OBB	2.21	0.40
5:AS:35:ILE:O	5:AS:39:VAL:HG23	2.21	0.40
5:AU:35:ILE:O	5:AU:38:ILE:HG22	2.20	0.40
6:AX:21:PHE:CD2	14:AX:102:CRT:C14	3.04	0.40
6:AX:40:TRP:HA	6:AX:45:TRP:CZ3	2.57	0.40
5:A1:46:TRP:CZ2	9:A1:102:BCL:CHC	3.03	0.40
6:A2:20:ILE:HG21	14:A2:102:CRT:H83	2.04	0.40
9:A2:101:BCL:O1D	9:A2:101:BCL:H2A	2.20	0.40
9:A3:103:BCL:HMD3	6:A4:36:HIS:NE2	2.37	0.40
5:A5:20:VAL:HB	9:A7:103:BCL:H202	2.02	0.40
6:A6:37:LEU:HD23	6:A6:37:LEU:C	2.46	0.40
1:BC:85:LEU:HD22	1:BC:89:GLU:CG	2.45	0.40
1:BC:151:THR:HG21	1:BC:323:MET:CB	2.51	0.40
1:BC:153:TYR:O	1:BC:157:ARG:CG	2.69	0.40
1:BC:236:MET:CE	7:BC:503:HEM:ND	2.84	0.40
2:BL:86:MET:CE	2:BL:96:GLN:HB2	2.51	0.40
2:BL:94:LEU:C	2:BL:94:LEU:HD23	2.47	0.40
2:BL:139:VAL:HG21	2:BL:258:LEU:HD22	2.03	0.40
2:BL:172:GLN:O	2:BL:172:GLN:HG3	2.21	0.40
2:BL:231:TYR:CE1	2:BL:233:ILE:HA	2.55	0.40
3:BM:104:LEU:HD21	3:BM:169:GLY:HA3	2.03	0.40
3:BM:122:LEU:O	3:BM:126:ILE:HD12	2.21	0.40
3:BM:187:ALA:O	3:BM:191:ILE:HG12	2.21	0.40
9:BM:402:BCL:H151	15:BQ:101:PEF:H453	2.02	0.40
4:BH:106:PRO:O	4:BH:107:MET:C	2.64	0.40
14:BA:102:CRT:H20	14:BA:102:CRT:H181	1.92	0.40
5:BD:38:ILE:O	5:BD:38:ILE:HG12	2.21	0.40
14:BF:103:CRT:H2M3	5:BK:36:HIS:HB3	2.03	0.40
5:BK:31:LEU:O	5:BK:34:LEU:HB3	2.22	0.40
9:BK:102:BCL:O1D	9:BK:102:BCL:C2A	2.68	0.40
5:BU:9:TYR:HB2	6:BV:15:LYS:CD	2.52	0.40
5:BU:49:ASP:OD2	5:BU:50:ASN:OD1	2.38	0.40
14:BU:103:CRT:H31	9:BY:102:BCL:H3A	2.03	0.40
5:BW:7:ASN:C	5:BW:9:TYR:N	2.78	0.40
5:B1:42:THR:O	5:B1:43:ASP:C	2.65	0.40
6:B4:28:TRP:HA	6:B4:31:LEU:HB2	2.02	0.40
9:B0:102:BCL:HBA1	9:B0:102:BCL:H3A	1.85	0.40
1:AC:205:ASP:CB	1:AC:304:ARG:HE	2.29	0.40
1:AC:248:THR:OG1	1:AC:249:PHE:N	2.54	0.40
2:AL:49:LEU:CD2	5:A9:37:MET:HG2	2.51	0.40
2:AL:53:GLY:O	2:AL:54:ALA:C	2.64	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:AL:101:CYS:C	2:AL:103:ALA:N	2.80	0.40
2:AL:247:LEU:HD23	2:AL:247:LEU:HA	1.86	0.40
9:AL:301:BCL:H161	9:AL:303:BCL:H161	2.03	0.40
11:AL:304:UQ8:H12A	11:AL:304:UQ8:H15	1.90	0.40
3:AM:135:LYS:C	3:AM:137:ALA:H	2.30	0.40
3:AM:177:PHE:CZ	14:AM:406:CRT:H25	2.57	0.40
3:AM:265:ILE:HA	3:AM:265:ILE:HD12	1.66	0.40
4:AH:136:MET:HE3	4:AH:172:VAL:CG1	2.46	0.40
4:AH:159:LEU:HD23	4:AH:213:ALA:C	2.47	0.40
5:AA:53:VAL:O	5:AA:54:SER:C	2.63	0.40
5:AF:36:HIS:CE1	9:AF:102:BCL:NA	2.89	0.40
5:AI:9:TYR:CA	6:AJ:18:HIS:CE1	2.79	0.40
5:AI:50:ASN:OD1	5:AI:51:ILE:N	2.50	0.40
5:AK:29:ILE:HB	9:AK:102:BCL:C4	2.47	0.40
5:AK:31:LEU:HA	5:AK:34:LEU:HB3	2.03	0.40
5:AQ:35:ILE:C	5:AQ:37:MET:N	2.80	0.40
5:AU:18:ARG:CD	5:AU:18:ARG:N	2.82	0.40
5:AU:43:ASP:HA	5:AW:47:LEU:O	2.21	0.40
5:AW:8:LEU:O	5:AW:11:ILE:CG1	2.70	0.40
6:AX:29:PHE:CE1	9:AX:101:BCL:H11	2.56	0.40
5:AY:51:ILE:CA	5:AY:52:PRO:C	2.94	0.40
6:AZ:26:TYR:O	6:AZ:27:ALA:C	2.64	0.40
5:A1:38:ILE:HG23	5:A1:39:VAL:H	1.87	0.40
5:A5:44:LEU:O	5:A5:46:TRP:N	2.47	0.40
9:A6:101:BCL:NB	9:A7:103:BCL:CMB	2.85	0.40
14:A7:102:CRT:H291	9:A7:103:BCL:O2A	2.21	0.40
6:A8:22:MET:O	6:A8:26:TYR:CD2	2.71	0.40
6:A8:30:GLY:O	6:A8:33:VAL:HG12	2.21	0.40
6:A8:30:GLY:C	6:A8:32:VAL:N	2.76	0.40
1:BC:293:ALA:HA	1:BC:296:LYS:HD2	2.03	0.40
2:BL:3:MET:HE1	2:BL:8:LYS:HG3	2.03	0.40
2:BL:44:LEU:O	2:BL:46:GLY:N	2.48	0.40
2:BL:126:VAL:CB	2:BL:127:PRO:HD3	2.44	0.40
2:BL:171:TYR:O	2:BL:174:LEU:N	2.44	0.40
2:BL:240:ARG:NH1	3:BM:7:ILE:O	2.54	0.40
3:BM:247:ARG:HG3	3:BM:247:ARG:H	1.67	0.40
3:BM:284:ILE:HG12	9:BM:402:BCL:CED	2.52	0.40
4:BH:153:GLY:H	4:BH:167:VAL:HG23	1.85	0.40
4:BH:173:ASP:OD2	4:BH:176:GLU:HG2	2.21	0.40
9:BA:101:BCL:CBC	9:BB:101:BCL:HHD	2.50	0.40
9:BB:101:BCL:HBB3	9:BD:102:BCL:CHC	2.52	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:BF:46:TRP:CD1	5:BF:46:TRP:C	2.99	0.40
5:BK:38:ILE:HG23	5:BK:39:VAL:HG23	2.03	0.40
5:BS:17:PRO:HA	5:BS:20:VAL:CG2	2.52	0.40
5:BU:8:LEU:C	5:BU:10:LYS:N	2.77	0.40
5:BU:35:ILE:O	5:BU:38:ILE:CG2	2.67	0.40
5:BU:44:LEU:HD13	6:BV:43:ARG:CD	2.51	0.40
6:BX:28:TRP:HE3	6:BX:31:LEU:HD12	1.86	0.40
5:B1:46:TRP:CH2	9:B1:102:BCL:HBC3	2.56	0.40
5:B3:5:ASN:HA	5:B3:8:LEU:HG	2.03	0.40
5:B3:53:VAL:O	5:B3:54:SER:CB	2.69	0.40
5:B7:13:LEU:O	6:B8:7:THR:CA	2.70	0.40
5:B7:42:THR:CB	5:B9:48:ASP:OD1	2.70	0.40
5:B7:46:TRP:CH2	9:B7:103:BCL:H2C	2.56	0.40
6:B8:10:THR:HG22	6:B8:11:ASP:H	1.85	0.40
6:B8:24:SER:O	6:B8:25:MET:C	2.64	0.40
6:B8:30:GLY:O	6:B8:33:VAL:HG12	2.22	0.40
5:B9:12:TRP:HZ2	6:B0:18:HIS:HD1	1.62	0.40

There are no symmetry-related clashes.

5.3 Torsion angles (i)

5.3.1 Protein backbone (i)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	AC	315/404 (78%)	195 (62%)	75 (24%)	45 (14%)	0	1
1	BC	315/404 (78%)	192 (61%)	87 (28%)	36 (11%)	0	1
2	AL	278/281 (99%)	137 (49%)	97 (35%)	44 (16%)	0	0
2	BL	278/281 (99%)	149 (54%)	97 (35%)	32 (12%)	0	1
3	AM	317/325 (98%)	162 (51%)	94 (30%)	61 (19%)	0	0
3	BM	317/325 (98%)	188 (59%)	85 (27%)	44 (14%)	0	1
4	AH	256/259 (99%)	162 (63%)	73 (28%)	21 (8%)	0	3

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
4	BH	256/259 (99%)	166 (65%)	66 (26%)	24 (9%)	0 2
5	A1	56/61 (92%)	43 (77%)	9 (16%)	4 (7%)	1 4
5	A3	55/61 (90%)	45 (82%)	7 (13%)	3 (6%)	1 8
5	A5	54/61 (88%)	42 (78%)	10 (18%)	2 (4%)	2 15
5	A7	49/61 (80%)	35 (71%)	11 (22%)	3 (6%)	1 7
5	A9	58/61 (95%)	47 (81%)	10 (17%)	1 (2%)	7 32
5	AA	46/61 (75%)	33 (72%)	11 (24%)	2 (4%)	2 12
5	AD	55/61 (90%)	40 (73%)	12 (22%)	3 (6%)	1 8
5	AF	57/61 (93%)	41 (72%)	13 (23%)	3 (5%)	1 9
5	AI	57/61 (93%)	47 (82%)	6 (10%)	4 (7%)	1 4
5	AK	56/61 (92%)	44 (79%)	10 (18%)	2 (4%)	2 16
5	AO	57/61 (93%)	46 (81%)	6 (10%)	5 (9%)	0 3
5	AQ	55/61 (90%)	36 (66%)	17 (31%)	2 (4%)	2 16
5	AS	57/61 (93%)	46 (81%)	7 (12%)	4 (7%)	1 4
5	AU	58/61 (95%)	43 (74%)	11 (19%)	4 (7%)	1 5
5	AW	58/61 (95%)	45 (78%)	8 (14%)	5 (9%)	0 3
5	AY	58/61 (95%)	43 (74%)	11 (19%)	4 (7%)	1 5
5	B1	52/61 (85%)	34 (65%)	11 (21%)	7 (14%)	0 1
5	B3	58/61 (95%)	38 (66%)	16 (28%)	4 (7%)	1 5
5	B5	49/61 (80%)	35 (71%)	10 (20%)	4 (8%)	0 3
5	B7	52/61 (85%)	40 (77%)	10 (19%)	2 (4%)	2 15
5	B9	49/61 (80%)	33 (67%)	14 (29%)	2 (4%)	2 13
5	BA	53/61 (87%)	29 (55%)	20 (38%)	4 (8%)	1 4
5	BD	43/61 (70%)	32 (74%)	8 (19%)	3 (7%)	1 4
5	BF	54/61 (88%)	42 (78%)	11 (20%)	1 (2%)	6 30
5	BI	48/61 (79%)	32 (67%)	13 (27%)	3 (6%)	1 6
5	BK	58/61 (95%)	43 (74%)	12 (21%)	3 (5%)	1 9
5	BO	57/61 (93%)	46 (81%)	11 (19%)	0	100 100
5	BQ	57/61 (93%)	41 (72%)	14 (25%)	2 (4%)	3 16
5	BS	57/61 (93%)	45 (79%)	9 (16%)	3 (5%)	1 9
5	BU	56/61 (92%)	43 (77%)	12 (21%)	1 (2%)	6 31

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
5	BW	56/61 (92%)	38 (68%)	14 (25%)	4 (7%)	1 4
5	BY	52/61 (85%)	30 (58%)	16 (31%)	6 (12%)	0 1
6	A0	38/47 (81%)	30 (79%)	7 (18%)	1 (3%)	4 23
6	A2	38/47 (81%)	29 (76%)	8 (21%)	1 (3%)	4 23
6	A4	38/47 (81%)	31 (82%)	6 (16%)	1 (3%)	4 23
6	A6	38/47 (81%)	35 (92%)	3 (8%)	0	100 100
6	A8	38/47 (81%)	27 (71%)	10 (26%)	1 (3%)	4 23
6	AB	38/47 (81%)	33 (87%)	5 (13%)	0	100 100
6	AE	38/47 (81%)	33 (87%)	5 (13%)	0	100 100
6	AG	38/47 (81%)	32 (84%)	5 (13%)	1 (3%)	4 23
6	AJ	38/47 (81%)	29 (76%)	9 (24%)	0	100 100
6	AN	38/47 (81%)	35 (92%)	3 (8%)	0	100 100
6	AP	38/47 (81%)	28 (74%)	10 (26%)	0	100 100
6	AR	38/47 (81%)	31 (82%)	7 (18%)	0	100 100
6	AT	38/47 (81%)	30 (79%)	5 (13%)	3 (8%)	1 3
6	AV	38/47 (81%)	35 (92%)	2 (5%)	1 (3%)	4 23
6	AX	38/47 (81%)	32 (84%)	6 (16%)	0	100 100
6	AZ	38/47 (81%)	29 (76%)	9 (24%)	0	100 100
6	B0	38/47 (81%)	30 (79%)	8 (21%)	0	100 100
6	B2	38/47 (81%)	31 (82%)	7 (18%)	0	100 100
6	B4	38/47 (81%)	32 (84%)	6 (16%)	0	100 100
6	B6	38/47 (81%)	35 (92%)	3 (8%)	0	100 100
6	B8	38/47 (81%)	27 (71%)	10 (26%)	1 (3%)	4 23
6	BB	38/47 (81%)	33 (87%)	5 (13%)	0	100 100
6	BE	38/47 (81%)	32 (84%)	6 (16%)	0	100 100
6	BG	38/47 (81%)	32 (84%)	6 (16%)	0	100 100
6	BJ	38/47 (81%)	29 (76%)	9 (24%)	0	100 100
6	BN	38/47 (81%)	35 (92%)	3 (8%)	0	100 100
6	BP	38/47 (81%)	25 (66%)	12 (32%)	1 (3%)	4 23
6	BR	38/47 (81%)	31 (82%)	7 (18%)	0	100 100
6	BT	38/47 (81%)	30 (79%)	5 (13%)	3 (8%)	1 3

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
6	BV	38/47 (81%)	36 (95%)	2 (5%)	0	100 100
6	BX	38/47 (81%)	32 (84%)	6 (16%)	0	100 100
6	BZ	38/47 (81%)	31 (82%)	7 (18%)	0	100 100
All	All	5285/5994 (88%)	3628 (69%)	1236 (23%)	421 (8%)	1 3

All (421) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	AC	64	ALA
1	AC	70	PRO
1	AC	97	VAL
1	AC	138	ASN
1	AC	154	THR
1	AC	173	LYS
1	AC	188	LYS
1	AC	253	THR
1	AC	262	SER
1	AC	287	LEU
1	AC	317	PRO
2	AL	11	ARG
2	AL	58	PRO
2	AL	141	VAL
2	AL	204	LEU
2	AL	210	GLN
2	AL	216	LYS
2	AL	222	ASN
2	AL	277	GLU
3	AM	6	ASN
3	AM	10	ALA
3	AM	24	PRO
3	AM	81	TRP
3	AM	93	LEU
3	AM	196	LEU
3	AM	234	GLU
3	AM	235	ILE
3	AM	241	ARG
3	AM	248	ALA
3	AM	258	PHE
4	AH	5	ILE
4	AH	18	ALA
4	AH	36	ARG

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Mol	Chain	Res	Type
4	AH	55	VAL
4	AH	250	ALA
5	AD	41	SER
5	AD	44	LEU
5	AI	44	LEU
5	AO	46	TRP
5	AO	49	ASP
5	AQ	46	TRP
5	AQ	49	ASP
5	AS	54	SER
5	AU	43	ASP
5	AU	60	LYS
5	AW	43	ASP
5	A1	50	ASN
6	A2	12	ASP
5	A3	43	ASP
5	A3	47	LEU
5	A9	43	ASP
1	BC	70	PRO
1	BC	97	VAL
1	BC	157	ARG
1	BC	173	LYS
1	BC	188	LYS
1	BC	233	PHE
1	BC	253	THR
1	BC	262	SER
1	BC	287	LEU
1	BC	317	PRO
2	BL	11	ARG
2	BL	58	PRO
2	BL	120	LEU
2	BL	141	VAL
2	BL	189	PHE
2	BL	210	GLN
2	BL	244	PHE
2	BL	256	CYS
3	BM	6	ASN
3	BM	10	ALA
3	BM	16	PRO
3	BM	24	PRO
3	BM	34	PRO
3	BM	81	TRP

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Mol	Chain	Res	Type
3	BM	93	LEU
3	BM	162	PHE
3	BM	248	ALA
3	BM	308	PRO
4	BH	36	ARG
4	BH	250	ALA
5	BD	44	LEU
5	BI	43	ASP
5	BI	53	VAL
5	BK	54	SER
5	BW	43	ASP
5	BW	51	ILE
5	BY	49	ASP
5	B1	54	SER
5	B3	3	THR
5	B3	4	MET
5	B3	43	ASP
5	B3	53	VAL
5	B5	46	TRP
5	B9	43	ASP
1	AC	22	GLY
1	AC	157	ARG
1	AC	206	GLN
1	AC	266	ARG
1	AC	278	ASP
1	AC	294	SER
1	AC	303	LEU
1	AC	304	ARG
2	AL	8	LYS
2	AL	12	VAL
2	AL	81	SER
2	AL	102	ALA
2	AL	120	LEU
2	AL	139	VAL
2	AL	169	VAL
2	AL	182	HIS
2	AL	189	PHE
2	AL	200	GLY
2	AL	215	VAL
2	AL	242	GLY
2	AL	244	PHE
2	AL	256	CYS

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Mol	Chain	Res	Type
3	AM	4	TYR
3	AM	16	PRO
3	AM	34	PRO
3	AM	78	SER
3	AM	80	HIS
3	AM	103	GLY
3	AM	111	GLU
3	AM	132	ARG
3	AM	173	LYS
3	AM	176	PRO
3	AM	192	ARG
3	AM	194	GLY
3	AM	244	ALA
3	AM	249	ALA
3	AM	266	HIS
3	AM	269	ALA
3	AM	270	TRP
3	AM	280	ALA
3	AM	281	GLY
4	AH	112	GLY
4	AH	132	LYS
4	AH	146	GLU
4	AH	171	TRP
4	AH	172	VAL
4	AH	186	VAL
4	AH	204	LYS
4	AH	241	ALA
4	AH	254	ARG
5	AF	46	TRP
5	AF	49	ASP
5	AK	46	TRP
5	AU	9	TYR
5	AW	46	TRP
5	AY	43	ASP
5	AY	46	TRP
5	AY	53	VAL
5	A1	43	ASP
5	A7	46	TRP
6	A8	42	TYR
1	BC	22	GLY
1	BC	54	GLN
1	BC	64	ALA

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Mol	Chain	Res	Type
1	BC	108	ASN
1	BC	154	THR
1	BC	185	TYR
1	BC	258	ASP
1	BC	278	ASP
1	BC	279	ILE
2	BL	83	GLY
2	BL	165	TRP
2	BL	168	ASN
2	BL	182	HIS
2	BL	193	CYS
2	BL	222	ASN
2	BL	260	SER
3	BM	9	THR
3	BM	65	LEU
3	BM	78	SER
3	BM	80	HIS
3	BM	103	GLY
3	BM	132	ARG
3	BM	175	VAL
3	BM	176	PRO
3	BM	216	PHE
3	BM	230	GLY
3	BM	235	ILE
3	BM	244	ALA
3	BM	245	ALA
3	BM	249	ALA
3	BM	262	MET
3	BM	287	SER
3	BM	290	VAL
4	BH	24	PHE
4	BH	38	GLY
4	BH	53	VAL
4	BH	55	VAL
4	BH	112	GLY
4	BH	146	GLU
4	BH	177	PRO
4	BH	186	VAL
4	BH	204	LYS
4	BH	253	GLU
4	BH	254	ARG
5	BA	5	ASN

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Mol	Chain	Res	Type
5	BA	43	ASP
5	BA	53	VAL
5	BF	46	TRP
5	BI	46	TRP
5	BK	46	TRP
5	BQ	54	SER
5	BS	54	SER
5	BU	43	ASP
5	BW	5	ASN
5	BY	43	ASP
5	BY	46	TRP
5	B1	43	ASP
5	B1	49	ASP
5	B1	60	LYS
5	B5	3	THR
5	B5	47	LEU
6	B8	42	TYR
1	AC	108	ASN
1	AC	115	ASN
1	AC	185	TYR
1	AC	233	PHE
1	AC	258	ASP
1	AC	260	THR
2	AL	117	CYS
2	AL	172	GLN
2	AL	199	HIS
2	AL	219	GLU
2	AL	262	PRO
2	AL	275	TRP
3	AM	9	THR
3	AM	45	ASP
3	AM	79	VAL
3	AM	175	VAL
3	AM	205	SER
3	AM	216	PHE
3	AM	290	VAL
3	AM	308	PRO
4	AH	38	GLY
4	AH	129	GLY
4	AH	177	PRO
4	AH	253	GLU
5	AA	43	ASP

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Mol	Chain	Res	Type
5	AA	46	TRP
5	AI	49	ASP
5	AO	50	ASN
5	AO	52	PRO
6	AV	42	TYR
5	AW	47	LEU
5	AY	47	LEU
5	A5	46	TRP
6	A0	8	GLY
1	BC	113	PRO
1	BC	115	ASN
1	BC	138	ASN
1	BC	206	GLN
1	BC	266	ARG
1	BC	300	GLY
2	BL	149	GLY
2	BL	219	GLU
2	BL	245	LEU
3	BM	45	ASP
3	BM	91	PHE
3	BM	111	GLU
3	BM	143	SER
3	BM	191	ILE
3	BM	192	ARG
3	BM	205	SER
3	BM	223	ILE
3	BM	277	VAL
5	BA	55	TYR
5	BD	9	TYR
5	BY	53	VAL
5	B1	50	ASN
5	B7	46	TRP
5	B9	6	ALA
1	AC	40	MET
1	AC	74	GLU
1	AC	191	ALA
1	AC	279	ILE
1	AC	300	GLY
1	AC	324	ALA
2	AL	174	LEU
2	AL	193	CYS
2	AL	197	SER

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Mol	Chain	Res	Type
2	AL	245	LEU
2	AL	272	TRP
3	AM	91	PHE
3	AM	100	PRO
3	AM	147	SER
3	AM	162	PHE
3	AM	201	PHE
3	AM	239	THR
3	AM	245	ALA
3	AM	260	VAL
3	AM	262	MET
3	AM	265	ILE
3	AM	298	ALA
4	AH	136	MET
4	AH	163	VAL
5	AD	60	LYS
5	AI	46	TRP
5	AS	60	LYS
6	AT	12	ASP
6	AT	42	TYR
6	AT	45	TRP
5	AW	52	PRO
6	A4	44	PRO
1	BC	236	MET
1	BC	273	ILE
2	BL	5	SER
2	BL	46	GLY
2	BL	81	SER
2	BL	181	ALA
2	BL	262	PRO
3	BM	164	ARG
3	BM	171	TRP
3	BM	260	VAL
4	BH	18	ALA
4	BH	107	MET
4	BH	129	GLY
4	BH	132	LYS
4	BH	171	TRP
5	BD	43	ASP
5	BQ	53	VAL
5	BS	60	LYS
6	BT	12	ASP

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Mol	Chain	Res	Type
6	BT	42	TYR
6	BT	45	TRP
5	B7	41	SER
1	AC	47	ARG
1	AC	54	GLN
1	AC	113	PRO
1	AC	192	TYR
1	AC	239	ILE
1	AC	292	PRO
2	AL	46	GLY
2	AL	59	THR
3	AM	148	TRP
3	AM	164	ARG
3	AM	197	TYR
5	AF	45	ASN
6	AG	45	TRP
5	AI	4	MET
5	AO	51	ILE
5	AS	44	LEU
5	AW	10	LYS
5	A1	47	LEU
5	A1	51	ILE
1	BC	47	ARG
1	BC	74	GLU
1	BC	125	VAL
1	BC	149	GLY
1	BC	294	SER
2	BL	12	VAL
2	BL	43	THR
2	BL	82	TYR
2	BL	93	GLY
2	BL	169	VAL
2	BL	174	LEU
2	BL	272	TRP
3	BM	20	GLY
3	BM	266	HIS
6	BP	31	LEU
5	BY	50	ASN
5	BY	54	SER
5	B1	55	TYR
5	B5	50	ASN
1	AC	149	GLY

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Mol	Chain	Res	Type
1	AC	273	ILE
2	AL	92	GLY
2	AL	101	CYS
2	AL	121	GLY
2	AL	250	ALA
3	AM	90	PHE
3	AM	102	TYR
3	AM	135	LYS
3	AM	179	ILE
3	AM	264	SER
5	AU	11	ILE
5	A3	53	VAL
5	A7	49	ASP
1	BC	40	MET
1	BC	292	PRO
3	BM	7	ILE
4	BH	209	VAL
4	BH	242	TYR
1	AC	198	ASP
2	AL	83	GLY
2	AL	159	ILE
2	AL	279	PRO
3	AM	20	GLY
5	AS	53	VAL
1	BC	234	GLY
1	BC	239	ILE
2	BL	139	VAL
3	BM	179	ILE
4	BH	163	VAL
1	AC	103	PRO
1	AC	125	VAL
2	AL	32	VAL
4	AH	209	VAL
2	BL	32	VAL
3	BM	28	LEU
4	BH	172	VAL
5	B1	53	VAL
1	AC	28	PRO
3	AM	282	ILE
5	AK	53	VAL
4	BH	4	GLY
5	BW	52	PRO

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Mol	Chain	Res	Type
1	AC	87	VAL
1	AC	302	PRO
2	AL	47	VAL
3	AM	277	VAL
5	A7	17	PRO
4	BH	158	GLY
3	AM	28	LEU
5	A5	52	PRO
5	BK	52	PRO
5	BS	52	PRO

5.3.2 Protein sidechains (i)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	AC	265/317 (84%)	234 (88%)	31 (12%)	5 23
1	BC	265/317 (84%)	240 (91%)	25 (9%)	8 32
2	AL	228/229 (100%)	193 (85%)	35 (15%)	2 13
2	BL	228/229 (100%)	207 (91%)	21 (9%)	8 33
3	AM	256/261 (98%)	214 (84%)	42 (16%)	2 12
3	BM	256/261 (98%)	221 (86%)	35 (14%)	3 17
4	AH	210/211 (100%)	194 (92%)	16 (8%)	12 41
4	BH	210/211 (100%)	191 (91%)	19 (9%)	9 34
5	A1	48/56 (86%)	44 (92%)	4 (8%)	10 37
5	A3	47/56 (84%)	41 (87%)	6 (13%)	4 19
5	A5	48/56 (86%)	44 (92%)	4 (8%)	10 37
5	A7	48/56 (86%)	40 (83%)	8 (17%)	2 12
5	A9	50/56 (89%)	45 (90%)	5 (10%)	7 29
5	AA	44/56 (79%)	38 (86%)	6 (14%)	3 17
5	AD	47/56 (84%)	43 (92%)	4 (8%)	10 36
5	AF	49/56 (88%)	44 (90%)	5 (10%)	7 29

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
5	AI	49/56 (88%)	43 (88%)	6 (12%)	5 21
5	AK	48/56 (86%)	43 (90%)	5 (10%)	7 28
5	AO	49/56 (88%)	43 (88%)	6 (12%)	5 21
5	AQ	47/56 (84%)	44 (94%)	3 (6%)	16 48
5	AS	49/56 (88%)	45 (92%)	4 (8%)	10 37
5	AU	50/56 (89%)	44 (88%)	6 (12%)	5 22
5	AW	50/56 (89%)	44 (88%)	6 (12%)	5 22
5	AY	50/56 (89%)	43 (86%)	7 (14%)	3 16
5	B1	45/56 (80%)	43 (96%)	2 (4%)	25 60
5	B3	50/56 (89%)	47 (94%)	3 (6%)	17 50
5	B5	48/56 (86%)	45 (94%)	3 (6%)	16 48
5	B7	45/56 (80%)	39 (87%)	6 (13%)	4 18
5	B9	48/56 (86%)	46 (96%)	2 (4%)	26 61
5	BA	50/56 (89%)	45 (90%)	5 (10%)	7 29
5	BD	43/56 (77%)	42 (98%)	1 (2%)	44 74
5	BF	48/56 (86%)	45 (94%)	3 (6%)	16 48
5	BI	46/56 (82%)	41 (89%)	5 (11%)	6 26
5	BK	50/56 (89%)	43 (86%)	7 (14%)	3 16
5	BO	49/56 (88%)	42 (86%)	7 (14%)	3 16
5	BQ	50/56 (89%)	46 (92%)	4 (8%)	11 38
5	BS	49/56 (88%)	48 (98%)	1 (2%)	48 76
5	BU	50/56 (89%)	44 (88%)	6 (12%)	5 22
5	BW	48/56 (86%)	42 (88%)	6 (12%)	4 20
5	BY	45/56 (80%)	40 (89%)	5 (11%)	6 25
6	A0	33/39 (85%)	21 (64%)	12 (36%)	0 1
6	A2	33/39 (85%)	28 (85%)	5 (15%)	3 14
6	A4	33/39 (85%)	27 (82%)	6 (18%)	2 9
6	A6	33/39 (85%)	29 (88%)	4 (12%)	5 21
6	A8	33/39 (85%)	25 (76%)	8 (24%)	1 4
6	AB	33/39 (85%)	27 (82%)	6 (18%)	2 9
6	AE	33/39 (85%)	31 (94%)	2 (6%)	17 49

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
6	AG	33/39 (85%)	24 (73%)	9 (27%)	0 2
6	AJ	33/39 (85%)	28 (85%)	5 (15%)	3 14
6	AN	33/39 (85%)	26 (79%)	7 (21%)	1 6
6	AP	33/39 (85%)	26 (79%)	7 (21%)	1 6
6	AR	33/39 (85%)	27 (82%)	6 (18%)	2 9
6	AT	33/39 (85%)	27 (82%)	6 (18%)	2 9
6	AV	33/39 (85%)	27 (82%)	6 (18%)	2 9
6	AX	33/39 (85%)	29 (88%)	4 (12%)	5 21
6	AZ	33/39 (85%)	28 (85%)	5 (15%)	3 14
6	B0	33/39 (85%)	26 (79%)	7 (21%)	1 6
6	B2	33/39 (85%)	29 (88%)	4 (12%)	5 21
6	B4	33/39 (85%)	28 (85%)	5 (15%)	3 14
6	B6	33/39 (85%)	29 (88%)	4 (12%)	5 21
6	B8	33/39 (85%)	24 (73%)	9 (27%)	0 2
6	BB	33/39 (85%)	25 (76%)	8 (24%)	1 4
6	BE	33/39 (85%)	31 (94%)	2 (6%)	17 49
6	BG	33/39 (85%)	29 (88%)	4 (12%)	5 21
6	BJ	33/39 (85%)	30 (91%)	3 (9%)	9 33
6	BN	33/39 (85%)	28 (85%)	5 (15%)	3 14
6	BP	33/39 (85%)	30 (91%)	3 (9%)	9 33
6	BR	33/39 (85%)	28 (85%)	5 (15%)	3 14
6	BT	33/39 (85%)	28 (85%)	5 (15%)	3 14
6	BV	33/39 (85%)	31 (94%)	2 (6%)	17 49
6	BX	33/39 (85%)	28 (85%)	5 (15%)	3 14
6	BZ	33/39 (85%)	31 (94%)	2 (6%)	17 49
All	All	4511/5076 (89%)	3965 (88%)	546 (12%)	5 21

All (546) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	AC	20	LEU
1	AC	30	THR
1	AC	54	GLN

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Mol	Chain	Res	Type
1	AC	56	ASN
1	AC	82	LEU
1	AC	90	PHE
1	AC	95	VAL
1	AC	99	THR
1	AC	104	LYS
1	AC	112	VAL
1	AC	115	ASN
1	AC	125	VAL
1	AC	128	ARG
1	AC	135	ARG
1	AC	150	VAL
1	AC	167	VAL
1	AC	174	TYR
1	AC	178	LEU
1	AC	181	THR
1	AC	190	VAL
1	AC	212	ILE
1	AC	223	PRO
1	AC	243	LEU
1	AC	247	CYS
1	AC	254	ARG
1	AC	274	ARG
1	AC	276	VAL
1	AC	283	TYR
1	AC	285	TRP
1	AC	302	PRO
1	AC	317	PRO
2	AL	5	SER
2	AL	37	VAL
2	AL	52	TRP
2	AL	68	TYR
2	AL	71	TRP
2	AL	82	TYR
2	AL	86	MET
2	AL	90	THR
2	AL	117	CYS
2	AL	126	VAL
2	AL	138	LEU
2	AL	145	PRO
2	AL	156	PRO
2	AL	162	HIS

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Mol	Chain	Res	Type
2	AL	166	VAL
2	AL	175	HIS
2	AL	187	SER
2	AL	192	ASN
2	AL	196	LEU
2	AL	204	LEU
2	AL	208	ASN
2	AL	217	THR
2	AL	222	ASN
2	AL	223	THR
2	AL	238	ILE
2	AL	247	LEU
2	AL	248	SER
2	AL	252	TRP
2	AL	255	VAL
2	AL	256	CYS
2	AL	257	ILE
2	AL	266	ARG
2	AL	276	LEU
2	AL	280	LEU
2	AL	281	TRP
3	AM	4	TYR
3	AM	5	GLN
3	AM	21	VAL
3	AM	34	PRO
3	AM	43	ILE
3	AM	58	THR
3	AM	63	PHE
3	AM	65	LEU
3	AM	85	GLN
3	AM	91	PHE
3	AM	102	TYR
3	AM	104	LEU
3	AM	114	TRP
3	AM	122	LEU
3	AM	132	ARG
3	AM	133	THR
3	AM	138	GLU
3	AM	148	TRP
3	AM	157	TYR
3	AM	159	VAL
3	AM	162	PHE

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Mol	Chain	Res	Type
3	AM	165	PRO
3	AM	171	TRP
3	AM	177	PHE
3	AM	179	ILE
3	AM	182	HIS
3	AM	183	LEU
3	AM	186	THR
3	AM	199	ASN
3	AM	214	LEU
3	AM	215	LEU
3	AM	234	GLU
3	AM	240	HIS
3	AM	246	GLU
3	AM	247	ARG
3	AM	250	LEU
3	AM	254	TRP
3	AM	259	ASN
3	AM	265	ILE
3	AM	275	LEU
3	AM	276	THR
3	AM	307	TYR
4	AH	16	ILE
4	AH	19	PHE
4	AH	22	PHE
4	AH	27	ILE
4	AH	31	ARG
4	AH	69	LEU
4	AH	128	GLU
4	AH	133	ILE
4	AH	140	LYS
4	AH	148	ASP
4	AH	155	THR
4	AH	170	VAL
4	AH	182	LEU
4	AH	197	ILE
4	AH	223	PRO
4	AH	248	LEU
5	AA	8	LEU
5	AA	18	ARG
5	AA	37	MET
5	AA	42	THR
5	AA	49	ASP

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Mol	Chain	Res	Type
5	AA	50	ASN
6	AB	9	LEU
6	AB	20	ILE
6	AB	23	GLN
6	AB	32	VAL
6	AB	34	ILE
6	AB	40	TRP
5	AD	7	ASN
5	AD	12	TRP
5	AD	14	ILE
5	AD	29	ILE
6	AE	21	PHE
6	AE	23	GLN
5	AF	12	TRP
5	AF	27	PHE
5	AF	40	LEU
5	AF	41	SER
5	AF	43	ASP
6	AG	13	GLU
6	AG	15	LYS
6	AG	17	PHE
6	AG	20	ILE
6	AG	21	PHE
6	AG	31	LEU
6	AG	34	ILE
6	AG	37	LEU
6	AG	38	LEU
5	AI	8	LEU
5	AI	9	TYR
5	AI	18	ARG
5	AI	38	ILE
5	AI	47	LEU
5	AI	55	TYR
6	AJ	13	GLU
6	AJ	21	PHE
6	AJ	33	VAL
6	AJ	34	ILE
6	AJ	41	LEU
5	AK	9	TYR
5	AK	12	TRP
5	AK	18	ARG
5	AK	29	ILE

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Mol	Chain	Res	Type
5	AK	44	LEU
6	AN	10	THR
6	AN	13	GLU
6	AN	20	ILE
6	AN	29	PHE
6	AN	33	VAL
6	AN	34	ILE
6	AN	41	LEU
5	AO	5	ASN
5	AO	8	LEU
5	AO	13	LEU
5	AO	18	ARG
5	AO	29	ILE
5	AO	47	LEU
6	AP	20	ILE
6	AP	21	PHE
6	AP	25	MET
6	AP	34	ILE
6	AP	37	LEU
6	AP	38	LEU
6	AP	41	LEU
5	AQ	9	TYR
5	AQ	51	ILE
5	AQ	55	TYR
6	AR	13	GLU
6	AR	20	ILE
6	AR	21	PHE
6	AR	29	PHE
6	AR	34	ILE
6	AR	38	LEU
5	AS	19	ARG
5	AS	29	ILE
5	AS	33	LEU
5	AS	55	TYR
6	AT	13	GLU
6	AT	15	LYS
6	AT	16	GLU
6	AT	20	ILE
6	AT	34	ILE
6	AT	40	TRP
5	AU	2	PHE
5	AU	9	TYR

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Mol	Chain	Res	Type
5	AU	18	ARG
5	AU	38	ILE
5	AU	47	LEU
5	AU	55	TYR
6	AV	13	GLU
6	AV	20	ILE
6	AV	32	VAL
6	AV	37	LEU
6	AV	38	LEU
6	AV	41	LEU
5	AW	7	ASN
5	AW	12	TRP
5	AW	16	ASP
5	AW	19	ARG
5	AW	33	LEU
5	AW	37	MET
6	AX	12	ASP
6	AX	20	ILE
6	AX	34	ILE
6	AX	36	HIS
5	AY	2	PHE
5	AY	5	ASN
5	AY	13	LEU
5	AY	16	ASP
5	AY	18	ARG
5	AY	29	ILE
5	AY	48	ASP
6	AZ	25	MET
6	AZ	34	ILE
6	AZ	36	HIS
6	AZ	37	LEU
6	AZ	43	ARG
5	A1	8	LEU
5	A1	12	TRP
5	A1	29	ILE
5	A1	38	ILE
6	A2	13	GLU
6	A2	20	ILE
6	A2	25	MET
6	A2	29	PHE
6	A2	41	LEU
5	A3	9	TYR

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Mol	Chain	Res	Type
5	A3	11	ILE
5	A3	19	ARG
5	A3	38	ILE
5	A3	47	LEU
5	A3	56	GLN
6	A4	13	GLU
6	A4	20	ILE
6	A4	23	GLN
6	A4	33	VAL
6	A4	34	ILE
6	A4	37	LEU
5	A5	9	TYR
5	A5	20	VAL
5	A5	29	ILE
5	A5	47	LEU
6	A6	20	ILE
6	A6	34	ILE
6	A6	37	LEU
6	A6	40	TRP
5	A7	7	ASN
5	A7	9	TYR
5	A7	18	ARG
5	A7	24	ILE
5	A7	27	PHE
5	A7	28	GLN
5	A7	29	ILE
5	A7	44	LEU
6	A8	16	GLU
6	A8	20	ILE
6	A8	21	PHE
6	A8	23	GLN
6	A8	25	MET
6	A8	32	VAL
6	A8	33	VAL
6	A8	34	ILE
5	A9	18	ARG
5	A9	19	ARG
5	A9	44	LEU
5	A9	55	TYR
5	A9	56	GLN
6	A0	16	GLU
6	A0	17	PHE

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Mol	Chain	Res	Type
6	A0	20	ILE
6	A0	21	PHE
6	A0	23	GLN
6	A0	25	MET
6	A0	32	VAL
6	A0	33	VAL
6	A0	34	ILE
6	A0	36	HIS
6	A0	37	LEU
6	A0	40	TRP
1	BC	20	LEU
1	BC	21	LEU
1	BC	30	THR
1	BC	54	GLN
1	BC	58	PRO
1	BC	90	PHE
1	BC	98	THR
1	BC	104	LYS
1	BC	115	ASN
1	BC	128	ARG
1	BC	130	MET
1	BC	150	VAL
1	BC	167	VAL
1	BC	174	TYR
1	BC	190	VAL
1	BC	212	ILE
1	BC	243	LEU
1	BC	247	CYS
1	BC	254	ARG
1	BC	273	ILE
1	BC	274	ARG
1	BC	283	TYR
1	BC	285	TRP
1	BC	302	PRO
1	BC	317	PRO
2	BL	52	TRP
2	BL	71	TRP
2	BL	82	TYR
2	BL	86	MET
2	BL	90	THR
2	BL	111	LEU
2	BL	116	ILE

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Mol	Chain	Res	Type
2	BL	117	CYS
2	BL	145	PRO
2	BL	172	GLN
2	BL	174	LEU
2	BL	192	ASN
2	BL	201	SER
2	BL	208	ASN
2	BL	222	ASN
2	BL	247	LEU
2	BL	252	TRP
2	BL	266	ARG
2	BL	273	ASN
2	BL	276	LEU
2	BL	281	TRP
3	BM	5	GLN
3	BM	16	PRO
3	BM	21	VAL
3	BM	27	ASN
3	BM	34	PRO
3	BM	63	PHE
3	BM	65	LEU
3	BM	102	TYR
3	BM	114	TRP
3	BM	126	ILE
3	BM	132	ARG
3	BM	138	GLU
3	BM	148	TRP
3	BM	155	PHE
3	BM	162	PHE
3	BM	171	TRP
3	BM	177	PHE
3	BM	179	ILE
3	BM	182	HIS
3	BM	183	LEU
3	BM	186	THR
3	BM	214	LEU
3	BM	218	MET
3	BM	222	THR
3	BM	229	PHE
3	BM	234	GLU
3	BM	235	ILE
3	BM	246	GLU

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Mol	Chain	Res	Type
3	BM	254	TRP
3	BM	255	THR
3	BM	259	ASN
3	BM	267	ARG
3	BM	275	LEU
3	BM	276	THR
3	BM	307	TYR
4	BH	13	GLN
4	BH	16	ILE
4	BH	22	PHE
4	BH	27	ILE
4	BH	29	TYR
4	BH	31	ARG
4	BH	54	LYS
4	BH	69	LEU
4	BH	123	CYS
4	BH	128	GLU
4	BH	140	LYS
4	BH	148	ASP
4	BH	155	THR
4	BH	182	LEU
4	BH	185	GLU
4	BH	186	VAL
4	BH	197	ILE
4	BH	225	LEU
4	BH	235	GLU
5	BA	8	LEU
5	BA	18	ARG
5	BA	37	MET
5	BA	42	THR
5	BA	55	TYR
6	BB	7	THR
6	BB	9	LEU
6	BB	20	ILE
6	BB	23	GLN
6	BB	26	TYR
6	BB	32	VAL
6	BB	34	ILE
6	BB	40	TRP
5	BD	12	TRP
6	BE	16	GLU
6	BE	23	GLN

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Mol	Chain	Res	Type
5	BF	4	MET
5	BF	12	TRP
5	BF	49	ASP
6	BG	13	GLU
6	BG	21	PHE
6	BG	23	GLN
6	BG	24	SER
5	BI	8	LEU
5	BI	9	TYR
5	BI	18	ARG
5	BI	47	LEU
5	BI	55	TYR
6	BJ	13	GLU
6	BJ	33	VAL
6	BJ	34	ILE
5	BK	9	TYR
5	BK	12	TRP
5	BK	18	ARG
5	BK	29	ILE
5	BK	44	LEU
5	BK	48	ASP
5	BK	55	TYR
6	BN	13	GLU
6	BN	24	SER
6	BN	29	PHE
6	BN	33	VAL
6	BN	34	ILE
5	BO	7	ASN
5	BO	9	TYR
5	BO	21	LEU
5	BO	43	ASP
5	BO	47	LEU
5	BO	51	ILE
5	BO	55	TYR
6	BP	21	PHE
6	BP	25	MET
6	BP	40	TRP
5	BQ	9	TYR
5	BQ	43	ASP
5	BQ	45	ASN
5	BQ	55	TYR
6	BR	20	ILE

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Mol	Chain	Res	Type
6	BR	21	PHE
6	BR	23	GLN
6	BR	29	PHE
6	BR	38	LEU
5	BS	55	TYR
6	BT	13	GLU
6	BT	20	ILE
6	BT	21	PHE
6	BT	29	PHE
6	BT	40	TRP
5	BU	9	TYR
5	BU	12	TRP
5	BU	18	ARG
5	BU	35	ILE
5	BU	47	LEU
5	BU	55	TYR
6	BV	13	GLU
6	BV	20	ILE
5	BW	7	ASN
5	BW	12	TRP
5	BW	14	ILE
5	BW	22	VAL
5	BW	37	MET
5	BW	55	TYR
6	BX	24	SER
6	BX	34	ILE
6	BX	36	HIS
6	BX	38	LEU
6	BX	44	PRO
5	BY	9	TYR
5	BY	18	ARG
5	BY	29	ILE
5	BY	37	MET
5	BY	56	GLN
6	BZ	29	PHE
6	BZ	36	HIS
5	B1	9	TYR
5	B1	18	ARG
6	B2	20	ILE
6	B2	24	SER
6	B2	25	MET
6	B2	41	LEU

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Mol	Chain	Res	Type
5	B3	45	ASN
5	B3	47	LEU
5	B3	56	GLN
6	B4	13	GLU
6	B4	20	ILE
6	B4	33	VAL
6	B4	34	ILE
6	B4	37	LEU
5	B5	9	TYR
5	B5	29	ILE
5	B5	47	LEU
6	B6	20	ILE
6	B6	34	ILE
6	B6	37	LEU
6	B6	40	TRP
5	B7	9	TYR
5	B7	18	ARG
5	B7	24	ILE
5	B7	28	GLN
5	B7	44	LEU
5	B7	47	LEU
6	B8	16	GLU
6	B8	20	ILE
6	B8	21	PHE
6	B8	23	GLN
6	B8	25	MET
6	B8	32	VAL
6	B8	33	VAL
6	B8	34	ILE
6	B8	41	LEU
5	B9	2	PHE
5	B9	44	LEU
6	B0	17	PHE
6	B0	21	PHE
6	B0	25	MET
6	B0	32	VAL
6	B0	34	ILE
6	B0	36	HIS
6	B0	40	TRP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (124) such sidechains are listed below:

Mol	Chain	Res	Type
1	AC	45	ASN
1	AC	54	GLN
1	AC	57	GLN
1	AC	77	GLN
1	AC	80	GLN
1	AC	138	ASN
1	AC	159	ASN
1	AC	206	GLN
1	AC	228	GLN
1	AC	238	ASN
1	AC	252	ASN
1	AC	322	GLN
2	AL	96	GLN
2	AL	192	ASN
2	AL	222	ASN
2	AL	273	ASN
3	AM	85	GLN
3	AM	101	GLN
3	AM	195	ASN
3	AM	199	ASN
3	AM	240	HIS
3	AM	315	ASN
4	AH	7	HIS
4	AH	88	ASN
4	AH	189	ASN
4	AH	218	HIS
4	AH	227	ASN
5	AA	28	GLN
6	AB	18	HIS
6	AB	23	GLN
5	AD	5	ASN
5	AD	7	ASN
5	AD	28	GLN
6	AE	18	HIS
5	AF	7	ASN
5	AF	56	GLN
6	AG	18	HIS
6	AG	23	GLN
5	AI	7	ASN
5	AI	28	GLN
5	AK	5	ASN
5	AO	5	ASN
5	AO	7	ASN

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Mol	Chain	Res	Type
5	AO	28	GLN
6	AP	18	HIS
5	AS	28	GLN
5	AS	45	ASN
6	AT	23	GLN
5	AW	5	ASN
5	AW	7	ASN
5	AW	56	GLN
6	AZ	18	HIS
6	AZ	23	GLN
5	A1	45	ASN
5	A1	56	GLN
5	A3	28	GLN
5	A3	45	ASN
5	A3	56	GLN
6	A4	23	GLN
5	A5	45	ASN
5	A5	56	GLN
5	A7	5	ASN
5	A7	7	ASN
5	A7	28	GLN
5	A9	7	ASN
5	A9	28	GLN
5	A9	56	GLN
1	BC	54	GLN
1	BC	57	GLN
1	BC	77	GLN
1	BC	80	GLN
1	BC	159	ASN
1	BC	184	ASN
1	BC	206	GLN
1	BC	238	ASN
1	BC	322	GLN
2	BL	96	GLN
2	BL	177	HIS
2	BL	192	ASN
2	BL	273	ASN
3	BM	5	GLN
3	BM	27	ASN
3	BM	47	GLN
3	BM	80	HIS
3	BM	195	ASN

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Mol	Chain	Res	Type
3	BM	199	ASN
3	BM	259	ASN
3	BM	315	ASN
4	BH	13	GLN
4	BH	88	ASN
4	BH	189	ASN
4	BH	218	HIS
4	BH	227	ASN
5	BA	7	ASN
6	BB	18	HIS
6	BB	23	GLN
6	BE	18	HIS
5	BF	7	ASN
5	BF	56	GLN
6	BG	18	HIS
5	BI	28	GLN
5	BK	28	GLN
5	BK	45	ASN
5	BK	56	GLN
5	BO	5	ASN
5	BO	7	ASN
5	BS	45	ASN
6	BT	23	GLN
5	BU	5	ASN
5	BU	28	GLN
5	BU	56	GLN
5	BW	7	ASN
6	BZ	23	GLN
5	B1	28	GLN
5	B1	56	GLN
5	B3	45	ASN
5	B3	56	GLN
6	B4	23	GLN
5	B5	28	GLN
6	B6	23	GLN
5	B7	56	GLN
5	B9	5	ASN
5	B9	7	ASN
5	B9	45	ASN

5.3.3 RNA

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates

There are no oligosaccharides in this entry.

5.6 Ligand geometry

Of 169 ligands modelled in this entry, 36 are monoatomic - leaving 133 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
9	BCL	B9	102	-	69,74,74	1.56	11 (15%)	79,115,115	2.36	24 (30%)
16	PO4	AH	302	-	4,4,4	1.69	0	6,6,6	0.44	0
9	BCL	AD	102	-	69,74,74	1.53	11 (15%)	79,115,115	2.30	23 (29%)
9	BCL	BQ	103	-	69,74,74	1.67	17 (24%)	79,115,115	2.19	18 (22%)
14	CRT	A5	103	-	43,43,43	1.23	5 (11%)	48,54,54	1.80	17 (35%)
14	CRT	BG	102	-	43,43,43	1.25	3 (6%)	48,54,54	1.82	16 (33%)
15	PEF	BM	407	-	18,18,46	3.17	7 (38%)	21,23,51	2.02	6 (28%)
7	HEM	AC	504	1	50,50,50	1.90	14 (28%)	67,82,82	1.47	9 (13%)
7	HEM	AC	503	1	50,50,50	1.84	14 (28%)	67,82,82	1.37	9 (13%)
13	MQ8	AM	405	-	54,54,54	1.14	5 (9%)	67,69,69	1.55	12 (17%)
7	HEM	BC	502	1	50,50,50	1.88	17 (34%)	67,82,82	1.46	11 (16%)
9	BCL	AM	401	-	69,74,74	1.32	8 (11%)	79,115,115	2.25	23 (29%)
14	CRT	AB	102	-	43,43,43	1.25	4 (9%)	48,54,54	1.99	17 (35%)
9	BCL	BX	101	-	69,74,74	1.83	20 (28%)	79,115,115	2.41	23 (29%)
16	PO4	BH	301	-	4,4,4	1.68	0	6,6,6	0.44	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\|>2$	Counts	RMSZ	$\#\|Z\|>2$
9	BCL	BT	101	-	69,74,74	1.64	14 (20%)	79,115,115	2.26	28 (35%)
7	HEM	AC	501	1	50,50,50	1.87	13 (26%)	67,82,82	1.49	7 (10%)
11	UQ8	BL	304	-	53,53,53	1.44	3 (5%)	66,67,67	1.63	14 (21%)
9	BCL	AM	402	-	69,74,74	1.32	9 (13%)	79,115,115	2.33	23 (29%)
9	BCL	AW	101	-	69,74,74	1.83	14 (20%)	79,115,115	3.10	28 (35%)
9	BCL	BJ	101	-	69,74,74	1.70	16 (23%)	79,115,115	2.16	24 (30%)
9	BCL	BY	102	-	69,74,74	1.51	12 (17%)	79,115,115	2.27	21 (26%)
14	CRT	B2	102	-	43,43,43	1.94	9 (20%)	48,54,54	2.14	17 (35%)
16	PO4	AM	410	-	4,4,4	1.73	0	6,6,6	0.47	0
9	BCL	A8	101	-	69,74,74	1.91	19 (27%)	79,115,115	2.87	28 (35%)
9	BCL	BF	102	-	69,74,74	1.77	14 (20%)	79,115,115	2.39	22 (27%)
9	BCL	A0	102	-	69,74,74	2.04	15 (21%)	79,115,115	2.93	26 (32%)
9	BCL	AV	102	-	69,74,74	1.79	16 (23%)	79,115,115	2.16	26 (32%)
9	BCL	AL	301	-	69,74,74	1.52	15 (21%)	79,115,115	2.18	23 (29%)
14	CRT	BV	102	-	43,43,43	1.54	10 (23%)	48,54,54	1.87	13 (27%)
9	BCL	AF	102	-	69,74,74	1.57	13 (18%)	79,115,115	2.24	24 (30%)
14	CRT	AW	102	-	43,43,43	1.54	6 (13%)	48,54,54	1.85	20 (41%)
9	BCL	BI	102	-	69,74,74	1.60	13 (18%)	79,115,115	2.30	22 (27%)
9	BCL	BL	303	-	69,74,74	1.33	10 (14%)	79,115,115	2.28	22 (27%)
9	BCL	BG	101	-	69,74,74	1.67	13 (18%)	79,115,115	2.35	23 (29%)
9	BCL	A5	102	-	69,74,74	1.88	17 (24%)	79,115,115	2.58	26 (32%)
14	CRT	AJ	102	-	43,43,43	1.28	4 (9%)	48,54,54	2.07	16 (33%)
14	CRT	BN	102	-	43,43,43	1.32	7 (16%)	48,54,54	1.97	16 (33%)
9	BCL	BK	102	-	69,74,74	1.57	12 (17%)	79,115,115	2.25	23 (29%)
14	CRT	AM	406	-	43,43,43	1.38	3 (6%)	48,54,54	1.64	10 (20%)
14	CRT	AT	102	-	43,43,43	1.28	4 (9%)	48,54,54	1.90	14 (29%)
14	CRT	BU	103	-	43,43,43	1.60	6 (13%)	48,54,54	2.38	16 (33%)
9	BCL	BD	102	-	69,74,74	1.83	17 (24%)	79,115,115	2.32	23 (29%)
14	CRT	BF	103	-	43,43,43	1.28	4 (9%)	48,54,54	2.38	17 (35%)
7	HEM	AC	502	1	50,50,50	1.93	16 (32%)	67,82,82	1.43	9 (13%)
9	BCL	AL	303	-	69,74,74	1.35	10 (14%)	79,115,115	2.28	20 (25%)
15	PEF	AH	301	-	18,18,46	3.17	7 (38%)	21,23,51	2.02	6 (28%)
9	BCL	BL	301	-	69,74,74	1.53	13 (18%)	79,115,115	2.18	23 (29%)
14	CRT	AX	102	-	43,43,43	1.60	9 (20%)	48,54,54	2.27	14 (29%)
14	CRT	AG	102	-	43,43,43	1.25	4 (9%)	48,54,54	1.87	17 (35%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\|>2$	Counts	RMSZ	$\#\|Z\|>2$
9	BCL	A9	102	-	69,74,74	1.64	15 (21%)	79,115,115	2.80	28 (35%)
9	BCL	AK	102	-	69,74,74	1.53	13 (18%)	79,115,115	2.34	24 (30%)
15	PEF	AM	408	-	13,13,46	2.98	5 (38%)	15,16,51	1.18	1 (6%)
15	PEF	BQ	101	-	46,46,46	2.19	7 (15%)	49,51,51	1.40	6 (12%)
9	BCL	AO	102	-	69,74,74	1.56	14 (20%)	79,115,115	2.24	26 (32%)
9	BCL	BM	402	-	69,74,74	1.42	13 (18%)	79,115,115	2.44	23 (29%)
9	BCL	AA	101	-	69,74,74	1.54	14 (20%)	79,115,115	2.21	21 (26%)
10	BPH	AL	302	-	59,70,70	1.58	6 (10%)	59,101,101	1.67	10 (16%)
9	BCL	AR	101	-	69,74,74	1.66	16 (23%)	79,115,115	2.31	24 (30%)
9	BCL	AS	103	-	69,74,74	1.66	14 (20%)	79,115,115	2.32	25 (31%)
9	BCL	A6	101	-	69,74,74	2.13	13 (18%)	79,115,115	2.45	27 (34%)
16	PO4	A3	101	-	4,4,4	1.73	1 (25%)	6,6,6	0.46	0
10	BPH	AM	403	-	59,70,70	1.65	7 (11%)	59,101,101	1.88	9 (15%)
9	BCL	BB	101	-	69,74,74	1.85	14 (20%)	79,115,115	2.51	25 (31%)
14	CRT	BO	103	-	43,43,43	1.24	3 (6%)	48,54,54	1.89	15 (31%)
9	BCL	AG	101	-	69,74,74	1.63	13 (18%)	79,115,115	2.24	24 (30%)
9	BCL	B8	101	-	69,74,74	1.64	15 (21%)	79,115,115	2.57	27 (34%)
14	CRT	AP	102	-	43,43,43	1.49	7 (16%)	48,54,54	1.87	16 (33%)
9	BCL	B2	101	-	69,74,74	1.66	13 (18%)	79,115,115	2.30	25 (31%)
9	BCL	AP	101	-	69,74,74	1.70	13 (18%)	79,115,115	2.42	26 (32%)
9	BCL	AE	101	-	69,74,74	1.75	14 (20%)	79,115,115	2.45	26 (32%)
7	HEM	BC	504	1	50,50,50	1.97	15 (30%)	67,82,82	1.44	8 (11%)
14	CRT	BP	102	-	43,43,43	1.60	7 (16%)	48,54,54	1.75	15 (31%)
9	BCL	AI	102	-	69,74,74	1.59	12 (17%)	79,115,115	2.25	23 (29%)
9	BCL	A3	104	-	69,74,74	1.66	15 (21%)	79,115,115	2.80	26 (32%)
7	HEM	BC	501	1	50,50,50	1.95	14 (28%)	67,82,82	1.43	8 (11%)
9	BCL	B0	102	-	69,74,74	1.49	14 (20%)	79,115,115	2.36	27 (34%)
13	MQ8	BM	405	-	54,54,54	1.19	7 (12%)	67,69,69	1.37	11 (16%)
9	BCL	BA	101	-	69,74,74	1.47	11 (15%)	79,115,115	2.21	21 (26%)
9	BCL	B1	102	-	69,74,74	1.49	14 (20%)	79,115,115	2.24	20 (25%)
9	BCL	BU	102	-	69,74,74	1.67	10 (14%)	79,115,115	2.29	21 (26%)
9	BCL	AT	101	-	69,74,74	1.78	18 (26%)	79,115,115	2.26	25 (31%)
9	BCL	BE	101	-	69,74,74	1.98	18 (26%)	79,115,115	2.14	23 (29%)
14	CRT	BW	103	-	43,43,43	1.50	7 (16%)	48,54,54	1.75	16 (33%)
9	BCL	A3	103	-	69,74,74	1.84	14 (20%)	79,115,115	2.29	23 (29%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\|>2$	Counts	RMSZ	$\#\|Z\|>2$
14	CRT	A0	101	-	43,43,43	1.36	6 (13%)	48,54,54	2.01	13 (27%)
9	BCL	AQ	102	-	69,74,74	1.56	15 (21%)	79,115,115	2.26	20 (25%)
9	BCL	BM	401	-	69,74,74	1.30	9 (13%)	79,115,115	2.26	25 (31%)
9	BCL	BN	101	-	69,74,74	1.57	13 (18%)	79,115,115	2.17	25 (31%)
9	BCL	A1	102	-	69,74,74	2.61	20 (28%)	79,115,115	2.53	24 (30%)
14	CRT	B1	103	-	43,43,43	1.23	3 (6%)	48,54,54	2.07	18 (37%)
11	UQ8	AL	304	-	53,53,53	1.49	3 (5%)	66,67,67	1.74	14 (21%)
9	BCL	B5	102	-	69,74,74	1.58	12 (17%)	79,115,115	2.27	24 (30%)
14	CRT	AS	104	-	43,43,43	1.26	2 (4%)	48,54,54	2.23	15 (31%)
9	BCL	BZ	101	-	69,74,74	1.60	12 (17%)	79,115,115	2.27	27 (34%)
9	BCL	B6	101	-	69,74,74	1.66	13 (18%)	79,115,115	2.33	28 (35%)
9	BCL	AN	101	-	69,74,74	1.67	14 (20%)	79,115,115	2.38	25 (31%)
9	BCL	AU	102	-	69,74,74	1.83	17 (24%)	79,115,115	2.73	30 (37%)
9	BCL	BO	102	-	69,74,74	1.53	13 (18%)	79,115,115	2.31	21 (26%)
14	CRT	BA	102	-	43,43,43	1.25	2 (4%)	48,54,54	1.73	14 (29%)
10	BPH	BM	403	-	59,70,70	1.68	5 (8%)	59,101,101	1.94	9 (15%)
9	BCL	B7	103	-	69,74,74	1.62	13 (18%)	79,115,115	2.37	24 (30%)
14	CRT	AN	102	-	43,43,43	1.25	4 (9%)	48,54,54	1.93	18 (37%)
9	BCL	BV	101	-	69,74,74	1.58	12 (17%)	79,115,115	2.28	29 (36%)
14	CRT	B5	103	-	43,43,43	1.26	4 (9%)	48,54,54	1.74	14 (29%)
9	BCL	BW	102	-	69,74,74	1.54	13 (18%)	79,115,115	2.29	23 (29%)
14	CRT	BB	102	-	43,43,43	1.34	4 (9%)	48,54,54	2.01	18 (37%)
15	PEF	AM	407	-	18,18,46	3.17	7 (38%)	21,23,51	2.02	6 (28%)
15	PEF	AS	101	-	46,46,46	2.20	7 (15%)	49,51,51	1.40	6 (12%)
14	CRT	A7	102	-	43,43,43	3.02	6 (13%)	48,54,54	3.75	14 (29%)
14	CRT	A2	102	-	43,43,43	1.50	8 (18%)	48,54,54	1.86	14 (29%)
14	CRT	AR	102	-	43,43,43	1.22	3 (6%)	48,54,54	1.89	15 (31%)
9	BCL	A7	103	-	69,74,74	2.12	16 (23%)	79,115,115	2.42	25 (31%)
14	CRT	A1	103	-	43,43,43	1.37	6 (13%)	48,54,54	2.46	17 (35%)
14	CRT	BS	103	-	43,43,43	1.24	2 (4%)	48,54,54	1.84	16 (33%)
7	HEM	BC	503	1	50,50,50	1.88	15 (30%)	67,82,82	1.38	10 (14%)
14	CRT	B7	102	-	43,43,43	1.57	6 (13%)	48,54,54	2.06	13 (27%)
14	CRT	AA	102	-	43,43,43	1.26	2 (4%)	48,54,54	1.73	12 (25%)
14	CRT	B0	101	-	43,43,43	1.59	10 (23%)	48,54,54	1.99	15 (31%)
9	BCL	AX	101	-	69,74,74	2.16	15 (21%)	79,115,115	2.63	24 (30%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\|>2$	Counts	RMSZ	$\#\|Z\|>2$
9	BCL	BS	102	-	69,74,74	1.38	10 (14%)	79,115,115	2.35	22 (27%)
9	BCL	A2	101	-	69,74,74	1.68	16 (23%)	79,115,115	2.48	22 (27%)
9	BCL	BP	101	-	69,74,74	1.77	19 (27%)	79,115,115	2.47	27 (34%)
9	BCL	B4	101	-	69,74,74	1.65	17 (24%)	79,115,115	2.40	26 (32%)
9	BCL	AB	101	-	69,74,74	1.63	15 (21%)	79,115,115	2.39	24 (30%)
9	BCL	AJ	101	-	69,74,74	1.56	13 (18%)	79,115,115	2.35	23 (29%)
10	BPH	BL	302	-	59,70,70	1.56	8 (13%)	59,101,101	1.65	9 (15%)
9	BCL	BQ	104	-	69,74,74	1.61	13 (18%)	79,115,115	2.20	25 (31%)
9	BCL	AZ	101	-	69,74,74	2.70	19 (27%)	79,115,115	3.07	24 (30%)
9	BCL	AY	102	-	69,74,74	1.85	15 (21%)	79,115,115	2.37	23 (29%)
9	BCL	B3	102	-	69,74,74	1.45	10 (14%)	79,115,115	2.27	23 (29%)
15	PEF	AM	409	-	46,46,46	2.20	7 (15%)	49,51,51	1.33	6 (12%)
14	CRT	BM	406	-	43,43,43	1.30	4 (9%)	48,54,54	1.68	13 (27%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
9	BCL	B9	102	-	-	14/41/137/137	-
9	BCL	AD	102	-	-	17/41/137/137	-
9	BCL	BQ	103	-	-	15/41/137/137	-
14	CRT	A5	103	-	-	1/51/51/51	-
14	CRT	BG	102	-	-	2/51/51/51	-
15	PEF	BM	407	-	-	11/20/20/50	-
7	HEM	AC	504	1	-	6/14/54/54	-
7	HEM	AC	503	1	-	9/14/54/54	-
13	MQ8	AM	405	-	-	10/47/67/67	0/2/2/2
7	HEM	BC	502	1	-	5/14/54/54	-
9	BCL	AM	401	-	-	12/41/137/137	-
14	CRT	AB	102	-	-	5/51/51/51	-
9	BCL	BX	101	-	-	12/41/137/137	-
9	BCL	BT	101	-	-	6/41/137/137	-
7	HEM	AC	501	1	-	7/14/54/54	-
11	UQ8	BL	304	-	-	4/51/75/75	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
9	BCL	AM	402	-	-	13/41/137/137	-
9	BCL	AW	101	-	-	16/41/137/137	-
9	BCL	BJ	101	-	-	7/41/137/137	-
9	BCL	BY	102	-	-	15/41/137/137	-
14	CRT	B2	102	-	-	2/51/51/51	-
9	BCL	A8	101	-	-	8/41/137/137	-
9	BCL	BF	102	-	-	13/41/137/137	-
9	BCL	A0	102	-	-	15/41/137/137	-
9	BCL	AV	102	-	-	13/41/137/137	-
9	BCL	AL	301	-	-	9/41/137/137	-
14	CRT	BV	102	-	-	2/51/51/51	-
9	BCL	AF	102	-	-	9/41/137/137	-
14	CRT	AW	102	-	-	0/51/51/51	-
9	BCL	BI	102	-	-	14/41/137/137	-
9	BCL	BL	303	-	-	11/41/137/137	-
9	BCL	BG	101	-	-	11/41/137/137	-
9	BCL	A5	102	-	-	14/41/137/137	-
14	CRT	AJ	102	-	-	1/51/51/51	-
14	CRT	BN	102	-	-	1/51/51/51	-
9	BCL	BK	102	-	-	12/41/137/137	-
14	CRT	AM	406	-	-	0/51/51/51	-
14	CRT	AT	102	-	-	2/51/51/51	-
14	CRT	BU	103	-	-	2/51/51/51	-
9	BCL	BD	102	-	-	13/41/137/137	-
14	CRT	BF	103	-	-	1/51/51/51	-
7	HEM	AC	502	1	-	9/14/54/54	-
9	BCL	AL	303	-	-	12/41/137/137	-
15	PEF	AH	301	-	-	9/20/20/50	-
9	BCL	BL	301	-	-	11/41/137/137	-
14	CRT	AX	102	-	-	1/51/51/51	-
14	CRT	AG	102	-	-	1/51/51/51	-
9	BCL	A9	102	-	-	19/41/137/137	-
9	BCL	AK	102	-	-	15/41/137/137	-
15	PEF	AM	408	-	-	8/13/13/50	-
15	PEF	BQ	101	-	-	22/50/50/50	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
9	BCL	AO	102	-	-	17/41/137/137	-
9	BCL	BM	402	-	-	13/41/137/137	-
9	BCL	AA	101	-	-	17/41/137/137	-
10	BPH	AL	302	-	-	12/37/105/105	0/5/6/6
9	BCL	AR	101	-	-	14/41/137/137	-
9	BCL	AS	103	-	-	17/41/137/137	-
9	BCL	A6	101	-	-	18/41/137/137	-
10	BPH	AM	403	-	-	9/37/105/105	0/5/6/6
9	BCL	BB	101	-	-	9/41/137/137	-
14	CRT	BO	103	-	-	1/51/51/51	-
9	BCL	AG	101	-	-	12/41/137/137	-
9	BCL	B8	101	-	-	11/41/137/137	-
14	CRT	AP	102	-	-	2/51/51/51	-
9	BCL	B2	101	-	-	11/41/137/137	-
9	BCL	AP	101	-	-	8/41/137/137	-
9	BCL	AE	101	-	-	9/41/137/137	-
7	HEM	BC	504	1	-	9/14/54/54	-
14	CRT	BP	102	-	-	2/51/51/51	-
9	BCL	AI	102	-	-	13/41/137/137	-
9	BCL	A3	104	-	-	16/41/137/137	-
7	HEM	BC	501	1	-	6/14/54/54	-
9	BCL	B0	102	-	-	12/41/137/137	-
13	MQ8	BM	405	-	-	14/47/67/67	0/2/2/2
9	BCL	BA	101	-	-	12/41/137/137	-
9	BCL	B1	102	-	-	14/41/137/137	-
9	BCL	BU	102	-	-	15/41/137/137	-
9	BCL	AT	101	-	-	9/41/137/137	-
9	BCL	BE	101	-	-	14/41/137/137	-
14	CRT	BW	103	-	-	0/51/51/51	-
9	BCL	A3	103	-	-	17/41/137/137	-
14	CRT	A0	101	-	-	2/51/51/51	-
9	BCL	AQ	102	-	-	17/41/137/137	-
9	BCL	BM	401	-	-	12/41/137/137	-
9	BCL	BN	101	-	-	12/41/137/137	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
9	BCL	A1	102	-	-	16/41/137/137	-
14	CRT	B1	103	-	-	2/51/51/51	-
11	UQ8	AL	304	-	-	9/51/75/75	0/1/1/1
9	BCL	B5	102	-	-	13/41/137/137	-
14	CRT	AS	104	-	-	2/51/51/51	-
9	BCL	BZ	101	-	-	13/41/137/137	-
9	BCL	B6	101	-	-	17/41/137/137	-
9	BCL	AN	101	-	-	13/41/137/137	-
9	BCL	AU	102	-	-	18/41/137/137	-
9	BCL	BO	102	-	-	15/41/137/137	-
14	CRT	BA	102	-	-	1/51/51/51	-
10	BPH	BM	403	-	-	9/37/105/105	0/5/6/6
9	BCL	B7	103	-	-	11/41/137/137	-
14	CRT	AN	102	-	-	1/51/51/51	-
9	BCL	BV	101	-	-	15/41/137/137	-
14	CRT	B5	103	-	-	1/51/51/51	-
9	BCL	BW	102	-	-	15/41/137/137	-
14	CRT	BB	102	-	-	2/51/51/51	-
15	PEF	AM	407	-	-	14/20/20/50	-
15	PEF	AS	101	-	-	27/50/50/50	-
14	CRT	A7	102	-	-	4/51/51/51	-
14	CRT	A2	102	-	-	1/51/51/51	-
14	CRT	AR	102	-	-	1/51/51/51	-
9	BCL	A7	103	-	-	13/41/137/137	-
14	CRT	A1	103	-	-	2/51/51/51	-
14	CRT	BS	103	-	-	2/51/51/51	-
7	HEM	BC	503	1	-	10/14/54/54	-
14	CRT	B7	102	-	-	2/51/51/51	-
14	CRT	AA	102	-	-	1/51/51/51	-
14	CRT	B0	101	-	-	5/51/51/51	-
9	BCL	AX	101	-	-	18/41/137/137	-
9	BCL	BS	102	-	-	15/41/137/137	-
9	BCL	A2	101	-	-	15/41/137/137	-
9	BCL	BP	101	-	-	6/41/137/137	-
9	BCL	B4	101	-	-	7/41/137/137	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
9	BCL	AB	101	-	-	10/41/137/137	-
9	BCL	AJ	101	-	-	11/41/137/137	-
10	BPH	BL	302	-	-	12/37/105/105	0/5/6/6
9	BCL	BQ	104	-	-	10/41/137/137	-
9	BCL	AZ	101	-	-	18/41/137/137	-
9	BCL	AY	102	-	-	17/41/137/137	-
9	BCL	B3	102	-	-	14/41/137/137	-
15	PEF	AM	409	-	-	19/50/50/50	-
14	CRT	BM	406	-	-	0/51/51/51	-

All (1387) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	AZ	101	BCL	O2A-C1	-18.12	0.97	1.46
9	A1	102	BCL	O2A-C1	15.94	1.89	1.46
14	A7	102	CRT	C21-C20	12.93	1.71	1.36
14	A7	102	CRT	C21-C22	12.54	1.82	1.43
9	A6	101	BCL	O2A-C1	12.28	1.79	1.46
9	AX	101	BCL	O2A-C1	11.74	1.77	1.46
9	A0	102	BCL	O2A-C1	11.17	1.76	1.46
9	A7	103	BCL	O2A-C1	10.71	1.75	1.46
10	BM	403	BPH	C1D-C2D	9.68	1.50	1.39
10	AM	403	BPH	C1D-C2D	9.15	1.49	1.39
10	AL	302	BPH	C1D-C2D	8.71	1.49	1.39
9	A5	102	BCL	O2A-C1	-8.44	1.23	1.46
15	AM	409	PEF	O4-C10	8.30	1.47	1.22
15	AS	101	PEF	O4-C10	8.28	1.47	1.22
15	BQ	101	PEF	O4-C10	8.27	1.47	1.22
10	BL	302	BPH	C1D-C2D	7.97	1.48	1.39
9	BE	101	BCL	CAA-CBA	-7.85	1.29	1.52
9	AY	102	BCL	CAA-C2A	7.83	1.68	1.54
9	A3	103	BCL	CAA-C2A	7.41	1.67	1.54
11	AL	304	UQ8	C43-C44	7.39	1.54	1.32
11	BL	304	UQ8	C43-C44	7.36	1.54	1.32
9	AW	101	BCL	O2A-C1	7.33	1.65	1.46
15	AM	407	PEF	O4-C10	7.30	1.47	1.21
15	AH	301	PEF	O4-C10	7.29	1.47	1.21
15	AM	408	PEF	O4-C10	7.29	1.47	1.21
15	BM	407	PEF	O4-C10	7.27	1.47	1.21
9	BB	101	BCL	MG-NA	7.05	2.23	2.06

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	A8	101	BCL	MG-NA	6.28	2.21	2.06
15	AH	301	PEF	O2-C10	6.14	1.48	1.35
14	B2	102	CRT	C4-C1	6.13	1.63	1.52
15	AM	407	PEF	O2-C10	6.13	1.48	1.35
15	BM	407	PEF	O2-C10	6.11	1.48	1.35
9	B6	101	BCL	CAA-CBA	-6.05	1.34	1.52
9	AD	102	BCL	O2D-CGD	6.02	1.48	1.33
14	BU	103	CRT	C4-C1	6.00	1.63	1.52
15	AS	101	PEF	O5-C30	5.99	1.40	1.22
15	AM	409	PEF	O5-C30	5.98	1.40	1.22
15	BQ	101	PEF	O5-C30	5.96	1.40	1.22
9	A2	101	BCL	O2A-C1	-5.78	1.30	1.46
9	A7	103	BCL	CAA-C2A	5.75	1.64	1.54
9	BV	101	BCL	CAA-CBA	-5.63	1.36	1.52
15	AM	409	PEF	P-O1P	5.62	1.70	1.50
15	AS	101	PEF	P-O1P	5.62	1.70	1.50
15	AM	408	PEF	P-O1P	5.59	1.70	1.50
15	AH	301	PEF	P-O1P	5.59	1.70	1.50
14	B2	102	CRT	C4-C5	5.59	1.58	1.50
15	BQ	101	PEF	P-O1P	5.58	1.70	1.50
9	B2	101	BCL	O2D-CGD	5.58	1.46	1.33
15	BM	407	PEF	P-O1P	5.57	1.70	1.50
15	AM	407	PEF	P-O1P	5.56	1.70	1.50
9	AT	101	BCL	CAA-CBA	-5.50	1.36	1.52
9	AU	102	BCL	O2D-CGD	5.43	1.46	1.33
14	AX	102	CRT	C4-C1	5.43	1.62	1.52
9	BD	102	BCL	O2D-CGD	5.40	1.46	1.33
9	A3	103	BCL	O2A-CGA	5.39	1.49	1.33
9	AF	102	BCL	O2D-CGD	5.38	1.46	1.33
15	AH	301	PEF	O5-C30	5.37	1.40	1.21
15	AM	407	PEF	O5-C30	5.36	1.40	1.21
11	BL	304	UQ8	C41-C42	-5.35	1.36	1.53
9	AV	102	BCL	CAA-CBA	-5.35	1.36	1.52
15	BM	407	PEF	O5-C30	5.32	1.40	1.21
9	A1	102	BCL	CAA-C2A	5.31	1.63	1.54
9	B5	102	BCL	O2D-CGD	5.25	1.46	1.33
15	AS	101	PEF	O2-C10	5.23	1.49	1.34
9	BU	102	BCL	O2D-CGD	5.22	1.46	1.33
11	AL	304	UQ8	C41-C42	-5.22	1.36	1.53
9	AU	102	BCL	CAA-C2A	5.22	1.63	1.54
9	BG	101	BCL	MG-NA	5.22	2.18	2.06
15	AM	409	PEF	O2-C10	5.22	1.49	1.34

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
15	BQ	101	PEF	O2-C10	5.20	1.49	1.34
9	BZ	101	BCL	MG-NA	5.19	2.18	2.06
9	BX	101	BCL	O2D-CGD	5.17	1.45	1.33
9	BM	402	BCL	O2D-CGD	5.15	1.45	1.33
9	A8	101	BCL	C5-C3	5.15	1.61	1.51
9	AM	402	BCL	O2D-CGD	5.14	1.45	1.33
9	AO	102	BCL	O2D-CGD	5.12	1.45	1.33
9	BP	101	BCL	O2D-CGD	5.08	1.45	1.33
9	AS	103	BCL	O2D-CGD	5.08	1.45	1.33
9	AL	301	BCL	O2D-CGD	5.08	1.45	1.33
9	AI	102	BCL	CAA-C2A	5.08	1.63	1.54
7	BC	501	HEM	C3C-C4C	5.07	1.54	1.46
9	BO	102	BCL	O2D-CGD	5.06	1.45	1.33
9	B9	102	BCL	O2D-CGD	5.06	1.45	1.33
9	AQ	102	BCL	O2D-CGD	5.03	1.45	1.33
9	BZ	101	BCL	O2D-CGD	5.02	1.45	1.33
9	AB	101	BCL	O2A-CGA	5.01	1.47	1.33
9	B3	102	BCL	O2D-CGD	5.00	1.45	1.33
9	BT	101	BCL	O2D-CGD	4.98	1.45	1.33
9	A6	101	BCL	O2D-CGD	4.96	1.45	1.33
9	BL	301	BCL	O2D-CGD	4.95	1.45	1.33
9	AA	101	BCL	O2D-CGD	4.93	1.45	1.33
9	AG	101	BCL	CAA-CBA	-4.92	1.38	1.52
9	AU	102	BCL	O2A-CGA	4.91	1.47	1.33
9	BT	101	BCL	MG-NC	4.90	2.17	2.06
9	BN	101	BCL	O2D-CGD	4.90	1.45	1.33
9	A3	104	BCL	O2D-CGD	4.90	1.45	1.33
9	B7	103	BCL	CAA-C2A	4.90	1.63	1.54
14	BP	102	CRT	C4-C1	4.88	1.61	1.52
9	AZ	101	BCL	O2D-CGD	4.87	1.45	1.33
9	BB	101	BCL	O2D-CGD	4.85	1.45	1.33
9	BK	102	BCL	O2D-CGD	4.85	1.45	1.33
9	BE	101	BCL	O2D-CGD	4.85	1.45	1.33
9	AP	101	BCL	CAA-CBA	-4.84	1.38	1.52
9	AJ	101	BCL	O2D-CGD	4.83	1.45	1.33
9	A3	103	BCL	O2D-CGD	4.82	1.45	1.33
9	AI	102	BCL	O2D-CGD	4.82	1.45	1.33
9	BI	102	BCL	O2D-CGD	4.82	1.45	1.33
9	AS	103	BCL	CAA-C2A	4.82	1.62	1.54
9	AV	102	BCL	O2D-CGD	4.81	1.45	1.33
9	A7	103	BCL	O2D-CGD	4.79	1.45	1.33
9	AE	101	BCL	O2A-CGA	4.79	1.47	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	A0	102	BCL	O2D-CGD	4.78	1.45	1.33
9	B7	103	BCL	O2D-CGD	4.78	1.45	1.33
9	BQ	103	BCL	O2D-CGD	4.77	1.45	1.33
9	BS	102	BCL	O2D-CGD	4.77	1.45	1.33
9	A9	102	BCL	CAA-C2A	4.77	1.62	1.54
9	BA	101	BCL	O2D-CGD	4.77	1.44	1.33
9	AN	101	BCL	O2D-CGD	4.75	1.44	1.33
9	AT	101	BCL	O2D-CGD	4.75	1.44	1.33
9	BL	303	BCL	O2D-CGD	4.74	1.44	1.33
9	A1	102	BCL	O2A-CGA	4.74	1.47	1.33
9	AW	101	BCL	CAA-C2A	4.72	1.62	1.54
9	AM	401	BCL	O2D-CGD	4.72	1.44	1.33
9	AK	102	BCL	O2D-CGD	4.71	1.44	1.33
9	AR	101	BCL	O2D-CGD	4.71	1.44	1.33
9	BF	102	BCL	MG-NA	4.71	2.17	2.06
9	BU	102	BCL	MG-NA	-4.70	1.95	2.06
7	AC	501	HEM	C3C-C4C	4.70	1.54	1.46
9	BV	101	BCL	O2D-CGD	4.69	1.44	1.33
9	BW	102	BCL	CAA-C2A	4.68	1.62	1.54
9	BW	102	BCL	O2D-CGD	4.67	1.44	1.33
9	BP	101	BCL	CAA-CBA	-4.67	1.38	1.52
14	AW	102	CRT	C4-C1	4.66	1.61	1.52
9	BD	102	BCL	C3C-C4C	-4.66	1.45	1.51
9	BF	102	BCL	O2A-CGA	4.65	1.46	1.33
9	BE	101	BCL	O2A-C1	-4.64	1.33	1.46
9	BX	101	BCL	O2A-CGA	4.64	1.46	1.33
9	BI	102	BCL	O2A-CGA	4.61	1.46	1.33
9	B6	101	BCL	O2D-CGD	4.61	1.44	1.33
9	BY	102	BCL	O2D-CGD	4.60	1.44	1.33
9	AB	101	BCL	O2D-CGD	4.60	1.44	1.33
7	AC	503	HEM	CAC-C3C	-4.58	1.35	1.47
9	AE	101	BCL	C5-C3	4.58	1.60	1.51
9	AY	102	BCL	O2D-CGD	4.57	1.44	1.33
9	AW	101	BCL	O2D-CGD	4.56	1.44	1.33
9	BU	102	BCL	CAA-C2A	4.54	1.62	1.54
9	BO	102	BCL	MG-NA	4.54	2.17	2.06
9	AO	102	BCL	O2A-CGA	4.53	1.46	1.33
9	A8	101	BCL	O2D-CGD	4.51	1.44	1.33
9	B1	102	BCL	O2D-CGD	4.51	1.44	1.33
9	B7	103	BCL	O2A-CGA	4.51	1.46	1.33
7	BC	503	HEM	CAC-C3C	-4.51	1.35	1.47
9	AI	102	BCL	O2A-CGA	4.51	1.46	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	AG	101	BCL	O2D-CGD	4.48	1.44	1.33
9	B8	101	BCL	MG-NA	4.48	2.16	2.06
9	AD	102	BCL	O2A-CGA	4.47	1.46	1.33
9	BQ	104	BCL	O2D-CGD	4.47	1.44	1.33
9	B9	102	BCL	CAA-C2A	4.46	1.62	1.54
9	B2	101	BCL	CAA-CBA	-4.45	1.39	1.52
7	BC	501	HEM	CAC-C3C	-4.45	1.35	1.47
9	AF	102	BCL	O2A-CGA	4.44	1.46	1.33
9	BY	102	BCL	O2A-CGA	4.42	1.46	1.33
9	AN	101	BCL	CAA-CBA	-4.42	1.39	1.52
7	AC	502	HEM	C3C-C4C	4.41	1.53	1.46
9	AX	101	BCL	CAA-CBA	-4.41	1.39	1.52
7	BC	504	HEM	C1B-NB	-4.39	1.32	1.40
9	B5	102	BCL	MG-NC	4.39	2.16	2.06
9	B9	102	BCL	O2A-CGA	4.39	1.46	1.33
9	BU	102	BCL	O2A-CGA	4.37	1.46	1.33
9	A9	102	BCL	O2D-CGD	4.37	1.44	1.33
9	AR	101	BCL	CAA-CBA	-4.37	1.39	1.52
14	A2	102	CRT	C4-C1	4.37	1.60	1.52
9	A5	102	BCL	CAA-C2A	4.37	1.62	1.54
9	BJ	101	BCL	O2D-CGD	4.36	1.44	1.33
9	AL	303	BCL	O2D-CGD	4.36	1.43	1.33
9	AS	103	BCL	CMA-C3A	4.35	1.62	1.53
9	BT	101	BCL	CAA-CBA	-4.34	1.39	1.52
7	BC	501	HEM	C1B-NB	-4.34	1.32	1.40
9	B8	101	BCL	O2D-CGD	4.33	1.43	1.33
7	AC	504	HEM	C3C-C4C	4.33	1.53	1.46
14	BW	103	CRT	C4-C1	4.32	1.60	1.52
15	AM	409	PEF	O3-C30	4.32	1.45	1.33
14	B7	102	CRT	C4-C1	4.31	1.60	1.52
9	BL	301	BCL	O2A-CGA	4.31	1.45	1.33
15	BQ	101	PEF	O3-C30	4.31	1.45	1.33
9	AS	103	BCL	O2A-CGA	4.30	1.45	1.33
9	BF	102	BCL	CAA-C2A	4.29	1.61	1.54
9	AX	101	BCL	O2D-CGD	4.29	1.43	1.33
15	AS	101	PEF	O3-C30	4.28	1.45	1.33
14	AP	102	CRT	C4-C1	4.28	1.60	1.52
9	BA	101	BCL	O2A-CGA	4.28	1.45	1.33
7	BC	502	HEM	C1B-NB	-4.27	1.32	1.40
9	BJ	101	BCL	CAA-CBA	-4.27	1.40	1.52
14	BP	102	CRT	C4-C5	4.26	1.56	1.50
7	BC	503	HEM	C3B-C2B	4.25	1.45	1.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	B0	102	BCL	O2D-CGD	4.25	1.43	1.33
9	BP	101	BCL	MG-NC	4.25	2.16	2.06
9	A5	102	BCL	O2D-CGD	4.25	1.43	1.33
9	BB	101	BCL	O2A-CGA	4.24	1.45	1.33
9	AE	101	BCL	C2-C3	4.24	1.42	1.33
7	BC	504	HEM	CAC-C3C	-4.22	1.36	1.47
9	A1	102	BCL	O2D-CGD	4.22	1.43	1.33
9	BX	101	BCL	C2-C3	4.21	1.42	1.33
9	BP	101	BCL	O2A-CGA	4.21	1.45	1.33
9	B2	101	BCL	O2A-CGA	4.19	1.45	1.33
9	AE	101	BCL	O2D-CGD	4.19	1.43	1.33
9	BQ	104	BCL	CAA-CBA	-4.19	1.40	1.52
14	B0	101	CRT	C19-C17	4.18	1.45	1.35
9	B5	102	BCL	O2A-CGA	4.15	1.45	1.33
13	AM	405	MQ8	C3-C2	4.15	1.42	1.35
9	BF	102	BCL	C1B-NB	4.14	1.43	1.37
9	AP	101	BCL	O2A-CGA	4.14	1.45	1.33
9	AJ	101	BCL	CAA-CBA	-4.13	1.40	1.52
10	AM	403	BPH	C4D-CHA	-4.12	1.33	1.39
9	B4	101	BCL	O2D-CGD	4.12	1.43	1.33
9	B6	101	BCL	O2A-CGA	4.12	1.45	1.33
9	AK	102	BCL	O2A-CGA	4.12	1.45	1.33
9	AY	102	BCL	O2A-CGA	4.10	1.45	1.33
9	B9	102	BCL	MG-NC	4.10	2.16	2.06
9	B1	102	BCL	O2A-CGA	4.09	1.45	1.33
9	A3	104	BCL	CAA-C2A	4.09	1.61	1.54
9	AE	101	BCL	CAA-C2A	4.09	1.61	1.54
9	BM	401	BCL	O2A-CGA	4.08	1.45	1.33
9	A1	102	BCL	C3B-C4B	4.08	1.49	1.41
9	BG	101	BCL	O2A-CGA	4.08	1.45	1.33
9	AL	303	BCL	O2A-CGA	4.07	1.45	1.33
9	AL	301	BCL	O2A-CGA	4.07	1.45	1.33
9	BK	102	BCL	O2A-CGA	4.05	1.45	1.33
9	AG	101	BCL	O2A-CGA	4.05	1.45	1.33
14	AW	102	CRT	C4-C5	4.05	1.56	1.50
9	BF	102	BCL	O2D-CGD	4.04	1.43	1.33
7	BC	504	HEM	C4C-NC	-4.04	1.32	1.39
14	B7	102	CRT	C4-C5	4.03	1.56	1.50
9	AV	102	BCL	C5-C3	4.03	1.59	1.51
9	B8	101	BCL	O2A-CGA	4.03	1.45	1.33
9	B8	101	BCL	C5-C3	4.02	1.59	1.51
7	AC	501	HEM	C1B-NB	-4.01	1.33	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	BK	102	BCL	MG-NC	4.01	2.15	2.06
14	B2	102	CRT	C19-C17	4.00	1.45	1.35
9	AZ	101	BCL	O2A-CGA	4.00	1.45	1.33
9	B3	102	BCL	O2A-CGA	3.98	1.45	1.33
9	BE	101	BCL	O1A-CGA	3.98	1.34	1.22
9	BM	401	BCL	O2D-CGD	3.98	1.43	1.33
9	AU	102	BCL	C3B-C4B	3.97	1.49	1.41
9	BL	303	BCL	O2A-CGA	3.97	1.44	1.33
9	AA	101	BCL	O2A-CGA	3.95	1.44	1.33
9	A6	101	BCL	MG-NA	3.95	2.15	2.06
9	AX	101	BCL	CMB-C2B	3.93	1.58	1.50
9	B4	101	BCL	O2A-CGA	3.93	1.44	1.33
9	AP	101	BCL	O2D-CGD	3.93	1.42	1.33
7	AC	502	HEM	C4C-NC	-3.90	1.32	1.39
9	BL	301	BCL	MG-NA	3.90	2.15	2.06
10	BM	403	BPH	C4D-CHA	-3.89	1.33	1.39
9	AV	102	BCL	C6-C7	3.88	1.68	1.52
9	BG	101	BCL	O2D-CGD	3.88	1.42	1.33
9	AM	401	BCL	O2A-CGA	3.87	1.44	1.33
7	AC	504	HEM	CAC-C3C	-3.86	1.37	1.47
9	AQ	102	BCL	O2A-CGA	3.86	1.44	1.33
7	AC	504	HEM	C1B-NB	-3.86	1.33	1.40
9	AU	102	BCL	C11-C10	3.86	1.68	1.52
7	BC	503	HEM	C1B-NB	-3.86	1.33	1.40
9	A3	104	BCL	MG-NA	3.84	2.15	2.06
7	BC	502	HEM	C3C-C4C	3.84	1.52	1.46
9	BV	101	BCL	O2A-CGA	3.83	1.44	1.33
9	A2	101	BCL	CAA-CBA	-3.83	1.41	1.52
9	BD	102	BCL	CAA-C2A	3.80	1.61	1.54
9	BD	102	BCL	O2A-CGA	3.80	1.44	1.33
9	B4	101	BCL	C5-C3	3.79	1.59	1.51
9	A0	102	BCL	MG-NC	3.79	2.15	2.06
10	BL	302	BPH	C4D-CHA	-3.78	1.33	1.39
7	AC	501	HEM	CAC-C3C	-3.78	1.37	1.47
9	A5	102	BCL	O2A-CGA	3.78	1.44	1.33
14	AP	102	CRT	C4-C5	3.78	1.55	1.50
9	AJ	101	BCL	O2A-CGA	3.77	1.44	1.33
7	AC	502	HEM	C1B-NB	-3.76	1.33	1.40
9	A9	102	BCL	O2A-CGA	3.75	1.44	1.33
9	AT	101	BCL	C5-C3	3.75	1.59	1.51
9	BQ	103	BCL	MG-NA	-3.75	1.97	2.06
9	AR	101	BCL	O2A-CGA	3.74	1.44	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	AP	101	BCL	CMB-C2B	3.74	1.58	1.50
9	AB	101	BCL	C5-C3	3.73	1.59	1.51
9	BB	101	BCL	C2-C3	3.73	1.41	1.33
9	AM	402	BCL	O2A-CGA	3.72	1.44	1.33
9	B7	103	BCL	MG-NC	3.72	2.15	2.06
9	BZ	101	BCL	O2A-CGA	3.72	1.44	1.33
9	BY	102	BCL	CAA-C2A	3.71	1.60	1.54
9	BA	101	BCL	CAA-C2A	3.71	1.60	1.54
7	AC	504	HEM	C3B-C2B	3.71	1.44	1.37
9	AV	102	BCL	C2-C3	3.71	1.41	1.33
9	B8	101	BCL	CAA-CBA	-3.70	1.41	1.52
9	BX	101	BCL	C5-C3	3.70	1.58	1.51
7	BC	502	HEM	O2A-CGA	-3.69	1.18	1.30
9	BM	402	BCL	O2A-CGA	3.69	1.44	1.33
14	A2	102	CRT	C4-C5	3.69	1.55	1.50
9	AR	101	BCL	MG-NA	3.68	2.15	2.06
9	A3	104	BCL	C5-C3	3.67	1.58	1.51
9	BQ	104	BCL	MG-NC	3.67	2.15	2.06
9	AN	101	BCL	O2A-CGA	3.66	1.44	1.33
9	AL	301	BCL	MG-NA	3.65	2.15	2.06
9	BQ	104	BCL	O2A-CGA	3.64	1.44	1.33
7	AC	504	HEM	C4C-NC	-3.63	1.32	1.39
9	BW	102	BCL	MG-NC	3.63	2.14	2.06
9	A3	103	BCL	CMB-C2B	3.63	1.58	1.50
9	BN	101	BCL	C5-C3	3.62	1.58	1.51
9	BW	102	BCL	O2A-CGA	3.61	1.43	1.33
9	BK	102	BCL	C3B-C4B	3.60	1.48	1.41
9	AY	102	BCL	C3B-C4B	3.60	1.48	1.41
9	BG	101	BCL	CAA-CBA	-3.60	1.42	1.52
9	BJ	101	BCL	C2-C3	3.59	1.41	1.33
9	BO	102	BCL	O2A-CGA	3.59	1.43	1.33
9	AX	101	BCL	C2-C3	3.58	1.41	1.33
9	BN	101	BCL	CAA-CBA	-3.58	1.42	1.52
9	A1	102	BCL	CMB-C2B	3.58	1.58	1.50
9	B4	101	BCL	MG-NC	3.58	2.14	2.06
9	BB	101	BCL	C5-C3	3.56	1.58	1.51
9	AS	103	BCL	C2-C3	3.56	1.41	1.33
9	B3	102	BCL	MG-NC	3.56	2.14	2.06
9	B4	101	BCL	CAA-CBA	-3.56	1.42	1.52
9	BT	101	BCL	MG-NA	-3.56	1.97	2.06
14	BW	103	CRT	C4-C5	3.55	1.55	1.50
9	B0	102	BCL	MG-NC	3.55	2.14	2.06

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	BI	102	BCL	CAA-C2A	3.55	1.60	1.54
9	AE	101	BCL	MG-NC	3.55	2.14	2.06
9	A7	103	BCL	MG-NC	3.55	2.14	2.06
9	BQ	103	BCL	C3B-C2B	-3.54	1.30	1.42
9	BJ	101	BCL	O2A-CGA	3.53	1.43	1.33
9	AR	101	BCL	CMB-C2B	3.53	1.58	1.50
9	AG	101	BCL	C2-C3	3.53	1.41	1.33
9	A5	102	BCL	MG-NC	3.53	2.14	2.06
9	BG	101	BCL	C5-C3	3.53	1.58	1.51
9	A2	101	BCL	C2-C3	3.52	1.41	1.33
9	BU	102	BCL	MG-NC	3.51	2.14	2.06
9	A8	101	BCL	CMB-C2B	3.50	1.58	1.50
9	A7	103	BCL	CMB-C2B	3.49	1.57	1.50
9	BD	102	BCL	O1D-CGD	3.49	1.30	1.21
9	BF	102	BCL	MG-NC	3.49	2.14	2.06
9	A3	103	BCL	C5-C3	3.49	1.58	1.51
9	BK	102	BCL	CAA-C2A	3.49	1.60	1.54
7	BC	502	HEM	CAC-C3C	-3.49	1.38	1.47
9	AN	101	BCL	MG-ND	-3.48	1.98	2.05
9	AF	102	BCL	C2-C3	3.48	1.41	1.33
9	BQ	103	BCL	O2A-CGA	3.48	1.43	1.33
7	BC	502	HEM	C4C-NC	-3.48	1.33	1.39
14	BV	102	CRT	C19-C17	3.45	1.43	1.35
9	A5	102	BCL	CMB-C2B	3.44	1.57	1.50
9	BL	301	BCL	O2D-CED	-3.44	1.37	1.45
9	A9	102	BCL	CMB-C2B	3.44	1.57	1.50
7	AC	501	HEM	C1D-ND	-3.43	1.32	1.38
9	AG	101	BCL	MG-NC	3.43	2.14	2.06
9	BG	101	BCL	C2-C3	3.43	1.41	1.33
9	A8	101	BCL	CAA-C2A	3.42	1.60	1.54
9	AO	102	BCL	CMB-C2B	3.42	1.57	1.50
9	AN	101	BCL	CMB-C2B	3.42	1.57	1.50
9	AQ	102	BCL	MG-NC	3.41	2.14	2.06
7	BC	503	HEM	C1D-C2D	3.41	1.51	1.44
7	BC	504	HEM	C3C-C4C	3.41	1.52	1.46
7	AC	502	HEM	O2A-CGA	-3.41	1.19	1.30
7	AC	503	HEM	C3B-C2B	3.41	1.44	1.37
9	BT	101	BCL	O2A-CGA	3.41	1.43	1.33
9	AX	101	BCL	C5-C3	3.40	1.58	1.51
9	AW	101	BCL	O2A-CGA	3.40	1.43	1.33
9	A3	103	BCL	MG-NC	3.39	2.14	2.06
9	BJ	101	BCL	C5-C3	3.39	1.58	1.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	B6	101	BCL	MG-NC	3.39	2.14	2.06
7	AC	502	HEM	CAC-C3C	-3.39	1.38	1.47
9	AJ	101	BCL	MG-NC	3.38	2.14	2.06
7	AC	501	HEM	C1D-C2D	3.38	1.51	1.44
9	BW	102	BCL	C3C-C4C	3.37	1.55	1.51
9	AL	301	BCL	O2D-CED	-3.37	1.37	1.45
9	AO	102	BCL	C2-C3	3.37	1.40	1.33
9	BI	102	BCL	C2-C3	3.37	1.40	1.33
10	AL	302	BPH	C4D-CHA	-3.37	1.34	1.39
9	A0	102	BCL	CAA-C2A	3.36	1.60	1.54
9	AT	101	BCL	MG-NA	3.35	2.14	2.06
9	AP	101	BCL	MG-NC	3.35	2.14	2.06
9	B9	102	BCL	C2-C3	3.34	1.40	1.33
9	BI	102	BCL	C2C-C3C	-3.34	1.45	1.54
13	BM	405	MQ8	C3-C2	3.34	1.41	1.35
9	BD	102	BCL	C1B-NB	3.33	1.42	1.37
9	BY	102	BCL	C3B-C4B	3.32	1.47	1.41
9	BN	101	BCL	C2-C3	3.32	1.40	1.33
9	AY	102	BCL	CMB-C2B	3.32	1.57	1.50
9	A7	103	BCL	O2A-CGA	3.32	1.43	1.33
9	A1	102	BCL	C2-C3	3.32	1.40	1.33
9	BP	101	BCL	C2-C3	3.31	1.40	1.33
9	BQ	103	BCL	CAA-CBA	-3.31	1.43	1.52
9	A9	102	BCL	MG-NC	3.31	2.14	2.06
7	AC	504	HEM	C1D-C2D	3.31	1.51	1.44
9	AQ	102	BCL	CMB-C2B	3.30	1.57	1.50
9	AJ	101	BCL	C5-C3	3.30	1.58	1.51
9	B2	101	BCL	C2-C3	3.29	1.40	1.33
7	BC	504	HEM	O2A-CGA	-3.28	1.20	1.30
7	AC	501	HEM	C4C-NC	-3.28	1.33	1.39
7	BC	504	HEM	C1D-C2D	3.28	1.51	1.44
9	B0	102	BCL	O2A-CGA	3.28	1.42	1.33
9	B6	101	BCL	C2-C3	3.28	1.40	1.33
9	AA	101	BCL	CAA-C2A	3.28	1.60	1.54
7	AC	503	HEM	O2A-CGA	-3.27	1.20	1.30
9	AA	101	BCL	O2D-CED	-3.27	1.38	1.45
9	A8	101	BCL	O2A-CGA	3.27	1.42	1.33
9	BD	102	BCL	C3B-C4B	3.27	1.47	1.41
9	BI	102	BCL	C5-C3	3.27	1.58	1.51
9	A3	104	BCL	CMB-C2B	3.27	1.57	1.50
7	AC	503	HEM	C3C-C4C	3.26	1.51	1.46
7	BC	504	HEM	CHB-C4A	-3.26	1.32	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	BF	102	BCL	C2C-C3C	-3.26	1.45	1.54
7	BC	502	HEM	C1D-ND	-3.25	1.32	1.38
9	AT	101	BCL	MG-NC	3.25	2.14	2.06
9	AY	102	BCL	C2C-C3C	-3.25	1.45	1.54
9	BK	102	BCL	C2-C3	3.25	1.40	1.33
9	A6	101	BCL	O2A-CGA	3.25	1.42	1.33
9	AB	101	BCL	C2-C3	3.25	1.40	1.33
7	BC	503	HEM	C3C-C4C	3.24	1.51	1.46
7	AC	502	HEM	C1D-ND	-3.24	1.32	1.38
9	B1	102	BCL	CAA-C2A	3.24	1.60	1.54
9	AY	102	BCL	C2-C3	3.23	1.40	1.33
9	AM	402	BCL	CAA-C2A	3.23	1.60	1.54
9	AD	102	BCL	CAA-C2A	3.23	1.60	1.54
14	AX	102	CRT	C19-C17	3.23	1.43	1.35
9	BF	102	BCL	C3C-C4C	3.23	1.55	1.51
15	AM	408	PEF	O2-C10	3.23	1.48	1.33
13	BM	405	MQ8	C11-C3	3.23	1.57	1.51
9	A0	102	BCL	O2A-CGA	3.23	1.42	1.33
9	A8	101	BCL	C6-C7	3.21	1.65	1.52
9	BP	101	BCL	MG-ND	-3.21	1.99	2.05
9	AK	102	BCL	CMB-C2B	3.21	1.57	1.50
14	BB	102	CRT	C25-C23	-3.21	1.39	1.46
14	BB	102	CRT	C30-C28	-3.21	1.39	1.46
14	B0	101	CRT	C30-C28	-3.21	1.39	1.46
9	BQ	103	BCL	CBD-CGD	-3.20	1.42	1.52
14	AB	102	CRT	C16-C17	-3.20	1.39	1.46
9	BU	102	BCL	CMB-C2B	3.19	1.57	1.50
14	A1	103	CRT	C16-C17	-3.19	1.39	1.46
9	AP	101	BCL	C2-C3	3.18	1.40	1.33
7	AC	503	HEM	C1D-C2D	3.18	1.51	1.44
9	BV	101	BCL	C2-C3	3.18	1.40	1.33
9	BE	101	BCL	CMB-C2B	3.18	1.57	1.50
9	BE	101	BCL	MG-NC	3.17	2.13	2.06
9	AV	102	BCL	C6-C5	3.17	1.63	1.52
9	BD	102	BCL	C5-C3	3.17	1.57	1.51
9	BE	101	BCL	C16-C17	3.17	1.65	1.52
14	B2	102	CRT	C22-C23	3.16	1.43	1.35
9	BX	101	BCL	MG-NC	3.16	2.13	2.06
9	BN	101	BCL	O2A-CGA	3.15	1.42	1.33
9	A6	101	BCL	CMB-C2B	3.15	1.57	1.50
13	BM	405	MQ8	C11-C12	3.15	1.55	1.50
13	AM	405	MQ8	C11-C3	3.15	1.57	1.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	B1	102	BCL	C2C-C3C	-3.14	1.46	1.54
9	BQ	104	BCL	CMB-C2B	3.14	1.57	1.50
9	BJ	101	BCL	MG-ND	-3.14	1.99	2.05
7	AC	504	HEM	C1D-ND	-3.14	1.32	1.38
9	AQ	102	BCL	C2-C3	3.14	1.40	1.33
7	BC	501	HEM	O2A-CGA	-3.14	1.20	1.30
14	AB	102	CRT	C25-C23	-3.14	1.39	1.46
9	B7	103	BCL	C2-C3	3.13	1.40	1.33
9	BZ	101	BCL	C5-C3	3.13	1.57	1.51
9	AP	101	BCL	C2C-C3C	-3.12	1.46	1.54
10	BL	302	BPH	C3D-C2D	-3.12	1.34	1.39
10	AL	302	BPH	C4D-ND	-3.11	1.33	1.38
9	BI	102	BCL	CMB-C2B	3.11	1.57	1.50
15	AM	407	PEF	P-O2P	3.10	1.69	1.55
7	AC	502	HEM	C1D-C2D	3.10	1.50	1.44
9	A3	104	BCL	O2A-CGA	3.10	1.42	1.33
15	AM	409	PEF	P-O2P	3.10	1.69	1.55
9	AU	102	BCL	CMB-C2B	3.10	1.57	1.50
9	A2	101	BCL	C3B-C2B	-3.09	1.32	1.42
15	AM	408	PEF	P-O2P	3.09	1.69	1.55
7	BC	503	HEM	O2A-CGA	-3.09	1.20	1.30
9	BJ	101	BCL	CMB-C2B	3.09	1.57	1.50
7	AC	504	HEM	CHB-C4A	-3.08	1.32	1.39
9	BD	102	BCL	CMA-C3A	3.08	1.59	1.53
15	BQ	101	PEF	P-O2P	3.08	1.69	1.55
9	BZ	101	BCL	CAA-CBA	-3.08	1.43	1.52
9	AY	102	BCL	O2A-C1	3.08	1.54	1.46
15	AH	301	PEF	P-O2P	3.08	1.69	1.55
15	AS	101	PEF	P-O2P	3.08	1.69	1.55
9	BQ	103	BCL	C2-C3	3.08	1.40	1.33
9	BK	102	BCL	CMB-C2B	3.07	1.57	1.50
9	A2	101	BCL	C2C-C3C	-3.07	1.46	1.54
14	BV	102	CRT	C22-C23	3.07	1.42	1.35
9	AK	102	BCL	CAA-C2A	3.07	1.59	1.54
9	AX	101	BCL	O2A-CGA	3.07	1.42	1.33
15	BM	407	PEF	P-O2P	3.06	1.69	1.55
9	AA	101	BCL	C2-C3	3.06	1.40	1.33
9	B3	102	BCL	CAA-C2A	3.06	1.59	1.54
7	AC	503	HEM	C1B-NB	-3.06	1.34	1.40
9	AJ	101	BCL	O2D-CED	-3.05	1.38	1.45
9	BO	102	BCL	CAA-C2A	3.05	1.59	1.54
9	BP	101	BCL	MG-NA	3.05	2.13	2.06

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
14	B2	102	CRT	C14-C12	3.05	1.42	1.35
9	BL	301	BCL	CAA-C2A	3.04	1.59	1.54
10	AL	302	BPH	C3B-C2B	3.04	1.44	1.39
9	B5	102	BCL	MG-NA	3.04	2.13	2.06
9	AU	102	BCL	CAA-CBA	-3.04	1.43	1.52
9	A0	102	BCL	CMB-C2B	3.03	1.57	1.50
9	BI	102	BCL	MG-NA	-3.03	1.99	2.06
9	BE	101	BCL	C3B-C2B	-3.03	1.32	1.42
9	A3	103	BCL	MG-NA	3.03	2.13	2.06
9	BQ	103	BCL	C3B-C4B	3.03	1.47	1.41
9	AS	103	BCL	MG-NC	3.03	2.13	2.06
9	BG	101	BCL	CMB-C2B	3.02	1.57	1.50
10	BL	302	BPH	C4D-ND	-3.02	1.34	1.38
9	B1	102	BCL	CHB-C4A	3.01	1.45	1.37
9	A6	101	BCL	MG-NC	3.00	2.13	2.06
9	A9	102	BCL	C2-C3	3.00	1.40	1.33
9	AG	101	BCL	C5-C3	3.00	1.57	1.51
9	AI	102	BCL	MG-NC	3.00	2.13	2.06
9	AA	101	BCL	MG-NC	3.00	2.13	2.06
9	BX	101	BCL	CHB-C4A	2.99	1.45	1.37
9	AR	101	BCL	C2C-C3C	-2.99	1.46	1.54
7	BC	504	HEM	CHD-C4C	-2.99	1.32	1.38
9	AU	102	BCL	MG-NC	2.99	2.13	2.06
9	AV	102	BCL	O2A-CGA	2.98	1.42	1.33
9	BM	402	BCL	MG-NA	2.98	2.13	2.06
9	AL	301	BCL	CAA-C2A	2.98	1.59	1.54
9	AG	101	BCL	C2C-C3C	-2.98	1.46	1.54
9	AS	103	BCL	C1B-NB	2.98	1.41	1.37
14	B7	102	CRT	C14-C12	2.98	1.42	1.35
9	BQ	104	BCL	C2-C3	2.97	1.39	1.33
9	BB	101	BCL	C3B-C4B	2.97	1.47	1.41
9	BA	101	BCL	C2-C3	2.97	1.39	1.33
9	A8	101	BCL	C7-C8	2.97	1.67	1.52
9	B5	102	BCL	CAA-CBA	-2.97	1.44	1.52
14	B0	101	CRT	C14-C12	2.97	1.42	1.35
9	A2	101	BCL	O1A-CGA	2.96	1.31	1.22
9	BY	102	BCL	C5-C3	2.96	1.57	1.51
9	AT	101	BCL	C4-C3	2.96	1.57	1.50
9	BX	101	BCL	CMB-C2B	2.96	1.56	1.50
7	BC	502	HEM	C3B-C2B	2.96	1.43	1.37
9	AV	102	BCL	MG-NC	2.96	2.13	2.06
9	BF	102	BCL	CMB-C2B	2.96	1.56	1.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	BC	503	HEM	CHB-C4A	-2.96	1.32	1.39
9	BK	102	BCL	C5-C3	2.95	1.57	1.51
14	A1	103	CRT	C4-C1	2.95	1.57	1.52
9	BS	102	BCL	CMB-C2B	2.95	1.56	1.50
9	A6	101	BCL	CAA-C2A	2.95	1.59	1.54
9	AU	102	BCL	CMD-C2D	2.95	1.56	1.50
9	BJ	101	BCL	C2C-C3C	-2.95	1.46	1.54
9	B2	101	BCL	C2C-C3C	-2.95	1.46	1.54
9	BF	102	BCL	C2-C3	2.94	1.39	1.33
9	A9	102	BCL	MG-NA	2.94	2.13	2.06
14	AX	102	CRT	C14-C12	2.93	1.42	1.35
9	AD	102	BCL	C3B-C4B	2.93	1.46	1.41
9	BX	101	BCL	C2C-C3C	-2.93	1.46	1.54
9	AF	102	BCL	C5-C3	2.93	1.57	1.51
9	B4	101	BCL	C2-C3	2.93	1.39	1.33
9	AB	101	BCL	CAA-CBA	-2.93	1.44	1.52
9	B3	102	BCL	C2C-C3C	-2.93	1.46	1.54
9	BL	303	BCL	C3B-C4B	2.93	1.46	1.41
9	BD	102	BCL	CBD-CGD	2.92	1.61	1.52
9	AJ	101	BCL	C2C-C3C	-2.92	1.46	1.54
9	BS	102	BCL	O2A-CGA	2.92	1.41	1.33
9	A0	102	BCL	C2C-C3C	-2.92	1.46	1.54
9	AV	102	BCL	C4-C3	2.92	1.57	1.50
9	AN	101	BCL	MG-NC	2.91	2.13	2.06
9	AX	101	BCL	MG-NC	2.91	2.13	2.06
9	B2	101	BCL	MG-NC	2.91	2.13	2.06
9	AP	101	BCL	C5-C3	2.91	1.57	1.51
9	A1	102	BCL	MG-NC	2.91	2.13	2.06
9	AX	101	BCL	CAA-C2A	2.90	1.59	1.54
9	AO	102	BCL	C2C-C3C	-2.90	1.46	1.54
9	AW	101	BCL	C3A-C2A	-2.90	1.46	1.54
9	AJ	101	BCL	C2-C3	2.90	1.39	1.33
9	BT	101	BCL	C2A-C1A	2.90	1.58	1.52
9	BO	102	BCL	C2-C3	2.89	1.39	1.33
9	BO	102	BCL	C2C-C3C	-2.89	1.46	1.54
9	BX	101	BCL	CAA-CBA	-2.89	1.44	1.52
9	AD	102	BCL	MG-NC	2.89	2.13	2.06
9	AF	102	BCL	CAA-CBA	-2.89	1.44	1.52
7	AC	503	HEM	C2A-C3A	-2.88	1.31	1.38
9	AK	102	BCL	C3B-C4B	2.88	1.46	1.41
9	B0	102	BCL	O2D-CED	-2.88	1.38	1.45
9	BU	102	BCL	C2-C3	2.88	1.39	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	BC	502	HEM	C1D-C2D	2.87	1.50	1.44
9	BY	102	BCL	MG-NC	2.87	2.13	2.06
9	BX	101	BCL	C1D-ND	-2.87	1.34	1.37
9	AQ	102	BCL	C3B-C4B	2.87	1.46	1.41
9	A2	101	BCL	O2D-CED	-2.87	1.38	1.45
14	B2	102	CRT	C6-C5	2.87	1.40	1.32
9	AD	102	BCL	C2-C3	2.86	1.39	1.33
9	AZ	101	BCL	C3B-C4B	2.86	1.46	1.41
9	AT	101	BCL	C2C-C3C	-2.86	1.46	1.54
9	BE	101	BCL	CBA-CGA	-2.86	1.42	1.50
9	BN	101	BCL	MG-ND	-2.86	2.00	2.05
9	A8	101	BCL	C3A-C2A	-2.86	1.46	1.54
9	BG	101	BCL	C3B-C4B	2.86	1.46	1.41
7	BC	501	HEM	C3B-C2B	2.85	1.43	1.37
9	AF	102	BCL	MG-NC	2.85	2.13	2.06
9	BS	102	BCL	MG-NC	2.85	2.13	2.06
9	AE	101	BCL	C2C-C3C	-2.85	1.46	1.54
9	A3	104	BCL	MG-NC	2.84	2.13	2.06
9	B1	102	BCL	MG-NC	2.84	2.13	2.06
9	AS	103	BCL	CHB-C1B	2.84	1.45	1.39
9	BS	102	BCL	C2-C3	2.84	1.39	1.33
14	AM	406	CRT	C37-C38	2.83	1.57	1.52
9	AU	102	BCL	C3B-C2B	-2.83	1.32	1.42
7	BC	501	HEM	C1D-C2D	2.83	1.50	1.44
9	B9	102	BCL	CMB-C2B	2.83	1.56	1.50
9	A2	101	BCL	C3B-C4B	2.83	1.46	1.41
9	AY	102	BCL	CAA-CBA	-2.83	1.44	1.52
9	AG	101	BCL	O1A-CGA	2.83	1.30	1.22
9	A7	103	BCL	CHB-C4A	2.83	1.44	1.37
9	AO	102	BCL	C3A-C2A	-2.82	1.46	1.54
9	BE	101	BCL	C19-C18	2.82	1.66	1.51
9	B7	103	BCL	CMB-C2B	2.82	1.56	1.50
9	BM	402	BCL	CAA-C2A	2.82	1.59	1.54
9	BN	101	BCL	C3B-C4B	2.82	1.46	1.41
7	AC	502	HEM	CHA-C1A	-2.82	1.33	1.39
9	BG	101	BCL	O2D-CED	-2.82	1.39	1.45
9	AB	101	BCL	O2D-CED	-2.82	1.39	1.45
9	BQ	104	BCL	O2D-CED	-2.82	1.39	1.45
9	A7	103	BCL	C2C-C3C	-2.82	1.46	1.54
9	AT	101	BCL	O1A-CGA	2.81	1.30	1.22
9	A9	102	BCL	C2C-C3C	-2.81	1.46	1.54
7	BC	504	HEM	C1D-ND	-2.81	1.33	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	A5	102	BCL	C2-C3	2.81	1.39	1.33
9	A3	103	BCL	C2C-C3C	-2.81	1.46	1.54
9	AP	101	BCL	MG-ND	-2.81	2.00	2.05
9	AD	102	BCL	C5-C3	2.81	1.57	1.51
9	AR	101	BCL	C2-C3	2.81	1.39	1.33
9	B1	102	BCL	CMB-C2B	2.81	1.56	1.50
14	A0	101	CRT	C4-C1	2.80	1.57	1.52
9	A1	102	BCL	CHB-C4A	2.80	1.44	1.37
10	AL	302	BPH	C3D-C2D	-2.80	1.34	1.39
9	AK	102	BCL	MG-NC	2.80	2.12	2.06
9	A2	101	BCL	O2D-CGD	2.80	1.40	1.33
9	AI	102	BCL	C5-C3	2.80	1.57	1.51
9	AZ	101	BCL	O1A-CGA	2.79	1.30	1.22
9	BL	301	BCL	C5-C3	2.79	1.57	1.51
7	BC	501	HEM	CHB-C4A	-2.79	1.33	1.39
9	BA	101	BCL	C5-C3	2.79	1.57	1.51
9	AW	101	BCL	CMB-C2B	2.79	1.56	1.50
9	B3	102	BCL	CMB-C2B	2.78	1.56	1.50
9	BD	102	BCL	CMB-C2B	2.78	1.56	1.50
9	B8	101	BCL	MG-NC	2.78	2.12	2.06
9	BO	102	BCL	CMB-C2B	2.78	1.56	1.50
9	AE	101	BCL	O2D-CED	-2.77	1.39	1.45
9	A6	101	BCL	C2C-C3C	-2.77	1.47	1.54
9	B3	102	BCL	C5-C3	2.77	1.57	1.51
9	BU	102	BCL	C2C-C3C	-2.77	1.47	1.54
9	BO	102	BCL	C3B-C4B	2.77	1.46	1.41
9	AT	101	BCL	C2-C3	2.76	1.39	1.33
10	BM	403	BPH	C3D-C2D	-2.76	1.34	1.39
9	AM	401	BCL	CAA-C2A	2.76	1.59	1.54
9	AI	102	BCL	C2-C3	2.76	1.39	1.33
9	BP	101	BCL	CMB-C2B	2.76	1.56	1.50
9	A0	102	BCL	C3B-C4B	2.75	1.46	1.41
9	AT	101	BCL	MG-ND	-2.75	2.00	2.05
9	BD	102	BCL	C2C-C3C	-2.75	1.47	1.54
9	B7	103	BCL	C5-C3	2.75	1.57	1.51
7	BC	503	HEM	C4A-NA	-2.75	1.34	1.39
9	AZ	101	BCL	CAA-C2A	2.75	1.59	1.54
9	AA	101	BCL	CMB-C2B	2.75	1.56	1.50
14	B0	101	CRT	C9-C7	2.74	1.42	1.35
9	B0	102	BCL	CAA-CBA	-2.74	1.44	1.52
14	B2	102	CRT	C15-C16	2.74	1.41	1.34
14	B1	103	CRT	C25-C23	-2.74	1.40	1.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
14	BB	102	CRT	C16-C17	-2.73	1.40	1.46
9	AI	102	BCL	CMB-C2B	2.73	1.56	1.50
9	A3	104	BCL	C2C-C3C	-2.73	1.47	1.54
7	AC	501	HEM	O2A-CGA	-2.73	1.21	1.30
9	BZ	101	BCL	C2-C3	2.73	1.39	1.33
9	BI	102	BCL	MG-NC	2.73	2.12	2.06
9	AZ	101	BCL	CMB-C2B	2.73	1.56	1.50
9	B8	101	BCL	CMB-C2B	2.73	1.56	1.50
14	BV	102	CRT	C15-C16	2.73	1.41	1.34
9	BV	101	BCL	C5-C3	2.73	1.56	1.51
9	BT	101	BCL	C2-C3	2.72	1.39	1.33
9	BA	101	BCL	MG-NC	2.72	2.12	2.06
9	B5	102	BCL	CMB-C2B	2.72	1.56	1.50
7	BC	504	HEM	C3B-C2B	2.72	1.42	1.37
9	AP	101	BCL	C1B-C2B	2.72	1.49	1.43
9	AL	301	BCL	C5-C3	2.71	1.56	1.51
9	B2	101	BCL	O1A-CGA	2.71	1.30	1.22
9	B9	102	BCL	C2C-C3C	-2.71	1.47	1.54
9	AQ	102	BCL	CAA-CBA	-2.70	1.44	1.52
9	BA	101	BCL	CMB-C2B	2.70	1.56	1.50
9	BB	101	BCL	CAA-CBA	-2.70	1.44	1.52
9	AF	102	BCL	CMB-C2B	2.70	1.56	1.50
9	BP	101	BCL	C2C-C3C	-2.70	1.47	1.54
10	AM	403	BPH	C3D-C2D	-2.70	1.35	1.39
9	AF	102	BCL	O2D-CED	-2.70	1.39	1.45
9	AB	101	BCL	C2C-C3C	-2.69	1.47	1.54
9	AI	102	BCL	C2C-C3C	-2.69	1.47	1.54
10	BM	403	BPH	C4D-ND	-2.69	1.34	1.38
9	AW	101	BCL	C2C-C3C	-2.69	1.47	1.54
9	B6	101	BCL	C2C-C3C	-2.69	1.47	1.54
9	B0	102	BCL	MG-ND	-2.68	2.00	2.05
14	AB	102	CRT	C30-C28	-2.68	1.40	1.46
9	AE	101	BCL	CMB-C2B	2.68	1.56	1.50
9	BV	101	BCL	CMB-C2B	2.68	1.56	1.50
9	AZ	101	BCL	MG-NC	2.68	2.12	2.06
9	A8	101	BCL	C2C-C3C	-2.68	1.47	1.54
7	BC	504	HEM	C1A-C2A	-2.67	1.39	1.44
9	BZ	101	BCL	CMB-C2B	2.67	1.56	1.50
9	B2	101	BCL	C5-C3	2.67	1.56	1.51
9	BS	102	BCL	CAA-C2A	2.67	1.58	1.54
9	AM	401	BCL	C2C-C3C	-2.67	1.47	1.54
10	BL	302	BPH	CHA-CBD	2.67	1.54	1.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	AC	502	HEM	CMB-C2B	-2.66	1.45	1.50
9	AN	101	BCL	C3B-C4B	2.66	1.46	1.41
9	AZ	101	BCL	CMC-C2C	2.66	1.58	1.53
9	BQ	103	BCL	C4B-NB	-2.66	1.34	1.37
9	AA	101	BCL	C2C-C3C	-2.66	1.47	1.54
14	BV	102	CRT	C14-C12	2.66	1.42	1.35
9	BV	101	BCL	MG-NA	2.66	2.12	2.06
7	BC	501	HEM	C1D-ND	-2.66	1.33	1.38
9	B0	102	BCL	CMB-C2B	2.66	1.56	1.50
9	AL	301	BCL	C2-C3	2.66	1.39	1.33
10	AM	403	BPH	C4D-ND	-2.65	1.34	1.38
7	BC	503	HEM	C1A-C2A	-2.65	1.39	1.44
9	BE	101	BCL	MG-ND	-2.65	2.00	2.05
9	AL	301	BCL	C2C-C3C	-2.65	1.47	1.54
9	AW	101	BCL	C2-C3	2.65	1.39	1.33
9	B0	102	BCL	C5-C3	2.65	1.56	1.51
9	AL	303	BCL	C2C-C3C	-2.65	1.47	1.54
9	AL	303	BCL	CAA-C2A	2.65	1.58	1.54
14	BG	102	CRT	C30-C28	-2.65	1.40	1.46
9	BQ	104	BCL	O1A-CGA	2.65	1.30	1.22
9	BD	102	BCL	C2-C3	2.65	1.39	1.33
9	BX	101	BCL	MG-ND	-2.64	2.00	2.05
9	B7	103	BCL	C3B-C4B	2.64	1.46	1.41
9	BE	101	BCL	O2D-CED	-2.64	1.39	1.45
9	AB	101	BCL	CMB-C2B	2.63	1.56	1.50
9	BY	102	BCL	C2C-C3C	-2.63	1.47	1.54
9	BJ	101	BCL	C3B-C2B	-2.63	1.33	1.42
9	BE	101	BCL	C2C-C3C	-2.62	1.47	1.54
14	B7	102	CRT	C15-C16	2.62	1.41	1.34
14	BU	103	CRT	C8-C7	2.62	1.56	1.50
7	AC	502	HEM	C3B-C2B	2.62	1.42	1.37
9	B4	101	BCL	MG-ND	-2.62	2.00	2.05
9	AT	101	BCL	O2A-CGA	2.62	1.41	1.33
9	AY	102	BCL	C4-C3	2.61	1.57	1.50
9	BJ	101	BCL	C3B-C4B	2.61	1.46	1.41
9	AD	102	BCL	C2C-C3C	-2.61	1.47	1.54
9	BF	102	BCL	C5-C3	2.61	1.56	1.51
9	AJ	101	BCL	CMB-C2B	2.61	1.56	1.50
9	AZ	101	BCL	MG-ND	-2.61	2.00	2.05
14	A2	102	CRT	C14-C12	2.61	1.41	1.35
9	B4	101	BCL	CMB-C2B	2.60	1.56	1.50
9	B4	101	BCL	C2C-C3C	-2.60	1.47	1.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	BY	102	BCL	C2-C3	2.60	1.39	1.33
7	AC	504	HEM	C1A-C2A	-2.60	1.39	1.44
9	B4	101	BCL	CAA-C2A	2.60	1.58	1.54
9	B1	102	BCL	CMA-C3A	2.60	1.58	1.53
7	AC	501	HEM	CHB-C4A	-2.60	1.33	1.39
9	A5	102	BCL	C2C-C3C	-2.60	1.47	1.54
9	AO	102	BCL	C3B-C2B	-2.60	1.33	1.42
9	BX	101	BCL	O1A-CGA	2.60	1.30	1.22
9	B5	102	BCL	C3B-C4B	2.60	1.46	1.41
14	BP	102	CRT	C14-C12	2.60	1.41	1.35
9	BQ	104	BCL	C3B-C4B	2.60	1.46	1.41
9	BL	301	BCL	C2-C3	2.60	1.39	1.33
15	AM	407	PEF	O3-C30	2.59	1.45	1.33
9	BG	101	BCL	C1B-NB	2.59	1.41	1.37
9	A7	103	BCL	CAA-CBA	-2.59	1.45	1.52
9	BI	102	BCL	C3B-C2B	-2.59	1.33	1.42
9	AX	101	BCL	CHB-C4A	2.59	1.44	1.37
9	AF	102	BCL	C4-C3	2.58	1.57	1.50
9	BM	401	BCL	CAA-C2A	2.58	1.58	1.54
14	BN	102	CRT	C4-C5	2.58	1.54	1.50
9	B0	102	BCL	C3B-C4B	2.58	1.46	1.41
9	AR	101	BCL	O2D-CED	-2.58	1.39	1.45
9	B5	102	BCL	C2C-C3C	-2.58	1.47	1.54
7	AC	504	HEM	O2A-CGA	-2.58	1.22	1.30
15	AH	301	PEF	O3-C30	2.58	1.45	1.33
9	AB	101	BCL	MG-NA	2.58	2.12	2.06
9	BV	101	BCL	O2D-CED	-2.57	1.39	1.45
9	AW	101	BCL	CMA-C3A	2.57	1.58	1.53
9	AX	101	BCL	C2C-C3C	-2.57	1.47	1.54
9	AB	101	BCL	MG-NC	2.57	2.12	2.06
9	B4	101	BCL	O2D-CED	-2.57	1.39	1.45
9	A3	104	BCL	C3B-C4B	2.56	1.46	1.41
9	A6	101	BCL	C3B-C4B	2.56	1.46	1.41
9	BV	101	BCL	MG-ND	-2.56	2.00	2.05
14	A2	102	CRT	C19-C17	2.56	1.41	1.35
7	AC	501	HEM	C3B-C2B	2.56	1.42	1.37
9	A5	102	BCL	C4-C3	2.56	1.57	1.50
9	B6	101	BCL	C3B-C4B	2.56	1.46	1.41
9	A5	102	BCL	C3B-C4B	2.56	1.46	1.41
9	AN	101	BCL	C2C-C3C	-2.56	1.47	1.54
9	BB	101	BCL	MG-ND	-2.56	2.00	2.05
9	A9	102	BCL	O2A-C1	2.55	1.53	1.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	BN	101	BCL	CMC-C2C	2.55	1.58	1.53
9	AU	102	BCL	C2C-C3C	-2.55	1.47	1.54
9	AM	401	BCL	C3B-C2B	-2.55	1.33	1.42
15	BM	407	PEF	O3-C30	2.55	1.45	1.33
9	BL	303	BCL	C2C-C3C	-2.55	1.47	1.54
9	AF	102	BCL	C2C-C3C	-2.55	1.47	1.54
9	AT	101	BCL	CHB-C1B	2.55	1.45	1.39
9	AG	101	BCL	CMB-C2B	2.54	1.56	1.50
9	AX	101	BCL	C3B-C4B	2.54	1.46	1.41
9	B8	101	BCL	C2C-C3C	-2.54	1.47	1.54
9	BB	101	BCL	C4D-CHA	2.54	1.47	1.38
9	BL	301	BCL	C2C-C3C	-2.54	1.47	1.54
9	B7	103	BCL	C2C-C3C	-2.54	1.47	1.54
9	A6	101	BCL	C2-C3	2.54	1.38	1.33
9	B0	102	BCL	C2C-C3C	-2.54	1.47	1.54
9	A8	101	BCL	O2D-CED	-2.54	1.39	1.45
9	AQ	102	BCL	C4-C3	2.54	1.56	1.50
9	BE	101	BCL	C16-C15	2.54	1.62	1.52
14	B1	103	CRT	C16-C17	-2.54	1.40	1.46
9	AL	303	BCL	C3B-C4B	2.53	1.46	1.41
7	BC	504	HEM	C2A-C3A	-2.53	1.32	1.38
9	BX	101	BCL	CMA-C3A	2.53	1.58	1.53
14	BO	103	CRT	C16-C17	-2.53	1.40	1.46
9	A3	104	BCL	O2D-CED	-2.53	1.39	1.45
7	BC	501	HEM	CHA-C4D	-2.53	1.33	1.38
9	AW	101	BCL	C3B-C2B	-2.52	1.33	1.42
9	AT	101	BCL	C2A-C1A	2.52	1.57	1.52
14	BO	103	CRT	C25-C23	-2.52	1.40	1.46
9	B4	101	BCL	C2A-C1A	2.52	1.57	1.52
9	BE	101	BCL	C1-C2	-2.52	1.42	1.49
14	B7	102	CRT	C19-C17	2.52	1.41	1.35
9	AA	101	BCL	C4-C3	2.51	1.56	1.50
9	BJ	101	BCL	C4-C3	2.51	1.56	1.50
9	AT	101	BCL	C3B-C2B	-2.51	1.34	1.42
9	AS	103	BCL	C2C-C3C	-2.51	1.47	1.54
9	A5	102	BCL	CAA-CBA	-2.51	1.45	1.52
7	BC	504	HEM	C4A-NA	-2.51	1.34	1.39
9	AG	101	BCL	O2D-CED	-2.51	1.39	1.45
9	A3	103	BCL	CHB-C4A	2.50	1.44	1.37
9	BY	102	BCL	CMB-C2B	2.50	1.55	1.50
9	AK	102	BCL	CMA-C3A	2.50	1.58	1.53
9	AK	102	BCL	C2C-C3C	-2.50	1.47	1.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
14	BU	103	CRT	C16-C17	-2.50	1.40	1.46
14	B0	101	CRT	C15-C16	2.50	1.41	1.34
9	B1	102	BCL	C3B-C2B	-2.50	1.34	1.42
9	A9	102	BCL	CAA-CBA	-2.50	1.45	1.52
9	AV	102	BCL	C2C-C3C	-2.49	1.47	1.54
9	B0	102	BCL	C2-C3	2.49	1.38	1.33
9	AW	101	BCL	MG-NC	2.49	2.12	2.06
9	A0	102	BCL	C2-C3	2.49	1.38	1.33
9	A1	102	BCL	C2C-C3C	-2.49	1.47	1.54
9	A2	101	BCL	O2A-CGA	2.49	1.40	1.33
9	BM	401	BCL	C3B-C2B	-2.49	1.34	1.42
7	BC	502	HEM	CHA-C1A	-2.49	1.33	1.39
9	AW	101	BCL	C4-C3	2.48	1.56	1.50
9	AM	402	BCL	CMB-C2B	2.48	1.55	1.50
9	A1	102	BCL	CMA-C3A	2.48	1.58	1.53
9	BJ	101	BCL	O2D-CED	-2.48	1.39	1.45
9	B4	101	BCL	MG-NA	2.48	2.12	2.06
9	BQ	103	BCL	CHB-C1B	-2.48	1.33	1.39
14	AB	102	CRT	C11-C12	-2.48	1.40	1.46
9	AQ	102	BCL	C3B-C2B	-2.48	1.34	1.42
9	B2	101	BCL	CMB-C2B	2.48	1.55	1.50
9	BA	101	BCL	CMA-C3A	2.48	1.58	1.53
14	AN	102	CRT	C16-C17	-2.48	1.40	1.46
9	BQ	103	BCL	CMB-C2B	2.48	1.55	1.50
9	A8	101	BCL	C10-C8	2.48	1.64	1.52
9	B4	101	BCL	C3B-C4B	2.47	1.46	1.41
9	BA	101	BCL	C2C-C3C	-2.47	1.47	1.54
9	B8	101	BCL	C3B-C4B	2.47	1.46	1.41
9	A1	102	BCL	C4-C3	2.47	1.56	1.50
9	A3	103	BCL	C3B-C4B	2.47	1.46	1.41
9	A7	103	BCL	C3A-C2A	-2.47	1.47	1.54
9	BM	402	BCL	C2C-C3C	-2.47	1.47	1.54
9	AU	102	BCL	CMC-C2C	2.46	1.58	1.53
9	BE	101	BCL	C3B-C4B	2.46	1.46	1.41
7	BC	501	HEM	C3B-C4B	-2.46	1.40	1.44
9	B2	101	BCL	MG-ND	-2.46	2.00	2.05
9	B6	101	BCL	CMB-C2B	2.46	1.55	1.50
9	BX	101	BCL	C3B-C4B	2.46	1.46	1.41
14	BV	102	CRT	C37-C36	2.46	1.53	1.50
9	BN	101	BCL	C4-C3	2.45	1.56	1.50
9	BF	102	BCL	C3B-C4B	2.45	1.46	1.41
9	AF	102	BCL	CHB-C4A	2.45	1.43	1.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	A8	101	BCL	MG-NC	2.45	2.12	2.06
9	AT	101	BCL	C3B-C4B	2.45	1.45	1.41
14	AN	102	CRT	C25-C23	-2.45	1.40	1.46
9	AR	101	BCL	CHB-C4A	2.45	1.43	1.37
9	A1	102	BCL	CHB-C1B	2.45	1.44	1.39
14	AP	102	CRT	C19-C17	2.45	1.41	1.35
14	B7	102	CRT	C22-C23	2.44	1.41	1.35
9	AE	101	BCL	CHB-C4A	2.44	1.43	1.37
9	BK	102	BCL	C2C-C3C	-2.44	1.47	1.54
9	B1	102	BCL	MG-ND	-2.44	2.00	2.05
9	AM	402	BCL	MG-NC	2.44	2.12	2.06
9	AA	101	BCL	C5-C3	2.44	1.56	1.51
9	B9	102	BCL	CHB-C4A	2.44	1.43	1.37
7	AC	501	HEM	O2D-CGD	-2.44	1.22	1.30
14	BU	103	CRT	C25-C23	-2.44	1.40	1.46
14	B1	103	CRT	C30-C28	-2.44	1.40	1.46
9	AP	101	BCL	C3B-C4B	2.43	1.45	1.41
7	AC	502	HEM	FE-NB	2.43	2.02	1.94
14	BP	102	CRT	C19-C17	2.43	1.41	1.35
9	BL	303	BCL	C4-C3	2.43	1.56	1.50
9	AQ	102	BCL	C2C-C3C	-2.43	1.47	1.54
9	BO	102	BCL	CMA-C3A	2.43	1.58	1.53
9	AR	101	BCL	MG-NC	2.43	2.12	2.06
9	A9	102	BCL	O2D-CED	-2.43	1.39	1.45
9	AG	101	BCL	CHB-C4A	2.43	1.43	1.37
9	BQ	103	BCL	CMD-C2D	2.43	1.55	1.50
9	A1	102	BCL	CAA-CBA	-2.42	1.45	1.52
9	BW	102	BCL	CMC-C2C	2.42	1.58	1.53
9	A0	102	BCL	O2D-CED	-2.42	1.39	1.45
9	A0	102	BCL	CHB-C4A	2.42	1.43	1.37
9	BD	102	BCL	CHB-C1B	2.42	1.44	1.39
9	BT	101	BCL	CMB-C2B	2.42	1.55	1.50
14	BF	103	CRT	C25-C23	-2.42	1.40	1.46
14	BG	102	CRT	C25-C23	-2.42	1.40	1.46
9	AA	101	BCL	C3B-C4B	2.42	1.45	1.41
9	BU	102	BCL	C3B-C4B	2.41	1.45	1.41
9	BS	102	BCL	C2C-C3C	-2.41	1.47	1.54
9	AN	101	BCL	C5-C3	2.41	1.56	1.51
9	AX	101	BCL	O1A-CGA	2.41	1.29	1.22
9	BP	101	BCL	C3B-C4B	2.41	1.45	1.41
9	AQ	102	BCL	CAA-C2A	2.41	1.58	1.54
9	BN	101	BCL	CMB-C2B	2.41	1.55	1.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
14	A1	103	CRT	C30-C28	-2.41	1.40	1.46
9	BP	101	BCL	CHB-C1B	2.41	1.44	1.39
9	AB	101	BCL	O1A-CGA	2.40	1.29	1.22
13	BM	405	MQ8	C10-C5	2.40	1.44	1.40
9	BW	102	BCL	C5-C3	2.40	1.56	1.51
9	BQ	103	BCL	C2C-C3C	-2.40	1.48	1.54
9	AN	101	BCL	C2-C3	2.40	1.38	1.33
14	B0	101	CRT	C4-C5	2.40	1.53	1.50
9	BS	102	BCL	CAA-CBA	-2.40	1.45	1.52
14	A1	103	CRT	C6-C5	2.40	1.38	1.32
7	BC	502	HEM	C2A-C3A	-2.40	1.32	1.38
9	AK	102	BCL	C2-C3	2.39	1.38	1.33
9	AZ	101	BCL	CHD-C1D	-2.39	1.33	1.38
14	AP	102	CRT	C14-C12	2.39	1.41	1.35
9	AK	102	BCL	C3B-C2B	-2.39	1.34	1.42
14	A1	103	CRT	C25-C23	-2.39	1.40	1.46
9	BN	101	BCL	C6-C5	2.39	1.60	1.52
14	AT	102	CRT	C19-C17	2.39	1.41	1.35
14	BP	102	CRT	C22-C23	2.39	1.41	1.35
9	AV	102	BCL	CMB-C2B	2.38	1.55	1.50
9	BZ	101	BCL	CHB-C4A	2.38	1.43	1.37
9	A8	101	BCL	C4-C3	2.38	1.56	1.50
7	AC	503	HEM	C4D-ND	-2.38	1.36	1.40
14	AR	102	CRT	C16-C17	-2.38	1.40	1.46
9	B6	101	BCL	O1A-CGA	2.38	1.29	1.22
7	AC	502	HEM	CHB-C4A	-2.38	1.34	1.39
7	AC	502	HEM	C1A-C2A	-2.38	1.39	1.44
14	AS	104	CRT	C22-C23	2.37	1.41	1.35
9	AJ	101	BCL	C3B-C4B	2.37	1.45	1.41
9	BI	102	BCL	CMA-C3A	2.37	1.58	1.53
9	BG	101	BCL	O1A-CGA	2.37	1.29	1.22
14	BO	103	CRT	C30-C28	-2.37	1.40	1.46
9	BW	102	BCL	C2-C3	2.37	1.38	1.33
14	BP	102	CRT	C25-C23	-2.37	1.40	1.46
9	A7	103	BCL	C3B-C4B	2.37	1.45	1.41
9	BV	101	BCL	C2C-C3C	-2.36	1.48	1.54
7	BC	502	HEM	C4D-ND	-2.36	1.36	1.40
9	BI	102	BCL	C4-C3	2.36	1.56	1.50
14	AR	102	CRT	C25-C23	-2.36	1.40	1.46
9	AT	101	BCL	CHB-C4A	2.36	1.43	1.37
9	BJ	101	BCL	MG-NC	2.36	2.11	2.06
9	A8	101	BCL	C3B-C4B	2.36	1.45	1.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	BJ	101	BCL	CMA-C3A	2.36	1.58	1.53
9	A9	102	BCL	CHB-C4A	2.36	1.43	1.37
9	AY	102	BCL	CHB-C4A	2.36	1.43	1.37
9	AB	101	BCL	CAA-C2A	2.35	1.58	1.54
9	BM	402	BCL	MG-NC	2.35	2.11	2.06
9	BS	102	BCL	CMA-C3A	2.35	1.58	1.53
9	B3	102	BCL	CHB-C4A	2.35	1.43	1.37
9	AK	102	BCL	C5-C3	2.35	1.56	1.51
14	A1	103	CRT	C4-C5	2.34	1.53	1.50
9	BM	401	BCL	C2C-C3C	-2.34	1.48	1.54
9	A3	103	BCL	C1B-C2B	2.34	1.48	1.43
9	AQ	102	BCL	C5-C3	2.34	1.56	1.51
14	BV	102	CRT	C9-C7	2.34	1.41	1.35
9	A2	101	BCL	CAA-C2A	2.34	1.58	1.54
9	AG	101	BCL	C3B-C4B	2.34	1.45	1.41
14	AR	102	CRT	C30-C28	-2.33	1.41	1.46
9	AL	301	BCL	CMB-C2B	2.33	1.55	1.50
7	BC	502	HEM	C4A-NA	-2.33	1.35	1.39
10	AM	403	BPH	C1A-C2A	2.33	1.54	1.51
9	BL	301	BCL	C3B-C2B	-2.32	1.34	1.42
9	BS	102	BCL	MG-NB	-2.32	2.01	2.05
14	AW	102	CRT	C25-C23	-2.32	1.41	1.46
14	B0	101	CRT	C35-C33	-2.32	1.41	1.46
14	B2	102	CRT	C21-C20	2.32	1.42	1.36
9	A7	103	BCL	CMA-C3A	2.32	1.58	1.53
9	AV	102	BCL	C7-C8	2.32	1.64	1.52
9	AP	101	BCL	O1A-CGA	2.32	1.29	1.22
9	BP	101	BCL	C5-C3	2.32	1.56	1.51
9	BZ	101	BCL	C2C-C3C	-2.32	1.48	1.54
9	B5	102	BCL	C2-C3	2.32	1.38	1.33
14	AG	102	CRT	C25-C23	-2.32	1.41	1.46
9	BK	102	BCL	C4-C3	2.32	1.56	1.50
9	BV	101	BCL	O1A-CGA	2.31	1.29	1.22
14	AX	102	CRT	C30-C28	-2.31	1.41	1.46
9	BQ	103	BCL	CMA-C3A	2.31	1.58	1.53
9	BZ	101	BCL	O2D-CED	-2.31	1.40	1.45
7	AC	503	HEM	CHD-C4C	-2.31	1.33	1.38
14	BA	102	CRT	C19-C17	2.31	1.41	1.35
9	AN	101	BCL	C4-C3	2.31	1.56	1.50
9	BW	102	BCL	C3B-C2B	-2.31	1.34	1.42
9	AO	102	BCL	C4-C3	2.31	1.56	1.50
9	AA	101	BCL	CHB-C4A	2.30	1.43	1.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
10	AM	403	BPH	C3A-C2A	-2.30	1.52	1.54
9	AI	102	BCL	C4-C3	2.30	1.56	1.50
9	BM	401	BCL	C2-C3	2.30	1.38	1.33
9	AY	102	BCL	O1D-CGD	2.30	1.27	1.21
9	B4	101	BCL	CMC-C2C	2.30	1.58	1.53
9	B5	102	BCL	CHB-C4A	2.30	1.43	1.37
9	BM	401	BCL	C3B-C4B	2.30	1.45	1.41
9	AM	402	BCL	C2C-C3C	-2.30	1.48	1.54
14	AN	102	CRT	C30-C28	-2.30	1.41	1.46
9	BB	101	BCL	C4-C3	2.30	1.56	1.50
14	BM	406	CRT	C37-C36	2.30	1.53	1.50
9	A9	102	BCL	C4-C3	2.30	1.56	1.50
9	BZ	101	BCL	MG-ND	-2.30	2.01	2.05
9	BU	102	BCL	CMC-C2C	2.30	1.58	1.53
9	AI	102	BCL	C3B-C4B	2.30	1.45	1.41
7	BC	503	HEM	FE-ND	2.29	2.02	1.94
13	BM	405	MQ8	C6-C5	2.29	1.43	1.39
9	AM	401	BCL	C2-C3	2.29	1.38	1.33
9	A5	102	BCL	O2D-CED	-2.29	1.40	1.45
9	AZ	101	BCL	C1D-C2D	-2.29	1.40	1.45
9	BM	402	BCL	C3B-C4B	2.29	1.45	1.41
9	AZ	101	BCL	C2-C3	2.29	1.38	1.33
7	BC	503	HEM	C4C-NC	-2.29	1.35	1.39
9	B7	103	BCL	CHB-C4A	2.29	1.43	1.37
9	BG	101	BCL	CMC-C2C	2.29	1.58	1.53
9	BQ	104	BCL	C2C-C3C	-2.28	1.48	1.54
9	AD	102	BCL	CMB-C2B	2.28	1.55	1.50
9	A2	101	BCL	C4B-NB	-2.28	1.34	1.37
7	AC	502	HEM	C3B-C4B	-2.28	1.40	1.44
9	AS	103	BCL	C4-C3	2.28	1.56	1.50
9	AX	101	BCL	CHB-C1B	2.28	1.44	1.39
7	BC	504	HEM	C3B-C4B	-2.28	1.40	1.44
14	A0	101	CRT	C6-C5	2.28	1.38	1.32
9	BL	303	BCL	C2-C3	2.28	1.38	1.33
9	AL	301	BCL	MG-NC	2.28	2.11	2.06
9	B6	101	BCL	C4-C3	2.28	1.56	1.50
9	AO	102	BCL	CAA-C2A	2.28	1.58	1.54
7	AC	501	HEM	C4D-ND	-2.28	1.36	1.40
9	B1	102	BCL	C3B-C4B	2.27	1.45	1.41
9	B0	102	BCL	CAA-C2A	2.27	1.58	1.54
7	BC	504	HEM	CAD-C3D	2.27	1.57	1.51
7	BC	501	HEM	C4C-NC	-2.27	1.35	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
14	AJ	102	CRT	C22-C23	2.27	1.41	1.35
9	BY	102	BCL	C4-C3	2.27	1.56	1.50
9	BW	102	BCL	C2C-C3C	-2.27	1.48	1.54
14	BN	102	CRT	C30-C28	-2.27	1.41	1.46
9	AR	101	BCL	CHB-C1B	2.27	1.44	1.39
9	BG	101	BCL	C4-C3	2.27	1.56	1.50
9	AK	102	BCL	CAA-CBA	-2.27	1.46	1.52
9	B4	101	BCL	C4-C3	2.27	1.56	1.50
9	AV	102	BCL	CMA-C3A	2.27	1.58	1.53
9	BD	102	BCL	C4-C3	2.27	1.56	1.50
9	B7	103	BCL	C4-C3	2.27	1.56	1.50
7	BC	501	HEM	C4D-ND	-2.27	1.36	1.40
7	BC	503	HEM	C4B-NB	2.27	1.43	1.38
9	B1	102	BCL	C2-C3	2.27	1.38	1.33
14	A0	101	CRT	C27-C28	2.27	1.41	1.35
9	B5	102	BCL	CHB-C1B	2.26	1.44	1.39
14	AX	102	CRT	C27-C28	2.26	1.41	1.35
14	A5	103	CRT	C25-C23	-2.26	1.41	1.46
9	BM	402	BCL	C3B-C2B	-2.26	1.34	1.42
9	A3	103	BCL	C4-C3	2.26	1.56	1.50
9	AZ	101	BCL	C2A-C1A	2.26	1.57	1.52
9	BT	101	BCL	C2C-C3C	-2.26	1.48	1.54
9	BY	102	BCL	CHB-C4A	2.26	1.43	1.37
9	A2	101	BCL	MG-ND	-2.25	2.01	2.05
9	A5	102	BCL	MG-NA	2.25	2.11	2.06
14	AX	102	CRT	C4-C5	2.25	1.53	1.50
9	A8	101	BCL	C2-C3	2.25	1.38	1.33
9	AF	102	BCL	MG-NA	2.25	2.11	2.06
9	A3	104	BCL	CHB-C4A	2.25	1.43	1.37
9	A3	104	BCL	C2-C3	2.25	1.38	1.33
9	AR	101	BCL	C3B-C2B	-2.25	1.34	1.42
14	BW	103	CRT	C19-C17	2.25	1.41	1.35
10	BM	403	BPH	C3B-C2B	2.25	1.43	1.39
9	AF	102	BCL	C3B-C4B	2.25	1.45	1.41
9	A0	102	BCL	CHB-C1B	2.25	1.44	1.39
9	B2	101	BCL	C4-C3	2.25	1.56	1.50
9	AV	102	BCL	CHB-C4A	2.25	1.43	1.37
9	AY	102	BCL	MG-NC	2.25	2.11	2.06
9	AL	303	BCL	C3B-C2B	-2.25	1.34	1.42
9	B9	102	BCL	C4-C3	2.24	1.56	1.50
9	BM	401	BCL	MG-NC	2.24	2.11	2.06
9	BN	101	BCL	MG-NC	2.24	2.11	2.06

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	AR	101	BCL	MG-ND	-2.24	2.01	2.05
9	AR	101	BCL	C3C-C4C	-2.24	1.48	1.51
9	AQ	102	BCL	O1D-CGD	2.24	1.26	1.21
14	A7	102	CRT	C22-C23	2.24	1.41	1.35
9	BQ	103	BCL	C4-C3	2.24	1.56	1.50
9	BB	101	BCL	MG-NC	2.24	2.11	2.06
9	AG	101	BCL	CBA-CGA	-2.24	1.44	1.50
9	AL	301	BCL	C4-C3	2.23	1.56	1.50
14	BN	102	CRT	C19-C17	2.23	1.41	1.35
9	BA	101	BCL	C4-C3	2.23	1.56	1.50
9	BP	101	BCL	O2D-CED	-2.23	1.40	1.45
14	A2	102	CRT	C15-C16	2.23	1.40	1.34
14	BF	103	CRT	C4-C5	2.23	1.53	1.50
9	A7	103	BCL	C2-C3	2.23	1.38	1.33
14	AX	102	CRT	C26-C25	2.23	1.40	1.34
9	AU	102	BCL	C4-C3	2.23	1.56	1.50
9	BP	101	BCL	C4D-CHA	2.23	1.46	1.38
14	AA	102	CRT	C19-C17	2.23	1.40	1.35
14	A7	102	CRT	C19-C17	2.23	1.40	1.35
7	AC	503	HEM	CHB-C4A	-2.23	1.34	1.39
14	A5	103	CRT	C16-C17	-2.23	1.41	1.46
9	AE	101	BCL	C4-C3	2.22	1.56	1.50
9	A1	102	BCL	C3B-C2B	-2.22	1.35	1.42
9	A7	103	BCL	C1B-C2B	2.22	1.48	1.43
9	B2	101	BCL	C3B-C2B	-2.22	1.35	1.42
7	BC	502	HEM	C1A-C2A	-2.22	1.40	1.44
9	AJ	101	BCL	CHB-C4A	2.22	1.43	1.37
9	BX	101	BCL	C3D-C2D	-2.22	1.32	1.39
11	BL	304	UQ8	C6-C1	2.22	1.39	1.35
9	A1	102	BCL	CBA-CGA	2.22	1.57	1.50
9	BV	101	BCL	C4-C3	2.22	1.56	1.50
14	BU	103	CRT	C30-C28	-2.22	1.41	1.46
9	B8	101	BCL	C2-C3	2.22	1.38	1.33
9	AM	402	BCL	C3B-C2B	-2.22	1.35	1.42
7	BC	503	HEM	C2A-C3A	-2.22	1.33	1.38
7	AC	503	HEM	CHA-C4D	-2.22	1.34	1.38
14	AP	102	CRT	C22-C23	2.22	1.40	1.35
9	BT	101	BCL	O2D-CED	-2.21	1.40	1.45
9	BX	101	BCL	CHB-C1B	2.21	1.44	1.39
9	BL	301	BCL	CMA-C3A	2.21	1.57	1.53
9	BM	402	BCL	CMB-C2B	2.21	1.55	1.50
13	AM	405	MQ8	C10-C5	2.21	1.44	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
14	AM	406	CRT	C22-C23	2.21	1.40	1.35
9	AE	101	BCL	C3B-C4B	2.21	1.45	1.41
14	AX	102	CRT	C22-C23	2.21	1.40	1.35
9	AJ	101	BCL	C4-C3	2.21	1.56	1.50
9	AX	101	BCL	CBD-CGD	-2.20	1.45	1.52
9	A0	102	BCL	MG-ND	-2.20	2.01	2.05
7	AC	504	HEM	C4A-NA	-2.20	1.35	1.39
14	BW	103	CRT	C16-C17	-2.20	1.41	1.46
14	A2	102	CRT	C22-C23	2.20	1.40	1.35
9	A3	104	BCL	MG-ND	-2.20	2.01	2.05
14	AW	102	CRT	C16-C17	-2.20	1.41	1.46
10	AL	302	BPH	CHA-CBD	2.20	1.54	1.51
7	AC	503	HEM	C4A-NA	-2.20	1.35	1.39
9	AM	401	BCL	MG-NC	2.20	2.11	2.06
14	AP	102	CRT	C25-C23	-2.19	1.41	1.46
7	BC	501	HEM	C1A-C2A	-2.19	1.40	1.44
9	AL	303	BCL	C4-C3	2.19	1.56	1.50
9	BL	303	BCL	C3B-C2B	-2.19	1.35	1.42
9	AY	102	BCL	CMA-C3A	2.19	1.57	1.53
9	AZ	101	BCL	C3B-C2B	-2.19	1.35	1.42
14	AN	102	CRT	C11-C12	-2.19	1.41	1.46
9	B9	102	BCL	C5-C3	2.19	1.55	1.51
9	BO	102	BCL	MG-ND	-2.19	2.01	2.05
9	BZ	101	BCL	CMC-C2C	2.18	1.57	1.53
9	AN	101	BCL	C3C-C4C	-2.18	1.48	1.51
9	AD	102	BCL	C4-C3	2.18	1.56	1.50
9	B6	101	BCL	MG-NA	2.18	2.11	2.06
14	AA	102	CRT	C22-C23	2.18	1.40	1.35
14	BS	103	CRT	C16-C17	-2.18	1.41	1.46
9	BW	102	BCL	C4-C3	2.18	1.56	1.50
14	BA	102	CRT	C22-C23	2.18	1.40	1.35
9	A2	101	BCL	CHB-C4A	2.17	1.43	1.37
9	AV	102	BCL	O2D-CED	-2.17	1.40	1.45
9	BJ	101	BCL	CMC-C2C	2.17	1.57	1.53
14	AW	102	CRT	C14-C12	2.17	1.40	1.35
7	AC	501	HEM	CMA-C3A	2.17	1.55	1.50
14	B0	101	CRT	C22-C23	2.17	1.40	1.35
9	A5	102	BCL	CMA-C3A	2.17	1.57	1.53
9	AL	303	BCL	C2-C3	2.17	1.38	1.33
9	A6	101	BCL	CMA-C3A	2.17	1.57	1.53
9	BX	101	BCL	C4-C3	2.17	1.56	1.50
9	B6	101	BCL	C5-C3	2.17	1.55	1.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	BO	102	BCL	C3B-C2B	-2.17	1.35	1.42
9	BT	101	BCL	MG-ND	-2.17	2.01	2.05
9	BQ	104	BCL	C4-C3	2.17	1.56	1.50
9	AO	102	BCL	MG-NA	2.17	2.11	2.06
9	AW	101	BCL	CHB-C4A	2.17	1.43	1.37
14	BN	102	CRT	C25-C23	-2.16	1.41	1.46
10	BL	302	BPH	C3D-C4D	2.16	1.44	1.41
9	AT	101	BCL	CMB-C2B	2.16	1.55	1.50
9	AB	101	BCL	C3B-C4B	2.16	1.45	1.41
7	BC	502	HEM	CHB-C4A	-2.16	1.34	1.39
9	AU	102	BCL	CHB-C4A	2.16	1.43	1.37
9	AU	102	BCL	C2-C3	2.16	1.38	1.33
7	BC	503	HEM	C4D-ND	-2.16	1.36	1.40
9	B2	101	BCL	C3B-C4B	2.16	1.45	1.41
9	A0	102	BCL	C4-C3	2.16	1.56	1.50
9	BP	101	BCL	CHB-C4A	2.16	1.43	1.37
14	BN	102	CRT	C15-C16	2.16	1.40	1.34
14	BN	102	CRT	C22-C23	2.16	1.40	1.35
14	BM	406	CRT	C25-C23	-2.16	1.41	1.46
9	AL	303	BCL	CMB-C2B	2.15	1.55	1.50
9	B6	101	BCL	CHB-C4A	2.15	1.43	1.37
9	BW	102	BCL	CMB-C2B	2.15	1.55	1.50
9	AM	401	BCL	CMB-C2B	2.15	1.55	1.50
9	AA	101	BCL	CMA-C3A	2.15	1.57	1.53
9	BM	401	BCL	C4-C3	2.15	1.56	1.50
9	BK	102	BCL	C3B-C2B	-2.15	1.35	1.42
14	AM	406	CRT	C19-C17	2.15	1.40	1.35
9	BL	301	BCL	C3B-C4B	2.15	1.45	1.41
14	A5	103	CRT	C30-C28	-2.15	1.41	1.46
9	BL	303	BCL	CMB-C2B	2.15	1.55	1.50
14	B2	102	CRT	C26-C25	2.15	1.40	1.34
9	BB	101	BCL	CAA-C2A	2.15	1.58	1.54
9	AN	101	BCL	O2D-CED	-2.14	1.40	1.45
9	AT	101	BCL	O2D-CED	-2.14	1.40	1.45
9	B4	101	BCL	O1A-CGA	2.14	1.28	1.22
14	BV	102	CRT	C21-C20	2.14	1.42	1.36
9	B3	102	BCL	C2-C3	2.14	1.38	1.33
9	AS	103	BCL	CHB-C4A	2.14	1.43	1.37
14	BN	102	CRT	C14-C12	2.14	1.40	1.35
9	BJ	101	BCL	C4B-NB	-2.14	1.35	1.37
7	BC	502	HEM	C3B-C4B	-2.14	1.40	1.44
15	AM	409	PEF	P-O4P	2.14	1.67	1.59

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
14	AS	104	CRT	C30-C28	-2.14	1.41	1.46
9	BI	102	BCL	O2D-CED	-2.14	1.40	1.45
9	AR	101	BCL	C3B-C4B	2.14	1.45	1.41
14	AJ	102	CRT	C25-C23	-2.14	1.41	1.46
14	BP	102	CRT	C15-C16	2.14	1.40	1.34
9	AQ	102	BCL	CMC-C2C	2.13	1.57	1.53
9	BL	301	BCL	CAA-CBA	-2.13	1.46	1.52
9	AE	101	BCL	O1A-CGA	2.13	1.28	1.22
9	A3	103	BCL	C3B-C2B	-2.13	1.35	1.42
9	AU	102	BCL	C1D-C2D	2.13	1.49	1.45
9	AD	102	BCL	CMA-C3A	2.13	1.57	1.53
9	BT	101	BCL	C4-C3	2.13	1.55	1.50
9	B1	102	BCL	CMD-C2D	2.13	1.55	1.50
14	BS	103	CRT	C19-C17	2.13	1.40	1.35
14	AG	102	CRT	C16-C17	-2.13	1.41	1.46
9	BO	102	BCL	CMC-C2C	2.13	1.57	1.53
9	BQ	103	BCL	CAA-C2A	2.13	1.58	1.54
9	AM	402	BCL	CMD-C2D	2.13	1.55	1.50
9	AY	102	BCL	C1B-C2B	2.13	1.48	1.43
14	AT	102	CRT	C16-C17	-2.12	1.41	1.46
9	BT	101	BCL	O1A-CGA	2.12	1.28	1.22
15	BQ	101	PEF	P-O4P	2.12	1.67	1.59
14	BV	102	CRT	C26-C25	2.12	1.40	1.34
9	BD	102	BCL	MG-NC	2.12	2.11	2.06
9	BP	101	BCL	C1B-NB	2.12	1.40	1.37
9	BE	101	BCL	CBB-CAB	2.12	1.54	1.50
11	AL	304	UQ8	C6-C1	2.12	1.39	1.35
9	AL	301	BCL	CMA-C3A	2.12	1.57	1.53
9	AO	102	BCL	CMA-C3A	2.11	1.57	1.53
9	BF	102	BCL	O2D-CED	-2.11	1.40	1.45
9	BP	101	BCL	C4-C3	2.11	1.55	1.50
9	AA	101	BCL	O1D-CGD	2.11	1.26	1.21
9	AK	102	BCL	O2D-CED	-2.11	1.40	1.45
7	AC	502	HEM	C3C-C2C	2.11	1.41	1.37
14	B5	103	CRT	C19-C17	2.11	1.40	1.35
9	AM	402	BCL	C3B-C4B	2.11	1.45	1.41
10	BL	302	BPH	O2D-CGD	2.11	1.38	1.33
13	AM	405	MQ8	C8-C9	2.11	1.42	1.38
9	A5	102	BCL	CHB-C4A	2.11	1.43	1.37
9	BQ	104	BCL	C4B-NB	2.11	1.40	1.37
15	AS	101	PEF	P-O4P	2.11	1.67	1.59
9	BL	303	BCL	CAA-C2A	2.11	1.57	1.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	A2	101	BCL	CMB-C2B	2.11	1.55	1.50
9	A6	101	BCL	O2D-CED	-2.11	1.40	1.45
9	BM	402	BCL	CMD-C2D	2.11	1.55	1.50
9	A2	101	BCL	C4-C3	2.11	1.55	1.50
9	B8	101	BCL	C6-C5	2.11	1.59	1.52
14	BW	103	CRT	C37-C36	2.11	1.53	1.50
9	BT	101	BCL	C1D-C2D	-2.11	1.41	1.45
9	AP	101	BCL	CHB-C1B	2.10	1.44	1.39
9	BK	102	BCL	CMA-C3A	2.10	1.57	1.53
9	A8	101	BCL	C1B-C2B	2.10	1.48	1.43
9	AL	301	BCL	C3B-C2B	-2.10	1.35	1.42
9	B8	101	BCL	O2D-CED	-2.10	1.40	1.45
9	AW	101	BCL	C3C-C4C	2.10	1.54	1.51
9	AL	301	BCL	CAA-CBA	-2.10	1.46	1.52
9	AL	303	BCL	MG-ND	-2.10	2.01	2.05
9	A1	102	BCL	CMC-C2C	2.10	1.57	1.53
9	BP	101	BCL	CHD-C1D	-2.10	1.34	1.38
9	AO	102	BCL	CAA-CBA	-2.10	1.46	1.52
9	AQ	102	BCL	C3A-C2A	-2.10	1.48	1.54
14	AX	102	CRT	C10-C11	2.09	1.40	1.34
14	AW	102	CRT	C37-C36	2.09	1.53	1.50
9	AS	103	BCL	C1D-ND	-2.09	1.35	1.37
9	B7	103	BCL	MG-NA	2.09	2.11	2.06
9	AB	101	BCL	C4-C3	2.09	1.55	1.50
9	B8	101	BCL	C3A-C2A	-2.09	1.48	1.54
15	AM	407	PEF	P-O4P	2.09	1.67	1.59
9	BX	101	BCL	CBB-CAB	2.09	1.54	1.50
14	BF	103	CRT	C30-C28	-2.09	1.41	1.46
7	BC	503	HEM	CHB-C1B	2.09	1.42	1.38
9	A6	101	BCL	O1A-CGA	2.09	1.28	1.22
9	A8	101	BCL	CMA-C3A	2.08	1.57	1.53
9	B5	102	BCL	C4-C3	2.08	1.55	1.50
15	AH	301	PEF	P-O4P	2.08	1.67	1.59
9	AO	102	BCL	MG-NC	2.08	2.11	2.06
14	BB	102	CRT	C11-C12	-2.08	1.41	1.46
9	AN	101	BCL	CHB-C4A	2.08	1.42	1.37
14	A0	101	CRT	C22-C23	2.08	1.40	1.35
14	B5	103	CRT	C22-C23	2.08	1.40	1.35
14	AG	102	CRT	C11-C12	-2.08	1.41	1.46
15	BM	407	PEF	P-O4P	2.08	1.67	1.59
9	BB	101	BCL	C3B-C2B	-2.08	1.35	1.42
15	AM	408	PEF	P-O4P	2.08	1.67	1.59

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	A3	104	BCL	C2A-C1A	2.07	1.56	1.52
14	AJ	102	CRT	C16-C17	-2.07	1.41	1.46
14	BF	103	CRT	C14-C12	2.07	1.40	1.35
14	A0	101	CRT	C16-C17	-2.07	1.41	1.46
14	AG	102	CRT	C30-C28	-2.07	1.41	1.46
9	BX	101	BCL	CMC-C2C	2.07	1.57	1.53
9	BD	102	BCL	CHD-C1D	-2.07	1.34	1.38
9	AI	102	BCL	CHB-C4A	2.07	1.42	1.37
9	BF	102	BCL	C4-C3	2.07	1.55	1.50
14	AJ	102	CRT	C30-C28	-2.07	1.41	1.46
7	AC	504	HEM	CHD-C4C	-2.07	1.34	1.38
14	AT	102	CRT	C22-C23	2.07	1.40	1.35
9	BL	301	BCL	MG-NC	2.07	2.11	2.06
7	AC	504	HEM	C2A-C3A	-2.07	1.33	1.38
7	BC	502	HEM	CHA-C4D	-2.06	1.34	1.38
9	B9	102	BCL	CMA-C3A	2.06	1.57	1.53
9	AZ	101	BCL	C4-C3	2.06	1.55	1.50
13	AM	405	MQ8	C6-C5	2.06	1.43	1.39
14	BV	102	CRT	C4-C5	2.06	1.53	1.50
9	AZ	101	BCL	CAA-CBA	-2.06	1.46	1.52
9	BE	101	BCL	CMA-C3A	2.06	1.57	1.53
9	BM	402	BCL	C4-C3	2.06	1.55	1.50
9	B0	102	BCL	CMC-C2C	2.06	1.57	1.53
14	AT	102	CRT	C4-C5	2.06	1.53	1.50
9	BN	101	BCL	C2C-C3C	-2.06	1.48	1.54
9	B1	102	BCL	C4-C3	2.06	1.55	1.50
14	BV	102	CRT	C4-C1	2.06	1.56	1.52
9	A5	102	BCL	C3B-C2B	-2.06	1.35	1.42
9	B7	103	BCL	CMA-C3A	2.06	1.57	1.53
7	BC	501	HEM	C1C-C2C	2.05	1.49	1.45
9	BW	102	BCL	C3B-C4B	2.05	1.45	1.41
9	BO	102	BCL	C4-C3	2.05	1.55	1.50
7	AC	502	HEM	C4D-ND	-2.05	1.36	1.40
14	BU	103	CRT	C14-C12	2.05	1.40	1.35
7	AC	504	HEM	CAD-C3D	2.05	1.56	1.51
14	BW	103	CRT	C25-C23	-2.05	1.41	1.46
14	A5	103	CRT	C4-C5	2.05	1.53	1.50
9	A7	103	BCL	O2D-CED	-2.05	1.40	1.45
9	BB	101	BCL	C1D-C2D	-2.05	1.41	1.45
14	BG	102	CRT	C22-C23	2.04	1.40	1.35
9	AJ	101	BCL	C3B-C2B	-2.04	1.35	1.42
9	A9	102	BCL	CMA-C3A	2.04	1.57	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	BM	402	BCL	CAA-CBA	-2.04	1.46	1.52
9	A1	102	BCL	C3A-C4A	2.04	1.58	1.51
9	A8	101	BCL	CHB-C4A	2.04	1.42	1.37
9	A9	102	BCL	C1B-C2B	2.04	1.47	1.43
9	A7	103	BCL	O1A-CGA	2.04	1.28	1.22
9	A5	102	BCL	C1B-C2B	2.04	1.47	1.43
7	AC	503	HEM	O2D-CGD	-2.04	1.24	1.30
9	B8	101	BCL	C7-C8	2.04	1.62	1.52
9	BQ	103	BCL	MG-ND	-2.04	2.01	2.05
9	AV	102	BCL	MG-NA	2.04	2.11	2.06
14	A2	102	CRT	C30-C28	-2.04	1.41	1.46
9	AB	101	BCL	CHB-C4A	2.03	1.42	1.37
9	A1	102	BCL	C1B-C2B	2.03	1.47	1.43
14	B0	101	CRT	C4-C1	2.03	1.56	1.52
9	AO	102	BCL	C5-C3	2.03	1.55	1.51
9	AS	103	BCL	O2D-CED	-2.03	1.40	1.45
14	A5	103	CRT	C4-C1	2.03	1.56	1.52
14	AP	102	CRT	C15-C16	2.03	1.40	1.34
7	AC	501	HEM	C3B-C4B	-2.03	1.41	1.44
9	AR	101	BCL	O1A-CGA	2.03	1.28	1.22
14	B0	101	CRT	C25-C23	-2.03	1.41	1.46
9	B3	102	BCL	CMA-C3A	2.03	1.57	1.53
9	BP	101	BCL	O1A-CGA	2.03	1.28	1.22
14	B5	103	CRT	C4-C1	2.03	1.56	1.52
9	BL	303	BCL	CMA-C3A	2.03	1.57	1.53
14	BW	103	CRT	C22-C23	2.03	1.40	1.35
14	A2	102	CRT	C25-C23	-2.02	1.41	1.46
9	A1	102	BCL	MG-NA	2.02	2.11	2.06
9	BA	101	BCL	C3B-C4B	2.02	1.45	1.41
9	A3	103	BCL	CBA-CGA	2.02	1.56	1.50
9	BQ	104	BCL	MG-ND	-2.02	2.01	2.05
9	BY	102	BCL	MG-NA	-2.02	2.01	2.06
9	A0	102	BCL	C3B-C2B	-2.02	1.35	1.42
10	AM	403	BPH	C3B-C2B	2.02	1.43	1.39
13	BM	405	MQ8	C8-C9	2.02	1.42	1.38
9	AL	301	BCL	C3B-C4B	2.02	1.45	1.41
9	AS	103	BCL	C5-C3	2.02	1.55	1.51
9	B8	101	BCL	C4-C3	2.02	1.55	1.50
9	BM	402	BCL	C2-C3	2.02	1.37	1.33
9	B0	102	BCL	C4-C3	2.02	1.55	1.50
9	AU	102	BCL	MG-ND	-2.02	2.01	2.05
9	AZ	101	BCL	O2D-CED	-2.02	1.40	1.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
14	BM	406	CRT	C22-C23	2.02	1.40	1.35
14	A0	101	CRT	C25-C23	-2.02	1.41	1.46
9	AI	102	BCL	O2D-CED	-2.01	1.40	1.45
7	BC	502	HEM	C3C-C2C	2.01	1.41	1.37
9	AZ	101	BCL	CMA-C3A	2.01	1.57	1.53
14	B5	103	CRT	C30-C28	-2.01	1.41	1.46
9	A3	104	BCL	CHB-C1B	2.01	1.43	1.39
14	A7	102	CRT	C26-C25	2.01	1.39	1.34
10	BL	302	BPH	O2A-CGA	2.01	1.39	1.33
9	BX	101	BCL	CBD-CGD	2.01	1.58	1.52
14	A7	102	CRT	C15-C16	2.01	1.39	1.34
9	AE	101	BCL	MG-ND	-2.01	2.01	2.05
13	BM	405	MQ8	C3-C4	2.01	1.52	1.47
14	BM	406	CRT	C19-C17	2.00	1.40	1.35
16	A3	101	PO4	P-O4	-2.00	1.48	1.54
7	AC	503	HEM	C3C-C2C	2.00	1.41	1.37

All (2456) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
14	A7	102	CRT	C20-C21-C22	-18.13	86.41	123.52
9	AW	101	BCL	O2A-C1-C2	17.78	176.53	108.11
9	A0	102	BCL	O2A-C1-C2	15.01	165.88	108.11
9	AZ	101	BCL	C1-O2A-CGA	14.84	152.58	116.65
9	A3	104	BCL	O2A-C1-C2	13.20	158.90	108.11
14	A7	102	CRT	C21-C22-C23	-11.15	111.64	127.28
9	B8	101	BCL	C1-C2-C3	11.15	144.47	126.20
9	AZ	101	BCL	O2A-C1-C2	10.87	149.94	108.11
9	BM	402	BCL	C1-C2-C3	10.62	143.59	126.20
9	AN	101	BCL	C1-C2-C3	10.45	143.32	126.20
9	A8	101	BCL	O2A-C1-C2	10.36	147.99	108.11
9	A8	101	BCL	C1-C2-C3	10.36	143.17	126.20
9	AX	101	BCL	O2A-C1-C2	10.32	147.82	108.11
9	AE	101	BCL	C1-C2-C3	10.24	142.97	126.20
9	B0	102	BCL	C1-C2-C3	10.21	142.93	126.20
14	A7	102	CRT	C21-C20-C19	-9.98	103.09	123.52
9	B4	101	BCL	C1-C2-C3	9.92	142.45	126.20
9	A9	102	BCL	O2A-C1-C2	9.77	145.70	108.11
9	BP	101	BCL	C1-C2-C3	9.74	142.16	126.20
9	AJ	101	BCL	C1-C2-C3	9.66	142.03	126.20
9	BX	101	BCL	C1-C2-C3	9.59	141.91	126.20
9	A5	102	BCL	O2A-C1-C2	9.51	144.71	108.11

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
9	A9	102	BCL	C1-O2A-CGA	-9.33	94.07	116.65
9	AP	101	BCL	C1-C2-C3	9.33	141.48	126.20
9	AU	102	BCL	O2A-C1-C2	9.24	143.66	108.11
9	B7	103	BCL	C1-C2-C3	9.17	141.22	126.20
9	A9	102	BCL	C1-C2-C3	9.15	141.19	126.20
10	AM	403	BPH	C4D-CHA-CBD	9.10	112.82	108.45
9	AA	101	BCL	C1-C2-C3	9.08	141.07	126.20
9	BV	101	BCL	C1-C2-C3	9.05	141.02	126.20
9	AB	101	BCL	C1-C2-C3	9.04	141.00	126.20
9	BB	101	BCL	C1-C2-C3	8.98	140.91	126.20
14	BF	103	CRT	C4-C5-C6	-8.97	112.32	124.91
9	A1	102	BCL	C1-C2-C3	8.93	140.83	126.20
9	B2	101	BCL	C1-C2-C3	8.89	140.76	126.20
9	BE	101	BCL	C1-C2-C3	8.81	140.63	126.20
10	BM	403	BPH	C4D-CHA-CBD	8.80	112.68	108.45
9	A2	101	BCL	C1-C2-C3	8.80	140.61	126.20
9	AX	101	BCL	C1-C2-C3	8.69	140.44	126.20
9	A6	101	BCL	C1-C2-C3	8.68	140.42	126.20
14	AS	104	CRT	C37-C36-C35	-8.68	112.73	124.91
9	A0	102	BCL	C1-C2-C3	8.64	140.36	126.20
9	AF	102	BCL	C1-C2-C3	8.64	140.35	126.20
9	BY	102	BCL	C1-C2-C3	8.61	140.30	126.20
9	AM	402	BCL	C1-C2-C3	8.57	140.23	126.20
9	A3	104	BCL	C1-C2-C3	8.56	140.22	126.20
9	B9	102	BCL	C1-C2-C3	8.50	140.12	126.20
9	BT	101	BCL	C1-C2-C3	8.48	140.09	126.20
9	BA	101	BCL	C1-C2-C3	8.46	140.06	126.20
9	B6	101	BCL	C1-C2-C3	8.45	140.04	126.20
9	BK	102	BCL	C1-C2-C3	8.40	139.97	126.20
9	A8	101	BCL	C1-O2A-CGA	8.40	136.98	116.65
9	BQ	103	BCL	C1-C2-C3	8.37	139.91	126.20
9	B5	102	BCL	C1-C2-C3	8.36	139.90	126.20
9	BS	102	BCL	C1-C2-C3	8.36	139.89	126.20
9	AM	401	BCL	C1-C2-C3	8.34	139.86	126.20
9	A5	102	BCL	C1-C2-C3	8.33	139.85	126.20
9	BI	102	BCL	C1-C2-C3	8.33	139.84	126.20
9	A2	101	BCL	O2A-C1-C2	8.26	139.91	108.11
9	AR	101	BCL	C1-C2-C3	8.23	139.69	126.20
9	BB	101	BCL	C4A-NA-C1A	8.23	110.43	106.68
9	AU	102	BCL	C1-C2-C3	8.18	139.60	126.20
9	AQ	102	BCL	C1-C2-C3	8.16	139.56	126.20
9	BP	101	BCL	C4A-NA-C1A	8.16	110.40	106.68

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
9	AS	103	BCL	C1-C2-C3	8.14	139.53	126.20
9	A7	103	BCL	C1-C2-C3	8.13	139.52	126.20
9	BM	401	BCL	C1-C2-C3	8.12	139.49	126.20
9	AY	102	BCL	C1-C2-C3	8.09	139.46	126.20
9	BU	102	BCL	C1-C2-C3	8.08	139.43	126.20
9	AL	303	BCL	C1-C2-C3	8.08	139.43	126.20
9	BF	102	BCL	C4A-NA-C1A	8.04	110.35	106.68
9	AG	101	BCL	C1-C2-C3	8.04	139.37	126.20
9	B3	102	BCL	C1-C2-C3	8.04	139.37	126.20
9	BL	303	BCL	C1-C2-C3	8.01	139.32	126.20
9	AD	102	BCL	C1-C2-C3	7.95	139.22	126.20
9	B1	102	BCL	C1-C2-C3	7.95	139.22	126.20
9	BD	102	BCL	C4A-NA-C1A	7.93	110.30	106.68
9	AI	102	BCL	C1-C2-C3	7.93	139.19	126.20
9	BQ	104	BCL	C1-C2-C3	7.91	139.15	126.20
9	AK	102	BCL	C1-C2-C3	7.78	138.94	126.20
9	AO	102	BCL	C1-C2-C3	7.72	138.85	126.20
9	BD	102	BCL	C1-C2-C3	7.71	138.82	126.20
9	BW	102	BCL	C1-C2-C3	7.67	138.77	126.20
9	AW	101	BCL	C1-C2-C3	7.62	138.69	126.20
9	BF	102	BCL	C1-C2-C3	7.60	138.64	126.20
9	BJ	101	BCL	C1-C2-C3	7.59	138.63	126.20
9	BG	101	BCL	C4A-NA-C1A	7.58	110.14	106.68
9	BG	101	BCL	C1-C2-C3	7.58	138.61	126.20
9	AT	101	BCL	C4A-NA-C1A	7.57	110.13	106.68
9	A5	102	BCL	C4A-NA-C1A	7.51	110.10	106.68
9	BO	102	BCL	C1-C2-C3	7.50	138.48	126.20
14	BU	103	CRT	C10-C9-C7	-7.41	116.88	127.28
14	B7	102	CRT	C4-C5-C6	-7.39	114.54	124.91
9	A3	103	BCL	C4A-NA-C1A	7.35	110.03	106.68
9	AP	101	BCL	C4A-NA-C1A	7.29	110.01	106.68
9	AK	102	BCL	C4A-NA-C1A	7.25	109.98	106.68
9	BZ	101	BCL	C1-C2-C3	7.25	138.07	126.20
9	AU	102	BCL	C4A-NA-C1A	7.22	109.97	106.68
9	B9	102	BCL	C4A-NA-C1A	7.17	109.95	106.68
9	AR	101	BCL	C4A-NA-C1A	7.15	109.94	106.68
14	BF	103	CRT	C37-C36-C35	-7.15	114.87	124.91
9	A2	101	BCL	C4A-NA-C1A	7.11	109.92	106.68
9	AZ	101	BCL	C1-C2-C3	7.09	137.82	126.20
14	B1	103	CRT	C37-C36-C35	-7.04	115.02	124.91
9	A3	103	BCL	C1-C2-C3	6.98	137.64	126.20
9	AT	101	BCL	C1-C2-C3	6.97	137.62	126.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
9	A1	102	BCL	C4A-NA-C1A	6.97	109.86	106.68
9	B3	102	BCL	C4A-NA-C1A	6.95	109.85	106.68
9	B5	102	BCL	C4A-NA-C1A	6.92	109.83	106.68
14	BV	102	CRT	C37-C36-C35	-6.91	115.22	124.91
9	BI	102	BCL	C4A-NA-C1A	6.89	109.82	106.68
9	AE	101	BCL	C4A-NA-C1A	6.89	109.82	106.68
14	AX	102	CRT	C10-C9-C7	-6.89	117.62	127.28
9	BO	102	BCL	C4-C3-C5	-6.81	103.40	115.23
9	AJ	101	BCL	C4A-NA-C1A	6.81	109.78	106.68
9	AG	101	BCL	C4A-NA-C1A	6.81	109.78	106.68
9	AX	101	BCL	C4A-NA-C1A	6.80	109.78	106.68
9	A9	102	BCL	C4A-NA-C1A	6.78	109.77	106.68
9	AD	102	BCL	C4A-NA-C1A	6.76	109.77	106.68
9	A0	102	BCL	C4A-NA-C1A	6.74	109.75	106.68
9	B7	103	BCL	C4A-NA-C1A	6.71	109.74	106.68
10	BL	302	BPH	C4D-CHA-CBD	6.69	111.67	108.45
14	B2	102	CRT	C4-C5-C6	-6.66	115.56	124.91
9	AS	103	BCL	C4-C3-C5	-6.63	103.72	115.23
9	B6	101	BCL	C4A-NA-C1A	6.62	109.70	106.68
14	A1	103	CRT	C10-C9-C7	-6.60	118.02	127.28
14	AT	102	CRT	C4-C5-C6	-6.58	115.68	124.91
9	BQ	104	BCL	C4A-NA-C1A	6.55	109.67	106.68
14	AJ	102	CRT	C37-C36-C35	-6.54	115.72	124.91
9	A7	103	BCL	C4A-NA-C1A	6.54	109.66	106.68
9	BU	102	BCL	C4A-NA-C1A	6.53	109.66	106.68
9	BL	301	BCL	C1-C2-C3	6.46	136.79	126.20
9	BM	402	BCL	C4A-NA-C1A	6.46	109.63	106.68
9	BX	101	BCL	C4-C3-C5	-6.45	104.03	115.23
9	BP	101	BCL	C4-C3-C5	-6.45	104.03	115.23
9	AB	101	BCL	C4A-NA-C1A	6.44	109.62	106.68
9	AQ	102	BCL	C4A-NA-C1A	6.44	109.62	106.68
9	AS	103	BCL	C4A-NA-C1A	6.44	109.62	106.68
9	BB	101	BCL	C4-C3-C5	-6.43	104.07	115.23
9	BS	102	BCL	C4-C3-C5	-6.42	104.09	115.23
9	BN	101	BCL	C1-C2-C3	6.41	136.70	126.20
9	AL	301	BCL	C1-C2-C3	6.41	136.70	126.20
9	AD	102	BCL	OBB-CAB-C3B	6.40	131.75	120.43
9	A6	101	BCL	C4A-NA-C1A	6.40	109.60	106.68
9	AL	301	BCL	C4A-NA-C1A	6.40	109.60	106.68
9	BV	101	BCL	C4A-NA-C1A	6.37	109.58	106.68
9	AG	101	BCL	C4-C3-C5	-6.36	104.19	115.23
9	BU	102	BCL	C4-C3-C5	-6.35	104.21	115.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
9	AP	101	BCL	C4-C3-C5	-6.34	104.22	115.23
9	BG	101	BCL	C4-C3-C5	-6.32	104.25	115.23
9	BL	301	BCL	C4A-NA-C1A	6.32	109.56	106.68
9	A2	101	BCL	C4-C3-C5	-6.31	104.27	115.23
9	AX	101	BCL	C4-C3-C5	-6.31	104.27	115.23
9	BZ	101	BCL	C4A-NA-C1A	6.30	109.56	106.68
9	AA	101	BCL	C4A-NA-C1A	6.28	109.54	106.68
10	AL	302	BPH	C4D-CHA-CBD	6.28	111.47	108.45
14	AJ	102	CRT	C4-C5-C6	-6.27	116.11	124.91
9	AZ	101	BCL	C4-C3-C5	-6.25	104.37	115.23
9	BK	102	BCL	C4A-NA-C1A	6.24	109.53	106.68
14	B7	102	CRT	C37-C36-C35	-6.22	116.17	124.91
9	B4	101	BCL	C4A-NA-C1A	6.21	109.51	106.68
9	AY	102	BCL	C4-C3-C5	-6.21	104.45	115.23
9	BY	102	BCL	C4A-NA-C1A	6.20	109.51	106.68
9	AB	101	BCL	C4-C3-C5	-6.20	104.47	115.23
9	BT	101	BCL	C4A-NA-C1A	6.20	109.51	106.68
9	BX	101	BCL	C4A-NA-C1A	6.19	109.50	106.68
9	AK	102	BCL	C4-C3-C5	-6.19	104.49	115.23
9	AE	101	BCL	C4-C3-C5	-6.17	104.52	115.23
9	BU	102	BCL	OBB-CAB-C3B	6.17	131.34	120.43
9	B6	101	BCL	C4-C3-C5	-6.17	104.52	115.23
9	BS	102	BCL	C4A-NA-C1A	6.13	109.48	106.68
9	B2	101	BCL	C4-C3-C5	-6.13	104.60	115.23
9	A1	102	BCL	OBB-CAB-C3B	6.12	131.25	120.43
9	AY	102	BCL	OBB-CAB-C3B	6.11	131.24	120.43
9	AD	102	BCL	C4-C3-C5	-6.10	104.63	115.23
14	AX	102	CRT	C15-C14-C12	-6.10	118.72	127.28
9	A2	101	BCL	OBB-CAB-C3B	6.10	131.21	120.43
9	BQ	103	BCL	C4-C3-C5	-6.10	104.64	115.23
9	A1	102	BCL	O2A-CGA-CBA	6.08	130.37	111.83
9	AO	102	BCL	C4-C3-C5	-6.07	104.69	115.23
9	AM	402	BCL	C4-C3-C5	-6.06	104.71	115.23
9	BD	102	BCL	OBB-CAB-C3B	6.05	131.14	120.43
9	A3	104	BCL	C4A-NA-C1A	6.04	109.43	106.68
9	AW	101	BCL	C4-C3-C5	-6.03	104.76	115.23
9	AZ	101	BCL	OBB-CAB-C3B	6.03	131.10	120.43
9	B1	102	BCL	OBB-CAB-C3B	6.02	131.08	120.43
9	BE	101	BCL	OBB-CAB-C3B	6.02	131.08	120.43
9	BY	102	BCL	OBB-CAB-C3B	6.00	131.03	120.43
9	BK	102	BCL	C4-C3-C5	-5.99	104.83	115.23
9	A7	103	BCL	O2A-C1-C2	5.99	131.15	108.11

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
9	AV	102	BCL	C4A-NA-C1A	5.98	109.41	106.68
9	BW	102	BCL	C4-C3-C5	-5.98	104.86	115.23
9	A1	102	BCL	C4-C3-C5	-5.97	104.86	115.23
9	AB	101	BCL	OBB-CAB-C3B	5.97	130.99	120.43
9	B1	102	BCL	C4A-NA-C1A	5.96	109.40	106.68
9	B9	102	BCL	C4-C3-C5	-5.95	104.89	115.23
9	BQ	103	BCL	OBB-CAB-C3B	5.95	130.96	120.43
9	AF	102	BCL	C4-C3-C5	-5.95	104.90	115.23
14	B0	101	CRT	C26-C27-C28	-5.95	118.93	127.28
9	A9	102	BCL	C4-C3-C5	-5.94	104.91	115.23
9	AJ	101	BCL	C4-C3-C5	-5.94	104.91	115.23
9	A6	101	BCL	C4-C3-C5	-5.94	104.92	115.23
10	BM	403	BPH	C2D-C1D-ND	-5.93	105.14	109.43
9	B0	102	BCL	C4A-NA-C1A	5.93	109.38	106.68
9	B1	102	BCL	C4-C3-C5	-5.92	104.95	115.23
9	BW	102	BCL	OBB-CAB-C3B	5.92	130.89	120.43
9	BI	102	BCL	OBB-CAB-C3B	5.91	130.89	120.43
9	A7	103	BCL	C4-C3-C5	-5.90	104.99	115.23
9	BN	101	BCL	C4A-NA-C1A	5.89	109.36	106.68
9	AN	101	BCL	C4A-NA-C1A	5.87	109.36	106.68
9	AA	101	BCL	OBB-CAB-C3B	5.87	130.81	120.43
9	AQ	102	BCL	C4-C3-C5	-5.87	105.04	115.23
9	AQ	102	BCL	OBB-CAB-C3B	5.84	130.76	120.43
9	BL	303	BCL	OBB-CAB-C3B	5.84	130.76	120.43
9	BJ	101	BCL	OBB-CAB-C3B	5.84	130.76	120.43
9	AR	101	BCL	C4-C3-C5	-5.83	105.10	115.23
9	BF	102	BCL	C4-C3-C5	-5.83	105.11	115.23
9	BO	102	BCL	C4A-NA-C1A	5.82	109.33	106.68
9	AU	102	BCL	OBB-CAB-C3B	5.82	130.72	120.43
14	AA	102	CRT	C37-C36-C35	-5.82	116.75	124.91
9	BQ	104	BCL	C4-C3-C5	-5.81	105.15	115.23
9	AY	102	BCL	C4A-NA-C1A	5.80	109.33	106.68
9	BG	101	BCL	OBB-CAB-C3B	5.80	130.69	120.43
11	BL	304	UQ8	C46-C44-C45	5.79	127.92	114.59
9	BW	102	BCL	O2D-CGD-CBD	5.79	121.36	111.23
9	AW	101	BCL	OBB-CAB-C3B	5.79	130.67	120.43
9	A1	102	BCL	C1-O2A-CGA	-5.78	102.65	116.65
9	B2	101	BCL	OBB-CAB-C3B	5.78	130.65	120.43
11	AL	304	UQ8	C46-C44-C45	5.78	127.88	114.59
9	B3	102	BCL	C4-C3-C5	-5.77	105.20	115.23
9	AL	303	BCL	O2D-CGD-CBD	5.77	121.32	111.23
9	AT	101	BCL	OBB-CAB-C3B	5.77	130.63	120.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
9	BK	102	BCL	OBB-CAB-C3B	5.77	130.63	120.43
9	AF	102	BCL	C4A-NA-C1A	5.76	109.31	106.68
9	AM	402	BCL	C4A-NA-C1A	5.76	109.31	106.68
9	BJ	101	BCL	C4-C3-C5	-5.76	105.23	115.23
9	BM	402	BCL	OBB-CAB-C3B	5.76	130.61	120.43
9	B8	101	BCL	C4A-NA-C1A	5.76	109.31	106.68
9	BU	102	BCL	O2D-CGD-CBD	5.75	121.29	111.23
9	BT	101	BCL	OBB-CAB-C3B	5.75	130.60	120.43
9	AM	401	BCL	C4-C3-C5	-5.75	105.25	115.23
10	AM	403	BPH	C2D-C1D-ND	-5.75	105.27	109.43
9	AL	303	BCL	OBB-CAB-C3B	5.75	130.59	120.43
9	BY	102	BCL	C4-C3-C5	-5.74	105.26	115.23
9	A3	104	BCL	C4-C3-C5	-5.74	105.27	115.23
9	AS	103	BCL	OBB-CAB-C3B	5.74	130.57	120.43
9	BD	102	BCL	C4-C3-C5	-5.73	105.28	115.23
9	A7	103	BCL	OBB-CAB-C3B	5.73	130.56	120.43
9	BL	301	BCL	OBB-CAB-C3B	5.73	130.56	120.43
14	BN	102	CRT	C4-C5-C6	-5.73	116.87	124.91
9	AL	301	BCL	OBB-CAB-C3B	5.72	130.54	120.43
9	AI	102	BCL	OBB-CAB-C3B	5.72	130.54	120.43
9	BE	101	BCL	C4A-NA-C1A	5.71	109.28	106.68
9	AK	102	BCL	OBB-CAB-C3B	5.71	130.53	120.43
9	A5	102	BCL	C4-C3-C5	-5.70	105.33	115.23
9	BI	102	BCL	C4-C3-C5	-5.70	105.33	115.23
14	BU	103	CRT	C5-C6-C7	-5.70	117.28	125.89
9	AV	102	BCL	OBB-CAB-C3B	5.70	130.51	120.43
9	AF	102	BCL	OBB-CAB-C3B	5.70	130.50	120.43
9	BF	102	BCL	OBB-CAB-C3B	5.70	130.50	120.43
9	AJ	101	BCL	OBB-CAB-C3B	5.69	130.50	120.43
9	B0	102	BCL	OBB-CAB-C3B	5.69	130.49	120.43
9	B7	103	BCL	OBB-CAB-C3B	5.68	130.48	120.43
9	BO	102	BCL	OBB-CAB-C3B	5.68	130.48	120.43
9	AZ	101	BCL	C4A-NA-C1A	5.68	109.27	106.68
9	AN	101	BCL	C4-C3-C5	-5.67	105.38	115.23
9	AM	401	BCL	C4A-NA-C1A	5.67	109.27	106.68
9	A5	102	BCL	OBB-CAB-C3B	5.67	130.45	120.43
9	BV	101	BCL	C4-C3-C5	-5.66	105.41	115.23
9	B0	102	BCL	C4-C3-C5	-5.65	105.42	115.23
9	BX	101	BCL	OBB-CAB-C3B	5.65	130.42	120.43
9	A0	102	BCL	OBB-CAB-C3B	5.65	130.42	120.43
9	B5	102	BCL	C4-C3-C5	-5.65	105.42	115.23
9	B3	102	BCL	OBB-CAB-C3B	5.64	130.41	120.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
14	AG	102	CRT	C37-C36-C35	-5.64	116.99	124.91
9	A3	103	BCL	OBB-CAB-C3B	5.64	130.40	120.43
9	BA	101	BCL	OBB-CAB-C3B	5.63	130.38	120.43
9	A0	102	BCL	C4-C3-C5	-5.62	105.47	115.23
9	AU	102	BCL	O2A-CGA-CBA	5.62	128.97	111.83
9	A6	101	BCL	OBB-CAB-C3B	5.62	130.37	120.43
9	AU	102	BCL	C4-C3-C5	-5.62	105.48	115.23
14	BU	103	CRT	C15-C14-C12	-5.61	119.41	127.28
9	AP	101	BCL	OBB-CAB-C3B	5.61	130.35	120.43
9	AI	102	BCL	C4A-NA-C1A	5.61	109.24	106.68
9	B6	101	BCL	OBB-CAB-C3B	5.61	130.34	120.43
9	BM	401	BCL	C4-C3-C5	-5.60	105.50	115.23
9	BW	102	BCL	C4A-NA-C1A	5.60	109.23	106.68
9	AG	101	BCL	OBB-CAB-C3B	5.60	130.33	120.43
9	AI	102	BCL	C4-C3-C5	-5.60	105.51	115.23
9	B4	101	BCL	OBB-CAB-C3B	5.60	130.33	120.43
9	B5	102	BCL	OBB-CAB-C3B	5.60	130.33	120.43
9	BL	303	BCL	O2D-CGD-CBD	5.59	121.01	111.23
14	AN	102	CRT	C37-C36-C35	-5.59	117.06	124.91
14	BA	102	CRT	C37-C36-C35	-5.59	117.06	124.91
9	AN	101	BCL	OBB-CAB-C3B	5.59	130.32	120.43
9	BM	402	BCL	C4-C3-C5	-5.59	105.53	115.23
9	BN	101	BCL	OBB-CAB-C3B	5.59	130.31	120.43
9	BZ	101	BCL	OBB-CAB-C3B	5.59	130.31	120.43
9	BX	101	BCL	CAA-C2A-C3A	-5.58	97.92	113.00
9	AM	402	BCL	OBB-CAB-C3B	5.58	130.29	120.43
9	A3	104	BCL	OBB-CAB-C3B	5.57	130.29	120.43
9	AW	101	BCL	C1-O2A-CGA	-5.57	103.17	116.65
9	BT	101	BCL	C4-C3-C5	-5.56	105.57	115.23
9	AO	102	BCL	OBB-CAB-C3B	5.56	130.26	120.43
9	A9	102	BCL	OBB-CAB-C3B	5.55	130.25	120.43
9	BM	401	BCL	C4A-NA-C1A	5.55	109.21	106.68
9	B9	102	BCL	OBB-CAB-C3B	5.54	130.23	120.43
9	B2	101	BCL	C4A-NA-C1A	5.54	109.21	106.68
9	BM	401	BCL	OBB-CAB-C3B	5.54	130.23	120.43
9	AM	401	BCL	OBB-CAB-C3B	5.53	130.21	120.43
9	AV	102	BCL	C1-O2A-CGA	5.53	130.04	116.65
14	BN	102	CRT	C37-C36-C35	-5.53	117.15	124.91
9	BN	101	BCL	C4-C3-C5	-5.53	105.64	115.23
9	BS	102	BCL	O2A-CGA-CBA	5.52	128.67	111.83
9	B7	103	BCL	C4-C3-C5	-5.51	105.66	115.23
9	B8	101	BCL	OBB-CAB-C3B	5.51	130.18	120.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
14	A0	101	CRT	C10-C9-C7	-5.51	119.55	127.28
9	BB	101	BCL	OBB-CAB-C3B	5.51	130.17	120.43
9	B1	102	BCL	O2D-CGD-CBD	5.51	120.86	111.23
9	AL	303	BCL	C4-C3-C5	-5.49	105.70	115.23
9	AX	101	BCL	OBB-CAB-C3B	5.49	130.13	120.43
9	A6	101	BCL	C1-O2A-CGA	-5.49	103.37	116.65
9	BP	101	BCL	OBB-CAB-C3B	5.48	130.13	120.43
9	A3	103	BCL	O2D-CGD-CBD	5.48	120.81	111.23
9	A8	101	BCL	OBB-CAB-C3B	5.48	130.12	120.43
9	B4	101	BCL	C4-C3-C5	-5.47	105.72	115.23
9	B8	101	BCL	CAA-C2A-C1A	5.47	129.91	111.97
9	A1	102	BCL	O2D-CGD-CBD	5.47	120.79	111.23
9	AR	101	BCL	OBB-CAB-C3B	5.46	130.08	120.43
9	BS	102	BCL	OBB-CAB-C3B	5.45	130.07	120.43
9	BA	101	BCL	C4A-NA-C1A	5.44	109.16	106.68
9	A1	102	BCL	O2A-CGA-O1A	-5.44	110.02	123.63
9	BL	303	BCL	C4-C3-C5	-5.44	105.79	115.23
9	BV	101	BCL	OBB-CAB-C3B	5.43	130.03	120.43
9	AV	102	BCL	C4-C3-C5	-5.42	105.82	115.23
9	BX	101	BCL	O2D-CGD-CBD	5.42	120.70	111.23
9	BA	101	BCL	C4-C3-C5	-5.41	105.84	115.23
9	BM	402	BCL	O2A-CGA-CBA	5.40	128.31	111.83
14	A1	103	CRT	C5-C6-C7	-5.40	117.73	125.89
9	AD	102	BCL	O2D-CGD-CBD	5.39	120.66	111.23
9	AY	102	BCL	O2A-CGA-CBA	5.38	128.24	111.83
9	B9	102	BCL	O2D-CGD-CBD	5.38	120.63	111.23
14	BS	103	CRT	C4-C5-C6	-5.38	117.36	124.91
9	BY	102	BCL	O2D-CGD-CBD	5.36	120.60	111.23
9	BQ	104	BCL	OBB-CAB-C3B	5.36	129.90	120.43
14	B5	103	CRT	C37-C36-C35	-5.36	117.39	124.91
9	B7	103	BCL	O2A-CGA-CBA	5.35	128.15	111.83
9	B3	102	BCL	O2D-CGD-CBD	5.34	120.57	111.23
9	AW	101	BCL	O2A-CGA-O1A	-5.34	110.26	123.63
9	BJ	101	BCL	C4A-NA-C1A	5.33	109.11	106.68
9	AE	101	BCL	CAA-C2A-C1A	5.33	129.45	111.97
9	BI	102	BCL	O2D-CGD-CBD	5.33	120.55	111.23
9	AE	101	BCL	OBB-CAB-C3B	5.32	129.83	120.43
9	AM	402	BCL	O2A-CGA-CBA	5.32	128.04	111.83
14	AX	102	CRT	C37-C36-C35	-5.30	117.47	124.91
9	B7	103	BCL	O2D-CGD-CBD	5.30	120.49	111.23
14	B2	102	CRT	C37-C36-C35	-5.29	117.49	124.91
9	AV	102	BCL	C1-C2-C3	5.28	134.85	126.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
9	BZ	101	BCL	C4-C3-C5	-5.28	106.06	115.23
9	AU	102	BCL	O2D-CGD-CBD	5.27	120.44	111.23
9	AA	101	BCL	C4-C3-C5	-5.27	106.09	115.23
14	A1	103	CRT	C37-C36-C35	-5.25	117.54	124.91
9	B6	101	BCL	O2D-CGD-CBD	5.25	120.41	111.23
9	AI	102	BCL	O2D-CGD-CBD	5.25	120.40	111.23
9	A8	101	BCL	C4A-NA-C1A	5.23	109.06	106.68
14	BU	103	CRT	C37-C36-C35	-5.22	117.58	124.91
9	BK	102	BCL	O2D-CGD-CBD	5.22	120.35	111.23
9	B8	101	BCL	O2D-CGD-CBD	5.21	120.34	111.23
9	BL	301	BCL	C4-C3-C5	-5.19	106.22	115.23
9	AI	102	BCL	O2A-CGA-CBA	5.19	127.64	111.83
14	A5	103	CRT	C37-C36-C35	-5.18	117.64	124.91
9	B8	101	BCL	C4-C3-C5	-5.18	106.25	115.23
14	A7	102	CRT	C4-C5-C6	-5.17	117.66	124.91
9	BL	303	BCL	C4A-NA-C1A	5.17	109.04	106.68
10	BL	302	BPH	C2D-C1D-ND	-5.15	105.71	109.43
9	B4	101	BCL	CAA-C2A-C1A	5.12	128.76	111.97
9	BN	101	BCL	O2D-CGD-CBD	5.12	120.18	111.23
9	B9	102	BCL	O2A-CGA-CBA	5.12	127.43	111.83
9	BO	102	BCL	O2D-CGD-CBD	5.10	120.15	111.23
9	AZ	101	BCL	CAA-C2A-C3A	-5.08	99.26	113.00
9	BQ	103	BCL	O2D-CGD-CBD	5.08	120.10	111.23
9	AL	301	BCL	O2A-CGA-CBA	5.07	127.31	111.83
9	BS	102	BCL	O2A-CGA-O1A	-5.07	110.94	123.63
9	AW	101	BCL	O2D-CGD-CBD	5.04	120.04	111.23
13	AM	405	MQ8	C11-C3-C4	-5.04	113.27	118.58
14	B0	101	CRT	C4-C5-C6	-5.03	117.84	124.91
9	BM	401	BCL	O2D-CGD-CBD	5.02	120.01	111.23
14	BS	103	CRT	C37-C36-C35	-5.01	117.87	124.91
9	A3	104	BCL	CAA-C2A-C1A	5.01	128.40	111.97
9	A5	102	BCL	O2D-CGD-CBD	5.01	119.98	111.23
9	BA	101	BCL	O2D-CGD-CBD	5.00	119.97	111.23
9	A3	104	BCL	O2A-CGA-O1A	-5.00	111.13	123.63
9	BS	102	BCL	O2D-CGD-CBD	5.00	119.96	111.23
9	AI	102	BCL	O2A-CGA-O1A	-4.98	111.17	123.63
10	AL	302	BPH	C2D-C1D-ND	-4.98	105.83	109.43
9	AM	401	BCL	O2D-CGD-CBD	4.97	119.92	111.23
9	BO	102	BCL	O2A-CGA-CBA	4.97	126.98	111.83
9	A0	102	BCL	C1-O2A-CGA	4.96	128.66	116.65
9	B9	102	BCL	O2A-CGA-O1A	-4.96	111.23	123.63
9	AM	402	BCL	O2D-CGD-CBD	4.95	119.89	111.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
9	B5	102	BCL	O2D-CGD-CBD	4.95	119.89	111.23
9	AX	101	BCL	CAA-C2A-C1A	4.95	128.20	111.97
9	AU	102	BCL	O2A-CGA-O1A	-4.94	111.27	123.63
9	A7	103	BCL	O2A-CGA-O1A	-4.94	111.28	123.63
9	BI	102	BCL	O2A-CGA-CBA	4.94	126.89	111.83
9	AK	102	BCL	O2D-CGD-CBD	4.93	119.86	111.23
9	BG	101	BCL	C1C-NC-C4C	4.93	108.93	106.68
9	BF	102	BCL	O2A-CGA-CBA	4.93	126.86	111.83
9	BW	102	BCL	O2A-CGA-CBA	4.93	126.86	111.83
9	AL	303	BCL	C4A-NA-C1A	4.92	108.92	106.68
9	AP	101	BCL	CAA-C2A-C1A	4.92	128.10	111.97
9	BB	101	BCL	O2D-CGD-CBD	4.92	119.83	111.23
9	AQ	102	BCL	O2D-CGD-CBD	4.91	119.82	111.23
14	AX	102	CRT	C4-C5-C6	-4.91	118.02	124.91
9	AL	301	BCL	C4-C3-C5	-4.91	106.71	115.23
9	B8	101	BCL	C1-O2A-CGA	4.90	128.51	116.65
9	BM	402	BCL	O2A-CGA-O1A	-4.89	111.40	123.63
9	BL	301	BCL	O2A-CGA-CBA	4.88	126.72	111.83
9	AO	102	BCL	O2D-CGD-CBD	4.88	119.76	111.23
9	BM	402	BCL	O2D-CGD-CBD	4.88	119.76	111.23
9	AW	101	BCL	O2A-CGA-CBA	4.87	126.68	111.83
9	B7	103	BCL	O2A-CGA-O1A	-4.86	111.46	123.63
9	BE	101	BCL	C4-C3-C5	-4.86	106.80	115.23
9	AW	101	BCL	C4A-NA-C1A	4.86	108.89	106.68
9	AK	102	BCL	O2A-CGA-CBA	4.85	126.62	111.83
9	AT	101	BCL	O2D-CGD-CBD	4.85	119.70	111.23
9	BG	101	BCL	O2D-CGD-CBD	4.85	119.70	111.23
9	A8	101	BCL	O2A-CGA-O1A	-4.84	111.51	123.63
9	A9	102	BCL	O2A-CGA-O1A	-4.84	111.52	123.63
9	BF	102	BCL	O2D-CGD-CBD	4.84	119.69	111.23
9	AL	301	BCL	O2A-CGA-O1A	-4.83	111.55	123.63
9	AY	102	BCL	O2A-CGA-O1A	-4.82	111.57	123.63
14	A2	102	CRT	C4-C5-C6	-4.82	118.15	124.91
9	A0	102	BCL	O2A-CGA-O1A	-4.82	111.58	123.63
9	BU	102	BCL	O2A-CGA-CBA	4.82	126.52	111.83
9	BW	102	BCL	O2A-CGA-O1A	-4.80	111.61	123.63
9	A3	103	BCL	O2A-CGA-CBA	4.79	126.44	111.83
14	BO	103	CRT	C37-C36-C35	-4.79	118.19	124.91
9	BO	102	BCL	O2A-CGA-O1A	-4.79	111.65	123.63
14	A2	102	CRT	C37-C36-C35	-4.79	118.19	124.91
9	AM	402	BCL	O2A-CGA-O1A	-4.78	111.66	123.63
14	AW	102	CRT	C37-C36-C35	-4.78	118.21	124.91

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
9	A7	103	BCL	O2A-CGA-CBA	4.77	126.38	111.83
9	A5	102	BCL	O2A-CGA-O1A	-4.77	111.71	123.63
9	AY	102	BCL	CBA-CAA-C2A	4.76	127.96	113.79
14	AP	102	CRT	C37-C36-C35	-4.76	118.23	124.91
9	AB	101	BCL	CAA-C2A-C1A	4.76	127.57	111.97
14	BU	103	CRT	C20-C19-C17	-4.76	120.61	127.28
9	AT	101	BCL	C4-C3-C5	-4.76	106.97	115.23
9	AA	101	BCL	O2A-CGA-CBA	4.74	126.28	111.83
9	B4	101	BCL	O2D-CGD-CBD	4.73	119.50	111.23
14	BG	102	CRT	C4-C5-C6	-4.72	118.28	124.91
9	AS	103	BCL	O2D-CGD-CBD	4.72	119.48	111.23
9	AN	101	BCL	O2D-CGD-CBD	4.71	119.46	111.23
9	AO	102	BCL	C4A-NA-C1A	4.70	108.82	106.68
9	AN	101	BCL	CAA-C2A-C1A	4.69	127.35	111.97
9	A3	103	BCL	O2A-CGA-O1A	-4.69	111.90	123.63
9	A8	101	BCL	O2A-CGA-CBA	4.69	126.13	111.83
9	A5	102	BCL	O2A-CGA-CBA	4.68	126.11	111.83
9	AK	102	BCL	O2A-CGA-O1A	-4.68	111.92	123.63
9	AS	103	BCL	O2A-CGA-O1A	-4.68	111.93	123.63
9	AF	102	BCL	O2D-CGD-CBD	4.67	119.40	111.23
9	BZ	101	BCL	CAA-C2A-C3A	-4.67	100.38	113.00
9	AR	101	BCL	CAA-C2A-C3A	-4.67	100.39	113.00
9	BI	102	BCL	O2A-CGA-O1A	-4.66	111.97	123.63
9	A8	101	BCL	C4-C3-C5	-4.65	107.15	115.23
9	B1	102	BCL	O2A-CGA-CBA	4.65	126.02	111.83
9	A6	101	BCL	O2A-CGA-O1A	-4.65	112.00	123.63
9	BL	301	BCL	O2A-CGA-O1A	-4.65	112.00	123.63
9	AN	101	BCL	CAA-C2A-C3A	-4.64	100.45	113.00
9	AS	103	BCL	O2A-CGA-CBA	4.64	125.97	111.83
14	B1	103	CRT	C4-C5-C6	-4.63	118.41	124.91
9	A0	102	BCL	CAA-C2A-C1A	4.63	127.15	111.97
9	AZ	101	BCL	CAA-C2A-C1A	4.62	127.13	111.97
9	BF	102	BCL	O2A-CGA-O1A	-4.61	112.09	123.63
9	BU	102	BCL	O2A-CGA-O1A	-4.61	112.09	123.63
9	BY	102	BCL	O2A-CGA-CBA	4.61	125.89	111.83
14	AX	102	CRT	C20-C19-C17	-4.60	120.83	127.28
9	B6	101	BCL	CAA-C2A-C1A	4.60	127.05	111.97
9	A3	103	BCL	C4-C3-C5	-4.59	107.25	115.23
9	A3	104	BCL	O2A-CGA-CBA	4.58	125.81	111.83
9	AA	101	BCL	O2A-CGA-O1A	-4.58	112.17	123.63
9	B2	101	BCL	O2D-CGD-CBD	4.58	119.23	111.23
9	BB	101	BCL	CAA-C2A-C1A	4.57	126.96	111.97

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
15	AM	407	PEF	C2-O2-C10	4.57	125.92	117.85
9	B1	102	BCL	O2A-CGA-O1A	-4.57	112.19	123.63
15	AH	301	PEF	C2-O2-C10	4.57	125.91	117.85
9	BL	303	BCL	C1C-NC-C4C	4.56	108.76	106.68
9	AB	101	BCL	CAA-C2A-C3A	-4.56	100.67	113.00
9	BB	101	BCL	CAA-C2A-C3A	-4.56	100.67	113.00
9	AB	101	BCL	C1-O2A-CGA	4.56	127.68	116.65
15	BM	407	PEF	C2-O2-C10	4.55	125.88	117.85
9	AV	102	BCL	O2D-CGD-CBD	4.54	119.16	111.23
9	A8	101	BCL	CAA-C2A-C1A	4.53	126.83	111.97
9	AP	101	BCL	O2D-CGD-CBD	4.53	119.14	111.23
7	AC	501	HEM	C3B-C4B-NB	4.52	112.72	109.47
9	BA	101	BCL	O2A-CGA-CBA	4.51	125.59	111.83
9	BQ	103	BCL	O2A-CGA-CBA	4.50	125.56	111.83
9	BK	102	BCL	O2A-CGA-O1A	-4.50	112.38	123.63
9	BQ	103	BCL	O2A-CGA-O1A	-4.49	112.39	123.63
7	AC	504	HEM	C3B-C4B-NB	4.48	112.69	109.47
9	A0	102	BCL	O2A-CGA-CBA	4.48	125.49	111.83
9	B4	101	BCL	CAA-C2A-C3A	-4.47	100.92	113.00
9	BT	101	BCL	O2D-CGD-CBD	4.46	119.03	111.23
15	AS	101	PEF	O3-C30-C31	4.46	125.43	111.83
9	BO	102	BCL	C5-C3-C2	4.46	131.17	121.17
14	AT	102	CRT	C37-C36-C35	-4.45	118.66	124.91
9	BM	401	BCL	O2A-CGA-CBA	4.45	125.40	111.83
9	BK	102	BCL	O2A-CGA-CBA	4.45	125.40	111.83
15	AM	409	PEF	O3-C30-C31	4.45	125.40	111.83
9	AL	303	BCL	O2A-CGA-O1A	-4.45	112.51	123.63
9	BM	401	BCL	C1C-NC-C4C	4.45	108.71	106.68
9	A2	101	BCL	O2A-CGA-O1A	-4.44	112.52	123.63
15	BQ	101	PEF	O3-C30-C31	4.44	125.36	111.83
9	BM	401	BCL	O2A-CGA-O1A	-4.44	112.53	123.63
9	BD	102	BCL	O2A-CGA-O1A	-4.43	112.54	123.63
9	AQ	102	BCL	O2A-CGA-CBA	4.43	125.35	111.83
14	AR	102	CRT	C37-C36-C35	-4.43	118.69	124.91
9	AL	303	BCL	C1C-NC-C4C	4.43	108.70	106.68
14	A1	103	CRT	C14-C15-C16	-4.43	110.36	123.20
9	AL	303	BCL	O2A-CGA-CBA	4.43	125.34	111.83
9	AQ	102	BCL	O2A-CGA-O1A	-4.42	112.56	123.63
9	BD	102	BCL	O2A-CGA-CBA	4.41	125.30	111.83
14	AG	102	CRT	C4-C5-C6	-4.41	118.72	124.91
9	BD	102	BCL	O2D-CGD-CBD	4.41	118.94	111.23
9	AO	102	BCL	O2A-CGA-O1A	-4.41	112.60	123.63

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
9	A6	101	BCL	CAA-C2A-C1A	4.41	126.42	111.97
9	BA	101	BCL	O2A-CGA-O1A	-4.41	112.61	123.63
9	AJ	101	BCL	CAA-C2A-C1A	4.40	126.40	111.97
9	BG	101	BCL	CAA-C2A-C3A	-4.40	101.12	113.00
14	AR	102	CRT	C5-C6-C7	-4.39	119.25	125.89
9	BL	303	BCL	O2A-CGA-O1A	-4.39	112.65	123.63
7	AC	504	HEM	C4C-C3C-C2C	-4.39	103.01	106.81
9	AD	102	BCL	O2A-CGA-O1A	-4.38	112.67	123.63
9	AM	401	BCL	O2A-CGA-O1A	-4.38	112.68	123.63
9	A9	102	BCL	O2A-CGA-CBA	4.37	125.17	111.83
9	AZ	101	BCL	C5-C3-C2	4.37	130.98	121.17
9	AG	101	BCL	O2D-CGD-CBD	4.37	118.87	111.23
9	A9	102	BCL	O2D-CGD-CBD	4.37	118.86	111.23
9	AD	102	BCL	O2A-CGA-CBA	4.36	125.13	111.83
9	BZ	101	BCL	CAA-C2A-C1A	4.36	126.25	111.97
9	A6	101	BCL	O2A-CGA-CBA	4.35	125.11	111.83
9	BL	303	BCL	O2A-CGA-CBA	4.35	125.11	111.83
9	AY	102	BCL	O2D-CGD-CBD	4.35	118.83	111.23
9	BY	102	BCL	O2A-CGA-O1A	-4.34	112.78	123.63
7	BC	502	HEM	C4C-C3C-C2C	-4.34	103.05	106.81
9	B5	102	BCL	O2A-CGA-CBA	4.33	125.04	111.83
9	BP	101	BCL	O2D-CGD-CBD	4.32	118.79	111.23
9	AM	401	BCL	O2A-CGA-CBA	4.32	125.00	111.83
9	BL	301	BCL	C1C-NC-C4C	4.31	108.64	106.68
9	A8	101	BCL	CBA-CAA-C2A	4.30	126.59	113.79
13	BM	405	MQ8	C12-C11-C3	4.30	122.67	112.08
9	BJ	101	BCL	O2D-CGD-CBD	4.29	118.72	111.23
9	AO	102	BCL	O2A-CGA-CBA	4.28	124.88	111.83
14	BM	406	CRT	C37-C36-C35	-4.27	118.91	124.91
9	BQ	103	BCL	C1C-NC-C4C	4.27	108.63	106.68
9	AR	101	BCL	CAA-C2A-C1A	4.27	125.96	111.97
7	AC	501	HEM	C4C-C3C-C2C	-4.26	103.12	106.81
14	A1	103	CRT	C6-C7-C9	4.26	125.71	119.01
7	BC	504	HEM	C4C-C3C-C2C	-4.25	103.13	106.81
9	AT	101	BCL	CAA-C2A-C1A	4.25	125.89	111.97
9	BG	101	BCL	C5-C3-C2	4.24	130.69	121.17
9	A0	102	BCL	CAA-C2A-C3A	-4.24	101.54	113.00
9	BV	101	BCL	O2D-CGD-CBD	4.24	118.64	111.23
9	AZ	101	BCL	O2A-CGA-O1A	-4.23	113.04	123.63
14	B2	102	CRT	C1-C4-C5	4.23	123.97	112.98
9	B0	102	BCL	CAA-C2A-C1A	4.23	125.84	111.97
9	AZ	101	BCL	O2A-CGA-CBA	4.21	124.66	111.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
9	A7	103	BCL	O2D-CGD-CBD	4.20	118.57	111.23
9	BZ	101	BCL	O2D-CGD-CBD	4.20	118.57	111.23
9	BQ	104	BCL	CAA-C2A-C3A	-4.19	101.67	113.00
7	AC	502	HEM	C4C-C3C-C2C	-4.19	103.18	106.81
15	AM	407	PEF	O2-C10-C11	4.18	118.55	111.09
9	A8	101	BCL	O2D-CGD-CBD	4.18	118.54	111.23
9	BX	101	BCL	CAA-C2A-C1A	4.18	125.67	111.97
9	AX	101	BCL	O2A-CGA-O1A	-4.17	113.19	123.63
15	AH	301	PEF	O2-C10-C11	4.17	118.53	111.09
13	AM	405	MQ8	C12-C11-C3	4.17	122.35	112.08
9	AE	101	BCL	C1-O2A-CGA	4.17	126.74	116.65
9	AJ	101	BCL	CAA-C2A-C3A	-4.16	101.76	113.00
9	AL	301	BCL	C1C-NC-C4C	4.15	108.57	106.68
15	BM	407	PEF	O2-C10-C11	4.14	118.48	111.09
9	AS	103	BCL	C5-C3-C2	4.14	130.46	121.17
14	AS	104	CRT	C21-C22-C23	-4.12	121.49	127.28
14	AB	102	CRT	C21-C20-C19	-4.12	115.08	123.52
9	AX	101	BCL	C5-C3-C2	4.12	130.41	121.17
9	AK	102	BCL	C5-C3-C2	4.11	130.40	121.17
7	BC	503	HEM	C4C-C3C-C2C	-4.11	103.25	106.81
9	B0	102	BCL	O2D-CGD-CBD	4.10	118.41	111.23
9	BP	101	BCL	CAA-C2A-C3A	-4.10	101.91	113.00
9	A2	101	BCL	CAA-C2A-C1A	4.10	125.42	111.97
9	A3	104	BCL	CAA-C2A-C3A	-4.10	101.92	113.00
9	AB	101	BCL	O2D-CGD-CBD	4.10	118.39	111.23
9	AG	101	BCL	C5-C3-C2	4.08	130.32	121.17
9	B3	102	BCL	O2A-CGA-CBA	4.07	124.25	111.83
9	B5	102	BCL	O2A-CGA-O1A	-4.07	113.44	123.63
9	AB	101	BCL	C5-C3-C2	4.06	130.28	121.17
7	AC	503	HEM	C4C-C3C-C2C	-4.05	103.30	106.81
9	BO	102	BCL	C1C-NC-C4C	4.05	108.53	106.68
9	B8	101	BCL	CBA-CAA-C2A	4.05	125.84	113.79
9	AD	102	BCL	C5-C3-C2	4.05	130.25	121.17
9	AU	102	BCL	C12-C11-C10	-4.04	95.16	113.28
9	A3	103	BCL	C5-C3-C2	4.04	130.23	121.17
11	AL	304	UQ8	C42-C41-C39	4.03	126.55	113.19
15	AM	409	PEF	O3-C30-O5	-4.03	113.54	123.63
14	A0	101	CRT	C26-C27-C28	-4.03	121.62	127.28
9	AK	102	BCL	C4D-CHA-C1A	4.02	126.04	121.24
15	AS	101	PEF	O3-C30-O5	-4.02	113.58	123.63
15	BQ	101	PEF	O3-C30-O5	-4.02	113.58	123.63
9	B8	101	BCL	CAA-C2A-C3A	-4.02	102.15	113.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
9	A8	101	BCL	C6-C5-C3	4.01	123.23	113.47
15	AS	101	PEF	O3-C3-C2	4.00	119.94	108.40
9	AE	101	BCL	CBA-CAA-C2A	4.00	125.70	113.79
15	BQ	101	PEF	O3-C3-C2	4.00	119.93	108.40
9	A6	101	BCL	CAA-C2A-C3A	-4.00	102.19	113.00
9	AE	101	BCL	O2D-CGD-CBD	4.00	118.22	111.23
10	BM	403	BPH	C1A-C2A-C3A	-4.00	99.03	102.84
9	BW	102	BCL	C5-C3-C2	4.00	130.14	121.17
9	A5	102	BCL	C4D-CHA-C1A	3.99	126.01	121.24
9	AQ	102	BCL	C4D-CHA-C1A	3.99	126.01	121.24
9	BB	101	BCL	C5-C3-C2	3.99	130.13	121.17
9	BT	101	BCL	CAA-C2A-C1A	3.99	125.06	111.97
15	AH	301	PEF	O3-C3-C2	3.99	119.89	108.40
14	AP	102	CRT	C4-C5-C6	-3.98	119.32	124.91
11	AL	304	UQ8	C35-C34-C36	3.98	122.14	115.23
9	AU	102	BCL	C11-C10-C8	-3.98	102.74	115.97
15	AM	407	PEF	O3-C3-C2	3.98	119.86	108.40
9	A1	102	BCL	C4D-CHA-C1A	3.98	125.99	121.24
9	A6	101	BCL	O2D-CGD-CBD	3.97	118.18	111.23
15	BM	407	PEF	O3-C3-C2	3.97	119.83	108.40
9	A3	103	BCL	C4D-CHA-C1A	3.97	125.97	121.24
9	A8	101	BCL	C1C-NC-C4C	3.95	108.48	106.68
9	A3	104	BCL	C5-C3-C2	3.95	130.04	121.17
9	BU	102	BCL	C5-C3-C2	3.95	130.04	121.17
9	B3	102	BCL	C5-C3-C2	3.95	130.04	121.17
9	AP	101	BCL	C5-C3-C2	3.95	130.03	121.17
10	AL	302	BPH	C1-C2-C3	3.94	132.66	126.20
14	BP	102	CRT	C37-C36-C35	-3.94	119.38	124.91
9	AX	101	BCL	C1-O2A-CGA	-3.94	107.11	116.65
9	BM	402	BCL	C1C-NC-C4C	3.94	108.47	106.68
7	AC	502	HEM	C3B-C4B-NB	3.94	112.30	109.47
9	BB	101	BCL	C1C-NC-C4C	3.93	108.47	106.68
9	BN	101	BCL	C5-C3-C2	3.93	130.00	121.17
9	AZ	101	BCL	O2D-CGD-CBD	3.93	118.09	111.23
9	AV	102	BCL	C5-C3-C2	3.92	129.98	121.17
9	AX	101	BCL	O2A-CGA-CBA	3.92	123.79	111.83
9	AX	101	BCL	O2D-CGD-CBD	3.91	118.06	111.23
14	AS	104	CRT	C10-C9-C7	-3.89	121.83	127.28
9	AM	402	BCL	C5-C3-C2	3.89	129.89	121.17
9	AG	101	BCL	CAA-C2A-C3A	-3.89	102.50	113.00
9	AJ	101	BCL	C1-O2A-CGA	3.88	126.04	116.65
14	A1	103	CRT	C21-C22-C23	-3.87	121.84	127.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
9	A9	102	BCL	C4D-CHA-C1A	3.87	125.86	121.24
7	BC	501	HEM	C3B-C4B-NB	3.87	112.25	109.47
9	BD	102	BCL	C5-C3-C2	3.87	129.86	121.17
9	B3	102	BCL	O2A-CGA-O1A	-3.87	113.95	123.63
9	AR	101	BCL	O2D-CGD-CBD	3.87	117.99	111.23
9	A9	102	BCL	CAA-C2A-C1A	3.87	124.65	111.97
14	A0	101	CRT	C31-C32-C33	-3.87	121.86	127.28
14	BO	103	CRT	C21-C22-C23	-3.86	121.86	127.28
14	BB	102	CRT	C4-C5-C6	-3.86	119.50	124.91
9	AU	102	BCL	C9-C8-C10	-3.86	97.52	111.27
9	AY	102	BCL	C5-C3-C2	3.85	129.82	121.17
14	AM	406	CRT	C31-C32-C33	-3.85	121.88	127.28
9	BZ	101	BCL	O2A-CGA-O1A	-3.85	114.00	123.63
14	A1	103	CRT	C8-C7-C9	-3.85	116.58	122.82
9	BM	401	BCL	CAA-C2A-C3A	-3.85	102.61	113.00
9	AF	102	BCL	O2A-CGA-O1A	-3.85	114.01	123.63
9	BJ	101	BCL	CAA-C2A-C3A	-3.84	102.61	113.00
7	BC	502	HEM	C3B-C4B-NB	3.84	112.23	109.47
11	BL	304	UQ8	C42-C41-C39	3.84	125.90	113.19
9	AX	101	BCL	CAA-C2A-C3A	-3.83	102.64	113.00
14	BO	103	CRT	C10-C9-C7	-3.83	121.91	127.28
14	AB	102	CRT	C37-C36-C35	-3.83	119.54	124.91
9	B1	102	BCL	C5-C3-C2	3.83	129.75	121.17
9	B6	101	BCL	CAA-C2A-C3A	-3.82	102.67	113.00
9	BP	101	BCL	CAA-C2A-C1A	3.82	124.49	111.97
9	BF	102	BCL	C3A-C2A-C1A	3.81	107.05	101.34
14	AB	102	CRT	C21-C22-C23	-3.81	121.93	127.28
9	A7	103	BCL	C4D-CHA-C1A	3.81	125.79	121.24
9	BL	301	BCL	C5-C3-C2	3.81	129.72	121.17
7	BC	504	HEM	C3B-C4B-NB	3.80	112.20	109.47
14	BB	102	CRT	C32-C31-C30	-3.80	112.20	123.20
9	BE	101	BCL	O2D-CGD-CBD	3.79	117.86	111.23
14	A0	101	CRT	C21-C22-C23	-3.79	121.96	127.28
14	BB	102	CRT	C21-C20-C19	-3.79	115.76	123.52
9	AM	401	BCL	C1C-NC-C4C	3.79	108.41	106.68
9	BV	101	BCL	CAA-C2A-C1A	3.78	124.38	111.97
9	BF	102	BCL	C5-C3-C2	3.78	129.65	121.17
9	A2	101	BCL	CAA-C2A-C3A	-3.77	102.81	113.00
9	A2	101	BCL	O2A-CGA-CBA	3.77	123.32	111.83
9	B8	101	BCL	CAA-CBA-CGA	-3.77	102.51	113.21
14	A0	101	CRT	C37-C36-C35	-3.76	119.63	124.91
9	BK	102	BCL	C5-C3-C2	3.76	129.60	121.17

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
9	A3	104	BCL	C6-C5-C3	3.76	122.62	113.47
9	BS	102	BCL	C5-C3-C2	3.75	129.59	121.17
7	BC	501	HEM	C4C-C3C-C2C	-3.75	103.56	106.81
14	BB	102	CRT	C21-C22-C23	-3.75	122.02	127.28
9	B6	101	BCL	C5-C3-C2	3.75	129.58	121.17
9	A2	101	BCL	C5-C3-C2	3.75	129.57	121.17
9	AW	101	BCL	C5-C3-C2	3.74	129.57	121.17
9	BO	102	BCL	C6-C5-C3	3.74	122.59	113.47
9	BQ	103	BCL	C5-C3-C2	3.74	129.57	121.17
9	AO	102	BCL	C5-C3-C2	3.74	129.56	121.17
9	BX	101	BCL	C5-C3-C2	3.73	129.55	121.17
9	BN	101	BCL	CAA-C2A-C3A	-3.73	102.91	113.00
9	AR	101	BCL	O2A-CGA-CBA	3.72	123.18	111.83
9	AE	101	BCL	C6-C5-C3	3.71	122.50	113.47
14	BG	102	CRT	C37-C36-C35	-3.71	119.70	124.91
9	A0	102	BCL	O2D-CGD-CBD	3.70	117.70	111.23
9	BJ	101	BCL	C5-C3-C2	3.70	129.48	121.17
9	B0	102	BCL	CAA-C2A-C3A	-3.70	103.00	113.00
9	B2	101	BCL	CAA-C2A-C1A	3.70	124.09	111.97
9	B2	101	BCL	C5-C3-C2	3.69	129.46	121.17
14	BW	103	CRT	C37-C36-C35	-3.69	119.73	124.91
9	B5	102	BCL	C4D-CHA-C1A	3.69	125.65	121.24
9	AF	102	BCL	C5-C3-C2	3.69	129.45	121.17
9	BQ	104	BCL	O2D-CGD-CBD	3.69	117.68	111.23
9	BP	101	BCL	C5-C3-C2	3.69	129.44	121.17
9	AU	102	BCL	C4D-CHA-C1A	3.69	125.64	121.24
10	BL	302	BPH	C1-C2-C3	3.68	132.23	126.20
9	AM	401	BCL	CAA-C2A-C3A	-3.68	103.04	113.00
14	BA	102	CRT	C4-C5-C6	-3.68	119.75	124.91
9	BD	102	BCL	C3A-C2A-C1A	3.68	106.84	101.34
14	AS	104	CRT	C35-C33-C32	-3.67	113.23	119.01
9	BY	102	BCL	C5-C3-C2	3.67	129.41	121.17
9	B2	101	BCL	C1-O2A-CGA	3.67	125.54	116.65
9	AW	101	BCL	CAA-C2A-C1A	3.67	124.00	111.97
9	AM	402	BCL	C1C-NC-C4C	3.67	108.35	106.68
9	BI	102	BCL	C5-C3-C2	3.67	129.40	121.17
9	AR	101	BCL	C5-C3-C2	3.66	129.39	121.17
9	AF	102	BCL	O2A-CGA-CBA	3.66	123.00	111.83
9	BX	101	BCL	O2D-CGD-O1D	-3.65	116.73	123.85
9	AW	101	BCL	CBA-CAA-C2A	3.65	124.67	113.79
9	AQ	102	BCL	C5-C3-C2	3.65	129.36	121.17
14	A0	101	CRT	C4-C5-C6	-3.65	119.79	124.91

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
9	BZ	101	BCL	C5-C3-C2	3.65	129.35	121.17
14	BN	102	CRT	C32-C31-C30	-3.64	112.65	123.20
9	BF	102	BCL	C4D-CHA-C1A	3.64	125.59	121.24
9	A7	103	BCL	C5-C3-C2	3.64	129.34	121.17
14	AR	102	CRT	C21-C22-C23	-3.64	122.17	127.28
9	AP	101	BCL	C2C-C3C-C4C	3.63	106.78	101.34
14	AR	102	CRT	C10-C9-C7	-3.62	122.20	127.28
9	B0	102	BCL	O2A-CGA-O1A	-3.62	114.57	123.63
9	AJ	101	BCL	C5-C3-C2	3.62	129.29	121.17
14	A7	102	CRT	C37-C36-C35	-3.62	119.83	124.91
9	B9	102	BCL	C5-C3-C2	3.62	129.29	121.17
9	BN	101	BCL	CAA-C2A-C1A	3.61	123.79	111.97
9	BL	303	BCL	CAA-C2A-C3A	-3.60	103.26	113.00
9	B8	101	BCL	C1C-NC-C4C	3.60	108.32	106.68
9	B5	102	BCL	C5-C3-C2	3.60	129.24	121.17
9	AL	301	BCL	C5-C3-C2	3.59	129.24	121.17
9	BP	101	BCL	CAB-C3B-C2B	3.59	135.21	127.74
14	AN	102	CRT	C21-C22-C23	-3.59	122.24	127.28
9	BQ	104	BCL	O2A-CGA-CBA	3.59	122.79	111.83
9	BZ	101	BCL	C1C-NC-C4C	3.58	108.31	106.68
9	BD	102	BCL	C4D-CHA-C1A	3.58	125.52	121.24
10	BM	403	BPH	OBB-CAB-C3B	3.58	125.97	119.99
9	A3	104	BCL	O2D-CGD-CBD	3.58	117.48	111.23
9	AI	102	BCL	C5-C3-C2	3.57	129.19	121.17
9	BF	102	BCL	C6-C5-C3	3.57	122.17	113.47
9	BQ	103	BCL	C4A-NA-C1A	3.57	108.31	106.68
9	A6	101	BCL	C5-C3-C2	3.55	129.14	121.17
9	A2	101	BCL	OBB-CAB-CBB	-3.55	111.83	119.77
14	B0	101	CRT	C31-C32-C33	-3.54	122.31	127.28
9	AY	102	BCL	C1-O2A-CGA	3.54	125.22	116.65
9	A7	103	BCL	C1-O2A-CGA	-3.54	108.08	116.65
14	A1	103	CRT	C13-C12-C11	3.53	123.49	118.09
14	BB	102	CRT	C30-C28-C27	-3.53	113.46	119.01
9	BQ	104	BCL	CAB-C3B-C2B	3.52	135.06	127.74
14	AS	104	CRT	C34-C33-C35	3.52	123.46	118.09
9	BZ	101	BCL	O2A-CGA-CBA	3.51	122.55	111.83
9	AE	101	BCL	C5-C3-C2	3.51	129.06	121.17
7	AC	504	HEM	C2B-C1B-NB	3.50	113.87	109.84
9	BN	101	BCL	C1-O2A-CGA	3.50	125.12	116.65
14	A1	103	CRT	C20-C19-C17	-3.50	122.37	127.28
7	BC	503	HEM	C2B-C1B-NB	3.50	113.86	109.84
9	A1	102	BCL	C5-C3-C2	3.49	129.01	121.17

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
9	A5	102	BCL	C5-C3-C2	3.49	129.01	121.17
9	BF	102	BCL	OBB-CAB-CBB	-3.49	111.96	119.77
9	AY	102	BCL	C4D-CHA-C1A	3.48	125.40	121.24
9	BD	102	BCL	O2D-CGD-O1D	-3.48	117.07	123.85
9	AU	102	BCL	C10-C8-C7	3.48	129.71	112.07
9	AJ	101	BCL	O2D-CGD-CBD	3.48	117.32	111.23
9	AM	401	BCL	C5-C3-C2	3.48	128.97	121.17
9	BE	101	BCL	C2A-C1A-CHA	3.47	129.89	123.87
11	AL	304	UQ8	C20-C19-C21	3.47	121.25	115.23
9	AU	102	BCL	CAA-C2A-C1A	3.46	123.31	111.97
9	AP	101	BCL	CAA-C2A-C3A	-3.46	103.65	113.00
14	B2	102	CRT	C8-C7-C6	3.46	123.37	118.09
9	A1	102	BCL	OBB-CAB-CBB	-3.45	112.04	119.77
9	BS	102	BCL	CAB-C3B-C2B	3.45	134.90	127.74
9	AR	101	BCL	O2A-CGA-O1A	-3.44	115.01	123.63
9	BV	101	BCL	C5-C3-C2	3.44	128.89	121.17
7	AC	501	HEM	C2B-C1B-NB	3.44	113.79	109.84
9	BW	102	BCL	O2D-CGD-O1D	-3.44	117.16	123.85
9	BP	101	BCL	C4D-CHA-C1A	3.44	125.34	121.24
9	BQ	104	BCL	C5-C3-C2	3.43	128.87	121.17
14	B0	101	CRT	C34-C33-C35	3.43	123.33	118.09
13	BM	405	MQ8	C11-C3-C4	-3.43	114.96	118.58
14	BG	102	CRT	C32-C31-C30	-3.43	113.26	123.20
9	BA	101	BCL	C5-C3-C2	3.43	128.86	121.17
9	AL	303	BCL	CAA-C2A-C3A	-3.43	103.74	113.00
14	A2	102	CRT	C3-C1-C2	-3.42	104.17	110.43
9	BE	101	BCL	C2C-C3C-C4C	3.42	106.46	101.34
9	BM	401	BCL	C5-C3-C2	3.42	128.84	121.17
9	AT	101	BCL	C5-C3-C2	3.42	128.84	121.17
9	AU	102	BCL	C5-C3-C2	3.41	128.83	121.17
15	AS	101	PEF	C2-O2-C10	3.41	125.96	117.80
9	B4	101	BCL	C5-C3-C2	3.41	128.82	121.17
9	AD	102	BCL	OBB-CAB-CBB	-3.41	112.14	119.77
9	A3	104	BCL	C1-O2A-CGA	3.41	124.90	116.65
15	BQ	101	PEF	C2-O2-C10	3.41	125.95	117.80
9	BV	101	BCL	C1C-NC-C4C	3.41	108.23	106.68
7	BC	501	HEM	C2B-C1B-NB	3.41	113.75	109.84
14	BU	103	CRT	C10-C11-C12	-3.40	117.03	126.36
14	BO	103	CRT	C5-C6-C7	-3.40	120.75	125.89
9	AF	102	BCL	C1C-NC-C4C	3.40	108.23	106.68
9	AW	101	BCL	C4D-CHA-C1A	3.40	125.30	121.24
9	B2	101	BCL	CAA-CBA-CGA	-3.40	103.55	113.21

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
9	BT	101	BCL	C5-C3-C2	3.40	128.79	121.17
9	AL	303	BCL	C5-C3-C2	3.40	128.79	121.17
9	BT	101	BCL	CAA-C2A-C3A	-3.39	103.83	113.00
15	AM	409	PEF	C2-O2-C10	3.39	125.90	117.80
14	AS	104	CRT	C21-C20-C19	-3.39	116.59	123.52
7	BC	504	HEM	C2B-C1B-NB	3.38	113.73	109.84
14	AW	102	CRT	C3-C1-C2	-3.38	104.25	110.43
9	B0	102	BCL	OBB-CAB-CBB	-3.38	112.20	119.77
9	BP	101	BCL	C1-O2A-CGA	3.38	124.83	116.65
9	B2	101	BCL	CAA-C2A-C3A	-3.38	103.87	113.00
9	B7	103	BCL	C4D-CHA-C1A	3.37	125.27	121.24
9	BT	101	BCL	O2A-CGA-O1A	-3.37	115.20	123.63
7	AC	503	HEM	C2B-C1B-NB	3.37	113.71	109.84
9	A6	101	BCL	C1C-NC-C4C	3.36	108.21	106.68
9	BG	101	BCL	C6-C5-C3	3.36	121.65	113.47
9	B9	102	BCL	C4D-CHA-C1A	3.36	125.25	121.24
9	AT	101	BCL	C2C-C3C-C4C	3.36	106.37	101.34
14	B7	102	CRT	C1-C4-C5	3.35	121.68	112.98
14	B0	101	CRT	C20-C19-C17	-3.35	122.58	127.28
9	BL	303	BCL	C5-C3-C2	3.35	128.68	121.17
9	BQ	103	BCL	C3A-C2A-C1A	3.35	106.35	101.34
10	BM	403	BPH	CMA-C3A-C4A	-3.34	107.41	114.61
14	AS	104	CRT	C32-C31-C30	-3.34	113.51	123.20
9	AZ	101	BCL	OBB-CAB-CBB	-3.34	112.28	119.77
14	AJ	102	CRT	C21-C22-C23	-3.34	122.59	127.28
9	A5	102	BCL	CAA-C2A-C1A	3.34	122.92	111.97
9	BB	101	BCL	O2A-CGA-O1A	-3.34	115.27	123.63
9	AP	101	BCL	C1-O2A-CGA	3.34	124.73	116.65
9	A0	102	BCL	C5-C3-C2	3.33	128.65	121.17
9	B0	102	BCL	C5-C3-C2	3.33	128.65	121.17
9	B0	102	BCL	O2A-CGA-CBA	3.33	122.00	111.83
14	BF	103	CRT	C32-C31-C30	-3.33	113.55	123.20
14	AN	102	CRT	C21-C20-C19	-3.32	116.72	123.52
9	BD	102	BCL	C1C-NC-C4C	3.32	108.19	106.68
7	BC	502	HEM	C2B-C1B-NB	3.32	113.65	109.84
9	AV	102	BCL	C1C-NC-C4C	3.31	108.19	106.68
14	BO	103	CRT	C21-C20-C19	-3.31	116.76	123.52
14	BB	102	CRT	C29-C28-C30	3.30	123.14	118.09
14	B0	101	CRT	C13-C12-C11	3.30	123.14	118.09
9	BZ	101	BCL	OBB-CAB-CBB	-3.30	112.37	119.77
9	AK	102	BCL	C2A-C1A-CHA	3.30	129.60	123.87
9	A7	103	BCL	CAA-C2A-C1A	3.30	122.79	111.97

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
14	AM	406	CRT	C37-C36-C35	-3.30	120.28	124.91
9	BO	102	BCL	OBB-CAB-CBB	-3.30	112.38	119.77
9	AR	101	BCL	C2C-C3C-C4C	3.30	106.28	101.34
9	BQ	104	BCL	OBB-CAB-CBB	-3.30	112.39	119.77
14	AJ	102	CRT	C32-C31-C30	-3.30	113.65	123.20
9	BJ	101	BCL	C6-C5-C3	3.30	121.50	113.47
9	AY	102	BCL	C2C-C3C-C4C	3.29	106.27	101.34
14	A1	103	CRT	C30-C28-C27	-3.28	113.84	119.01
9	B7	103	BCL	C5-C3-C2	3.28	128.53	121.17
15	AS	101	PEF	O2-C10-C11	3.28	118.58	111.48
9	BQ	104	BCL	CAA-C2A-C1A	3.28	122.72	111.97
14	A1	103	CRT	C11-C12-C14	-3.27	113.86	119.01
9	AN	101	BCL	C5-C3-C2	3.27	128.51	121.17
9	AN	101	BCL	O2A-CGA-O1A	-3.27	115.44	123.63
9	AO	102	BCL	C4D-CHA-C1A	3.27	125.14	121.24
9	A6	101	BCL	CBA-CAA-C2A	3.27	123.52	113.79
9	BV	101	BCL	CAA-C2A-C3A	-3.27	104.17	113.00
9	BB	101	BCL	C2A-C1A-CHA	3.27	129.53	123.87
9	BW	102	BCL	C2C-C3C-C4C	3.26	106.23	101.34
14	A1	103	CRT	C32-C31-C30	-3.26	113.75	123.20
9	A1	102	BCL	CAA-C2A-C1A	3.26	122.66	111.97
15	AM	409	PEF	O2-C10-C11	3.26	118.53	111.48
15	BQ	101	PEF	O2-C10-C11	3.26	118.53	111.48
9	B4	101	BCL	O2A-CGA-O1A	-3.25	115.49	123.63
14	B2	102	CRT	C13-C12-C11	3.25	123.05	118.09
9	BI	102	BCL	C6-C5-C3	3.25	121.38	113.47
9	AD	102	BCL	C4D-CHA-C1A	3.24	125.11	121.24
9	BA	101	BCL	C1C-NC-C4C	3.24	108.16	106.68
9	B7	103	BCL	OBB-CAB-CBB	-3.24	112.51	119.77
9	A8	101	BCL	C2C-C3C-C4C	3.24	106.19	101.34
9	AP	101	BCL	C4D-CHA-C1A	3.24	125.11	121.24
9	AL	303	BCL	OBB-CAB-CBB	-3.24	112.52	119.77
9	BL	303	BCL	OBB-CAB-CBB	-3.24	112.52	119.77
9	AU	102	BCL	OBB-CAB-CBB	-3.24	112.52	119.77
9	BK	102	BCL	OBB-CAB-CBB	-3.23	112.53	119.77
14	AB	102	CRT	C16-C17-C19	-3.23	113.92	119.01
9	BW	102	BCL	OBB-CAB-CBB	-3.23	112.53	119.77
14	A1	103	CRT	C21-C20-C19	-3.23	116.91	123.52
9	B1	102	BCL	C2C-C3C-C4C	3.23	106.18	101.34
14	AN	102	CRT	C32-C31-C30	-3.23	113.84	123.20
11	AL	304	UQ8	C42-C43-C44	-3.23	116.88	127.64
9	BD	102	BCL	C2C-C3C-C4C	3.22	106.17	101.34

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
9	BM	402	BCL	C4D-CHA-C1A	3.22	125.09	121.24
9	A3	103	BCL	OBB-CAB-CBB	-3.22	112.56	119.77
9	AY	102	BCL	OBB-CAB-CBB	-3.22	112.56	119.77
10	AM	403	BPH	OBB-CAB-C3B	3.22	125.37	119.99
9	B8	101	BCL	C6-C5-C3	3.22	121.30	113.47
14	BV	102	CRT	C14-C15-C16	-3.21	113.89	123.20
9	AS	103	BCL	C2B-C1B-NB	-3.21	107.00	110.33
9	BP	101	BCL	OBB-CAB-CBB	-3.21	112.58	119.77
9	BL	303	BCL	C2C-C3C-C4C	3.21	106.14	101.34
14	AR	102	CRT	C21-C20-C19	-3.20	116.96	123.52
9	AT	101	BCL	OBB-CAB-CBB	-3.20	112.60	119.77
14	BM	406	CRT	C20-C19-C17	-3.20	122.79	127.28
9	BN	101	BCL	OBB-CAB-CBB	-3.20	112.60	119.77
9	BY	102	BCL	O2D-CGD-O1D	-3.20	117.62	123.85
9	BP	101	BCL	C2C-C3C-C4C	3.20	106.13	101.34
9	A2	101	BCL	C4D-CHA-C1A	3.20	125.06	121.24
14	AN	102	CRT	C4-C5-C6	-3.19	120.43	124.91
14	AM	406	CRT	C10-C9-C7	-3.19	122.80	127.28
9	A3	103	BCL	CAA-C2A-C1A	3.19	122.43	111.97
14	B1	103	CRT	C10-C9-C7	-3.19	122.80	127.28
9	A6	101	BCL	C2C-C3C-C4C	3.19	106.12	101.34
9	AR	101	BCL	C4D-CHA-C1A	3.19	125.05	121.24
9	AO	102	BCL	C2C-C3C-C4C	3.19	106.11	101.34
9	AB	101	BCL	OBB-CAB-CBB	-3.19	112.63	119.77
9	AA	101	BCL	OBB-CAB-CBB	-3.19	112.63	119.77
7	AC	502	HEM	C2B-C1B-NB	3.19	113.50	109.84
14	BP	102	CRT	C14-C15-C16	-3.19	113.97	123.20
9	AL	303	BCL	C2C-C3C-C4C	3.19	106.11	101.34
9	BY	102	BCL	OBB-CAB-CBB	-3.19	112.64	119.77
9	B3	102	BCL	OBB-CAB-CBB	-3.18	112.65	119.77
9	BM	402	BCL	OBB-CAB-CBB	-3.18	112.65	119.77
9	BE	101	BCL	C5-C3-C2	3.18	128.31	121.17
9	AM	401	BCL	C4D-CHA-C1A	3.18	125.03	121.24
9	BS	102	BCL	CBA-CAA-C2A	3.18	123.24	113.79
9	BU	102	BCL	OBB-CAB-CBB	-3.18	112.66	119.77
9	BX	101	BCL	O2A-CGA-CBA	3.17	121.52	111.83
9	AB	101	BCL	C1C-NC-C4C	3.17	108.13	106.68
9	AY	102	BCL	CAA-C2A-C1A	3.17	122.37	111.97
14	AB	102	CRT	C14-C15-C16	-3.17	114.01	123.20
9	A1	102	BCL	O2D-CGD-O1D	-3.17	117.68	123.85
14	A1	103	CRT	C29-C28-C30	3.17	122.93	118.09
9	AN	101	BCL	O2A-CGA-CBA	3.17	121.50	111.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
9	BS	102	BCL	C4D-CHA-C1A	3.17	125.02	121.24
9	BJ	101	BCL	C2C-C3C-C4C	3.17	106.08	101.34
9	B5	102	BCL	OBB-CAB-CBB	-3.17	112.68	119.77
9	BS	102	BCL	OBB-CAB-CBB	-3.17	112.68	119.77
9	BX	101	BCL	OBB-CAB-CBB	-3.16	112.68	119.77
9	AM	402	BCL	OBB-CAB-CBB	-3.16	112.68	119.77
14	B1	103	CRT	C21-C20-C19	-3.16	117.05	123.52
9	B3	102	BCL	C4D-CHA-C1A	3.16	125.02	121.24
9	BP	101	BCL	C2A-C1A-CHA	3.16	129.36	123.87
9	AN	101	BCL	C2C-C3C-C4C	3.16	106.08	101.34
9	BT	101	BCL	OBB-CAB-CBB	-3.16	112.69	119.77
9	AP	101	BCL	CAB-C3B-C2B	3.16	134.31	127.74
9	AA	101	BCL	O2D-CGD-CBD	3.16	116.76	111.23
14	BP	102	CRT	C27-C26-C25	-3.16	114.04	123.20
9	A5	102	BCL	OBB-CAB-CBB	-3.16	112.69	119.77
9	AF	102	BCL	C2C-C3C-C4C	3.16	106.07	101.34
14	AW	102	CRT	C4-C5-C6	-3.16	120.48	124.91
9	BY	102	BCL	C3A-C2A-C1A	3.16	106.07	101.34
9	AI	102	BCL	C4D-CHA-C1A	3.16	125.01	121.24
9	BB	101	BCL	CHB-C1B-NB	3.16	128.78	124.05
14	BM	406	CRT	C31-C32-C33	-3.16	122.85	127.28
9	AL	301	BCL	OBB-CAB-CBB	-3.15	112.71	119.77
9	BT	101	BCL	O2A-CGA-CBA	3.15	121.45	111.83
9	B6	101	BCL	OBB-CAB-CBB	-3.15	112.71	119.77
14	AB	102	CRT	C27-C26-C25	-3.15	114.07	123.20
9	B2	101	BCL	C2C-C3C-C4C	3.15	106.06	101.34
14	BF	103	CRT	C21-C22-C23	-3.15	122.86	127.28
9	BE	101	BCL	CAA-C2A-C1A	3.15	122.30	111.97
9	BQ	104	BCL	O2A-CGA-O1A	-3.15	115.75	123.63
9	AQ	102	BCL	OBB-CAB-CBB	-3.15	112.72	119.77
9	BF	102	BCL	CHB-C1B-NB	3.15	128.77	124.05
9	AW	101	BCL	C3A-C2A-C1A	3.15	106.05	101.34
9	AO	102	BCL	CAA-C2A-C3A	-3.14	104.50	113.00
9	BL	301	BCL	C2C-C3C-C4C	3.14	106.05	101.34
9	B9	102	BCL	C2C-C3C-C4C	3.14	106.05	101.34
9	AT	101	BCL	C1-O2A-CGA	3.14	124.26	116.65
9	AU	102	BCL	C2C-C3C-C4C	3.14	106.04	101.34
9	AG	101	BCL	C6-C5-C3	3.14	121.12	113.47
14	AB	102	CRT	C18-C17-C16	3.14	122.88	118.09
7	AC	503	HEM	C3B-C4B-NB	3.14	111.72	109.47
14	BN	102	CRT	C29-C28-C30	3.14	122.88	118.09
7	BC	503	HEM	C3B-C4B-NB	3.14	111.72	109.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
9	AM	402	BCL	C4D-CHA-C1A	3.13	124.98	121.24
14	AT	102	CRT	C20-C19-C17	-3.13	122.88	127.28
9	A9	102	BCL	C5-C3-C2	3.13	128.20	121.17
9	AN	101	BCL	OBB-CAB-CBB	-3.13	112.75	119.77
9	AV	102	BCL	OBB-CAB-CBB	-3.13	112.75	119.77
9	BM	402	BCL	C5-C3-C2	3.13	128.20	121.17
9	AW	101	BCL	OBB-CAB-CBB	-3.13	112.75	119.77
9	B4	101	BCL	C2C-C3C-C4C	3.13	106.03	101.34
9	AM	402	BCL	C6-C5-C3	3.13	121.09	113.47
9	AV	102	BCL	C2C-C3C-C4C	3.13	106.03	101.34
9	A2	101	BCL	C1C-NC-C4C	3.13	108.11	106.68
9	BG	101	BCL	OBB-CAB-CBB	-3.13	112.77	119.77
10	BL	302	BPH	C2B-C1B-NB	-3.13	107.17	109.43
14	BF	103	CRT	C29-C28-C30	3.13	122.86	118.09
9	A2	101	BCL	O2D-CGD-CBD	3.12	116.69	111.23
14	B2	102	CRT	C14-C15-C16	-3.12	114.15	123.20
9	BM	402	BCL	C2C-C3C-C4C	3.12	106.02	101.34
14	AB	102	CRT	C29-C28-C30	3.12	122.86	118.09
9	B7	103	BCL	C6-C5-C3	3.12	121.07	113.47
9	AJ	101	BCL	OBB-CAB-CBB	-3.12	112.78	119.77
9	BY	102	BCL	C4D-CHA-C1A	3.12	124.97	121.24
9	A0	102	BCL	OBB-CAB-CBB	-3.12	112.78	119.77
9	AF	102	BCL	OBB-CAB-CBB	-3.12	112.78	119.77
9	B5	102	BCL	C2C-C3C-C4C	3.12	106.01	101.34
9	BO	102	BCL	C2C-C3C-C4C	3.12	106.01	101.34
9	BN	101	BCL	C1C-NC-C4C	3.12	108.10	106.68
14	BB	102	CRT	C18-C17-C16	3.12	122.85	118.09
9	BM	401	BCL	C4D-CHA-C1A	3.12	124.96	121.24
9	BD	102	BCL	OBB-CAB-CBB	-3.11	112.80	119.77
9	A7	103	BCL	C2C-C3C-C4C	3.11	106.00	101.34
9	B3	102	BCL	CAB-C3B-C2B	3.11	134.21	127.74
9	AM	401	BCL	C2C-C3C-C4C	3.11	105.99	101.34
9	B4	101	BCL	C6-C5-C3	3.11	121.04	113.47
9	AD	102	BCL	C2C-C3C-C4C	3.11	105.99	101.34
9	AA	101	BCL	C2C-C3C-C4C	3.11	105.99	101.34
9	AU	102	BCL	CAA-CBA-CGA	3.10	122.01	113.21
9	B4	101	BCL	C1-O2A-CGA	3.10	124.15	116.65
14	BB	102	CRT	C36-C35-C33	-3.10	121.21	125.89
9	BB	101	BCL	C1-O2A-CGA	3.10	124.15	116.65
9	AP	101	BCL	C4D-C3D-CAD	-3.10	104.74	108.11
9	BT	101	BCL	CHC-C4B-NB	3.10	128.69	124.05
9	B4	101	BCL	OBB-CAB-CBB	-3.09	112.84	119.77

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
9	BX	101	BCL	O2A-CGA-O1A	-3.09	115.89	123.63
9	AB	101	BCL	C2C-C3C-C4C	3.09	105.97	101.34
9	BV	101	BCL	OBB-CAB-CBB	-3.09	112.85	119.77
14	BB	102	CRT	C27-C26-C25	-3.09	114.25	123.20
9	AQ	102	BCL	C2A-C1A-CHA	3.09	129.23	123.87
9	B3	102	BCL	C2C-C3C-C4C	3.09	105.96	101.34
9	BB	101	BCL	C6-C5-C3	3.08	120.97	113.47
9	AI	102	BCL	C6-C5-C3	3.08	120.97	113.47
9	AT	101	BCL	C4D-CHA-C1A	3.08	124.92	121.24
9	BL	301	BCL	OBB-CAB-CBB	-3.08	112.87	119.77
14	BS	103	CRT	C20-C19-C17	-3.08	122.96	127.28
14	BB	102	CRT	C14-C15-C16	-3.08	114.28	123.20
9	B9	102	BCL	C2A-C1A-CHA	3.08	129.21	123.87
9	B9	102	BCL	CAB-C3B-C2B	3.08	134.13	127.74
9	BA	101	BCL	OBB-CAB-CBB	-3.07	112.88	119.77
11	AL	304	UQ8	C32-C33-C34	-3.07	120.59	127.62
9	B7	103	BCL	C2C-C3C-C4C	3.07	105.94	101.34
9	BJ	101	BCL	O2A-CGA-CBA	3.07	121.20	111.83
9	AP	101	BCL	OBB-CAB-CBB	-3.07	112.89	119.77
9	B4	101	BCL	O2A-CGA-CBA	3.07	121.20	111.83
9	AA	101	BCL	C4D-CHA-C1A	3.07	124.91	121.24
14	AA	102	CRT	C4-C5-C6	-3.07	120.60	124.91
9	BK	102	BCL	C2C-C3C-C4C	3.07	105.94	101.34
9	B8	101	BCL	C2C-C3C-C4C	3.07	105.94	101.34
9	BX	101	BCL	C2A-C1A-CHA	3.07	129.19	123.87
9	A6	101	BCL	OBB-CAB-CBB	-3.07	112.90	119.77
14	BP	102	CRT	C4-C5-C6	-3.07	120.61	124.91
9	AG	101	BCL	OBB-CAB-CBB	-3.07	112.90	119.77
9	BJ	101	BCL	C4D-CHA-C1A	3.07	124.90	121.24
10	BL	302	BPH	OBB-CAB-C3B	3.06	125.11	119.99
9	AL	301	BCL	C2C-C3C-C4C	3.06	105.93	101.34
9	A8	101	BCL	OBB-CAB-CBB	-3.06	112.91	119.77
9	AS	103	BCL	C2C-C3C-C4C	3.06	105.93	101.34
9	AM	402	BCL	C2C-C3C-C4C	3.06	105.93	101.34
9	A5	102	BCL	C2C-C3C-C4C	3.06	105.93	101.34
9	BQ	103	BCL	OBB-CAB-CBB	-3.06	112.91	119.77
11	BL	304	UQ8	C46-C44-C43	-3.06	113.47	122.66
9	AK	102	BCL	OBB-CAB-CBB	-3.06	112.91	119.77
9	AI	102	BCL	C2C-C3C-C4C	3.06	105.92	101.34
9	AB	101	BCL	C6-C5-C3	3.06	120.92	113.47
9	AZ	101	BCL	C4D-CHA-C1A	3.06	124.89	121.24
15	BM	407	PEF	O3-C30-C31	3.06	125.34	112.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
9	BI	102	BCL	C2C-C3C-C4C	3.06	105.92	101.34
14	A2	102	CRT	C1M-O1-C1	3.06	136.58	117.25
10	AM	403	BPH	C2B-C1B-NB	-3.05	107.23	109.43
9	AE	101	BCL	O2A-CGA-O1A	-3.05	116.00	123.63
9	AI	102	BCL	OBB-CAB-CBB	-3.05	112.94	119.77
9	A7	103	BCL	CMA-C3A-C2A	-3.05	102.20	113.98
9	A6	101	BCL	C4D-CHA-C1A	3.05	124.88	121.24
9	AQ	102	BCL	C2C-C3C-C4C	3.05	105.90	101.34
15	AH	301	PEF	O3-C30-C31	3.05	125.30	112.38
9	AE	101	BCL	C2A-C1A-CHA	3.05	129.15	123.87
13	BM	405	MQ8	C45-C43-C44	3.05	120.51	115.23
14	BW	103	CRT	C14-C15-C16	-3.04	114.38	123.20
9	BS	102	BCL	C2C-C3C-C4C	3.04	105.90	101.34
14	BF	103	CRT	C1-C4-C5	3.04	120.88	112.98
9	A9	102	BCL	C2C-C3C-C4C	3.04	105.90	101.34
9	BN	101	BCL	CAB-C3B-C2B	3.04	134.06	127.74
10	AM	403	BPH	O2D-CGD-CBD	-3.04	107.60	110.95
9	BF	102	BCL	C2A-C1A-CHA	3.04	129.15	123.87
9	AE	101	BCL	C2C-C3C-C4C	3.04	105.89	101.34
9	A3	104	BCL	C2C-C3C-C4C	3.04	105.89	101.34
15	AM	407	PEF	O3-C30-C31	3.04	125.28	112.38
9	AL	301	BCL	CBA-CAA-C2A	3.04	122.84	113.79
9	BN	101	BCL	O2A-CGA-O1A	-3.04	116.03	123.63
14	AP	102	CRT	C14-C15-C16	-3.04	114.40	123.20
9	A8	101	BCL	CAA-C2A-C3A	-3.04	104.79	113.00
14	AB	102	CRT	C30-C28-C27	-3.04	114.23	119.01
9	BT	101	BCL	C2C-C3C-C4C	3.03	105.88	101.34
9	AS	103	BCL	OBB-CAB-CBB	-3.03	112.97	119.77
9	BJ	101	BCL	CAA-C2A-C1A	3.03	121.92	111.97
9	B8	101	BCL	OBB-CAB-CBB	-3.03	112.98	119.77
9	A9	102	BCL	CAB-C3B-C2B	3.03	134.04	127.74
9	AO	102	BCL	OBB-CAB-CBB	-3.03	112.98	119.77
9	A8	101	BCL	CAB-C3B-C2B	3.03	134.04	127.74
9	B6	101	BCL	C2C-C3C-C4C	3.03	105.88	101.34
14	AJ	102	CRT	C21-C20-C19	-3.03	117.32	123.52
9	AT	101	BCL	CBA-CAA-C2A	3.03	122.80	113.79
9	AV	102	BCL	CAA-C2A-C1A	3.02	121.89	111.97
9	BZ	101	BCL	CAB-C3B-C2B	3.02	134.03	127.74
9	AF	102	BCL	C4D-CHA-C1A	3.02	124.85	121.24
9	BG	101	BCL	O2A-CGA-CBA	3.02	121.05	111.83
10	BL	302	BPH	C4C-C3C-C2C	-3.02	99.96	102.84
9	AJ	101	BCL	C6-C5-C3	3.02	120.82	113.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
9	AG	101	BCL	C2C-C3C-C4C	3.02	105.86	101.34
9	B2	101	BCL	OBB-CAB-CBB	-3.02	113.01	119.77
9	A7	103	BCL	OBB-CAB-CBB	-3.02	113.01	119.77
9	BE	101	BCL	C3A-C2A-C1A	3.02	105.86	101.34
9	BM	401	BCL	C2C-C3C-C4C	3.01	105.85	101.34
14	BV	102	CRT	C34-C33-C35	3.01	122.69	118.09
14	BN	102	CRT	C21-C22-C23	-3.01	123.05	127.28
14	B5	103	CRT	C4-C5-C6	-3.01	120.68	124.91
9	A3	103	BCL	C2C-C3C-C4C	3.01	105.85	101.34
9	AA	101	BCL	C5-C3-C2	3.01	127.92	121.17
14	AB	102	CRT	C32-C31-C30	-3.01	114.48	123.20
9	AJ	101	BCL	C2C-C3C-C4C	3.01	105.84	101.34
10	AL	302	BPH	OBB-CAB-C3B	3.01	125.01	119.99
9	AO	102	BCL	C4D-C3D-CAD	-3.00	104.84	108.11
9	BW	102	BCL	C4D-CHA-C1A	3.00	124.83	121.24
9	A6	101	BCL	CAB-C3B-C2B	3.00	133.98	127.74
9	BF	102	BCL	CAB-C3B-C2B	3.00	133.98	127.74
9	A3	104	BCL	OBB-CAB-CBB	-3.00	113.05	119.77
9	BM	402	BCL	CAA-C2A-C3A	-3.00	104.89	113.00
9	AS	103	BCL	C3A-C2A-C1A	3.00	105.83	101.34
9	AV	102	BCL	C4D-CHA-C1A	3.00	124.82	121.24
14	B7	102	CRT	C6-C7-C9	-3.00	114.30	119.01
14	B1	103	CRT	C27-C26-C25	-2.99	114.52	123.20
14	AP	102	CRT	C27-C26-C25	-2.99	114.53	123.20
9	B0	102	BCL	C2C-C3C-C4C	2.99	105.82	101.34
14	BF	103	CRT	C21-C20-C19	-2.99	117.40	123.52
9	A1	102	BCL	CAA-CBA-CGA	2.99	121.69	113.21
9	BA	101	BCL	C2C-C3C-C4C	2.99	105.81	101.34
14	B2	102	CRT	C5-C6-C7	2.98	130.40	125.89
9	BB	101	BCL	CHA-C1A-NA	-2.98	119.63	126.39
14	BW	103	CRT	C13-C12-C11	2.98	122.64	118.09
9	BF	102	BCL	C2C-C3C-C4C	2.98	105.81	101.34
9	B8	101	BCL	O2D-CGD-O1D	-2.98	118.04	123.85
9	AX	101	BCL	OBB-CAB-CBB	-2.98	113.09	119.77
14	A5	103	CRT	C21-C22-C23	-2.98	123.10	127.28
9	A3	103	BCL	O2D-CGD-O1D	-2.98	118.05	123.85
9	AL	303	BCL	O2D-CGD-O1D	-2.98	118.05	123.85
14	BG	102	CRT	C29-C28-C30	2.98	122.64	118.09
9	AR	101	BCL	C2A-C3A-C4A	2.98	106.68	101.87
9	AT	101	BCL	C2A-C1A-CHA	2.98	129.04	123.87
9	BQ	104	BCL	C2C-C3C-C4C	2.98	105.80	101.34
9	AU	102	BCL	C1C-NC-C4C	2.98	108.04	106.68

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
9	AV	102	BCL	CAA-C2A-C3A	-2.98	104.96	113.00
9	BZ	101	BCL	C6-C5-C3	2.97	120.71	113.47
9	BV	101	BCL	CAB-C3B-C2B	2.97	133.92	127.74
9	B8	101	BCL	C5-C3-C2	2.97	127.84	121.17
9	AK	102	BCL	C2C-C3C-C4C	2.97	105.79	101.34
9	A3	104	BCL	C1C-NC-C4C	2.97	108.03	106.68
9	A0	102	BCL	C2C-C3C-C4C	2.96	105.78	101.34
9	BL	303	BCL	C4D-CHA-C1A	2.96	124.78	121.24
9	AM	401	BCL	OBB-CAB-CBB	-2.96	113.14	119.77
14	B1	103	CRT	C14-C15-C16	-2.96	114.62	123.20
9	BY	102	BCL	C2C-C3C-C4C	2.96	105.77	101.34
9	A0	102	BCL	CBA-CAA-C2A	2.96	122.60	113.79
9	A1	102	BCL	C2C-C3C-C4C	2.96	105.77	101.34
9	A8	101	BCL	CMA-C3A-C2A	-2.96	102.55	113.98
9	AV	102	BCL	C6-C5-C3	2.95	120.66	113.47
9	AP	101	BCL	O2D-CGD-O1D	-2.95	118.10	123.85
9	AL	301	BCL	C2A-C1A-CHA	2.95	128.99	123.87
9	AT	101	BCL	CHA-C1A-NA	-2.95	119.71	126.39
14	AX	102	CRT	C10-C11-C12	-2.95	118.28	126.36
9	AZ	101	BCL	C2A-C3A-C4A	2.95	106.63	101.87
14	BM	406	CRT	C10-C9-C7	-2.94	123.15	127.28
14	B0	101	CRT	C38-C37-C36	-2.94	105.35	112.98
14	BP	102	CRT	C29-C28-C30	2.94	122.58	118.09
9	BF	102	BCL	C1C-NC-C4C	2.94	108.02	106.68
9	BW	102	BCL	C3A-C2A-C1A	2.94	105.74	101.34
9	B8	101	BCL	C2A-C1A-CHA	2.94	128.97	123.87
9	A3	104	BCL	CAB-C3B-C2B	2.94	133.84	127.74
9	A3	104	BCL	C4D-CHA-C1A	2.94	124.75	121.24
9	BX	101	BCL	C1-O2A-CGA	2.93	123.75	116.65
9	AR	101	BCL	OBB-CAB-CBB	-2.93	113.20	119.77
9	A7	103	BCL	CAB-C3B-C2B	2.93	133.83	127.74
9	B9	102	BCL	OBB-CAB-CBB	-2.92	113.22	119.77
9	AL	303	BCL	C4D-CHA-C1A	2.92	124.73	121.24
9	BB	101	BCL	C4D-C3D-CAD	-2.92	104.93	108.11
14	AG	102	CRT	C32-C31-C30	-2.92	114.74	123.20
14	B5	103	CRT	C32-C31-C30	-2.92	114.74	123.20
9	BI	102	BCL	C3A-C2A-C1A	2.92	105.71	101.34
9	BN	101	BCL	O2A-CGA-CBA	2.92	120.73	111.83
9	AN	101	BCL	C1-O2A-CGA	2.91	123.70	116.65
9	A5	102	BCL	CAB-C3B-C2B	2.91	133.79	127.74
14	BG	102	CRT	C30-C28-C27	-2.91	114.43	119.01
9	B1	102	BCL	OBB-CAB-CBB	-2.91	113.25	119.77

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
9	B7	103	BCL	C3A-C2A-C1A	2.91	105.70	101.34
9	B0	102	BCL	C2A-C1A-CHA	2.91	128.91	123.87
9	AJ	101	BCL	C4D-CHA-C1A	2.91	124.71	121.24
14	AM	406	CRT	C15-C14-C12	-2.91	123.20	127.28
14	BB	102	CRT	C16-C17-C19	-2.90	114.44	119.01
9	AX	101	BCL	CAB-C3B-C2B	2.90	133.78	127.74
9	BM	402	BCL	CHB-C1B-NB	2.90	128.40	124.05
9	BU	102	BCL	C2C-C3C-C4C	2.90	105.68	101.34
9	AR	101	BCL	C1C-NC-C4C	2.90	108.00	106.68
14	AS	104	CRT	C4-C5-C6	-2.90	120.85	124.91
9	AB	101	BCL	C2A-C1A-CHA	2.89	128.89	123.87
14	A5	103	CRT	C32-C31-C30	-2.89	114.82	123.20
14	AA	102	CRT	C20-C19-C17	-2.89	123.22	127.28
9	BG	101	BCL	CHB-C1B-NB	2.89	128.39	124.05
14	AN	102	CRT	C29-C28-C30	2.89	122.50	118.09
9	BQ	103	BCL	C4D-C3D-CAD	-2.89	104.97	108.11
9	BA	101	BCL	C6-C5-C3	2.89	120.50	113.47
9	AP	101	BCL	C6-C5-C3	2.89	120.50	113.47
9	BL	301	BCL	CBA-CAA-C2A	2.89	122.38	113.79
9	B6	101	BCL	C1-O2A-CGA	2.89	123.64	116.65
9	B8	101	BCL	CAB-C3B-C2B	2.88	133.74	127.74
9	BK	102	BCL	C4D-CHA-C1A	2.88	124.69	121.24
9	BX	101	BCL	CHA-C1A-NA	-2.88	119.86	126.39
9	A9	102	BCL	CMA-C3A-C2A	-2.88	102.84	113.98
9	BM	401	BCL	OBB-CAB-CBB	-2.88	113.32	119.77
9	AY	102	BCL	CAB-C3B-C2B	2.88	133.73	127.74
9	BZ	101	BCL	C2C-C3C-C4C	2.88	105.65	101.34
9	BT	101	BCL	C2A-C1A-CHA	2.88	128.86	123.87
14	BP	102	CRT	C13-C12-C11	2.88	122.48	118.09
9	AO	102	BCL	CAB-C3B-C2B	2.88	133.72	127.74
9	BD	102	BCL	CHB-C1B-NB	2.88	128.36	124.05
9	A2	101	BCL	C2C-C3C-C4C	2.88	105.65	101.34
9	AT	101	BCL	CHB-C1B-NB	2.88	128.36	124.05
14	BB	102	CRT	C34-C33-C35	2.88	122.48	118.09
9	AR	101	BCL	CHA-C1A-NA	-2.87	119.89	126.39
14	AN	102	CRT	C14-C15-C16	-2.87	114.89	123.20
14	BO	103	CRT	C32-C31-C30	-2.87	114.89	123.20
9	AW	101	BCL	C2C-C3C-C4C	2.87	105.64	101.34
9	BI	102	BCL	OBB-CAB-CBB	-2.87	113.35	119.77
14	BV	102	CRT	C29-C28-C30	2.87	122.47	118.09
14	A5	103	CRT	C20-C19-C17	-2.87	123.26	127.28
9	AA	101	BCL	C1C-NC-C4C	2.86	107.98	106.68

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
9	AX	101	BCL	C2C-C3C-C4C	2.86	105.63	101.34
9	BL	303	BCL	O2D-CGD-O1D	-2.86	118.28	123.85
9	BJ	101	BCL	O2A-CGA-O1A	-2.86	116.47	123.63
9	AG	101	BCL	CAA-C2A-C1A	2.86	121.34	111.97
9	AN	101	BCL	CAB-C3B-C2B	2.86	133.68	127.74
9	AI	102	BCL	C1C-NC-C4C	2.86	107.98	106.68
9	AG	101	BCL	C1-O2A-CGA	2.85	123.56	116.65
14	AP	102	CRT	C32-C31-C30	-2.85	114.93	123.20
9	AZ	101	BCL	CAB-C3B-C2B	2.85	133.67	127.74
9	AO	102	BCL	C1C-NC-C4C	2.85	107.98	106.68
9	AP	101	BCL	O2A-CGA-O1A	-2.85	116.50	123.63
14	BM	406	CRT	C5-C6-C7	-2.85	121.58	125.89
9	B3	102	BCL	C6-C5-C3	2.85	120.41	113.47
14	BA	102	CRT	C32-C31-C30	-2.85	114.95	123.20
14	BU	103	CRT	C32-C31-C30	-2.85	114.95	123.20
9	BJ	101	BCL	OBB-CAB-CBB	-2.84	113.40	119.77
9	B2	101	BCL	C6-C5-C3	2.84	120.40	113.47
9	BL	301	BCL	C6-C5-C3	2.84	120.39	113.47
9	AE	101	BCL	CAA-C2A-C3A	-2.84	105.32	113.00
14	AM	406	CRT	C26-C27-C28	-2.84	123.29	127.28
14	B1	103	CRT	C29-C28-C30	2.84	122.43	118.09
13	AM	405	MQ8	C21-C22-C23	-2.84	121.12	127.62
7	AC	501	HEM	C3B-C2B-C1B	-2.84	104.28	106.41
9	BE	101	BCL	OBB-CAB-CBB	-2.84	113.41	119.77
9	BS	102	BCL	C1C-NC-C4C	2.84	107.97	106.68
9	BV	101	BCL	C6-C5-C3	2.84	120.38	113.47
9	BL	301	BCL	C2A-C1A-CHA	2.84	128.79	123.87
14	AR	102	CRT	C29-C28-C30	2.83	122.42	118.09
14	AP	102	CRT	C1M-O1-C1	2.83	135.17	117.25
9	BQ	103	BCL	C2C-C3C-C4C	2.83	105.58	101.34
14	BS	103	CRT	C13-C12-C11	2.83	122.41	118.09
9	BV	101	BCL	C2C-C3C-C4C	2.83	105.58	101.34
9	AE	101	BCL	CAB-C3B-C2B	2.83	133.62	127.74
14	AW	102	CRT	C13-C12-C11	2.83	122.41	118.09
14	BW	103	CRT	C4-C5-C6	-2.83	120.94	124.91
9	AO	102	BCL	CHA-C1A-NA	-2.83	119.99	126.39
9	B7	103	BCL	C1C-NC-C4C	2.83	107.97	106.68
9	AD	102	BCL	C1C-NC-C4C	2.83	107.97	106.68
9	AS	103	BCL	CHB-C1B-NB	2.82	128.29	124.05
14	B1	103	CRT	C21-C22-C23	-2.82	123.32	127.28
9	AX	101	BCL	C4D-CHA-C1A	2.82	124.61	121.24
9	A9	102	BCL	OBB-CAB-CBB	-2.82	113.45	119.77

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
14	BO	103	CRT	C29-C28-C30	2.82	122.40	118.09
9	A3	103	BCL	CAB-C3B-C2B	2.82	133.60	127.74
14	AT	102	CRT	C20-C21-C22	-2.82	117.75	123.52
9	BY	102	BCL	C6-C5-C3	2.82	120.33	113.47
9	AL	301	BCL	C6-C5-C3	2.82	120.33	113.47
9	A5	102	BCL	CMA-C3A-C2A	-2.81	103.10	113.98
9	B8	101	BCL	O2A-CGA-O1A	-2.81	116.59	123.63
11	BL	304	UQ8	C42-C43-C44	-2.81	118.27	127.64
9	BM	401	BCL	CHB-C1B-NB	2.81	128.26	124.05
9	A8	101	BCL	C5-C3-C2	2.81	127.47	121.17
9	BB	101	BCL	OBB-CAB-CBB	-2.81	113.48	119.77
9	AT	101	BCL	C6-C5-C3	2.81	120.31	113.47
9	BX	101	BCL	C2C-C3C-C4C	2.81	105.55	101.34
14	AG	102	CRT	C21-C22-C23	-2.81	123.34	127.28
14	BA	102	CRT	C20-C19-C17	-2.81	123.34	127.28
14	B1	103	CRT	C13-C12-C11	2.81	122.38	118.09
9	B0	102	BCL	CHA-C1A-NA	-2.81	120.03	126.39
9	AD	102	BCL	C3A-C2A-C1A	2.81	105.54	101.34
9	AJ	101	BCL	O2A-CGA-O1A	-2.81	116.61	123.63
9	B4	101	BCL	O2D-CGD-O1D	-2.80	118.39	123.85
14	BU	103	CRT	C20-C21-C22	-2.80	117.78	123.52
14	A0	101	CRT	C8-C7-C9	-2.80	118.28	122.82
14	BP	102	CRT	C32-C31-C30	-2.80	115.08	123.20
14	AR	102	CRT	C13-C12-C11	2.80	122.36	118.09
7	AC	502	HEM	CHC-C1C-NC	2.80	127.50	124.45
9	B3	102	BCL	O2D-CGD-O1D	-2.80	118.40	123.85
14	BV	102	CRT	C5-C6-C7	-2.80	121.66	125.89
9	B6	101	BCL	CAB-C3B-C2B	2.80	133.56	127.74
9	BT	101	BCL	CHA-C1A-NA	-2.80	120.06	126.39
9	B3	102	BCL	C2A-C1A-CHA	2.79	128.72	123.87
9	BF	102	BCL	C2B-C1B-NB	-2.79	107.43	110.33
9	BX	101	BCL	CAB-C3B-C2B	2.79	133.55	127.74
9	A8	101	BCL	C4D-CHA-C1A	2.79	124.58	121.24
9	B7	103	BCL	CAB-C3B-C2B	2.79	133.54	127.74
14	BW	103	CRT	C10-C9-C7	-2.79	123.36	127.28
14	AT	102	CRT	C8-C7-C6	2.79	122.35	118.09
9	BU	102	BCL	O2D-CGD-O1D	-2.79	118.42	123.85
14	B7	102	CRT	C8-C7-C6	2.79	122.35	118.09
13	AM	405	MQ8	C45-C43-C44	2.79	120.07	115.23
14	B7	102	CRT	C32-C31-C30	-2.79	115.12	123.20
9	AV	102	BCL	CHB-C1B-NB	2.79	128.23	124.05
9	BK	102	BCL	C3A-C2A-C1A	2.79	105.51	101.34

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
14	BF	103	CRT	C8-C7-C6	2.79	122.34	118.09
9	BP	101	BCL	CHB-C1B-NB	2.79	128.23	124.05
9	AN	101	BCL	C1C-NC-C4C	2.79	107.95	106.68
9	B1	102	BCL	O2D-CGD-O1D	-2.79	118.43	123.85
9	B4	101	BCL	CHA-C1A-NA	-2.78	120.09	126.39
14	BO	103	CRT	C14-C15-C16	-2.78	115.14	123.20
9	BO	102	BCL	C4D-CHA-C1A	2.78	124.56	121.24
14	AJ	102	CRT	C29-C28-C30	2.78	122.34	118.09
14	AR	102	CRT	C14-C15-C16	-2.78	115.14	123.20
14	B5	103	CRT	C20-C19-C17	-2.78	123.38	127.28
9	BQ	104	BCL	C2A-C3A-C4A	2.78	106.36	101.87
9	B7	103	BCL	C2A-C1A-CHA	2.78	128.69	123.87
11	BL	304	UQ8	C30-C29-C31	2.77	120.04	115.23
10	AL	302	BPH	OBB-CAB-CBB	-2.77	114.28	120.19
14	AA	102	CRT	C32-C31-C30	-2.77	115.17	123.20
14	AP	102	CRT	C29-C28-C30	2.77	122.32	118.09
9	BO	102	BCL	CHB-C1B-NB	2.77	128.21	124.05
9	B6	101	BCL	C6-C5-C3	2.77	120.22	113.47
9	BV	101	BCL	CHB-C1B-NB	2.77	128.20	124.05
9	A0	102	BCL	C4D-CHA-C1A	2.77	124.55	121.24
9	BA	101	BCL	CAB-C3B-C2B	2.77	133.49	127.74
9	BJ	101	BCL	CHB-C1B-NB	2.77	128.20	124.05
14	A0	101	CRT	C21-C20-C19	-2.77	117.86	123.52
9	A8	101	BCL	C2B-C1B-NB	-2.77	107.46	110.33
9	B4	101	BCL	C1C-NC-C4C	2.76	107.94	106.68
14	BG	102	CRT	C27-C26-C25	-2.76	115.19	123.20
9	BI	102	BCL	C4D-CHA-C1A	2.76	124.54	121.24
14	AT	102	CRT	C14-C15-C16	-2.76	115.20	123.20
9	AF	102	BCL	C3A-C2A-C1A	2.76	105.47	101.34
9	AB	101	BCL	CAB-C3B-C2B	2.76	133.47	127.74
9	BV	101	BCL	O2A-CGA-O1A	-2.76	116.73	123.63
14	AP	102	CRT	C13-C12-C11	2.76	122.30	118.09
14	B1	103	CRT	C18-C17-C16	2.76	122.30	118.09
9	AQ	102	BCL	C4D-C3D-CAD	-2.76	105.11	108.11
14	AR	102	CRT	C32-C31-C30	-2.76	115.21	123.20
9	A8	101	BCL	CHA-C1A-NA	-2.76	120.15	126.39
7	AC	504	HEM	C3B-C2B-C1B	-2.76	104.34	106.41
14	A2	102	CRT	C14-C15-C16	-2.76	115.22	123.20
9	B5	102	BCL	CAB-C3B-C2B	2.75	133.47	127.74
9	AS	103	BCL	C4D-CHA-C1A	2.75	124.53	121.24
9	AW	101	BCL	C4D-C3D-CAD	-2.75	105.11	108.11
9	BL	301	BCL	CAA-C2A-C1A	2.75	121.00	111.97

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
9	BT	101	BCL	C4D-CHA-C1A	2.75	124.53	121.24
9	B6	101	BCL	O2D-CGD-O1D	-2.75	118.49	123.85
14	AN	102	CRT	C34-C33-C35	2.75	122.29	118.09
14	BU	103	CRT	C29-C28-C30	2.75	122.29	118.09
10	AL	302	BPH	C1A-C2A-C3A	-2.75	100.22	102.84
10	AL	302	BPH	C2B-C1B-NB	-2.75	107.44	109.43
9	BG	101	BCL	CAB-C3B-C2B	2.75	133.45	127.74
9	BQ	104	BCL	C4D-CHA-C1A	2.74	124.52	121.24
9	A3	103	BCL	C6-C5-C3	2.74	120.15	113.47
14	AB	102	CRT	C13-C12-C11	2.74	122.28	118.09
9	BL	301	BCL	CHB-C1B-NB	2.74	128.16	124.05
14	AG	102	CRT	C29-C28-C30	2.74	122.27	118.09
9	B8	101	BCL	CMA-C3A-C2A	-2.74	103.39	113.98
14	A5	103	CRT	C29-C28-C30	2.74	122.27	118.09
14	AW	102	CRT	C14-C15-C16	-2.74	115.27	123.20
9	A3	103	BCL	CHA-C1A-NA	-2.74	120.19	126.39
11	AL	304	UQ8	C46-C44-C43	-2.74	114.44	122.66
9	A0	102	BCL	CMA-C3A-C2A	-2.74	103.39	113.98
11	AL	304	UQ8	C30-C29-C31	2.74	119.98	115.23
9	AK	102	BCL	CMA-C3A-C2A	-2.74	103.40	113.98
9	A5	102	BCL	CBA-CAA-C2A	2.74	121.94	113.79
9	AG	101	BCL	C4D-CHA-C1A	2.74	124.51	121.24
9	BV	101	BCL	C3A-C2A-C1A	2.74	105.44	101.34
14	BM	406	CRT	C4-C5-C6	-2.74	121.07	124.91
9	AR	101	BCL	CAB-C3B-C2B	2.74	133.43	127.74
9	BL	303	BCL	CHB-C1B-NB	2.74	128.15	124.05
14	AJ	102	CRT	C8-C7-C6	2.74	122.27	118.09
14	AM	406	CRT	C4-C5-C6	-2.74	121.07	124.91
9	AE	101	BCL	OBB-CAB-CBB	-2.74	113.64	119.77
14	AJ	102	CRT	C34-C33-C35	2.73	122.26	118.09
9	B7	103	BCL	O2D-CGD-O1D	-2.73	118.53	123.85
9	AO	102	BCL	C2A-C1A-CHA	2.73	128.60	123.87
9	AK	102	BCL	C6-C5-C3	2.73	120.12	113.47
9	BS	102	BCL	CHB-C1B-NB	2.73	128.14	124.05
14	BN	102	CRT	C21-C20-C19	-2.73	117.94	123.52
9	A0	102	BCL	CAB-C3B-C2B	2.73	133.41	127.74
14	A0	101	CRT	C14-C15-C16	-2.73	115.30	123.20
14	AT	102	CRT	C29-C28-C30	2.73	122.25	118.09
7	AC	502	HEM	C2A-C1A-NA	2.73	113.18	110.15
9	B5	102	BCL	CHB-C1B-NB	2.72	128.14	124.05
9	AM	402	BCL	CHB-C1B-NB	2.72	128.13	124.05
9	BB	101	BCL	O2A-CGA-CBA	2.72	120.13	111.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
9	AS	103	BCL	C1C-NC-C4C	2.72	107.92	106.68
10	AM	403	BPH	C1A-C2A-C3A	-2.72	100.25	102.84
9	B4	101	BCL	CAB-C3B-C2B	2.72	133.39	127.74
9	A3	103	BCL	CBA-CAA-C2A	2.72	121.88	113.79
9	AW	101	BCL	CAB-C3B-C2B	2.72	133.39	127.74
9	BN	101	BCL	C6-C5-C3	2.72	120.09	113.47
9	A0	102	BCL	CHA-C1A-NA	-2.72	120.24	126.39
14	AT	102	CRT	C13-C12-C11	2.72	122.24	118.09
9	AK	102	BCL	CAB-C3B-C2B	2.72	133.38	127.74
9	A6	101	BCL	CMA-C3A-C2A	-2.71	103.49	113.98
9	BE	101	BCL	CHB-C1B-NB	2.71	128.12	124.05
9	BG	101	BCL	O2A-CGA-O1A	-2.71	116.84	123.63
9	A7	103	BCL	C3A-C2A-C1A	2.71	105.40	101.34
9	B4	101	BCL	CHB-C1B-NB	2.71	128.11	124.05
14	AJ	102	CRT	C14-C15-C16	-2.71	115.35	123.20
9	AF	102	BCL	C6-C5-C3	2.71	120.07	113.47
9	AL	301	BCL	C3A-C2A-C1A	2.71	105.39	101.34
9	AO	102	BCL	CAA-C2A-C1A	2.71	120.84	111.97
9	BG	101	BCL	C2C-C3C-C4C	2.71	105.39	101.34
9	BJ	101	BCL	C1-O2A-CGA	2.71	123.20	116.65
9	BU	102	BCL	CAB-C3B-C2B	2.70	133.36	127.74
10	BM	403	BPH	C2B-C1B-NB	-2.70	107.47	109.43
9	AK	102	BCL	CHA-C1A-NA	-2.70	120.27	126.39
9	AO	102	BCL	CMA-C3A-C2A	-2.70	103.53	113.98
14	AN	102	CRT	C18-C17-C16	2.70	122.22	118.09
9	AV	102	BCL	O2D-CGD-O1D	-2.70	118.59	123.85
14	AN	102	CRT	C30-C28-C27	-2.70	114.76	119.01
9	BB	101	BCL	C2B-C1B-NB	-2.70	107.53	110.33
9	BZ	101	BCL	CHB-C1B-NB	2.70	128.10	124.05
9	AN	101	BCL	C2A-C3A-C4A	2.70	106.23	101.87
9	B8	101	BCL	CHB-C1B-NB	2.70	128.10	124.05
14	BN	102	CRT	C35-C33-C32	-2.70	114.77	119.01
14	A0	101	CRT	C18-C17-C16	2.70	122.21	118.09
9	BG	101	BCL	O2D-CGD-O1D	-2.70	118.60	123.85
9	AA	101	BCL	C3A-C2A-C1A	2.70	105.38	101.34
9	BN	101	BCL	CHB-C1B-NB	2.70	128.09	124.05
9	B8	101	BCL	C3A-C2A-C1A	2.69	105.38	101.34
7	BC	502	HEM	C2A-C1A-NA	2.69	113.14	110.15
14	AB	102	CRT	C4-C5-C6	-2.69	121.13	124.91
9	A9	102	BCL	C2B-C1B-NB	-2.69	107.53	110.33
7	BC	501	HEM	C3B-C2B-C1B	-2.69	104.39	106.41
9	A6	101	BCL	CHA-C1A-NA	-2.69	120.30	126.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
7	BC	504	HEM	C2A-C1A-NA	2.69	113.14	110.15
9	AK	102	BCL	CBA-CAA-C2A	2.69	121.80	113.79
14	B1	103	CRT	C32-C31-C30	-2.69	115.41	123.20
9	BW	102	BCL	CHB-C1B-NB	2.69	128.08	124.05
9	AL	303	BCL	CHB-C1B-NB	2.69	128.08	124.05
14	A5	103	CRT	C15-C14-C12	-2.69	123.51	127.28
9	BK	102	BCL	C6-C5-C3	2.68	120.01	113.47
14	AS	104	CRT	C5-C6-C7	-2.68	121.84	125.89
14	BO	103	CRT	C13-C12-C11	2.68	122.19	118.09
9	A8	101	BCL	CHB-C1B-NB	2.68	128.07	124.05
9	BV	101	BCL	CMA-C3A-C2A	-2.68	103.62	113.98
9	BP	101	BCL	C4D-C3D-CAD	-2.68	105.19	108.11
14	BN	102	CRT	C30-C28-C27	-2.68	114.80	119.01
9	A7	103	BCL	CHA-C1A-NA	-2.68	120.33	126.39
14	B5	103	CRT	C8-C7-C6	2.68	122.18	118.09
9	A8	101	BCL	C2A-C1A-CHA	2.68	128.51	123.87
9	AX	101	BCL	C6-C5-C3	2.68	119.99	113.47
14	BS	103	CRT	C29-C28-C30	2.68	122.18	118.09
9	BA	101	BCL	O2D-CGD-O1D	-2.68	118.64	123.85
9	AF	102	BCL	C2B-C1B-NB	-2.67	107.55	110.33
9	AX	101	BCL	C2A-C3A-C4A	2.67	106.19	101.87
9	BS	102	BCL	CHC-C4B-NB	2.67	128.06	124.05
9	A9	102	BCL	CHA-C1A-NA	-2.67	120.34	126.39
9	BQ	104	BCL	CHA-C1A-NA	-2.67	120.34	126.39
9	B2	101	BCL	C2A-C1A-CHA	2.67	128.50	123.87
9	BS	102	BCL	C2B-C1B-NB	-2.67	107.56	110.33
14	BM	406	CRT	C26-C27-C28	-2.67	123.53	127.28
9	AO	102	BCL	C6-C5-C3	2.67	119.97	113.47
9	B5	102	BCL	CMA-C3A-C2A	-2.67	103.66	113.98
9	BA	101	BCL	C4D-CHA-C1A	2.67	124.43	121.24
9	BA	101	BCL	C3A-C2A-C1A	2.67	105.34	101.34
14	BF	103	CRT	C5-C6-C7	2.67	129.92	125.89
9	A3	104	BCL	CMA-C3A-C2A	-2.67	103.67	113.98
9	AN	101	BCL	CAC-C3C-C4C	-2.67	106.67	112.58
9	AL	301	BCL	CAA-C2A-C1A	2.67	120.71	111.97
14	AP	102	CRT	C3-C1-C2	-2.67	105.55	110.43
9	A5	102	BCL	CHA-C1A-NA	-2.66	120.36	126.39
14	BW	103	CRT	C1M-O1-C1	2.66	134.08	117.25
9	AA	101	BCL	C6-C5-C3	2.66	119.95	113.47
7	BC	504	HEM	C3B-C2B-C1B	-2.66	104.41	106.41
7	AC	503	HEM	C2A-C1A-NA	2.66	113.10	110.15
9	B2	101	BCL	CHB-C1B-NB	2.66	128.03	124.05

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
7	AC	503	HEM	C3B-C2B-C1B	-2.66	104.42	106.41
9	BV	101	BCL	O2A-CGA-CBA	2.66	119.94	111.83
9	AL	301	BCL	CHB-C1B-NB	2.65	128.03	124.05
9	B2	101	BCL	C4D-CHA-C1A	2.65	124.41	121.24
14	B5	103	CRT	C29-C28-C30	2.65	122.14	118.09
9	BD	102	BCL	C2B-C1B-NB	-2.65	107.58	110.33
9	BP	101	BCL	C2B-C1B-NB	-2.65	107.58	110.33
9	AJ	101	BCL	CAB-C3B-C2B	2.65	133.25	127.74
9	BV	101	BCL	C1-O2A-CGA	2.65	123.07	116.65
9	A7	103	BCL	C4D-C3D-CAD	-2.65	105.23	108.11
9	AM	401	BCL	CHB-C1B-NB	2.65	128.03	124.05
14	B2	102	CRT	C11-C12-C14	-2.65	114.84	119.01
9	BV	101	BCL	C2B-C1B-NB	-2.65	107.58	110.33
9	B0	102	BCL	C1C-NC-C4C	2.65	107.89	106.68
9	AB	101	BCL	CHA-C1A-NA	-2.65	120.40	126.39
9	A3	104	BCL	CHA-C1A-NA	-2.65	120.40	126.39
14	B2	102	CRT	C35-C33-C32	-2.65	114.85	119.01
9	B3	102	BCL	CHA-C1A-NA	-2.64	120.40	126.39
9	AG	101	BCL	CAB-C3B-C2B	2.64	133.24	127.74
14	BF	103	CRT	C34-C33-C35	2.64	122.13	118.09
9	B0	102	BCL	C1-O2A-CGA	2.64	123.05	116.65
14	AR	102	CRT	C18-C17-C16	2.64	122.13	118.09
9	AK	102	BCL	O2D-CGD-O1D	-2.64	118.70	123.85
9	A9	102	BCL	CBA-CAA-C2A	2.64	121.65	113.79
14	AB	102	CRT	C8-C7-C6	2.64	122.12	118.09
9	AE	101	BCL	CHA-C1A-NA	-2.64	120.42	126.39
9	AF	102	BCL	CAB-C3B-C2B	2.64	133.22	127.74
10	AM	403	BPH	C4C-C3C-C2C	-2.64	100.33	102.84
9	AY	102	BCL	CMA-C3A-C2A	-2.64	103.79	113.98
9	B5	102	BCL	C1C-NC-C4C	2.63	107.88	106.68
9	BI	102	BCL	C2A-C1A-CHA	2.63	128.44	123.87
9	B9	102	BCL	C3A-C2A-C1A	2.63	105.28	101.34
14	BV	102	CRT	C10-C9-C7	-2.63	123.58	127.28
14	BF	103	CRT	C35-C33-C32	-2.63	114.87	119.01
9	A3	104	BCL	CHB-C1B-NB	2.63	127.99	124.05
14	BO	103	CRT	C18-C17-C16	2.63	122.10	118.09
9	AN	101	BCL	C6-C5-C3	2.63	119.87	113.47
7	AC	501	HEM	CHC-C1C-NC	2.63	127.31	124.45
9	A3	104	BCL	C2A-C3A-C4A	2.63	106.11	101.87
14	BN	102	CRT	C34-C33-C35	2.63	122.10	118.09
14	A7	102	CRT	C13-C12-C11	2.63	122.10	118.09
14	BP	102	CRT	C30-C28-C27	-2.62	114.88	119.01

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
9	B4	101	BCL	C2A-C1A-CHA	2.62	128.42	123.87
9	AT	101	BCL	CAA-CBA-CGA	-2.62	105.77	113.21
9	AT	101	BCL	C4D-C3D-CAD	-2.62	105.26	108.11
9	BL	301	BCL	C2B-C1B-NB	-2.62	107.61	110.33
9	B2	101	BCL	CAB-C3B-C2B	2.62	133.18	127.74
13	AM	405	MQ8	C19-C18-C20	2.62	119.77	115.23
9	AM	402	BCL	CAA-C2A-C3A	-2.62	105.93	113.00
9	AS	103	BCL	CHC-C4B-NB	2.61	127.97	124.05
9	AD	102	BCL	CHB-C1B-NB	2.61	127.97	124.05
9	BZ	101	BCL	CHA-C1A-NA	-2.61	120.48	126.39
14	BV	102	CRT	C32-C31-C30	-2.61	115.64	123.20
9	B9	102	BCL	C6-C5-C3	2.61	119.82	113.47
9	BB	101	BCL	O2D-CGD-O1D	-2.61	118.77	123.85
9	AR	101	BCL	C2B-C1B-NB	-2.61	107.62	110.33
13	BM	405	MQ8	C21-C22-C23	-2.61	121.65	127.62
9	B4	101	BCL	C4D-CHA-C1A	2.61	124.35	121.24
9	AA	101	BCL	CAB-C3B-C2B	2.61	133.16	127.74
9	AU	102	BCL	CBA-CAA-C2A	2.61	121.55	113.79
9	B5	102	BCL	C6-C5-C3	2.61	119.82	113.47
9	BM	401	BCL	O2D-CGD-O1D	-2.61	118.78	123.85
14	A2	102	CRT	C13-C12-C11	2.61	122.07	118.09
9	AK	102	BCL	CHB-C1B-NB	2.60	127.96	124.05
9	BK	102	BCL	C2A-C1A-CHA	2.60	128.39	123.87
9	AI	102	BCL	O2D-CGD-O1D	-2.60	118.78	123.85
9	BP	101	BCL	O2A-CGA-O1A	-2.60	117.12	123.63
9	BJ	101	BCL	O2D-CGD-O1D	-2.60	118.78	123.85
9	B0	102	BCL	CAB-C3B-C2B	2.60	133.15	127.74
14	A5	103	CRT	C8-C7-C6	2.60	122.06	118.09
9	AI	102	BCL	C3A-C2A-C1A	2.60	105.23	101.34
14	BU	103	CRT	C34-C33-C35	2.60	122.06	118.09
9	AV	102	BCL	CAB-C3B-C2B	2.60	133.14	127.74
9	BO	102	BCL	C2A-C1A-CHA	2.60	128.38	123.87
14	AX	102	CRT	C32-C31-C30	-2.60	115.68	123.20
9	B0	102	BCL	C6-C5-C3	2.60	119.79	113.47
9	BV	101	BCL	C2A-C1A-CHA	2.59	128.37	123.87
14	AG	102	CRT	C8-C7-C6	2.59	122.05	118.09
14	AS	104	CRT	C18-C17-C16	2.59	122.05	118.09
14	BS	103	CRT	C9-C10-C11	-2.59	115.69	123.20
14	BS	103	CRT	C32-C31-C30	-2.59	115.69	123.20
9	AZ	101	BCL	CHA-C1A-NA	-2.59	120.53	126.39
7	AC	504	HEM	C2A-C1A-NA	2.59	113.02	110.15
9	AQ	102	BCL	CHB-C1B-NB	2.59	127.93	124.05

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
14	BF	103	CRT	C18-C17-C16	2.59	122.04	118.09
9	BI	102	BCL	C2B-C1B-NB	-2.59	107.64	110.33
9	AZ	101	BCL	C2C-C3C-C4C	2.59	105.21	101.34
14	BG	102	CRT	C36-C35-C33	-2.59	121.98	125.89
9	A7	103	BCL	C2B-C1B-NB	-2.59	107.64	110.33
9	BX	101	BCL	C6-C5-C3	2.58	119.76	113.47
9	AQ	102	BCL	C6-C5-C3	2.58	119.76	113.47
14	BS	103	CRT	C14-C15-C16	-2.58	115.71	123.20
9	BG	101	BCL	C2A-C1A-CHA	2.58	128.35	123.87
9	B9	102	BCL	CMA-C3A-C2A	-2.58	104.00	113.98
14	BS	103	CRT	C8-C7-C6	2.58	122.03	118.09
7	BC	502	HEM	CHC-C1C-NC	2.58	127.26	124.45
9	A7	103	BCL	C2A-C1A-CHA	2.58	128.35	123.87
14	AN	102	CRT	C10-C9-C7	-2.58	123.66	127.28
9	BO	102	BCL	CHA-C1A-NA	-2.58	120.55	126.39
9	AP	101	BCL	O2A-CGA-CBA	2.58	119.70	111.83
9	B5	102	BCL	C2A-C1A-CHA	2.58	128.34	123.87
9	B5	102	BCL	CHA-C1A-NA	-2.58	120.55	126.39
9	AQ	102	BCL	CHA-C1A-NA	-2.58	120.56	126.39
9	AI	102	BCL	CAB-C3B-C2B	2.58	133.10	127.74
14	A5	103	CRT	C13-C12-C11	2.58	122.02	118.09
9	B5	102	BCL	C2B-C1B-NB	-2.57	107.66	110.33
9	A1	102	BCL	CHB-C1B-NB	2.57	127.91	124.05
9	A3	103	BCL	C2A-C1A-CHA	2.57	128.33	123.87
14	BW	103	CRT	C8-C7-C6	2.57	122.02	118.09
9	AL	301	BCL	C2B-C1B-NB	-2.57	107.66	110.33
9	AM	402	BCL	CAB-C3B-C2B	2.57	133.09	127.74
9	AT	101	BCL	CMA-C3A-C2A	-2.57	104.03	113.98
14	AG	102	CRT	C13-C12-C11	2.57	122.02	118.09
9	B3	102	BCL	C2B-C1B-NB	-2.57	107.66	110.33
9	B7	103	BCL	C2B-C1B-NB	-2.57	107.66	110.33
14	B0	101	CRT	C35-C33-C32	-2.57	114.97	119.01
9	AR	101	BCL	CHB-C1B-NB	2.57	127.90	124.05
14	AJ	102	CRT	C35-C33-C32	-2.57	114.97	119.01
9	BP	101	BCL	C6-C5-C3	2.57	119.72	113.47
9	A7	103	BCL	CBA-CAA-C2A	2.57	121.43	113.79
7	AC	504	HEM	C4B-C3B-C2B	-2.56	104.92	107.28
9	AQ	102	BCL	CAB-C3B-C2B	2.56	133.07	127.74
9	BA	101	BCL	CHB-C1B-NB	2.56	127.89	124.05
9	AV	102	BCL	CMA-C3A-C2A	-2.56	104.08	113.98
14	B2	102	CRT	C34-C33-C35	2.56	122.00	118.09
14	BA	102	CRT	C27-C26-C25	-2.56	115.78	123.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
14	BB	102	CRT	C35-C33-C32	-2.56	114.98	119.01
14	BO	103	CRT	C30-C28-C27	-2.56	114.98	119.01
7	BC	503	HEM	C4B-C3B-C2B	-2.56	104.93	107.28
9	AA	101	BCL	C2B-C1B-NB	-2.56	107.67	110.33
7	AC	504	HEM	CHC-C1C-NC	2.56	127.24	124.45
14	A7	102	CRT	C29-C28-C30	2.56	121.99	118.09
9	BE	101	BCL	O2A-CGA-O1A	-2.56	117.23	123.63
14	AT	102	CRT	C32-C31-C30	-2.56	115.80	123.20
10	AL	302	BPH	C4C-C3C-C2C	-2.55	100.41	102.84
9	B8	101	BCL	CHA-C1A-NA	-2.55	120.61	126.39
9	B6	101	BCL	CHB-C1B-NB	2.55	127.88	124.05
9	AP	101	BCL	C2B-C1B-NB	-2.55	107.68	110.33
9	BS	102	BCL	C3A-C2A-C1A	2.55	105.16	101.34
9	B7	103	BCL	CMA-C3A-C2A	-2.55	104.13	113.98
9	AX	101	BCL	C2B-C1B-NB	-2.55	107.68	110.33
9	AG	101	BCL	O2A-CGA-CBA	2.55	119.60	111.83
9	A3	103	BCL	CMA-C3A-C2A	-2.55	104.14	113.98
9	AY	102	BCL	C4D-C3D-CAD	-2.55	105.34	108.11
9	A5	102	BCL	CHB-C1B-NB	2.55	127.87	124.05
9	AJ	101	BCL	O2A-CGA-CBA	2.54	119.59	111.83
9	B9	102	BCL	CHA-C1A-NA	-2.54	120.63	126.39
9	AZ	101	BCL	C6-C5-C3	2.54	119.66	113.47
9	A9	102	BCL	C4D-C3D-CAD	-2.54	105.34	108.11
14	A0	101	CRT	C5-C6-C7	-2.54	122.05	125.89
14	AJ	102	CRT	C13-C12-C11	2.54	121.97	118.09
9	B6	101	BCL	CHA-C1A-NA	-2.54	120.64	126.39
9	A3	104	BCL	C2B-C1B-NB	-2.54	107.69	110.33
9	AL	301	BCL	CAB-C3B-C2B	2.54	133.01	127.74
9	A5	102	BCL	C4D-C3D-CAD	-2.54	105.35	108.11
15	AM	409	PEF	O3-C3-C2	2.54	115.71	108.40
9	AS	103	BCL	C2A-C1A-CHA	2.53	128.26	123.87
14	AJ	102	CRT	C18-C17-C16	2.53	121.96	118.09
14	AN	102	CRT	C27-C26-C25	-2.53	115.86	123.20
9	BP	101	BCL	CHA-C1A-NA	-2.53	120.66	126.39
9	A5	102	BCL	O2D-CGD-O1D	-2.53	118.92	123.85
14	AA	102	CRT	C29-C28-C30	2.53	121.95	118.09
9	AP	101	BCL	CHA-C1A-NA	-2.53	120.67	126.39
14	AP	102	CRT	C9-C10-C11	-2.53	115.88	123.20
13	AM	405	MQ8	C24-C23-C25	2.53	119.62	115.23
9	BW	102	BCL	C4D-C3D-CAD	-2.53	105.36	108.11
9	BS	102	BCL	O2D-CGD-O1D	-2.53	118.93	123.85
9	BP	101	BCL	O2A-CGA-CBA	2.53	119.54	111.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
9	AJ	101	BCL	C2A-C1A-CHA	2.53	128.25	123.87
14	AG	102	CRT	C27-C26-C25	-2.53	115.88	123.20
14	B7	102	CRT	C29-C28-C30	2.52	121.94	118.09
9	B2	101	BCL	O2D-CGD-O1D	-2.52	118.94	123.85
9	B7	103	BCL	CHB-C1B-NB	2.52	127.83	124.05
9	BE	101	BCL	CHA-C1A-NA	-2.52	120.68	126.39
9	AT	101	BCL	C2B-C1B-NB	-2.52	107.71	110.33
14	AX	102	CRT	C15-C16-C17	-2.52	119.46	126.36
9	AN	101	BCL	CMA-C3A-C2A	-2.52	104.25	113.98
9	BB	101	BCL	C2C-C3C-C4C	2.52	105.11	101.34
14	B1	103	CRT	C24-C23-C25	2.52	121.93	118.09
9	AW	101	BCL	O2D-CGD-O1D	-2.51	118.95	123.85
9	B6	101	BCL	C2A-C1A-CHA	2.51	128.23	123.87
9	BU	102	BCL	C6-C5-C3	2.51	119.59	113.47
9	A1	102	BCL	CHA-C1A-NA	-2.51	120.70	126.39
9	A3	103	BCL	C2B-C1B-NB	-2.51	107.72	110.33
9	A6	101	BCL	CHB-C1B-NB	2.51	127.81	124.05
9	B3	102	BCL	CHB-C1B-NB	2.51	127.81	124.05
9	AL	301	BCL	CHA-C1A-NA	-2.51	120.71	126.39
9	B9	102	BCL	C2B-C1B-NB	-2.51	107.72	110.33
9	A6	101	BCL	C2A-C1A-CHA	2.51	128.22	123.87
14	A7	102	CRT	C20-C19-C17	-2.51	123.76	127.28
14	AM	406	CRT	C5-C6-C7	-2.50	122.11	125.89
9	BZ	101	BCL	C2A-C3A-C4A	2.50	105.92	101.87
9	B6	101	BCL	O2A-CGA-O1A	-2.50	117.36	123.63
14	A0	101	CRT	C13-C12-C11	2.50	121.91	118.09
9	BK	102	BCL	O2D-CGD-O1D	-2.50	118.98	123.85
9	BI	102	BCL	CHB-C1B-NB	2.50	127.80	124.05
9	BG	101	BCL	C1-O2A-CGA	2.50	122.70	116.65
9	BD	102	BCL	C3C-C2C-C1C	2.50	105.91	101.87
9	BZ	101	BCL	C1-O2A-CGA	2.50	122.70	116.65
14	B1	103	CRT	C30-C28-C27	-2.50	115.08	119.01
14	A2	102	CRT	C32-C31-C30	-2.50	115.96	123.20
14	BG	102	CRT	C8-C7-C6	2.50	121.91	118.09
14	B2	102	CRT	C10-C9-C7	-2.50	123.78	127.28
9	AT	101	BCL	O2D-CGD-O1D	-2.50	118.99	123.85
9	BT	101	BCL	C2B-C1B-NB	-2.49	107.74	110.33
9	BL	301	BCL	C3A-C2A-C1A	2.49	105.07	101.34
9	AV	102	BCL	CHA-C1A-NA	-2.49	120.75	126.39
9	AL	303	BCL	CAC-C3C-C4C	-2.49	107.05	112.58
11	AL	304	UQ8	C15-C14-C16	2.49	119.55	115.23
9	BL	301	BCL	O2D-CGD-CBD	2.49	115.59	111.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
9	AB	101	BCL	O2D-CGD-O1D	-2.49	119.00	123.85
9	BI	102	BCL	O2D-CGD-O1D	-2.49	119.00	123.85
9	B6	101	BCL	CAA-CBA-CGA	-2.49	106.14	113.21
13	AM	405	MQ8	C39-C38-C40	2.49	119.55	115.23
14	BW	103	CRT	C20-C19-C17	-2.49	123.79	127.28
9	AQ	102	BCL	CMA-C3A-C2A	-2.49	104.36	113.98
9	BN	101	BCL	C2C-C3C-C4C	2.49	105.06	101.34
9	AX	101	BCL	CHB-C1B-NB	2.49	127.78	124.05
9	B0	102	BCL	CHB-C1B-NB	2.49	127.78	124.05
15	BM	407	PEF	C3-C2-C1	-2.48	105.99	111.78
9	AJ	101	BCL	CHB-C1B-NB	2.48	127.78	124.05
9	AA	101	BCL	CHB-C1B-NB	2.48	127.77	124.05
9	AS	103	BCL	CBA-CAA-C2A	2.48	121.18	113.79
9	B3	102	BCL	CMA-C3A-C2A	-2.48	104.39	113.98
9	B9	102	BCL	O2D-CGD-O1D	-2.48	119.02	123.85
9	A9	102	BCL	CAA-C2A-C3A	-2.48	106.30	113.00
9	AG	101	BCL	O2D-CGD-O1D	-2.48	119.02	123.85
14	BP	102	CRT	C18-C17-C16	2.48	121.87	118.09
9	A6	101	BCL	C2B-C1B-NB	-2.48	107.76	110.33
9	B9	102	BCL	CHB-C1B-NB	2.48	127.76	124.05
9	BZ	101	BCL	C2B-C1B-NB	-2.48	107.76	110.33
9	BM	402	BCL	O2D-CGD-O1D	-2.48	119.03	123.85
14	BB	102	CRT	C37-C36-C35	-2.47	121.44	124.91
14	B2	102	CRT	C32-C31-C30	-2.47	116.04	123.20
9	AU	102	BCL	C4D-C3D-CAD	-2.47	105.42	108.11
9	AT	101	BCL	O2A-CGA-O1A	-2.47	117.45	123.63
14	AS	104	CRT	C29-C28-C30	2.47	121.86	118.09
9	AL	303	BCL	CMA-C3A-C2A	-2.47	104.44	113.98
14	AT	102	CRT	C24-C23-C25	2.47	121.86	118.09
9	AL	303	BCL	CAB-C3B-C2B	2.47	132.87	127.74
9	BL	301	BCL	CHA-C1A-NA	-2.47	120.81	126.39
9	AY	102	BCL	CHA-C1A-NA	-2.46	120.81	126.39
9	B6	101	BCL	C1C-NC-C4C	2.46	107.80	106.68
7	AC	503	HEM	C1D-C2D-C3D	-2.46	104.39	106.98
9	A9	102	BCL	CHB-C1B-NB	2.46	127.74	124.05
14	AR	102	CRT	C30-C28-C27	-2.46	115.14	119.01
9	BU	102	BCL	C3A-C2A-C1A	2.46	105.03	101.34
14	AT	102	CRT	C9-C10-C11	-2.46	116.07	123.20
9	BJ	101	BCL	C1C-NC-C4C	2.46	107.80	106.68
9	B6	101	BCL	O2A-CGA-CBA	2.46	119.34	111.83
7	AC	501	HEM	C2A-C1A-NA	2.46	112.88	110.15
9	AX	101	BCL	CHA-C1A-NA	-2.46	120.82	126.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
14	BU	103	CRT	C1M-O1-C1	2.46	132.80	117.25
9	AF	102	BCL	CHB-C1B-NB	2.46	127.74	124.05
14	B5	103	CRT	C21-C22-C23	-2.46	123.83	127.28
14	AP	102	CRT	C18-C17-C16	2.46	121.84	118.09
9	B8	101	BCL	C2B-C1B-NB	-2.46	107.78	110.33
9	A2	101	BCL	CHA-C1A-NA	-2.45	120.83	126.39
9	A3	103	BCL	C4D-C3D-CAD	-2.45	105.44	108.11
9	BX	101	BCL	C1C-NC-C4C	2.45	107.80	106.68
14	AW	102	CRT	C20-C19-C17	-2.45	123.84	127.28
9	BU	102	BCL	C2A-C1A-CHA	2.45	128.12	123.87
9	AE	101	BCL	C4D-CHA-C1A	2.45	124.17	121.24
14	AS	104	CRT	C36-C35-C33	2.45	129.59	125.89
14	BA	102	CRT	C29-C28-C30	2.45	121.83	118.09
9	BZ	101	BCL	O2D-CGD-O1D	-2.45	119.08	123.85
9	AM	401	BCL	CAB-C3B-C2B	2.45	132.83	127.74
14	AP	102	CRT	C24-C23-C25	2.45	121.83	118.09
10	BM	403	BPH	C6-C5-C3	2.45	119.42	113.47
13	AM	405	MQ8	C11-C12-C13	-2.44	122.62	126.83
14	BN	102	CRT	C8-C7-C6	2.44	121.82	118.09
14	BW	103	CRT	C29-C28-C30	2.44	121.82	118.09
9	AT	101	BCL	C2A-C3A-C4A	2.44	105.82	101.87
9	B7	103	BCL	CHA-C1A-NA	-2.44	120.86	126.39
9	AG	101	BCL	CHB-C1B-NB	2.44	127.71	124.05
15	AM	409	PEF	C3-C2-C1	-2.44	106.09	111.78
9	A5	102	BCL	C2A-C1A-CHA	2.44	128.10	123.87
14	B1	103	CRT	C16-C17-C19	-2.44	115.17	119.01
9	BW	102	BCL	CAB-C3B-C2B	2.44	132.81	127.74
14	AJ	102	CRT	C30-C28-C27	-2.44	115.17	119.01
9	A1	102	BCL	C4D-C3D-CAD	-2.44	105.46	108.11
9	BQ	104	BCL	C2A-C1A-CHA	2.44	128.09	123.87
14	BS	103	CRT	C20-C21-C22	-2.44	118.54	123.52
9	A9	102	BCL	C2A-C1A-CHA	2.44	128.09	123.87
9	B4	101	BCL	C2A-C3A-C4A	2.43	105.80	101.87
15	AS	101	PEF	C3-C2-C1	-2.43	106.11	111.78
9	AE	101	BCL	CHB-C1B-NB	2.43	127.70	124.05
9	AJ	101	BCL	CHA-C1A-NA	-2.43	120.88	126.39
9	A0	102	BCL	C2A-C3A-C4A	2.43	105.80	101.87
14	BO	103	CRT	C27-C26-C25	-2.43	116.15	123.20
9	AK	102	BCL	C4D-C3D-CAD	-2.43	105.47	108.11
15	AM	407	PEF	C3-C2-C1	-2.43	106.12	111.78
9	A2	101	BCL	CHB-C1B-NB	2.43	127.69	124.05
9	AD	102	BCL	C6-C5-C3	2.43	119.38	113.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
14	AW	102	CRT	C29-C28-C30	2.43	121.80	118.09
14	BA	102	CRT	C13-C12-C11	2.43	121.80	118.09
14	AP	102	CRT	C30-C28-C27	-2.43	115.19	119.01
9	AW	101	BCL	CMA-C3A-C2A	-2.43	104.60	113.98
9	BL	301	BCL	C4D-CHA-C1A	2.43	124.14	121.24
9	AE	101	BCL	C2B-C1B-NB	-2.43	107.81	110.33
13	AM	405	MQ8	C36-C37-C38	-2.43	122.07	127.62
9	BE	101	BCL	O2D-CGD-O1D	-2.42	119.13	123.85
9	AI	102	BCL	C2B-C1B-NB	-2.42	107.81	110.33
9	B1	102	BCL	CHA-C1A-NA	-2.42	120.90	126.39
9	A3	103	BCL	CHB-C1B-NB	2.42	127.69	124.05
9	BT	101	BCL	C6-C5-C3	2.42	119.37	113.47
14	BV	102	CRT	C18-C17-C16	2.42	121.79	118.09
9	AG	101	BCL	C2A-C1A-CHA	2.42	128.07	123.87
14	AA	102	CRT	C34-C33-C35	2.42	121.79	118.09
9	B0	102	BCL	C4D-CHA-C1A	2.42	124.13	121.24
9	B1	102	BCL	CMA-C3A-C2A	-2.42	104.62	113.98
9	AL	303	BCL	C6-C5-C3	2.42	119.36	113.47
7	BC	502	HEM	C3B-C2B-C1B	-2.42	104.59	106.41
9	AI	102	BCL	CHB-C1B-NB	2.42	127.68	124.05
9	B8	101	BCL	C4D-CHA-C1A	2.42	124.13	121.24
9	AR	101	BCL	C6-C5-C3	2.42	119.36	113.47
14	BG	102	CRT	C13-C12-C11	2.42	121.78	118.09
9	BL	301	BCL	CAB-C3B-C2B	2.42	132.76	127.74
9	A0	102	BCL	C2A-C1A-CHA	2.42	128.06	123.87
9	AZ	101	BCL	CHB-C1B-NB	2.42	127.67	124.05
9	B0	102	BCL	O2D-CGD-O1D	-2.42	119.15	123.85
15	BQ	101	PEF	C3-C2-C1	-2.42	106.15	111.78
9	AB	101	BCL	CHB-C1B-NB	2.41	127.67	124.05
14	AA	102	CRT	C13-C12-C11	2.41	121.77	118.09
7	BC	501	HEM	C2A-C1A-NA	2.41	112.83	110.15
9	AM	402	BCL	O2D-CGD-O1D	-2.41	119.15	123.85
9	A2	101	BCL	CMA-C3A-C2A	-2.41	104.66	113.98
15	AH	301	PEF	C3-C2-C1	-2.41	106.17	111.78
14	AR	102	CRT	C27-C26-C25	-2.41	116.22	123.20
14	BN	102	CRT	C27-C26-C25	-2.41	116.22	123.20
9	B0	102	BCL	CAC-C3C-C4C	-2.41	107.24	112.58
11	BL	304	UQ8	C20-C19-C21	2.41	119.40	115.23
7	AC	502	HEM	CBC-CAC-C3C	2.40	139.54	127.53
11	BL	304	UQ8	C40-C39-C41	2.40	119.40	115.23
9	B3	102	BCL	CHC-C4B-NB	2.40	127.65	124.05
9	BQ	104	BCL	C2B-C1B-NB	-2.40	107.83	110.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
9	B2	101	BCL	CBA-CAA-C2A	2.40	120.94	113.79
9	AM	401	BCL	O2D-CGD-O1D	-2.40	119.17	123.85
9	AA	101	BCL	CMA-C3A-C2A	-2.40	104.70	113.98
9	BX	101	BCL	CMA-C3A-C2A	-2.40	104.70	113.98
14	AB	102	CRT	C24-C23-C25	2.40	121.75	118.09
9	AE	101	BCL	O2D-CGD-O1D	-2.40	119.18	123.85
9	AG	101	BCL	CHA-C1A-NA	-2.40	120.96	126.39
9	AP	101	BCL	C2A-C1A-CHA	2.40	128.03	123.87
14	BS	103	CRT	C11-C12-C14	-2.40	115.24	119.01
9	A1	102	BCL	CMA-C3A-C2A	-2.40	104.71	113.98
9	AZ	101	BCL	CMA-C3A-C2A	-2.40	104.71	113.98
9	AN	101	BCL	CHA-C1A-NA	-2.40	120.97	126.39
9	AA	101	BCL	C2A-C1A-CHA	2.40	128.02	123.87
14	BM	406	CRT	C15-C14-C12	-2.40	123.92	127.28
9	BM	402	BCL	CAB-C3B-C2B	2.39	132.72	127.74
9	AQ	102	BCL	C3A-C2A-C1A	2.39	104.93	101.34
14	AS	104	CRT	C14-C15-C16	-2.39	116.26	123.20
9	BL	303	BCL	C6-C5-C3	2.39	119.30	113.47
11	BL	304	UQ8	C15-C14-C16	2.39	119.38	115.23
14	BU	103	CRT	C13-C12-C14	-2.39	118.94	122.82
9	BD	102	BCL	C6-C5-C3	2.39	119.29	113.47
14	AW	102	CRT	C24-C23-C25	2.39	121.74	118.09
14	AW	102	CRT	C1M-O1-C1	2.39	132.37	117.25
7	BC	503	HEM	C1D-C2D-C3D	-2.39	104.47	106.98
11	AL	304	UQ8	C7-C8-C9	-2.39	122.72	126.83
9	BY	102	BCL	CAB-C3B-C2B	2.39	132.70	127.74
9	AB	101	BCL	C2B-C1B-NB	-2.39	107.85	110.33
9	A0	102	BCL	C2B-C1B-NB	-2.39	107.85	110.33
9	AF	102	BCL	CBA-CAA-C2A	2.39	120.89	113.79
14	BG	102	CRT	C21-C22-C23	-2.39	123.93	127.28
10	AM	403	BPH	OBD-CAD-C3D	-2.38	124.17	127.89
14	A7	102	CRT	C14-C15-C16	-2.38	116.29	123.20
9	BB	101	BCL	CAC-C3C-C4C	-2.38	107.29	112.58
14	AG	102	CRT	C34-C33-C35	2.38	121.73	118.09
13	AM	405	MQ8	C34-C33-C35	2.38	119.36	115.23
9	BT	101	BCL	CHB-C1B-NB	2.38	127.62	124.05
9	B6	101	BCL	C2B-C1B-NB	-2.38	107.86	110.33
14	BN	102	CRT	C13-C12-C11	2.38	121.72	118.09
9	AD	102	BCL	O2D-CGD-O1D	-2.38	119.21	123.85
14	BS	103	CRT	C27-C26-C25	-2.38	116.31	123.20
9	B5	102	BCL	O2D-CGD-O1D	-2.38	119.22	123.85
9	B2	101	BCL	CHA-C1A-NA	-2.38	121.01	126.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
9	AB	101	BCL	O2A-CGA-O1A	-2.38	117.68	123.63
9	AU	102	BCL	CHA-C1A-NA	-2.38	121.01	126.39
14	AG	102	CRT	C30-C28-C27	-2.38	115.27	119.01
9	BZ	101	BCL	CMA-C3A-C2A	-2.38	104.80	113.98
14	A2	102	CRT	C29-C28-C30	2.38	121.72	118.09
9	AK	102	BCL	C2B-C1B-NB	-2.37	107.86	110.33
9	AS	103	BCL	CMB-C2B-C1B	-2.37	121.81	125.42
7	BC	503	HEM	C3B-C2B-C1B	-2.37	104.63	106.41
9	BV	101	BCL	O2D-CGD-O1D	-2.37	119.23	123.85
14	A5	103	CRT	C27-C26-C25	-2.37	116.33	123.20
9	AS	103	BCL	O2D-CGD-O1D	-2.37	119.24	123.85
9	BQ	103	BCL	O1D-CGD-CBD	-2.37	119.84	124.52
9	BW	102	BCL	CHC-C4B-NB	2.37	127.60	124.05
9	B0	102	BCL	CMA-C3A-C2A	-2.37	104.83	113.98
9	A1	102	BCL	C1C-NC-C4C	2.37	107.76	106.68
9	AN	101	BCL	O2D-CGD-O1D	-2.36	119.25	123.85
9	AL	301	BCL	C4D-CHA-C1A	2.36	124.06	121.24
9	AW	101	BCL	CHA-C1A-NA	-2.36	121.04	126.39
14	BV	102	CRT	C13-C12-C11	2.36	121.69	118.09
9	B6	101	BCL	C4D-CHA-C1A	2.36	124.06	121.24
14	AG	102	CRT	C14-C15-C16	-2.36	116.36	123.20
7	BC	503	HEM	C2A-C1A-NA	2.36	112.77	110.15
9	BO	102	BCL	O1D-CGD-CBD	-2.36	119.87	124.52
9	BZ	101	BCL	C2A-C1A-CHA	2.36	127.96	123.87
14	BU	103	CRT	C21-C22-C23	-2.36	123.97	127.28
14	AG	102	CRT	C21-C20-C19	-2.36	118.70	123.52
9	AW	101	BCL	C2B-C1B-NB	-2.36	107.88	110.33
9	BG	101	BCL	CAA-C2A-C1A	2.35	119.69	111.97
9	BT	101	BCL	O2D-CGD-O1D	-2.35	119.27	123.85
14	AP	102	CRT	C8-C7-C6	2.35	121.68	118.09
14	A2	102	CRT	C34-C33-C35	2.35	121.68	118.09
9	B0	102	BCL	C2B-C1B-NB	-2.35	107.89	110.33
9	AE	101	BCL	CMA-C3A-C2A	-2.35	104.90	113.98
9	BO	102	BCL	C3A-C2A-C1A	2.35	104.86	101.34
9	AE	101	BCL	O2A-CGA-CBA	2.35	118.99	111.83
14	B1	103	CRT	C34-C33-C35	2.35	121.67	118.09
9	BA	101	BCL	CMA-C3A-C2A	-2.35	104.91	113.98
14	AX	102	CRT	C13-C12-C14	-2.35	119.02	122.82
9	BM	402	BCL	CMA-C3A-C2A	-2.35	104.92	113.98
9	BL	303	BCL	CAB-C3B-C2B	2.34	132.61	127.74
9	AI	102	BCL	CMA-C3A-C2A	-2.34	104.92	113.98
14	BW	103	CRT	C3-C1-C2	-2.34	106.14	110.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
14	BP	102	CRT	C24-C23-C25	2.34	121.67	118.09
14	AB	102	CRT	C34-C33-C35	2.34	121.67	118.09
9	BG	101	BCL	C2B-C1B-NB	-2.34	107.90	110.33
14	AT	102	CRT	C27-C26-C25	-2.34	116.41	123.20
9	AX	101	BCL	C3C-C2C-C1C	2.34	105.65	101.87
10	AL	302	BPH	O2D-CGD-CBD	-2.34	108.38	110.95
9	BM	402	BCL	C2B-C1B-NB	-2.34	107.90	110.33
9	A6	101	BCL	O2A-C1-C2	2.34	117.11	108.11
9	BO	102	BCL	CAB-C3B-C2B	2.34	132.60	127.74
9	AW	101	BCL	C2A-C1A-CHA	2.33	127.92	123.87
9	AT	101	BCL	CAB-C3B-C2B	2.33	132.59	127.74
14	BB	102	CRT	C38-C37-C36	-2.33	106.92	112.98
9	AM	401	BCL	C6-C5-C3	2.33	119.15	113.47
14	A7	102	CRT	C32-C31-C30	-2.33	116.44	123.20
9	A1	102	BCL	C3C-C2C-C1C	2.33	105.64	101.87
11	BL	304	UQ8	C7-C8-C9	-2.33	122.81	126.83
9	A5	102	BCL	C2B-C1B-NB	-2.33	107.91	110.33
9	B2	101	BCL	C4D-C3D-CAD	-2.33	105.57	108.11
9	BF	102	BCL	C3C-C2C-C1C	2.33	105.64	101.87
9	A3	104	BCL	CAC-C3C-C4C	-2.33	107.41	112.58
14	A1	103	CRT	C27-C26-C25	-2.33	116.44	123.20
9	AS	103	BCL	C6-C5-C3	2.33	119.15	113.47
14	AN	102	CRT	C8-C7-C6	2.33	121.65	118.09
14	BS	103	CRT	C24-C23-C25	2.33	121.65	118.09
9	AF	102	BCL	O2D-CGD-O1D	-2.33	119.31	123.85
9	BT	101	BCL	C4D-C3D-CAD	-2.33	105.58	108.11
9	AB	101	BCL	CMA-C3A-C2A	-2.33	104.98	113.98
9	B2	101	BCL	O2A-CGA-O1A	-2.32	117.81	123.63
9	BN	101	BCL	O2D-CGD-O1D	-2.32	119.32	123.85
14	BO	103	CRT	C34-C33-C35	2.32	121.64	118.09
9	BG	101	BCL	C2A-C3A-C4A	2.32	105.62	101.87
9	B4	101	BCL	CAC-C3C-C4C	-2.32	107.43	112.58
9	BE	101	BCL	CBB-CAB-C3B	-2.32	115.50	120.40
9	BE	101	BCL	C1C-NC-C4C	2.32	107.74	106.68
9	AR	101	BCL	CMA-C3A-C2A	-2.32	105.01	113.98
9	BB	101	BCL	CMA-C3A-C2A	-2.32	105.02	113.98
9	BJ	101	BCL	C4D-C3D-CAD	-2.32	105.59	108.11
14	BU	103	CRT	C30-C28-C27	-2.32	115.36	119.01
14	AN	102	CRT	C13-C12-C11	2.32	121.63	118.09
9	BB	101	BCL	CAB-C3B-C2B	2.32	132.56	127.74
9	BB	101	BCL	C4D-CHA-C1A	2.32	124.01	121.24
9	BX	101	BCL	C2B-C1B-NB	-2.32	107.92	110.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
9	BP	101	BCL	C3A-C2A-C1A	2.31	104.81	101.34
9	AV	102	BCL	C2B-C1B-NB	-2.31	107.93	110.33
9	BE	101	BCL	C2B-C1B-NB	-2.31	107.93	110.33
9	AF	102	BCL	CMA-C3A-C2A	-2.31	105.04	113.98
14	AW	102	CRT	C26-C27-C28	-2.31	124.04	127.28
9	BM	401	BCL	CAB-C3B-C2B	2.31	132.54	127.74
14	B2	102	CRT	C29-C28-C30	2.31	121.62	118.09
14	B2	102	CRT	C27-C26-C25	-2.31	116.50	123.20
9	B4	101	BCL	CMA-C3A-C2A	-2.31	105.06	113.98
14	AN	102	CRT	C16-C17-C19	-2.31	115.38	119.01
14	AR	102	CRT	C34-C33-C35	2.31	121.61	118.09
9	BT	101	BCL	CAB-C3B-C2B	2.31	132.53	127.74
9	B2	101	BCL	CMA-C3A-C2A	-2.31	105.07	113.98
14	AW	102	CRT	C21-C22-C23	-2.30	124.05	127.28
9	BV	101	BCL	CAA-CBA-CGA	-2.30	106.67	113.21
9	AU	102	BCL	O2D-CGD-O1D	-2.30	119.36	123.85
9	AZ	101	BCL	CAC-C3C-C4C	-2.30	107.48	112.58
9	AQ	102	BCL	O2D-CGD-O1D	-2.30	119.37	123.85
9	AG	101	BCL	C2B-C1B-NB	-2.30	107.94	110.33
9	BF	102	BCL	O2D-CGD-O1D	-2.30	119.37	123.85
9	AS	103	BCL	CHA-C1A-NA	-2.30	121.19	126.39
15	AM	408	PEF	O3P-C1-C2	2.30	117.58	109.45
9	B4	101	BCL	C2B-C1B-NB	-2.30	107.94	110.33
9	AG	101	BCL	CMA-C3A-C2A	-2.30	105.11	113.98
9	B1	102	BCL	C2A-C1A-CHA	2.30	127.85	123.87
14	A5	103	CRT	C30-C28-C27	-2.29	115.40	119.01
7	BC	501	HEM	CBC-CAC-C3C	2.29	138.98	127.53
9	BI	102	BCL	CHC-C4B-NB	2.29	127.49	124.05
14	BG	102	CRT	C14-C15-C16	-2.29	116.56	123.20
9	AP	101	BCL	C2A-C3A-C4A	2.29	105.57	101.87
9	A1	102	BCL	C2B-C1B-NB	-2.29	107.95	110.33
9	A0	102	BCL	CHB-C1B-NB	2.29	127.49	124.05
9	AU	102	BCL	CMA-C3A-C2A	-2.29	105.13	113.98
9	BN	101	BCL	CHA-C1A-NA	-2.29	121.21	126.39
14	AB	102	CRT	C5-C6-C7	-2.29	122.43	125.89
9	AY	102	BCL	C1C-NC-C4C	2.29	107.72	106.68
14	BV	102	CRT	C21-C22-C23	-2.29	124.07	127.28
9	BW	102	BCL	C6-C5-C3	2.29	119.04	113.47
9	AX	101	BCL	CMA-C3A-C2A	-2.29	105.14	113.98
9	BA	101	BCL	C2A-C1A-CHA	2.28	127.83	123.87
9	B6	101	BCL	CMA-C3A-C2A	-2.28	105.15	113.98
9	AD	102	BCL	CMA-C3A-C2A	-2.28	105.16	113.98

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
14	B5	103	CRT	C13-C12-C11	2.28	121.58	118.09
9	BN	101	BCL	C2A-C3A-C4A	2.28	105.56	101.87
14	BG	102	CRT	C9-C10-C11	-2.28	116.59	123.20
9	BN	101	BCL	CMA-C3A-C2A	-2.28	105.17	113.98
9	BJ	101	BCL	C2A-C1A-CHA	2.28	127.82	123.87
9	AV	102	BCL	C3C-C2C-C1C	2.28	105.55	101.87
14	BW	103	CRT	C24-C23-C25	2.28	121.57	118.09
9	AO	102	BCL	C2B-C1B-NB	-2.28	107.96	110.33
14	AJ	102	CRT	C27-C26-C25	-2.28	116.60	123.20
9	BZ	101	BCL	CHC-C4B-NB	2.28	127.47	124.05
9	BL	303	BCL	CMA-C3A-C2A	-2.28	105.18	113.98
14	AJ	102	CRT	C9-C10-C11	-2.28	116.60	123.20
14	AM	406	CRT	C36-C35-C33	-2.28	122.45	125.89
14	AA	102	CRT	C27-C26-C25	-2.28	116.61	123.20
9	AJ	101	BCL	CMA-C3A-C2A	-2.28	105.18	113.98
14	AX	102	CRT	C11-C12-C14	2.28	122.59	119.01
14	BB	102	CRT	C13-C12-C11	2.27	121.56	118.09
14	A1	103	CRT	C3-C1-C2	-2.27	106.27	110.43
11	AL	304	UQ8	C10-C9-C11	2.27	119.17	115.23
11	BL	304	UQ8	C25-C24-C26	2.27	119.17	115.23
9	BD	102	BCL	CAB-C3B-C2B	2.27	132.46	127.74
14	BG	102	CRT	C34-C33-C35	2.27	121.56	118.09
9	A8	101	BCL	C4D-C3D-CAD	-2.27	105.64	108.11
9	B1	102	BCL	C2B-C1B-NB	-2.27	107.97	110.33
9	BT	101	BCL	CMA-C3A-C2A	-2.27	105.20	113.98
9	BZ	101	BCL	C4D-CHA-C1A	2.27	123.95	121.24
9	BY	102	BCL	C3C-C2C-C1C	2.27	105.54	101.87
14	AW	102	CRT	C32-C31-C30	-2.27	116.62	123.20
9	AJ	101	BCL	C2B-C1B-NB	-2.27	107.97	110.33
9	BX	101	BCL	CHB-C1B-NB	2.27	127.45	124.05
7	BC	503	HEM	C4A-C3A-C2A	2.27	109.41	106.82
14	AG	102	CRT	C18-C17-C16	2.27	121.55	118.09
14	AS	104	CRT	C13-C12-C11	2.27	121.55	118.09
9	AM	401	BCL	CMA-C3A-C2A	-2.27	105.22	113.98
14	BP	102	CRT	C1M-O1-C1	2.27	131.58	117.25
14	B5	103	CRT	C27-C26-C25	-2.27	116.64	123.20
9	BW	102	BCL	CMA-C3A-C2A	-2.26	105.23	113.98
9	B9	102	BCL	CHC-C4B-NB	2.26	127.44	124.05
9	AG	101	BCL	O2A-CGA-O1A	-2.26	117.97	123.63
9	AG	101	BCL	C3C-C2C-C1C	2.26	105.52	101.87
9	BL	303	BCL	CBA-CAA-C2A	-2.26	107.07	113.79
14	A5	103	CRT	C21-C20-C19	-2.26	118.89	123.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
7	AC	501	HEM	CBC-CAC-C3C	2.26	138.82	127.53
9	A9	102	BCL	CAC-C3C-C4C	-2.26	107.57	112.58
14	A5	103	CRT	C4-C5-C6	-2.26	121.74	124.91
14	B0	101	CRT	C21-C22-C23	-2.26	124.11	127.28
14	AG	102	CRT	C9-C10-C11	-2.26	116.66	123.20
9	BT	101	BCL	CAC-C3C-C4C	-2.26	107.57	112.58
9	AV	102	BCL	O2A-CGA-O1A	-2.26	117.98	123.63
14	BG	102	CRT	C21-C20-C19	-2.25	118.91	123.52
14	A7	102	CRT	C24-C23-C25	2.25	121.53	118.09
14	B0	101	CRT	C9-C10-C11	-2.25	116.67	123.20
9	AD	102	BCL	C2B-C1B-NB	-2.25	107.99	110.33
14	AP	102	CRT	C6-C7-C9	-2.25	115.47	119.01
9	BD	102	BCL	CBA-CAA-C2A	2.25	120.49	113.79
9	AO	102	BCL	CHB-C1B-NB	2.25	127.43	124.05
9	AO	102	BCL	O2D-CGD-O1D	-2.25	119.47	123.85
14	BN	102	CRT	C18-C17-C16	2.25	121.53	118.09
14	AP	102	CRT	C21-C20-C19	-2.25	118.92	123.52
14	B7	102	CRT	C34-C33-C35	2.25	121.52	118.09
7	BC	504	HEM	CHC-C1C-NC	2.25	126.90	124.45
9	AZ	101	BCL	C1C-NC-C4C	2.25	107.70	106.68
9	B1	102	BCL	CBB-CAB-C3B	-2.25	115.67	120.40
10	AM	403	BPH	C6-C5-C3	2.25	118.94	113.47
14	B2	102	CRT	C3-C1-C2	-2.25	106.32	110.43
9	BF	102	BCL	CHC-C4B-NB	2.24	127.42	124.05
7	BC	504	HEM	CBC-CAC-C3C	2.24	138.74	127.53
14	BG	102	CRT	C18-C17-C16	2.24	121.52	118.09
9	AM	402	BCL	CMA-C3A-C2A	-2.24	105.31	113.98
9	AD	102	BCL	CAB-C3B-C2B	2.24	132.40	127.74
9	AJ	101	BCL	C3C-C2C-C1C	2.24	105.49	101.87
7	BC	502	HEM	CBC-CAC-C3C	2.24	138.73	127.53
9	A0	102	BCL	CAC-C3C-C4C	-2.24	107.61	112.58
9	BU	102	BCL	C3C-C2C-C1C	2.24	105.49	101.87
14	AX	102	CRT	C8-C7-C9	-2.24	119.18	122.82
14	AM	406	CRT	C13-C12-C11	2.24	121.51	118.09
9	AL	301	BCL	O2D-CGD-CBD	2.24	115.15	111.23
9	AN	101	BCL	C2B-C1B-NB	-2.24	108.01	110.33
14	B5	103	CRT	C9-C10-C11	-2.24	116.72	123.20
9	BQ	103	BCL	CHB-C1B-NB	2.24	127.40	124.05
9	BK	102	BCL	C1C-NC-C4C	2.24	107.70	106.68
9	AV	102	BCL	C2A-C3A-C4A	2.23	105.48	101.87
13	AM	405	MQ8	C41-C42-C43	-2.23	122.51	127.62
14	BU	103	CRT	C15-C16-C17	-2.23	120.24	126.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
9	B6	101	BCL	C2A-C3A-C4A	2.23	105.47	101.87
14	BV	102	CRT	C35-C33-C32	-2.23	115.50	119.01
9	AD	102	BCL	O1D-CGD-CBD	-2.23	120.12	124.52
14	AG	102	CRT	C24-C23-C25	2.23	121.49	118.09
9	BQ	104	BCL	CHC-C4B-NB	2.23	127.39	124.05
14	AW	102	CRT	C8-C7-C6	2.23	121.49	118.09
11	BL	304	UQ8	C35-C34-C36	2.23	119.09	115.23
7	BC	501	HEM	C4A-C3A-C2A	2.23	109.37	106.82
9	A9	102	BCL	C1C-NC-C4C	2.22	107.69	106.68
14	BV	102	CRT	C30-C28-C27	-2.22	115.51	119.01
9	BQ	104	BCL	CHB-C1B-NB	2.22	127.38	124.05
9	B0	102	BCL	CBA-CAA-C2A	2.22	120.41	113.79
9	A6	101	BCL	CAC-C3C-C4C	-2.22	107.65	112.58
9	A2	101	BCL	C2A-C3A-C4A	2.22	105.46	101.87
9	BT	101	BCL	CHC-C4B-C3B	-2.22	122.90	131.72
9	A3	103	BCL	C3C-C2C-C1C	2.22	105.45	101.87
9	BM	401	BCL	C6-C5-C3	2.22	118.87	113.47
9	AE	101	BCL	C3A-C2A-C1A	2.22	104.66	101.34
9	A7	103	BCL	CHB-C1B-NB	2.22	127.38	124.05
14	AW	102	CRT	C10-C9-C7	-2.22	124.17	127.28
14	BF	103	CRT	C14-C15-C16	-2.22	116.78	123.20
13	BM	405	MQ8	C19-C18-C20	2.22	119.08	115.23
14	BP	102	CRT	C9-C10-C11	-2.22	116.78	123.20
9	BK	102	BCL	CBA-CAA-C2A	2.22	120.39	113.79
9	BO	102	BCL	CBA-CAA-C2A	2.22	120.39	113.79
14	A2	102	CRT	C21-C22-C23	-2.22	124.17	127.28
7	AC	503	HEM	C4A-C3A-C2A	2.22	109.36	106.82
9	AB	101	BCL	C4D-CHA-C1A	2.21	123.89	121.24
14	AN	102	CRT	C35-C33-C32	-2.21	115.53	119.01
14	A5	103	CRT	C24-C23-C25	2.21	121.47	118.09
14	AX	102	CRT	C18-C17-C19	-2.21	119.23	122.82
9	BY	102	BCL	CMA-C3A-C2A	-2.21	105.44	113.98
9	BM	402	BCL	CHA-C1A-NA	-2.21	121.39	126.39
9	BW	102	BCL	C2A-C1A-CHA	2.21	127.70	123.87
9	AM	401	BCL	C2A-C3A-C4A	2.21	105.44	101.87
9	AL	301	BCL	CMA-C3A-C2A	-2.21	105.44	113.98
9	AI	102	BCL	C3C-C2C-C1C	2.20	105.43	101.87
14	AW	102	CRT	C21-C20-C19	-2.20	119.01	123.52
9	BK	102	BCL	CHB-C1B-NB	2.20	127.35	124.05
14	AX	102	CRT	C20-C21-C22	-2.20	119.01	123.52
9	BK	102	BCL	C3C-C2C-C1C	2.20	105.43	101.87
9	AA	101	BCL	CHA-C1A-NA	-2.20	121.41	126.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
9	AW	101	BCL	CHB-C1B-NB	2.20	127.35	124.05
9	BQ	104	BCL	C6-C5-C3	2.20	118.83	113.47
13	AM	405	MQ8	C26-C27-C28	-2.20	122.59	127.62
9	BP	101	BCL	CMA-C3A-C2A	-2.20	105.47	113.98
9	BI	102	BCL	CBB-CAB-C3B	-2.20	115.76	120.40
9	AN	101	BCL	C4D-CHA-C1A	2.20	123.87	121.24
14	BW	103	CRT	C27-C26-C25	-2.20	116.83	123.20
9	BX	101	BCL	CHC-C4B-NB	2.20	127.35	124.05
14	AW	102	CRT	C1-C4-C5	2.20	118.69	112.98
7	BC	504	HEM	CBD-CAD-C3D	2.20	118.61	112.53
10	AL	302	BPH	CMD-C2D-C3D	2.20	129.08	124.68
14	AT	102	CRT	C11-C12-C14	-2.20	115.55	119.01
14	BN	102	CRT	C14-C15-C16	-2.20	116.83	123.20
9	AO	102	BCL	C3A-C2A-C1A	2.20	104.63	101.34
14	A5	103	CRT	C9-C10-C11	-2.20	116.83	123.20
7	AC	503	HEM	CHC-C1C-NC	2.20	126.85	124.45
9	BW	102	BCL	CBA-CAA-C2A	2.20	120.33	113.79
7	BC	502	HEM	C4A-C3A-C2A	2.20	109.33	106.82
9	AM	401	BCL	C2B-C1B-NB	-2.20	108.05	110.33
9	BN	101	BCL	C2A-C1A-CHA	2.20	127.68	123.87
9	AY	102	BCL	C3A-C2A-C1A	2.20	104.63	101.34
9	AF	102	BCL	CHA-C1A-NA	-2.19	121.42	126.39
9	AU	102	BCL	O1D-CGD-CBD	-2.19	120.19	124.52
9	BI	102	BCL	CAB-C3B-C2B	2.19	132.30	127.74
9	B6	101	BCL	CAC-C3C-C4C	-2.19	107.72	112.58
14	A7	102	CRT	C27-C26-C25	-2.19	116.85	123.20
9	BV	101	BCL	CHA-C1A-NA	-2.19	121.43	126.39
9	B5	102	BCL	CHC-C4B-NB	2.19	127.34	124.05
9	BQ	104	BCL	O2D-CGD-O1D	-2.19	119.58	123.85
14	BM	406	CRT	C24-C23-C25	2.19	121.43	118.09
7	AC	502	HEM	C4A-C3A-C2A	2.19	109.32	106.82
14	BP	102	CRT	C21-C20-C19	-2.19	119.05	123.52
9	A8	101	BCL	O2D-CGD-O1D	-2.18	119.59	123.85
9	BT	101	BCL	C3C-C2C-C1C	2.18	105.40	101.87
9	A2	101	BCL	CAB-C3B-C2B	2.18	132.28	127.74
14	BW	103	CRT	C32-C31-C30	-2.18	116.87	123.20
14	B5	103	CRT	C30-C28-C27	-2.18	115.57	119.01
14	BU	103	CRT	C35-C33-C32	-2.18	115.58	119.01
9	AM	401	BCL	CAA-CBA-CGA	2.18	119.41	113.21
9	AD	102	BCL	C3C-C2C-C1C	2.18	105.39	101.87
9	A8	101	BCL	CAC-C3C-C4C	-2.18	107.75	112.58
13	BM	405	MQ8	C24-C23-C25	2.18	119.01	115.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
14	AB	102	CRT	C10-C9-C7	-2.18	124.22	127.28
14	BO	103	CRT	C16-C17-C19	-2.18	115.58	119.01
14	AR	102	CRT	C16-C17-C19	-2.18	115.59	119.01
9	AS	103	BCL	C3C-C2C-C1C	2.18	105.38	101.87
9	BM	402	BCL	O2A-C1-C2	2.17	116.48	108.11
9	B0	102	BCL	C2A-C3A-C4A	2.17	105.38	101.87
14	B0	101	CRT	C36-C35-C33	-2.17	122.61	125.89
9	AT	101	BCL	CHC-C4B-NB	2.17	127.31	124.05
10	BM	403	BPH	C1-C2-C3	2.17	129.76	126.20
9	AF	102	BCL	C2A-C1A-CHA	2.17	127.64	123.87
14	B1	103	CRT	C8-C7-C9	-2.17	119.30	122.82
9	A0	102	BCL	C4D-C3D-CAD	-2.17	105.75	108.11
9	BL	303	BCL	CAC-C3C-C4C	-2.17	107.76	112.58
14	AA	102	CRT	C24-C23-C25	2.17	121.40	118.09
14	BA	102	CRT	C8-C7-C6	2.17	121.40	118.09
13	BM	405	MQ8	C41-C42-C43	-2.17	122.66	127.62
9	AP	101	BCL	CAC-C3C-C4C	-2.17	107.77	112.58
14	AW	102	CRT	C20-C21-C22	-2.17	119.09	123.52
14	AA	102	CRT	C21-C22-C23	-2.16	124.24	127.28
9	B1	102	BCL	CAA-C2A-C3A	-2.16	107.15	113.00
11	BL	304	UQ8	C4M-O4-C4	-2.16	108.86	116.47
9	AP	101	BCL	CHB-C1B-NB	2.16	127.30	124.05
7	BC	501	HEM	CHC-C1C-NC	2.16	126.81	124.45
9	BL	301	BCL	CMA-C3A-C2A	-2.16	105.62	113.98
9	BJ	101	BCL	CBB-CAB-C3B	-2.16	115.84	120.40
9	B1	102	BCL	C3A-C2A-C1A	2.16	104.58	101.34
9	BO	102	BCL	C4D-C3D-CAD	-2.16	105.76	108.11
9	AS	103	BCL	CAB-C3B-C2B	2.16	132.23	127.74
9	AX	101	BCL	C4D-C3D-CAD	-2.16	105.76	108.11
14	BM	406	CRT	C13-C12-C11	2.16	121.39	118.09
9	A5	102	BCL	CAC-C3C-C4C	-2.16	107.79	112.58
14	BA	102	CRT	C20-C21-C22	-2.16	119.10	123.52
14	AA	102	CRT	C8-C7-C6	2.16	121.38	118.09
10	BM	403	BPH	OBB-CAB-CBB	-2.16	115.59	120.19
9	BS	102	BCL	CHA-C1A-NA	-2.16	121.51	126.39
11	AL	304	UQ8	C4M-O4-C4	-2.16	108.89	116.47
14	B0	101	CRT	C31-C30-C28	-2.16	120.45	126.36
14	BF	103	CRT	C13-C12-C11	2.15	121.38	118.09
9	BK	102	BCL	CMA-C3A-C2A	-2.15	105.66	113.98
9	BP	101	BCL	O2D-CGD-O1D	-2.15	119.66	123.85
9	BV	101	BCL	C4D-CHA-C1A	2.15	123.81	121.24
9	B4	101	BCL	CHC-C4B-NB	2.15	127.28	124.05

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
14	A2	102	CRT	C35-C33-C32	-2.15	115.63	119.01
9	AU	102	BCL	CHB-C1B-NB	2.15	127.27	124.05
14	AW	102	CRT	C34-C33-C35	2.15	121.37	118.09
14	B5	103	CRT	C24-C23-C25	2.15	121.37	118.09
14	BP	102	CRT	C3-C1-C2	-2.15	106.50	110.43
9	AM	402	BCL	C2B-C1B-NB	-2.15	108.10	110.33
9	BD	102	BCL	CMA-C3A-C2A	-2.15	105.68	113.98
9	AM	402	BCL	CHC-C4B-NB	2.14	127.27	124.05
9	AE	101	BCL	CHC-C4B-NB	2.14	127.27	124.05
9	BG	101	BCL	CHA-C1A-NA	-2.14	121.54	126.39
9	BK	102	BCL	CHA-C1A-NA	-2.14	121.54	126.39
9	B5	102	BCL	C3C-C2C-C1C	2.14	105.33	101.87
9	BU	102	BCL	O1D-CGD-CBD	-2.14	120.29	124.52
14	BA	102	CRT	C24-C23-C25	2.14	121.36	118.09
9	AU	102	BCL	CAA-C2A-C3A	-2.14	107.21	113.00
9	BM	401	BCL	CAC-C3C-C4C	-2.14	107.84	112.58
9	B9	102	BCL	C4D-C3D-CAD	-2.14	105.78	108.11
14	BF	103	CRT	C24-C23-C25	2.14	121.36	118.09
14	B7	102	CRT	C13-C12-C11	2.14	121.36	118.09
9	BI	102	BCL	C3C-C2C-C1C	2.14	105.32	101.87
14	BG	102	CRT	C24-C23-C25	2.14	121.35	118.09
14	BA	102	CRT	C21-C22-C23	-2.14	124.28	127.28
9	BM	401	BCL	CMA-C3A-C2A	-2.14	105.72	113.98
14	BF	103	CRT	C27-C26-C25	-2.14	117.01	123.20
10	BL	302	BPH	OBB-CAB-CBB	-2.13	115.64	120.19
9	A7	103	BCL	O2D-CGD-O1D	-2.13	119.69	123.85
14	B0	101	CRT	C26-C25-C23	-2.13	120.51	126.36
9	AN	101	BCL	CHB-C1B-NB	2.13	127.25	124.05
9	BP	101	BCL	CMB-C2B-C1B	-2.13	122.17	125.42
9	B3	102	BCL	C3A-C2A-C1A	2.13	104.53	101.34
9	AA	101	BCL	C3C-C2C-C1C	2.13	105.31	101.87
14	BA	102	CRT	C14-C15-C16	-2.13	117.02	123.20
9	AL	301	BCL	C3C-C2C-C1C	2.13	105.31	101.87
13	BM	405	MQ8	C29-C28-C30	2.13	118.92	115.23
7	BC	502	HEM	C4B-C3B-C2B	-2.13	105.32	107.28
9	A3	104	BCL	C4D-C3D-CAD	-2.13	105.80	108.11
9	BQ	104	BCL	C3C-C2C-C1C	2.13	105.31	101.87
9	BP	101	BCL	CHC-C4B-NB	2.13	127.24	124.05
7	AC	503	HEM	CBC-CAC-C3C	2.13	138.15	127.53
9	AY	102	BCL	C2A-C1A-CHA	2.13	127.56	123.87
14	AW	102	CRT	C18-C17-C16	2.12	121.33	118.09
14	BW	103	CRT	C20-C21-C22	-2.12	119.17	123.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
9	BA	101	BCL	CHA-C1A-NA	-2.12	121.58	126.39
9	AB	101	BCL	CHC-C4B-NB	2.12	127.23	124.05
9	A6	101	BCL	C2A-C3A-C4A	2.12	105.30	101.87
9	B9	102	BCL	O1D-CGD-CBD	-2.12	120.33	124.52
9	BM	401	BCL	C2A-C3A-C4A	2.12	105.29	101.87
9	AU	102	BCL	C2B-C1B-NB	-2.12	108.13	110.33
14	A2	102	CRT	C27-C26-C25	-2.12	117.06	123.20
9	AM	402	BCL	CHA-C1A-NA	-2.12	121.60	126.39
14	AG	102	CRT	C20-C19-C17	-2.12	124.31	127.28
9	AE	101	BCL	C3C-C2C-C1C	2.12	105.29	101.87
9	B3	102	BCL	C3C-C2C-C1C	2.12	105.29	101.87
9	B5	102	BCL	C4D-C3D-CAD	-2.11	105.81	108.11
9	B6	101	BCL	C3C-C2C-C1C	2.11	105.28	101.87
9	AB	101	BCL	CBA-CAA-C2A	2.11	120.08	113.79
9	BY	102	BCL	C2A-C1A-CHA	2.11	127.53	123.87
14	B1	103	CRT	C11-C12-C14	-2.11	115.69	119.01
9	BK	102	BCL	CAB-C3B-C2B	2.11	132.13	127.74
9	AK	102	BCL	C1C-NC-C4C	2.11	107.64	106.68
9	BS	102	BCL	CHC-C4B-C3B	-2.11	123.34	131.72
9	AW	101	BCL	OBD-CAD-C3D	2.11	133.35	128.42
13	BM	405	MQ8	C34-C33-C35	2.11	118.89	115.23
9	B7	103	BCL	C3C-C2C-C1C	2.11	105.27	101.87
14	A0	101	CRT	C6-C7-C9	2.11	122.32	119.01
9	BM	402	BCL	C2A-C1A-CHA	2.11	127.52	123.87
14	A5	103	CRT	C18-C17-C16	2.11	121.31	118.09
9	BU	102	BCL	CBB-CAB-C3B	-2.11	115.96	120.40
9	AF	102	BCL	C3C-C2C-C1C	2.11	105.27	101.87
9	B1	102	BCL	C4D-C3D-CAD	-2.11	105.82	108.11
9	A1	102	BCL	CBA-CAA-C2A	2.11	120.06	113.79
9	BU	102	BCL	C4D-CHA-C1A	2.11	123.75	121.24
9	B9	102	BCL	C3C-C2C-C1C	2.10	105.27	101.87
14	BB	102	CRT	C24-C23-C25	2.10	121.30	118.09
14	B5	103	CRT	C34-C33-C35	2.10	121.30	118.09
7	BC	503	HEM	CHC-C1C-NC	2.10	126.74	124.45
14	AW	102	CRT	C9-C10-C11	-2.10	117.11	123.20
14	A7	102	CRT	C9-C10-C11	-2.10	117.11	123.20
14	BN	102	CRT	C9-C10-C11	-2.10	117.11	123.20
9	AR	101	BCL	CHC-C4B-NB	2.10	127.20	124.05
14	A5	103	CRT	C20-C21-C22	-2.10	119.22	123.52
9	B2	101	BCL	CAC-C3C-C4C	-2.10	107.92	112.58
14	B7	102	CRT	C35-C33-C32	-2.10	115.71	119.01
9	B6	101	BCL	CHC-C4B-NB	2.10	127.20	124.05

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
9	A9	102	BCL	C3C-C2C-C1C	2.10	105.26	101.87
14	B7	102	CRT	C27-C26-C25	-2.10	117.12	123.20
9	BE	101	BCL	C4D-C3D-CAD	-2.10	105.83	108.11
14	BA	102	CRT	C9-C10-C11	-2.10	117.12	123.20
9	AI	102	BCL	CHA-C1A-NA	-2.10	121.64	126.39
14	AR	102	CRT	C24-C23-C25	2.10	121.29	118.09
14	B2	102	CRT	C24-C23-C25	2.10	121.29	118.09
9	A2	101	BCL	C2B-C1B-NB	-2.10	108.15	110.33
9	BA	101	BCL	CHC-C4B-NB	2.10	127.19	124.05
9	B8	101	BCL	CHC-C4B-NB	2.09	127.19	124.05
9	BE	101	BCL	C4D-CHA-C1A	2.09	123.74	121.24
9	AM	402	BCL	CAC-C3C-C4C	-2.09	107.94	112.58
9	AP	101	BCL	CHC-C4B-C3B	-2.09	123.41	131.72
9	AL	303	BCL	C4D-C3D-CAD	-2.09	105.83	108.11
14	BM	406	CRT	C18-C17-C16	2.09	121.28	118.09
9	A5	102	BCL	C1C-NC-C4C	2.09	107.63	106.68
14	BA	102	CRT	C34-C33-C35	2.09	121.28	118.09
14	A5	103	CRT	C34-C33-C35	2.09	121.28	118.09
14	BB	102	CRT	C10-C9-C7	-2.09	124.35	127.28
9	B0	102	BCL	CHC-C4B-NB	2.09	127.18	124.05
11	BL	304	UQ8	C10-C9-C11	2.09	118.85	115.23
9	BV	101	BCL	CAC-C3C-C4C	-2.09	107.95	112.58
14	AS	104	CRT	C16-C17-C19	-2.09	115.72	119.01
9	AO	102	BCL	CHC-C4B-NB	2.09	127.18	124.05
9	BD	102	BCL	CHC-C1C-NC	2.09	127.41	124.40
9	AU	102	BCL	C2A-C3A-C4A	2.09	105.24	101.87
9	BQ	104	BCL	CMA-C3A-C2A	-2.09	105.91	113.98
9	B5	102	BCL	C3A-C2A-C1A	2.09	104.46	101.34
9	BL	303	BCL	CAA-CBA-CGA	2.09	119.13	113.21
9	BL	301	BCL	CAC-C3C-C4C	-2.09	107.96	112.58
9	A0	102	BCL	C3C-C2C-C1C	2.09	105.24	101.87
9	BV	101	BCL	CHC-C4B-NB	2.08	127.17	124.05
14	AN	102	CRT	C24-C23-C25	2.08	121.27	118.09
10	BL	302	BPH	C1A-C2A-C3A	-2.08	100.86	102.84
9	AV	102	BCL	CAC-C3C-C4C	-2.08	107.96	112.58
7	BC	503	HEM	CBC-CAC-C3C	2.08	137.93	127.53
9	BW	102	BCL	C1C-NC-C4C	2.08	107.63	106.68
9	AM	401	BCL	CHC-C4B-NB	2.08	127.17	124.05
9	AR	101	BCL	O2D-CGD-O1D	-2.08	119.80	123.85
14	AN	102	CRT	C11-C12-C14	-2.08	115.74	119.01
9	BW	102	BCL	CHA-C1A-NA	-2.08	121.68	126.39
9	AP	101	BCL	CMA-C3A-C2A	-2.08	105.94	113.98

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
9	AI	102	BCL	C4D-C3D-CAD	-2.08	105.85	108.11
9	AG	101	BCL	CHC-C4B-NB	2.08	127.17	124.05
13	BM	405	MQ8	C26-C27-C28	-2.08	122.87	127.62
9	BZ	101	BCL	CAC-C3C-C4C	-2.08	107.97	112.58
7	BC	502	HEM	CAA-C2A-C1A	2.08	129.00	124.94
9	BQ	103	BCL	C6-C5-C3	2.08	118.52	113.47
9	A3	104	BCL	C3C-C2C-C1C	2.07	105.22	101.87
9	AS	103	BCL	CHD-C1D-ND	-2.07	121.88	124.80
15	AH	301	PEF	C3-O3-C30	2.07	122.20	117.08
14	B7	102	CRT	C18-C17-C16	2.07	121.25	118.09
9	BP	101	BCL	CAA-CBA-CGA	-2.07	107.32	113.21
9	BQ	103	BCL	C3C-C2C-C1C	2.07	105.22	101.87
9	A9	102	BCL	CHC-C4B-NB	2.07	127.16	124.05
7	AC	504	HEM	CAC-C3C-C4C	2.07	129.77	124.82
9	AK	102	BCL	CAC-C3C-C4C	-2.07	107.99	112.58
9	BM	401	BCL	CHA-C1A-NA	-2.07	121.70	126.39
9	BN	101	BCL	C3C-C2C-C1C	2.07	105.21	101.87
11	AL	304	UQ8	C7-C6-C1	-2.07	121.34	124.89
9	BJ	101	BCL	C2B-C1B-NB	-2.07	108.18	110.33
14	BM	406	CRT	C27-C26-C25	-2.07	117.20	123.20
9	BT	101	BCL	C2A-C3A-C4A	2.07	105.21	101.87
9	AZ	101	BCL	CHC-C4B-NB	2.07	127.15	124.05
9	B8	101	BCL	CAC-C3C-C4C	-2.07	108.00	112.58
14	A7	102	CRT	C18-C17-C16	2.07	121.24	118.09
9	BQ	104	BCL	CHC-C4B-C3B	-2.07	123.52	131.72
14	BN	102	CRT	C24-C23-C25	2.06	121.24	118.09
9	A7	103	BCL	C3C-C2C-C1C	2.06	105.20	101.87
9	BM	401	BCL	CHC-C4B-NB	2.06	127.14	124.05
9	B7	103	BCL	C4D-C3D-CAD	-2.06	105.87	108.11
9	BI	102	BCL	O1D-CGD-CBD	-2.06	120.45	124.52
15	AM	407	PEF	C3-O3-C30	2.06	122.17	117.08
14	A2	102	CRT	C24-C23-C25	2.06	121.24	118.09
14	BW	103	CRT	C21-C22-C23	-2.06	124.39	127.28
9	A6	101	BCL	O2D-CGD-O1D	-2.06	119.84	123.85
11	BL	304	UQ8	C3M-O3-C3	-2.06	109.23	116.47
9	BL	303	BCL	C4D-C3D-CAD	-2.06	105.87	108.11
9	BD	102	BCL	CBB-CAB-C3B	-2.06	116.06	120.40
9	B3	102	BCL	CHC-C4B-C3B	-2.06	123.55	131.72
9	A5	102	BCL	C3C-C2C-C1C	2.06	105.19	101.87
14	BM	406	CRT	C21-C22-C23	-2.06	124.39	127.28
14	BO	103	CRT	C24-C23-C25	2.06	121.23	118.09
14	BF	103	CRT	C30-C28-C27	-2.06	115.78	119.01

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
14	A1	103	CRT	C24-C23-C25	2.05	121.23	118.09
9	BG	101	BCL	CMA-C3A-C2A	-2.05	106.04	113.98
14	AT	102	CRT	C18-C17-C16	2.05	121.22	118.09
14	AM	406	CRT	C24-C23-C25	2.05	121.22	118.09
11	AL	304	UQ8	C3M-O3-C3	-2.05	109.27	116.47
9	BJ	101	BCL	CMA-C3A-C2A	-2.05	106.07	113.98
9	BZ	101	BCL	C4D-C3D-CAD	-2.05	105.88	108.11
14	B5	103	CRT	C20-C21-C22	-2.05	119.33	123.52
9	BU	102	BCL	CHC-C4B-NB	2.05	127.12	124.05
9	BN	101	BCL	O1D-CGD-CBD	-2.04	120.49	124.52
9	AY	102	BCL	O2D-CGD-O1D	-2.04	119.87	123.85
9	AM	401	BCL	C3C-C2C-C1C	2.04	105.17	101.87
9	AU	102	BCL	C3C-C2C-C1C	2.04	105.17	101.87
9	AI	102	BCL	CHC-C4B-NB	2.04	127.11	124.05
14	BP	102	CRT	C16-C17-C19	-2.04	115.80	119.01
9	BJ	101	BCL	CHA-C1A-NA	-2.04	121.78	126.39
9	AD	102	BCL	CBB-CAB-C3B	-2.04	116.11	120.40
9	AN	101	BCL	C4D-C3D-CAD	-2.04	105.89	108.11
9	BV	101	BCL	C4D-C3D-CAD	-2.04	105.89	108.11
14	B1	103	CRT	C35-C33-C32	-2.04	115.81	119.01
7	AC	502	HEM	C4B-C3B-C2B	-2.04	105.41	107.28
9	BE	101	BCL	O2A-CGA-CBA	2.04	118.04	111.83
9	B7	103	BCL	CHC-C4B-NB	2.04	127.10	124.05
9	A5	102	BCL	CAA-C2A-C3A	-2.03	107.50	113.00
9	BN	101	BCL	C2B-C1B-NB	-2.03	108.22	110.33
13	BM	405	MQ8	C39-C38-C40	2.03	118.76	115.23
9	AO	102	BCL	CHC-C4B-C3B	-2.03	123.64	131.72
7	AC	504	HEM	CBC-CAC-C3C	2.03	137.68	127.53
9	AW	101	BCL	CHC-C4B-NB	2.03	127.10	124.05
9	BK	102	BCL	C4D-C3D-CAD	-2.03	105.90	108.11
9	AY	102	BCL	C6-C5-C3	2.03	118.42	113.47
9	A1	102	BCL	CAA-C2A-C3A	-2.03	107.51	113.00
9	AV	102	BCL	CHC-C4B-NB	2.03	127.10	124.05
14	AJ	102	CRT	C16-C17-C19	-2.03	115.81	119.01
9	AF	102	BCL	CHC-C4B-NB	2.03	127.10	124.05
9	B1	102	BCL	CAB-C3B-C2B	2.03	131.96	127.74
9	A9	102	BCL	O2D-CGD-O1D	-2.03	119.89	123.85
15	BM	407	PEF	C3-O3-C30	2.03	122.09	117.08
9	AD	102	BCL	CHA-C1A-NA	-2.03	121.79	126.39
9	AR	101	BCL	C2A-C1A-CHA	2.03	127.39	123.87
14	B2	102	CRT	C26-C27-C28	-2.03	124.43	127.28
9	BQ	103	BCL	CBB-CAB-C3B	-2.03	116.13	120.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
9	AK	102	BCL	C3C-C2C-C1C	2.03	105.14	101.87
14	BS	103	CRT	C30-C28-C27	-2.03	115.82	119.01
9	A8	101	BCL	C3A-C2A-C1A	2.03	104.38	101.34
9	A6	101	BCL	C4D-C3D-CAD	-2.03	105.91	108.11
14	AX	102	CRT	C18-C17-C16	2.03	121.18	118.09
9	AW	101	BCL	C1C-NC-C4C	2.03	107.60	106.68
9	BU	102	BCL	CHA-C1A-NA	-2.03	121.81	126.39
14	BS	103	CRT	C18-C17-C16	2.02	121.18	118.09
9	BS	102	BCL	C3C-C2C-C1C	2.02	105.14	101.87
9	BE	101	BCL	CHC-C4B-NB	2.02	127.08	124.05
9	BY	102	BCL	CHB-C1B-NB	2.02	127.08	124.05
14	AA	102	CRT	C30-C28-C27	-2.02	115.83	119.01
9	BF	102	BCL	CHC-C4B-C3B	-2.02	123.70	131.72
9	BE	101	BCL	CAC-C3C-C4C	-2.02	108.11	112.58
9	BM	401	BCL	CAA-CBA-CGA	2.02	118.94	113.21
14	BV	102	CRT	C27-C26-C25	-2.02	117.36	123.20
14	B0	101	CRT	C29-C28-C30	2.02	121.17	118.09
9	B2	101	BCL	CHC-C4B-NB	2.01	127.07	124.05
14	AG	102	CRT	C11-C12-C14	-2.01	115.84	119.01
9	BM	402	BCL	CHC-C4B-NB	2.01	127.07	124.05
9	BY	102	BCL	C2B-C1B-NB	-2.01	108.24	110.33
9	AV	102	BCL	C2A-C1A-CHA	2.01	127.36	123.87
9	AM	402	BCL	C3C-C2C-C1C	2.01	105.12	101.87
14	B0	101	CRT	C27-C26-C25	-2.01	117.37	123.20
14	BS	103	CRT	C21-C22-C23	-2.01	124.46	127.28
9	BM	401	BCL	C2B-C1B-NB	-2.01	108.24	110.33
9	AW	101	BCL	C3C-C2C-C1C	2.01	105.11	101.87
9	BT	101	BCL	C1-O2A-CGA	2.01	121.51	116.65
9	BM	402	BCL	C4D-C3D-CAD	-2.01	105.92	108.11
9	BY	102	BCL	CHA-C1A-NA	-2.01	121.84	126.39
9	BM	401	BCL	C3C-C2C-C1C	2.01	105.11	101.87
7	BC	502	HEM	CHB-C1B-NB	-2.00	121.89	124.37
9	BV	101	BCL	C3C-C2C-C1C	2.00	105.11	101.87
9	AJ	101	BCL	CHC-C4B-NB	2.00	127.06	124.05
10	BL	302	BPH	CMD-C2D-C3D	2.00	128.69	124.68
14	A2	102	CRT	C9-C10-C11	-2.00	117.40	123.20
9	A2	101	BCL	C3C-C2C-C1C	2.00	105.10	101.87
14	B7	102	CRT	C21-C22-C23	-2.00	124.47	127.28
14	BW	103	CRT	C9-C10-C11	-2.00	117.40	123.20
7	AC	502	HEM	C1D-C2D-C3D	-2.00	104.88	106.98

There are no chirality outliers.

All (1250) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	AC	501	HEM	C2B-C3B-CAB-CBB
7	AC	501	HEM	C4B-C3B-CAB-CBB
7	AC	501	HEM	C2C-C3C-CAC-CBC
7	AC	502	HEM	C2B-C3B-CAB-CBB
7	AC	502	HEM	C4B-C3B-CAB-CBB
7	AC	502	HEM	C2C-C3C-CAC-CBC
7	AC	502	HEM	C4C-C3C-CAC-CBC
7	AC	503	HEM	C2B-C3B-CAB-CBB
7	AC	503	HEM	C4B-C3B-CAB-CBB
7	AC	503	HEM	C2C-C3C-CAC-CBC
7	AC	504	HEM	C2C-C3C-CAC-CBC
7	AC	504	HEM	C4C-C3C-CAC-CBC
7	BC	501	HEM	C2B-C3B-CAB-CBB
7	BC	501	HEM	C4B-C3B-CAB-CBB
7	BC	501	HEM	C2C-C3C-CAC-CBC
7	BC	502	HEM	C2B-C3B-CAB-CBB
7	BC	502	HEM	C4B-C3B-CAB-CBB
7	BC	502	HEM	C2C-C3C-CAC-CBC
7	BC	502	HEM	C4C-C3C-CAC-CBC
7	BC	503	HEM	C2B-C3B-CAB-CBB
7	BC	503	HEM	C4B-C3B-CAB-CBB
7	BC	503	HEM	C2C-C3C-CAC-CBC
7	BC	503	HEM	C4C-C3C-CAC-CBC
7	BC	504	HEM	C2B-C3B-CAB-CBB
7	BC	504	HEM	C4B-C3B-CAB-CBB
7	BC	504	HEM	C2C-C3C-CAC-CBC
7	BC	504	HEM	C4C-C3C-CAC-CBC
9	AL	303	BCL	CAD-CBD-CGD-O1D
9	AL	303	BCL	CAD-CBD-CGD-O2D
9	AM	402	BCL	CAD-CBD-CGD-O2D
9	AA	101	BCL	O2A-C1-C2-C3
9	AB	101	BCL	C2C-C3C-CAC-CBC
9	AB	101	BCL	C4C-C3C-CAC-CBC
9	AE	101	BCL	C2C-C3C-CAC-CBC
9	AE	101	BCL	C4C-C3C-CAC-CBC
9	AF	102	BCL	C4C-C3C-CAC-CBC
9	AG	101	BCL	C4C-C3C-CAC-CBC
9	AI	102	BCL	C4C-C3C-CAC-CBC
9	AK	102	BCL	C4C-C3C-CAC-CBC
9	AN	101	BCL	C2C-C3C-CAC-CBC
9	AN	101	BCL	C4C-C3C-CAC-CBC
9	AO	102	BCL	C4C-C3C-CAC-CBC

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Mol	Chain	Res	Type	Atoms
9	AQ	102	BCL	C4C-C3C-CAC-CBC
9	AS	103	BCL	C4C-C3C-CAC-CBC
9	AT	101	BCL	C4C-C3C-CAC-CBC
9	AU	102	BCL	C1A-C2A-CAA-CBA
9	AU	102	BCL	C2C-C3C-CAC-CBC
9	AU	102	BCL	C4C-C3C-CAC-CBC
9	AV	102	BCL	C4C-C3C-CAC-CBC
9	AV	102	BCL	O2A-C1-C2-C3
9	AW	101	BCL	C2-C1-O2A-CGA
9	AW	101	BCL	C4C-C3C-CAC-CBC
9	AX	101	BCL	C2C-C3C-CAC-CBC
9	AX	101	BCL	C4C-C3C-CAC-CBC
9	AY	102	BCL	C4C-C3C-CAC-CBC
9	AZ	101	BCL	C1A-C2A-CAA-CBA
9	AZ	101	BCL	C3A-C2A-CAA-CBA
9	AZ	101	BCL	C2C-C3C-CAC-CBC
9	AZ	101	BCL	C4C-C3C-CAC-CBC
9	A1	102	BCL	C4C-C3C-CAC-CBC
9	A2	101	BCL	C1A-C2A-CAA-CBA
9	A2	101	BCL	C2C-C3C-CAC-CBC
9	A2	101	BCL	C4C-C3C-CAC-CBC
9	A3	103	BCL	C4C-C3C-CAC-CBC
9	A3	104	BCL	C2-C1-O2A-CGA
9	A3	104	BCL	C2C-C3C-CAC-CBC
9	A3	104	BCL	C4C-C3C-CAC-CBC
9	A6	101	BCL	C2C-C3C-CAC-CBC
9	A6	101	BCL	C4C-C3C-CAC-CBC
9	A9	102	BCL	O2A-C1-C2-C3
9	A0	102	BCL	C2C-C3C-CAC-CBC
9	A0	102	BCL	C4C-C3C-CAC-CBC
9	BL	301	BCL	C2C-C3C-CAC-CBC
9	BL	301	BCL	C4C-C3C-CAC-CBC
9	BL	303	BCL	CAD-CBD-CGD-O1D
9	BL	303	BCL	CAD-CBD-CGD-O2D
9	BM	401	BCL	C2C-C3C-CAC-CBC
9	BM	401	BCL	C4C-C3C-CAC-CBC
9	BA	101	BCL	C4C-C3C-CAC-CBC
9	BB	101	BCL	C2C-C3C-CAC-CBC
9	BB	101	BCL	C4C-C3C-CAC-CBC
9	BE	101	BCL	C4C-C3C-CAC-CBC
9	BE	101	BCL	C11-C12-C13-C14
9	BF	102	BCL	C4C-C3C-CAC-CBC

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Mol	Chain	Res	Type	Atoms
9	BI	102	BCL	C4C-C3C-CAC-CBC
9	BJ	101	BCL	O2A-C1-C2-C3
9	BN	101	BCL	C4C-C3C-CAC-CBC
9	BN	101	BCL	O2A-C1-C2-C3
9	BO	102	BCL	C4C-C3C-CAC-CBC
9	BQ	103	BCL	C4C-C3C-CAC-CBC
9	BQ	104	BCL	C4C-C3C-CAC-CBC
9	BQ	104	BCL	O2A-C1-C2-C3
9	BS	102	BCL	C2C-C3C-CAC-CBC
9	BS	102	BCL	C4C-C3C-CAC-CBC
9	BT	101	BCL	C2C-C3C-CAC-CBC
9	BT	101	BCL	C4C-C3C-CAC-CBC
9	BU	102	BCL	C2C-C3C-CAC-CBC
9	BU	102	BCL	C4C-C3C-CAC-CBC
9	BV	101	BCL	C2C-C3C-CAC-CBC
9	BV	101	BCL	C4C-C3C-CAC-CBC
9	BX	101	BCL	C4C-C3C-CAC-CBC
9	BY	102	BCL	C4C-C3C-CAC-CBC
9	BZ	101	BCL	C2C-C3C-CAC-CBC
9	BZ	101	BCL	C4C-C3C-CAC-CBC
9	BZ	101	BCL	O2A-C1-C2-C3
9	B1	102	BCL	C4C-C3C-CAC-CBC
9	B2	101	BCL	C2C-C3C-CAC-CBC
9	B2	101	BCL	C4C-C3C-CAC-CBC
9	B3	102	BCL	C4C-C3C-CAC-CBC
9	B6	101	BCL	C2C-C3C-CAC-CBC
9	B6	101	BCL	C4C-C3C-CAC-CBC
9	B0	102	BCL	C2C-C3C-CAC-CBC
9	B0	102	BCL	C4C-C3C-CAC-CBC
11	AL	304	UQ8	C30-C29-C31-C32
11	AL	304	UQ8	C28-C29-C31-C32
13	AM	405	MQ8	C12-C11-C3-C2
13	AM	405	MQ8	C12-C11-C3-C4
13	BM	405	MQ8	C12-C11-C3-C4
14	AB	102	CRT	O1-C1-C4-C5
14	AB	102	CRT	C2-C1-C4-C5
14	AB	102	CRT	C3-C1-C4-C5
14	AB	102	CRT	C1-C4-C5-C6
14	AB	102	CRT	C35-C36-C37-C38
14	AG	102	CRT	C1-C4-C5-C6
14	AN	102	CRT	C35-C36-C37-C38
14	AP	102	CRT	C35-C36-C37-C38

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Mol	Chain	Res	Type	Atoms
14	AR	102	CRT	C35-C36-C37-C38
14	AS	104	CRT	C1-C4-C5-C6
14	AT	102	CRT	C35-C36-C37-C38
14	A1	103	CRT	C35-C36-C37-C38
14	A7	102	CRT	C35-C36-C37-C38
14	A0	101	CRT	C1-C4-C5-C6
14	BG	102	CRT	C1-C4-C5-C6
14	BN	102	CRT	C35-C36-C37-C38
14	BO	103	CRT	C35-C36-C37-C38
14	BP	102	CRT	C35-C36-C37-C38
14	BS	103	CRT	C35-C36-C37-C38
14	BU	103	CRT	C35-C36-C37-C38
14	BV	102	CRT	C1-C4-C5-C6
14	BV	102	CRT	C35-C36-C37-C38
14	B2	102	CRT	C1-C4-C5-C6
14	B2	102	CRT	C35-C36-C37-C38
14	B5	103	CRT	C35-C36-C37-C38
14	B7	102	CRT	C1-C4-C5-C6
14	B7	102	CRT	C35-C36-C37-C38
14	B0	101	CRT	O1-C1-C4-C5
14	B0	101	CRT	C2-C1-C4-C5
14	B0	101	CRT	C3-C1-C4-C5
14	B0	101	CRT	C1-C4-C5-C6
14	B0	101	CRT	C35-C36-C37-C38
15	AM	407	PEF	C11-C10-O2-C2
15	AM	407	PEF	O4-C10-O2-C2
15	AM	407	PEF	C1-O3P-P-O1P
15	AM	407	PEF	C1-O3P-P-O2P
15	AM	407	PEF	C1-O3P-P-O4P
15	AM	407	PEF	C4-O4P-P-O1P
15	AM	407	PEF	C4-O4P-P-O2P
15	AM	407	PEF	C4-O4P-P-O3P
15	AM	408	PEF	C5-C4-O4P-P
15	AM	408	PEF	C1-O3P-P-O1P
15	AM	408	PEF	C1-O3P-P-O2P
15	AM	408	PEF	C4-O4P-P-O3P
15	AM	409	PEF	C11-C10-O2-C2
15	AM	409	PEF	O4-C10-O2-C2
15	AH	301	PEF	C11-C10-O2-C2
15	AH	301	PEF	O4-C10-O2-C2
15	AH	301	PEF	C4-O4P-P-O1P
15	AS	101	PEF	C2-C1-O3P-P

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Mol	Chain	Res	Type	Atoms
15	AS	101	PEF	C11-C10-O2-C2
15	AS	101	PEF	C1-O3P-P-O1P
15	AS	101	PEF	C1-O3P-P-O2P
15	AS	101	PEF	C1-O3P-P-O4P
15	BM	407	PEF	C11-C10-O2-C2
15	BM	407	PEF	O4-C10-O2-C2
15	BM	407	PEF	C1-O3P-P-O2P
15	BM	407	PEF	C1-O3P-P-O4P
15	BM	407	PEF	C4-O4P-P-O2P
15	BM	407	PEF	C4-O4P-P-O3P
9	AL	303	BCL	C3-C5-C6-C7
9	BL	303	BCL	C3-C5-C6-C7
9	AZ	101	BCL	O1A-CGA-O2A-C1
9	A3	104	BCL	O1A-CGA-O2A-C1
9	A8	101	BCL	O1A-CGA-O2A-C1
9	A0	102	BCL	O1A-CGA-O2A-C1
15	AS	101	PEF	O4-C10-O2-C2
9	AF	102	BCL	C3-C5-C6-C7
9	AQ	102	BCL	C3-C5-C6-C7
9	A3	103	BCL	C3-C5-C6-C7
9	A6	101	BCL	C3-C5-C6-C7
9	A0	102	BCL	C3-C5-C6-C7
9	BW	102	BCL	C3-C5-C6-C7
10	AL	302	BPH	C3-C5-C6-C7
9	AZ	101	BCL	CBA-CGA-O2A-C1
9	A2	101	BCL	CBA-CGA-O2A-C1
9	A8	101	BCL	CBA-CGA-O2A-C1
9	A0	102	BCL	CBA-CGA-O2A-C1
11	AL	304	UQ8	C20-C19-C21-C22
11	AL	304	UQ8	C18-C19-C21-C22
9	AX	101	BCL	C3-C5-C6-C7
9	AZ	101	BCL	C3-C5-C6-C7
9	BO	102	BCL	C3-C5-C6-C7
9	A3	104	BCL	CBA-CGA-O2A-C1
9	A6	101	BCL	CBA-CGA-O2A-C1
15	AS	101	PEF	C31-C30-O3-C3
15	BQ	101	PEF	C31-C30-O3-C3
15	AM	407	PEF	C31-C30-O3-C3
9	AX	101	BCL	O1A-CGA-O2A-C1
9	A2	101	BCL	O1A-CGA-O2A-C1
9	A6	101	BCL	O1A-CGA-O2A-C1
15	AS	101	PEF	O5-C30-O3-C3

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Mol	Chain	Res	Type	Atoms
15	BQ	101	PEF	O5-C30-O3-C3
15	AH	301	PEF	O5-C30-O3-C3
15	AM	408	PEF	C11-C10-O2-C2
9	A2	101	BCL	C3-C5-C6-C7
9	A8	101	BCL	C3-C5-C6-C7
10	BL	302	BPH	C3-C5-C6-C7
15	BM	407	PEF	O5-C30-O3-C3
15	AH	301	PEF	C31-C30-O3-C3
9	AA	101	BCL	C3-C5-C6-C7
9	BU	102	BCL	C3-C5-C6-C7
9	AX	101	BCL	CBA-CGA-O2A-C1
11	AL	304	UQ8	C24-C26-C27-C28
11	BL	304	UQ8	C24-C26-C27-C28
13	AM	405	MQ8	C18-C20-C21-C22
13	AM	405	MQ8	C23-C25-C26-C27
13	AM	405	MQ8	C28-C30-C31-C32
7	AC	502	HEM	C3D-CAD-CBD-CGD
15	BM	407	PEF	C31-C30-O3-C3
9	AR	101	BCL	C11-C10-C8-C9
9	AV	102	BCL	C14-C13-C15-C16
10	AM	403	BPH	C14-C13-C15-C16
10	BM	403	BPH	C14-C13-C15-C16
15	AS	101	PEF	C36-C37-C38-C39
9	AA	101	BCL	C13-C15-C16-C17
9	AL	301	BCL	C15-C16-C17-C18
9	AU	102	BCL	C8-C10-C11-C12
9	BV	101	BCL	C15-C16-C17-C18
9	B3	102	BCL	C15-C16-C17-C18
15	AM	408	PEF	O4-C10-O2-C2
9	AL	303	BCL	C15-C16-C17-C18
9	B3	102	BCL	C8-C10-C11-C12
15	AM	407	PEF	O5-C30-O3-C3
15	BQ	101	PEF	C39-C40-C41-C42
11	BL	304	UQ8	C34-C36-C37-C38
13	BM	405	MQ8	C13-C15-C16-C17
13	BM	405	MQ8	C18-C20-C21-C22
13	BM	405	MQ8	C38-C40-C41-C42
9	AM	401	BCL	C13-C15-C16-C17
9	AD	102	BCL	C15-C16-C17-C18
15	AS	101	PEF	C30-C31-C32-C33
15	BQ	101	PEF	C30-C31-C32-C33
9	AL	303	BCL	C5-C6-C7-C8

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Mol	Chain	Res	Type	Atoms
9	AA	101	BCL	C8-C10-C11-C12
9	AF	102	BCL	C15-C16-C17-C18
9	AK	102	BCL	C15-C16-C17-C18
9	BQ	104	BCL	C15-C16-C17-C18
9	BU	102	BCL	C8-C10-C11-C12
9	B5	102	BCL	C8-C10-C11-C12
9	AE	101	BCL	C10-C11-C12-C13
9	AQ	102	BCL	C5-C6-C7-C8
9	A3	103	BCL	C5-C6-C7-C8
9	A3	103	BCL	C15-C16-C17-C18
9	BL	301	BCL	C15-C16-C17-C18
9	BA	101	BCL	C5-C6-C7-C8
9	BE	101	BCL	C10-C11-C12-C13
9	BS	102	BCL	C15-C16-C17-C18
9	BW	102	BCL	C15-C16-C17-C18
9	BX	101	BCL	C8-C10-C11-C12
9	BY	102	BCL	C13-C15-C16-C17
9	B1	102	BCL	C5-C6-C7-C8
9	B1	102	BCL	C15-C16-C17-C18
9	B9	102	BCL	C5-C6-C7-C8
10	BL	302	BPH	C15-C16-C17-C18
9	AG	101	BCL	C3-C5-C6-C7
9	AS	103	BCL	C3-C5-C6-C7
9	B5	102	BCL	C3-C5-C6-C7
9	AI	102	BCL	C15-C16-C17-C18
9	AK	102	BCL	C5-C6-C7-C8
9	AN	101	BCL	C10-C11-C12-C13
9	A2	101	BCL	C5-C6-C7-C8
9	BL	303	BCL	C5-C6-C7-C8
9	BF	102	BCL	C5-C6-C7-C8
9	BI	102	BCL	C5-C6-C7-C8
9	BK	102	BCL	C5-C6-C7-C8
9	BY	102	BCL	C15-C16-C17-C18
9	B3	102	BCL	C5-C6-C7-C8
9	B6	101	BCL	C15-C16-C17-C18
9	B7	103	BCL	C13-C15-C16-C17
9	B9	102	BCL	C15-C16-C17-C18
10	AL	302	BPH	C15-C16-C17-C18
9	BD	102	BCL	C15-C16-C17-C18
9	BQ	103	BCL	C5-C6-C7-C8
9	AO	102	BCL	C15-C16-C17-C18
9	AV	102	BCL	C13-C15-C16-C17

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Mol	Chain	Res	Type	Atoms
9	AV	102	BCL	C15-C16-C17-C18
9	A1	102	BCL	C15-C16-C17-C18
9	BF	102	BCL	C15-C16-C17-C18
9	AQ	102	BCL	C10-C11-C12-C13
9	B5	102	BCL	C5-C6-C7-C8
9	AY	102	BCL	CBA-CGA-O2A-C1
9	B5	102	BCL	C15-C16-C17-C18
9	BJ	101	BCL	C3-C5-C6-C7
9	AS	103	BCL	C8-C10-C11-C12
9	A3	103	BCL	C8-C10-C11-C12
9	A7	103	BCL	C5-C6-C7-C8
9	A9	102	BCL	C5-C6-C7-C8
9	BM	401	BCL	C13-C15-C16-C17
9	BM	402	BCL	C13-C15-C16-C17
9	BA	101	BCL	C15-C16-C17-C18
9	BN	101	BCL	C10-C11-C12-C13
9	BO	102	BCL	C15-C16-C17-C18
9	BY	102	BCL	C5-C6-C7-C8
9	B7	103	BCL	C5-C6-C7-C8
9	B7	103	BCL	C8-C10-C11-C12
9	AU	102	BCL	C5-C6-C7-C8
9	A9	102	BCL	C8-C10-C11-C12
9	AM	402	BCL	C5-C6-C7-C8
9	AS	103	BCL	C13-C15-C16-C17
9	AY	102	BCL	C5-C6-C7-C8
9	A1	102	BCL	C8-C10-C11-C12
9	A5	102	BCL	C5-C6-C7-C8
9	A5	102	BCL	C8-C10-C11-C12
9	B6	101	BCL	C5-C6-C7-C8
9	AW	101	BCL	C5-C6-C7-C8
9	A1	102	BCL	C5-C6-C7-C8
9	BU	102	BCL	C5-C6-C7-C8
9	B7	103	BCL	C15-C16-C17-C18
9	BK	102	BCL	C3-C5-C6-C7
9	AR	101	BCL	C16-C17-C18-C19
9	BV	101	BCL	C16-C17-C18-C20
9	BZ	101	BCL	C16-C17-C18-C20
10	BL	302	BPH	C16-C17-C18-C20
14	A7	102	CRT	C21-C22-C23-C24
7	BC	502	HEM	C3D-CAD-CBD-CGD
9	B6	101	BCL	C8-C10-C11-C12
15	BQ	101	PEF	C20-C21-C22-C23

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Mol	Chain	Res	Type	Atoms
9	AG	101	BCL	O2A-C1-C2-C3
9	AR	101	BCL	O2A-C1-C2-C3
9	BT	101	BCL	O2A-C1-C2-C3
10	AL	302	BPH	O2A-C1-C2-C3
10	BL	302	BPH	O2A-C1-C2-C3
9	AN	101	BCL	C16-C17-C18-C19
9	AZ	101	BCL	C16-C17-C18-C19
9	A6	101	BCL	C16-C17-C18-C19
9	A6	101	BCL	C16-C17-C18-C20
9	BN	101	BCL	C16-C17-C18-C20
9	B2	101	BCL	C16-C17-C18-C20
9	B8	101	BCL	C16-C17-C18-C20
10	BL	302	BPH	C16-C17-C18-C19
7	BC	504	HEM	C4D-C3D-CAD-CBD
14	A7	102	CRT	C21-C22-C23-C25
9	AJ	101	BCL	C8-C10-C11-C12
9	AU	102	BCL	C15-C16-C17-C18
9	BI	102	BCL	C15-C16-C17-C18
9	BK	102	BCL	C13-C15-C16-C17
9	A3	104	BCL	C3-C5-C6-C7
9	AN	101	BCL	C16-C17-C18-C20
9	AR	101	BCL	C16-C17-C18-C20
9	AX	101	BCL	C16-C17-C18-C20
9	AZ	101	BCL	C16-C17-C18-C20
9	A3	104	BCL	C16-C17-C18-C19
9	BI	102	BCL	C16-C17-C18-C20
9	BV	101	BCL	C16-C17-C18-C19
10	AL	302	BPH	C16-C17-C18-C20
10	BL	302	BPH	C5-C6-C7-C8
15	BQ	101	PEF	C36-C37-C38-C39
15	AS	101	PEF	C31-C32-C33-C34
15	BQ	101	PEF	C31-C32-C33-C34
9	A6	101	BCL	C5-C6-C7-C8
9	BV	101	BCL	C8-C10-C11-C12
7	BC	504	HEM	C2D-C3D-CAD-CBD
9	AV	102	BCL	C16-C17-C18-C20
9	AX	101	BCL	C16-C17-C18-C19
9	A3	104	BCL	C16-C17-C18-C20
9	A0	102	BCL	C16-C17-C18-C19
9	A0	102	BCL	C16-C17-C18-C20
9	BN	101	BCL	C16-C17-C18-C19
9	B2	101	BCL	C16-C17-C18-C19

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Mol	Chain	Res	Type	Atoms
9	B8	101	BCL	C16-C17-C18-C19
10	AL	302	BPH	C16-C17-C18-C19
9	AD	102	BCL	C5-C6-C7-C8
9	AQ	102	BCL	C11-C12-C13-C15
9	AS	103	BCL	C11-C12-C13-C15
9	AY	102	BCL	C11-C12-C13-C15
9	BD	102	BCL	C8-C10-C11-C12
9	AY	102	BCL	O1A-CGA-O2A-C1
9	AD	102	BCL	C3-C5-C6-C7
9	AK	102	BCL	C3-C5-C6-C7
15	AS	101	PEF	C13-C14-C15-C16
9	AS	103	BCL	C5-C6-C7-C8
9	BX	101	BCL	C15-C16-C17-C18
9	AV	102	BCL	C16-C17-C18-C19
9	A2	101	BCL	C16-C17-C18-C19
9	A2	101	BCL	C16-C17-C18-C20
9	B0	102	BCL	C16-C17-C18-C19
9	B0	102	BCL	C16-C17-C18-C20
15	AM	409	PEF	C17-C18-C19-C20
9	BQ	104	BCL	C10-C11-C12-C13
9	A1	102	BCL	C3-C5-C6-C7
9	B1	102	BCL	C3-C5-C6-C7
15	AS	101	PEF	C34-C35-C36-C37
15	BQ	101	PEF	C16-C17-C18-C19
9	AJ	101	BCL	C16-C17-C18-C20
9	AP	101	BCL	C16-C17-C18-C20
9	B6	101	BCL	C16-C17-C18-C20
9	AU	102	BCL	CBA-CGA-O2A-C1
15	AM	409	PEF	C31-C30-O3-C3
9	BL	301	BCL	C5-C6-C7-C8
9	B8	101	BCL	C10-C11-C12-C13
9	AY	102	BCL	C3-C5-C6-C7
7	BC	503	HEM	C2A-CAA-CBA-CGA
15	AS	101	PEF	C14-C15-C16-C17
9	A7	103	BCL	C15-C16-C17-C18
9	BO	102	BCL	C5-C6-C7-C8
9	AQ	102	BCL	C8-C10-C11-C12
9	B0	102	BCL	C13-C15-C16-C17
9	BK	102	BCL	C15-C16-C17-C18
9	AT	101	BCL	C15-C16-C17-C18
9	BA	101	BCL	C13-C15-C16-C17
15	AS	101	PEF	C33-C34-C35-C36

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Mol	Chain	Res	Type	Atoms
14	AA	102	CRT	C35-C36-C37-C38
14	A5	103	CRT	C35-C36-C37-C38
14	BB	102	CRT	C35-C36-C37-C38
14	BG	102	CRT	C35-C36-C37-C38
14	BS	103	CRT	C1-C4-C5-C6
14	BU	103	CRT	C1-C4-C5-C6
13	BM	405	MQ8	C12-C11-C3-C2
9	BA	101	BCL	C16-C17-C18-C20
9	BG	101	BCL	C16-C17-C18-C20
9	BX	101	BCL	C16-C17-C18-C19
9	BX	101	BCL	C16-C17-C18-C20
9	BZ	101	BCL	C16-C17-C18-C19
9	B6	101	BCL	C16-C17-C18-C19
9	AZ	101	BCL	C4-C3-C5-C6
9	A3	103	BCL	C4-C3-C5-C6
13	AM	405	MQ8	C29-C28-C30-C31
13	AM	405	MQ8	C27-C28-C30-C31
9	B6	101	BCL	C13-C15-C16-C17
15	AS	101	PEF	C41-C42-C43-C44
9	AM	402	BCL	C10-C11-C12-C13
9	AR	101	BCL	C10-C11-C12-C13
9	AW	101	BCL	C15-C16-C17-C18
9	BZ	101	BCL	C15-C16-C17-C18
9	B2	101	BCL	C15-C16-C17-C18
9	B7	103	BCL	C10-C11-C12-C13
15	AM	407	PEF	O3P-C1-C2-O2
15	AM	409	PEF	O3P-C1-C2-O2
15	AH	301	PEF	O3P-C1-C2-O2
9	AE	101	BCL	C5-C6-C7-C8
9	AG	101	BCL	C13-C15-C16-C17
9	AJ	101	BCL	C15-C16-C17-C18
9	AW	101	BCL	C8-C10-C11-C12
9	A5	102	BCL	C15-C16-C17-C18
9	A9	102	BCL	C15-C16-C17-C18
9	BW	102	BCL	C5-C6-C7-C8
7	AC	503	HEM	C4C-C3C-CAC-CBC
9	AP	101	BCL	C16-C17-C18-C19
9	AA	101	BCL	C5-C6-C7-C8
9	AY	102	BCL	C15-C16-C17-C18
9	BV	101	BCL	C5-C6-C7-C8
10	AM	403	BPH	C5-C6-C7-C8
15	BQ	101	PEF	C17-C18-C19-C20

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Mol	Chain	Res	Type	Atoms
9	AJ	101	BCL	C16-C17-C18-C19
9	BG	101	BCL	C16-C17-C18-C19
9	BT	101	BCL	C16-C17-C18-C19
13	BM	405	MQ8	C45-C43-C44-C46
7	BC	504	HEM	C3D-CAD-CBD-CGD
9	BM	402	BCL	C3-C5-C6-C7
9	AZ	101	BCL	C2-C3-C5-C6
9	AG	101	BCL	C15-C16-C17-C18
9	BQ	103	BCL	C15-C16-C17-C18
15	AM	409	PEF	O5-C30-O3-C3
9	AX	101	BCL	C1A-C2A-CAA-CBA
9	A1	102	BCL	C1A-C2A-CAA-CBA
9	AU	102	BCL	O1A-CGA-O2A-C1
9	AY	102	BCL	C8-C10-C11-C12
15	AM	409	PEF	O3P-C1-C2-C3
9	AM	402	BCL	C11-C12-C13-C15
9	AA	101	BCL	C11-C10-C8-C7
9	AA	101	BCL	C12-C13-C15-C16
9	AD	102	BCL	C11-C12-C13-C15
9	AE	101	BCL	C11-C12-C13-C15
9	AF	102	BCL	C12-C13-C15-C16
9	AI	102	BCL	C6-C7-C8-C10
9	AK	102	BCL	C11-C10-C8-C7
9	AK	102	BCL	C11-C12-C13-C15
9	AN	101	BCL	C12-C13-C15-C16
9	AO	102	BCL	C11-C12-C13-C15
9	AO	102	BCL	C12-C13-C15-C16
9	AQ	102	BCL	C12-C13-C15-C16
9	AU	102	BCL	C11-C12-C13-C15
9	AW	101	BCL	C11-C12-C13-C15
9	AW	101	BCL	C12-C13-C15-C16
9	AY	102	BCL	C6-C7-C8-C10
9	A1	102	BCL	C6-C7-C8-C10
9	A3	103	BCL	C6-C7-C8-C10
9	A3	103	BCL	C12-C13-C15-C16
9	A6	101	BCL	C11-C12-C13-C15
9	A7	103	BCL	C6-C7-C8-C10
9	A7	103	BCL	C11-C12-C13-C15
9	A8	101	BCL	C6-C7-C8-C10
9	A9	102	BCL	C6-C7-C8-C10
9	A9	102	BCL	C11-C12-C13-C15
9	A0	102	BCL	C6-C7-C8-C10

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Mol	Chain	Res	Type	Atoms
9	BM	401	BCL	C12-C13-C15-C16
9	BA	101	BCL	C12-C13-C15-C16
9	BB	101	BCL	C6-C7-C8-C10
9	BD	102	BCL	C11-C12-C13-C15
9	BG	101	BCL	C11-C12-C13-C15
9	BI	102	BCL	C11-C10-C8-C7
9	BK	102	BCL	C11-C10-C8-C7
9	BK	102	BCL	C12-C13-C15-C16
9	BN	101	BCL	C11-C12-C13-C15
9	BO	102	BCL	C11-C12-C13-C15
9	BQ	103	BCL	C11-C12-C13-C15
9	BQ	103	BCL	C12-C13-C15-C16
9	BX	101	BCL	C12-C13-C15-C16
9	B1	102	BCL	C11-C10-C8-C7
9	B2	101	BCL	C12-C13-C15-C16
9	B3	102	BCL	C12-C13-C15-C16
9	B4	101	BCL	C11-C12-C13-C15
9	B5	102	BCL	C11-C10-C8-C7
9	B6	101	BCL	C12-C13-C15-C16
9	B9	102	BCL	C11-C12-C13-C15
9	AP	101	BCL	C8-C10-C11-C12
9	AV	102	BCL	C2C-C3C-CAC-CBC
9	AY	102	BCL	C2C-C3C-CAC-CBC
9	BE	101	BCL	C2C-C3C-CAC-CBC
9	BF	102	BCL	C2C-C3C-CAC-CBC
9	BN	101	BCL	C2C-C3C-CAC-CBC
9	BQ	104	BCL	C2C-C3C-CAC-CBC
9	BX	101	BCL	C2C-C3C-CAC-CBC
9	B6	101	BCL	C4-C3-C5-C6
9	B6	101	BCL	C2-C3-C5-C6
15	BQ	101	PEF	C18-C19-C20-C21
9	AA	101	BCL	C11-C10-C8-C9
9	AA	101	BCL	C11-C12-C13-C14
9	AD	102	BCL	C6-C7-C8-C9
9	AI	102	BCL	C6-C7-C8-C9
9	AK	102	BCL	C14-C13-C15-C16
9	AN	101	BCL	C14-C13-C15-C16
9	AO	102	BCL	C14-C13-C15-C16
9	AQ	102	BCL	C11-C10-C8-C9
9	AW	101	BCL	C14-C13-C15-C16
9	A1	102	BCL	C6-C7-C8-C9
9	A1	102	BCL	C11-C12-C13-C14

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Mol	Chain	Res	Type	Atoms
9	A2	101	BCL	C11-C12-C13-C14
9	A3	103	BCL	C6-C7-C8-C9
9	A3	104	BCL	C14-C13-C15-C16
9	A6	101	BCL	C6-C7-C8-C9
9	A6	101	BCL	C11-C12-C13-C14
9	A8	101	BCL	C6-C7-C8-C9
9	BB	101	BCL	C6-C7-C8-C9
9	BF	102	BCL	C11-C10-C8-C9
9	BF	102	BCL	C14-C13-C15-C16
9	BI	102	BCL	C11-C10-C8-C9
9	BI	102	BCL	C14-C13-C15-C16
9	BJ	101	BCL	C11-C12-C13-C14
9	BN	101	BCL	C11-C10-C8-C9
9	BN	101	BCL	C11-C12-C13-C14
9	BO	102	BCL	C11-C10-C8-C9
9	BP	101	BCL	C11-C12-C13-C14
9	BQ	103	BCL	C11-C10-C8-C9
9	BU	102	BCL	C11-C10-C8-C9
9	BY	102	BCL	C14-C13-C15-C16
9	B1	102	BCL	C11-C12-C13-C14
9	B5	102	BCL	C11-C10-C8-C9
9	B5	102	BCL	C14-C13-C15-C16
9	B6	101	BCL	C14-C13-C15-C16
9	B7	103	BCL	C6-C7-C8-C9
9	B7	103	BCL	C11-C12-C13-C14
9	B0	102	BCL	C11-C12-C13-C14
10	AL	302	BPH	C14-C13-C15-C16
15	AS	101	PEF	C37-C38-C39-C40
15	BQ	101	PEF	C37-C38-C39-C40
9	AY	102	BCL	C13-C15-C16-C17
9	BO	102	BCL	C10-C11-C12-C13
15	AM	409	PEF	C1-C2-C3-O3
15	BM	407	PEF	C1-C2-C3-O3
9	A6	101	BCL	C8-C10-C11-C12
9	BU	102	BCL	C10-C11-C12-C13
9	AE	101	BCL	C16-C17-C18-C20
9	AK	102	BCL	C16-C17-C18-C19
9	A5	102	BCL	C16-C17-C18-C19
10	AM	403	BPH	C16-C17-C18-C20
9	BE	101	BCL	C3-C5-C6-C7
15	AS	101	PEF	C39-C40-C41-C42
9	A6	101	BCL	C4-C3-C5-C6

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Mol	Chain	Res	Type	Atoms
9	A3	103	BCL	C2-C3-C5-C6
9	A6	101	BCL	C2-C3-C5-C6
9	AL	301	BCL	C5-C6-C7-C8
10	AL	302	BPH	C5-C6-C7-C8
10	BM	403	BPH	C13-C15-C16-C17
9	AE	101	BCL	C16-C17-C18-C19
9	A5	102	BCL	C16-C17-C18-C20
9	BO	102	BCL	C13-C15-C16-C17
9	BI	102	BCL	C8-C10-C11-C12
10	AM	403	BPH	C13-C15-C16-C17
9	AM	402	BCL	O2A-C1-C2-C3
9	AT	101	BCL	O2A-C1-C2-C3
9	BE	101	BCL	O2A-C1-C2-C3
9	BG	101	BCL	O2A-C1-C2-C3
9	BV	101	BCL	O2A-C1-C2-C3
9	AL	301	BCL	C13-C15-C16-C17
9	A9	102	BCL	C13-C15-C16-C17
15	BQ	101	PEF	C12-C13-C14-C15
10	AM	403	BPH	C16-C17-C18-C19
15	AM	409	PEF	C21-C22-C23-C24
15	AM	409	PEF	C39-C40-C41-C42
9	A5	102	BCL	C3-C5-C6-C7
9	BY	102	BCL	C3-C5-C6-C7
9	AT	101	BCL	C10-C11-C12-C13
9	A1	102	BCL	C13-C15-C16-C17
9	BS	102	BCL	C8-C10-C11-C12
15	AM	409	PEF	C18-C19-C20-C21
9	BL	301	BCL	C13-C15-C16-C17
9	AS	103	BCL	C15-C16-C17-C18
9	AD	102	BCL	C4-C3-C5-C6
9	AX	101	BCL	C4-C3-C5-C6
9	A3	104	BCL	C4-C3-C5-C6
9	BO	102	BCL	C4-C3-C5-C6
13	BM	405	MQ8	C34-C33-C35-C36
9	AX	101	BCL	C2-C3-C5-C6
11	AL	304	UQ8	C38-C39-C41-C42
15	AM	409	PEF	C30-C31-C32-C33
9	AM	402	BCL	C11-C12-C13-C14
9	AA	101	BCL	C14-C13-C15-C16
9	AD	102	BCL	C11-C12-C13-C14
9	AE	101	BCL	C11-C12-C13-C14
9	AF	102	BCL	C14-C13-C15-C16

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Mol	Chain	Res	Type	Atoms
9	AK	102	BCL	C11-C10-C8-C9
9	AK	102	BCL	C11-C12-C13-C14
9	AO	102	BCL	C11-C12-C13-C14
9	AQ	102	BCL	C11-C12-C13-C14
9	AQ	102	BCL	C14-C13-C15-C16
9	AS	103	BCL	C11-C12-C13-C14
9	AU	102	BCL	C6-C7-C8-C9
9	AU	102	BCL	C11-C12-C13-C14
9	AW	101	BCL	C6-C7-C8-C9
9	AW	101	BCL	C11-C12-C13-C14
9	AX	101	BCL	C11-C12-C13-C14
9	AY	102	BCL	C6-C7-C8-C9
9	AY	102	BCL	C11-C12-C13-C14
9	A3	103	BCL	C14-C13-C15-C16
9	A7	103	BCL	C6-C7-C8-C9
9	A7	103	BCL	C11-C12-C13-C14
9	A7	103	BCL	C14-C13-C15-C16
9	A9	102	BCL	C6-C7-C8-C9
9	A9	102	BCL	C11-C12-C13-C14
9	A0	102	BCL	C6-C7-C8-C9
9	BM	401	BCL	C14-C13-C15-C16
9	BA	101	BCL	C14-C13-C15-C16
9	BD	102	BCL	C11-C12-C13-C14
9	BD	102	BCL	C14-C13-C15-C16
9	BG	101	BCL	C11-C12-C13-C14
9	BK	102	BCL	C11-C10-C8-C9
9	BK	102	BCL	C14-C13-C15-C16
9	BO	102	BCL	C11-C12-C13-C14
9	BQ	103	BCL	C11-C12-C13-C14
9	BQ	103	BCL	C14-C13-C15-C16
9	BS	102	BCL	C14-C13-C15-C16
9	BV	101	BCL	C14-C13-C15-C16
9	BX	101	BCL	C14-C13-C15-C16
9	B1	102	BCL	C11-C10-C8-C9
9	B2	101	BCL	C14-C13-C15-C16
9	B3	102	BCL	C11-C10-C8-C9
9	B3	102	BCL	C11-C12-C13-C14
9	B3	102	BCL	C14-C13-C15-C16
9	B4	101	BCL	C11-C12-C13-C14
9	B6	101	BCL	C11-C12-C13-C14
9	B9	102	BCL	C11-C12-C13-C14
10	AL	302	BPH	C11-C10-C8-C9

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Mol	Chain	Res	Type	Atoms
9	B0	102	BCL	C8-C10-C11-C12
9	AM	401	BCL	C5-C6-C7-C8
15	AM	407	PEF	O3P-C1-C2-C3
15	AH	301	PEF	O3P-C1-C2-C3
9	AA	101	BCL	C11-C12-C13-C15
9	AI	102	BCL	C11-C12-C13-C15
9	AJ	101	BCL	C11-C12-C13-C15
9	AJ	101	BCL	C12-C13-C15-C16
9	AK	102	BCL	C12-C13-C15-C16
9	AQ	102	BCL	C11-C10-C8-C7
9	AS	103	BCL	C12-C13-C15-C16
9	AU	102	BCL	C6-C7-C8-C10
9	AV	102	BCL	C11-C12-C13-C15
9	AW	101	BCL	C6-C7-C8-C10
9	AX	101	BCL	C11-C12-C13-C15
9	A1	102	BCL	C11-C12-C13-C15
9	A2	101	BCL	C11-C12-C13-C15
9	A3	104	BCL	C6-C7-C8-C10
9	A3	104	BCL	C11-C12-C13-C15
9	A3	104	BCL	C12-C13-C15-C16
9	A7	103	BCL	C12-C13-C15-C16
9	BL	301	BCL	C12-C13-C15-C16
9	BL	303	BCL	C11-C10-C8-C7
9	BL	303	BCL	C11-C12-C13-C15
9	BD	102	BCL	C12-C13-C15-C16
9	BE	101	BCL	C11-C10-C8-C7
9	BE	101	BCL	C11-C12-C13-C15
9	BF	102	BCL	C11-C10-C8-C7
9	BF	102	BCL	C12-C13-C15-C16
9	BI	102	BCL	C11-C12-C13-C15
9	BI	102	BCL	C12-C13-C15-C16
9	BJ	101	BCL	C11-C12-C13-C15
9	BO	102	BCL	C11-C10-C8-C7
9	BO	102	BCL	C12-C13-C15-C16
9	BP	101	BCL	C11-C12-C13-C15
9	BP	101	BCL	C12-C13-C15-C16
9	BS	102	BCL	C12-C13-C15-C16
9	BU	102	BCL	C11-C10-C8-C7
9	BU	102	BCL	C12-C13-C15-C16
9	BV	101	BCL	C12-C13-C15-C16
9	BW	102	BCL	C12-C13-C15-C16
9	BY	102	BCL	C12-C13-C15-C16

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Mol	Chain	Res	Type	Atoms
9	BZ	101	BCL	C11-C12-C13-C15
9	B1	102	BCL	C12-C13-C15-C16
9	B3	102	BCL	C11-C10-C8-C7
9	B3	102	BCL	C11-C12-C13-C15
9	B5	102	BCL	C12-C13-C15-C16
9	B6	101	BCL	C11-C12-C13-C15
9	B7	103	BCL	C6-C7-C8-C10
9	B0	102	BCL	C6-C7-C8-C10
10	AL	302	BPH	C11-C10-C8-C7
10	AL	302	BPH	C12-C13-C15-C16
9	A7	103	BCL	C8-C10-C11-C12
9	BI	102	BCL	C16-C17-C18-C19
9	AM	402	BCL	C4-C3-C5-C6
9	AG	101	BCL	C4-C3-C5-C6
9	AU	102	BCL	C3A-C2A-CAA-CBA
9	AX	101	BCL	C3A-C2A-CAA-CBA
9	A2	101	BCL	C3A-C2A-CAA-CBA
9	A5	102	BCL	C4-C3-C5-C6
9	BQ	103	BCL	C4-C3-C5-C6
9	B9	102	BCL	C4-C3-C5-C6
10	AL	302	BPH	C4-C3-C5-C6
11	AL	304	UQ8	C25-C24-C26-C27
9	AN	101	BCL	C8-C10-C11-C12
9	AV	102	BCL	C8-C10-C11-C12
15	AM	409	PEF	C12-C13-C14-C15
9	AF	102	BCL	C8-C10-C11-C12
9	BS	102	BCL	C13-C15-C16-C17
9	AG	101	BCL	C16-C17-C18-C19
15	AS	101	PEF	C1-C2-C3-O3
15	BQ	101	PEF	C1-C2-C3-O3
9	AG	101	BCL	C16-C17-C18-C20
9	BA	101	BCL	C16-C17-C18-C19
11	AL	304	UQ8	C40-C39-C41-C42
9	AS	103	BCL	CAA-CBA-CGA-O2A
9	AM	402	BCL	C2-C3-C5-C6
9	A5	102	BCL	C2-C3-C5-C6
9	BO	102	BCL	C2-C3-C5-C6
9	B9	102	BCL	C2-C3-C5-C6
11	AL	304	UQ8	C23-C24-C26-C27
9	B1	102	BCL	C8-C10-C11-C12
15	BQ	101	PEF	C22-C23-C24-C25
9	AD	102	BCL	C8-C10-C11-C12

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Mol	Chain	Res	Type	Atoms
9	AO	102	BCL	C13-C15-C16-C17
9	B1	102	BCL	C13-C15-C16-C17
9	BT	101	BCL	C16-C17-C18-C20
15	AS	101	PEF	O2-C2-C3-O3
15	BQ	101	PEF	O2-C2-C3-O3
9	B3	102	BCL	C3-C5-C6-C7
9	B9	102	BCL	C8-C10-C11-C12
9	AG	101	BCL	C2-C3-C5-C6
9	BQ	103	BCL	C2-C3-C5-C6
10	AL	302	BPH	C2-C3-C5-C6
13	BM	405	MQ8	C32-C33-C35-C36
9	AD	102	BCL	C11-C10-C8-C9
9	AJ	101	BCL	C11-C12-C13-C14
9	AJ	101	BCL	C14-C13-C15-C16
9	AS	103	BCL	C14-C13-C15-C16
9	AZ	101	BCL	C6-C7-C8-C9
9	A3	104	BCL	C6-C7-C8-C9
9	A3	104	BCL	C11-C12-C13-C14
9	BL	303	BCL	C11-C10-C8-C9
9	BU	102	BCL	C14-C13-C15-C16
9	BA	101	BCL	C3-C5-C6-C7
9	AL	303	BCL	C8-C10-C11-C12
11	BL	304	UQ8	C19-C21-C22-C23
13	AM	405	MQ8	C33-C35-C36-C37
9	AA	101	BCL	C4C-C3C-CAC-CBC
9	A8	101	BCL	C4C-C3C-CAC-CBC
9	A9	102	BCL	C4C-C3C-CAC-CBC
9	BK	102	BCL	C4C-C3C-CAC-CBC
9	BW	102	BCL	C4C-C3C-CAC-CBC
9	B9	102	BCL	C4C-C3C-CAC-CBC
9	AN	101	BCL	C4-C3-C5-C6
9	BY	102	BCL	CAA-CBA-CGA-O2A
9	AD	102	BCL	C2-C3-C5-C6
9	A3	104	BCL	C2-C3-C5-C6
9	BG	101	BCL	C13-C15-C16-C17
9	A7	103	BCL	C13-C15-C16-C17
13	BM	405	MQ8	C35-C36-C37-C38
9	BJ	101	BCL	C5-C6-C7-C8
9	BW	102	BCL	C8-C10-C11-C12
9	BX	101	BCL	C13-C15-C16-C17
9	AL	301	BCL	C6-C7-C8-C10
9	AL	301	BCL	C12-C13-C15-C16

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Mol	Chain	Res	Type	Atoms
9	AL	303	BCL	C11-C10-C8-C7
9	AA	101	BCL	C6-C7-C8-C10
9	AD	102	BCL	C12-C13-C15-C16
9	AO	102	BCL	C11-C10-C8-C7
9	AR	101	BCL	C6-C7-C8-C10
9	AT	101	BCL	C11-C12-C13-C15
9	AU	102	BCL	C12-C13-C15-C16
9	AY	102	BCL	C12-C13-C15-C16
9	AZ	101	BCL	C11-C12-C13-C15
9	A1	102	BCL	C12-C13-C15-C16
9	A2	101	BCL	C12-C13-C15-C16
9	A5	102	BCL	C6-C7-C8-C10
9	A8	101	BCL	C12-C13-C15-C16
9	A9	102	BCL	C12-C13-C15-C16
9	A0	102	BCL	C11-C12-C13-C15
9	BG	101	BCL	C6-C7-C8-C10
9	BQ	103	BCL	C11-C10-C8-C7
9	BV	101	BCL	C6-C7-C8-C10
9	BV	101	BCL	C11-C12-C13-C15
9	BY	102	BCL	C11-C10-C8-C7
9	BZ	101	BCL	C6-C7-C8-C10
9	BZ	101	BCL	C12-C13-C15-C16
9	B8	101	BCL	C11-C12-C13-C15
14	AP	102	CRT	C1-C4-C5-C6
14	AS	104	CRT	C35-C36-C37-C38
14	AX	102	CRT	C35-C36-C37-C38
14	A1	103	CRT	C1-C4-C5-C6
14	A2	102	CRT	C1-C4-C5-C6
14	A7	102	CRT	C1-C4-C5-C6
14	A0	101	CRT	C35-C36-C37-C38
14	BA	102	CRT	C35-C36-C37-C38
14	BB	102	CRT	C1-C4-C5-C6
14	BF	103	CRT	C1-C4-C5-C6
14	BP	102	CRT	C1-C4-C5-C6
14	B1	103	CRT	C1-C4-C5-C6
14	B1	103	CRT	C35-C36-C37-C38
9	BN	101	BCL	C8-C10-C11-C12
7	AC	501	HEM	C3D-CAD-CBD-CGD
7	BC	503	HEM	C3D-CAD-CBD-CGD
9	AO	102	BCL	C5-C6-C7-C8
9	B4	101	BCL	C16-C17-C18-C20
7	AC	504	HEM	C2B-C3B-CAB-CBB

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Mol	Chain	Res	Type	Atoms
9	AW	101	BCL	C3-C5-C6-C7
9	BQ	103	BCL	C3-C5-C6-C7
9	AU	102	BCL	O2A-C1-C2-C3
9	AK	102	BCL	C16-C17-C18-C20
9	B4	101	BCL	C16-C17-C18-C19
15	AS	101	PEF	O3P-C1-C2-O2
9	AF	102	BCL	C5-C6-C7-C8
9	AS	103	BCL	C10-C11-C12-C13
7	AC	501	HEM	C4C-C3C-CAC-CBC
7	BC	501	HEM	C4C-C3C-CAC-CBC
9	AJ	101	BCL	C4-C3-C5-C6
9	AN	101	BCL	C2-C3-C5-C6
13	BM	405	MQ8	C42-C43-C44-C46
15	AM	409	PEF	C5-C4-O4P-P
15	BQ	101	PEF	C5-C4-O4P-P
9	BE	101	BCL	C15-C16-C17-C18
9	BY	102	BCL	C10-C11-C12-C13
9	AY	102	BCL	C14-C13-C15-C16
9	A2	101	BCL	C14-C13-C15-C16
9	BE	101	BCL	C14-C13-C15-C16
9	BW	102	BCL	C14-C13-C15-C16
9	BY	102	BCL	C11-C10-C8-C9
9	BZ	101	BCL	C14-C13-C15-C16
15	AM	409	PEF	C16-C17-C18-C19
9	AX	101	BCL	C5-C6-C7-C8
9	BS	102	BCL	C10-C11-C12-C13
9	BM	402	BCL	C8-C10-C11-C12
9	BD	102	BCL	C13-C15-C16-C17
9	BQ	103	BCL	C8-C10-C11-C12
9	BB	101	BCL	C16-C17-C18-C20
9	B5	102	BCL	C16-C17-C18-C20
10	BM	403	BPH	C16-C17-C18-C20
9	AW	101	BCL	C13-C15-C16-C17
9	A9	102	BCL	C3-C5-C6-C7
9	BS	102	BCL	C3-C5-C6-C7
9	AK	102	BCL	C13-C15-C16-C17
9	AU	102	BCL	C13-C15-C16-C17
9	BM	401	BCL	C16-C17-C18-C19
9	AB	101	BCL	C3-C5-C6-C7
9	AU	102	BCL	C3-C5-C6-C7
9	AL	303	BCL	C11-C12-C13-C15
9	AM	401	BCL	C6-C7-C8-C10

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Mol	Chain	Res	Type	Atoms
9	AD	102	BCL	C6-C7-C8-C10
9	AN	101	BCL	C6-C7-C8-C10
9	AO	102	BCL	C6-C7-C8-C10
9	AT	101	BCL	C12-C13-C15-C16
9	AV	102	BCL	C12-C13-C15-C16
9	AZ	101	BCL	C12-C13-C15-C16
9	A6	101	BCL	C6-C7-C8-C10
9	A0	102	BCL	C12-C13-C15-C16
9	BL	301	BCL	C6-C7-C8-C10
9	BM	402	BCL	C6-C7-C8-C10
9	BG	101	BCL	C12-C13-C15-C16
9	B5	102	BCL	C11-C12-C13-C15
9	B7	103	BCL	C11-C12-C13-C15
9	B0	102	BCL	C11-C12-C13-C15
10	BL	302	BPH	C11-C10-C8-C7
9	AT	101	BCL	C2C-C3C-CAC-CBC
9	AW	101	BCL	C2C-C3C-CAC-CBC
9	A9	102	BCL	CBA-CGA-O2A-C1
9	BB	101	BCL	C16-C17-C18-C19
9	AL	301	BCL	C6-C7-C8-C9
9	AL	303	BCL	C11-C10-C8-C9
9	AD	102	BCL	C14-C13-C15-C16
9	AT	101	BCL	C11-C12-C13-C14
9	AU	102	BCL	C14-C13-C15-C16
9	A1	102	BCL	C14-C13-C15-C16
9	A3	103	BCL	C11-C12-C13-C14
9	A5	102	BCL	C6-C7-C8-C9
9	A9	102	BCL	C14-C13-C15-C16
9	BW	102	BCL	C11-C12-C13-C14
9	BZ	101	BCL	C6-C7-C8-C9
10	BM	403	BPH	C6-C7-C8-C9
9	AD	102	BCL	C13-C15-C16-C17
9	A9	102	BCL	O1A-CGA-O2A-C1
15	BM	407	PEF	O2-C2-C3-O3
9	B5	102	BCL	C16-C17-C18-C19
9	BF	102	BCL	C13-C15-C16-C17
9	AJ	101	BCL	C2-C3-C5-C6
15	BQ	101	PEF	C42-C43-C44-C45
9	AM	401	BCL	CAD-CBD-CGD-O2D
9	A3	103	BCL	CAD-CBD-CGD-O2D
9	BM	401	BCL	CAD-CBD-CGD-O2D
9	BM	402	BCL	CAD-CBD-CGD-O2D

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Mol	Chain	Res	Type	Atoms
9	BW	102	BCL	C13-C15-C16-C17
9	B5	102	BCL	C10-C11-C12-C13
9	B7	103	BCL	CAA-CBA-CGA-O2A
9	AM	401	BCL	CAD-CBD-CGD-O1D
9	AM	402	BCL	CAD-CBD-CGD-O1D
9	A3	103	BCL	CAD-CBD-CGD-O1D
9	BM	401	BCL	CAD-CBD-CGD-O1D
9	BM	402	BCL	CAD-CBD-CGD-O1D
9	BD	102	BCL	CHA-CBD-CGD-O1D
15	AM	408	PEF	C1-O3P-P-O4P
9	BM	401	BCL	C16-C17-C18-C20
9	AQ	102	BCL	C13-C15-C16-C17
9	BG	101	BCL	C8-C10-C11-C12
10	BL	302	BPH	C4-C3-C5-C6
13	BM	405	MQ8	C14-C13-C15-C16
9	B8	101	BCL	C2A-CAA-CBA-CGA
10	BM	403	BPH	C16-C17-C18-C19
15	BM	407	PEF	C3-C2-O2-C10
9	BQ	104	BCL	C13-C15-C16-C17
15	AS	101	PEF	O3P-C1-C2-C3
9	AL	301	BCL	C14-C13-C15-C16
9	AM	401	BCL	C11-C12-C13-C14
9	AA	101	BCL	C6-C7-C8-C9
9	AI	102	BCL	C11-C10-C8-C9
9	AO	102	BCL	C11-C10-C8-C9
9	AP	101	BCL	C11-C12-C13-C14
9	AQ	102	BCL	C6-C7-C8-C9
9	AR	101	BCL	C14-C13-C15-C16
9	AS	103	BCL	C11-C10-C8-C9
9	AZ	101	BCL	C11-C12-C13-C14
9	A0	102	BCL	C11-C12-C13-C14
9	BE	101	BCL	C11-C10-C8-C9
9	BK	102	BCL	C11-C12-C13-C14
9	BO	102	BCL	C14-C13-C15-C16
9	BP	101	BCL	C14-C13-C15-C16
9	BS	102	BCL	C11-C12-C13-C14
9	BV	101	BCL	C6-C7-C8-C9
9	BV	101	BCL	C11-C12-C13-C14
9	BZ	101	BCL	C11-C12-C13-C14
9	B8	101	BCL	C11-C12-C13-C14
9	B9	102	BCL	C11-C10-C8-C9
10	AM	403	BPH	C11-C10-C8-C9

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Mol	Chain	Res	Type	Atoms
10	BL	302	BPH	C11-C10-C8-C9
10	BL	302	BPH	C14-C13-C15-C16
9	AR	101	BCL	C11-C10-C8-C7
9	AX	101	BCL	C12-C13-C15-C16
9	AZ	101	BCL	C6-C7-C8-C10
9	B1	102	BCL	C11-C12-C13-C15
10	BL	302	BPH	C12-C13-C15-C16
15	AM	408	PEF	O3P-C1-C2-O2
9	AR	101	BCL	C4-C3-C5-C6
9	BB	101	BCL	C4-C3-C5-C6
15	AM	409	PEF	O2-C2-C3-O3
9	A5	102	BCL	C13-C15-C16-C17
11	BL	304	UQ8	C14-C16-C17-C18
14	AJ	102	CRT	C1-C4-C5-C6
14	AT	102	CRT	C1-C4-C5-C6
9	AG	101	BCL	C8-C10-C11-C12
9	BQ	104	BCL	C16-C17-C18-C20
9	BQ	104	BCL	C4-C3-C5-C6
10	BL	302	BPH	C2-C3-C5-C6
9	AF	102	BCL	C11-C10-C8-C9
9	AI	102	BCL	C11-C12-C13-C14
9	AR	101	BCL	C6-C7-C8-C9
9	AX	101	BCL	C6-C7-C8-C9
9	A1	102	BCL	C11-C10-C8-C9
9	BL	301	BCL	C6-C7-C8-C9
9	BL	301	BCL	C14-C13-C15-C16
9	BL	303	BCL	C11-C12-C13-C14
9	BA	101	BCL	C11-C10-C8-C9
9	BI	102	BCL	C11-C12-C13-C14
9	BW	102	BCL	C11-C10-C8-C9
9	B1	102	BCL	C14-C13-C15-C16
9	AM	401	BCL	C4-C3-C5-C6
9	AI	102	BCL	C4-C3-C5-C6
9	AR	101	BCL	C2-C3-C5-C6
9	BB	101	BCL	C2-C3-C5-C6
13	BM	405	MQ8	C43-C44-C46-C47
9	AI	102	BCL	C13-C15-C16-C17
7	AC	504	HEM	C4B-C3B-CAB-CBB
15	AM	409	PEF	C14-C15-C16-C17
9	AF	102	BCL	C11-C10-C8-C7
9	A1	102	BCL	C11-C10-C8-C7
9	A3	103	BCL	C11-C12-C13-C15

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Mol	Chain	Res	Type	Atoms
9	A5	102	BCL	C11-C10-C8-C7
9	BA	101	BCL	C11-C10-C8-C7
9	BK	102	BCL	C11-C12-C13-C15
9	BN	101	BCL	C6-C7-C8-C10
9	BS	102	BCL	C11-C12-C13-C15
9	B9	102	BCL	C11-C10-C8-C7
10	AM	403	BPH	C11-C10-C8-C7
10	AM	403	BPH	C12-C13-C15-C16
9	BI	102	BCL	C13-C15-C16-C17
7	BC	503	HEM	CAA-CBA-CGA-O2A
9	BQ	104	BCL	C16-C17-C18-C19
9	AB	101	BCL	C4-C3-C5-C6
9	A1	102	BCL	C3A-C2A-CAA-CBA
9	BP	101	BCL	C4-C3-C5-C6
9	AA	101	BCL	C15-C16-C17-C18
13	BM	405	MQ8	C12-C13-C15-C16
9	AO	102	BCL	C8-C10-C11-C12
9	B8	101	BCL	C13-C15-C16-C17
9	AA	101	BCL	C2-C1-O2A-CGA
9	BU	102	BCL	C2-C1-O2A-CGA
9	B9	102	BCL	C13-C15-C16-C17
9	AK	102	BCL	C4-C3-C5-C6
9	B3	102	BCL	C4-C3-C5-C6
9	AZ	101	BCL	C5-C6-C7-C8
9	B2	101	BCL	C2A-CAA-CBA-CGA
7	AC	503	HEM	CAA-CBA-CGA-O2A
7	BC	501	HEM	CAA-CBA-CGA-O1A
7	BC	503	HEM	CAA-CBA-CGA-O1A
9	AV	102	BCL	C11-C12-C13-C14
9	AW	101	BCL	C11-C10-C8-C9
9	A5	102	BCL	C11-C10-C8-C9
9	BD	102	BCL	C11-C10-C8-C9
9	BF	102	BCL	C11-C12-C13-C14
9	BS	102	BCL	C11-C10-C8-C9
9	BU	102	BCL	C6-C7-C8-C9
9	B6	101	BCL	C6-C7-C8-C9
9	B8	101	BCL	C11-C10-C8-C9
9	B0	102	BCL	C6-C7-C8-C9
10	AM	403	BPH	C6-C7-C8-C9
10	BM	403	BPH	C11-C10-C8-C9
9	BL	303	BCL	C15-C16-C17-C18
9	BF	102	BCL	C8-C10-C11-C12

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Mol	Chain	Res	Type	Atoms
9	BM	402	BCL	O2A-C1-C2-C3
15	AM	407	PEF	C3-C2-O2-C10
15	AH	301	PEF	C3-C2-O2-C10
15	AS	101	PEF	C1-C2-O2-C10
15	AM	409	PEF	C11-C12-C13-C14
15	BQ	101	PEF	C11-C12-C13-C14
9	AS	103	BCL	C4-C3-C5-C6
9	BU	102	BCL	CAA-CBA-CGA-O2A
9	BQ	104	BCL	C2-C3-C5-C6
15	AS	101	PEF	C12-C13-C14-C15
9	BQ	103	BCL	C13-C15-C16-C17
7	AC	503	HEM	CAD-CBD-CGD-O1D
7	AC	502	HEM	CAD-CBD-CGD-O1D
7	BC	501	HEM	CAA-CBA-CGA-O2A
9	A7	103	BCL	C3-C5-C6-C7
7	AC	501	HEM	CAA-CBA-CGA-O1A
7	AC	501	HEM	CAA-CBA-CGA-O2A
7	BC	504	HEM	CAD-CBD-CGD-O1D
9	BT	101	BCL	C4-C3-C5-C6
9	AM	401	BCL	C2-C3-C5-C6
9	AB	101	BCL	C2-C3-C5-C6
9	BP	101	BCL	C2-C3-C5-C6
7	AC	502	HEM	CAA-CBA-CGA-O2A
9	AM	401	BCL	C11-C12-C13-C15
9	AB	101	BCL	C11-C12-C13-C15
9	AD	102	BCL	C11-C10-C8-C7
9	AI	102	BCL	C11-C10-C8-C7
9	AI	102	BCL	C12-C13-C15-C16
9	AP	101	BCL	C11-C12-C13-C15
9	AQ	102	BCL	C6-C7-C8-C10
9	AR	101	BCL	C12-C13-C15-C16
9	AS	103	BCL	C11-C10-C8-C7
9	A6	101	BCL	C12-C13-C15-C16
9	BE	101	BCL	C12-C13-C15-C16
9	BN	101	BCL	C11-C10-C8-C7
9	BW	102	BCL	C11-C12-C13-C15
9	BY	102	BCL	C11-C12-C13-C15
9	B4	101	BCL	C12-C13-C15-C16
9	B8	101	BCL	C12-C13-C15-C16
9	AG	101	BCL	C2C-C3C-CAC-CBC
9	BI	102	BCL	C2C-C3C-CAC-CBC
9	AO	102	BCL	C4-C3-C5-C6

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Mol	Chain	Res	Type	Atoms
9	AP	101	BCL	C4-C3-C5-C6
9	BM	401	BCL	C4-C3-C5-C6
13	AM	405	MQ8	C45-C43-C44-C46
9	B1	102	BCL	CAA-CBA-CGA-O2A
9	BM	401	BCL	C5-C6-C7-C8
9	AI	102	BCL	C2-C3-C5-C6
9	BM	401	BCL	C2-C3-C5-C6
9	B3	102	BCL	C2-C3-C5-C6
9	AL	301	BCL	C2-C1-O2A-CGA
9	BJ	101	BCL	C2-C1-O2A-CGA
9	AL	303	BCL	C14-C13-C15-C16
9	AY	102	BCL	C11-C10-C8-C9
9	BM	402	BCL	C11-C12-C13-C14
15	AS	101	PEF	C35-C36-C37-C38
15	BQ	101	PEF	C35-C36-C37-C38
7	AC	502	HEM	CAA-CBA-CGA-O1A
7	AC	503	HEM	CAA-CBA-CGA-O1A
7	AC	503	HEM	CAD-CBD-CGD-O2D
9	BD	102	BCL	C4-C3-C5-C6
9	BE	101	BCL	C4-C3-C5-C6
9	BF	102	BCL	C4-C3-C5-C6
7	AC	504	HEM	CAD-CBD-CGD-O1D
9	AK	102	BCL	C2-C3-C5-C6
9	AG	101	BCL	C5-C6-C7-C8
9	AL	301	BCL	C4C-C3C-CAC-CBC
9	B8	101	BCL	C4C-C3C-CAC-CBC
9	B9	102	BCL	CAA-CBA-CGA-O2A
9	B1	102	BCL	C4-C3-C5-C6
9	B0	102	BCL	C4-C3-C5-C6
9	AM	401	BCL	C16-C17-C18-C19
9	AM	401	BCL	C16-C17-C18-C20
7	AC	502	HEM	CAD-CBD-CGD-O2D
9	AS	103	BCL	CAA-CBA-CGA-O1A
9	A3	103	BCL	CAA-CBA-CGA-O2A
7	BC	504	HEM	CAD-CBD-CGD-O2D
9	AX	101	BCL	C6-C7-C8-C10
9	BW	102	BCL	C11-C10-C8-C7
9	BK	102	BCL	CAA-CBA-CGA-O2A
9	AO	102	BCL	C6-C7-C8-C9
9	AT	101	BCL	C14-C13-C15-C16
9	A7	103	BCL	C11-C10-C8-C9
9	A9	102	BCL	C11-C10-C8-C9

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Mol	Chain	Res	Type	Atoms
9	BW	102	BCL	C6-C7-C8-C9
9	AN	101	BCL	C2A-CAA-CBA-CGA
9	BQ	103	BCL	C2-C1-O2A-CGA
9	AR	101	BCL	C3A-C2A-CAA-CBA
9	AD	102	BCL	CAA-CBA-CGA-O2A
9	AP	101	BCL	C2-C3-C5-C6
15	AM	407	PEF	C1-C2-O2-C10
15	AH	301	PEF	C1-C2-O2-C10
7	AC	504	HEM	CAD-CBD-CGD-O2D
7	BC	503	HEM	CAD-CBD-CGD-O2D
9	AB	101	BCL	C2A-CAA-CBA-CGA
9	AP	101	BCL	C2A-CAA-CBA-CGA
7	BC	503	HEM	CAD-CBD-CGD-O1D
9	BV	101	BCL	C4-C3-C5-C6
9	BZ	101	BCL	C13-C15-C16-C17
9	AZ	101	BCL	C14-C13-C15-C16
9	A0	102	BCL	C14-C13-C15-C16
9	BG	101	BCL	C6-C7-C8-C9
9	BY	102	BCL	CAA-CBA-CGA-O1A
7	AC	503	HEM	C2D-C3D-CAD-CBD
9	AA	101	BCL	C16-C17-C18-C19
9	AJ	101	BCL	C2A-CAA-CBA-CGA
9	AL	303	BCL	C12-C13-C15-C16
9	AW	101	BCL	C11-C10-C8-C7
9	AY	102	BCL	C11-C10-C8-C7
9	A7	103	BCL	C11-C10-C8-C7
9	BD	102	BCL	C11-C10-C8-C7
9	BF	102	BCL	C11-C12-C13-C15
9	BS	102	BCL	C11-C10-C8-C7
9	BU	102	BCL	C6-C7-C8-C10
9	B9	102	BCL	C12-C13-C15-C16
9	B0	102	BCL	C12-C13-C15-C16
10	BM	403	BPH	C11-C10-C8-C7
10	BM	403	BPH	C12-C13-C15-C16
9	AM	402	BCL	C2-C1-O2A-CGA
9	A5	102	BCL	C2-C1-O2A-CGA
9	BL	301	BCL	C2-C1-O2A-CGA
9	BL	303	BCL	C8-C10-C11-C12
9	B2	101	BCL	CAA-CBA-CGA-O2A
9	BM	402	BCL	C16-C17-C18-C19
9	AQ	102	BCL	C2-C3-C5-C6
9	A0	102	BCL	C2-C3-C5-C6

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Mol	Chain	Res	Type	Atoms
9	B6	101	BCL	C2A-CAA-CBA-CGA
9	B8	101	BCL	CAA-CBA-CGA-O2A
9	BX	101	BCL	C2A-CAA-CBA-CGA
9	AM	401	BCL	C6-C7-C8-C9
9	AN	101	BCL	C6-C7-C8-C9
9	BL	303	BCL	C14-C13-C15-C16
9	BG	101	BCL	C14-C13-C15-C16
9	B5	102	BCL	C11-C12-C13-C14
9	B6	101	BCL	C11-C10-C8-C9
9	BW	102	BCL	CAA-CBA-CGA-O2A
9	AB	101	BCL	C16-C17-C18-C19
9	AQ	102	BCL	C4-C3-C5-C6
9	B2	101	BCL	C4-C3-C5-C6
9	A2	101	BCL	C8-C10-C11-C12
9	A9	102	BCL	CAA-CBA-CGA-O2A
9	AB	101	BCL	C16-C17-C18-C20
9	B7	103	BCL	C16-C17-C18-C20
9	AE	101	BCL	CAA-CBA-CGA-O2A
9	BM	402	BCL	CAA-CBA-CGA-O2A
9	BU	102	BCL	C16-C17-C18-C19
9	A0	102	BCL	C4-C3-C5-C6
15	AM	409	PEF	C37-C38-C39-C40
9	AO	102	BCL	C2-C1-O2A-CGA
9	BA	101	BCL	C2-C1-O2A-CGA
9	BO	102	BCL	C2-C1-O2A-CGA
9	AM	402	BCL	C11-C10-C8-C7
9	A9	102	BCL	C11-C10-C8-C7
9	BM	402	BCL	C11-C10-C8-C7
9	BJ	101	BCL	C6-C7-C8-C10
9	BX	101	BCL	C6-C7-C8-C10
9	B4	101	BCL	C6-C7-C8-C10
9	AS	103	BCL	C2-C3-C5-C6
9	BD	102	BCL	C5-C6-C7-C8
9	AO	102	BCL	CAA-CBA-CGA-O2A
9	AQ	102	BCL	C15-C16-C17-C18
10	BM	403	BPH	C15-C16-C17-C18
13	BM	405	MQ8	C20-C21-C22-C23
9	AB	101	BCL	CAA-CBA-CGA-O2A
9	BM	402	BCL	C16-C17-C18-C20
13	AM	405	MQ8	C42-C43-C44-C46
15	BQ	101	PEF	C13-C14-C15-C16
9	BD	102	BCL	C3-C5-C6-C7

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Mol	Chain	Res	Type	Atoms
9	A9	102	BCL	CAA-CBA-CGA-O1A
9	AL	303	BCL	C11-C12-C13-C14
9	AX	101	BCL	C14-C13-C15-C16
9	A3	103	BCL	C11-C10-C8-C9
9	A6	101	BCL	C14-C13-C15-C16
9	A8	101	BCL	C14-C13-C15-C16
9	BY	102	BCL	C11-C12-C13-C14
9	BL	301	BCL	C8-C10-C11-C12
9	B3	102	BCL	CAA-CBA-CGA-O2A
9	B2	101	BCL	CAA-CBA-CGA-O1A
9	BM	402	BCL	CAA-CBA-CGA-O1A
9	AM	402	BCL	C13-C15-C16-C17
9	AO	102	BCL	C2-C3-C5-C6
9	AM	402	BCL	C3-C5-C6-C7
9	AR	101	BCL	C3-C5-C6-C7
9	AI	102	BCL	C10-C11-C12-C13
9	AD	102	BCL	CAD-CBD-CGD-O2D
9	BS	102	BCL	CAD-CBD-CGD-O2D
9	B9	102	BCL	CAD-CBD-CGD-O2D
9	BB	101	BCL	CAA-CBA-CGA-O2A
15	AS	101	PEF	O3-C30-C31-C32
9	A6	101	BCL	C2-C1-O2A-CGA
9	BE	101	BCL	C2-C1-O2A-CGA
9	BX	101	BCL	C2-C1-O2A-CGA
9	BY	102	BCL	C2-C1-O2A-CGA
9	B4	101	BCL	C2-C1-O2A-CGA
15	BQ	101	PEF	O3-C30-C31-C32
9	BW	102	BCL	CAA-CBA-CGA-O1A
15	AS	101	PEF	C10-C11-C12-C13
9	AV	102	BCL	CAA-CBA-CGA-O2A
9	BS	102	BCL	C5-C6-C7-C8
15	BQ	101	PEF	O5-C30-C31-C32

There are no ring outliers.

131 monomers are involved in 3136 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
9	B9	102	BCL	38	0
16	AH	302	PO4	1	0
9	AD	102	BCL	27	0
9	BQ	103	BCL	30	0
14	A5	103	CRT	29	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
14	BG	102	CRT	13	0
15	BM	407	PEF	2	0
7	AC	504	HEM	5	0
7	AC	503	HEM	13	0
13	AM	405	MQ8	10	0
7	BC	502	HEM	7	0
9	AM	401	BCL	15	0
14	AB	102	CRT	46	0
9	BX	101	BCL	33	0
16	BH	301	PO4	1	0
9	BT	101	BCL	22	0
7	AC	501	HEM	9	0
11	BL	304	UQ8	12	0
9	AM	402	BCL	31	0
9	AW	101	BCL	40	0
9	BJ	101	BCL	26	0
9	BY	102	BCL	35	0
14	B2	102	CRT	75	0
9	A8	101	BCL	46	0
9	BF	102	BCL	39	0
9	A0	102	BCL	69	0
9	AV	102	BCL	30	0
9	AL	301	BCL	40	0
14	BV	102	CRT	59	0
9	AF	102	BCL	38	0
14	AW	102	CRT	32	0
9	BI	102	BCL	50	0
9	BL	303	BCL	13	0
9	BG	101	BCL	46	0
9	A5	102	BCL	34	0
14	AJ	102	CRT	21	0
14	BN	102	CRT	15	0
9	BK	102	BCL	26	0
14	AM	406	CRT	13	0
14	AT	102	CRT	18	0
14	BU	103	CRT	61	0
9	BD	102	BCL	34	0
14	BF	103	CRT	19	0
7	AC	502	HEM	11	0
9	AL	303	BCL	21	0
15	AH	301	PEF	8	0
9	BL	301	BCL	33	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
14	AX	102	CRT	47	0
14	AG	102	CRT	11	0
9	A9	102	BCL	36	0
9	AK	102	BCL	58	0
15	AM	408	PEF	6	0
15	BQ	101	PEF	12	0
9	AO	102	BCL	50	0
9	BM	402	BCL	35	0
9	AA	101	BCL	42	0
10	AL	302	BPH	17	0
9	AR	101	BCL	34	0
9	AS	103	BCL	32	0
9	A6	101	BCL	27	0
10	AM	403	BPH	12	0
9	BB	101	BCL	47	0
14	BO	103	CRT	13	0
9	AG	101	BCL	36	0
9	B8	101	BCL	36	0
14	AP	102	CRT	24	0
9	B2	101	BCL	43	0
9	AP	101	BCL	37	0
9	AE	101	BCL	32	0
7	BC	504	HEM	9	0
14	BP	102	CRT	29	0
9	AI	102	BCL	36	0
9	A3	104	BCL	25	0
7	BC	501	HEM	7	0
9	B0	102	BCL	57	0
13	BM	405	MQ8	8	0
9	BA	101	BCL	38	0
9	B1	102	BCL	37	0
9	BU	102	BCL	38	0
9	AT	101	BCL	25	0
9	BE	101	BCL	46	0
14	BW	103	CRT	20	0
9	A3	103	BCL	42	0
14	A0	101	CRT	28	0
9	AQ	102	BCL	25	0
9	BM	401	BCL	14	0
9	BN	101	BCL	30	0
9	A1	102	BCL	52	0
14	B1	103	CRT	41	0

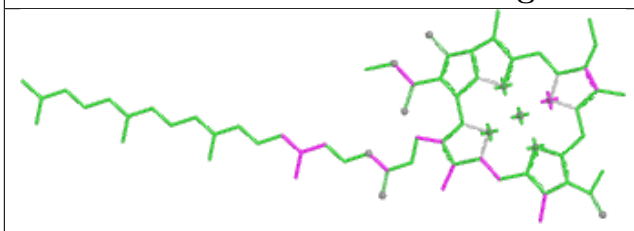
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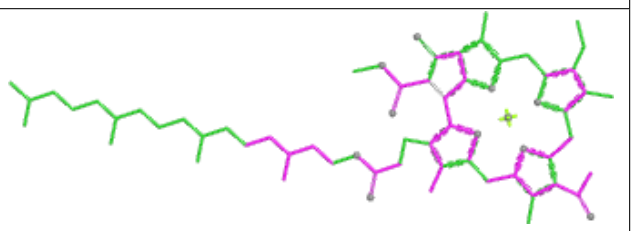
Mol	Chain	Res	Type	Clashes	Symm-Clashes
11	AL	304	UQ8	8	0
9	B5	102	BCL	20	0
14	AS	104	CRT	77	0
9	BZ	101	BCL	31	0
9	B6	101	BCL	24	0
9	AN	101	BCL	53	0
9	AU	102	BCL	46	0
9	BO	102	BCL	48	0
14	BA	102	CRT	26	0
10	BM	403	BPH	7	0
9	B7	103	BCL	39	0
14	AN	102	CRT	15	0
9	BV	101	BCL	23	0
14	B5	103	CRT	22	0
9	BW	102	BCL	42	0
14	BB	102	CRT	27	0
15	AM	407	PEF	3	0
15	AS	101	PEF	37	0
14	A7	102	CRT	46	0
14	A2	102	CRT	51	0
14	AR	102	CRT	20	0
9	A7	103	BCL	52	0
14	A1	103	CRT	29	0
14	BS	103	CRT	10	0
7	BC	503	HEM	11	0
14	B7	102	CRT	39	0
14	AA	102	CRT	24	0
14	B0	101	CRT	44	0
9	AX	101	BCL	33	0
9	BS	102	BCL	28	0
9	A2	101	BCL	28	0
9	BP	101	BCL	35	0
9	B4	101	BCL	20	0
9	AB	101	BCL	35	0
9	AJ	101	BCL	43	0
10	BL	302	BPH	11	0
9	BQ	104	BCL	34	0
9	AZ	101	BCL	41	0
9	AY	102	BCL	48	0
9	B3	102	BCL	47	0
15	AM	409	PEF	14	0
14	BM	406	CRT	8	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less then 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

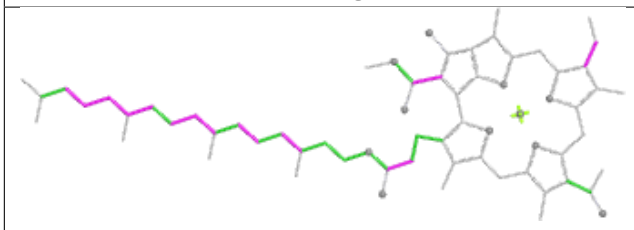
Ligand BCL B9 102



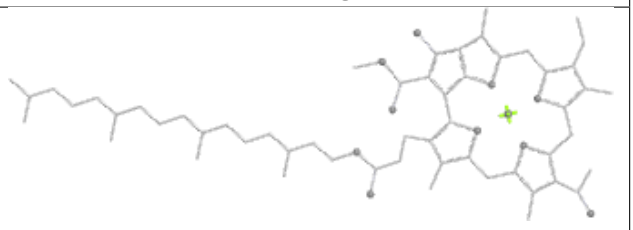
Bond lengths



Bond angles

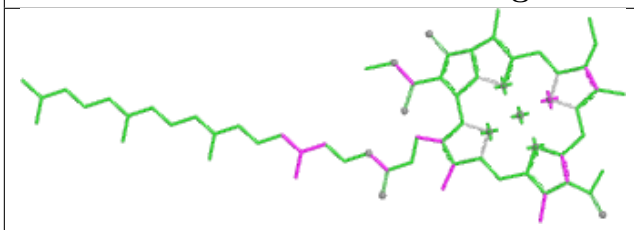


Torsions

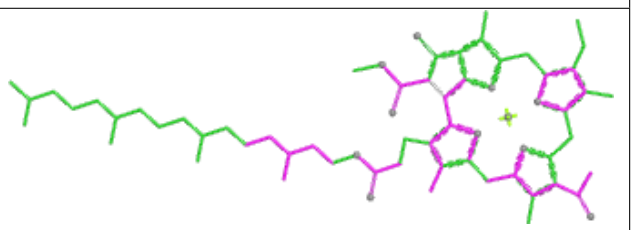


Rings

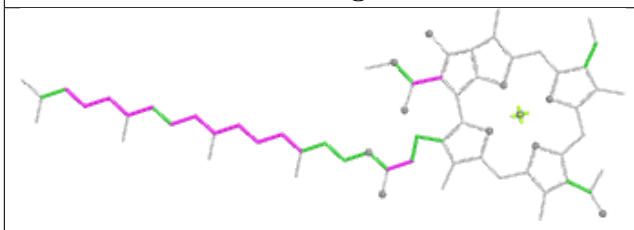
Ligand BCL AD 102



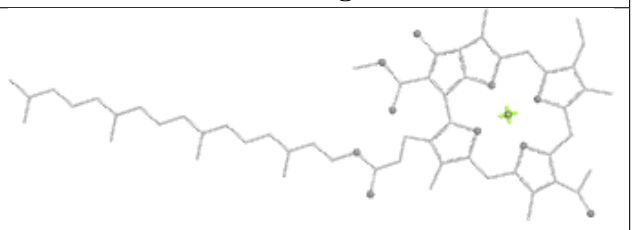
Bond lengths



Bond angles

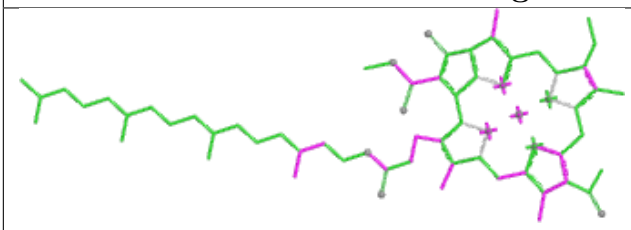


Torsions

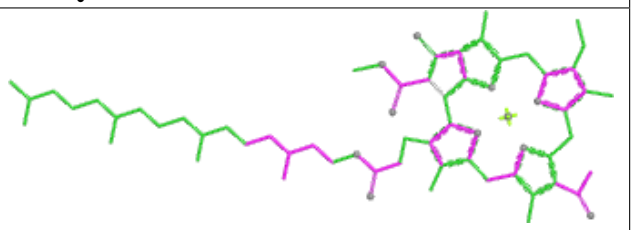


Rings

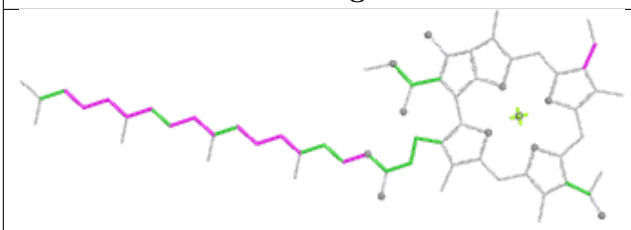
Ligand BCL BQ 103



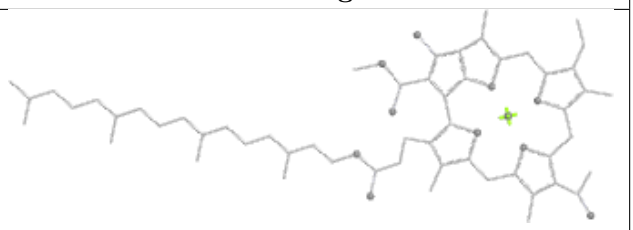
Bond lengths



Bond angles

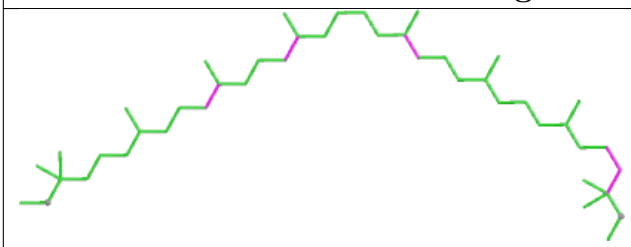


Torsions

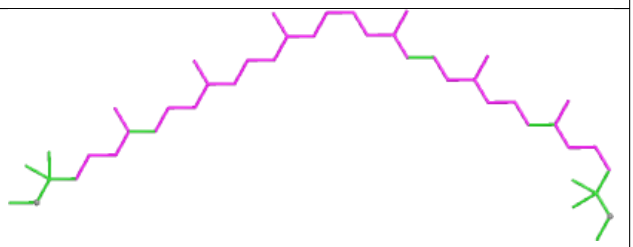


Rings

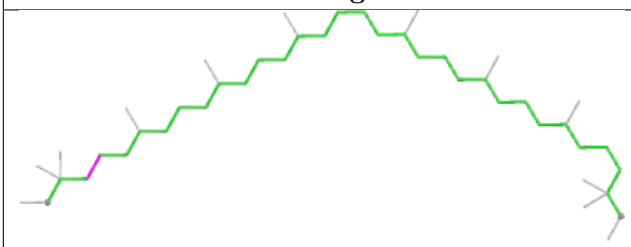
Ligand CRT A5 103



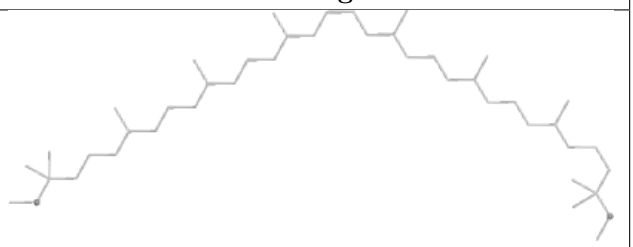
Bond lengths



Bond angles

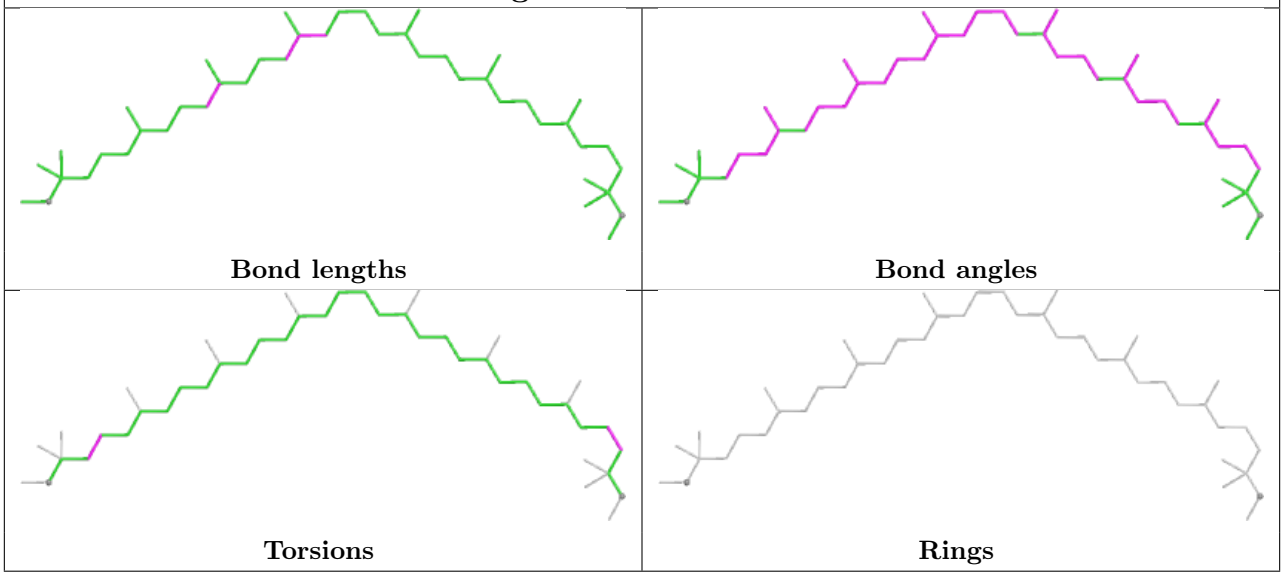


Torsions

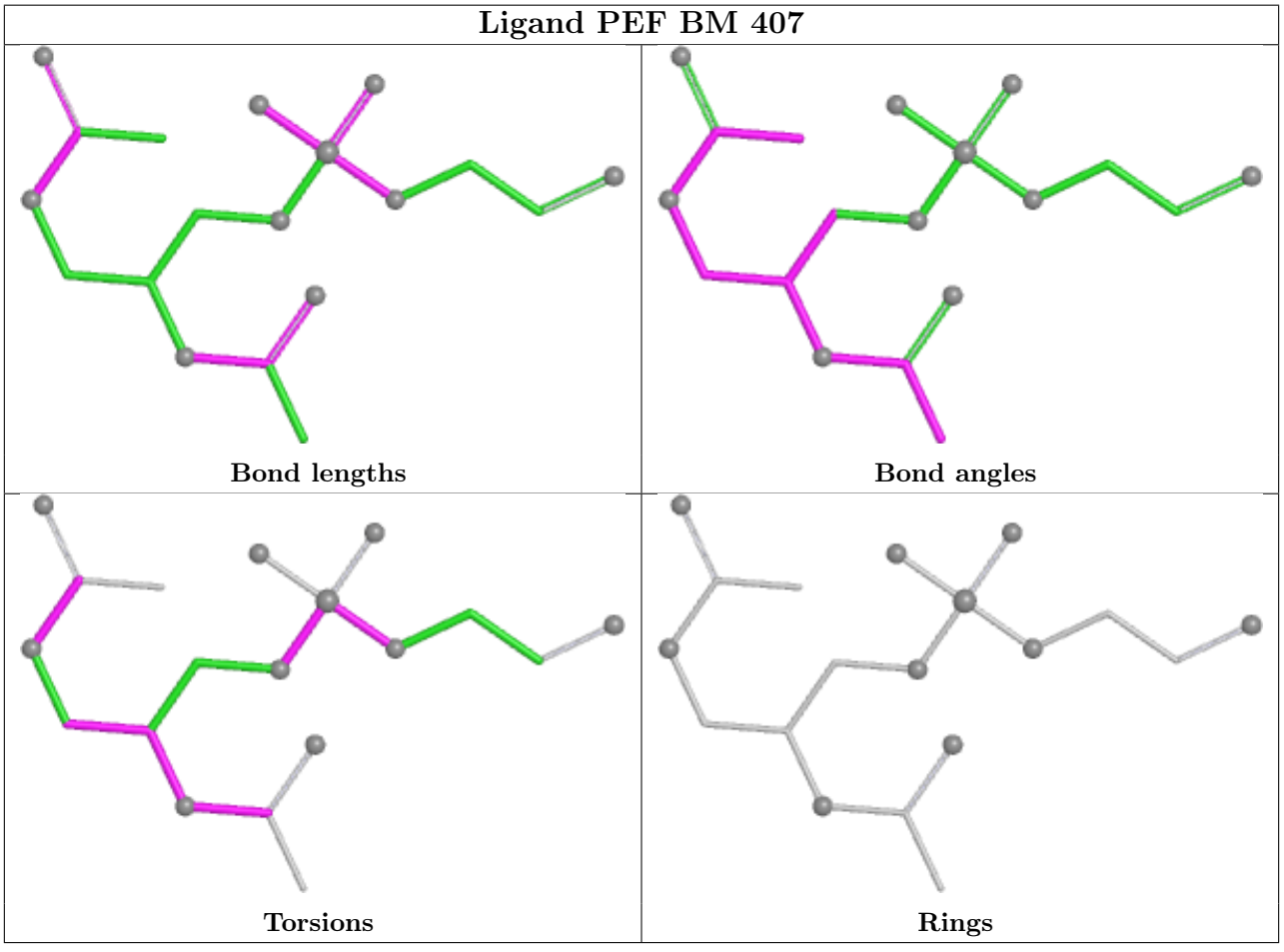


Rings

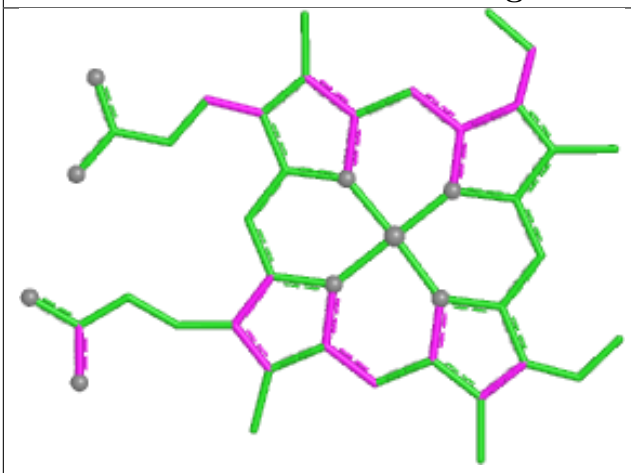
Ligand CRT BG 102



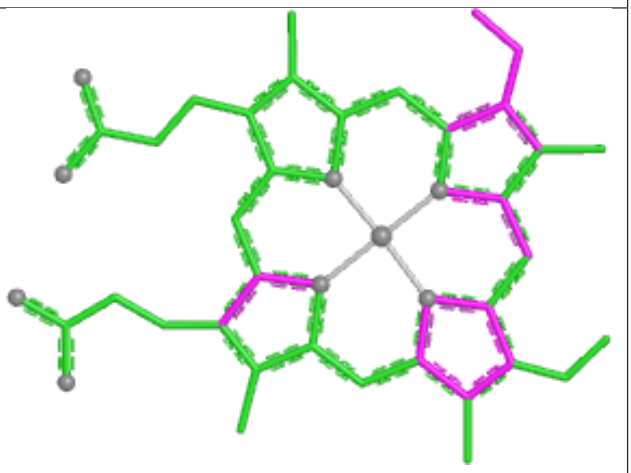
Ligand PEF BM 407



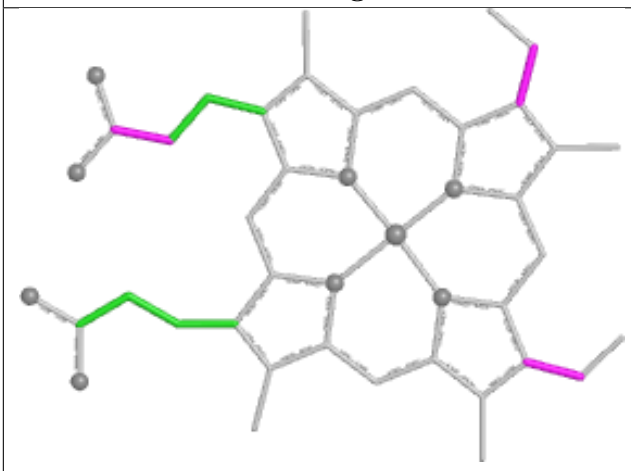
Ligand HEM AC 504



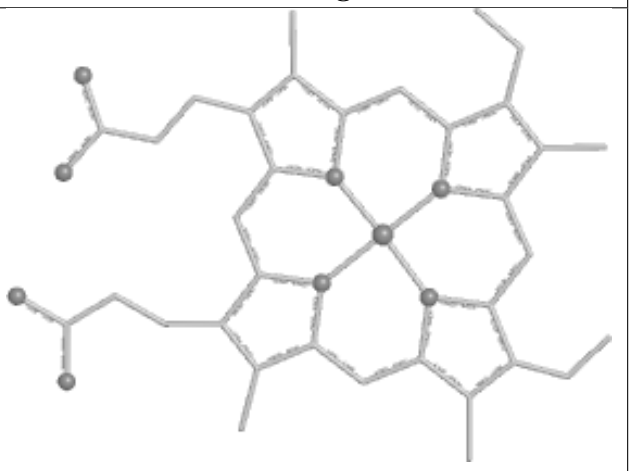
Bond lengths



Bond angles

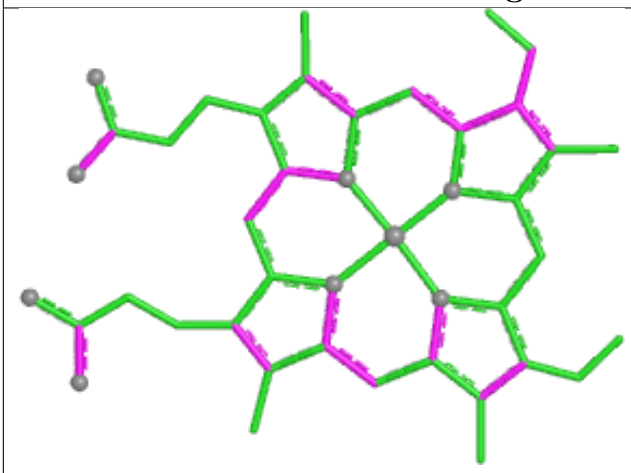


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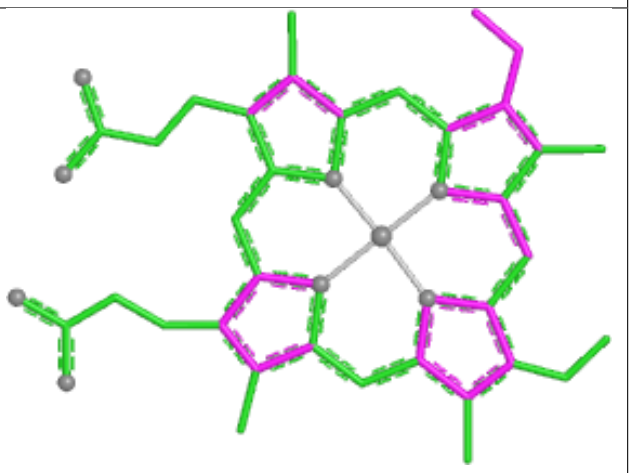


Rings

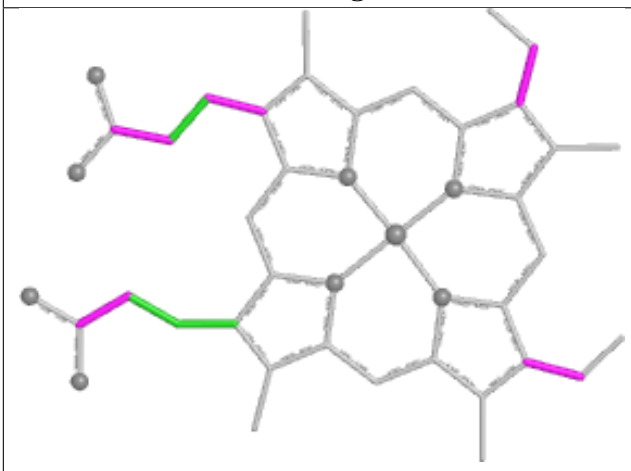
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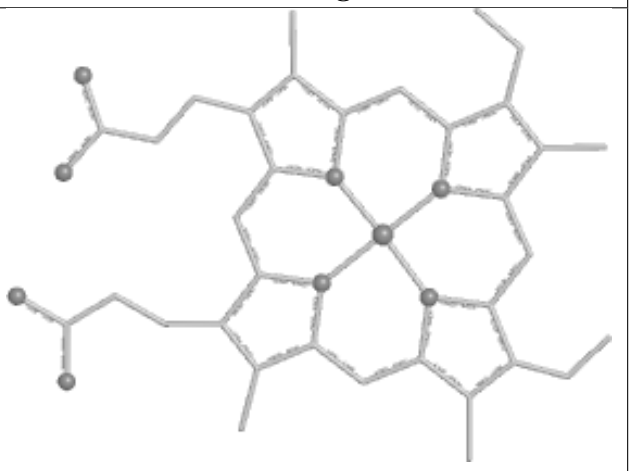
Bond lengths



Bond angles

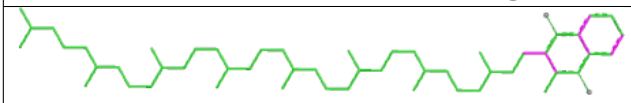


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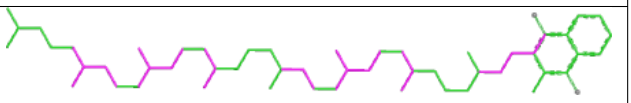


Rings

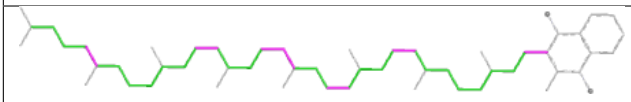
Ligand MQ8 AM 405



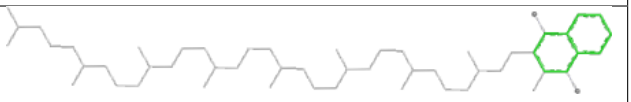
Bond lengths



Bond angles

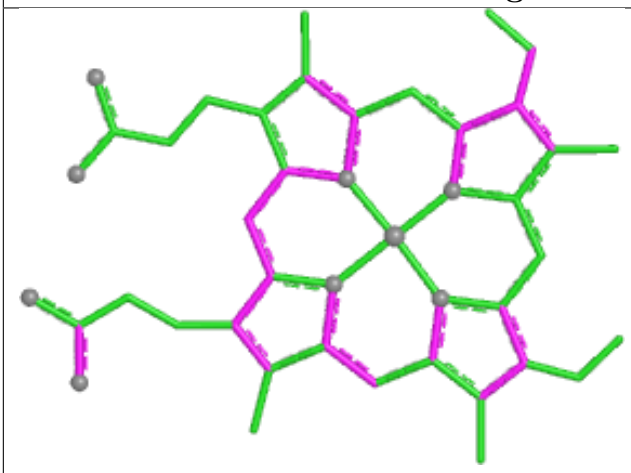


Torsions

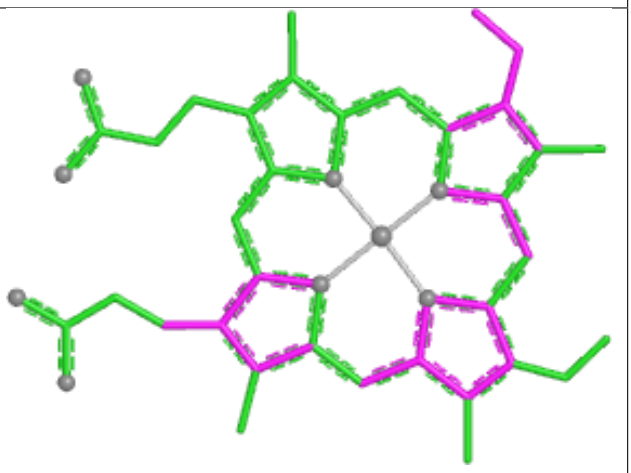


Rings

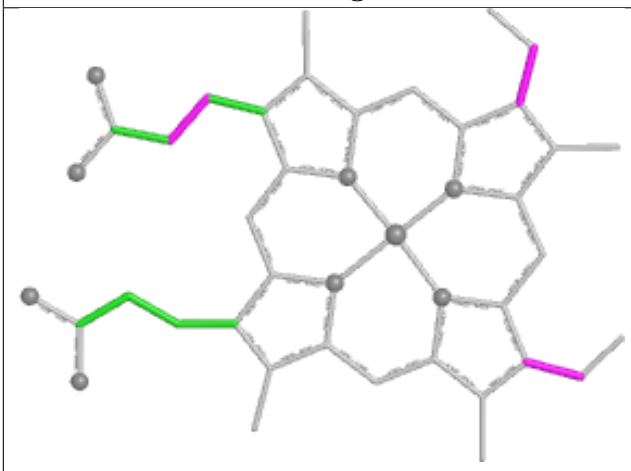
Ligand HEM BC 502



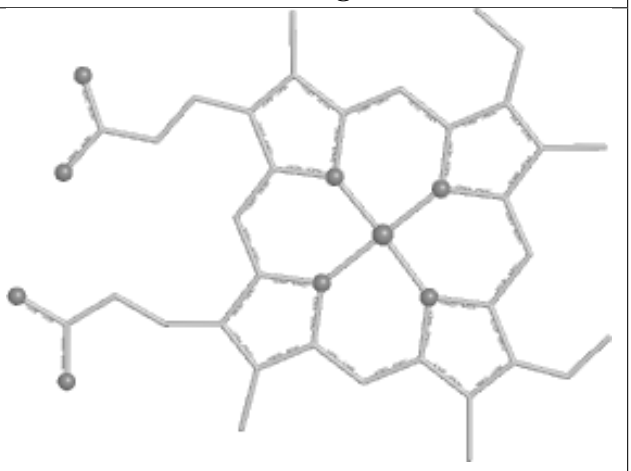
Bond lengths



Bond angles

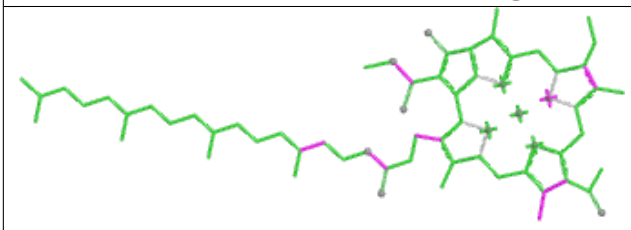


Torsions

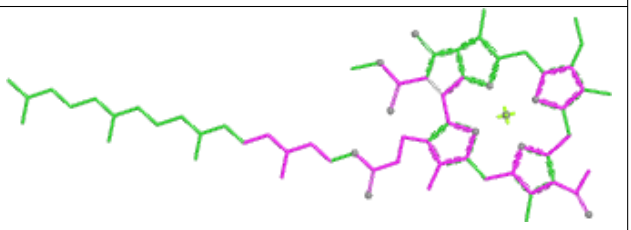


Rings

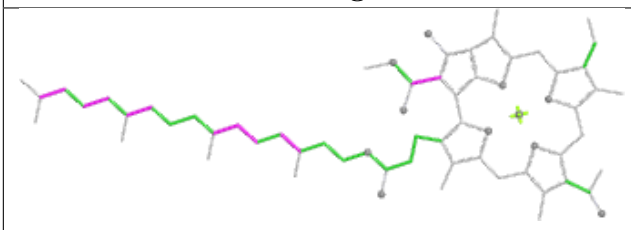
Ligand BCL AM 401



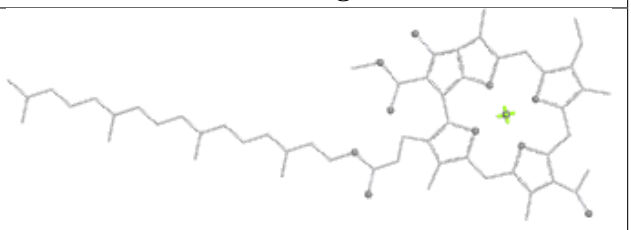
Bond lengths



Bond angles

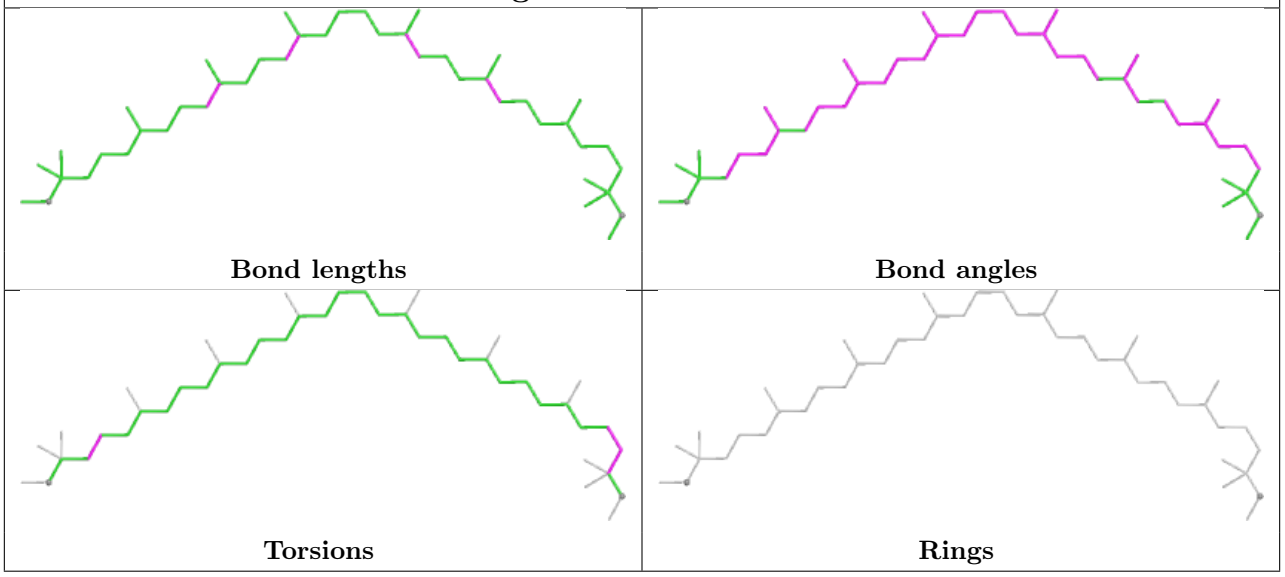


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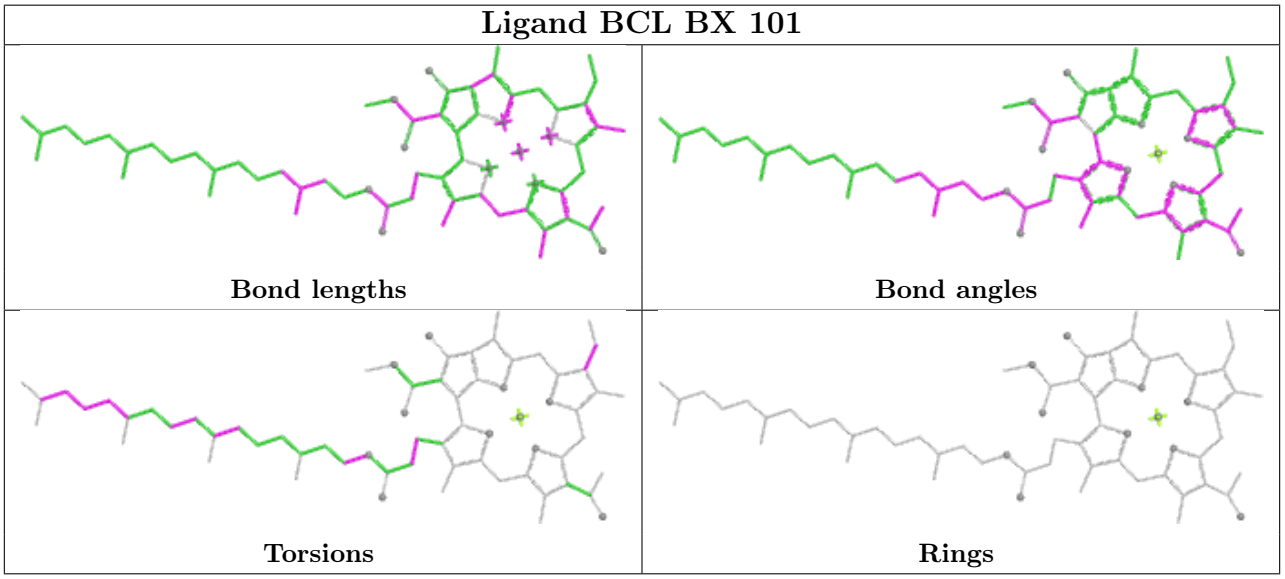


Rings

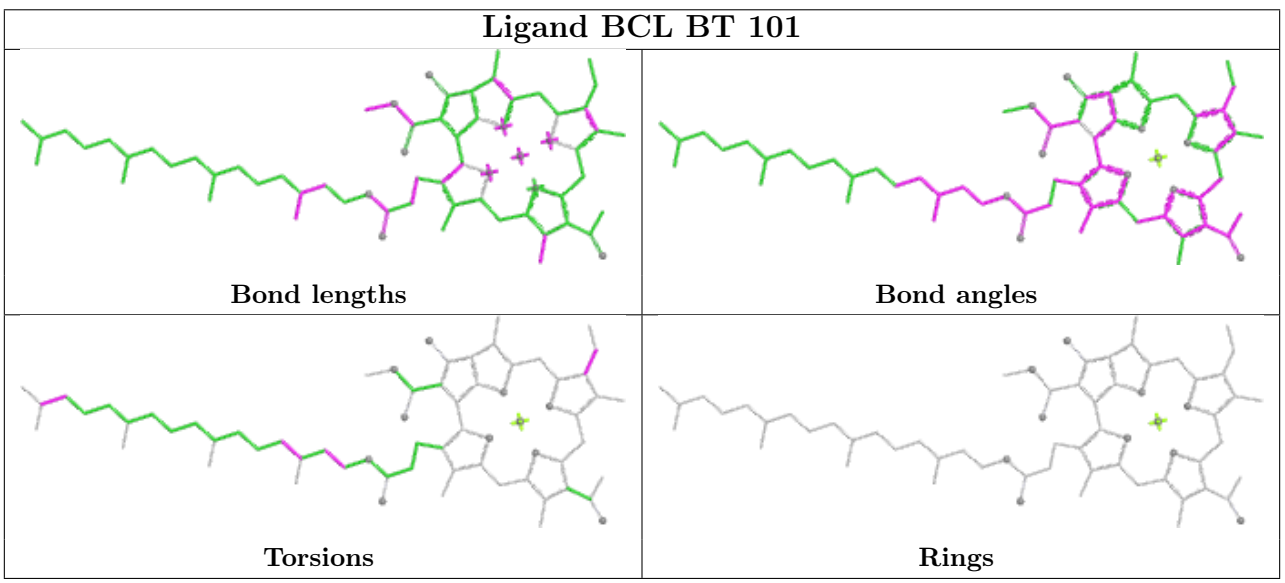
Ligand CRT AB 102



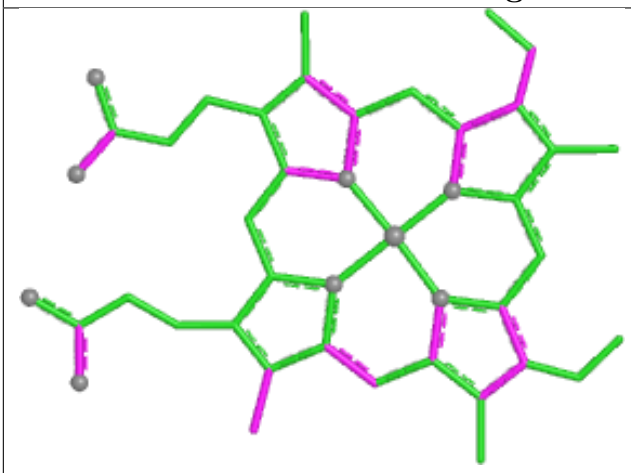
Ligand BCL BX 101



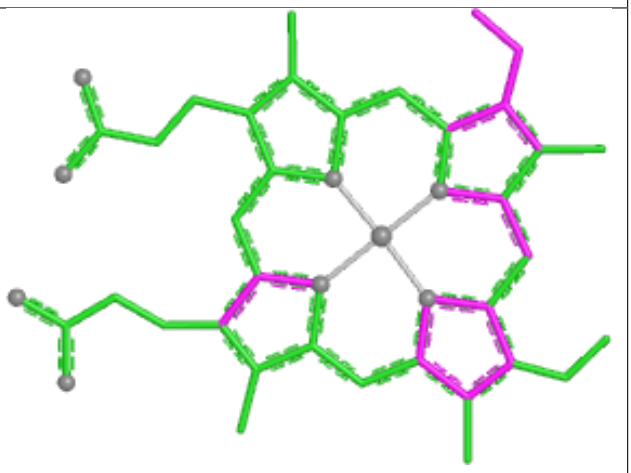
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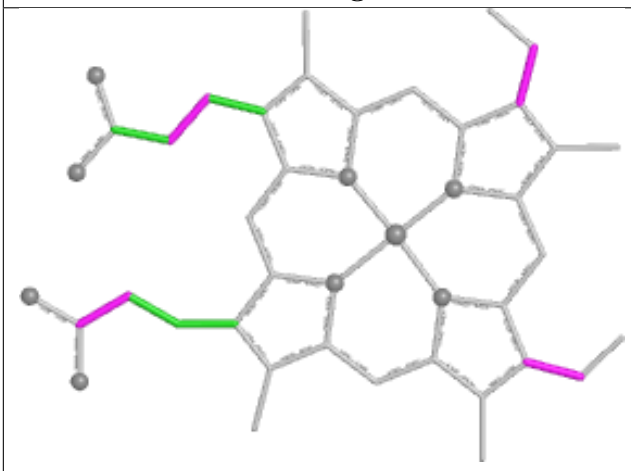
Ligand HEM AC 501



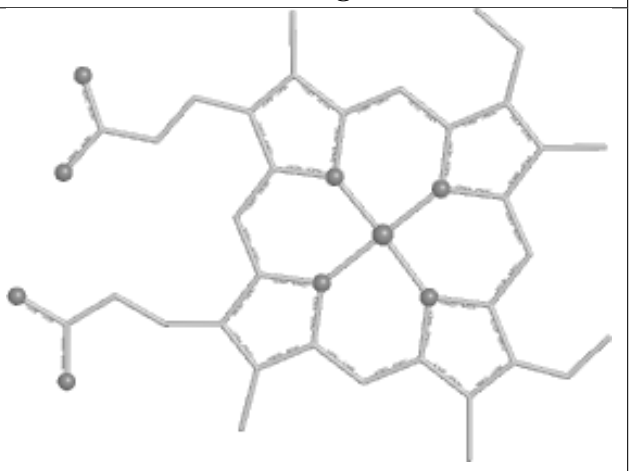
Bond lengths



Bond angles

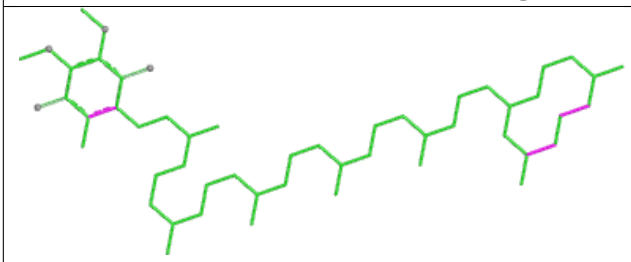


Torsions

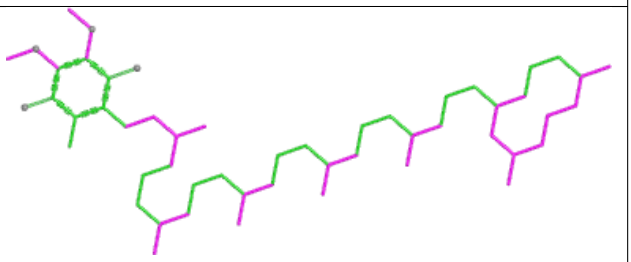


Rings

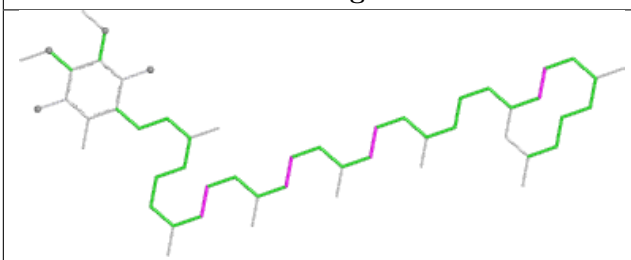
Ligand UQ8 BL 304



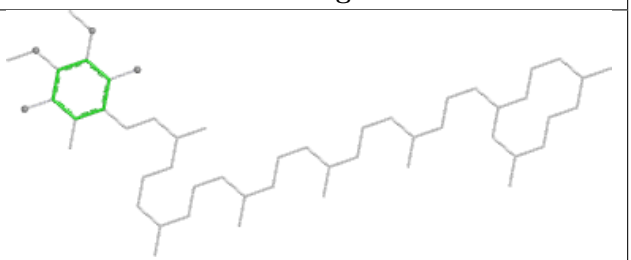
Bond lengths



Bond angles

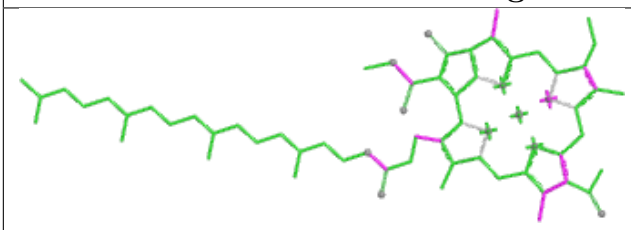


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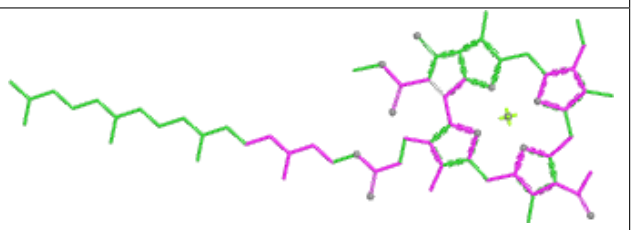


Rings

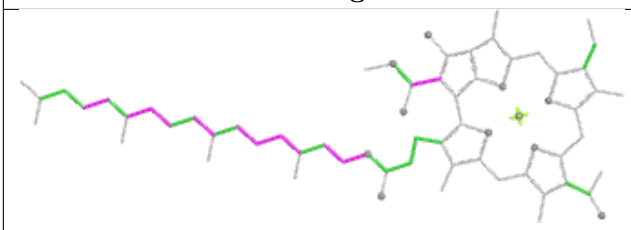
Ligand BCL AM 402



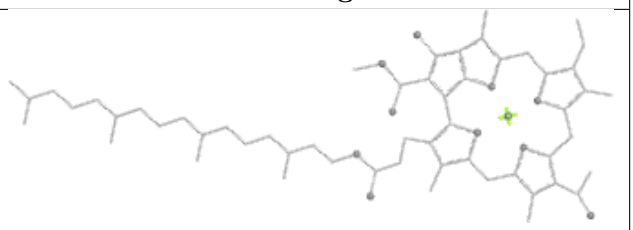
Bond lengths



Bond angles

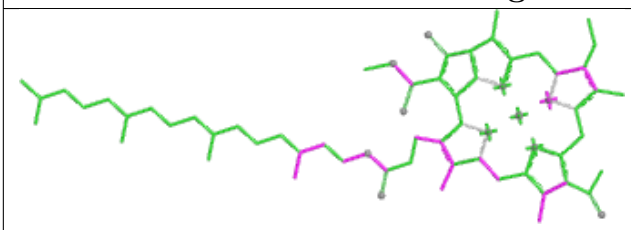


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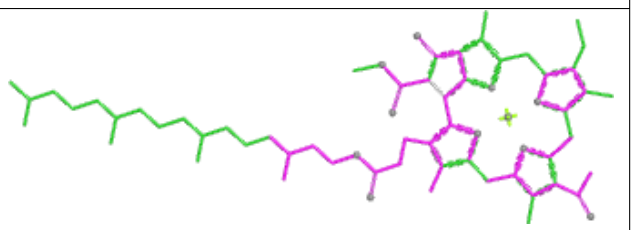


Rings

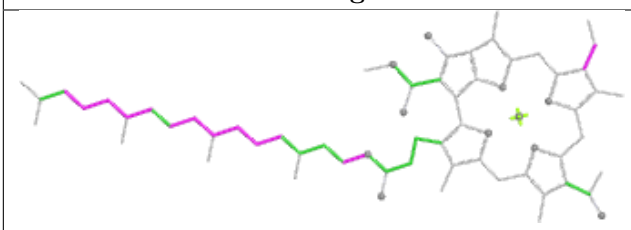
Ligand BCL AW 101



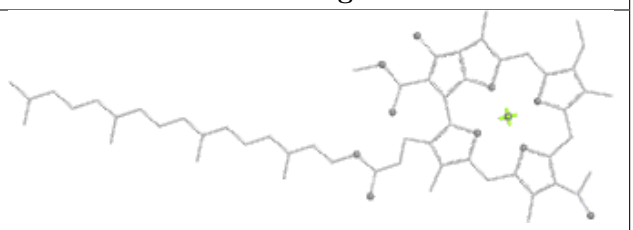
Bond lengths



Bond angles

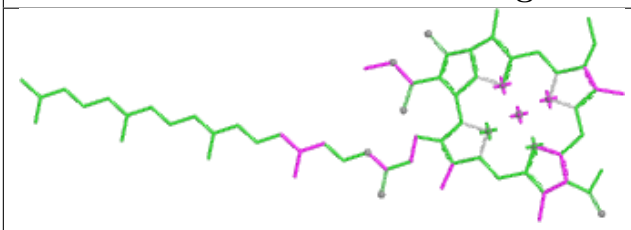


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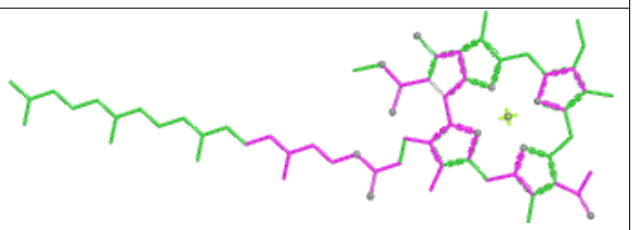


Rings

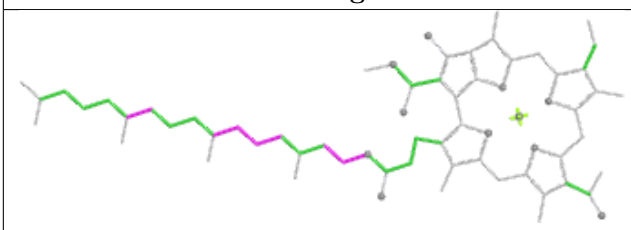
Ligand BCL BJ 101



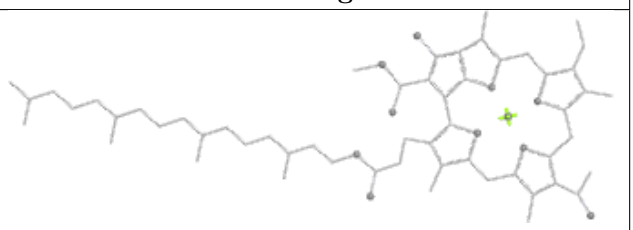
Bond lengths



Bond angles

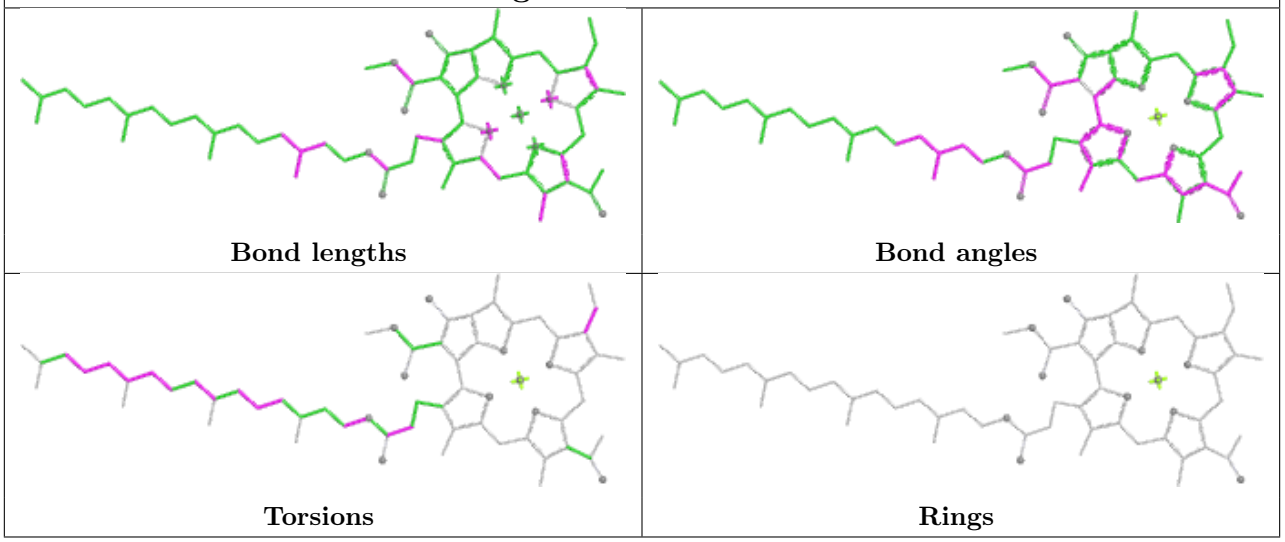


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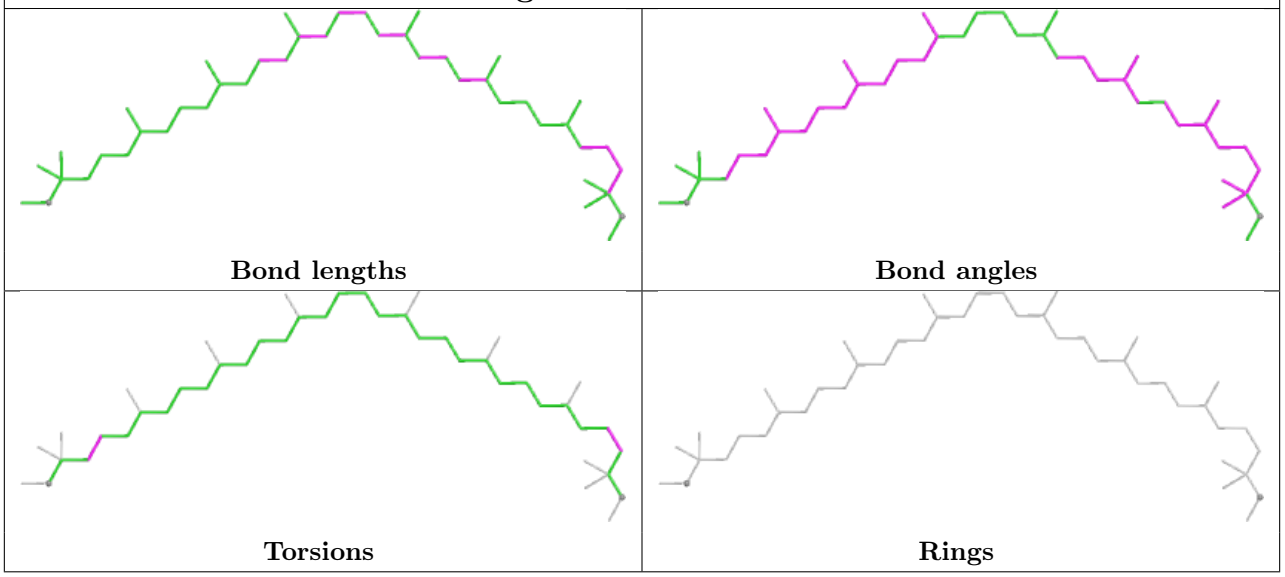


Rings

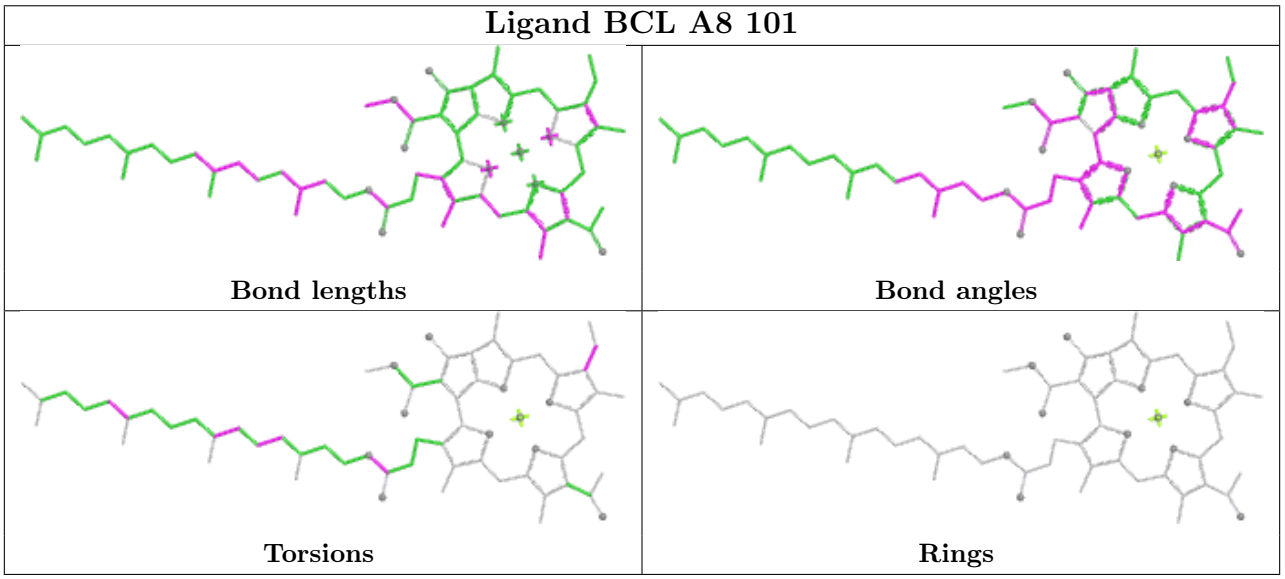
Ligand BCL BY 102



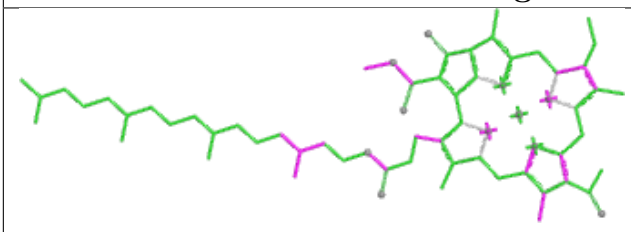
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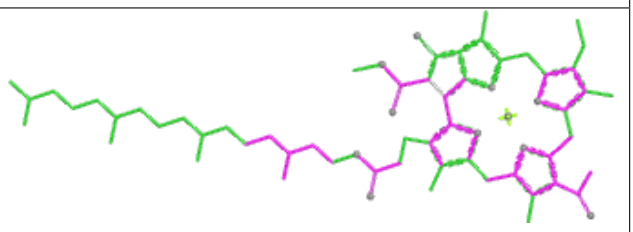
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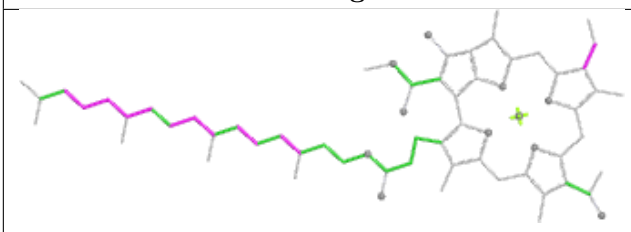
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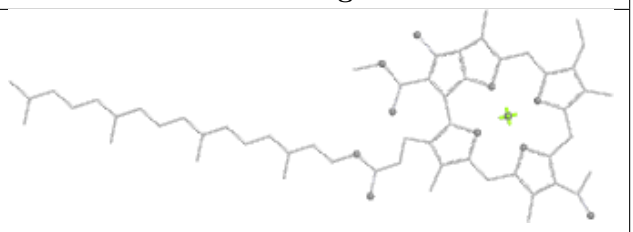
Bond lengths



Bond angles

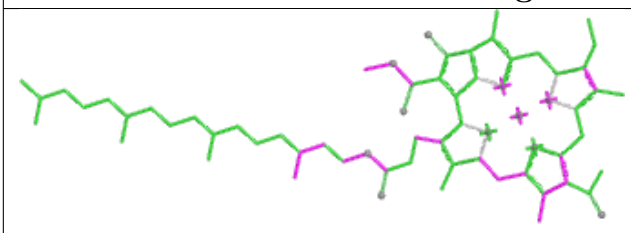


Torsions

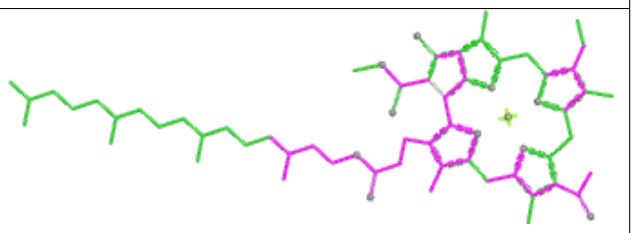


Rings

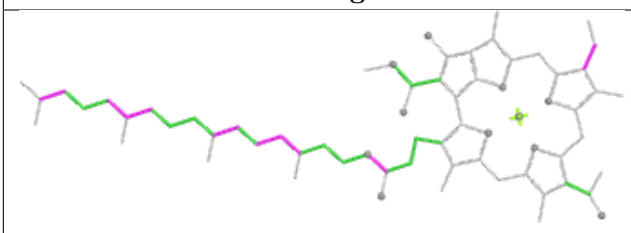
Ligand BCL A0 102



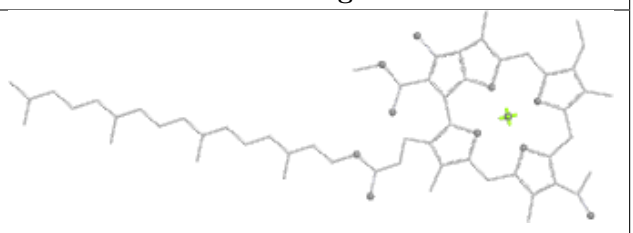
Bond lengths



Bond angles

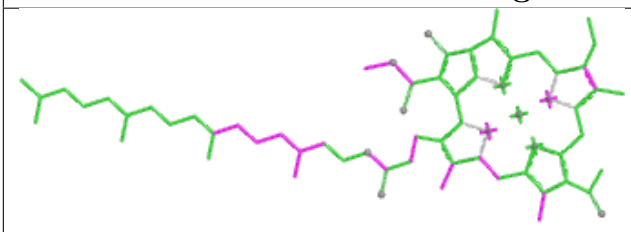


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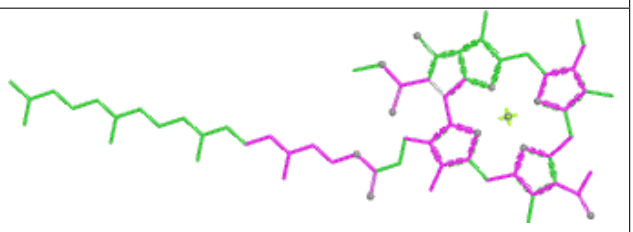


Rings

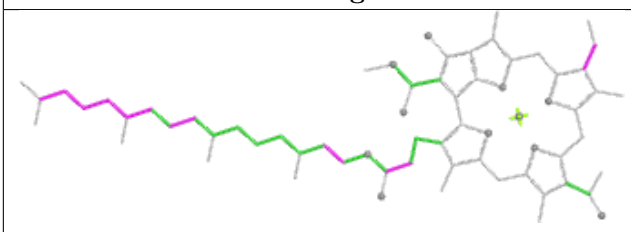
Ligand BCL AV 102



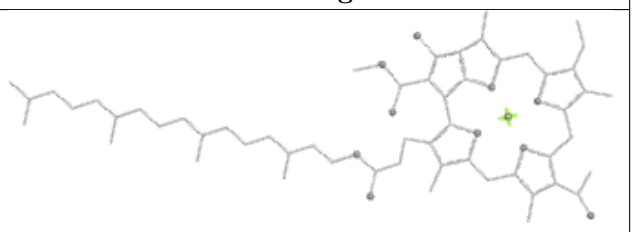
Bond lengths



Bond angles

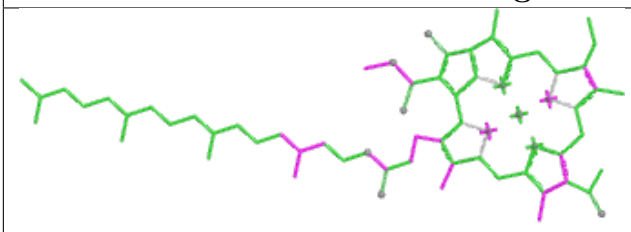


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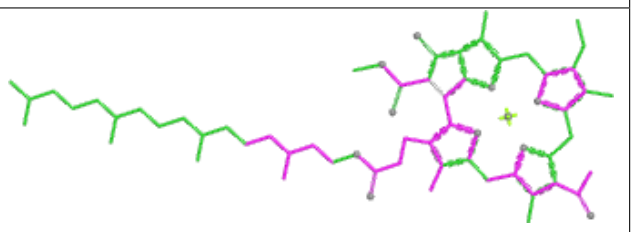


Rings

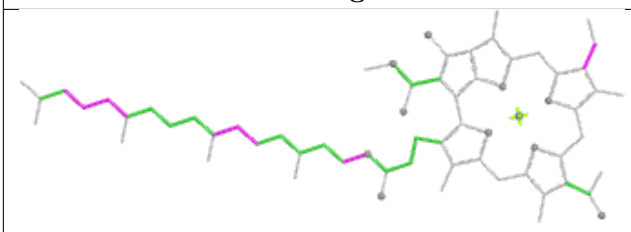
Ligand BCL AL 301



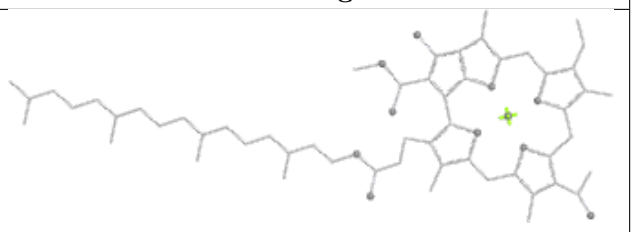
Bond lengths



Bond angles

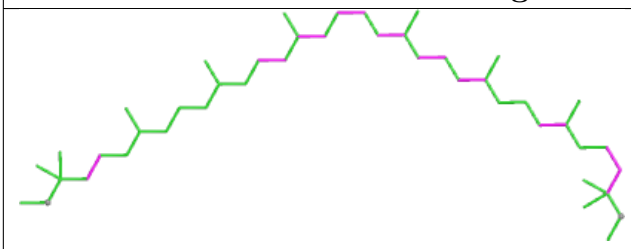


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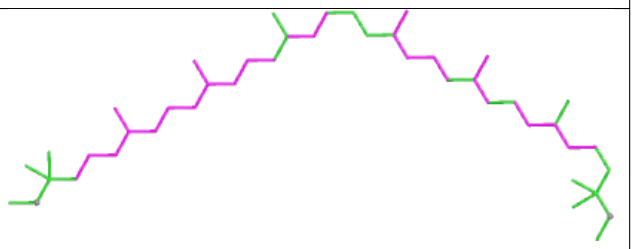


Rings

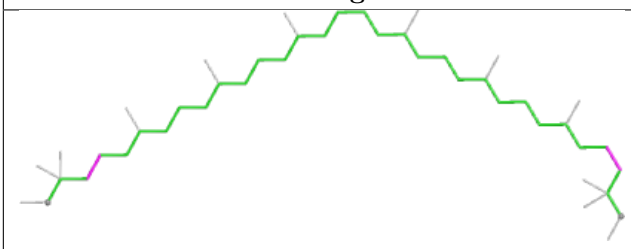
Ligand CRT BV 102



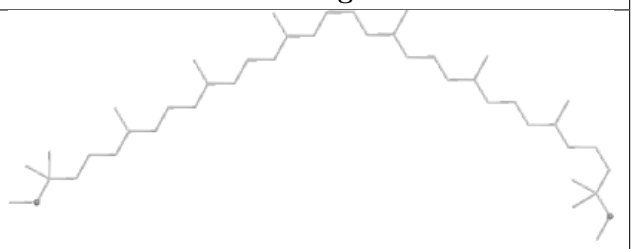
Bond lengths



Bond angles

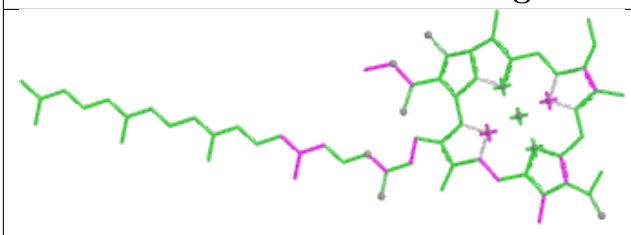


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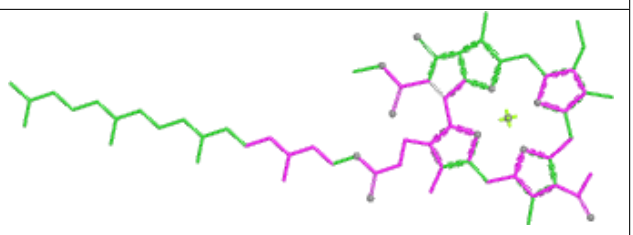


Rings

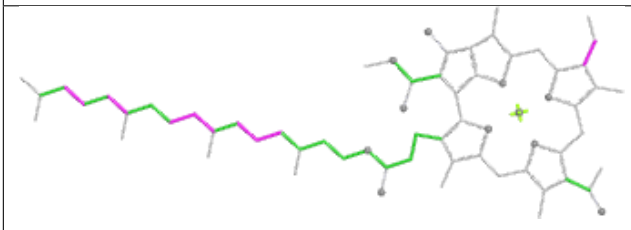
Ligand BCL AF 102



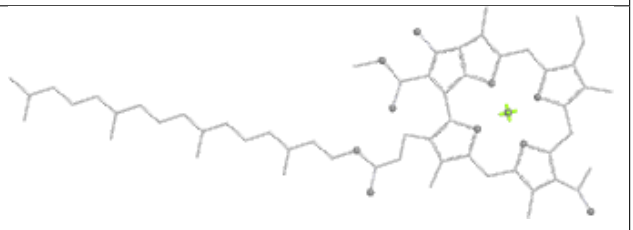
Bond lengths



Bond angles

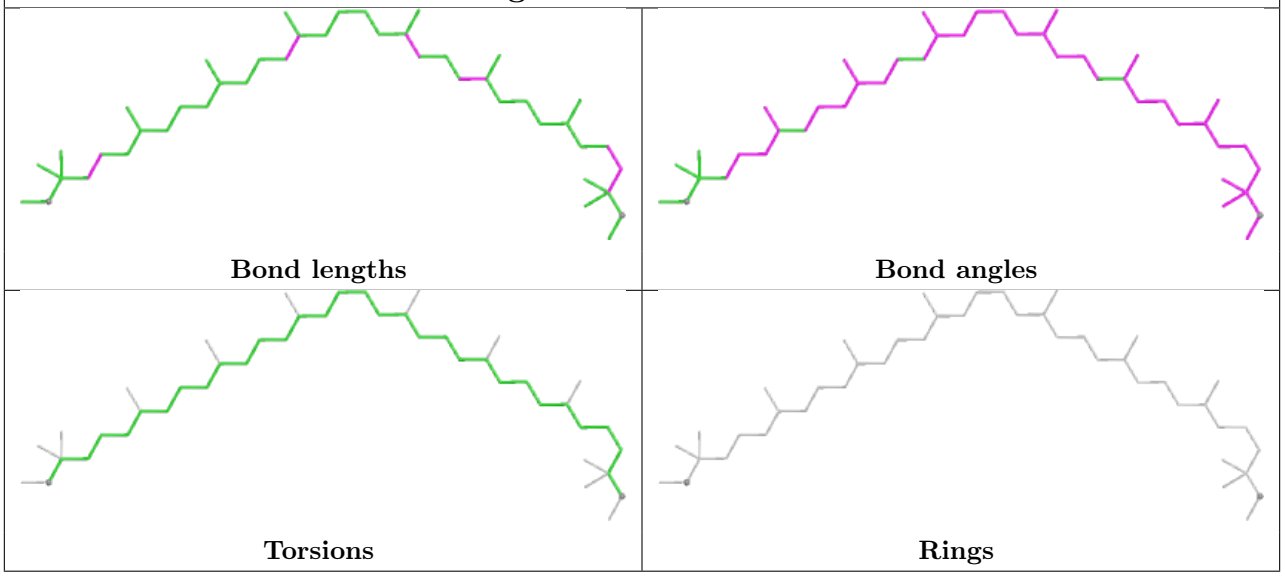


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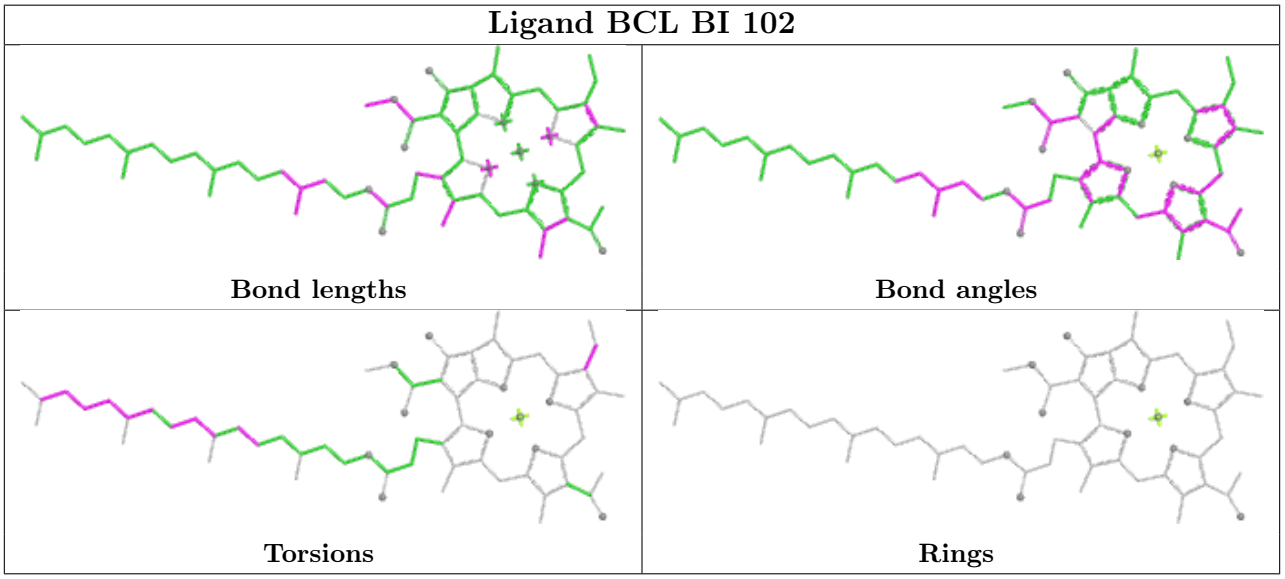


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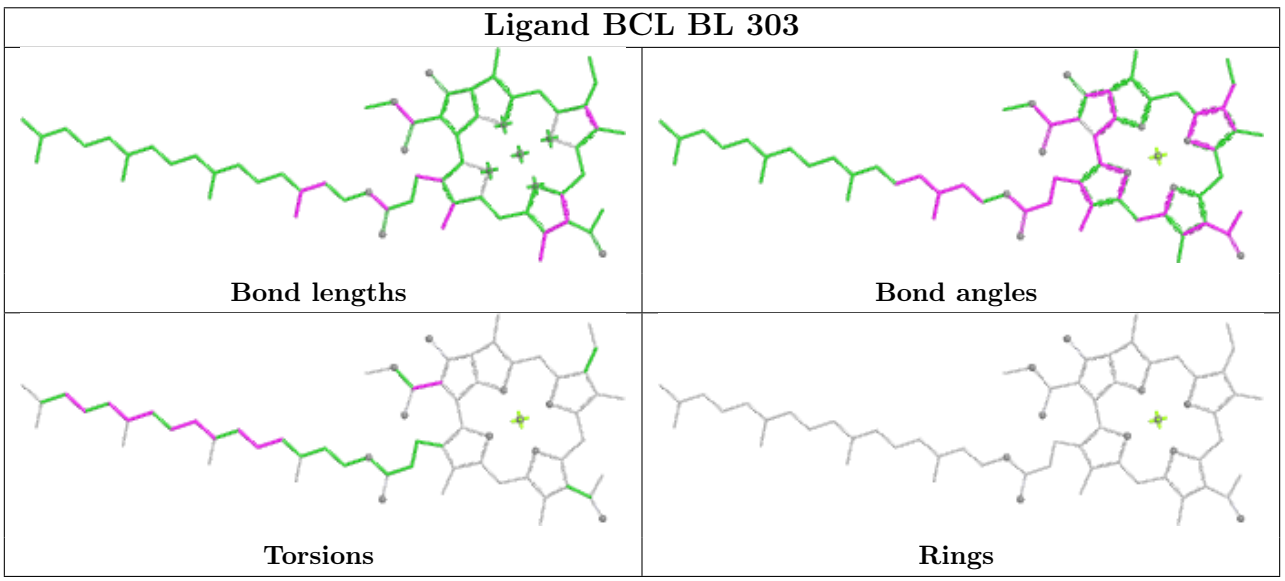
Ligand CRT AW 102



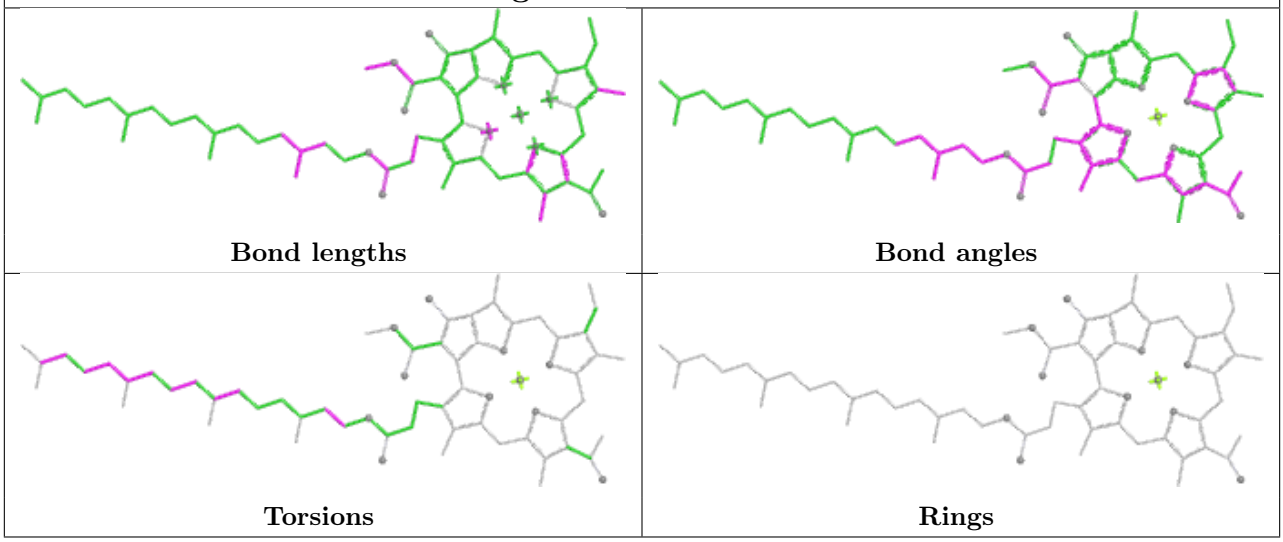
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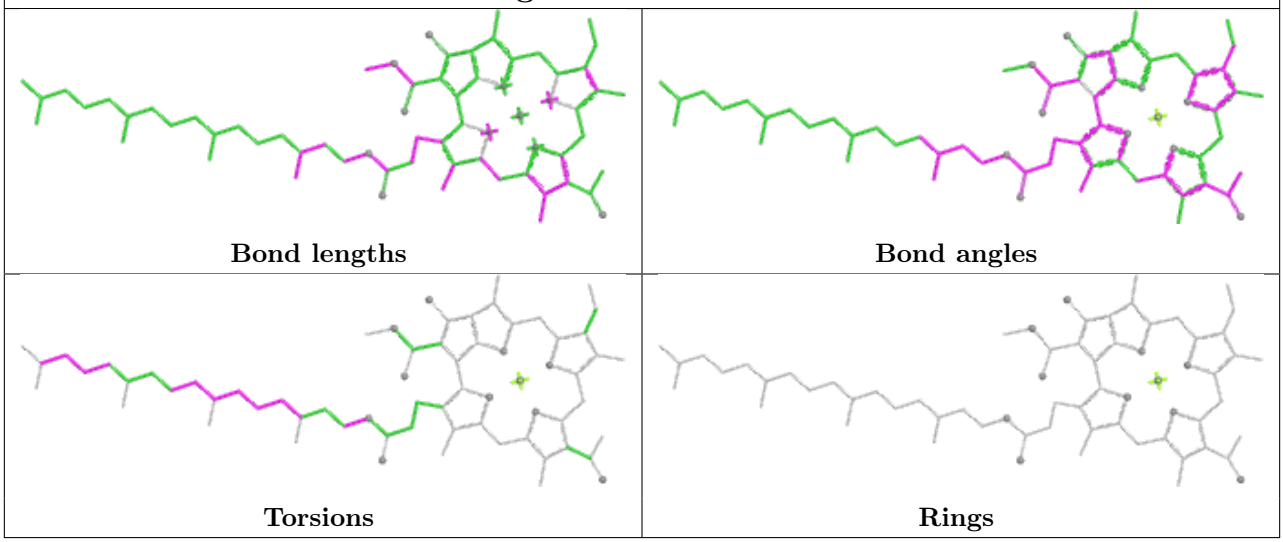
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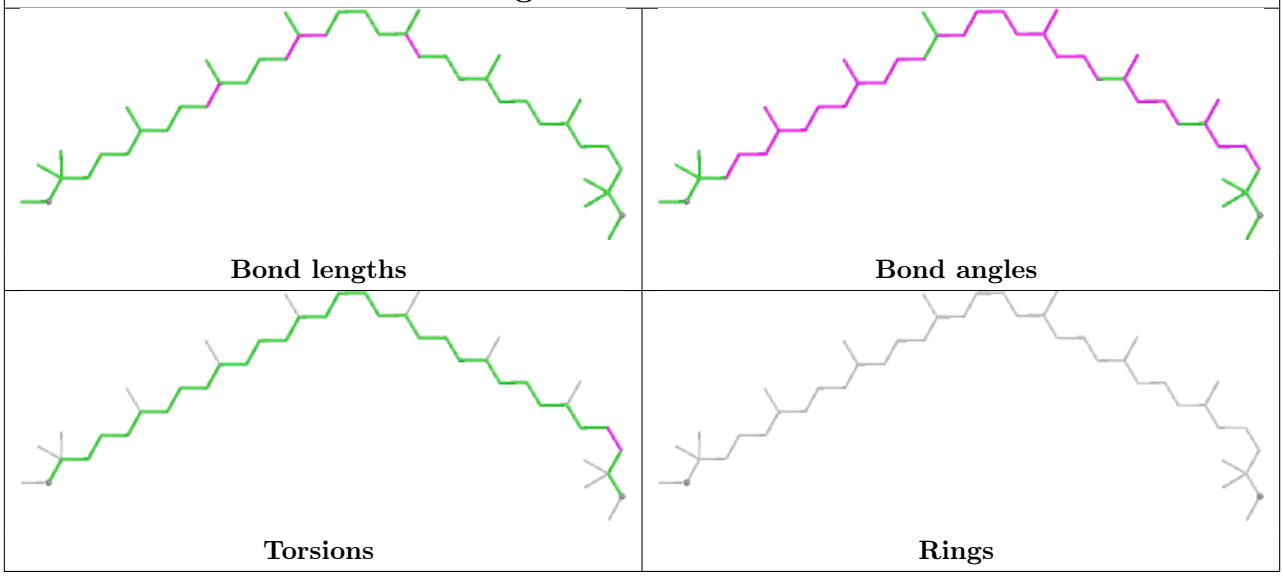
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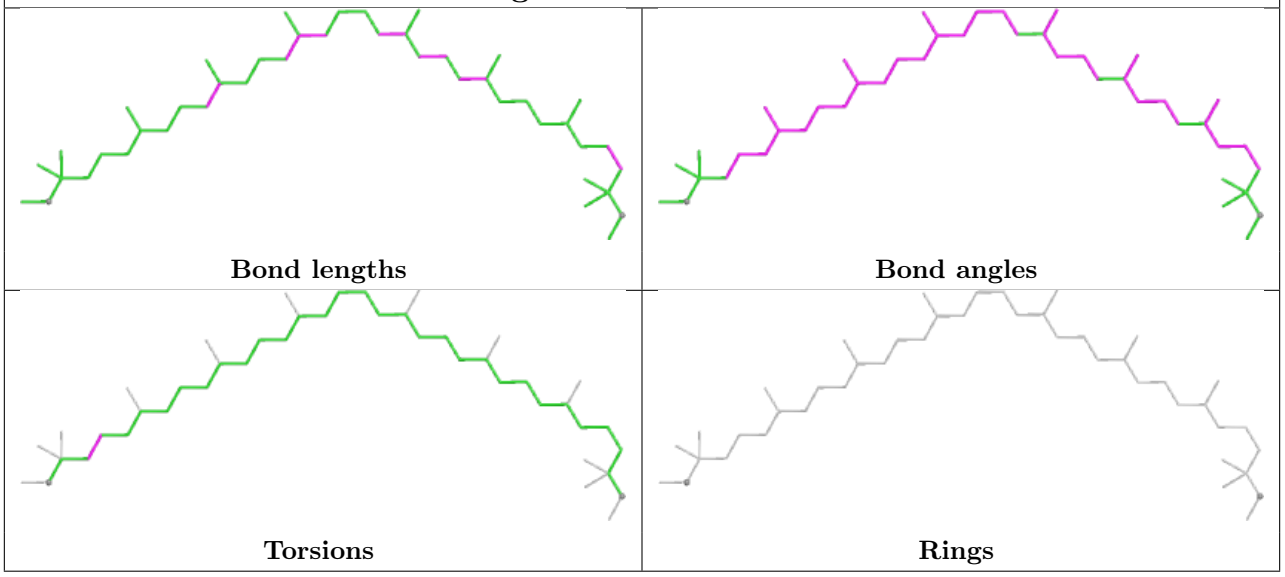
Ligand BCL A5 102



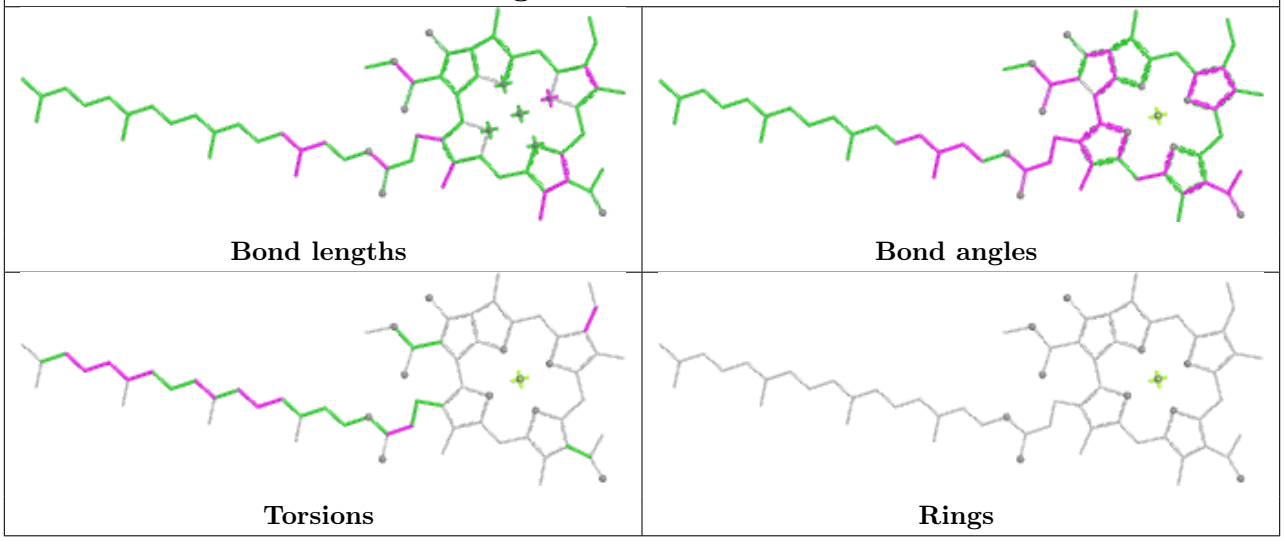
Ligand CRT AJ 102



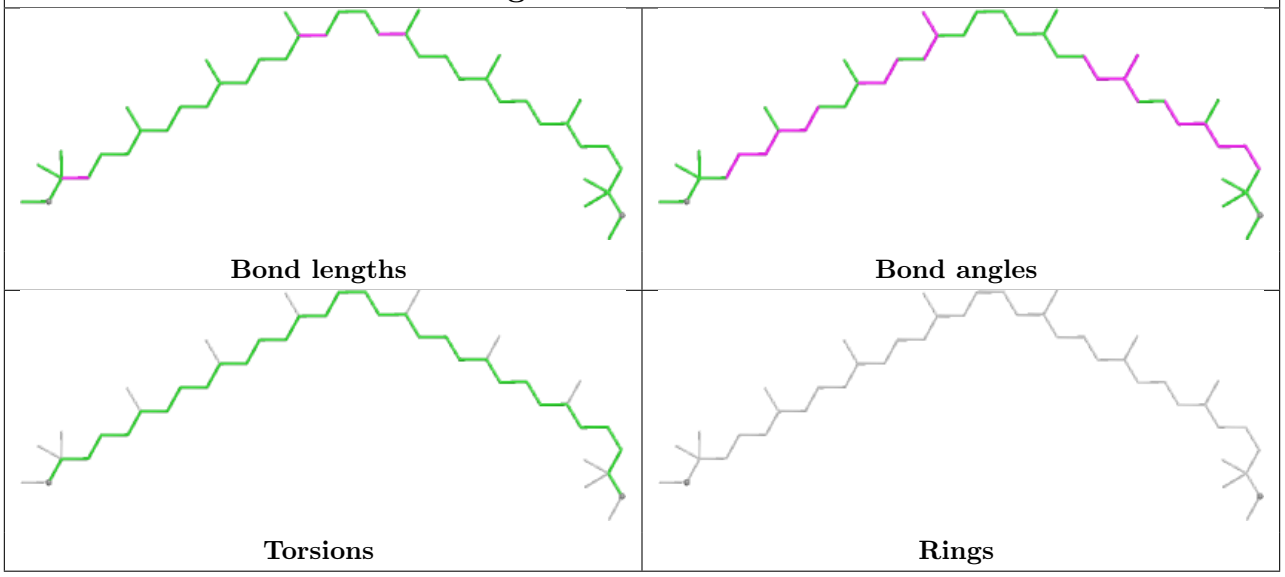
Ligand CRT BN 102



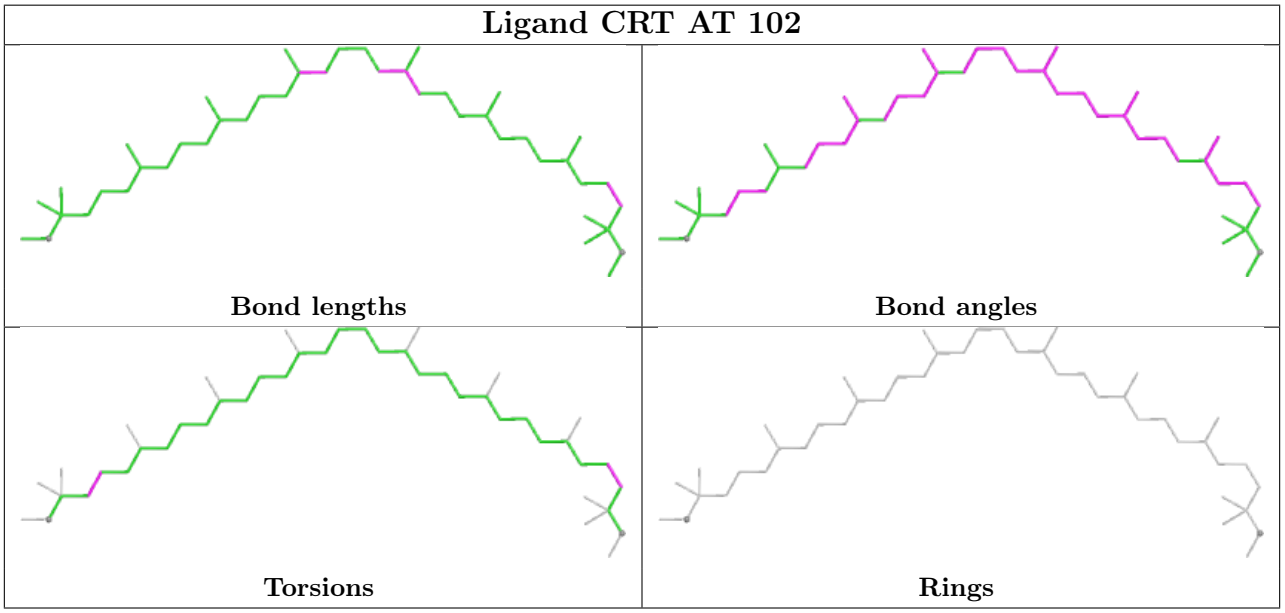
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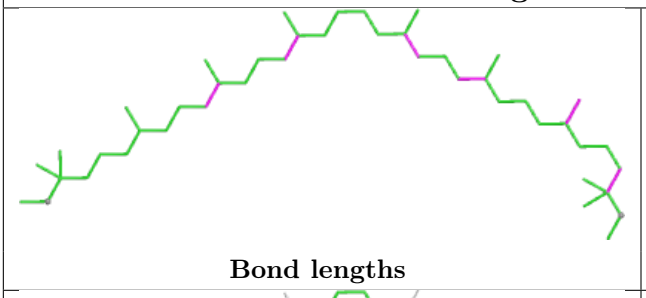
Ligand CRT AM 406



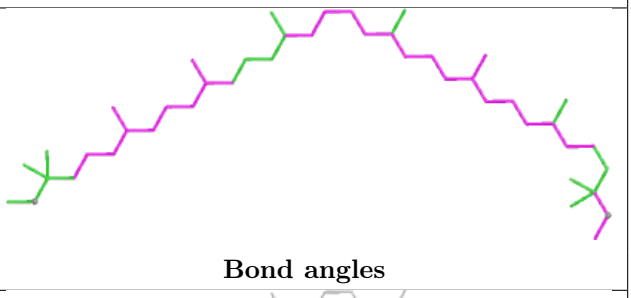
Ligand CRT AT 102



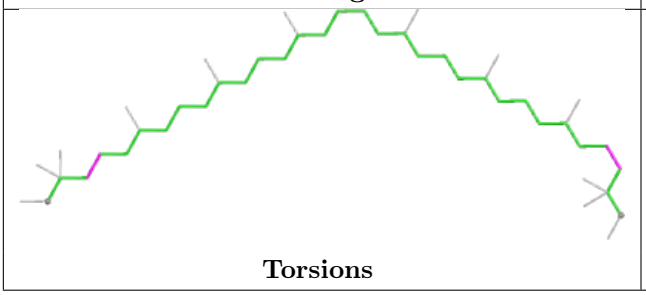
Ligand CRT BU 103



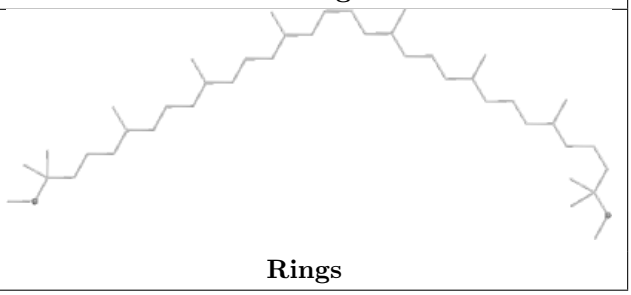
Bond lengths



Bond angles

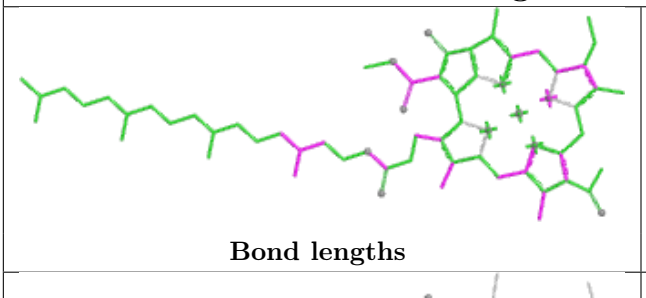


Torsions

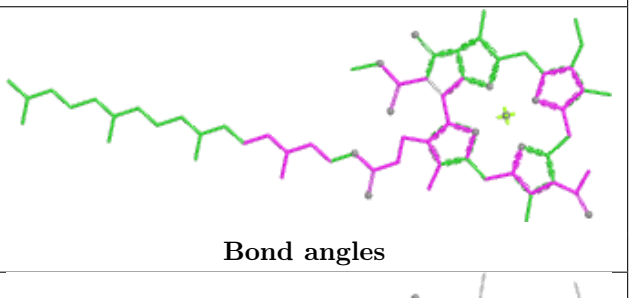


Rings

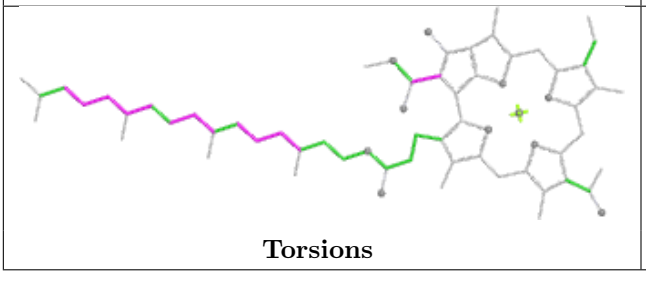
Ligand BCL BD 102



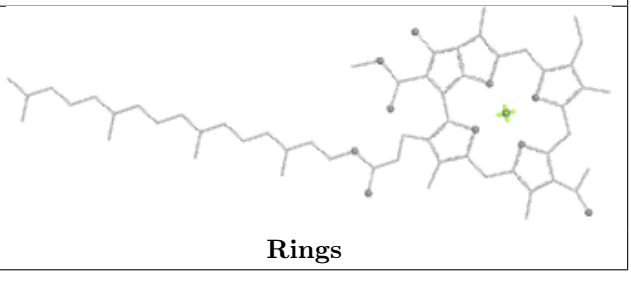
Bond lengths



Bond angles

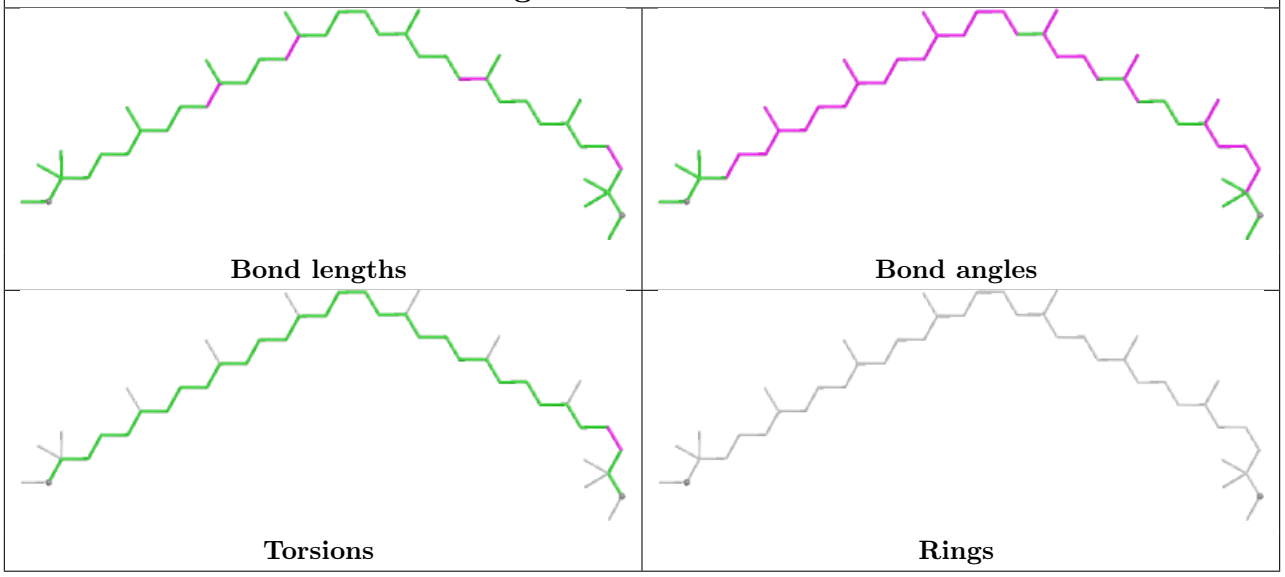


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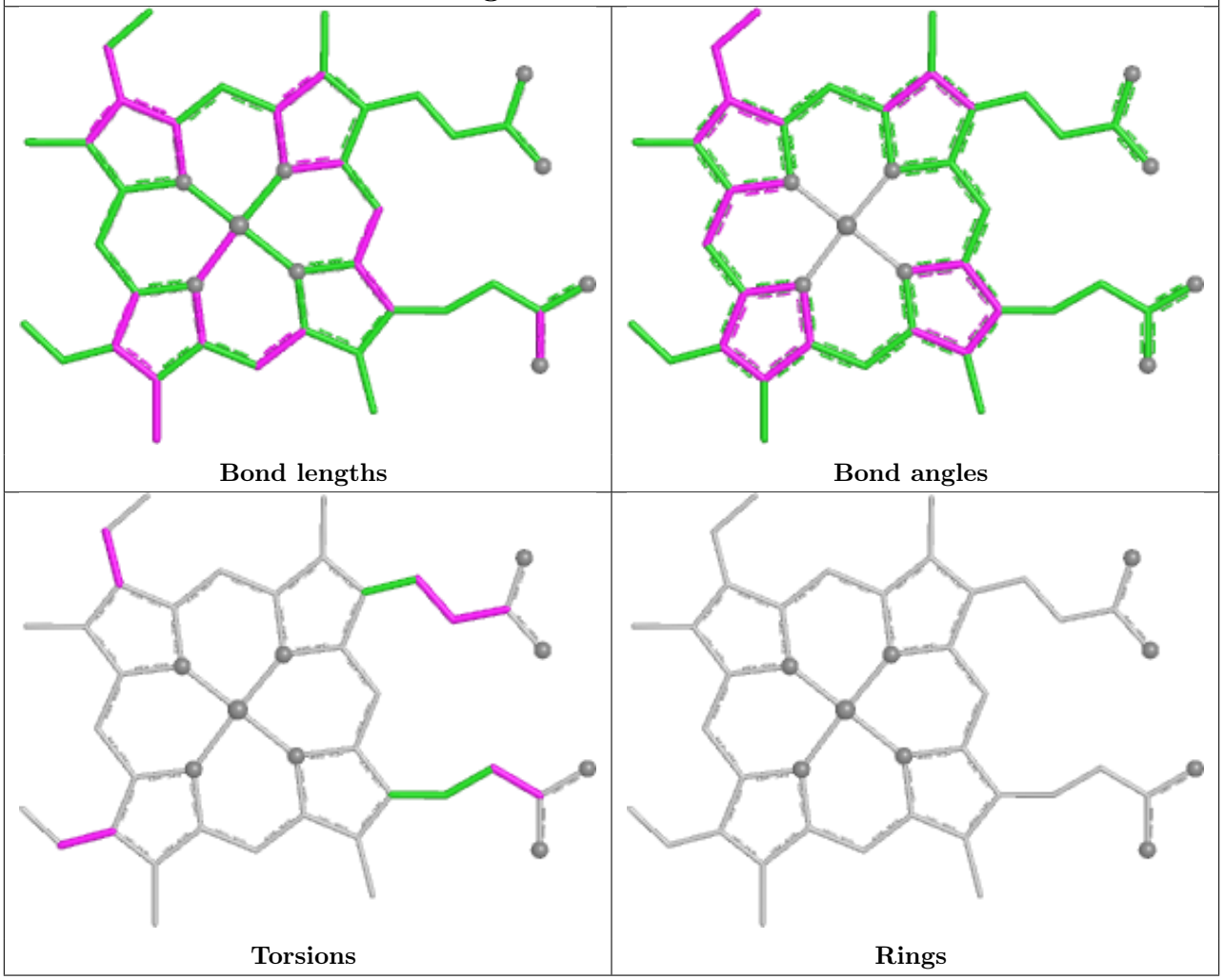


Rings

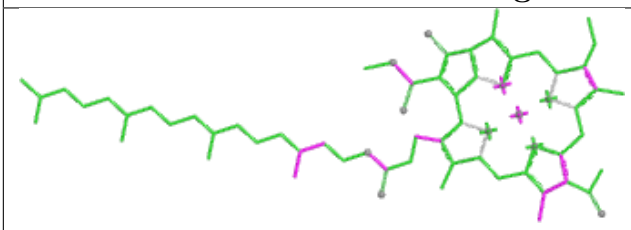
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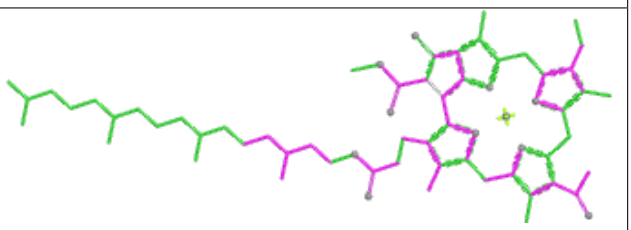
Ligand HEM AC 502



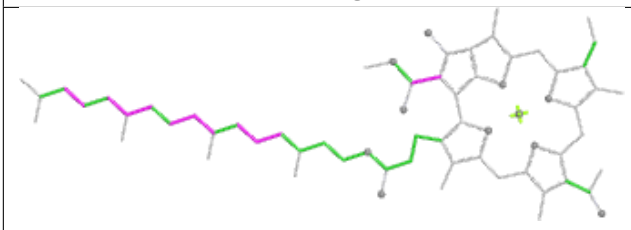
Ligand BCL AL 303



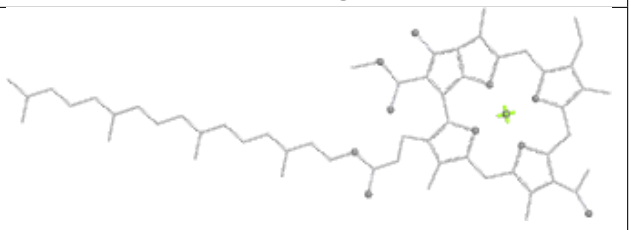
Bond lengths



Bond angles

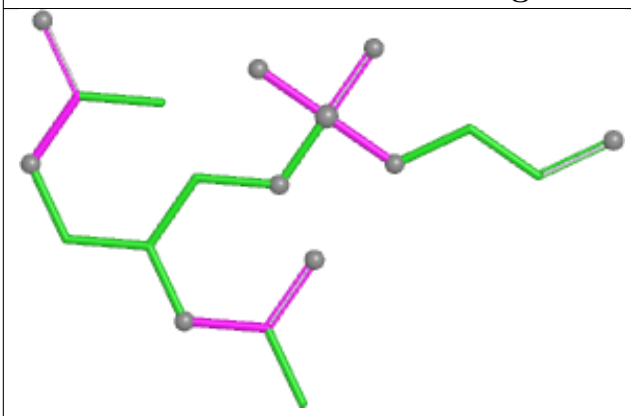


Torsions

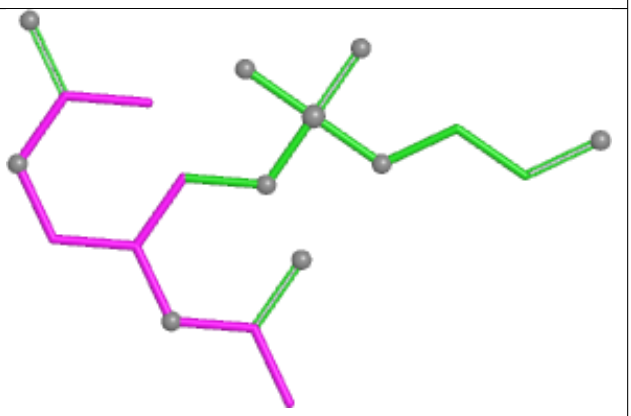


Rings

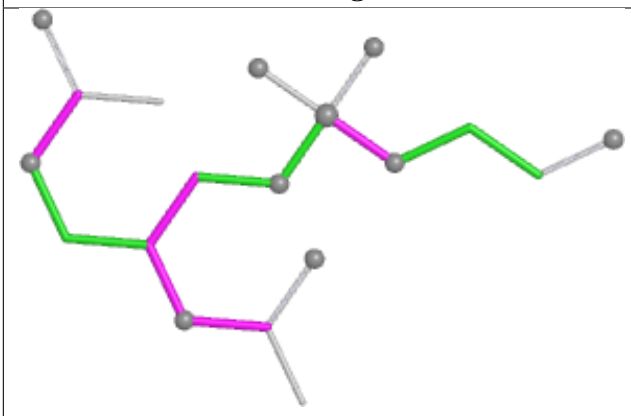
Ligand PEF AH 301



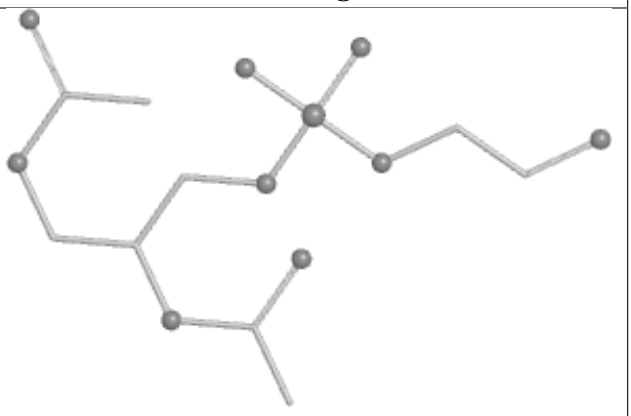
Bond lengths



Bond angles

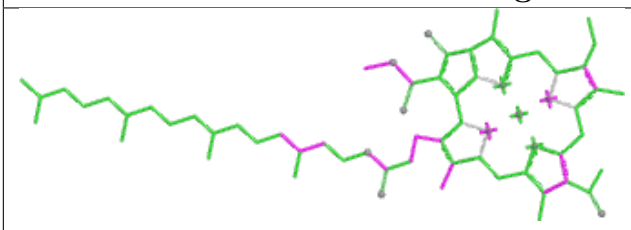


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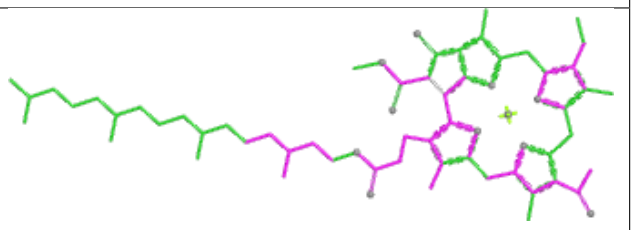


Rings

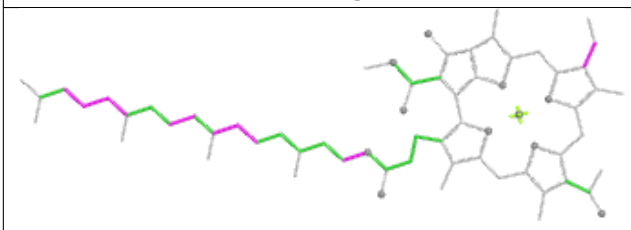
Ligand BCL BL 301



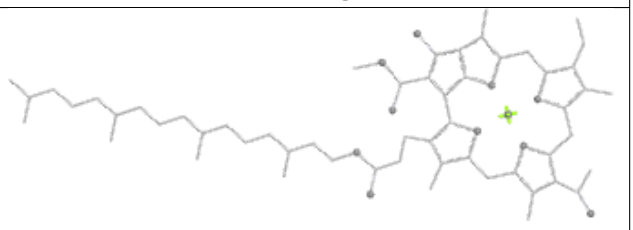
Bond lengths



Bond angles

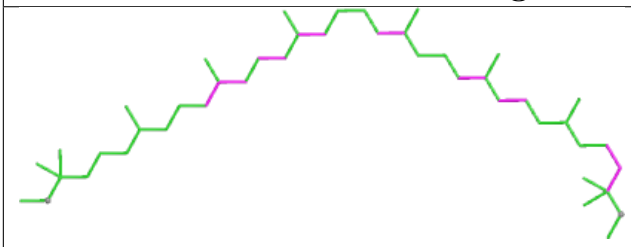


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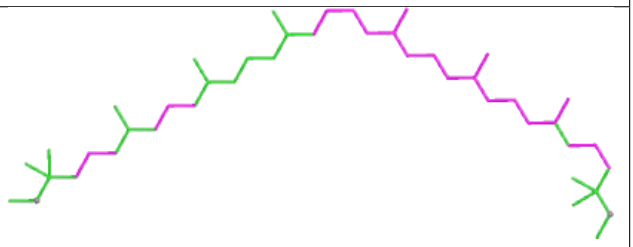


Rings

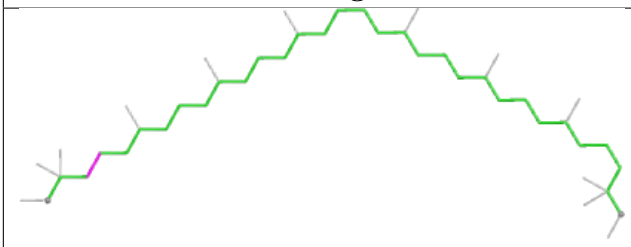
Ligand CRT AX 102



Bond lengths



Bond angles

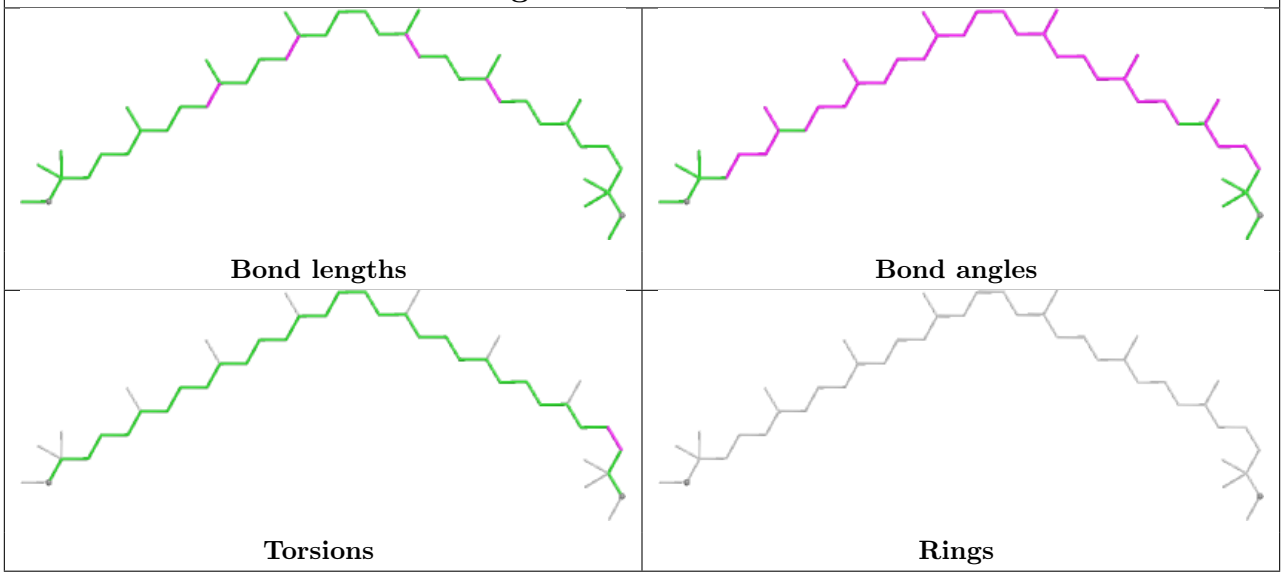


Torsions

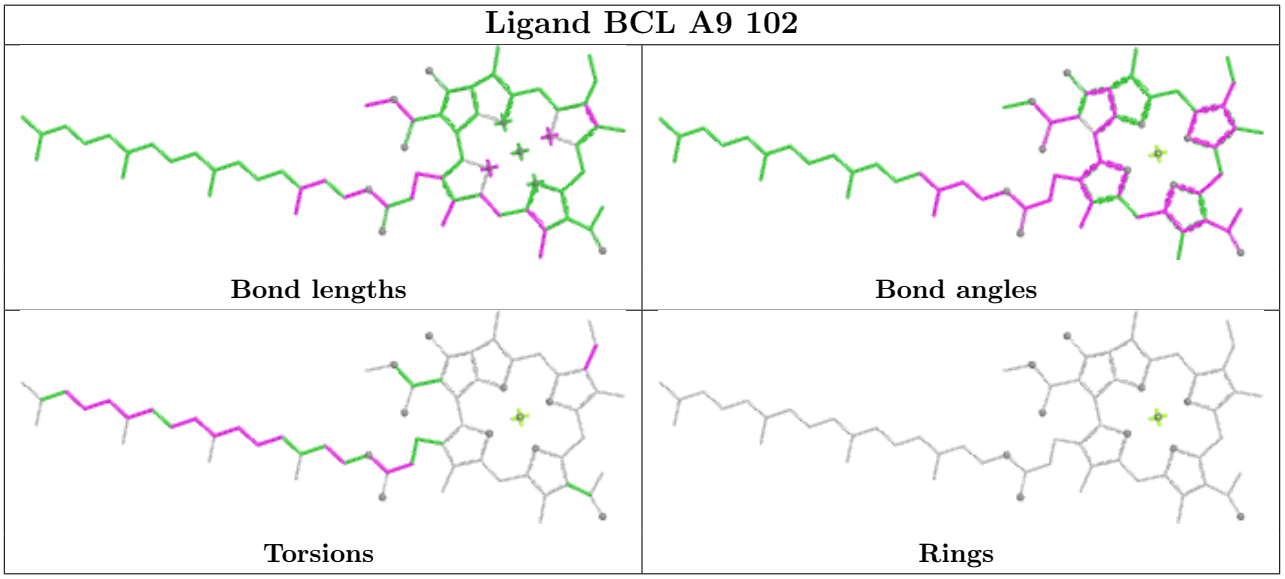


Rings

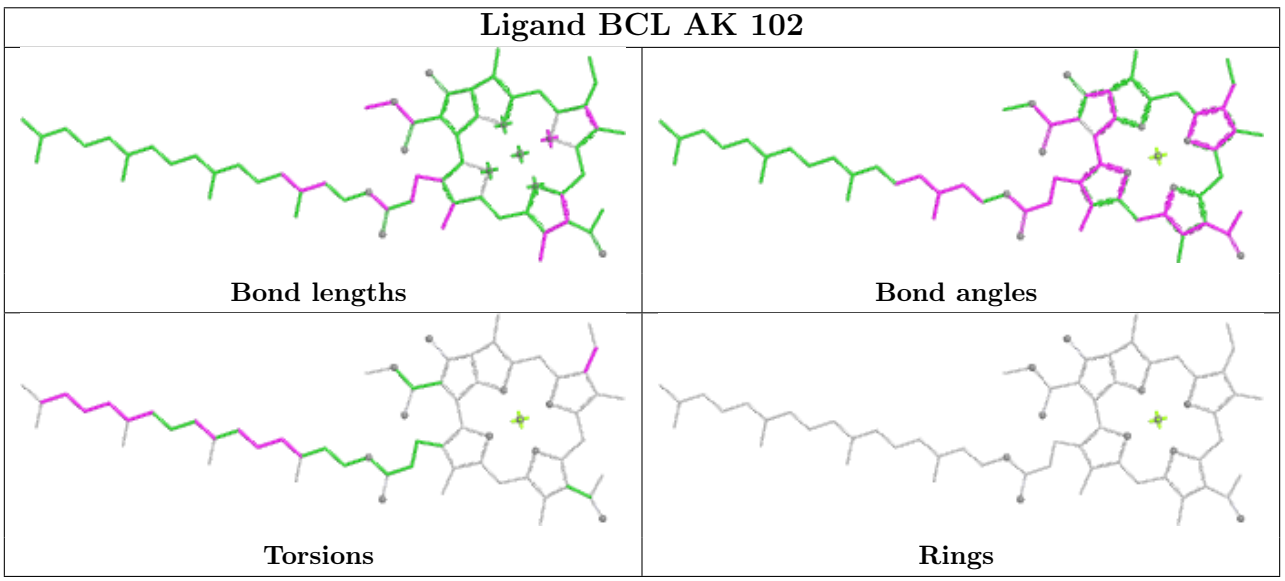
Ligand CRT AG 102



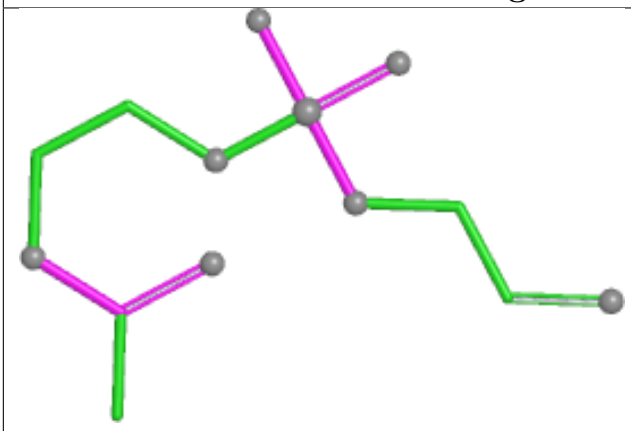
Ligand BCL A9 102



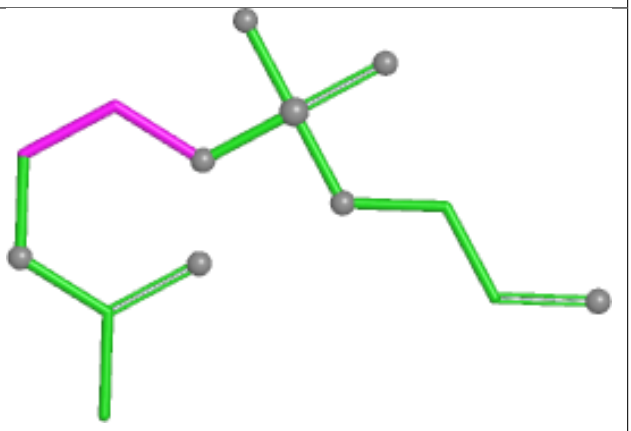
Ligand BCL AK 102



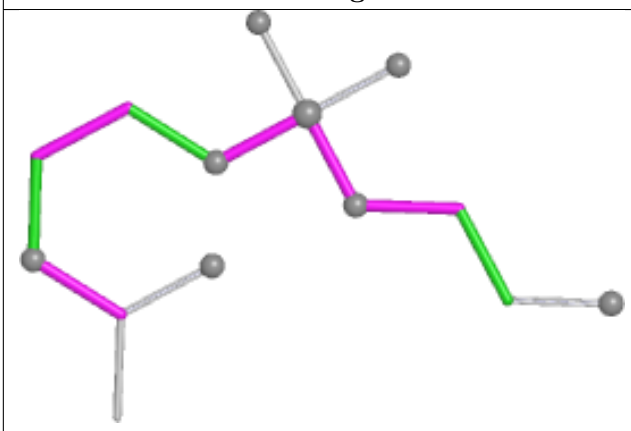
Ligand PEF AM 408



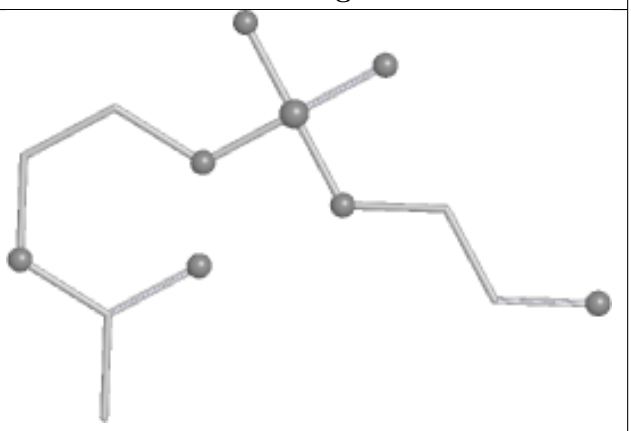
Bond lengths



Bond angles

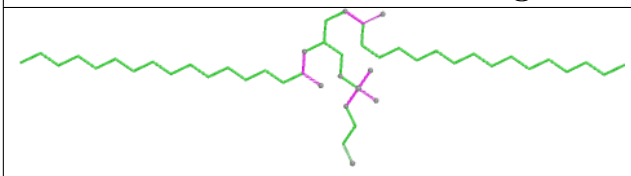


Torsions

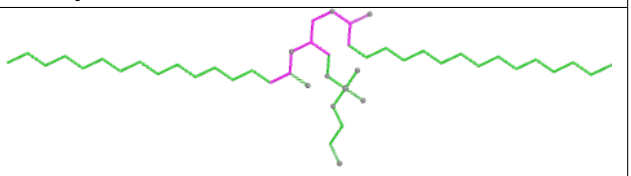


Rings

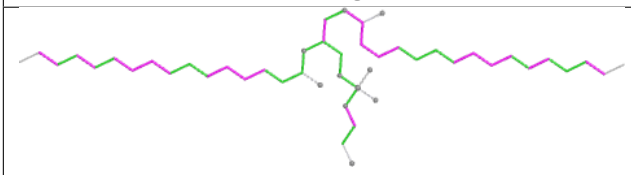
Ligand PEF BQ 101



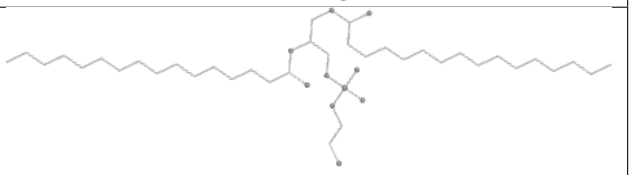
Bond lengths



Bond angles

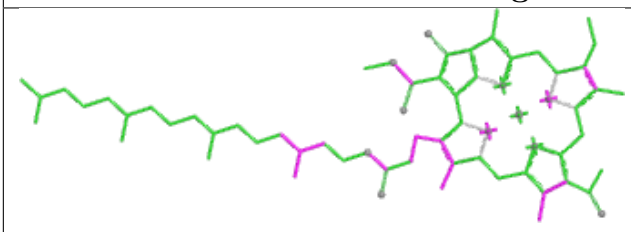


Torsions

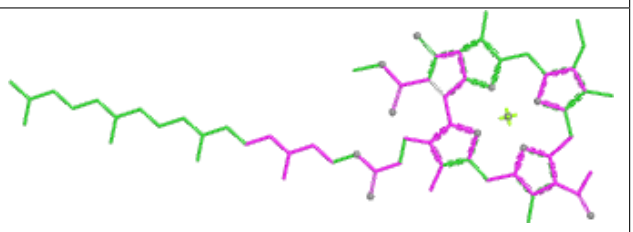


Rings

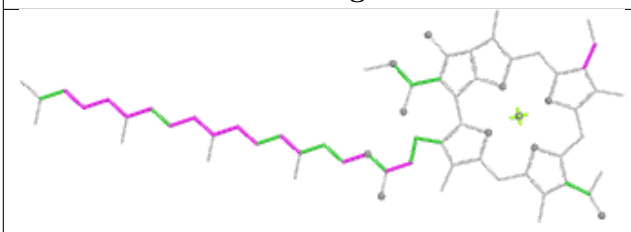
Ligand BCL AO 102



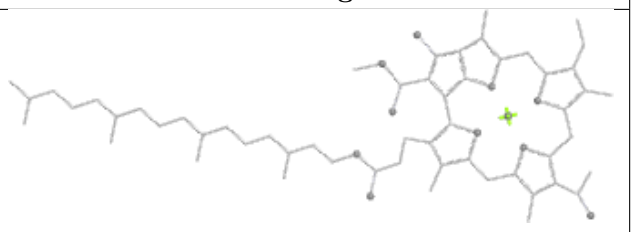
Bond lengths



Bond angles

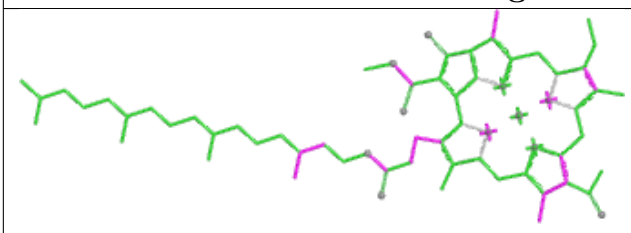


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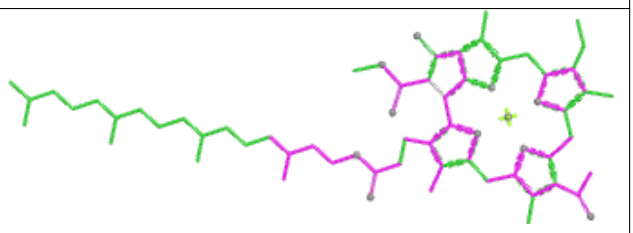


Rings

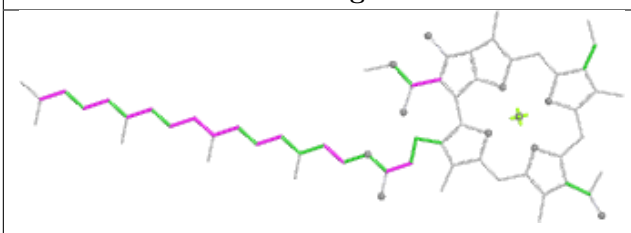
Ligand BCL BM 402



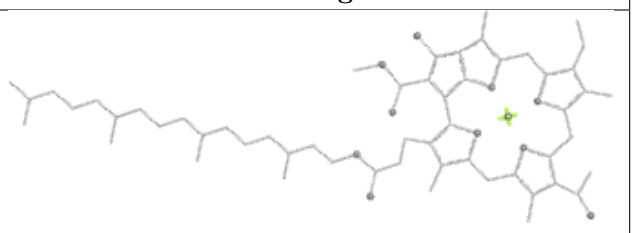
Bond lengths



Bond angles

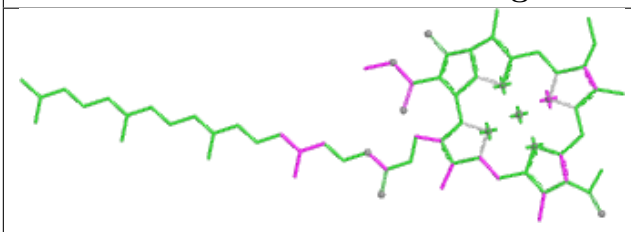


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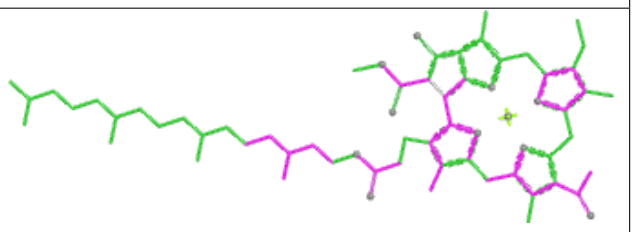


Rings

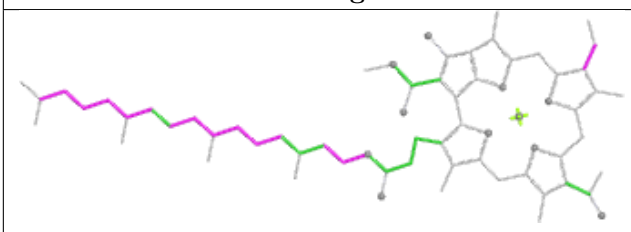
Ligand BCL AA 101



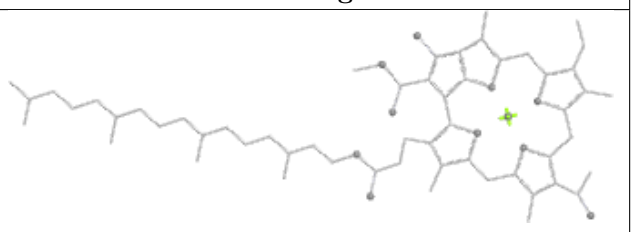
Bond lengths



Bond angles

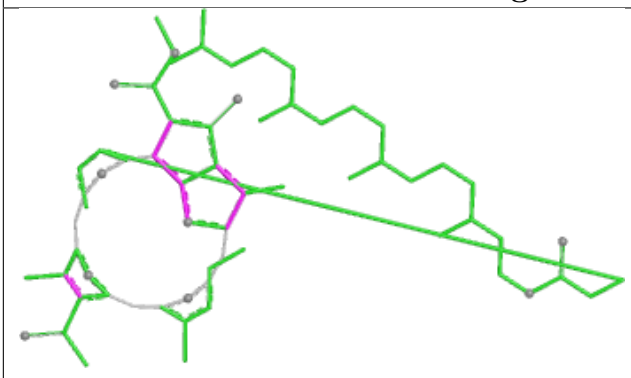


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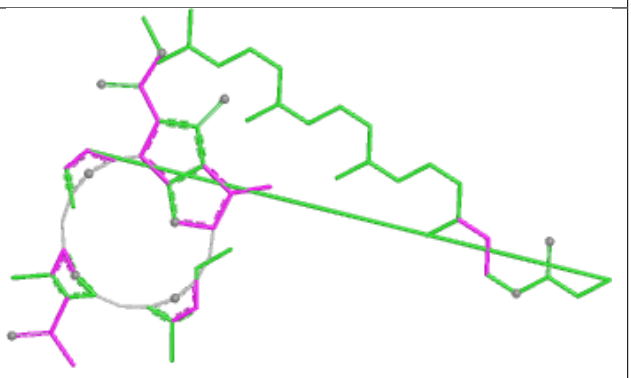


Rings

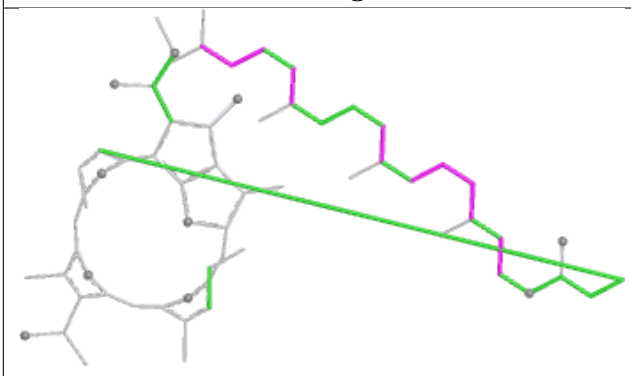
Ligand BPH AL 302



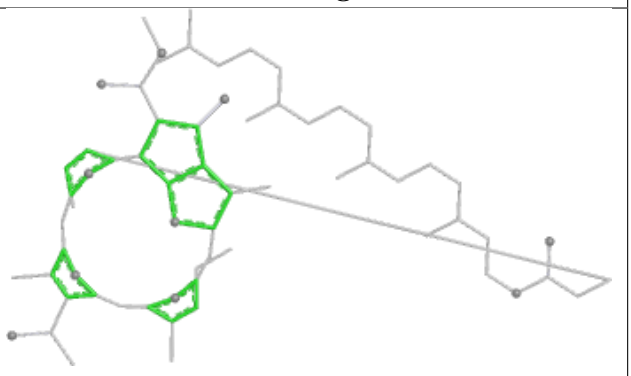
Bond lengths



Bond angles

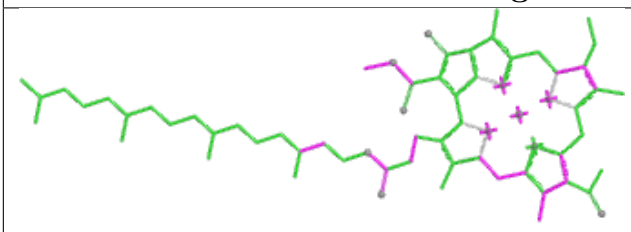


Torsions

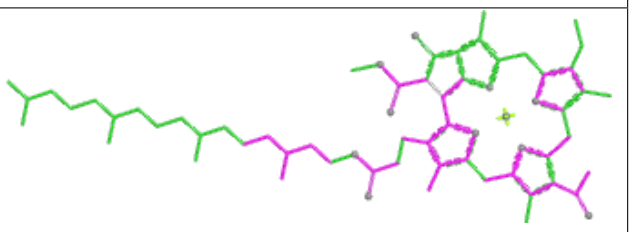


Rings

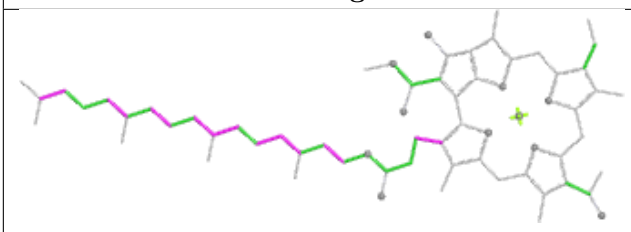
Ligand BCL AR 101



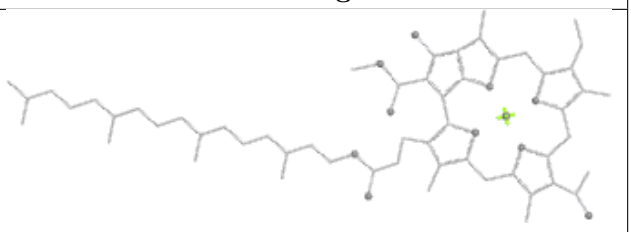
Bond lengths



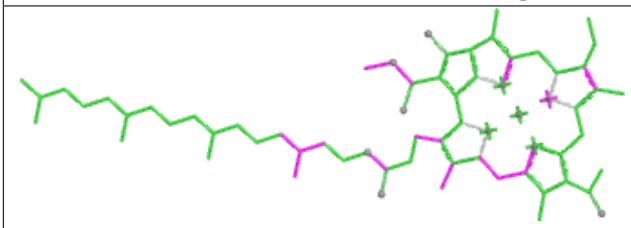
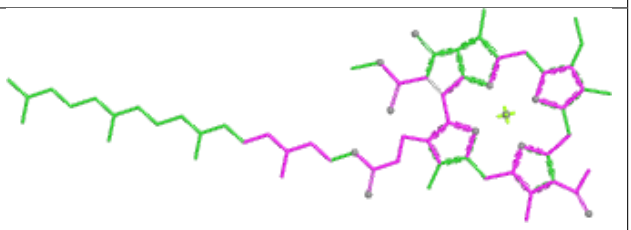
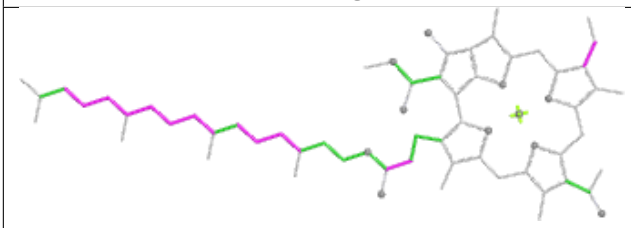
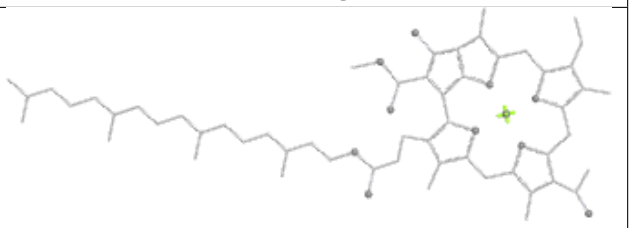
Bond angles

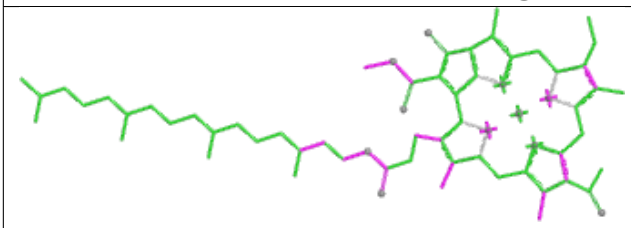
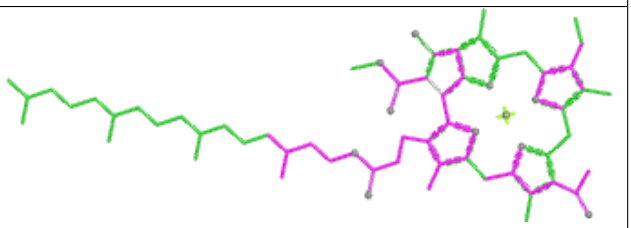
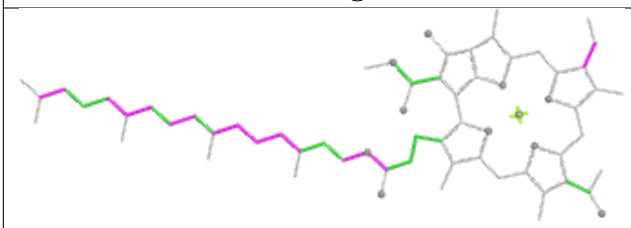
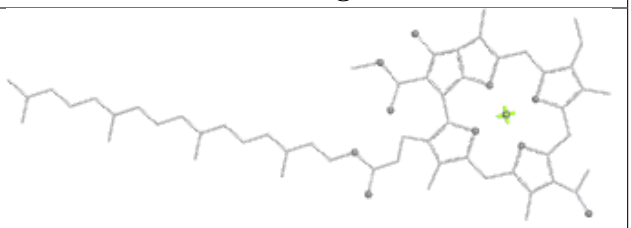


Torsions

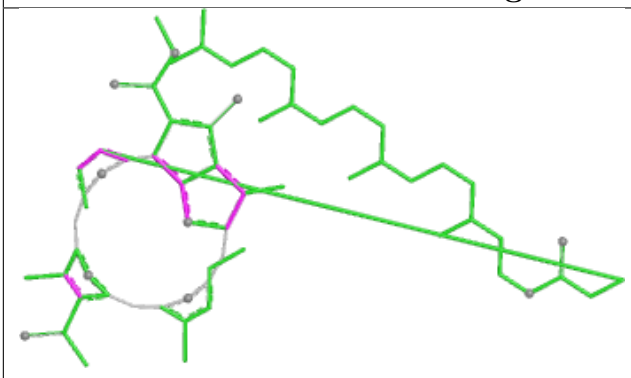


Rings

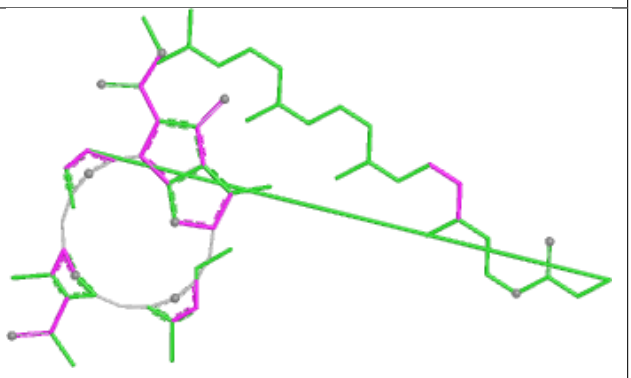
Ligand BCL AS 103	
 Bond lengths	 Bond angles
 Torsions	 Rings

Ligand BCL A6 101	
 Bond lengths	 Bond angles
 Torsions	 Rings

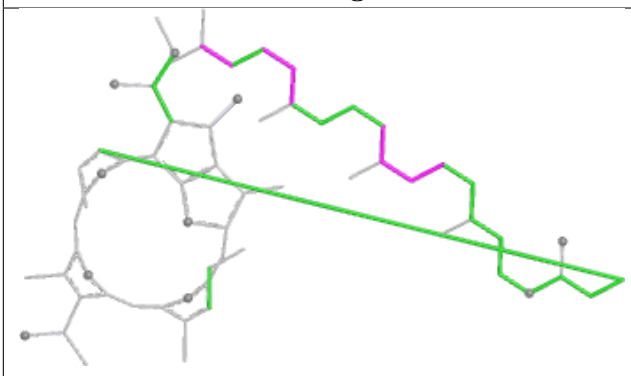
Ligand BPH AM 403



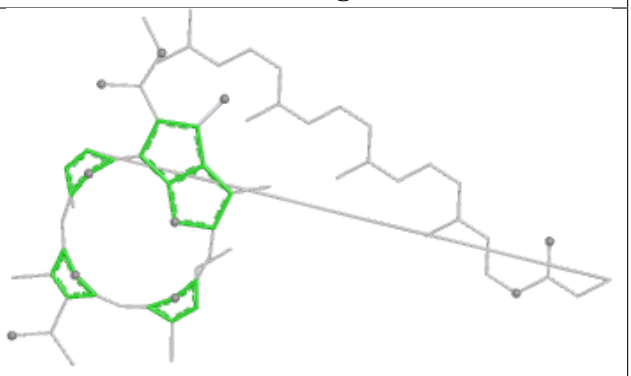
Bond lengths



Bond angles

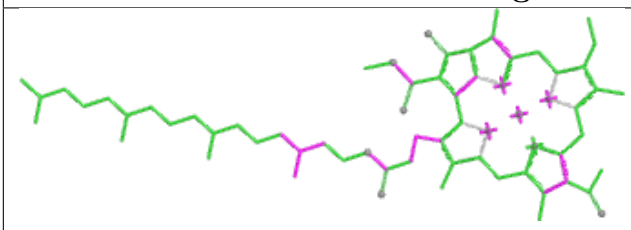


Torsions

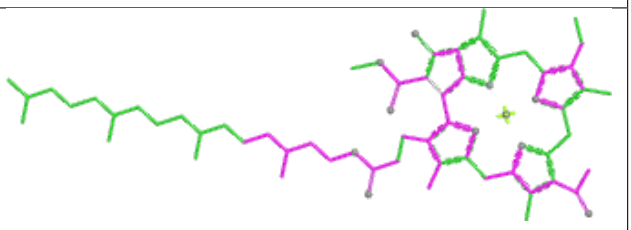


Rings

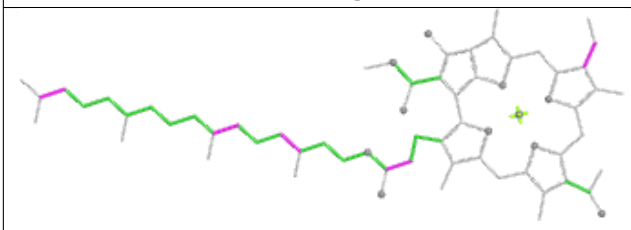
Ligand BCL BB 101



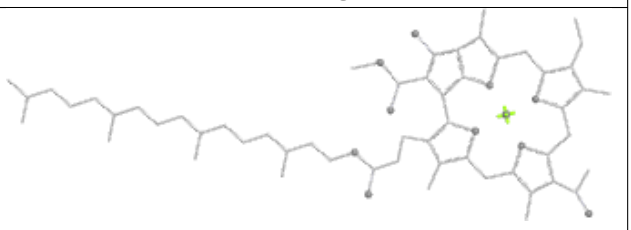
Bond lengths



Bond angles

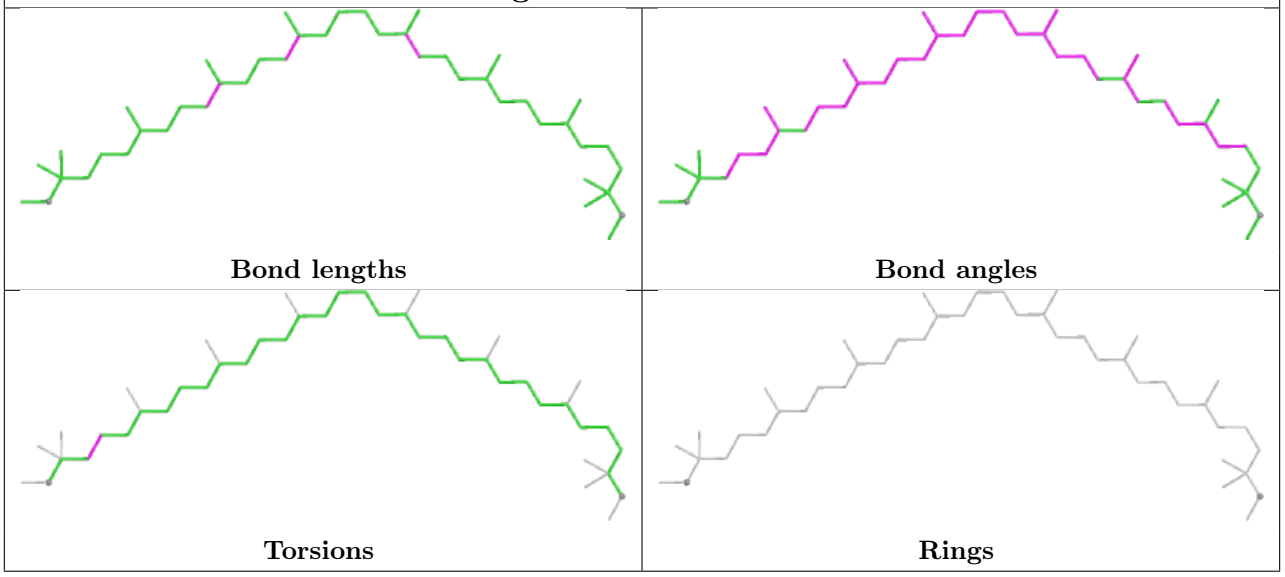


Torsions

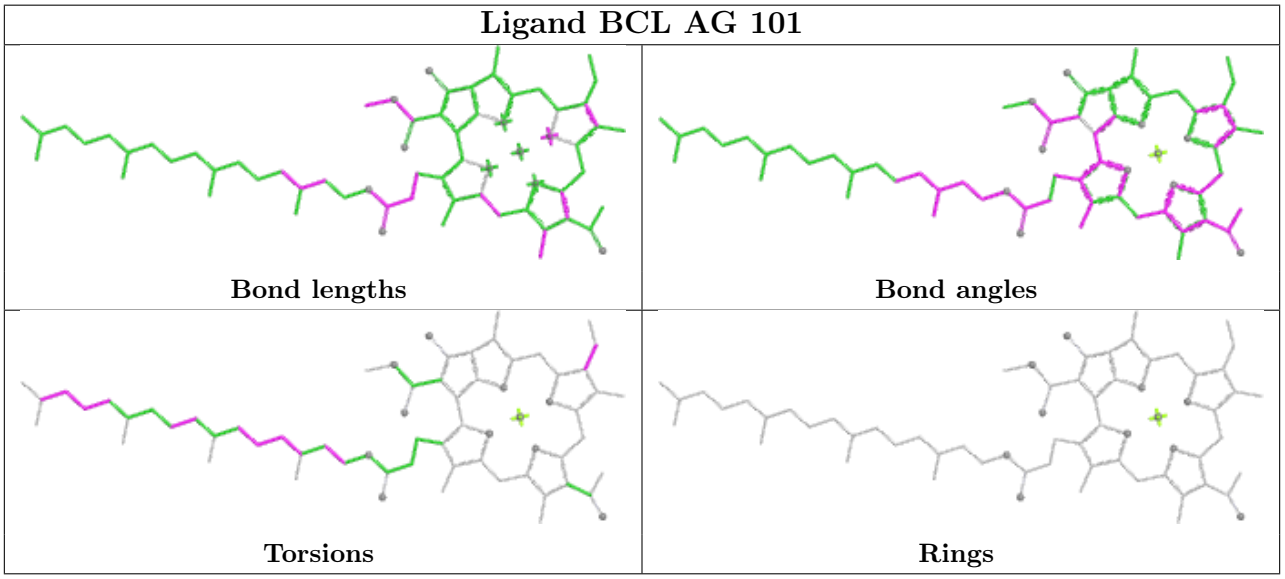


Rings

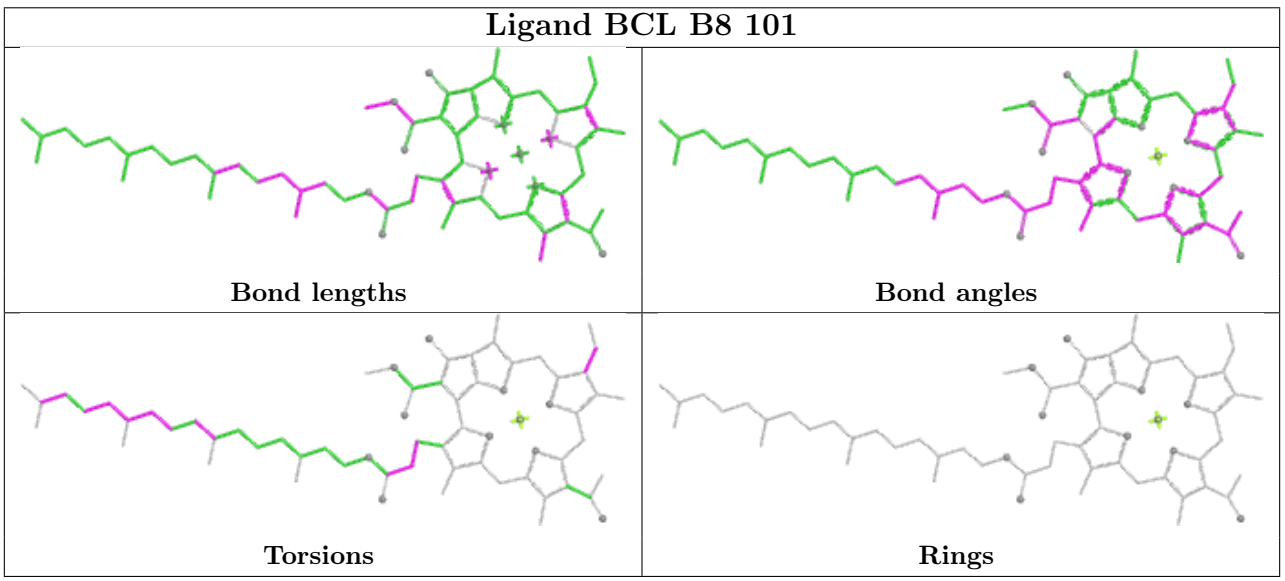
Ligand CRT BO 103



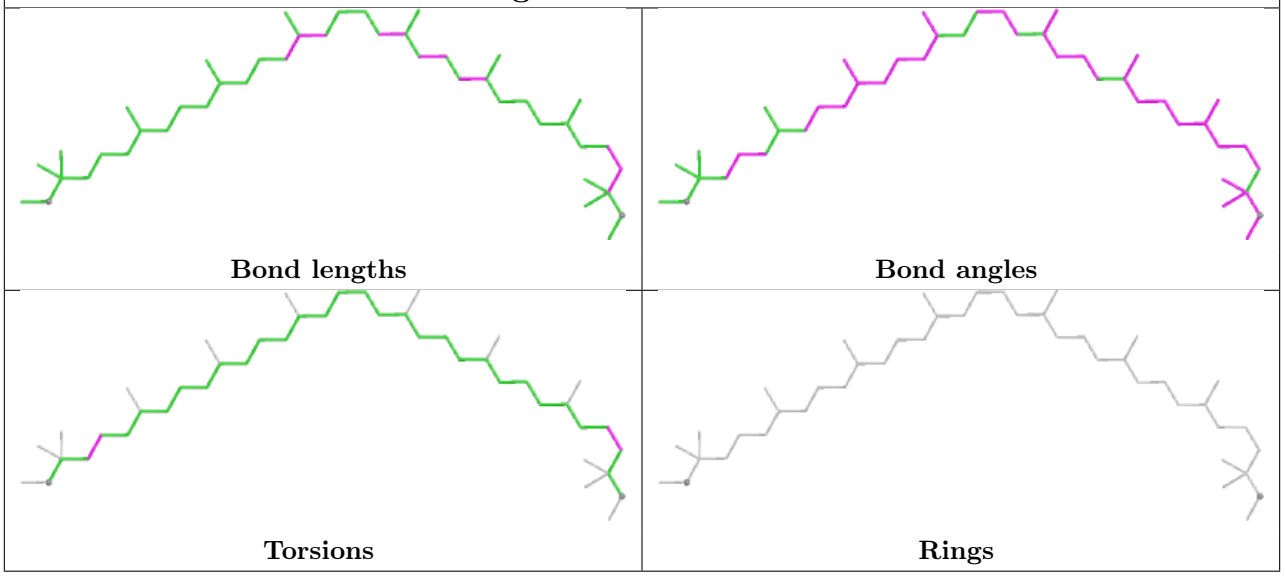
Ligand BCL AG 101



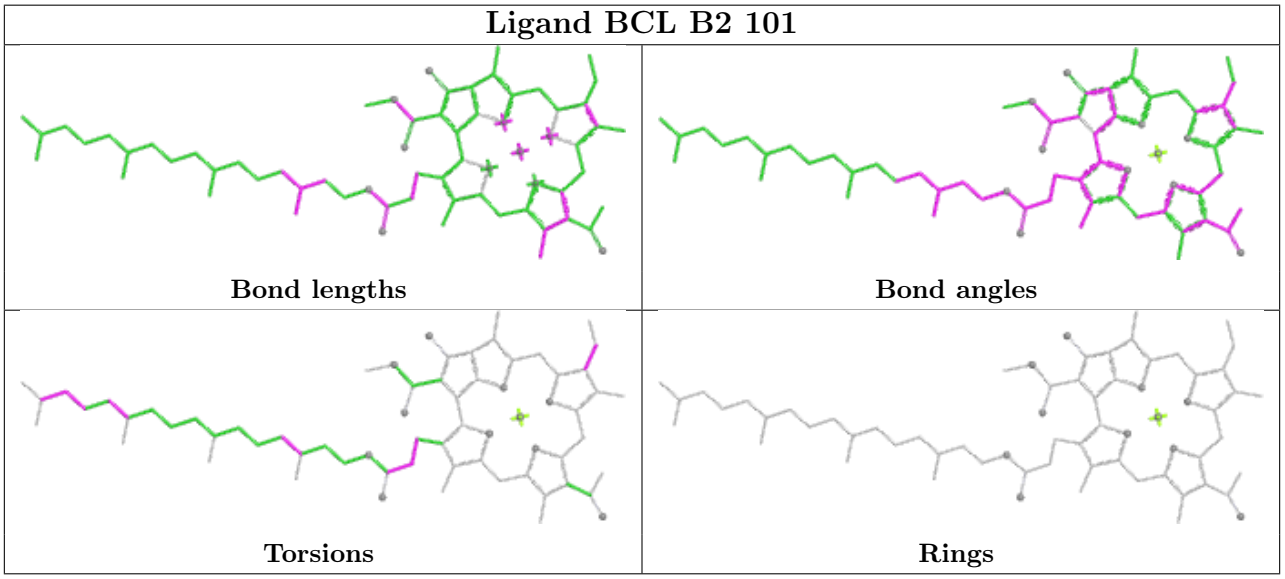
Ligand BCL B8 101



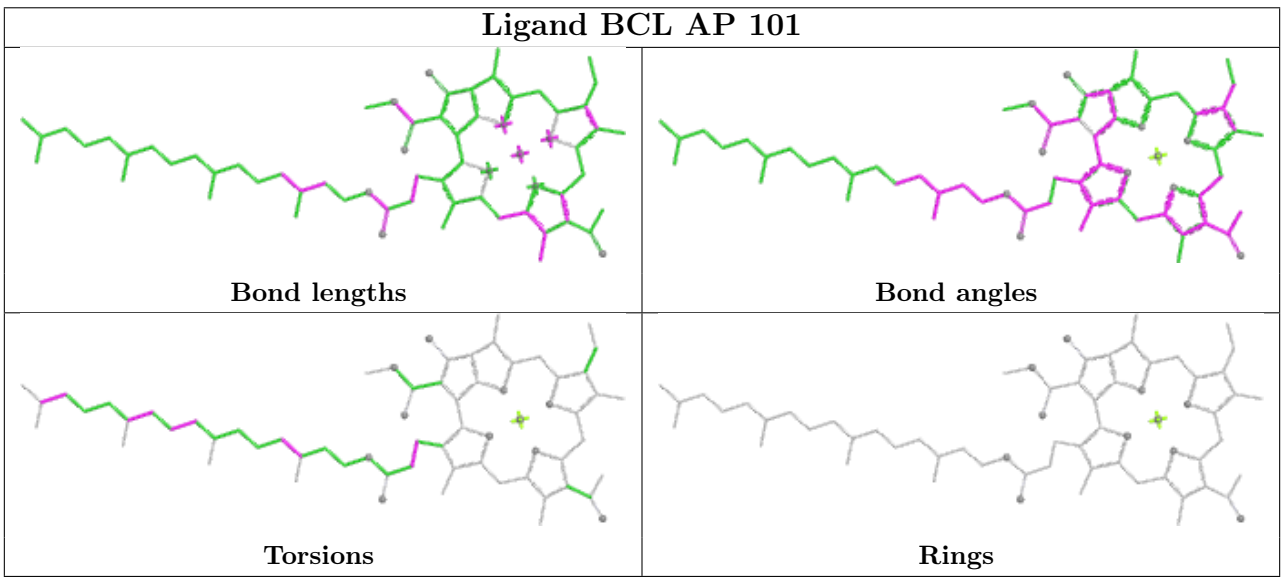
Ligand CRT AP 102



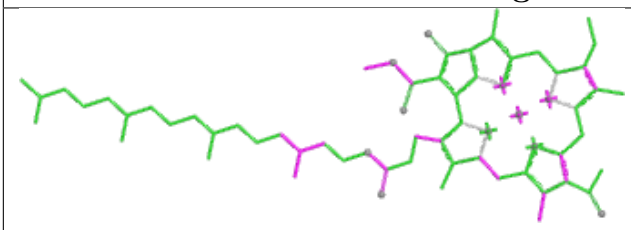
Ligand BCL B2 101



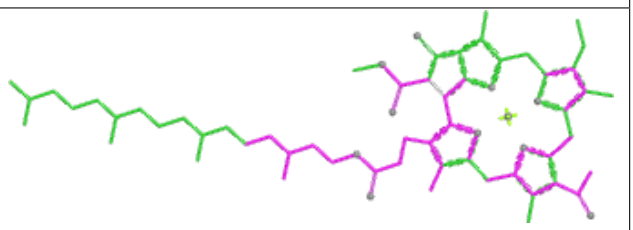
Ligand BCL AP 101



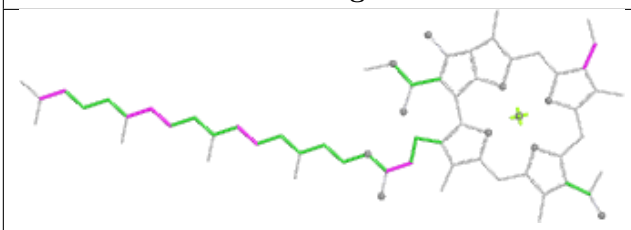
Ligand BCL AE 101



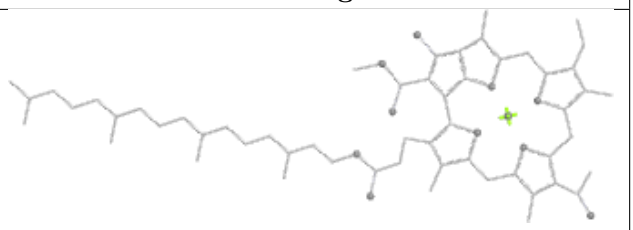
Bond lengths



Bond angles

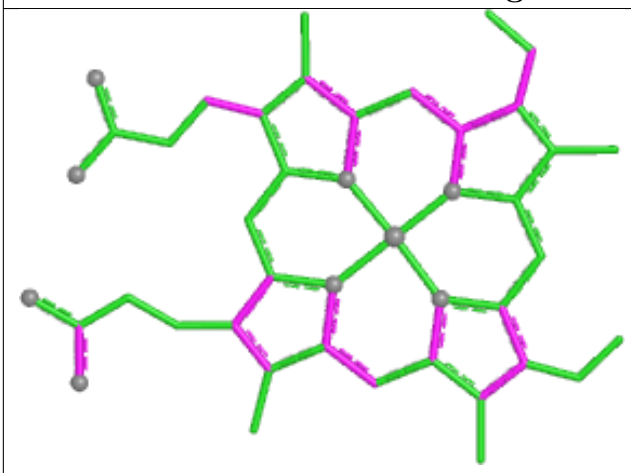


Torsions

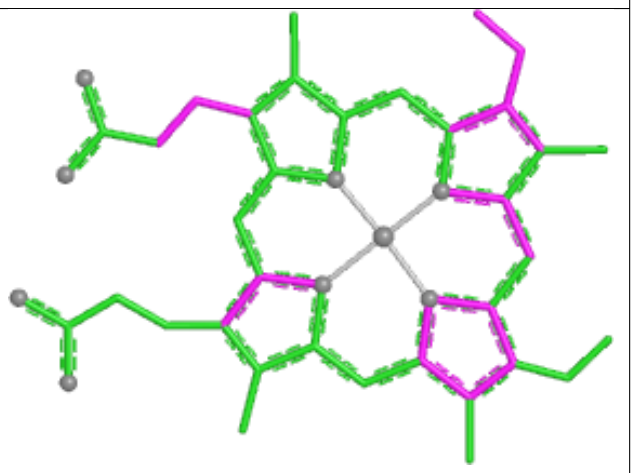


Rings

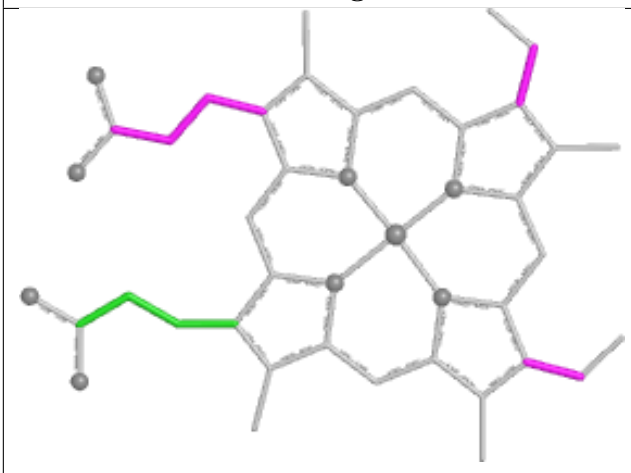
Ligand HEM BC 504



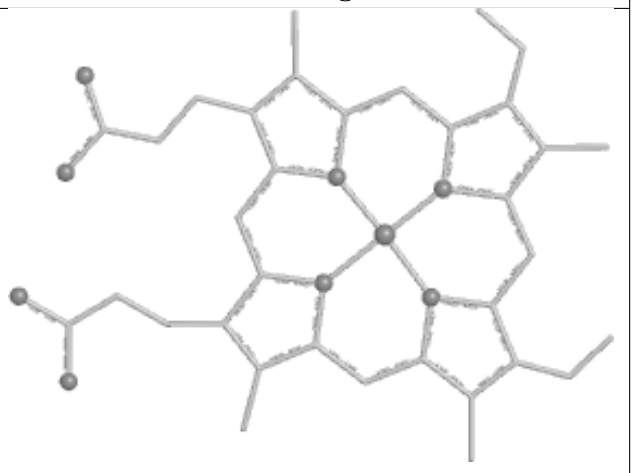
Bond lengths



Bond angles

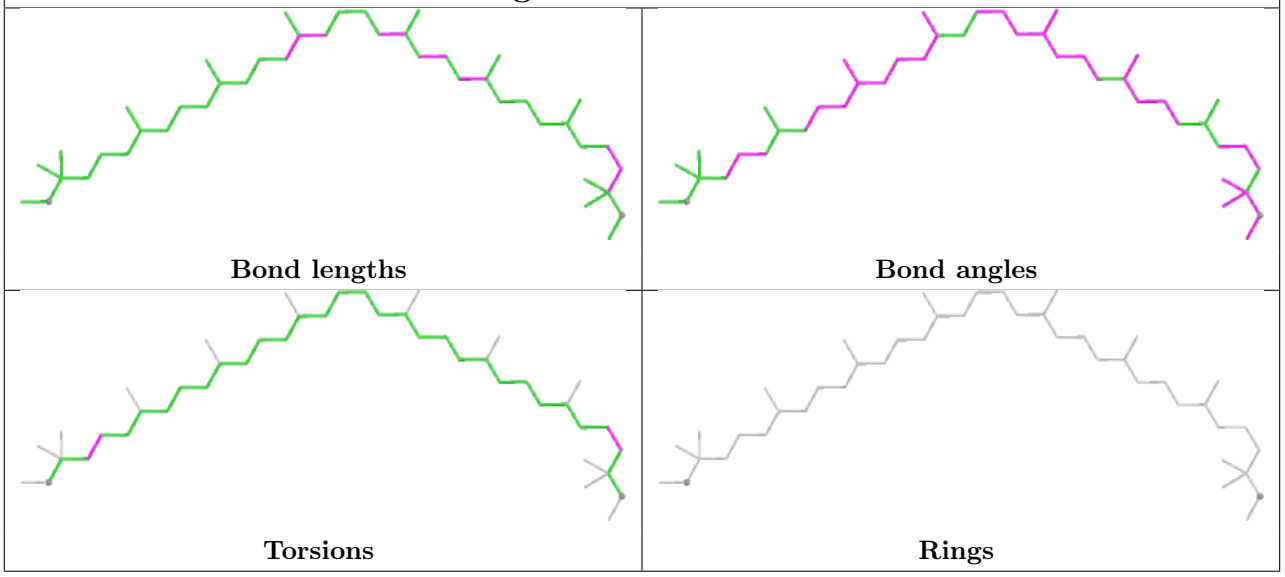


Torsions

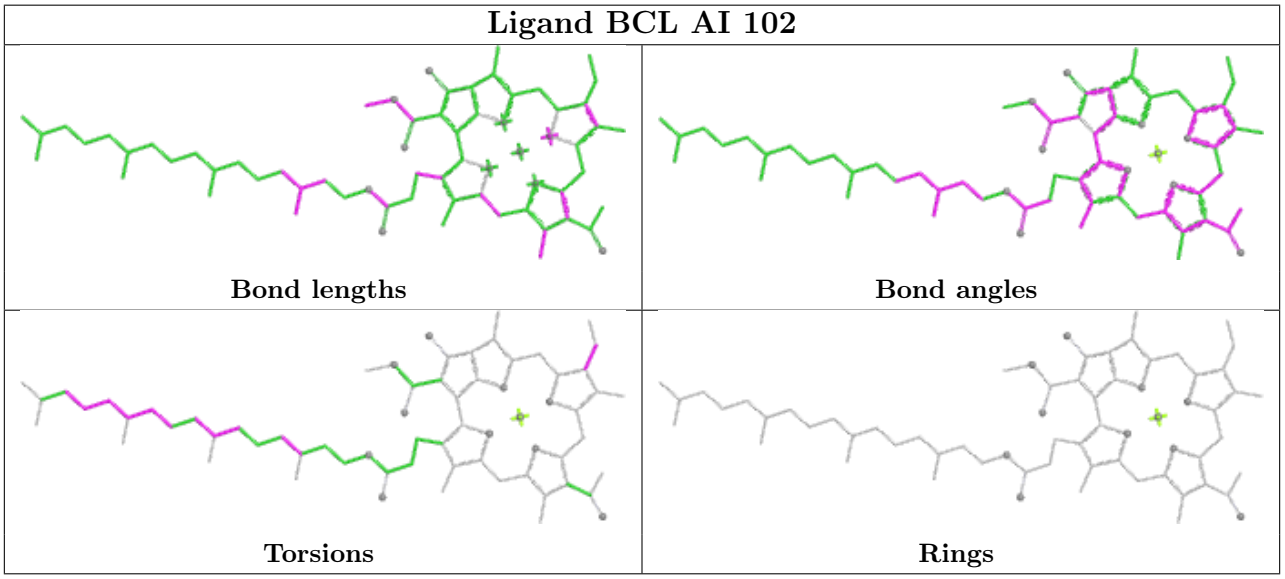


Rings

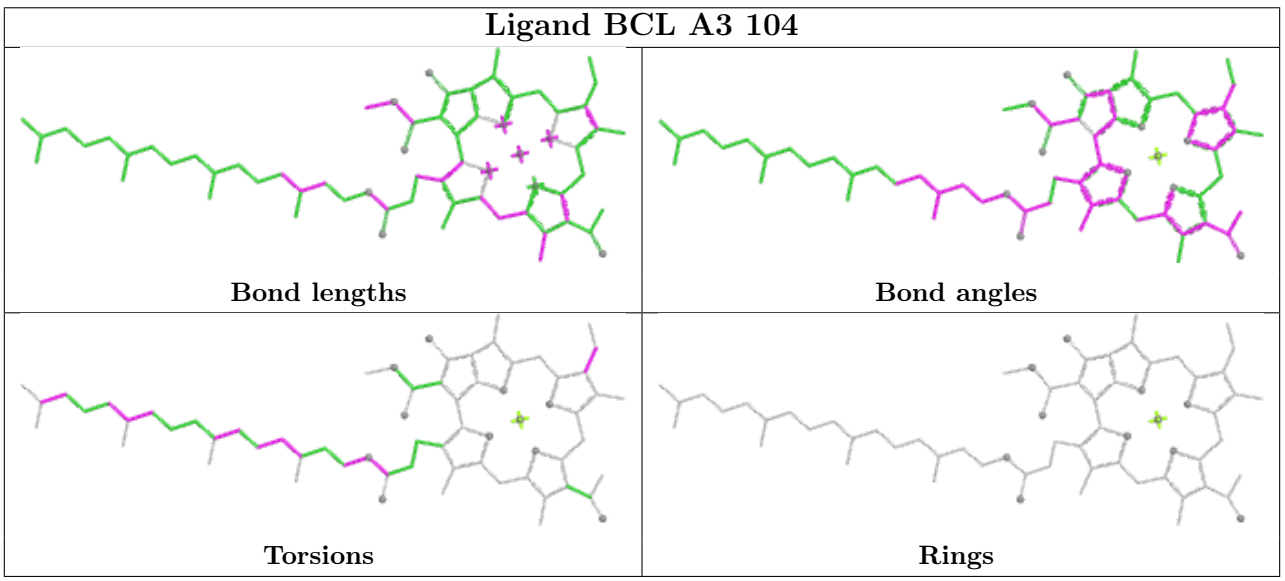
Ligand CRT BP 102



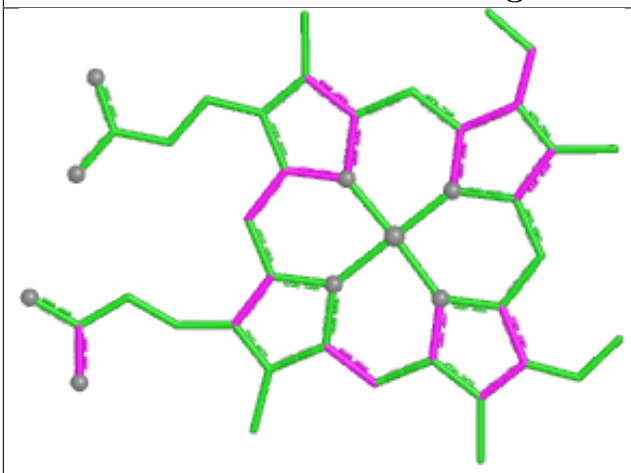
Ligand BCL AI 102



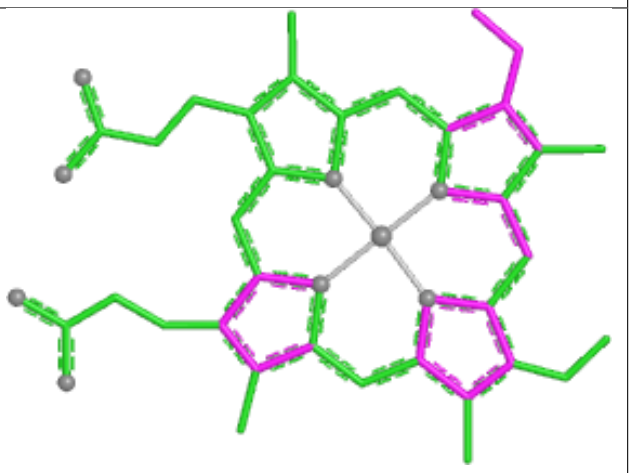
Ligand BCL A3 104



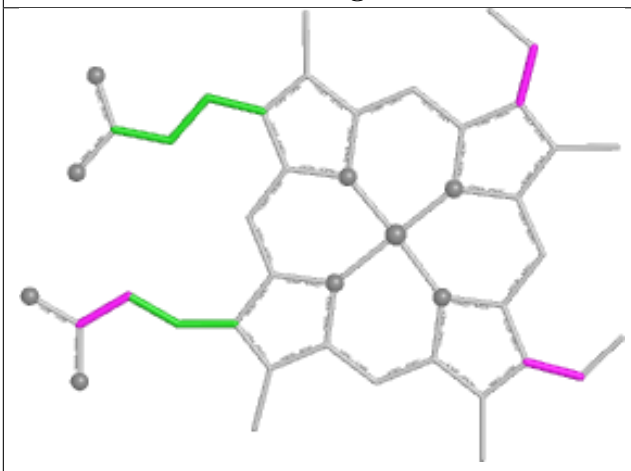
Ligand HEM BC 501



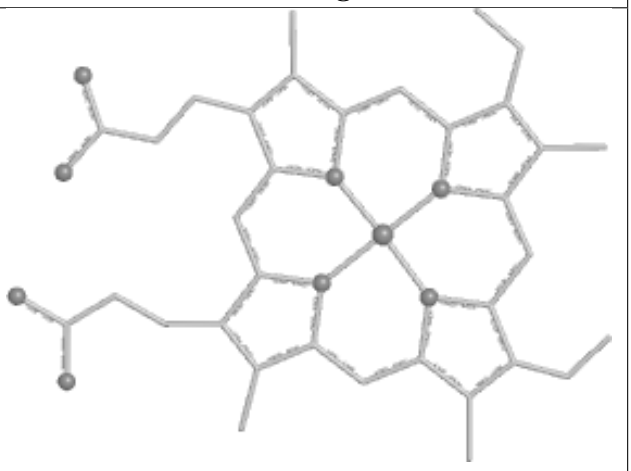
Bond lengths



Bond angles

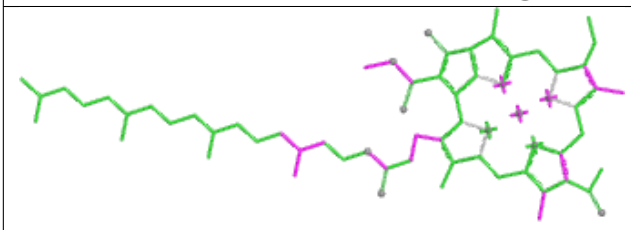


Torsions

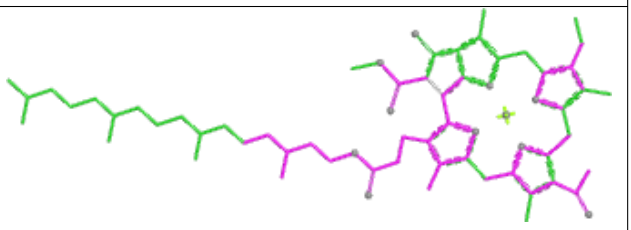


Rings

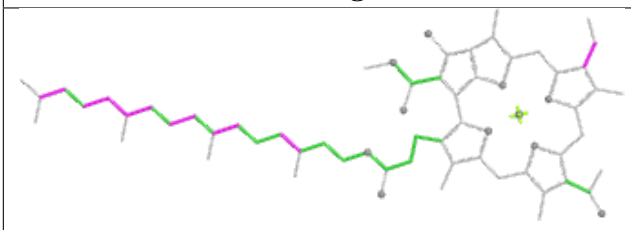
Ligand BCL B0 102



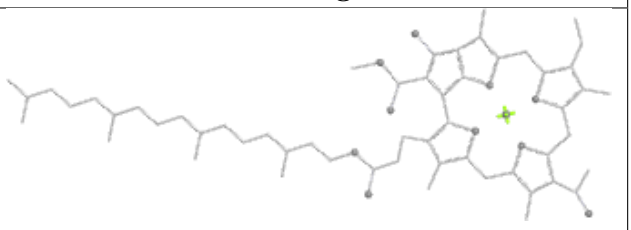
Bond lengths



Bond angles

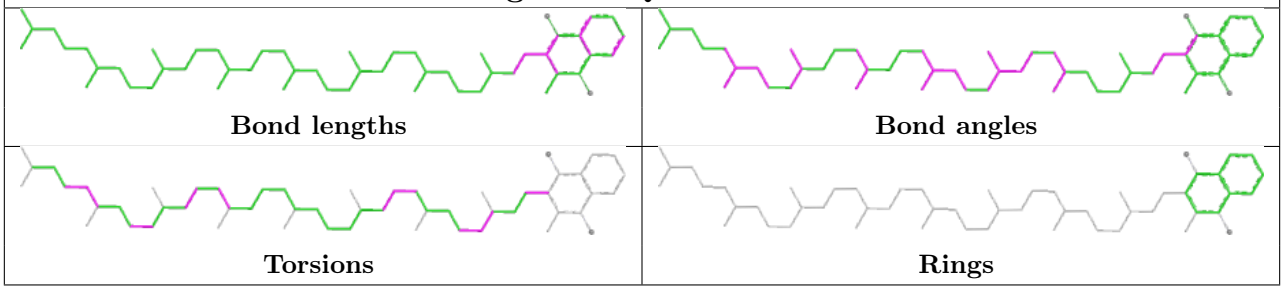


Torsions

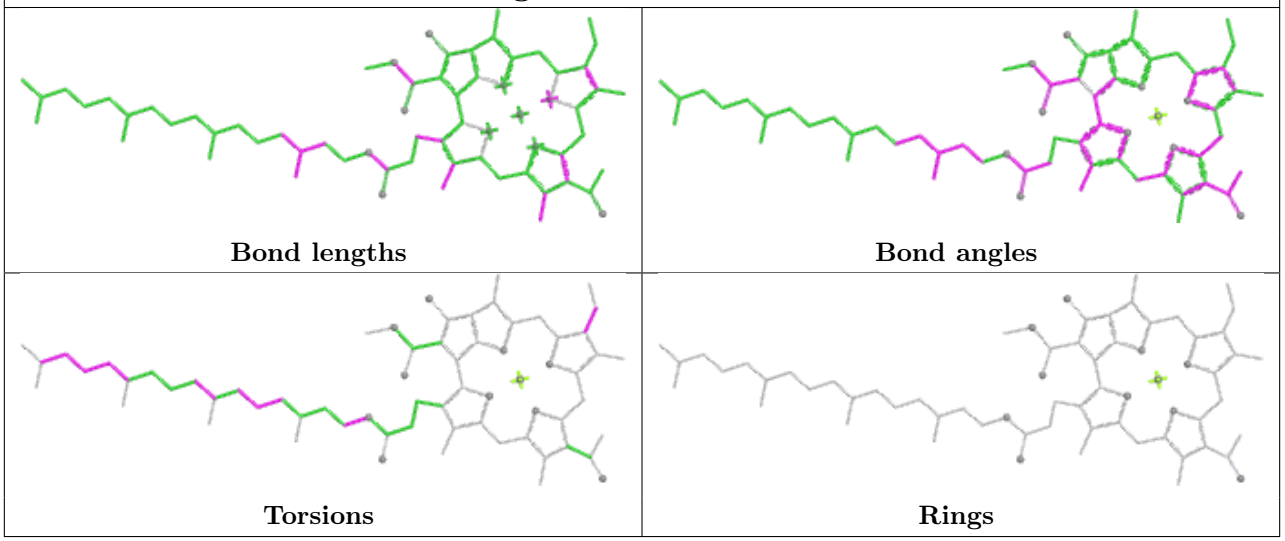


Rings

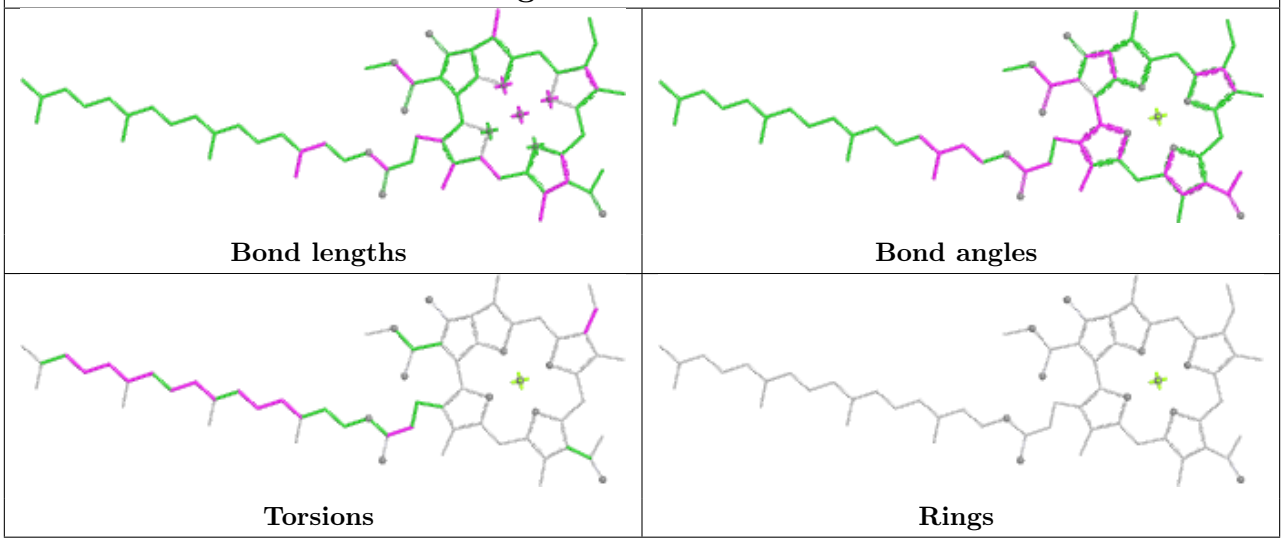
Ligand MQ8 BM 405



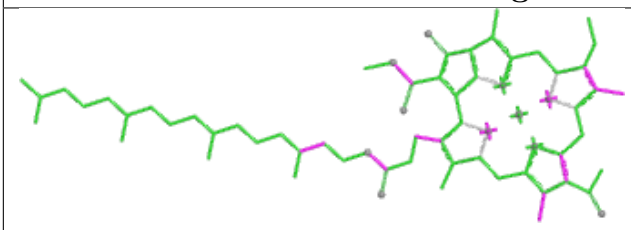
Ligand BCL BA 101



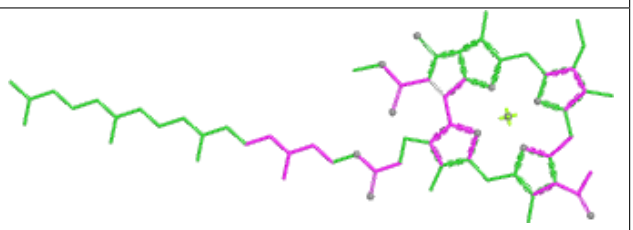
Ligand BCL B1 102



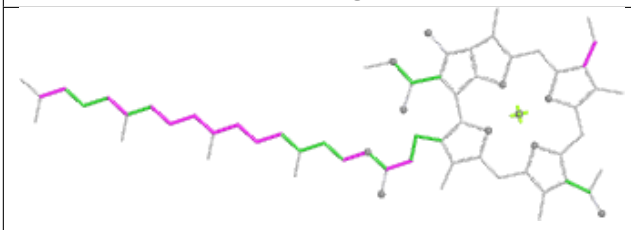
Ligand BCL BU 102



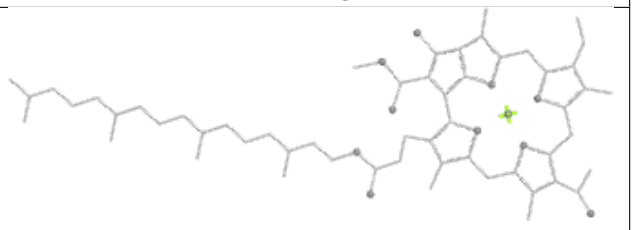
Bond lengths



Bond angles

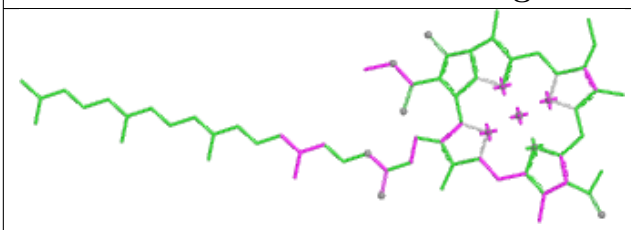


Torsions

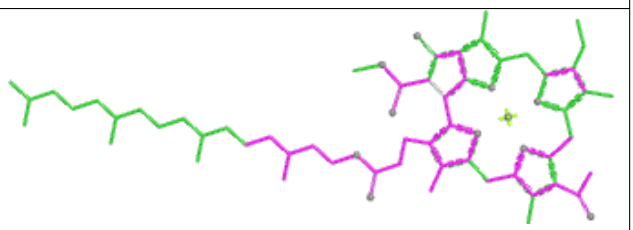


Rings

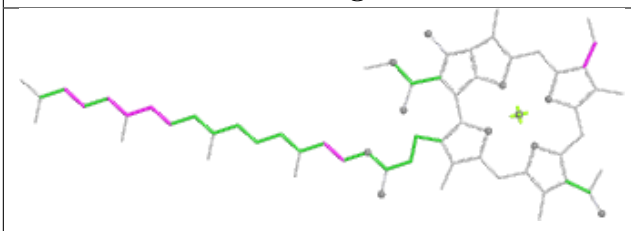
Ligand BCL AT 101



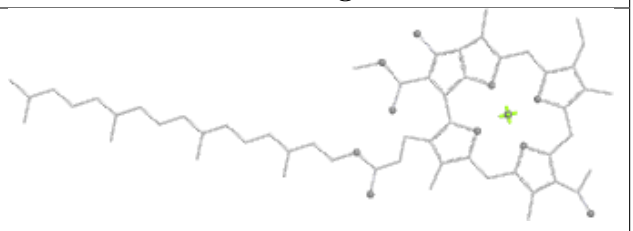
Bond lengths



Bond angles

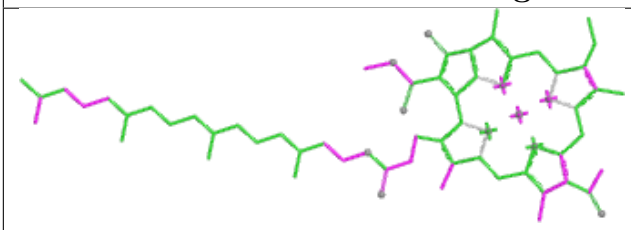


Torsions

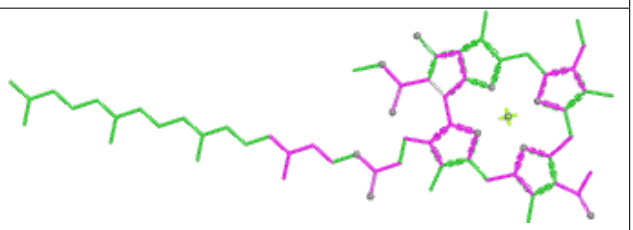


Rings

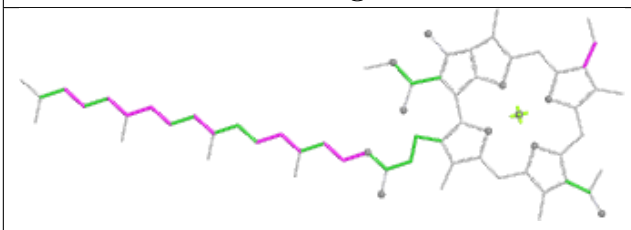
Ligand BCL BE 101



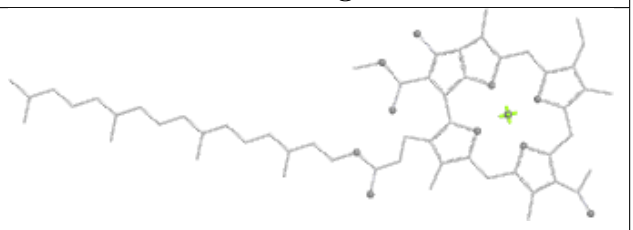
Bond lengths



Bond angles

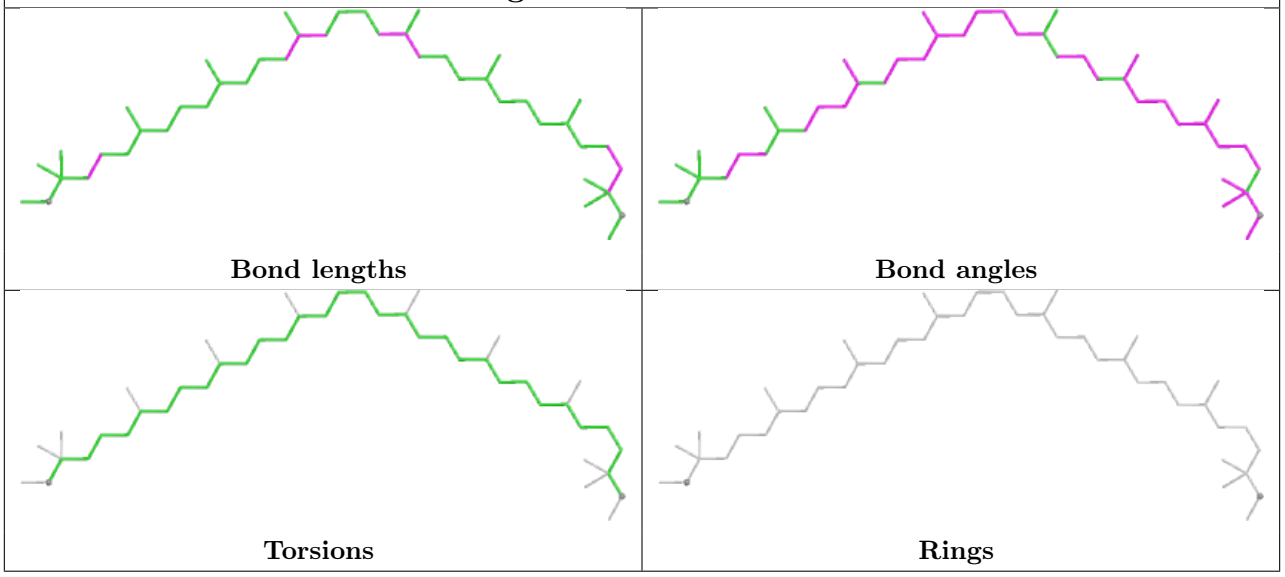


Torsions

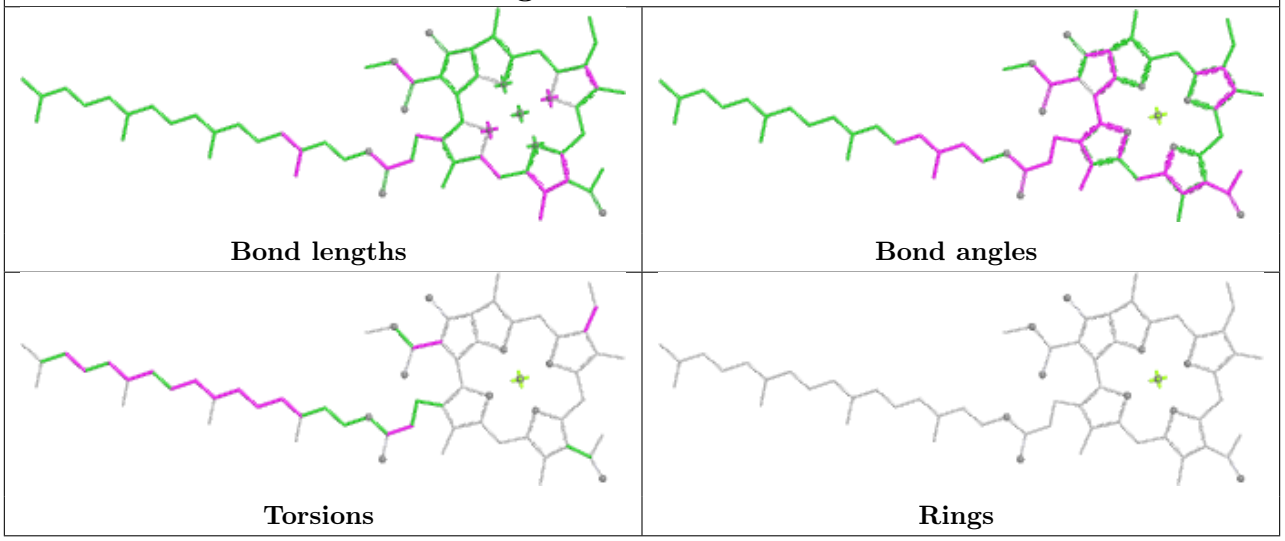


Rings

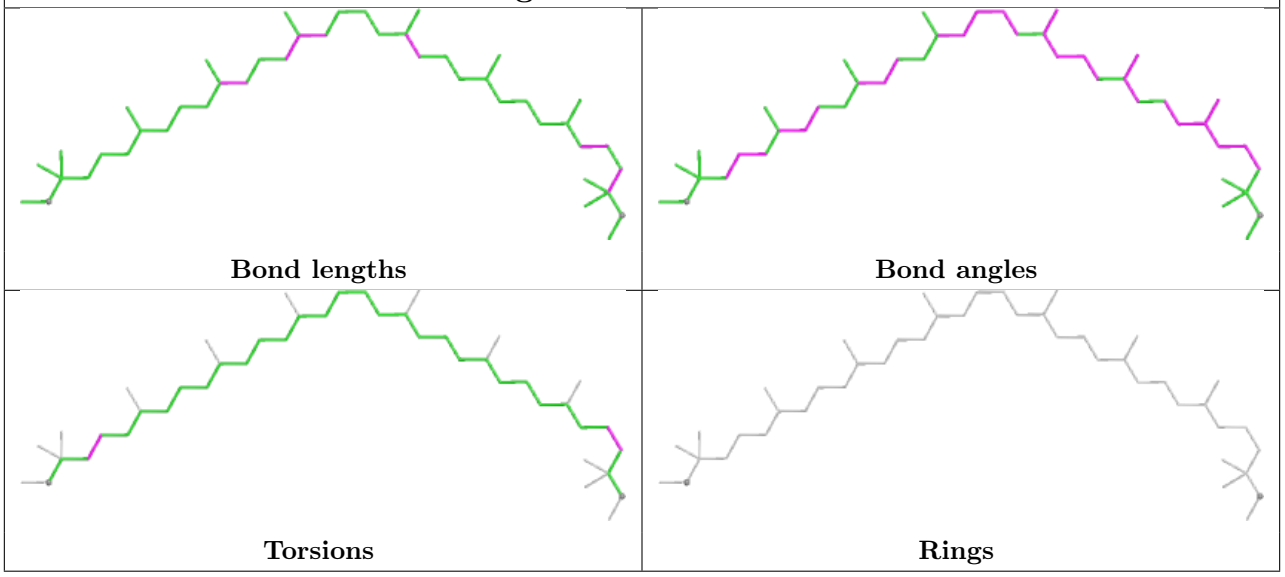
Ligand CRT BW 103



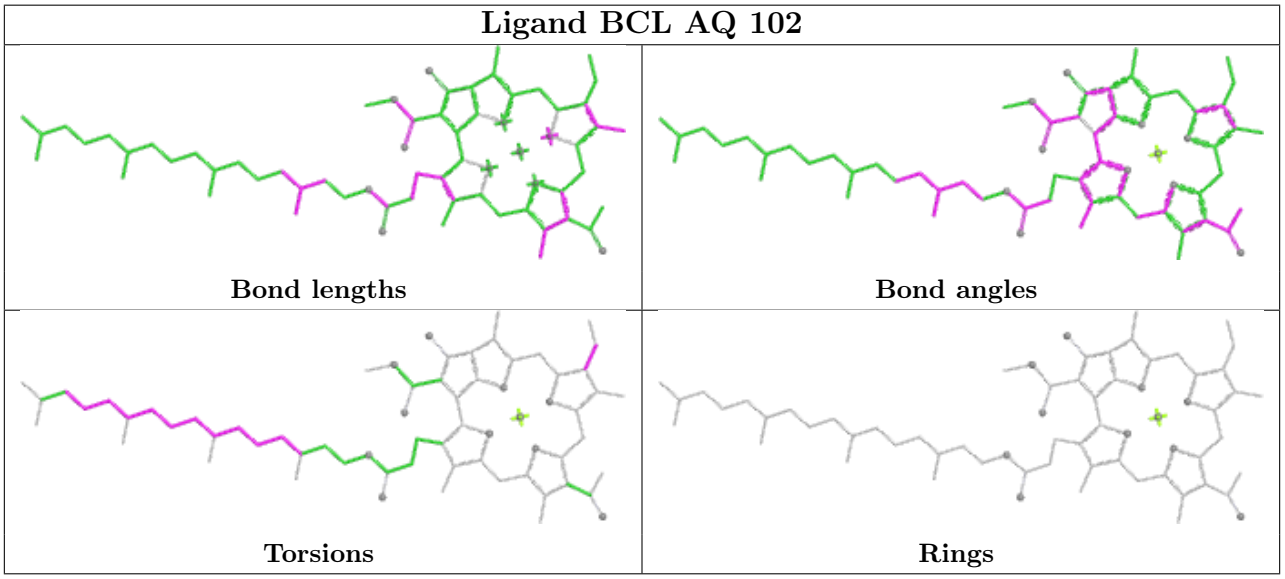
Ligand BCL A3 103



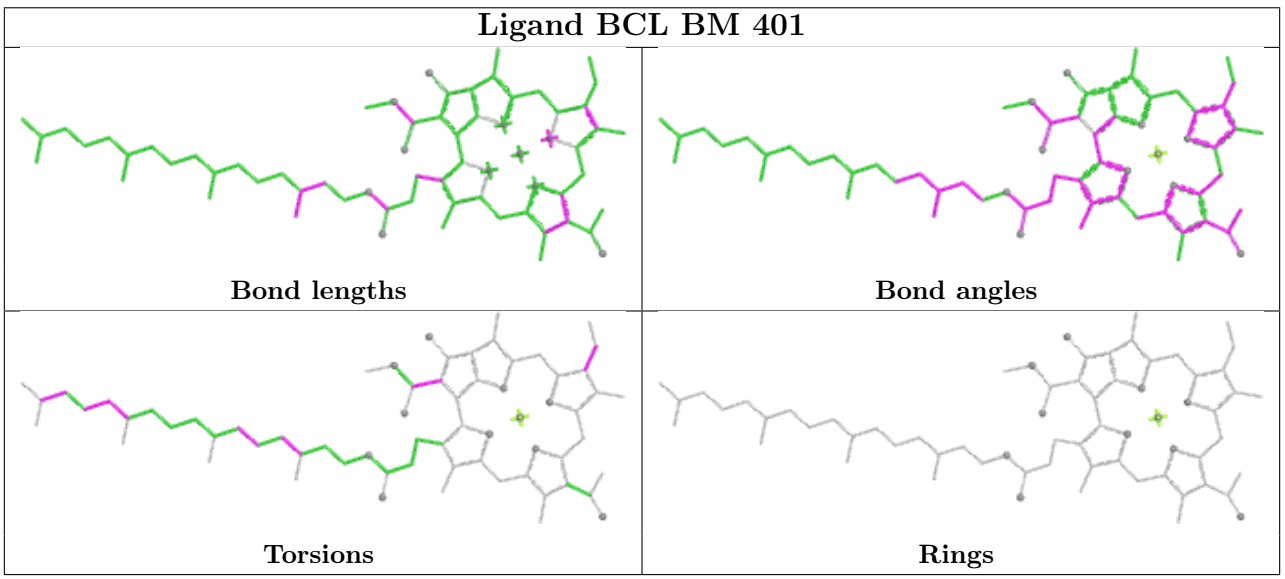
Ligand CRT A0 101



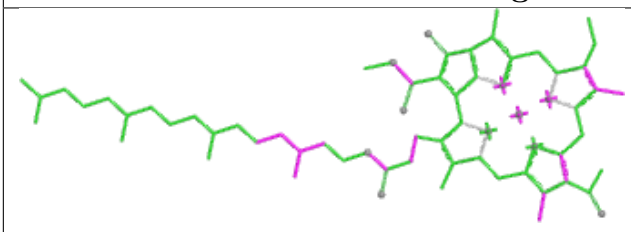
Ligand BCL AQ 102



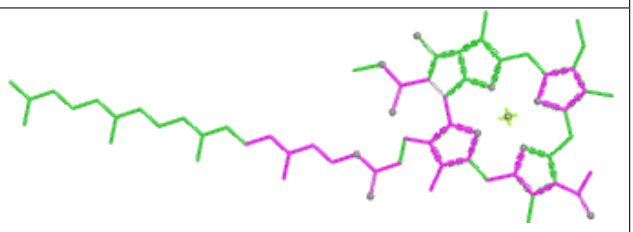
Ligand BCL BM 401



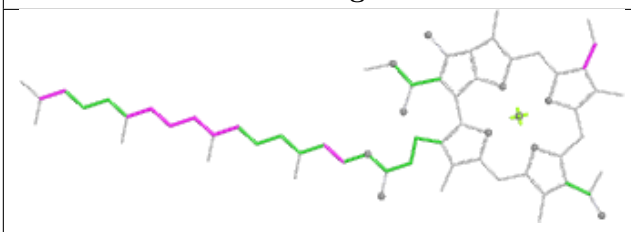
Ligand BCL BN 101



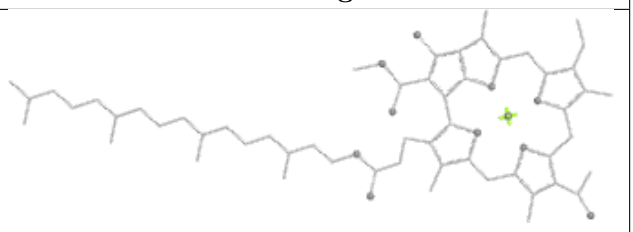
Bond lengths



Bond angles

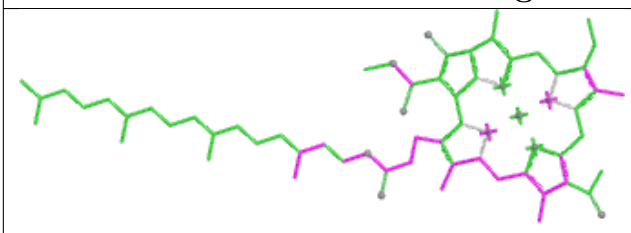


Torsions

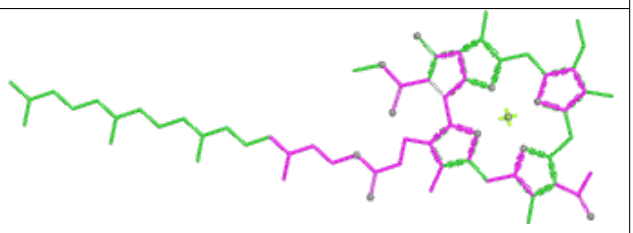


Rings

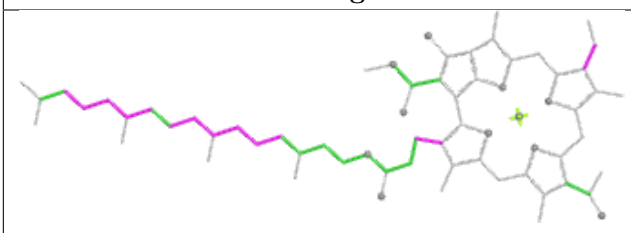
Ligand BCL A1 102



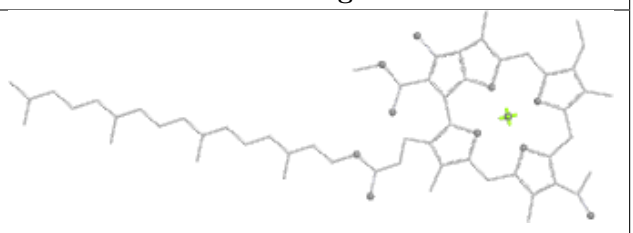
Bond lengths



Bond angles

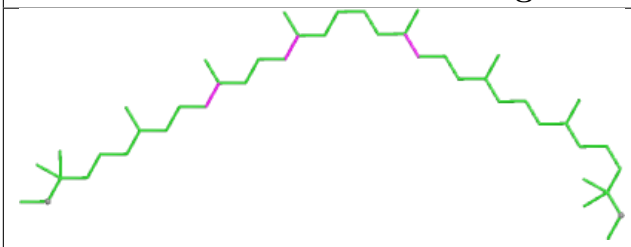


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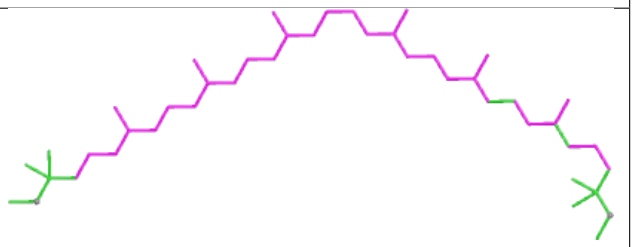


Rings

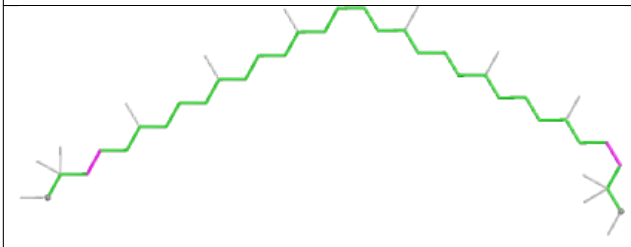
Ligand CRT B1 103



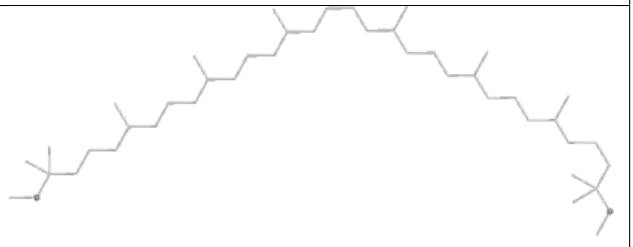
Bond lengths



Bond angles

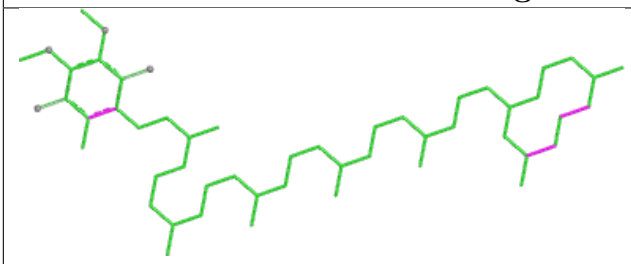


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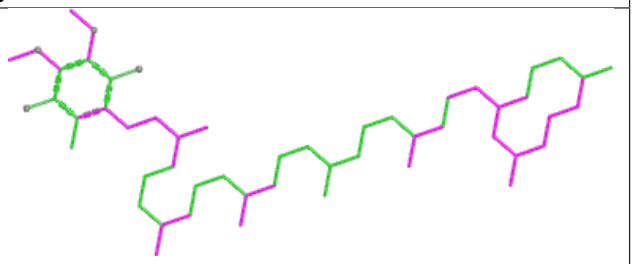


Rings

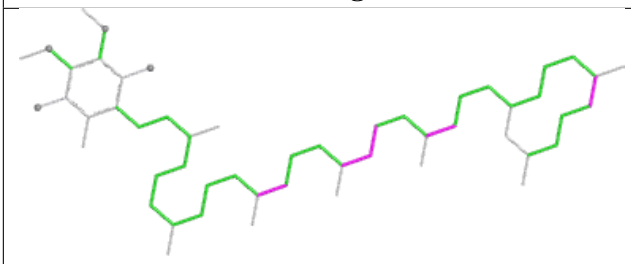
Ligand UQ8 AL 304



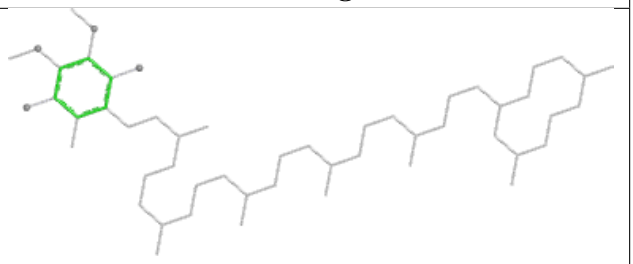
Bond lengths



Bond angles

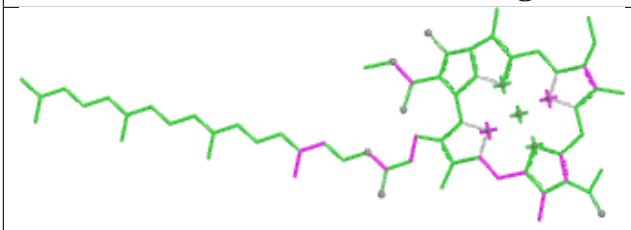


Torsions

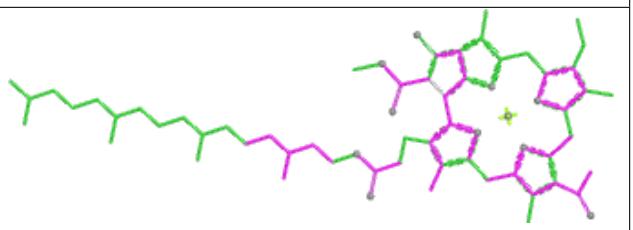


Rings

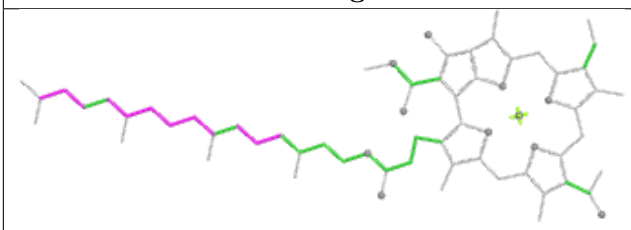
Ligand BCL B5 102



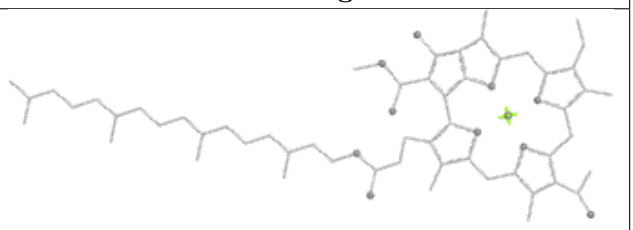
Bond lengths



Bond angles

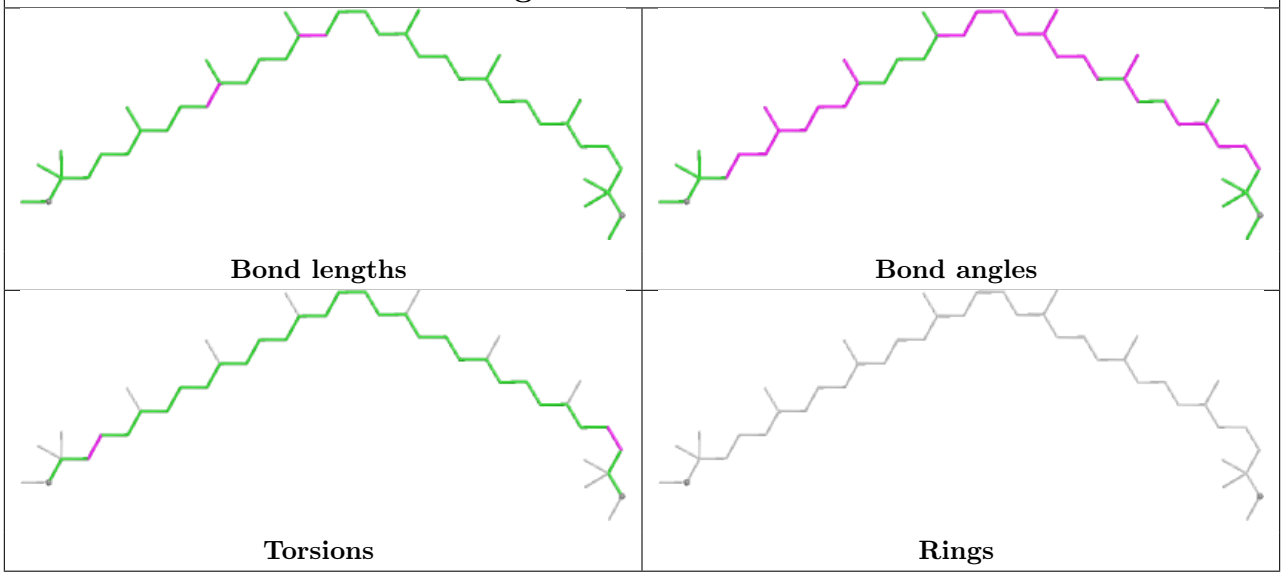


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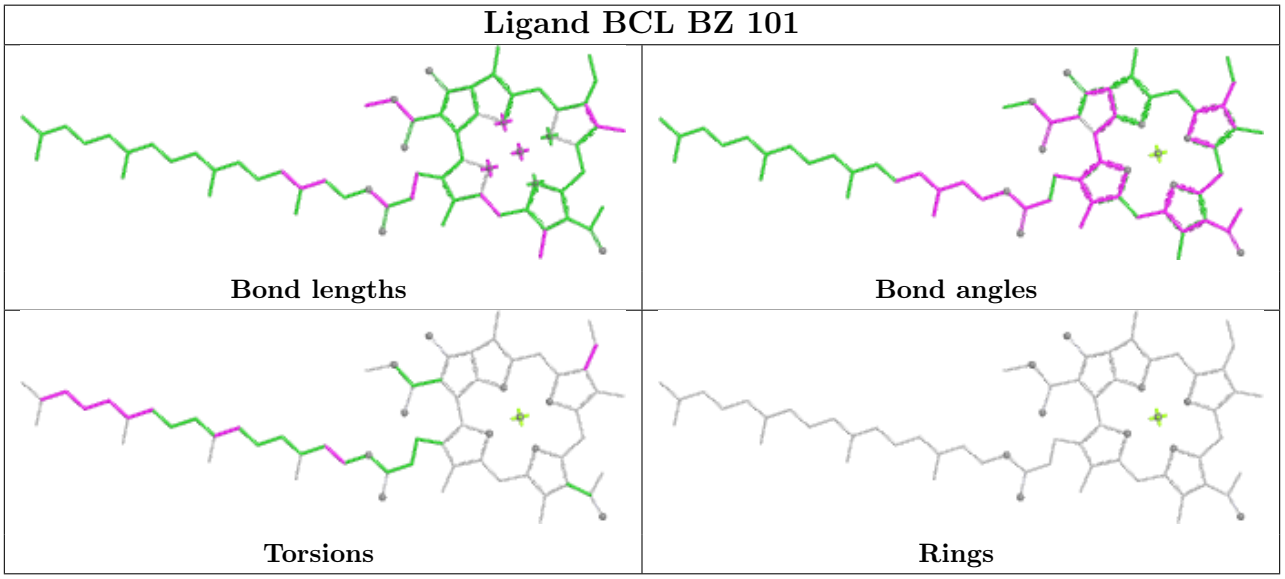


Rings

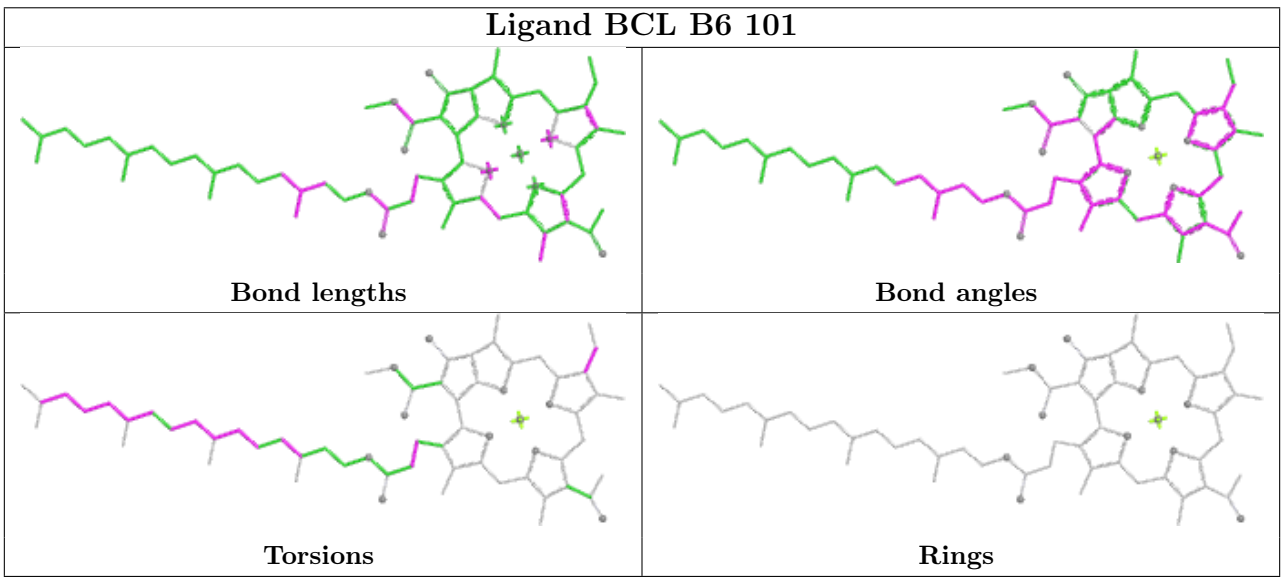
Ligand CRT AS 104



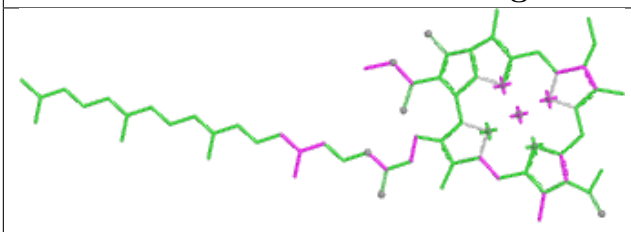
Ligand BCL BZ 101



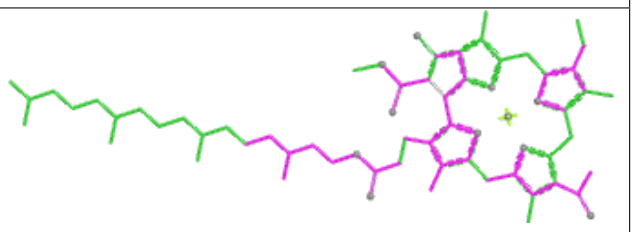
Ligand BCL B6 101



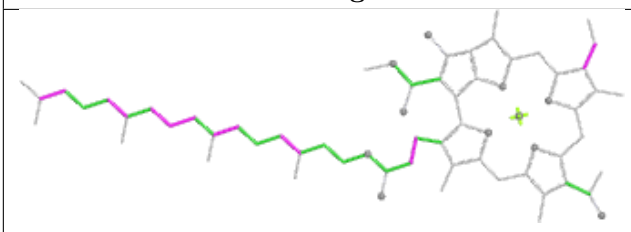
Ligand BCL AN 101



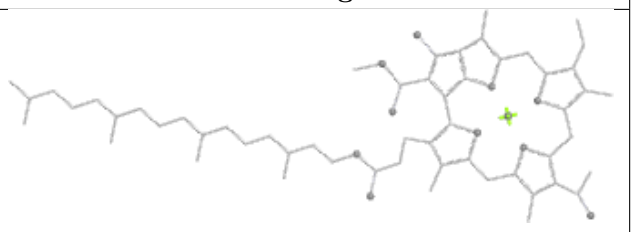
Bond lengths



Bond angles

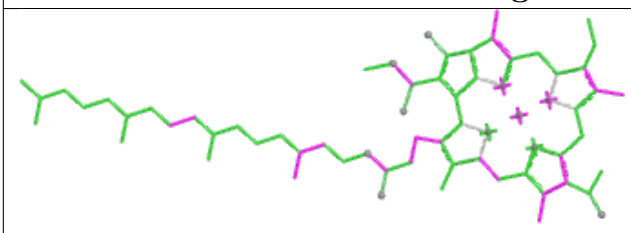


Torsions

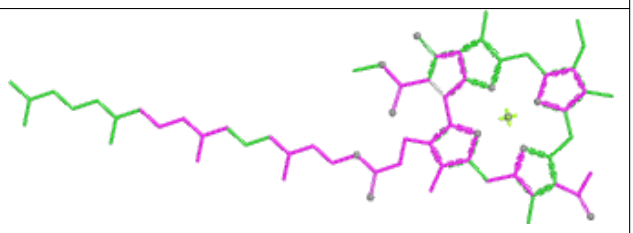


Rings

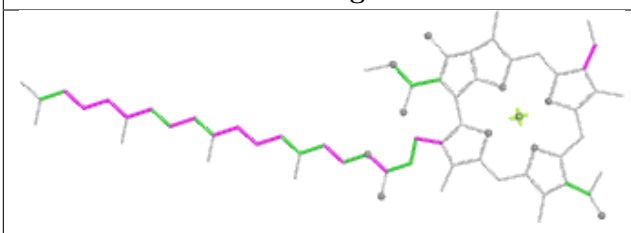
Ligand BCL AU 102



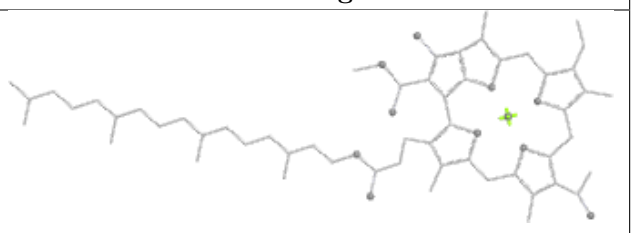
Bond lengths



Bond angles

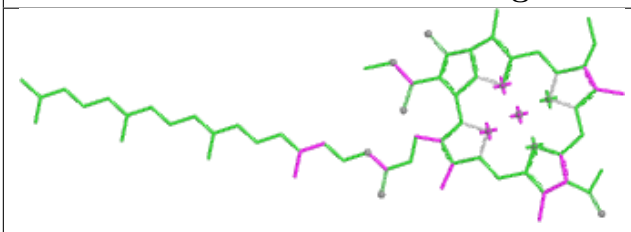


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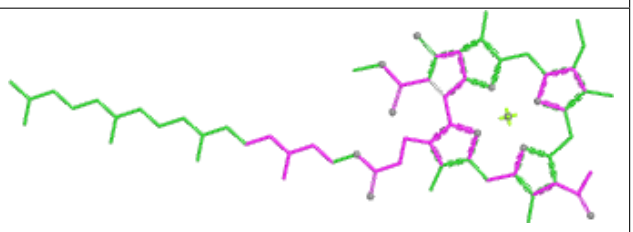


Rings

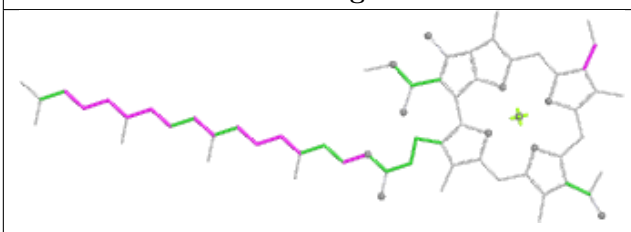
Ligand BCL BO 102



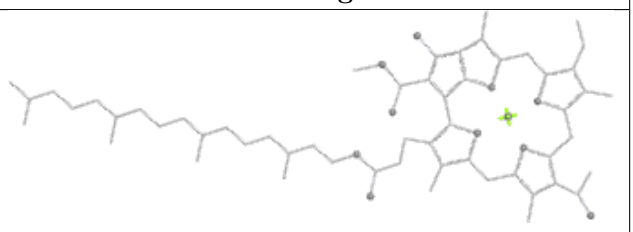
Bond lengths



Bond angles

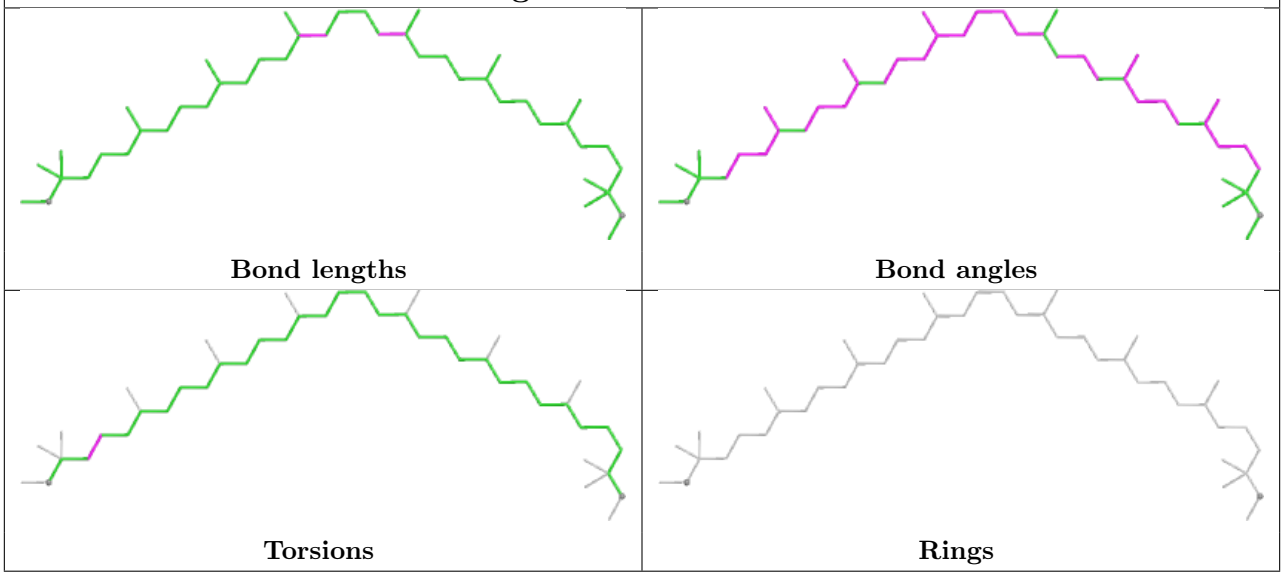


Torsions

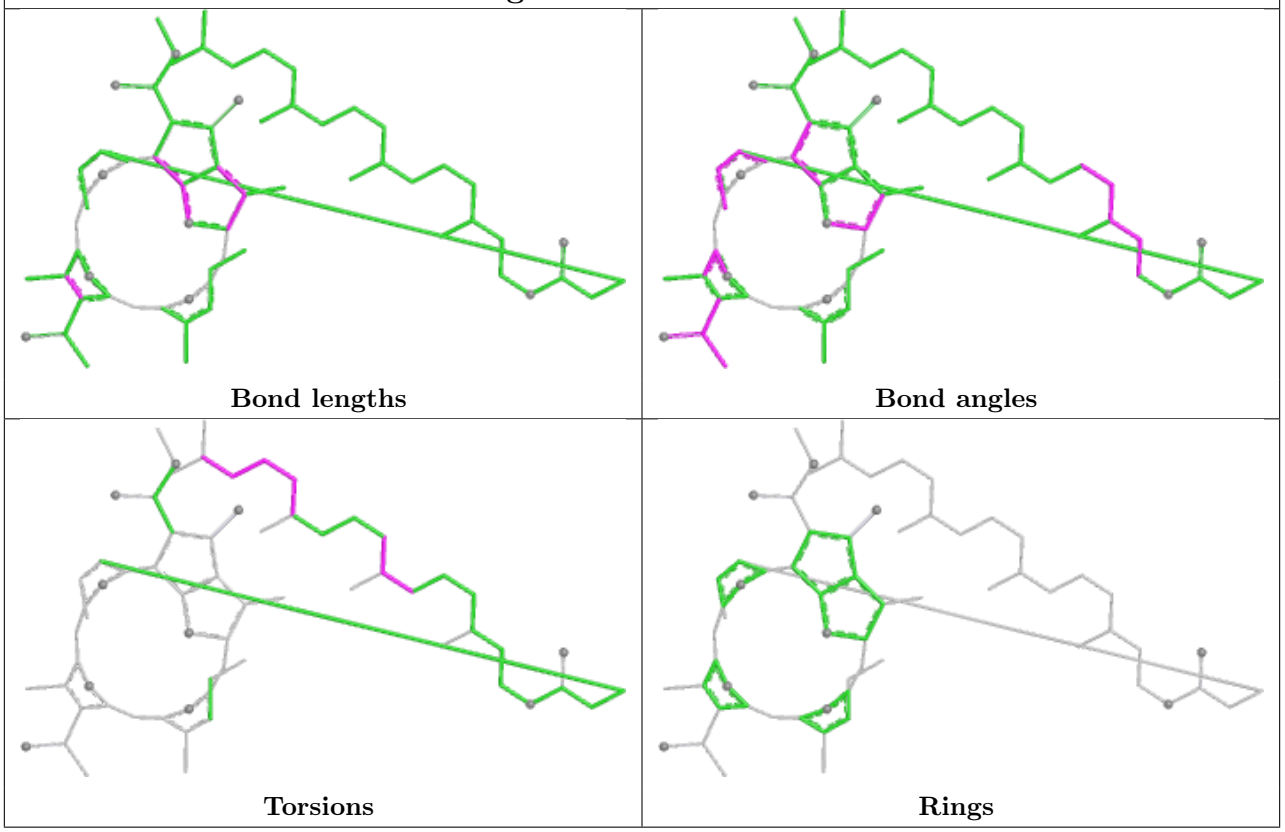


Rings

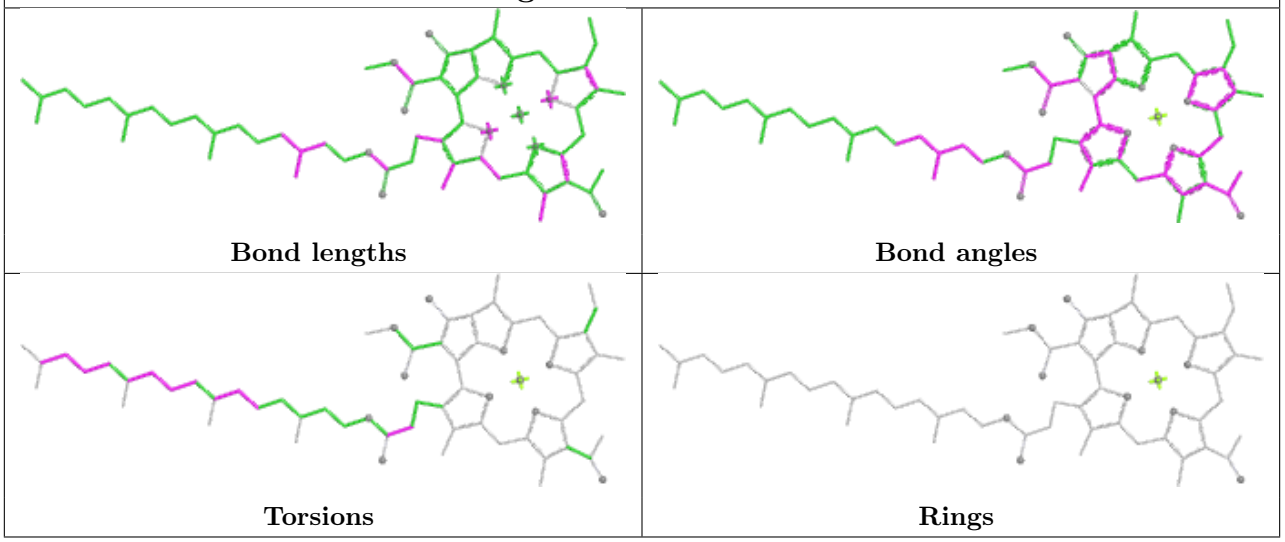
Ligand CRT BA 102



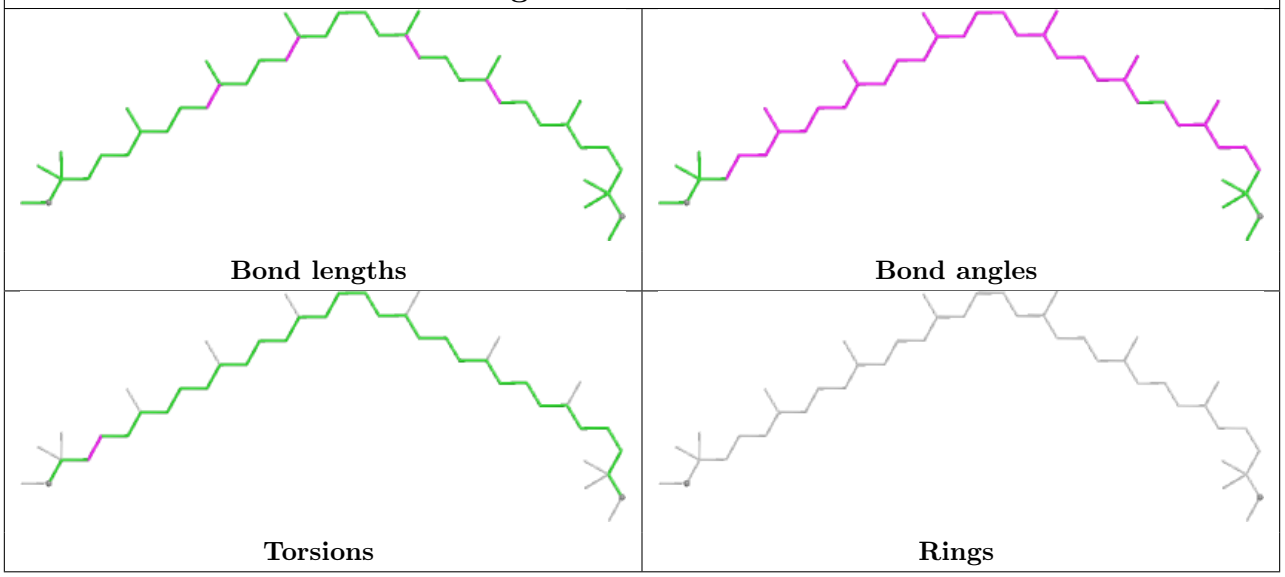
Ligand BPH BM 403



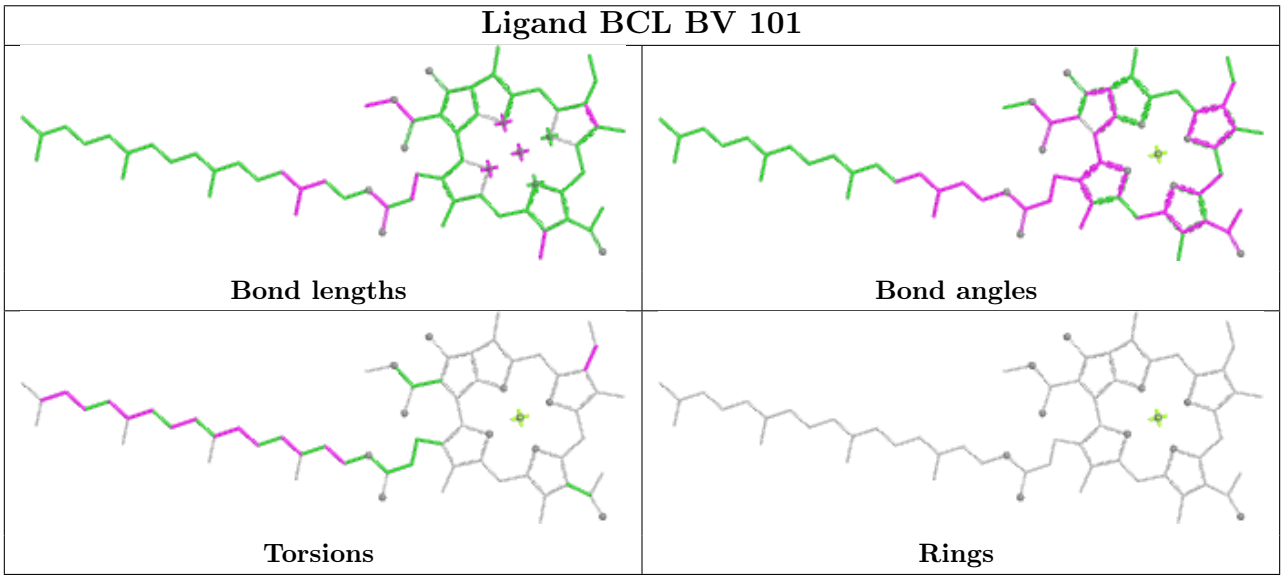
Ligand BCL B7 103



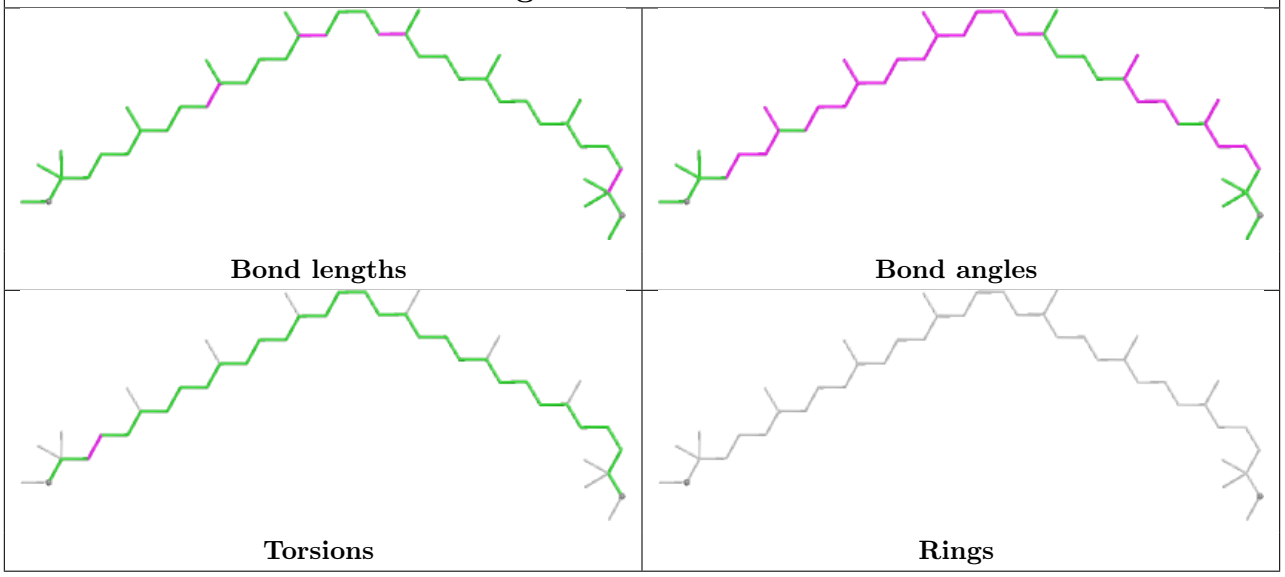
Ligand CRT AN 102



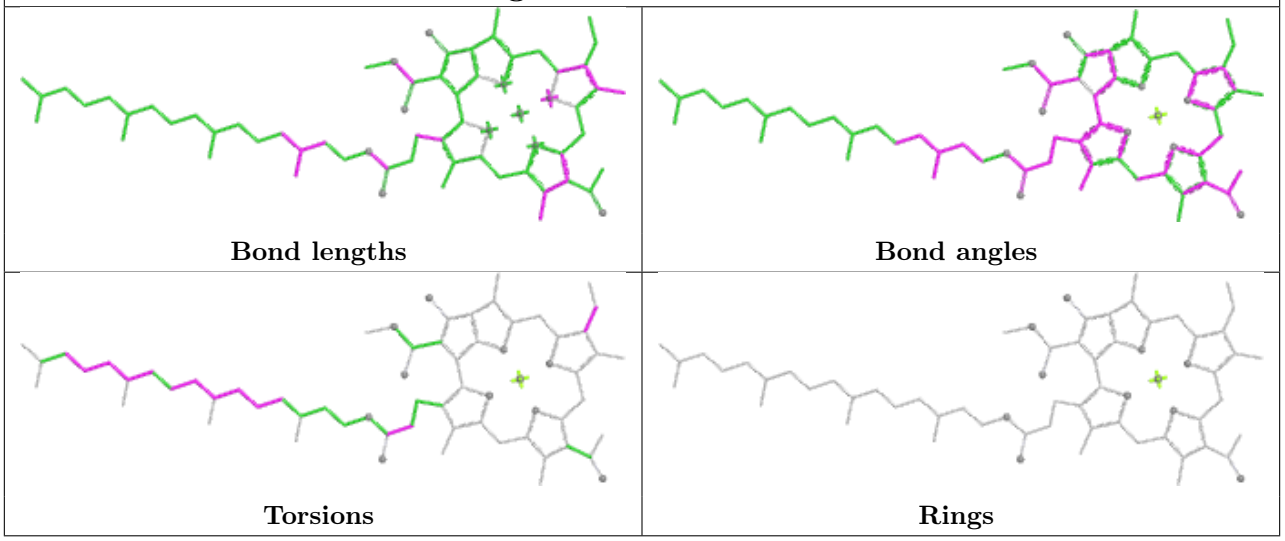
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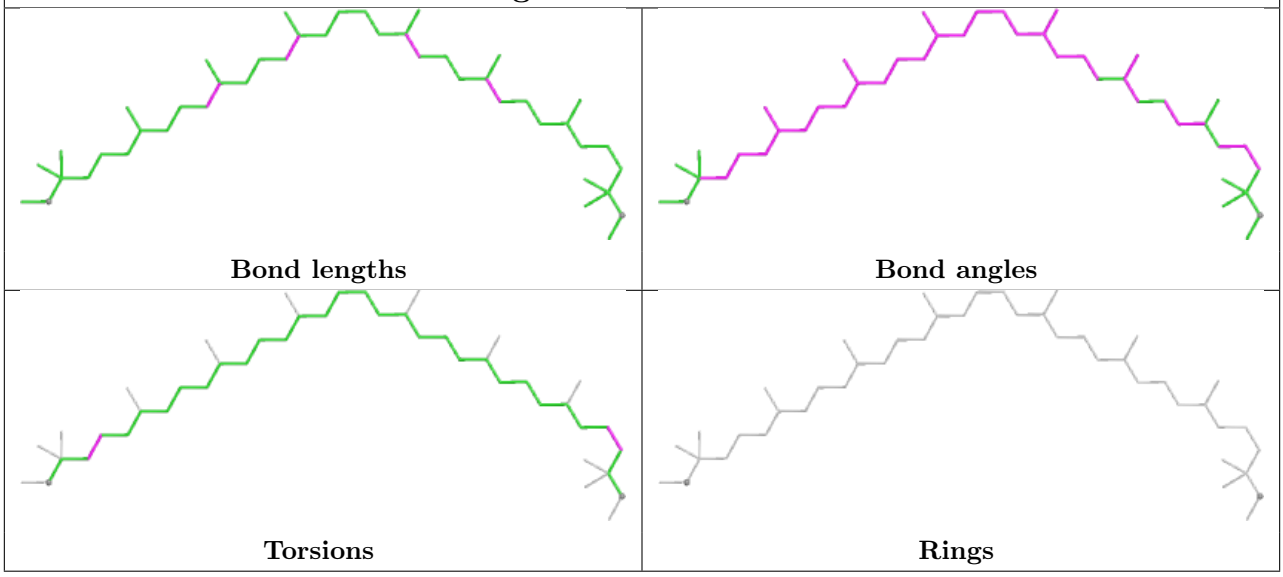
Ligand CRT B5 103



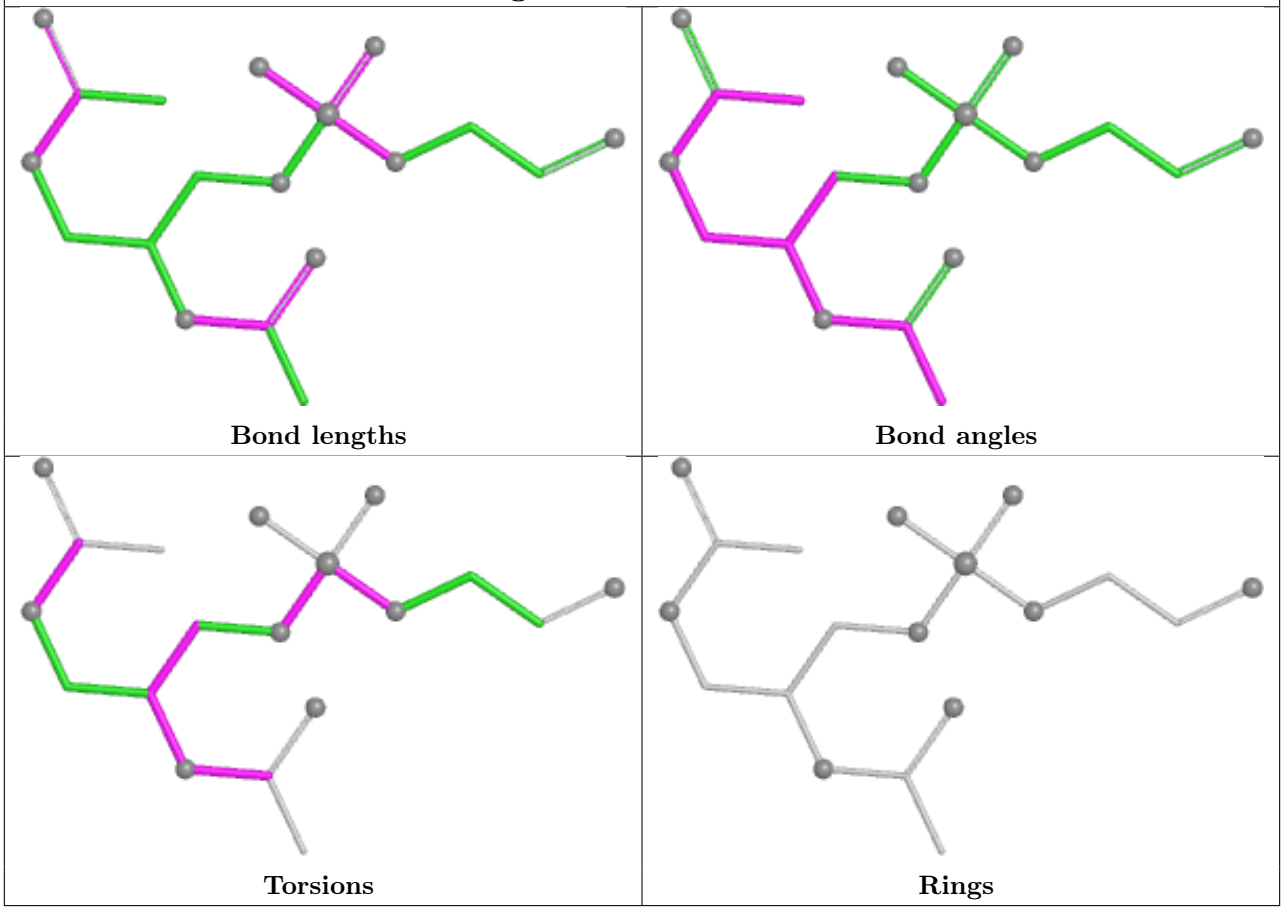
Ligand BCL BW 102



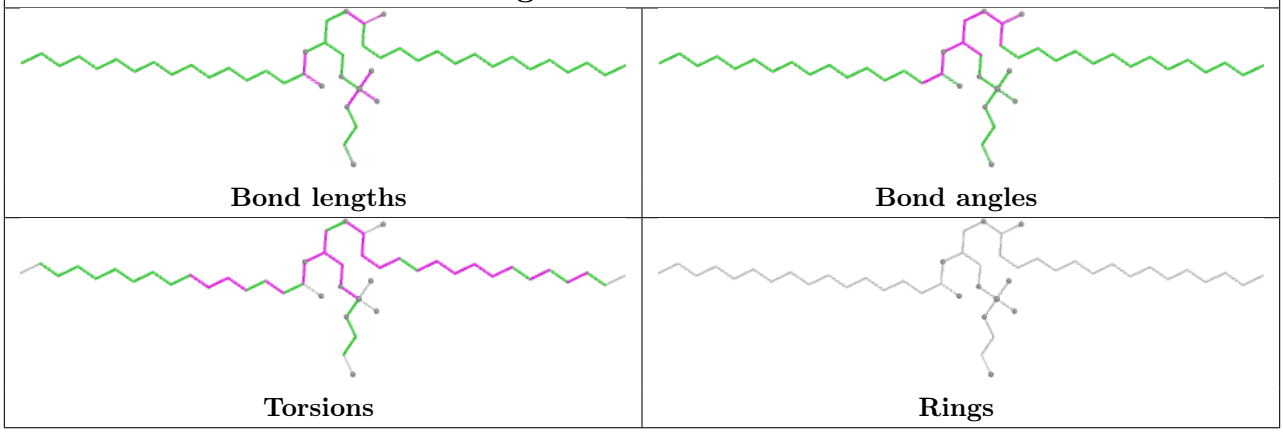
Ligand CRT BB 102



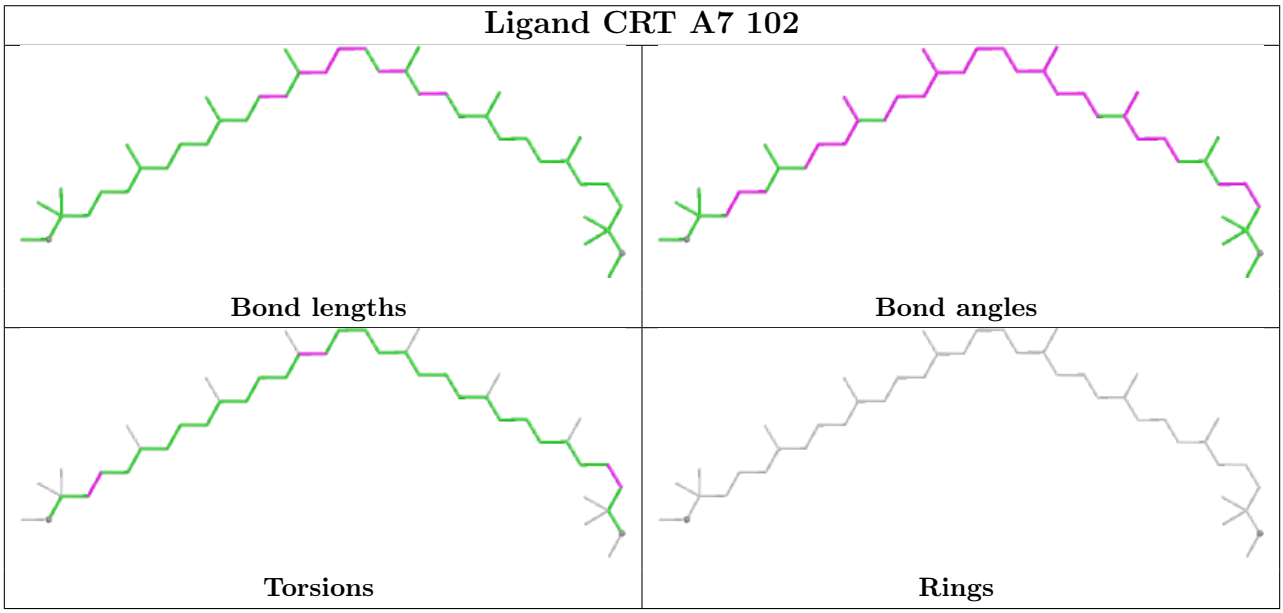
Ligand PEF AM 407



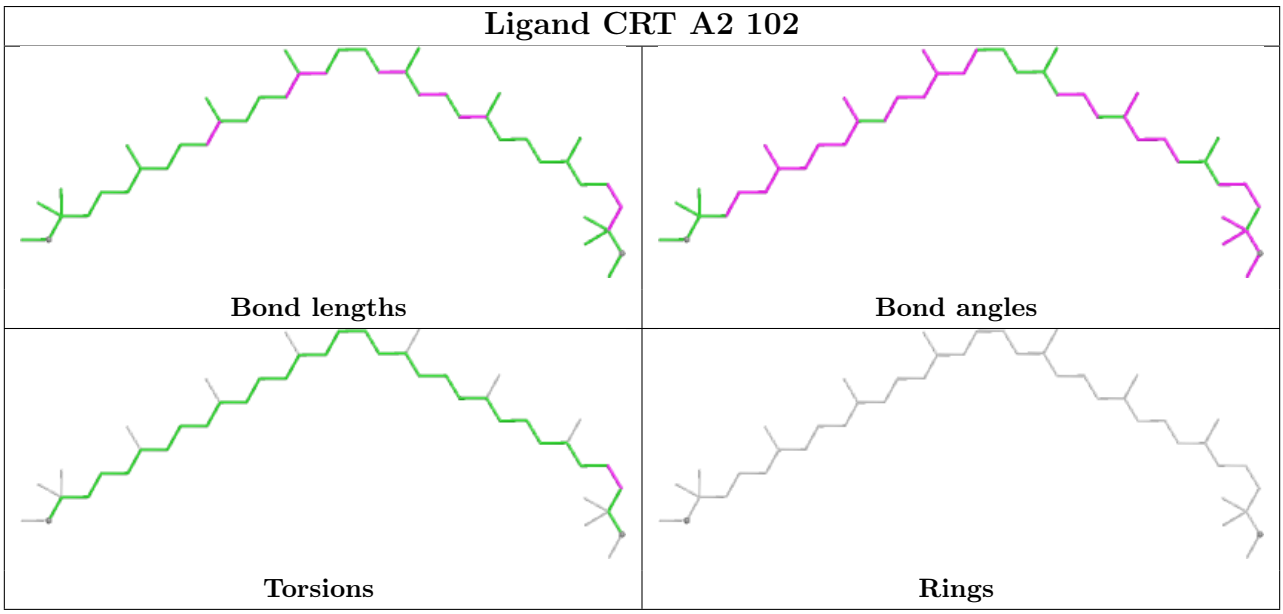
Ligand PEF AS 101



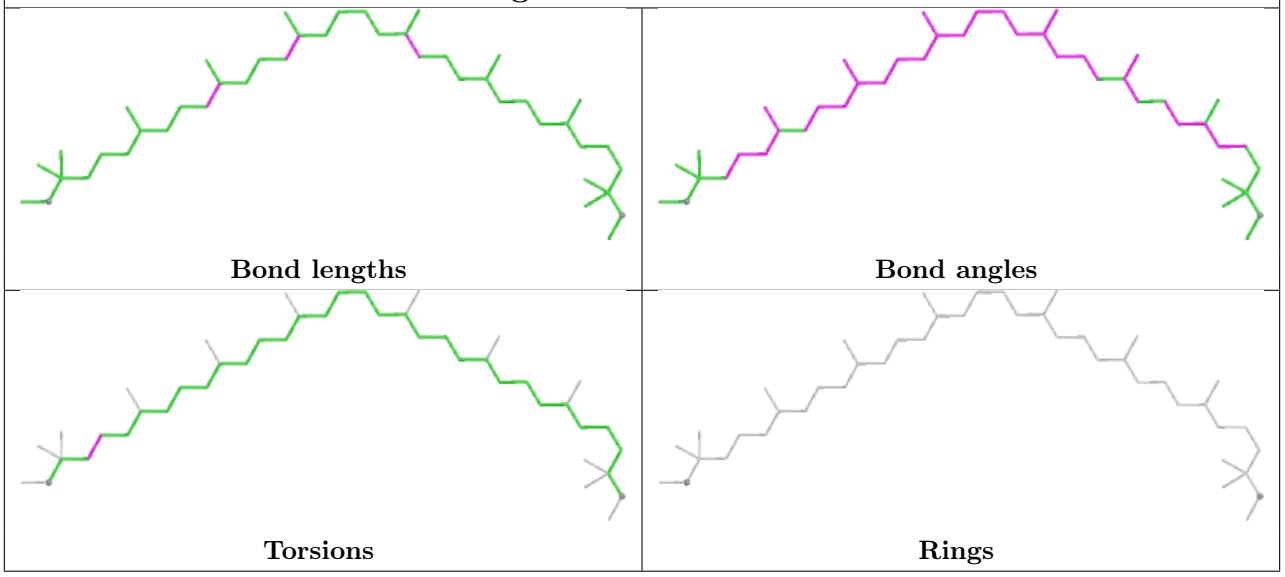
Ligand CRT A7 102



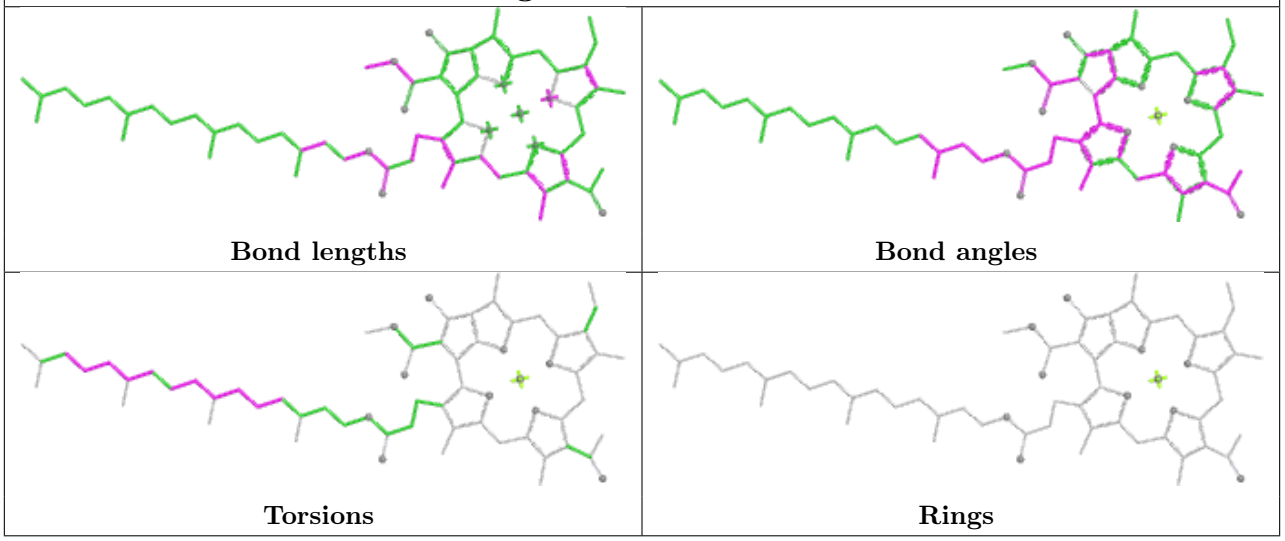
Ligand CRT A2 102



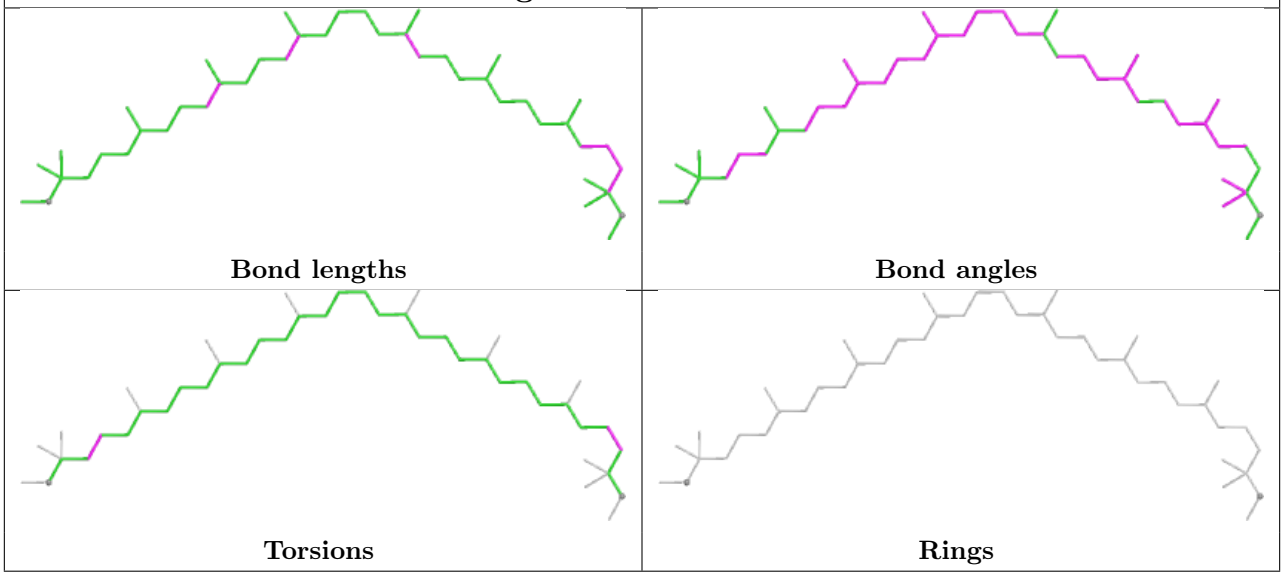
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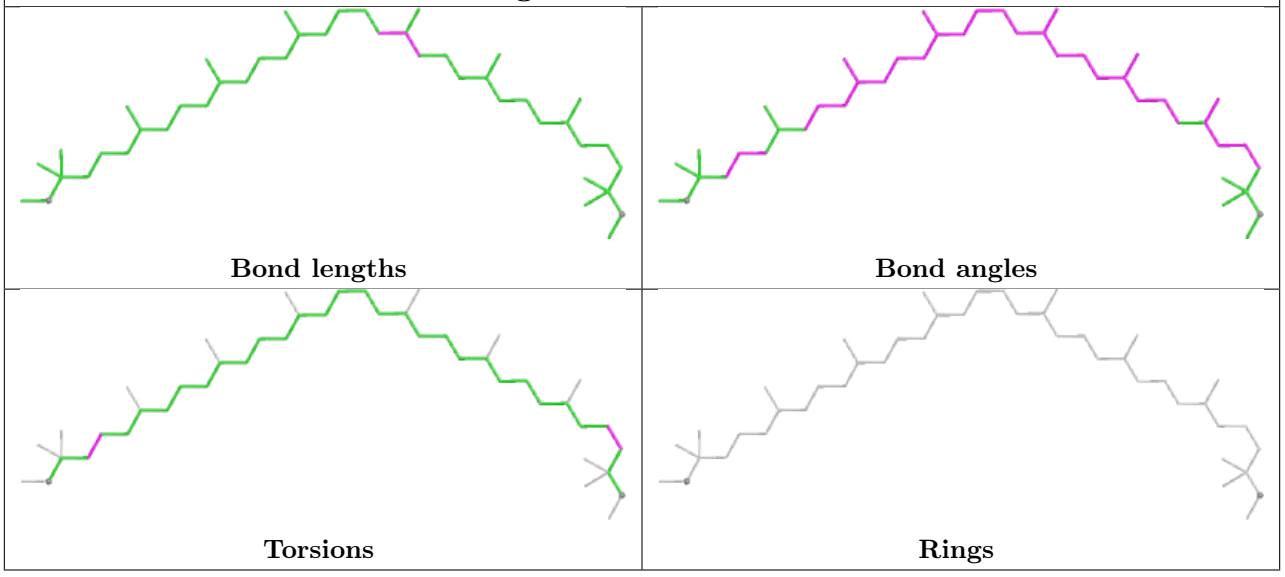
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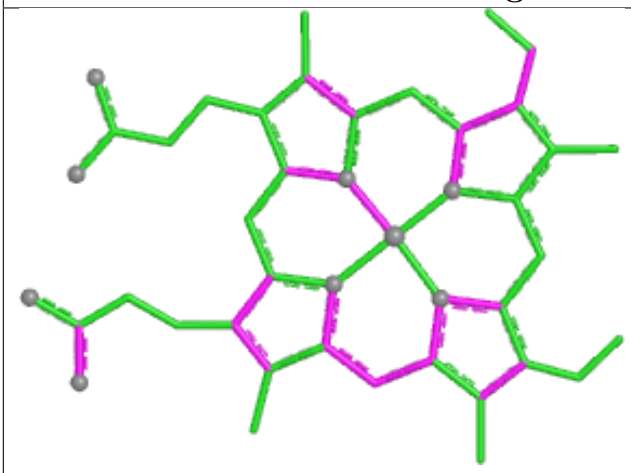
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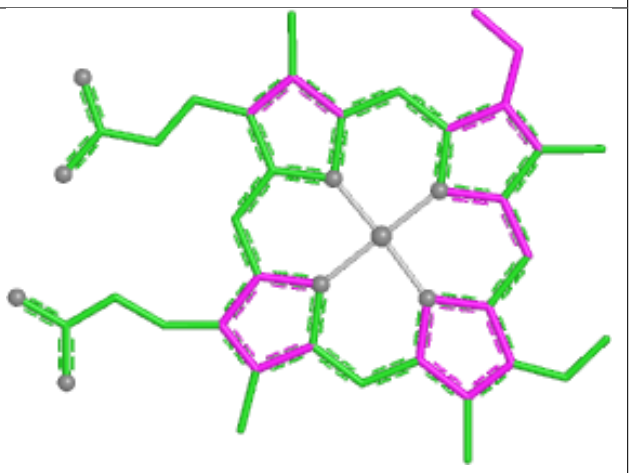
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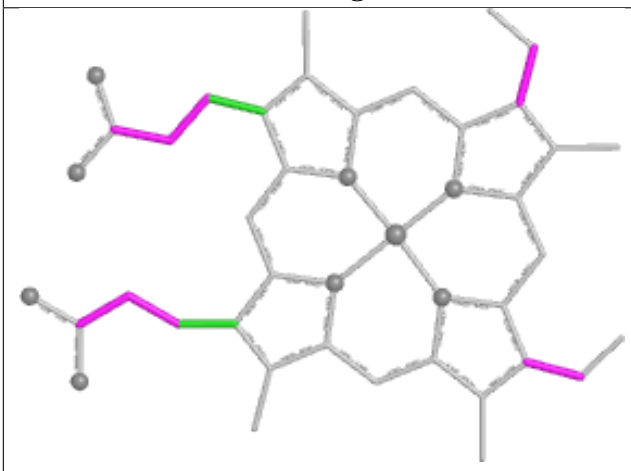
Ligand HEM BC 503



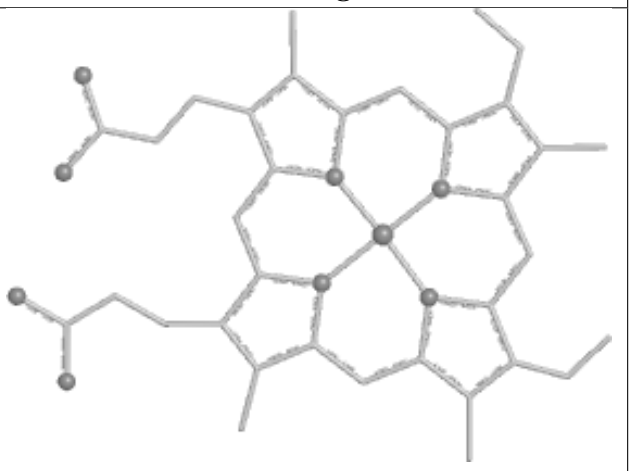
Bond lengths



Bond angles

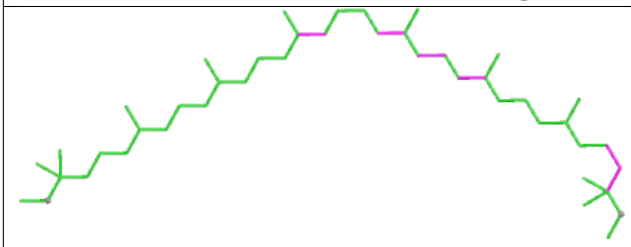


Torsions

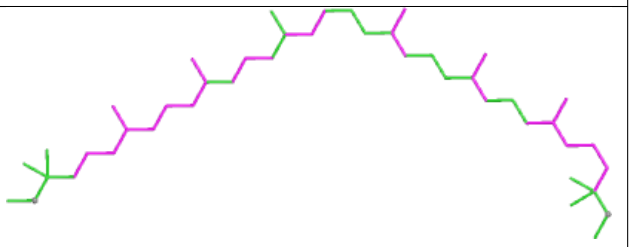


Rings

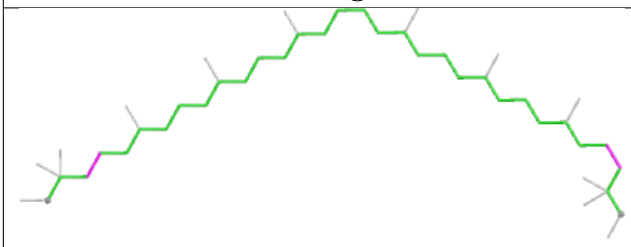
Ligand CRT B7 102



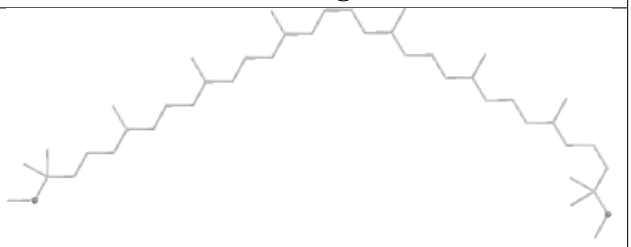
Bond lengths



Bond angles

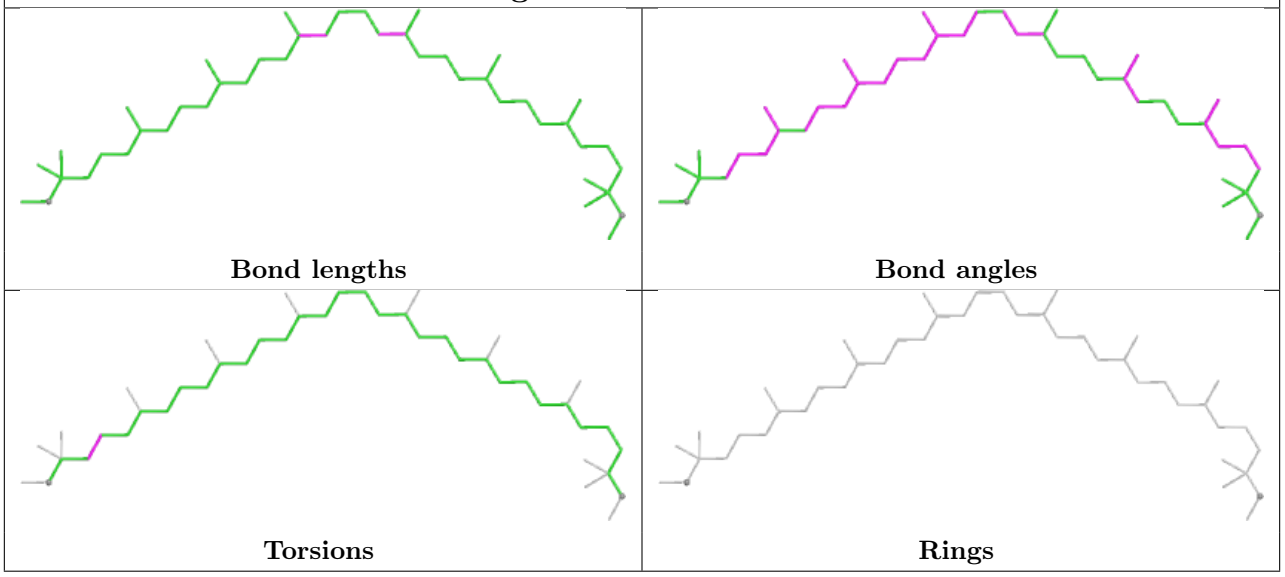


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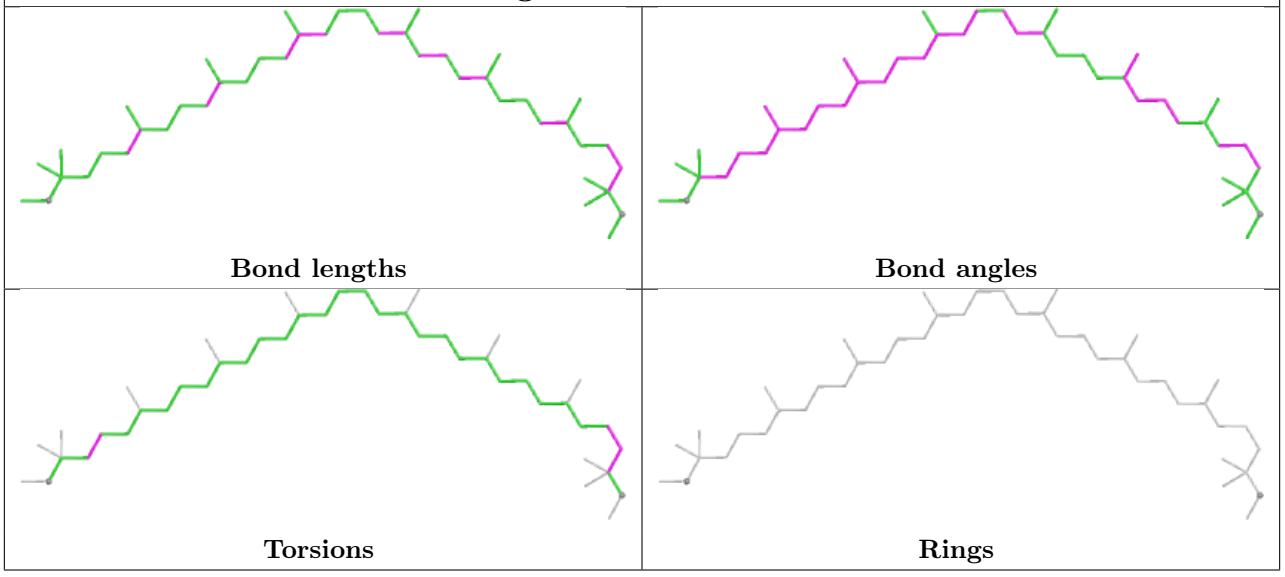


Rings

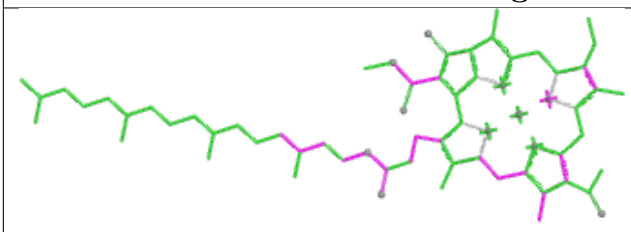
Ligand CRT AA 102



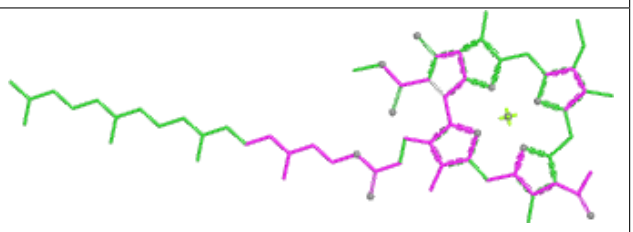
Ligand CRT B0 101



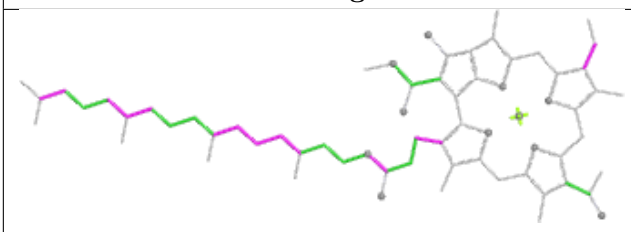
Ligand BCL AX 101



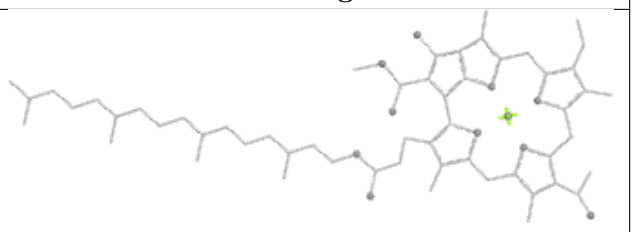
Bond lengths



Bond angles

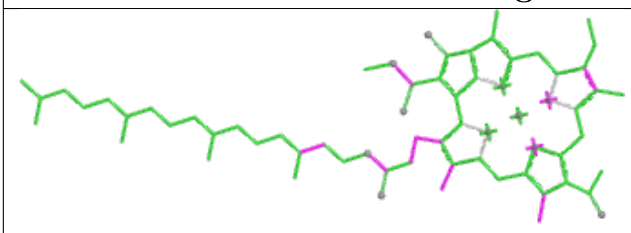


Torsions

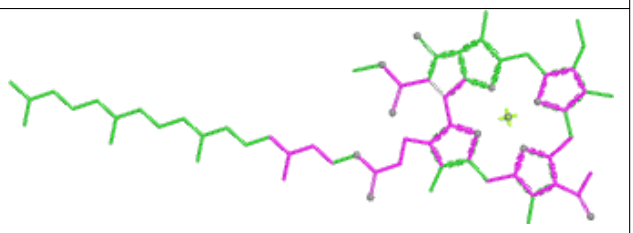


Rings

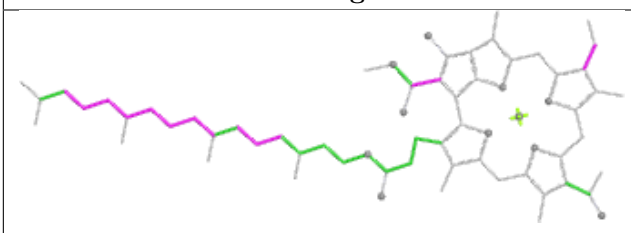
Ligand BCL BS 102



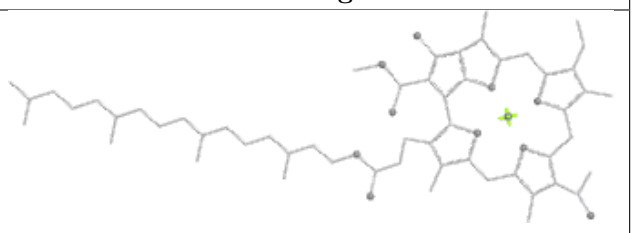
Bond lengths



Bond angles

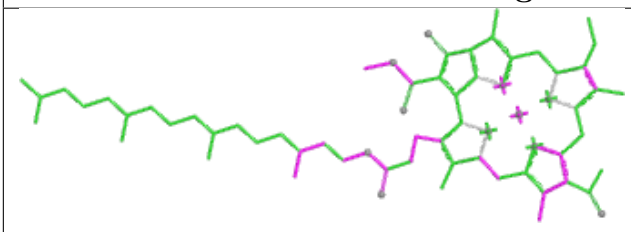


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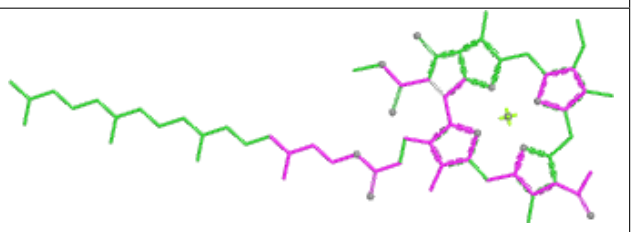


Rings

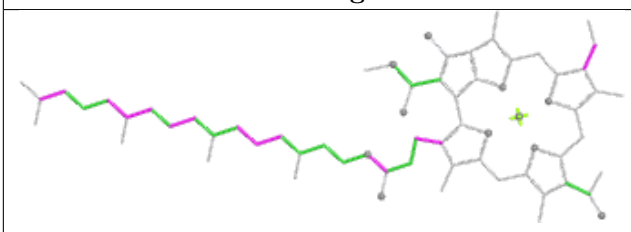
Ligand BCL A2 101



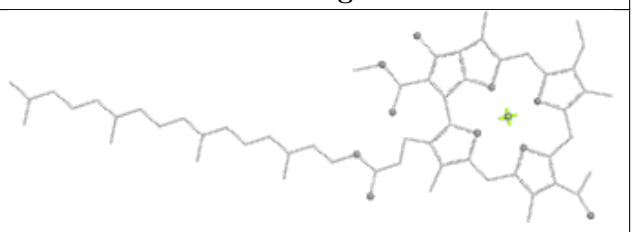
Bond lengths



Bond angles

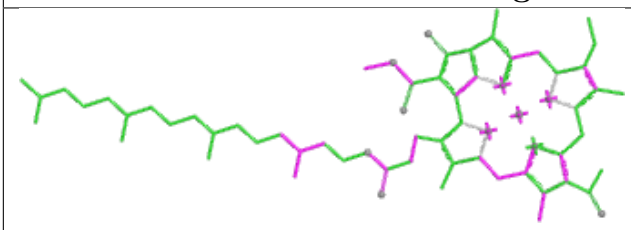


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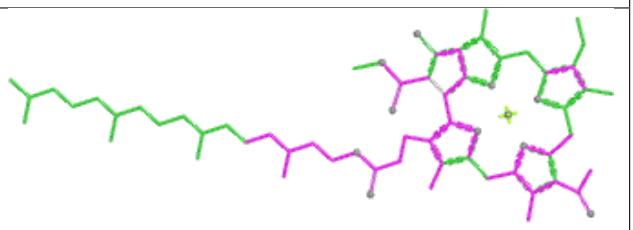


Rings

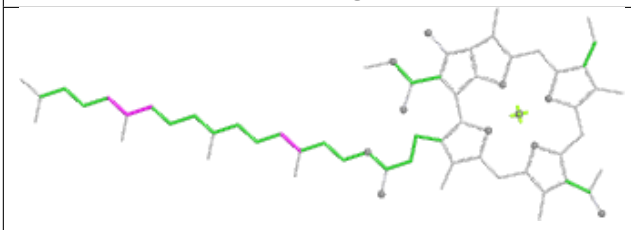
Ligand BCL BP 101



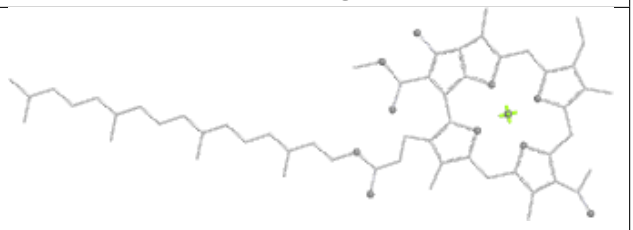
Bond lengths



Bond angles

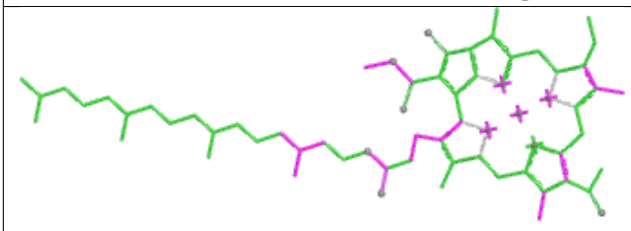


Torsions

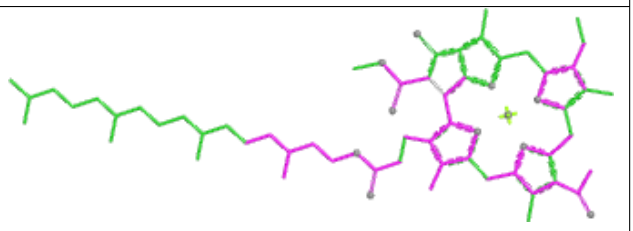


Rings

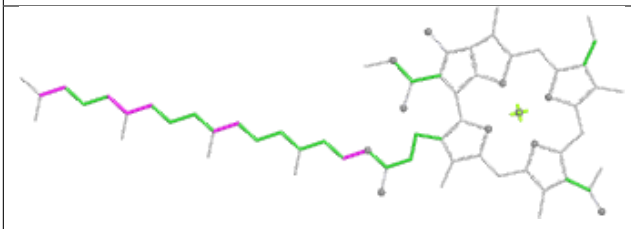
Ligand BCL B4 101



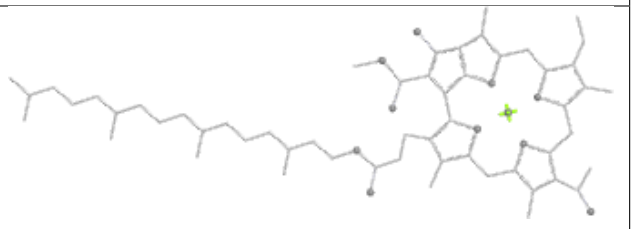
Bond lengths



Bond angles

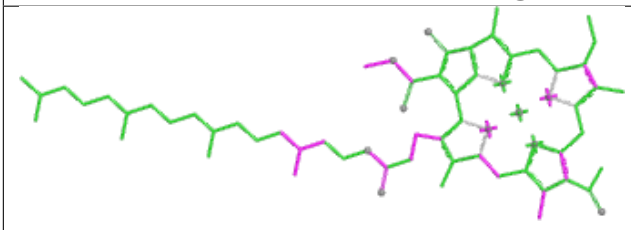


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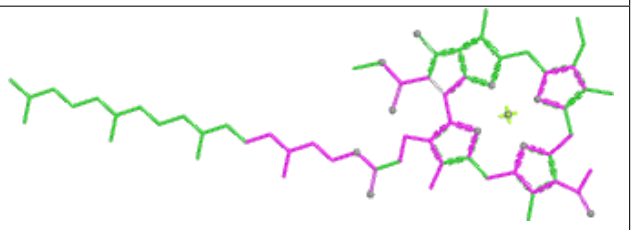


Rings

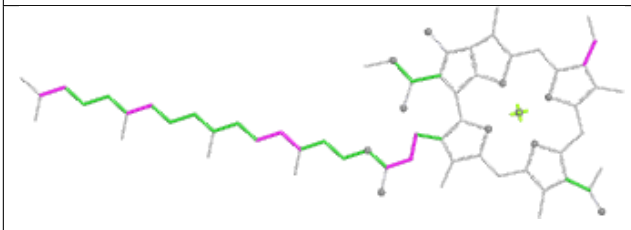
Ligand BCL AB 101



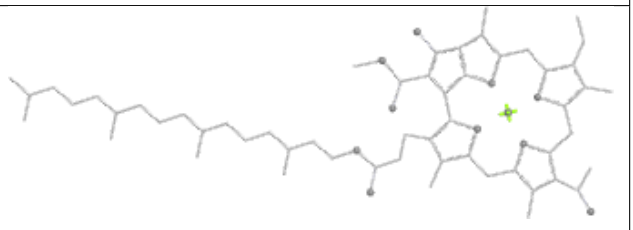
Bond lengths



Bond angles

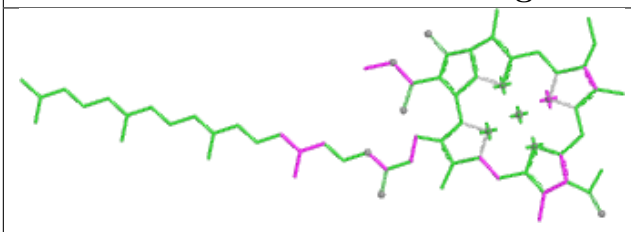


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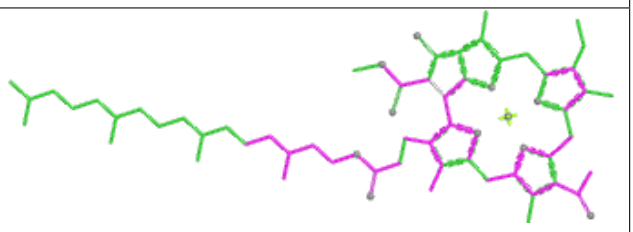


Rings

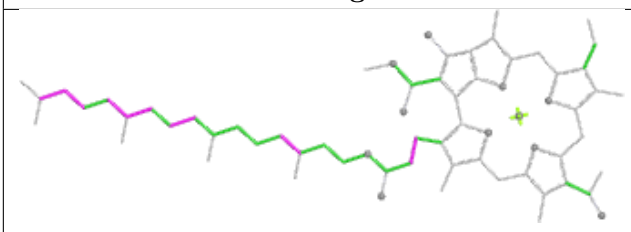
Ligand BCL AJ 101



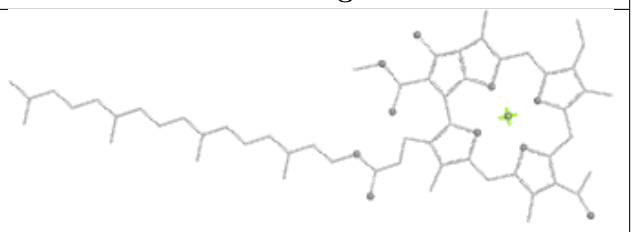
Bond lengths



Bond angles

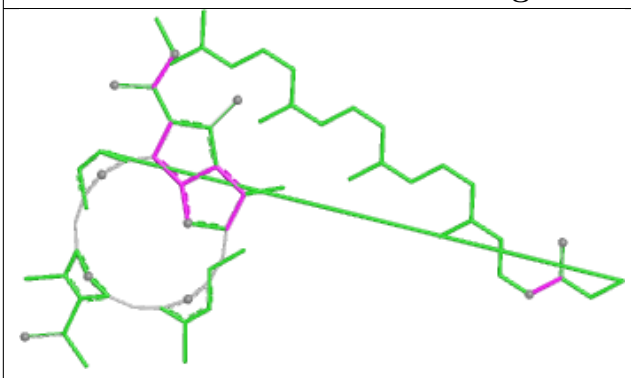


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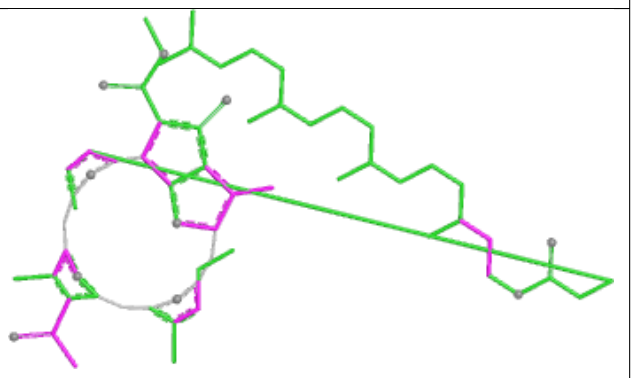


Rings

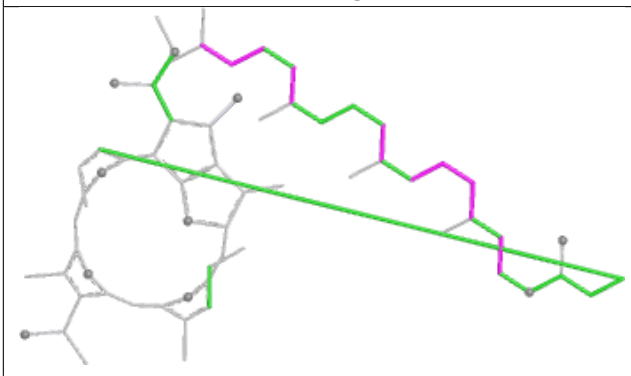
Ligand BPH BL 302



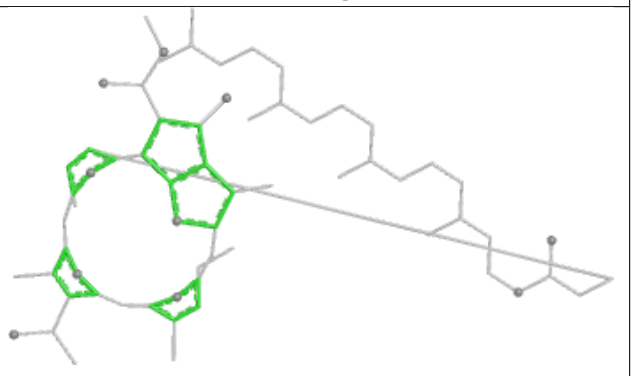
Bond lengths



Bond angles

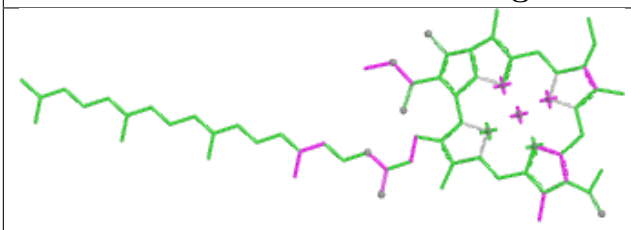


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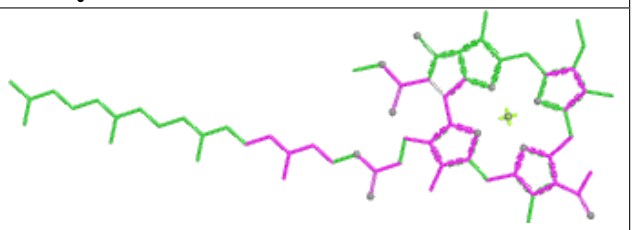


Rings

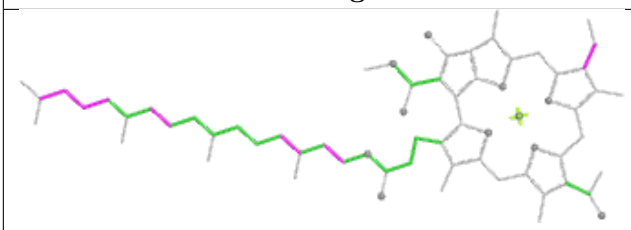
Ligand BCL BQ 104



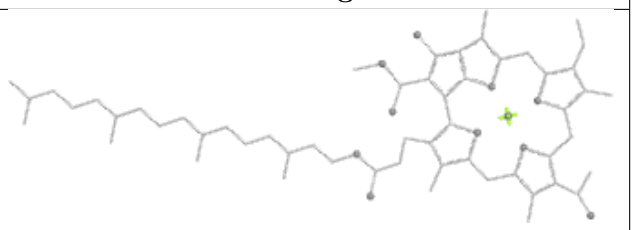
Bond lengths



Bond angles

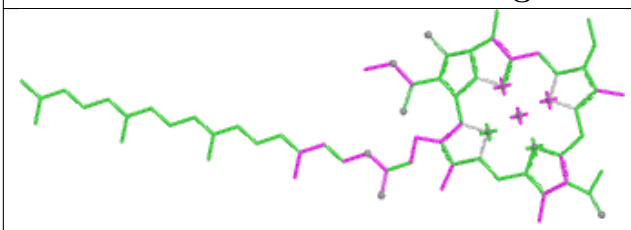


Torsions

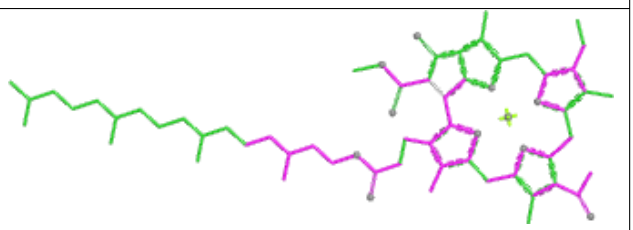


Rings

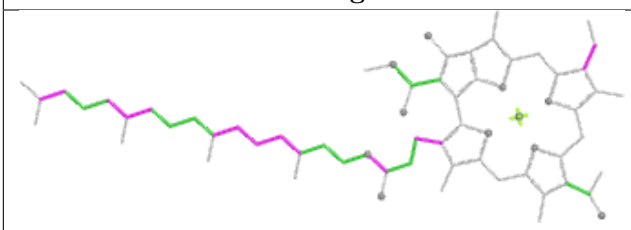
Ligand BCL AZ 101



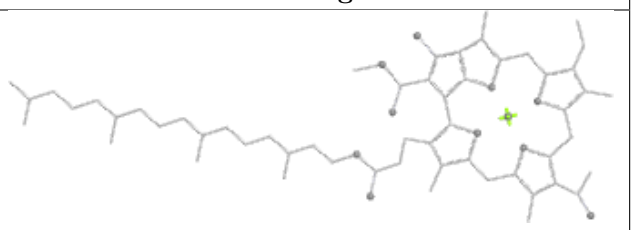
Bond lengths



Bond angles

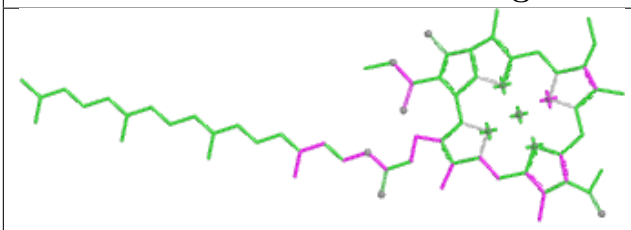


Torsions

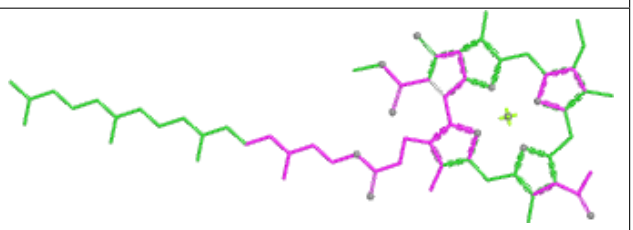


Rings

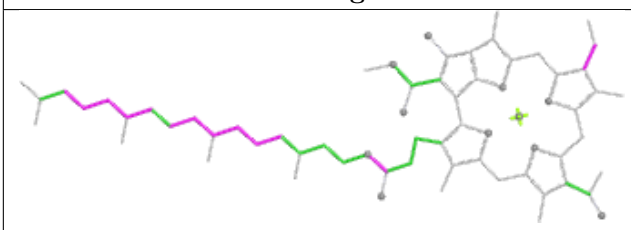
Ligand BCL AY 102



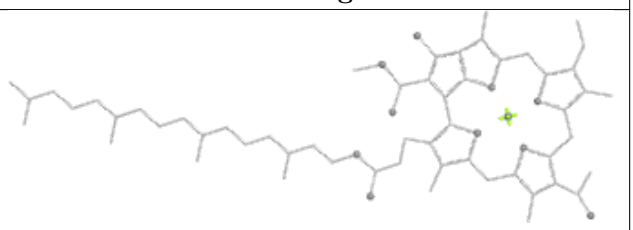
Bond lengths



Bond angles

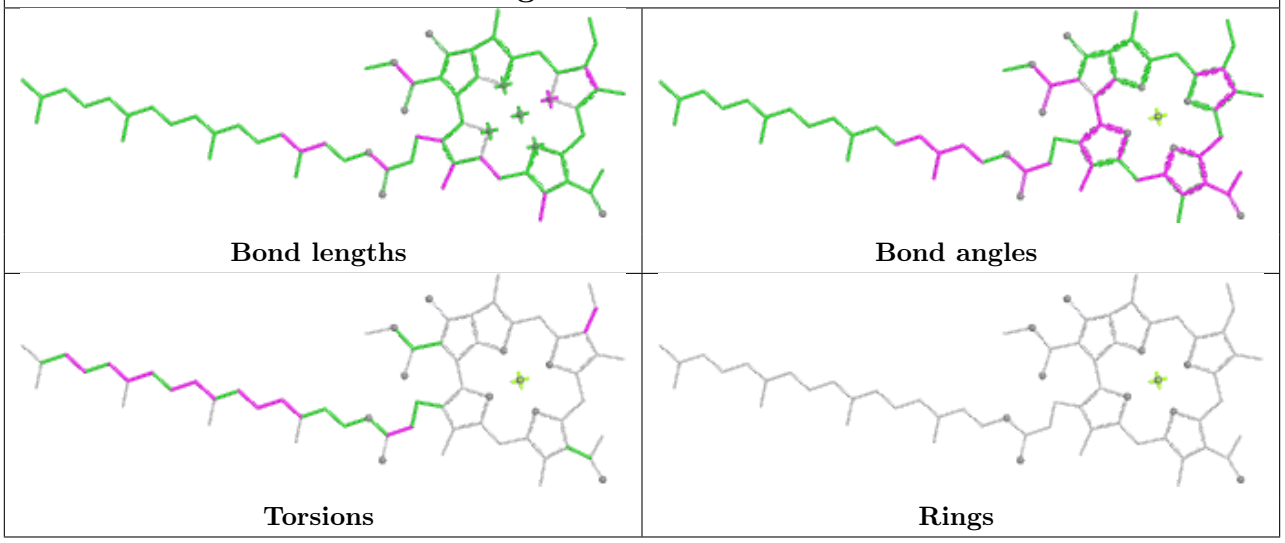


Torsions

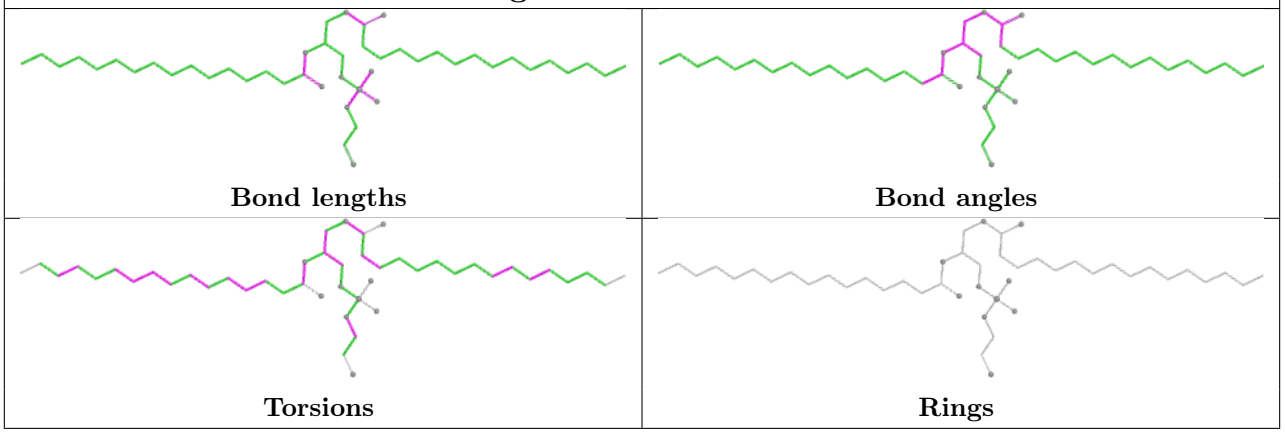


Rings

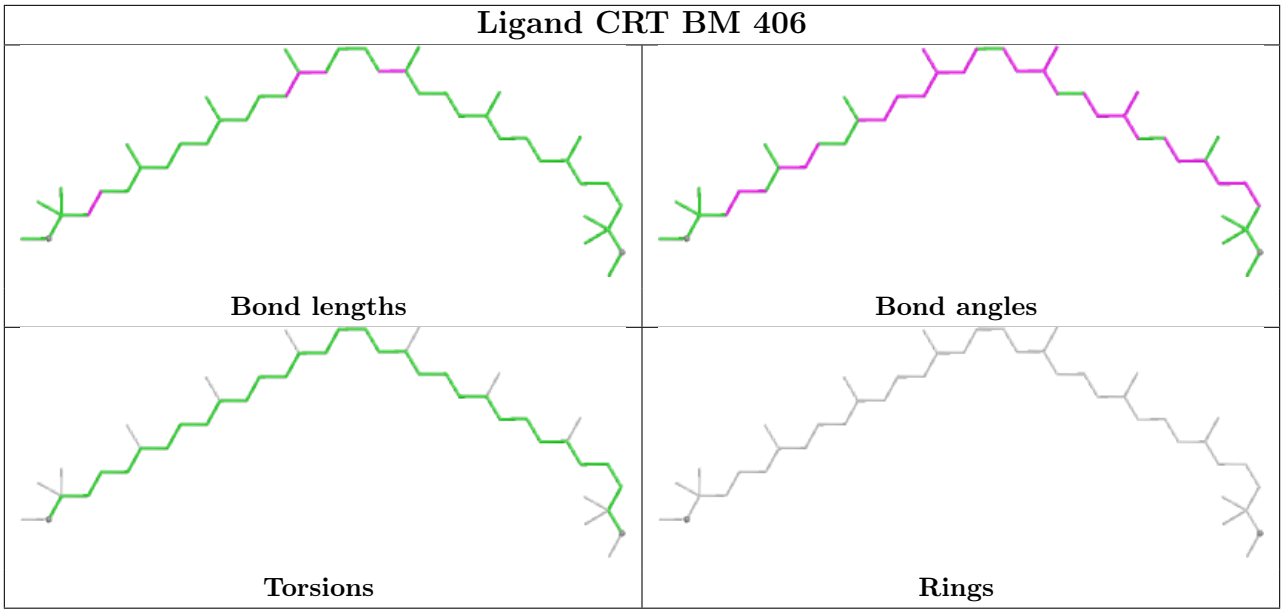
Ligand BCL B3 102



Ligand PEF AM 409



Ligand CRT BM 406



5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled '#RSRZ> 2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95^{th} percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled 'Q< 0.9' lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å^2)	Q<0.9
1	AC	317/404 (78%)	-0.60	4 (1%)	75	53	5, 84, 141, 197	1 (0%)
1	BC	317/404 (78%)	-0.48	3 (0%)	81	61	8, 91, 139, 165	1 (0%)
2	AL	280/281 (99%)	-0.66	1 (0%)	88	76	21, 58, 119, 145	0
2	BL	280/281 (99%)	-0.60	1 (0%)	88	76	33, 77, 136, 156	0
3	AM	319/325 (98%)	-0.61	2 (0%)	85	69	19, 65, 107, 121	0
3	BM	319/325 (98%)	-0.67	0	100	100	37, 80, 125, 184	0
4	AH	258/259 (99%)	-0.55	4 (1%)	70	47	46, 95, 149, 183	0
4	BH	258/259 (99%)	-0.57	1 (0%)	88	76	57, 104, 164, 183	0
5	A1	58/61 (95%)	-0.37	3 (5%)	33	17	73, 163, 300, 305	0
5	A3	57/61 (93%)	-0.55	0	100	100	117, 162, 318, 320	0
5	A5	56/61 (91%)	-0.40	3 (5%)	31	16	70, 165, 320, 321	0
5	A7	51/61 (83%)	-0.35	0	100	100	111, 148, 234, 251	0
5	A9	60/61 (98%)	-0.39	0	100	100	96, 151, 319, 319	0
5	AA	48/61 (78%)	-0.52	0	100	100	92, 144, 241, 257	0
5	AD	57/61 (93%)	-0.12	4 (7%)	22	12	97, 145, 222, 239	0
5	AF	59/61 (96%)	-0.31	1 (1%)	69	45	99, 130, 226, 235	0
5	AI	59/61 (96%)	-0.13	3 (5%)	33	17	81, 143, 236, 267	0
5	AK	58/61 (95%)	-0.57	0	100	100	81, 140, 226, 261	0
5	AO	59/61 (96%)	-0.19	2 (3%)	48	27	88, 158, 258, 262	0
5	AQ	57/61 (93%)	-0.56	2 (3%)	47	27	56, 135, 277, 280	0
5	AS	59/61 (96%)	-0.24	3 (5%)	33	17	86, 158, 300, 309	0
5	AU	60/61 (98%)	-0.04	4 (6%)	24	12	144, 167, 252, 254	0
5	AW	60/61 (98%)	-0.34	4 (6%)	24	12	68, 135, 239, 250	0
5	AY	60/61 (98%)	-0.24	2 (3%)	49	28	128, 152, 278, 284	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å^2)	Q<0.9
5	B1	54/61 (88%)	-0.45	0 100 100	88, 126, 237, 238	0
5	B3	60/61 (98%)	-0.22	2 (3%) 49 28	103, 151, 262, 263	0
5	B5	51/61 (83%)	-0.57	0 100 100	131, 174, 234, 237	0
5	B7	54/61 (88%)	-0.34	1 (1%) 66 43	120, 191, 255, 261	0
5	B9	51/61 (83%)	-0.58	0 100 100	101, 150, 240, 241	0
5	BA	55/61 (90%)	-0.30	3 (5%) 30 15	112, 161, 261, 269	0
5	BD	45/61 (73%)	-0.08	3 (6%) 24 12	135, 140, 226, 247	0
5	BF	56/61 (91%)	-0.26	4 (7%) 22 11	135, 168, 237, 251	0
5	BI	50/61 (81%)	-0.57	1 (2%) 65 41	107, 134, 223, 229	0
5	BK	60/61 (98%)	-0.35	2 (3%) 49 28	152, 166, 314, 318	0
5	BO	59/61 (96%)	-0.66	1 (1%) 69 45	76, 129, 292, 295	0
5	BQ	59/61 (96%)	-0.29	2 (3%) 48 27	150, 168, 266, 274	0
5	BS	59/61 (96%)	-0.46	2 (3%) 48 27	91, 159, 250, 253	0
5	BU	58/61 (95%)	-0.21	1 (1%) 69 45	109, 150, 280, 282	0
5	BW	58/61 (95%)	-0.42	1 (1%) 69 45	49, 114, 230, 232	0
5	BY	54/61 (88%)	-0.44	0 100 100	46, 95, 222, 230	0
6	A0	40/47 (85%)	-0.35	0 100 100	166, 177, 205, 220	0
6	A2	40/47 (85%)	-0.59	0 100 100	122, 146, 202, 211	0
6	A4	40/47 (85%)	-0.60	2 (5%) 34 18	147, 151, 221, 222	0
6	A6	40/47 (85%)	-0.63	1 (2%) 58 35	140, 155, 199, 213	0
6	A8	40/47 (85%)	-0.35	3 (7%) 20 10	129, 187, 225, 229	0
6	AB	40/47 (85%)	-0.39	1 (2%) 58 35	122, 162, 189, 190	0
6	AE	40/47 (85%)	-0.30	2 (5%) 34 18	120, 145, 169, 184	0
6	AG	40/47 (85%)	-0.63	0 100 100	74, 116, 140, 145	0
6	AJ	40/47 (85%)	-0.59	1 (2%) 58 35	118, 129, 154, 159	0
6	AN	40/47 (85%)	-0.69	1 (2%) 58 35	101, 121, 155, 161	0
6	AP	40/47 (85%)	-0.63	1 (2%) 58 35	106, 142, 245, 248	0
6	AR	40/47 (85%)	-0.30	2 (5%) 34 18	122, 157, 194, 199	0
6	AT	40/47 (85%)	-0.52	0 100 100	119, 148, 198, 205	0
6	AV	40/47 (85%)	-0.30	1 (2%) 58 35	130, 167, 222, 225	0
6	AX	40/47 (85%)	-0.23	0 100 100	157, 167, 194, 199	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å^2)	Q<0.9
6	AZ	40/47 (85%)	-0.52	0 100 100	111, 137, 238, 240	0
6	B0	40/47 (85%)	-0.35	2 (5%) 34 18	197, 208, 220, 221	0
6	B2	40/47 (85%)	-0.90	0 100 100	116, 127, 158, 162	0
6	B4	40/47 (85%)	-0.48	2 (5%) 34 18	134, 156, 191, 199	0
6	B6	40/47 (85%)	-0.73	0 100 100	115, 152, 213, 215	0
6	B8	40/47 (85%)	-0.59	0 100 100	123, 199, 232, 234	0
6	BB	40/47 (85%)	-0.78	0 100 100	155, 164, 228, 231	0
6	BE	40/47 (85%)	-0.73	0 100 100	148, 170, 194, 212	0
6	BG	40/47 (85%)	-0.13	4 (10%) 12 7	185, 204, 218, 220	0
6	BJ	40/47 (85%)	-0.56	0 100 100	199, 204, 208, 209	0
6	BN	40/47 (85%)	-0.66	0 100 100	152, 160, 194, 200	0
6	BP	40/47 (85%)	-0.90	1 (2%) 58 35	124, 149, 226, 230	0
6	BR	40/47 (85%)	-0.25	4 (10%) 12 7	138, 171, 206, 213	0
6	BT	40/47 (85%)	-0.64	0 100 100	132, 157, 227, 233	0
6	BV	40/47 (85%)	-0.75	1 (2%) 58 35	91, 143, 170, 174	0
6	BX	40/47 (85%)	-0.77	0 100 100	100, 127, 169, 172	0
6	BZ	40/47 (85%)	-0.72	0 100 100	121, 135, 178, 186	0
All	All	5429/5994 (90%)	-0.50	99 (1%) 67 44	5, 123, 236, 321	2 (0%)

All (99) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
5	BD	41	SER	8.9
5	AI	13	LEU	7.4
5	BS	5	ASN	5.8
5	AW	46	TRP	5.3
5	AS	48	ASP	5.0
6	BR	45	TRP	5.0
5	BQ	2	PHE	4.6
6	A8	19	ALA	4.5
6	BG	41	LEU	4.5
5	BD	37	MET	4.5
5	BW	40	LEU	4.4
5	AD	40	LEU	4.2
5	AD	39	VAL	4.2
5	A1	13	LEU	4.1

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Mol	Chain	Res	Type	RSRZ
1	AC	17	SER	3.9
5	AD	41	SER	3.9
5	B7	10	LYS	3.9
5	BD	40	LEU	3.9
5	B3	6	ALA	3.8
6	A4	45	TRP	3.7
5	AF	13	LEU	3.7
5	AW	47	LEU	3.6
6	BR	39	ALA	3.6
5	AU	39	VAL	3.6
5	BA	41	SER	3.5
6	BG	39	ALA	3.5
5	AO	57	ALA	3.4
5	AY	13	LEU	3.4
5	A5	39	VAL	3.4
6	AV	45	TRP	3.4
1	AC	171	GLY	3.3
5	A1	47	LEU	3.3
5	A5	5	ASN	3.2
6	B4	19	ALA	3.2
5	BF	13	LEU	3.2
5	BO	54	SER	3.2
5	AU	38	ILE	3.1
5	BF	44	LEU	3.1
5	B3	51	ILE	3.0
5	BA	6	ALA	3.0
5	BK	58	LEU	3.0
5	AW	40	LEU	2.9
5	A5	8	LEU	2.9
2	BL	151	TRP	2.9
5	AS	46	TRP	2.9
5	AI	14	ILE	2.8
5	A1	14	ILE	2.8
5	AD	42	THR	2.8
6	AN	16	GLU	2.7
6	BG	42	TYR	2.7
6	AR	11	ASP	2.7
5	BS	8	LEU	2.6
6	A8	21	PHE	2.6
4	AH	47	GLU	2.6
5	BK	57	ALA	2.6
6	A4	39	ALA	2.6

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Mol	Chain	Res	Type	RSRZ
6	B0	39	ALA	2.6
5	BA	42	THR	2.6
4	AH	39	TYR	2.6
4	AH	151	PRO	2.5
5	BQ	8	LEU	2.5
1	BC	38	VAL	2.5
5	AU	43	ASP	2.5
1	BC	181	THR	2.5
3	AM	289	THR	2.5
5	AW	39	VAL	2.5
5	AI	6	ALA	2.4
5	AQ	42	THR	2.4
3	AM	290	VAL	2.4
6	A8	17	PHE	2.4
5	BF	43	ASP	2.4
6	BP	11	ASP	2.4
6	AR	40	TRP	2.4
6	BR	40	TRP	2.4
6	AJ	39	ALA	2.4
6	BR	44	PRO	2.4
2	AL	154	GLY	2.3
5	BI	8	LEU	2.3
6	AB	28	TRP	2.2
6	AE	45	TRP	2.2
6	AP	11	ASP	2.2
6	A6	16	GLU	2.2
6	BV	12	ASP	2.2
6	AE	39	ALA	2.2
5	AO	8	LEU	2.2
6	B0	42	TYR	2.2
5	BF	45	ASN	2.2
6	B4	16	GLU	2.1
1	AC	58	PRO	2.1
4	BH	56	VAL	2.1
5	AS	47	LEU	2.1
5	BU	40	LEU	2.1
1	AC	191	ALA	2.1
1	BC	59	VAL	2.0
4	AH	56	VAL	2.0
5	AY	40	LEU	2.0
5	AU	46	TRP	2.0
5	AQ	13	LEU	2.0

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Mol	Chain	Res	Type	RSRZ
6	BG	9	LEU	2.0

6.2 Non-standard residues in protein, DNA, RNA chains (i)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates (i)

There are no oligosaccharides in this entry.

6.4 Ligands (i)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
8	CA	AO	101	1/1	0.94	0.07	267,267,267,267	0
16	PO4	A3	101	5/5	0.94	0.08	153,154,154,154	0
8	CA	B9	101	1/1	0.95	0.05	150,150,150,150	0
14	CRT	AN	102	44/44	0.95	0.15	133,135,140,140	0
14	CRT	A0	101	44/44	0.95	0.17	179,184,194,195	0
14	CRT	BV	102	44/44	0.95	0.19	208,214,218,220	0
15	PEF	BQ	101	47/47	0.95	0.17	121,145,161,161	0
8	CA	AS	102	1/1	0.95	0.06	190,190,190,190	0
14	CRT	AB	102	44/44	0.96	0.14	170,172,177,178	0
8	CA	A5	101	1/1	0.96	0.04	255,255,255,255	0
9	BCL	AM	402	66/66	0.96	0.11	40,48,81,101	0
14	CRT	BG	102	44/44	0.96	0.10	122,130,143,145	0
14	CRT	BP	102	44/44	0.96	0.13	149,151,153,154	0
9	BCL	A3	103	66/66	0.96	0.14	127,137,178,178	0
14	CRT	B1	103	44/44	0.96	0.15	205,224,238,240	0
9	BCL	BQ	104	66/66	0.96	0.14	181,191,253,254	0
11	UQ8	AL	304	53/53	0.96	0.14	74,107,143,151	0
9	BCL	AR	101	66/66	0.97	0.10	187,200,230,231	0
9	BCL	A2	101	66/66	0.97	0.14	122,133,201,203	0
8	CA	AK	101	1/1	0.97	0.06	216,216,216,216	0
9	BCL	A3	104	66/66	0.97	0.12	151,177,223,224	0
9	BCL	A6	101	66/66	0.97	0.12	238,254,268,272	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
9	BCL	A8	101	66/66	0.97	0.09	201,214,245,247	0
9	BCL	BL	301	66/66	0.97	0.09	36,53,94,100	0
9	BCL	BM	402	66/66	0.97	0.11	41,58,73,79	0
9	BCL	BA	101	66/66	0.97	0.10	206,224,237,238	0
9	BCL	BD	102	66/66	0.97	0.11	147,182,237,238	0
9	BCL	BF	102	66/66	0.97	0.12	143,168,206,210	0
9	BCL	BI	102	66/66	0.97	0.10	131,159,208,208	0
9	BCL	BJ	101	66/66	0.97	0.10	202,204,207,208	0
9	BCL	BP	101	66/66	0.97	0.09	156,173,208,209	0
8	CA	BD	101	1/1	0.97	0.04	236,236,236,236	0
9	BCL	BU	102	66/66	0.97	0.12	144,161,240,240	0
9	BCL	BW	102	66/66	0.97	0.15	129,145,213,213	0
9	BCL	BX	101	66/66	0.97	0.11	126,139,204,205	0
9	BCL	BZ	101	66/66	0.97	0.10	125,139,192,193	0
9	BCL	B8	101	66/66	0.97	0.09	226,245,252,258	0
9	BCL	B9	102	66/66	0.97	0.11	205,229,236,237	0
10	BPH	AM	403	65/65	0.97	0.10	43,61,150,153	0
10	BPH	BL	302	65/65	0.97	0.10	36,63,98,116	0
8	CA	BI	101	1/1	0.97	0.05	236,236,236,236	0
11	UQ8	BL	304	53/53	0.97	0.12	66,137,159,166	0
13	MQ8	AM	405	53/53	0.97	0.12	76,93,128,137	0
14	CRT	AA	102	44/44	0.97	0.13	164,164,167,167	0
8	CA	B7	101	1/1	0.97	0.05	259,259,259,259	0
14	CRT	AG	102	44/44	0.97	0.14	126,127,135,138	0
14	CRT	AJ	102	44/44	0.97	0.12	138,140,144,147	0
8	CA	AA	103	1/1	0.97	0.07	275,275,275,275	0
14	CRT	AP	102	44/44	0.97	0.13	137,140,147,148	0
14	CRT	AS	104	44/44	0.97	0.15	165,170,177,180	0
14	CRT	AW	102	44/44	0.97	0.16	127,136,144,147	0
14	CRT	AX	102	44/44	0.97	0.14	218,225,227,229	0
14	CRT	A5	103	44/44	0.97	0.16	167,174,185,187	0
14	CRT	A7	102	44/44	0.97	0.11	126,137,150,152	0
9	BCL	AL	301	66/66	0.97	0.10	38,50,92,96	0
14	CRT	BA	102	44/44	0.97	0.13	143,157,171,174	0
14	CRT	BB	102	44/44	0.97	0.15	169,185,201,205	0
14	CRT	BF	103	44/44	0.97	0.11	134,141,147,149	0
8	CA	AD	101	1/1	0.97	0.06	238,238,238,238	0
14	CRT	BN	102	44/44	0.97	0.11	134,135,140,141	0
14	CRT	BO	103	44/44	0.97	0.15	179,182,185,186	0
9	BCL	AB	101	66/66	0.97	0.11	124,158,224,224	0
14	CRT	BU	103	44/44	0.97	0.15	161,173,176,176	0
9	BCL	AE	101	66/66	0.97	0.12	135,156,211,212	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
9	BCL	AJ	101	66/66	0.97	0.11	134,165,191,193	0
14	CRT	B5	103	44/44	0.97	0.17	178,198,221,226	0
14	CRT	B0	101	44/44	0.97	0.19	202,220,238,242	0
15	PEF	AM	409	47/47	0.97	0.15	129,183,183,183	0
15	PEF	AS	101	47/47	0.97	0.12	178,225,225,225	0
9	BCL	AK	102	66/66	0.97	0.12	142,154,184,184	0
9	BCL	AN	101	66/66	0.97	0.10	106,126,157,171	0
9	BCL	BT	101	66/66	0.98	0.11	176,191,253,254	0
8	CA	AV	101	1/1	0.98	0.05	168,168,168,168	0
9	BCL	BV	101	66/66	0.98	0.14	172,194,218,220	0
9	BCL	AF	102	66/66	0.98	0.14	156,175,222,222	0
9	BCL	AG	101	66/66	0.98	0.11	105,132,177,178	0
9	BCL	BY	102	66/66	0.98	0.09	110,129,169,169	0
9	BCL	AI	102	66/66	0.98	0.11	148,163,190,191	0
9	BCL	B1	102	66/66	0.98	0.09	87,118,185,186	0
9	BCL	B2	101	66/66	0.98	0.11	136,152,203,204	0
9	BCL	B3	102	66/66	0.98	0.07	104,112,169,170	0
9	BCL	B4	101	66/66	0.98	0.10	149,166,212,214	0
9	BCL	B5	102	66/66	0.98	0.09	154,171,221,221	0
9	BCL	B6	101	66/66	0.98	0.12	238,251,261,264	0
9	BCL	B7	103	66/66	0.98	0.11	232,250,259,260	0
8	CA	A3	102	1/1	0.98	0.08	256,256,256,256	0
7	HEM	AC	502	43/43	0.98	0.09	53,68,81,88	0
9	BCL	B0	102	66/66	0.98	0.10	198,202,205,207	0
10	BPH	AL	302	65/65	0.98	0.09	31,54,75,84	0
8	CA	A9	101	1/1	0.98	0.04	268,268,268,268	0
8	CA	BA	103	1/1	0.98	0.05	283,283,283,283	0
10	BPH	BM	403	65/65	0.98	0.10	72,96,164,176	0
9	BCL	AS	103	66/66	0.98	0.12	206,220,229,233	0
9	BCL	AT	101	66/66	0.98	0.10	187,198,247,248	0
9	BCL	AV	102	66/66	0.98	0.11	183,194,249,250	0
13	MQ8	BM	405	53/53	0.98	0.11	68,99,166,168	0
14	CRT	AM	406	44/44	0.98	0.09	58,70,105,107	0
9	BCL	AX	101	66/66	0.98	0.12	157,171,252,255	0
9	BCL	AZ	101	66/66	0.98	0.12	120,142,217,218	0
7	HEM	AC	503	43/43	0.98	0.10	71,84,106,112	0
8	CA	AI	101	1/1	0.98	0.04	224,224,224,224	0
8	CA	BO	101	1/1	0.98	0.04	238,238,238,238	0
9	BCL	A5	102	66/66	0.98	0.11	152,159,199,200	0
14	CRT	AR	102	44/44	0.98	0.12	189,193,200,201	0
8	CA	BU	101	1/1	0.98	0.04	237,237,237,237	0
14	CRT	AT	102	44/44	0.98	0.14	200,205,209,210	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
9	BCL	A7	103	66/66	0.98	0.12	239,243,267,268	0
8	CA	BW	101	1/1	0.98	0.04	204,204,204,204	0
14	CRT	A1	103	44/44	0.98	0.10	110,123,135,138	0
14	CRT	A2	102	44/44	0.98	0.13	107,117,146,149	0
9	BCL	A9	102	66/66	0.98	0.14	213,227,251,252	0
9	BCL	A0	102	66/66	0.98	0.11	199,210,239,240	0
7	HEM	AC	504	43/43	0.98	0.09	39,76,101,104	0
9	BCL	BL	303	66/66	0.98	0.08	39,55,82,90	0
9	BCL	BM	401	66/66	0.98	0.08	33,56,137,146	0
7	HEM	BC	502	43/43	0.98	0.09	50,80,94,103	0
8	CA	AQ	101	1/1	0.98	0.05	207,207,207,207	0
9	BCL	AL	303	66/66	0.98	0.08	32,54,95,104	0
9	BCL	AM	401	66/66	0.98	0.09	19,34,93,96	0
9	BCL	BG	101	66/66	0.98	0.09	201,207,214,217	0
14	CRT	BS	103	44/44	0.98	0.12	193,197,199,200	0
7	HEM	BC	504	43/43	0.98	0.08	55,78,92,106	0
9	BCL	AA	101	66/66	0.98	0.16	235,252,257,260	0
14	CRT	BW	103	44/44	0.98	0.11	129,145,156,161	0
9	BCL	BK	102	66/66	0.98	0.10	194,257,267,269	0
14	CRT	B2	102	44/44	0.98	0.14	132,151,166,171	0
9	BCL	BN	101	66/66	0.98	0.10	153,170,192,193	0
14	CRT	B7	102	44/44	0.98	0.12	173,194,212,217	0
9	BCL	BO	102	66/66	0.98	0.09	153,180,189,193	0
15	PEF	AM	407	19/47	0.98	0.10	114,165,177,177	0
8	CA	AU	101	1/1	0.98	0.05	223,223,223,223	0
15	PEF	AH	301	19/47	0.98	0.08	128,144,161,167	0
9	BCL	BQ	103	66/66	0.98	0.13	168,191,225,226	0
15	PEF	BM	407	19/47	0.98	0.11	102,136,173,178	0
9	BCL	AD	102	66/66	0.98	0.14	179,196,260,263	0
9	BCL	BS	102	66/66	0.98	0.11	152,193,231,231	0
16	PO4	BH	301	5/5	0.98	0.07	108,109,109,110	0
8	CA	B5	101	1/1	0.99	0.03	226,226,226,226	0
9	BCL	BE	101	66/66	0.99	0.08	146,170,225,227	0
8	CA	A7	101	1/1	0.99	0.03	213,213,213,213	0
14	CRT	BM	406	44/44	0.99	0.08	69,74,96,109	0
8	CA	AF	101	1/1	0.99	0.05	254,254,254,254	0
9	BCL	AU	102	66/66	0.99	0.11	161,175,231,234	0
8	CA	BC	505	1/1	0.99	0.06	69,69,69,69	0
9	BCL	AW	101	66/66	0.99	0.10	143,191,243,243	0
7	HEM	BC	501	43/43	0.99	0.07	52,60,117,131	0
9	BCL	AY	102	66/66	0.99	0.10	133,140,195,196	0
8	CA	AC	505	1/1	0.99	0.09	30,30,30,30	0

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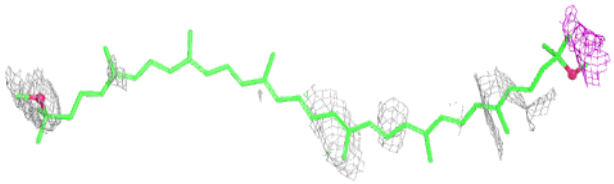
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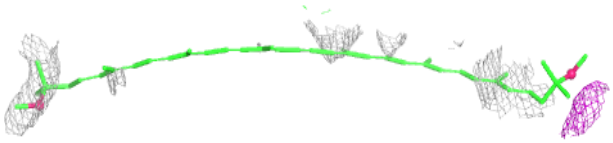
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
12	FE	AM	404	1/1	0.99	0.04	47,47,47,47	0
9	BCL	A1	102	66/66	0.99	0.10	96,135,194,194	0
8	CA	BF	101	1/1	0.99	0.03	276,276,276,276	0
8	CA	AY	101	1/1	0.99	0.06	197,197,197,197	0
8	CA	BK	101	1/1	0.99	0.05	275,275,275,275	0
8	CA	A1	101	1/1	0.99	0.05	212,212,212,212	0
8	CA	BQ	102	1/1	0.99	0.04	249,249,249,249	0
8	CA	BS	101	1/1	0.99	0.04	222,222,222,222	0
7	HEM	AC	501	43/43	0.99	0.08	86,86,86,96	0
7	HEM	BC	503	43/43	0.99	0.10	84,110,135,146	0
15	PEF	AM	408	14/47	0.99	0.07	12,90,119,125	0
8	CA	BY	101	1/1	0.99	0.03	148,148,148,148	0
8	CA	B1	101	1/1	0.99	0.06	184,184,184,184	0
8	CA	B3	101	1/1	0.99	0.07	222,222,222,222	0
9	BCL	AO	102	66/66	0.99	0.09	199,207,214,219	0
9	BCL	AP	101	66/66	0.99	0.09	107,127,189,190	0
16	PO4	AM	410	5/5	0.99	0.09	83,84,84,84	0
16	PO4	AH	302	5/5	0.99	0.04	102,103,104,104	0
9	BCL	AQ	102	66/66	0.99	0.08	187,198,210,221	0
9	BCL	BB	101	66/66	0.99	0.11	187,209,245,245	0
12	FE	BM	404	1/1	1.00	0.05	44,44,44,44	0

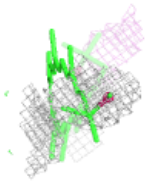
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around CRT AN 102:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

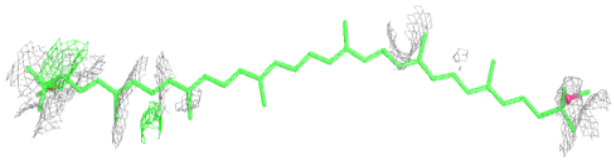


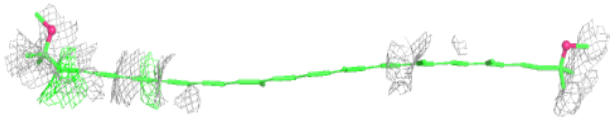




Electron density around CRT A0 101:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

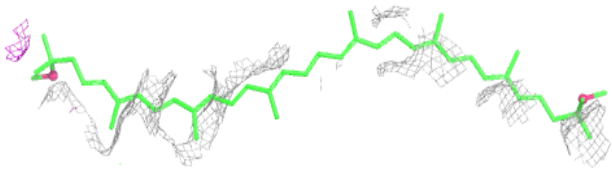


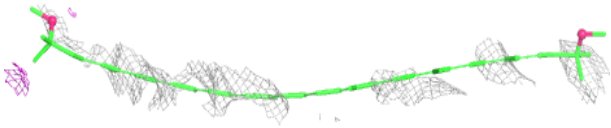




Electron density around CRT BV 102:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

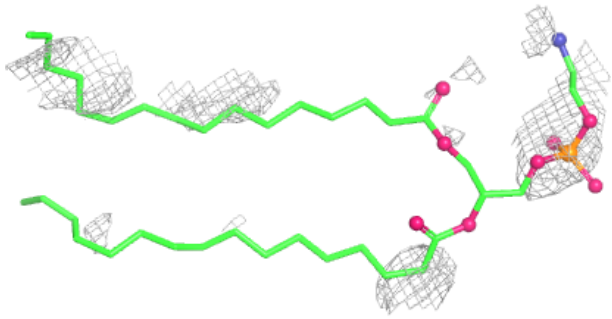


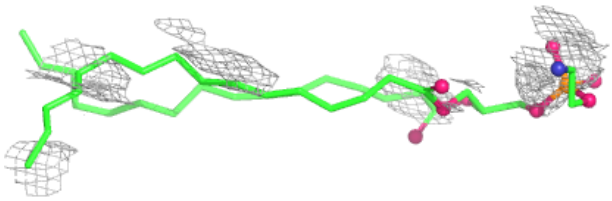


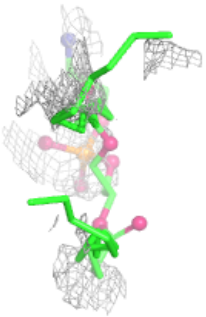


Electron density around PEF BQ 101:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

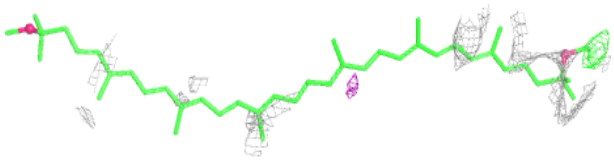




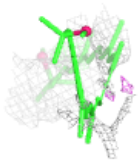


Electron density around CRT AB 102:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

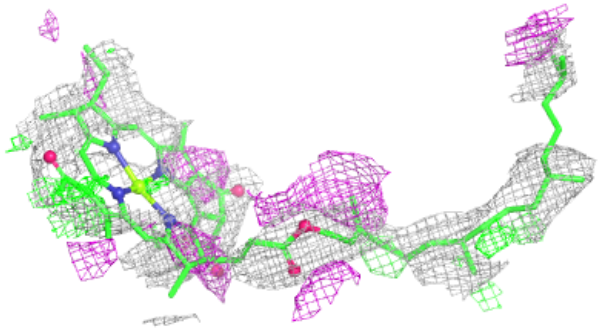


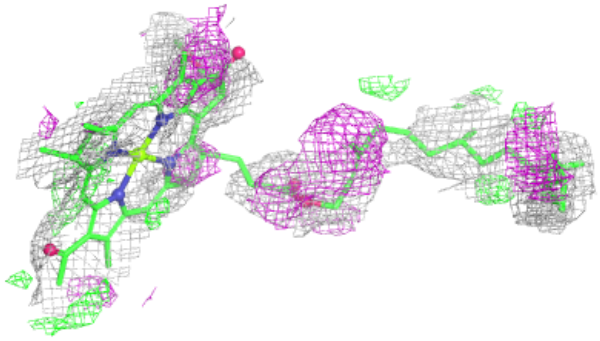


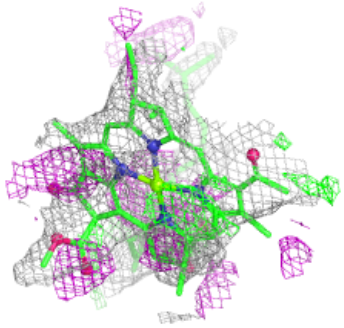


Electron density around BCL AM 402:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

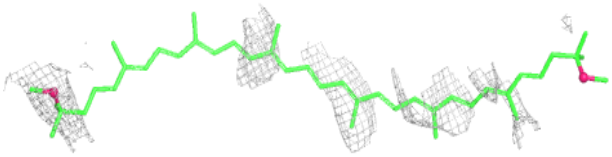


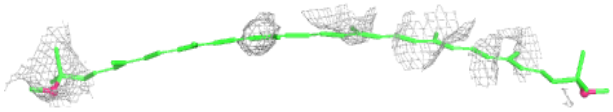


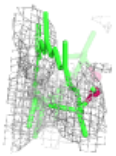


Electron density around CRT BG 102:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

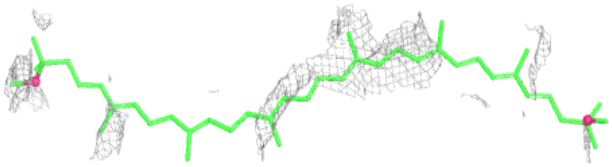


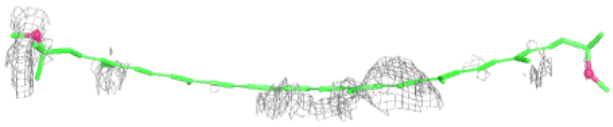


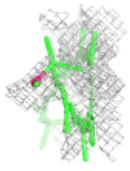


Electron density around CRT BP 102:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

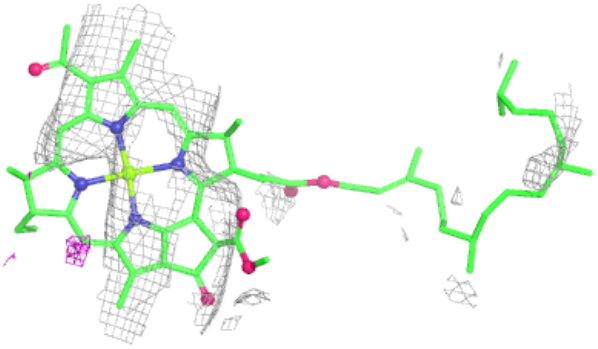


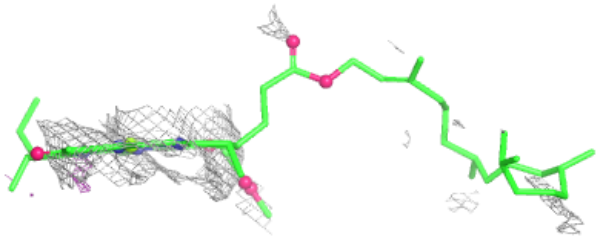


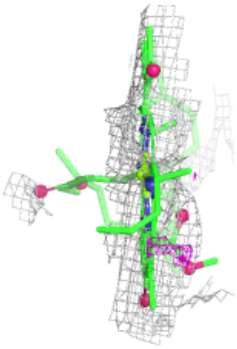


Electron density around BCL A3 103:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

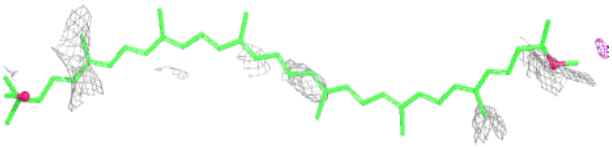


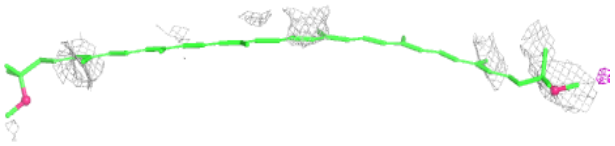


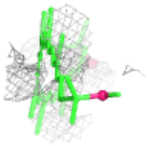


Electron density around CRT B1 103:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

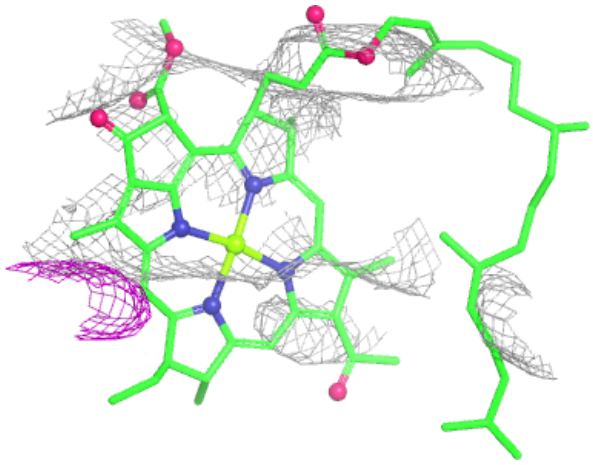


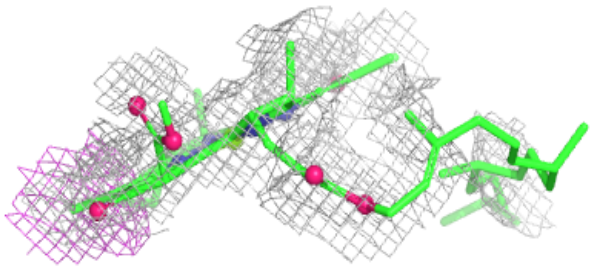


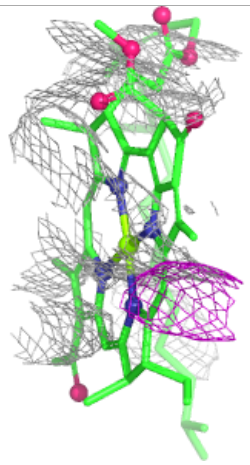


Electron density around BCL BQ 104:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

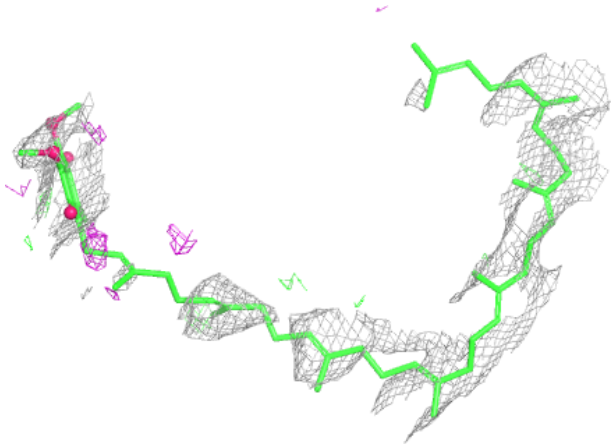


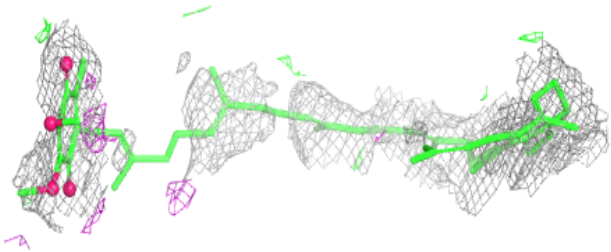


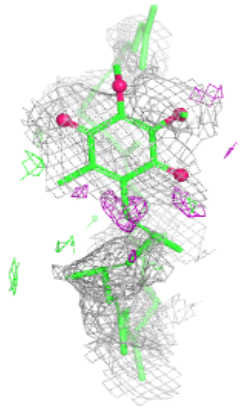


Electron density around UQ8 AL 304:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

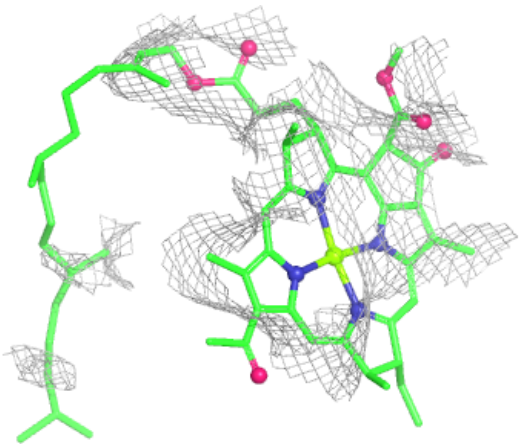


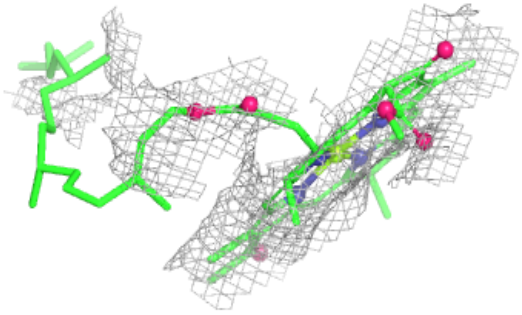


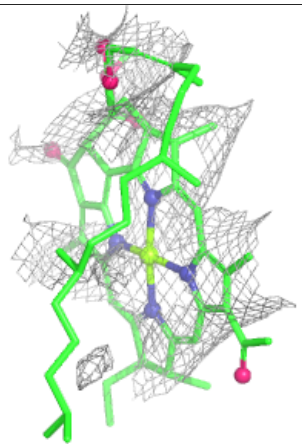


Electron density around BCL AR 101:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

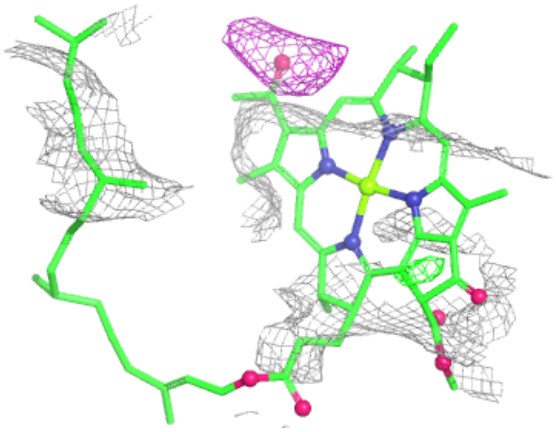


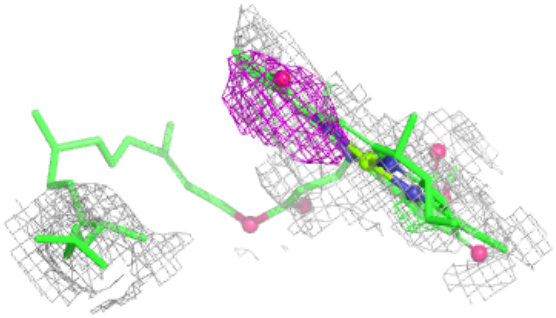


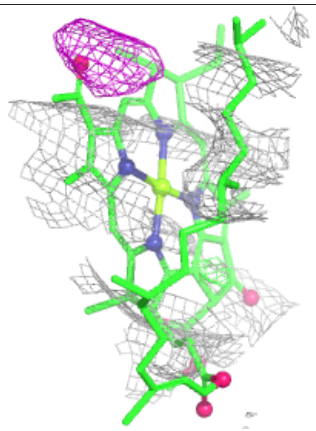


Electron density around BCL A2 101:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

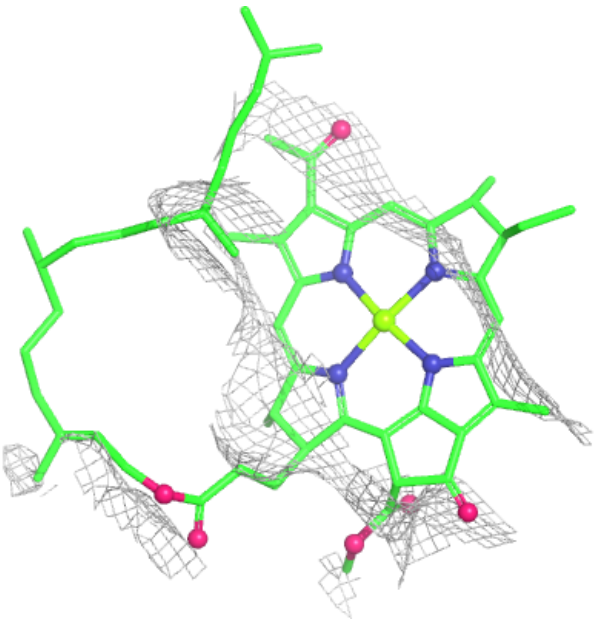


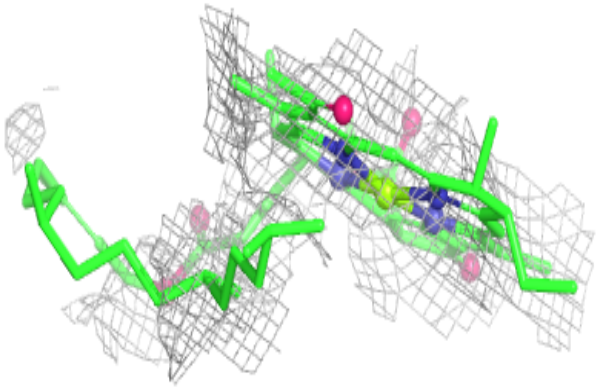


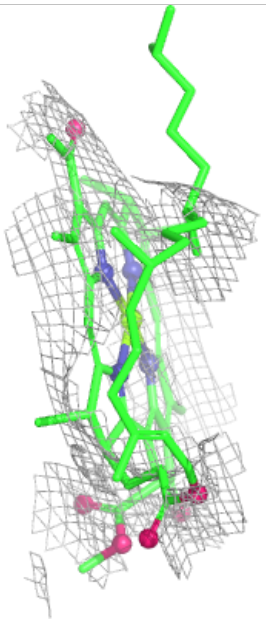


Electron density around BCL A3 104:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

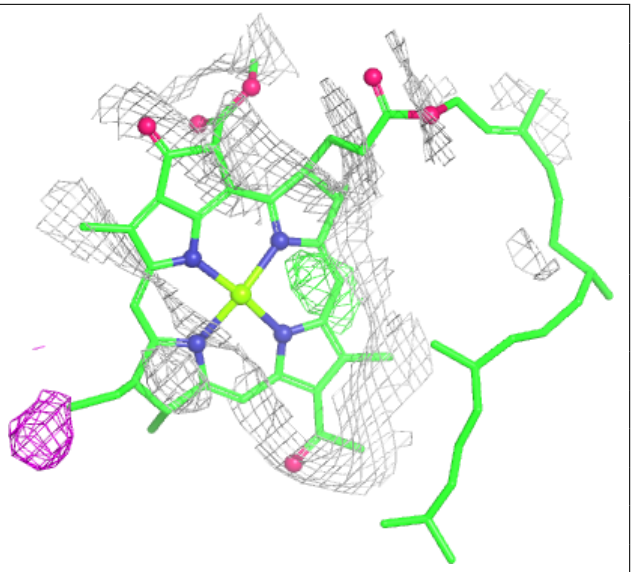


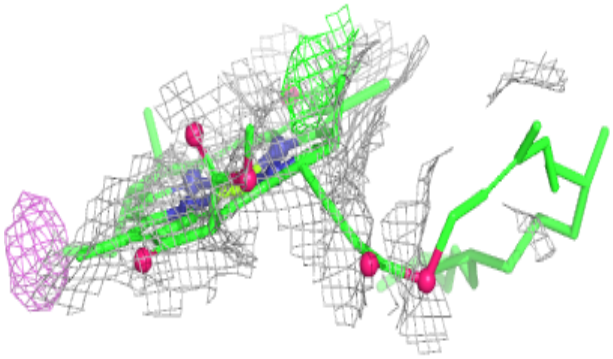


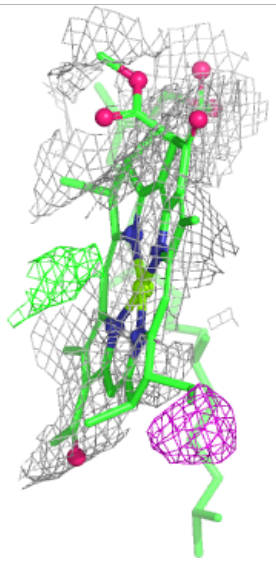


Electron density around BCL A6 101:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

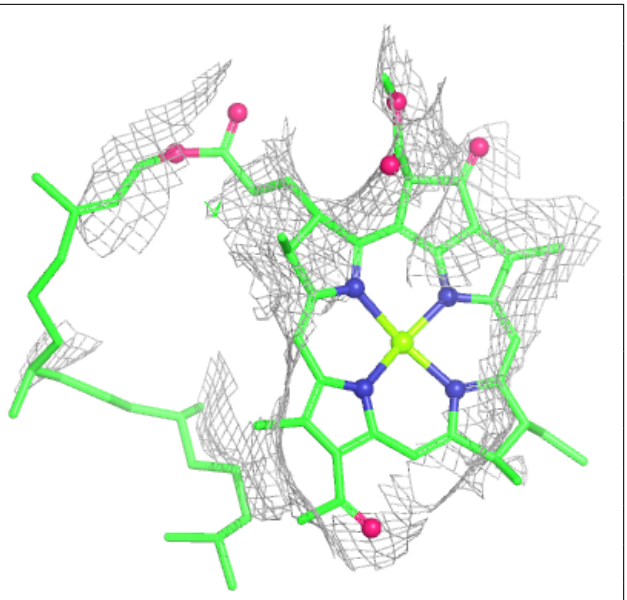


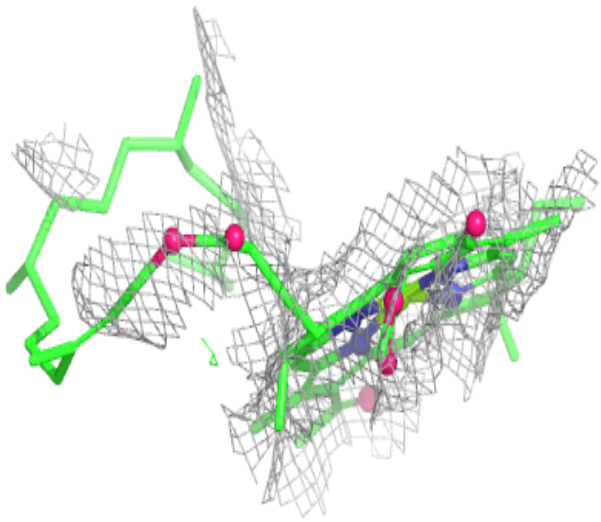


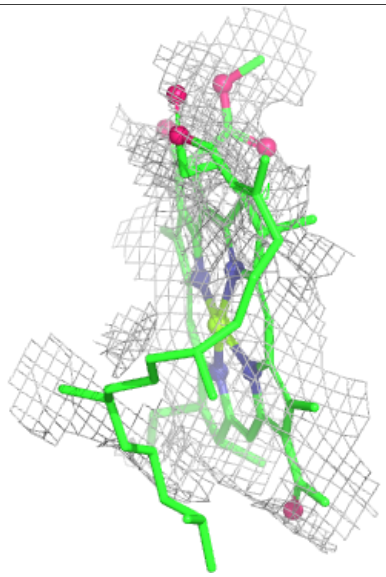


Electron density around BCL A8 101:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

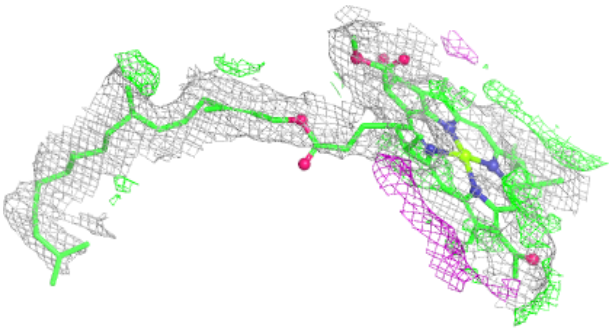


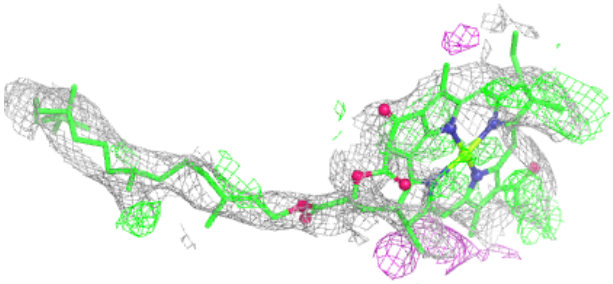


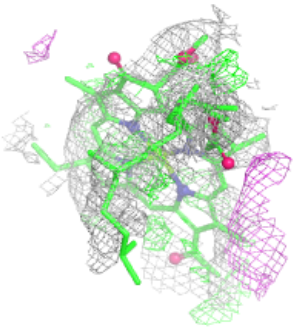


Electron density around BCL BL 301:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

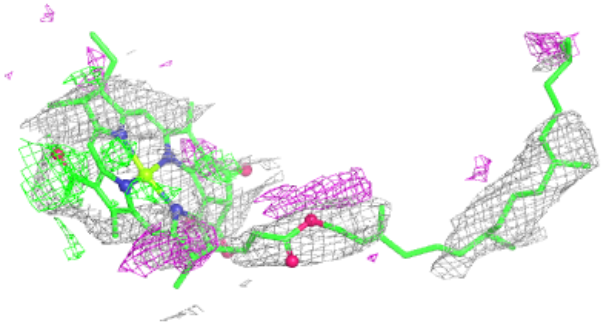


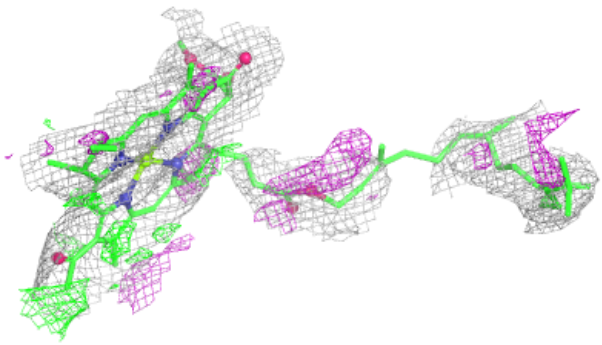


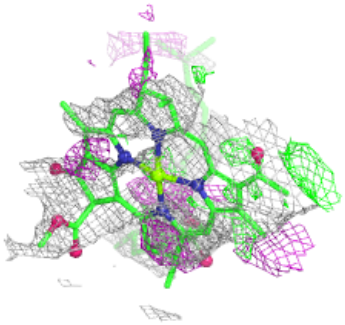


Electron density around BCL BM 402:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

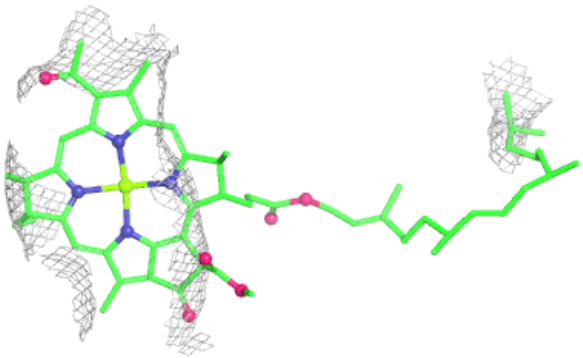


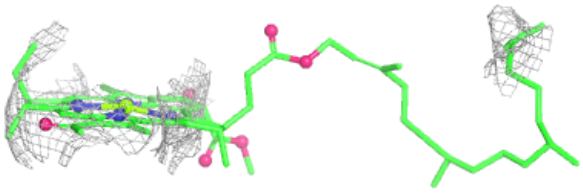


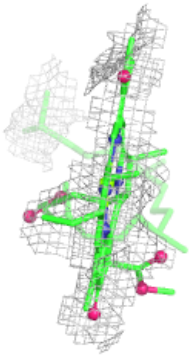


Electron density around BCL BA 101:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

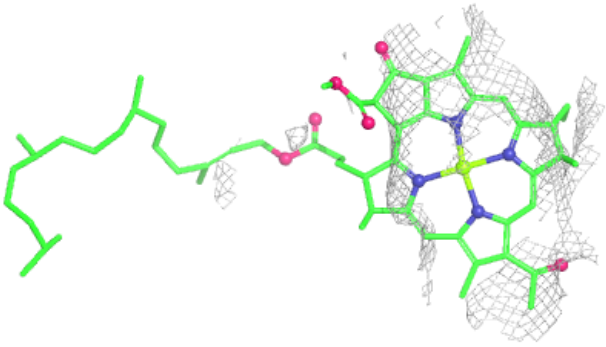


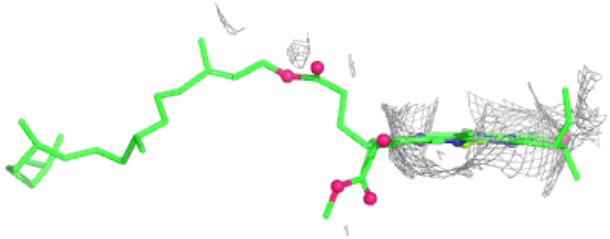




Electron density around BCL BD 102:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

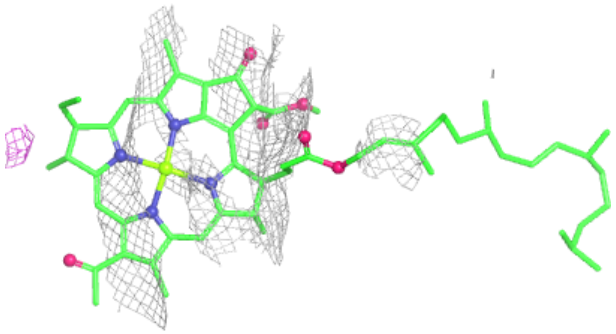


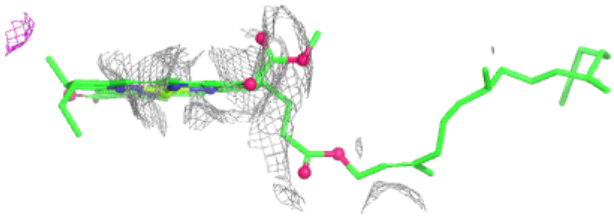


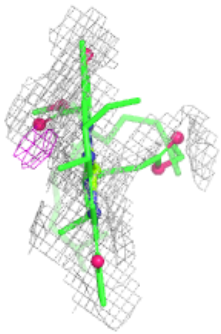


Electron density around BCL BF 102:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

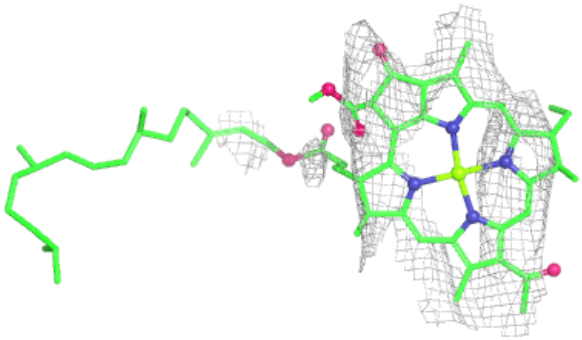


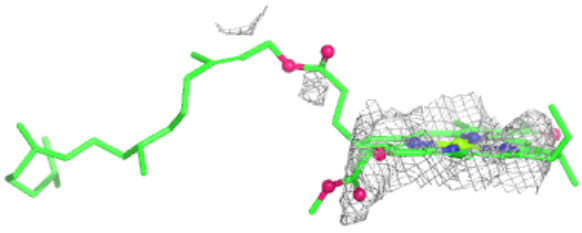


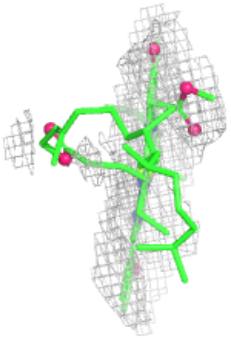


Electron density around BCL BI 102:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

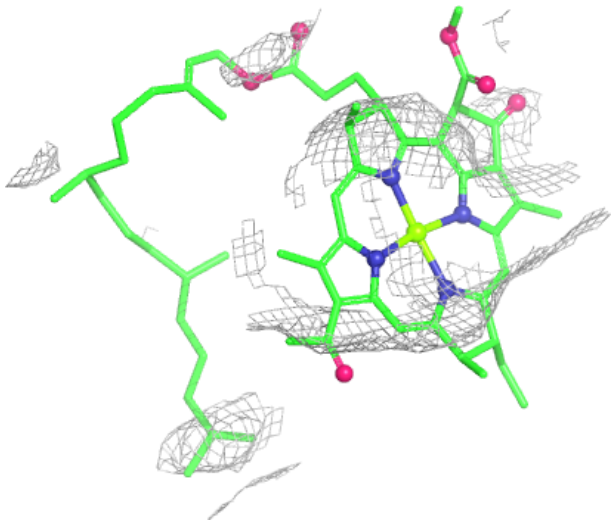


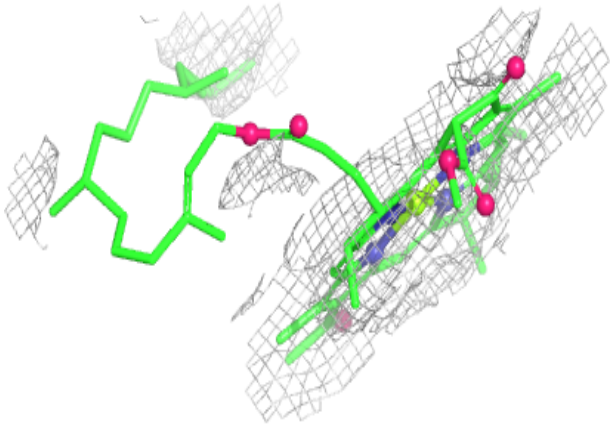


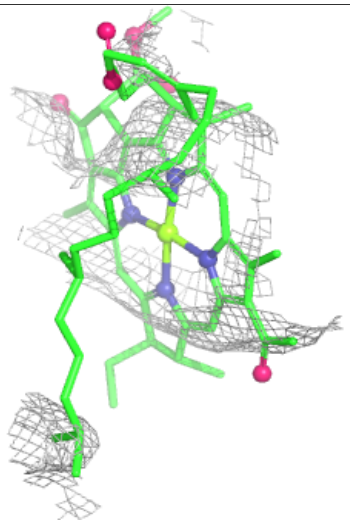


Electron density around BCL BJ 101:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

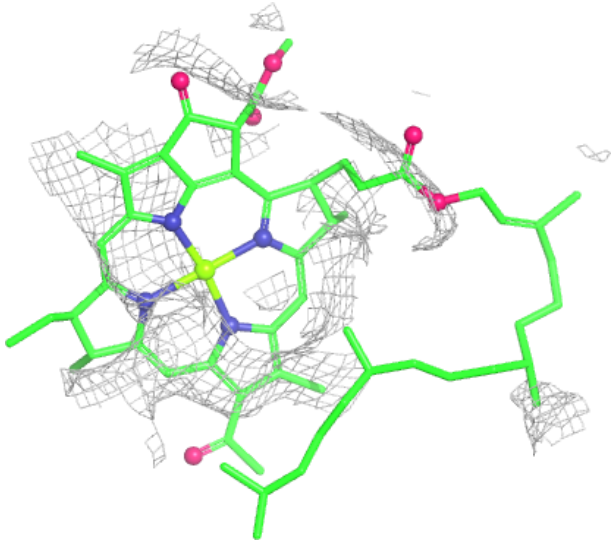


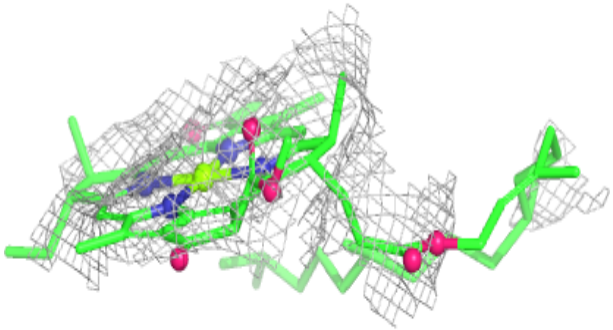




Electron density around BCL BP 101:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

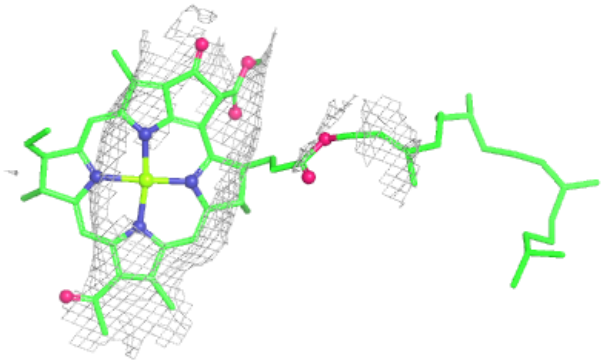


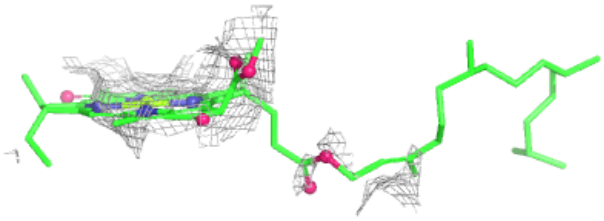


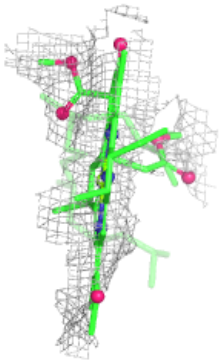


Electron density around BCL BU 102:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

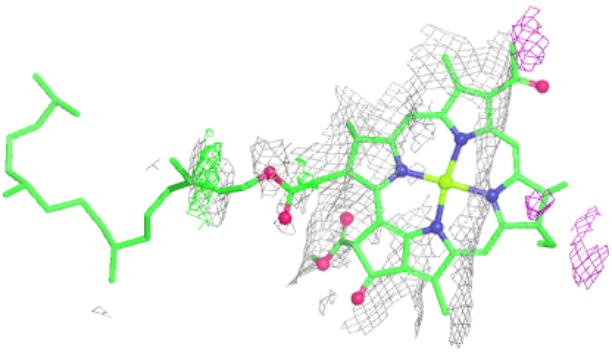


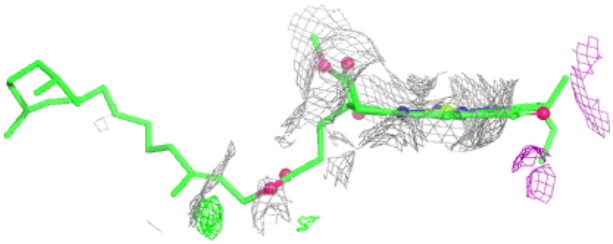


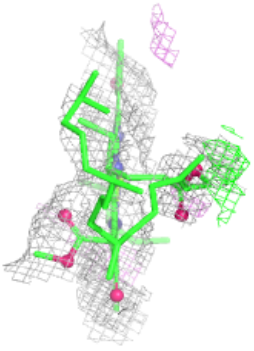


Electron density around BCL BW 102:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

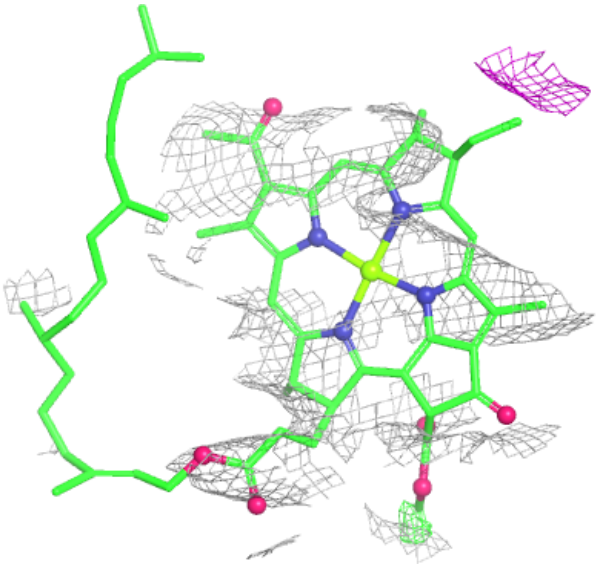


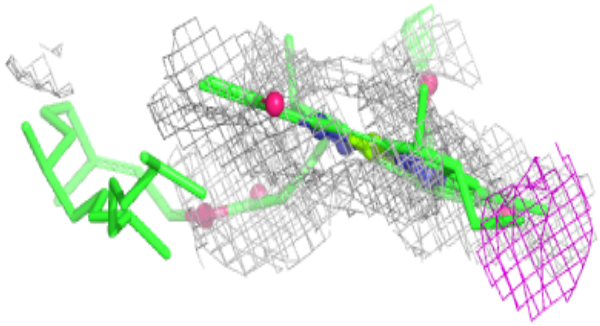


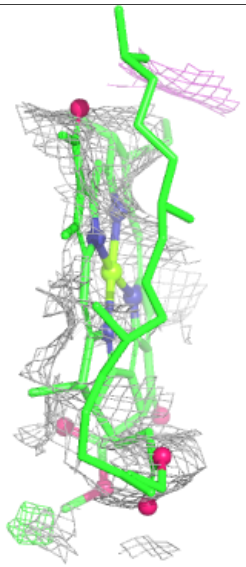


Electron density around BCL BX 101:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

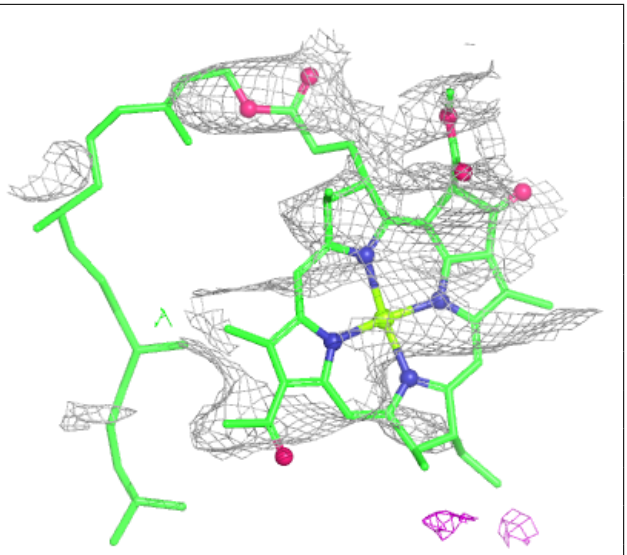


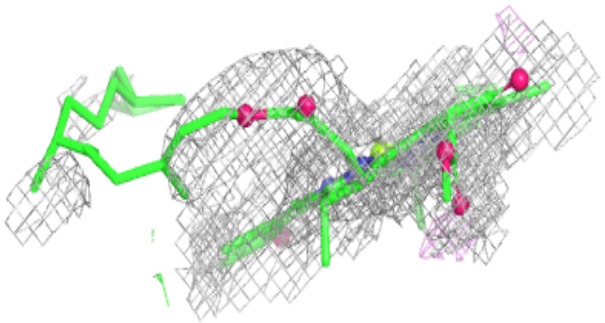


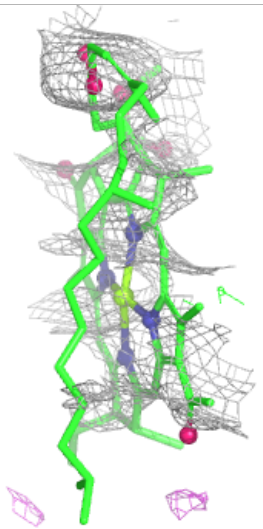


Electron density around BCL BZ 101:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

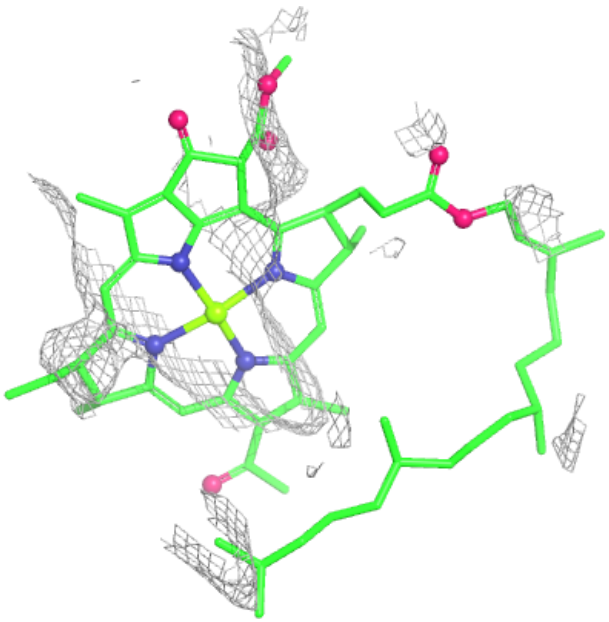


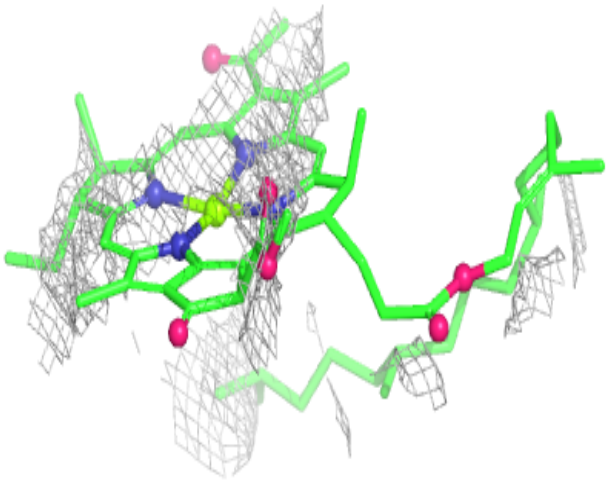


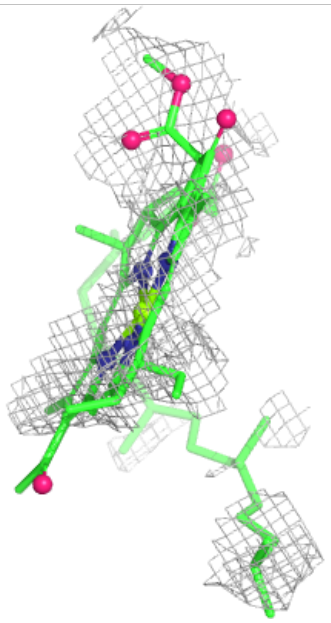


Electron density around BCL B8 101:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

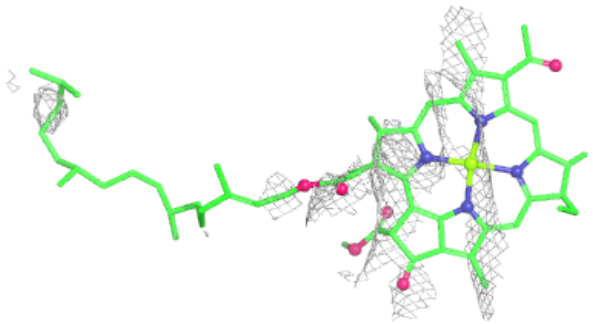


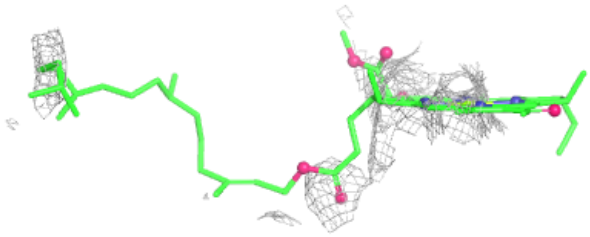


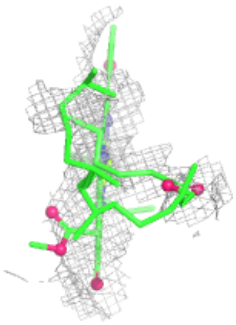


Electron density around BCL B9 102:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

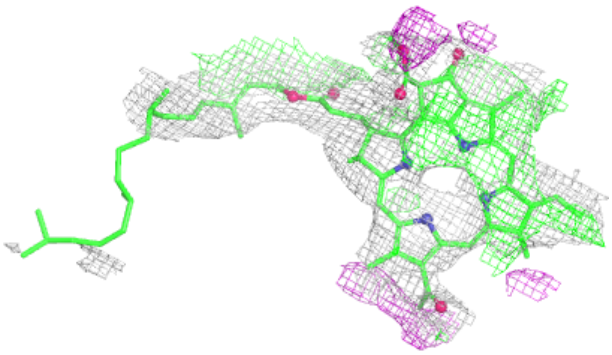


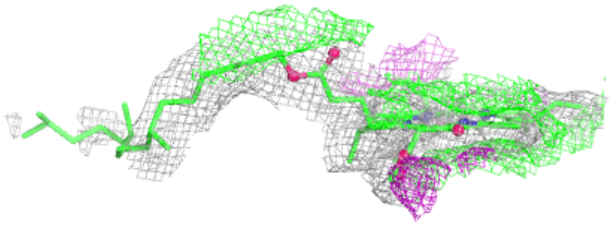




Electron density around BPH AM 403:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

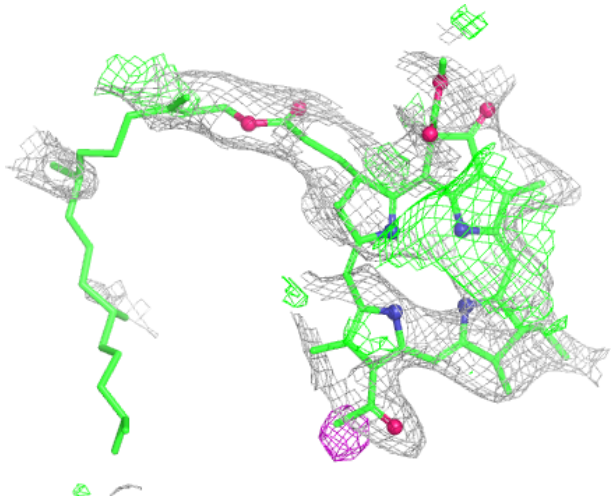


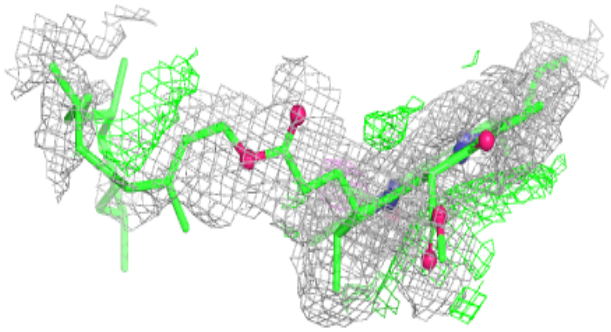


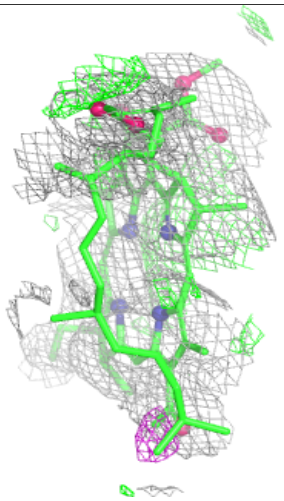


Electron density around BPH BL 302:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

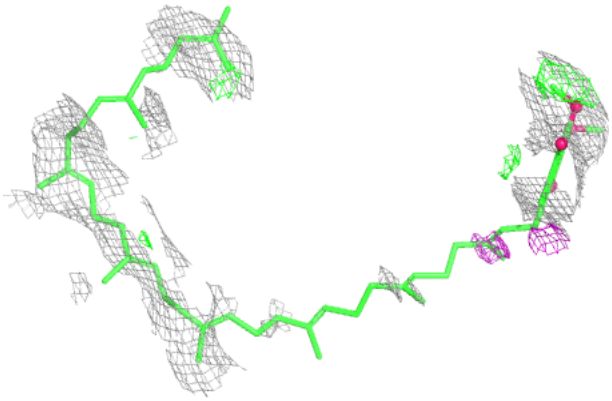


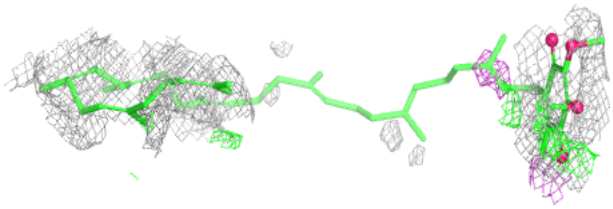


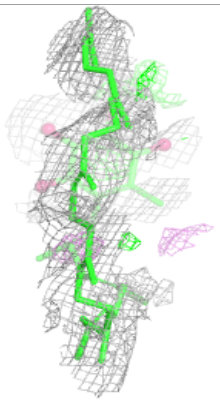


Electron density around UQ8 BL 304:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

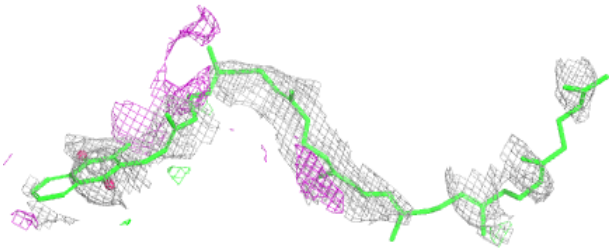


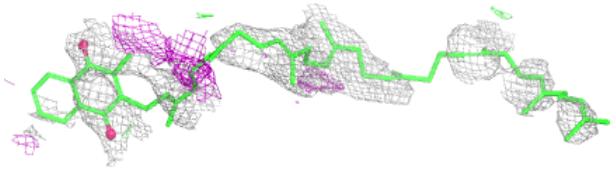


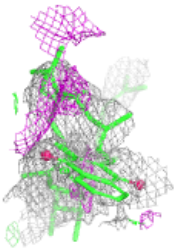


Electron density around MQ8 AM 405:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

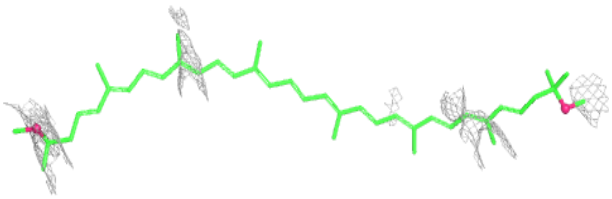


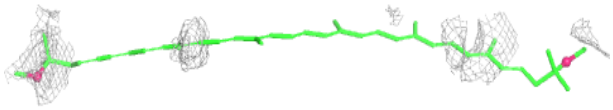


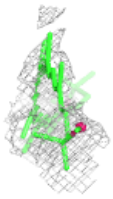


Electron density around CRT AA 102:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

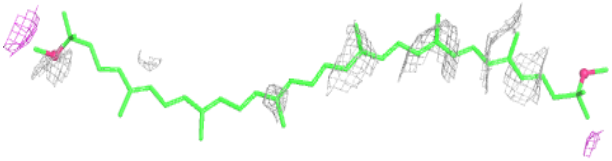


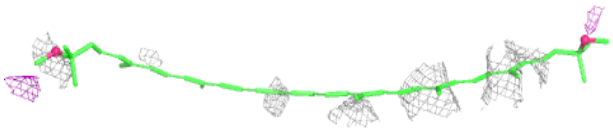


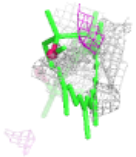


Electron density around CRT AG 102:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

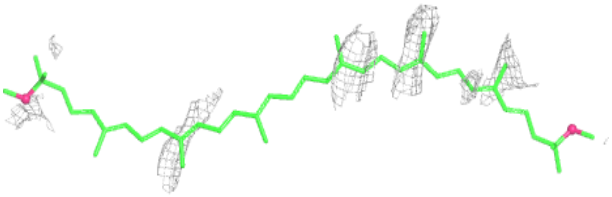


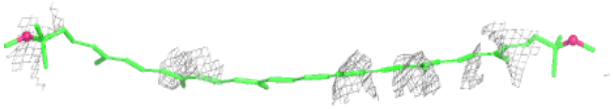




Electron density around CRT AJ 102:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

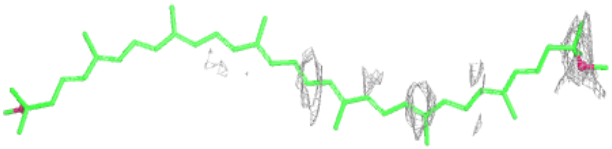


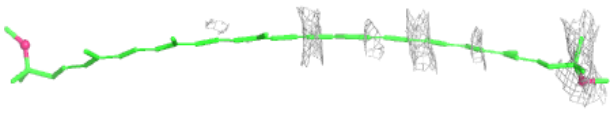


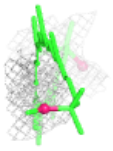


Electron density around CRT AP 102:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

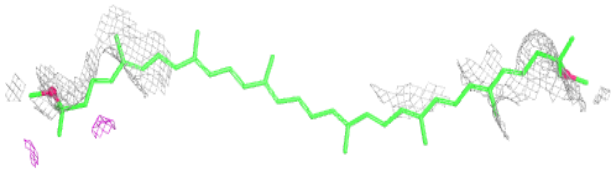


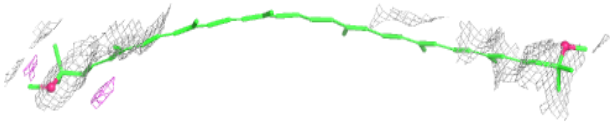


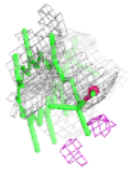


Electron density around CRT AS 104:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

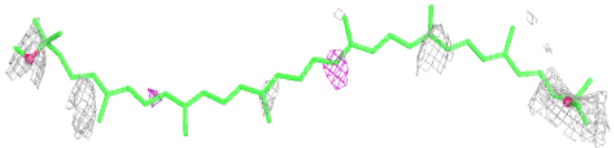






Electron density around CRT AW 102:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

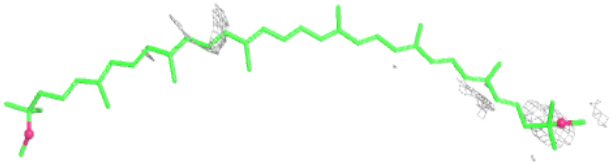




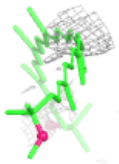


Electron density around CRT AX 102:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

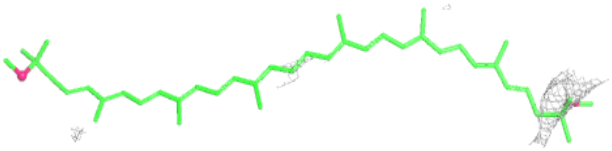


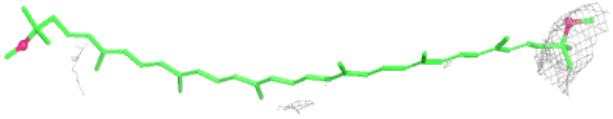


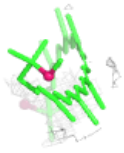


Electron density around CRT A5 103:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

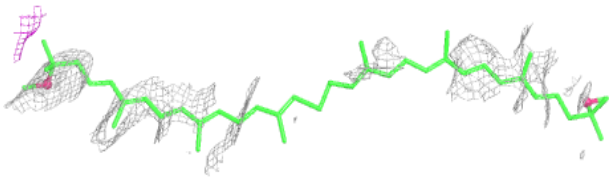


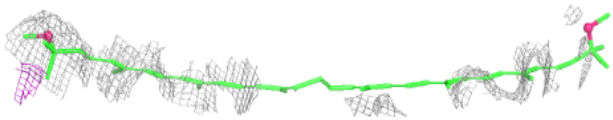


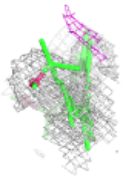


Electron density around CRT A7 102:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

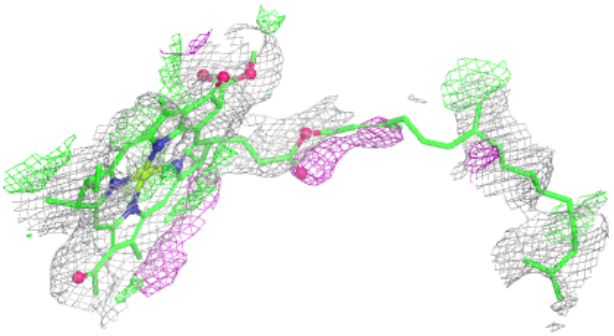


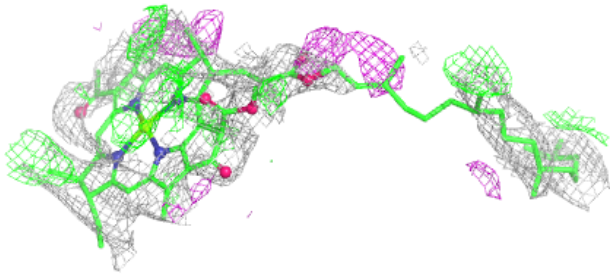


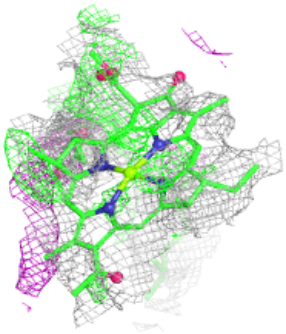


Electron density around BCL AL 301:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)





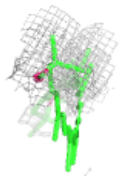


Electron density around CRT BA 102:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

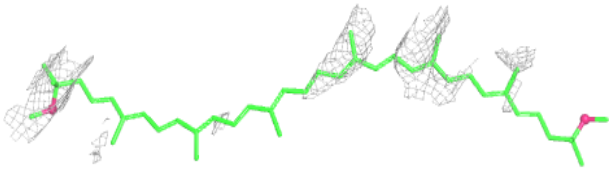


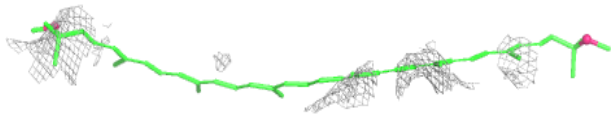




Electron density around CRT BB 102:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

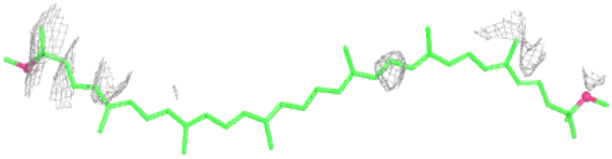


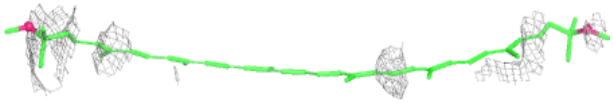




Electron density around CRT BF 103:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

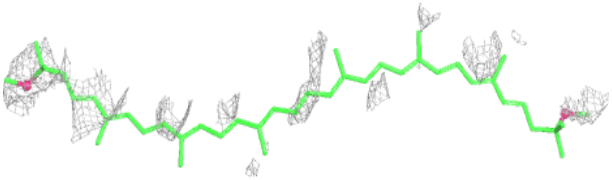


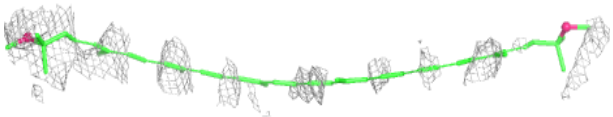




Electron density around CRT BN 102:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

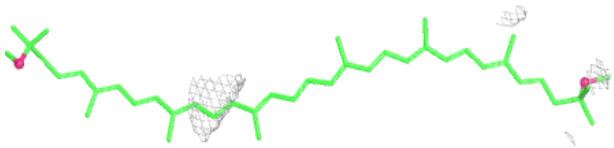


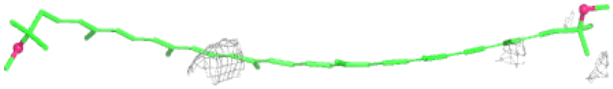




Electron density around CRT BO 103:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

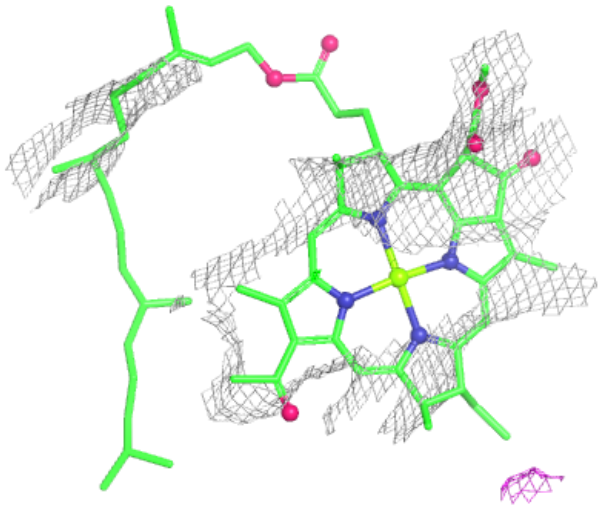


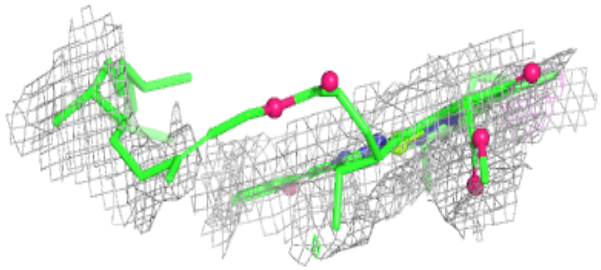




Electron density around BCL AB 101:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

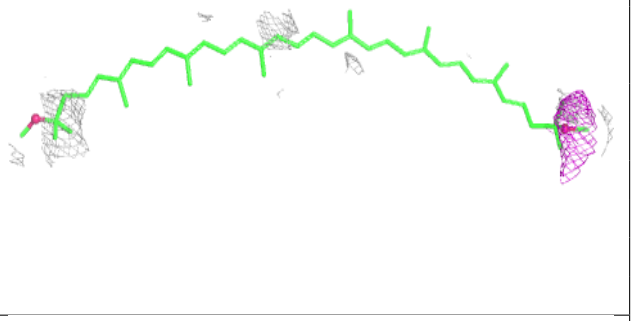


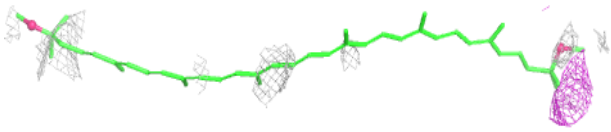


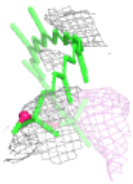


Electron density around CRT BU 103:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

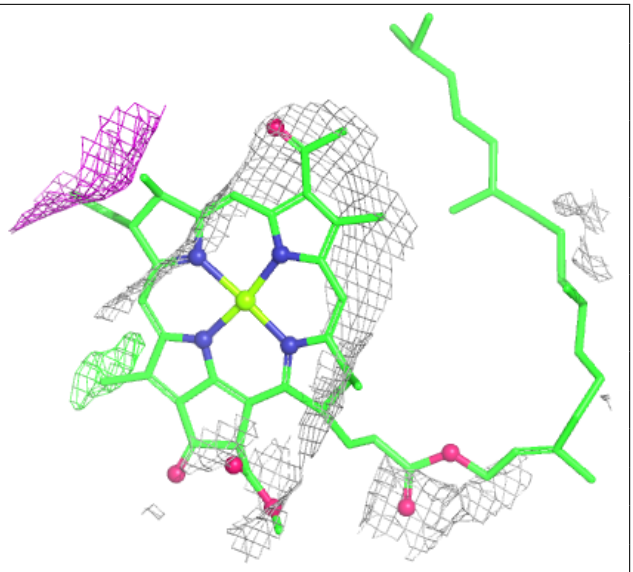


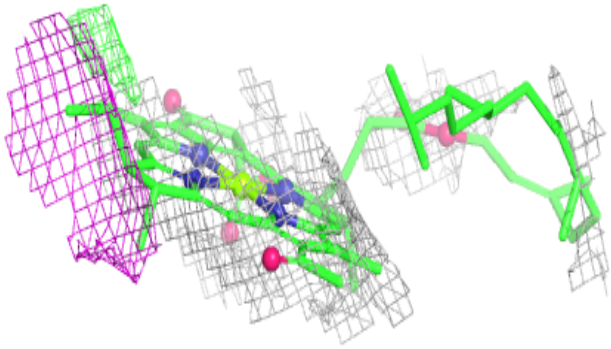


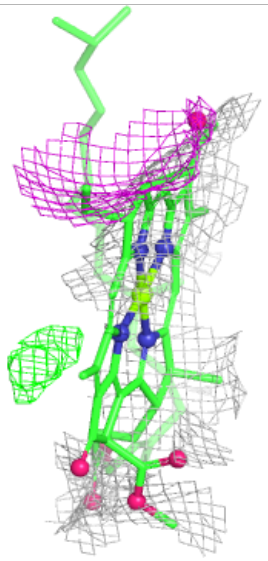


Electron density around BCL AE 101:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

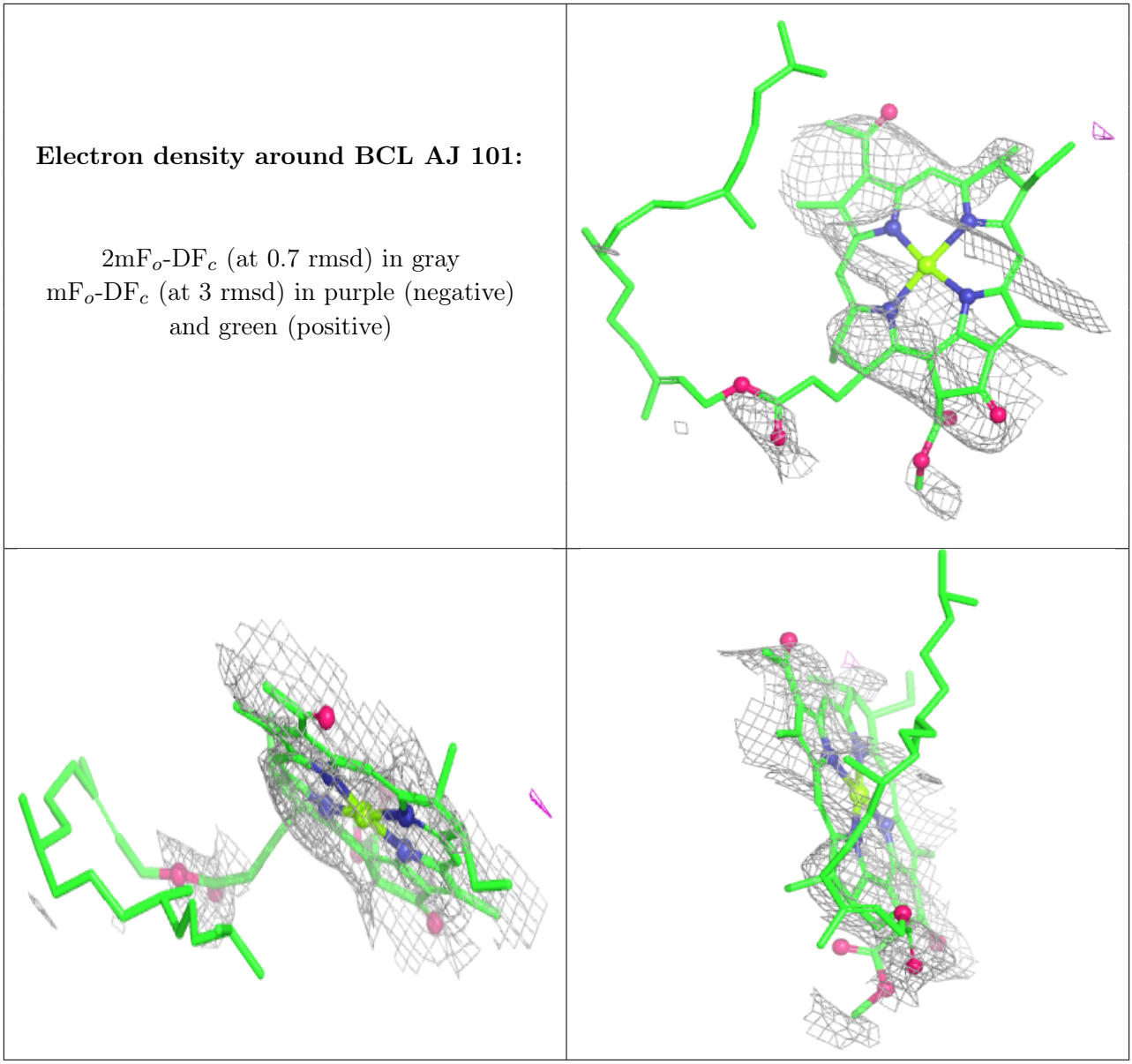






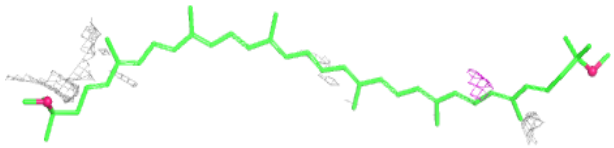
Electron density around BCL AJ 101:

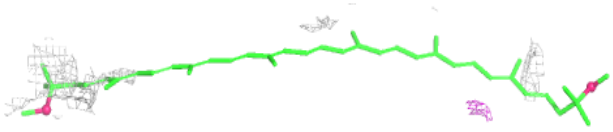
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

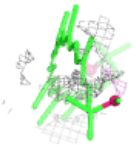


Electron density around CRT B5 103:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

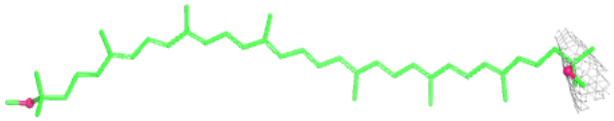


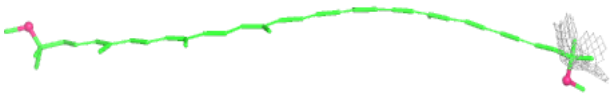




Electron density around CRT B0 101:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

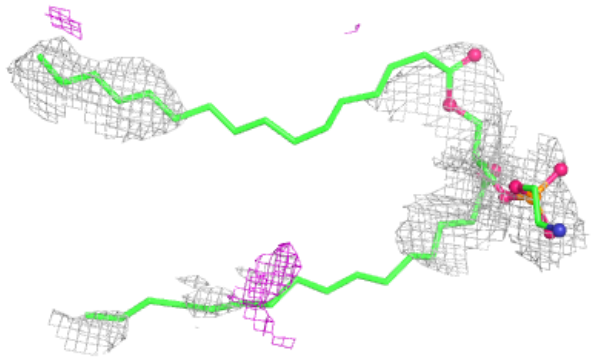


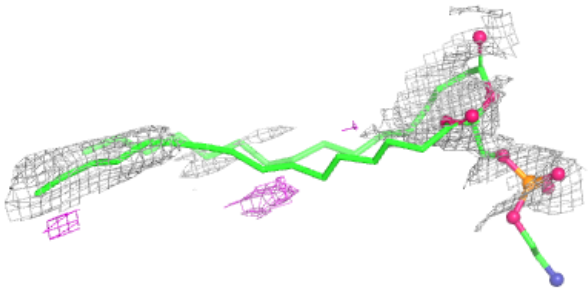


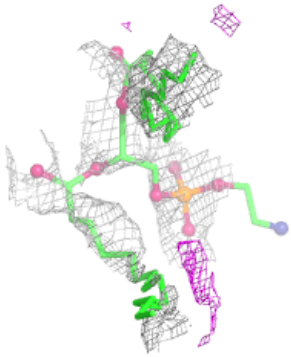


Electron density around PEF AM 409:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

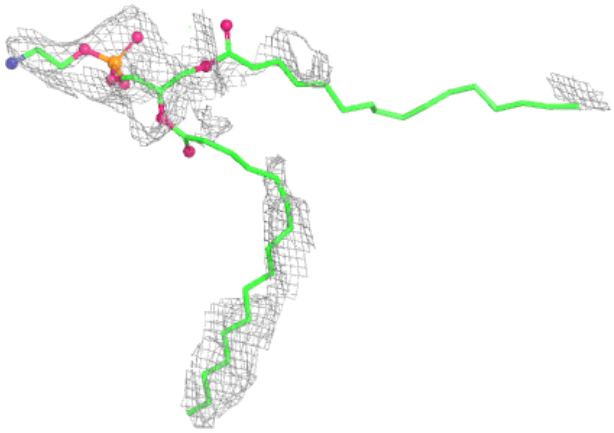


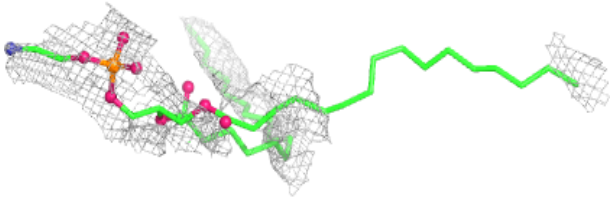




Electron density around PEF AS 101:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

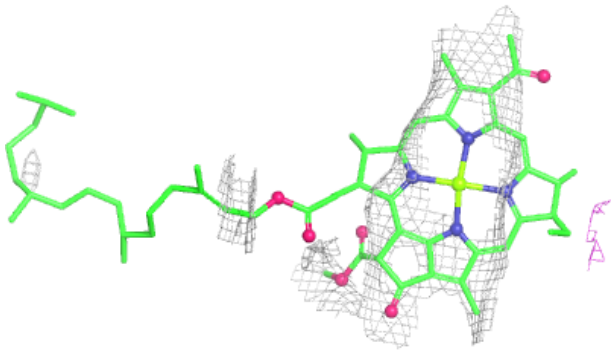


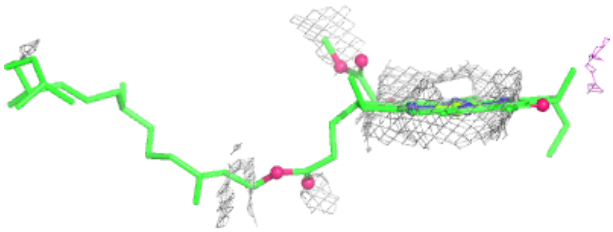




Electron density around BCL AK 102:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

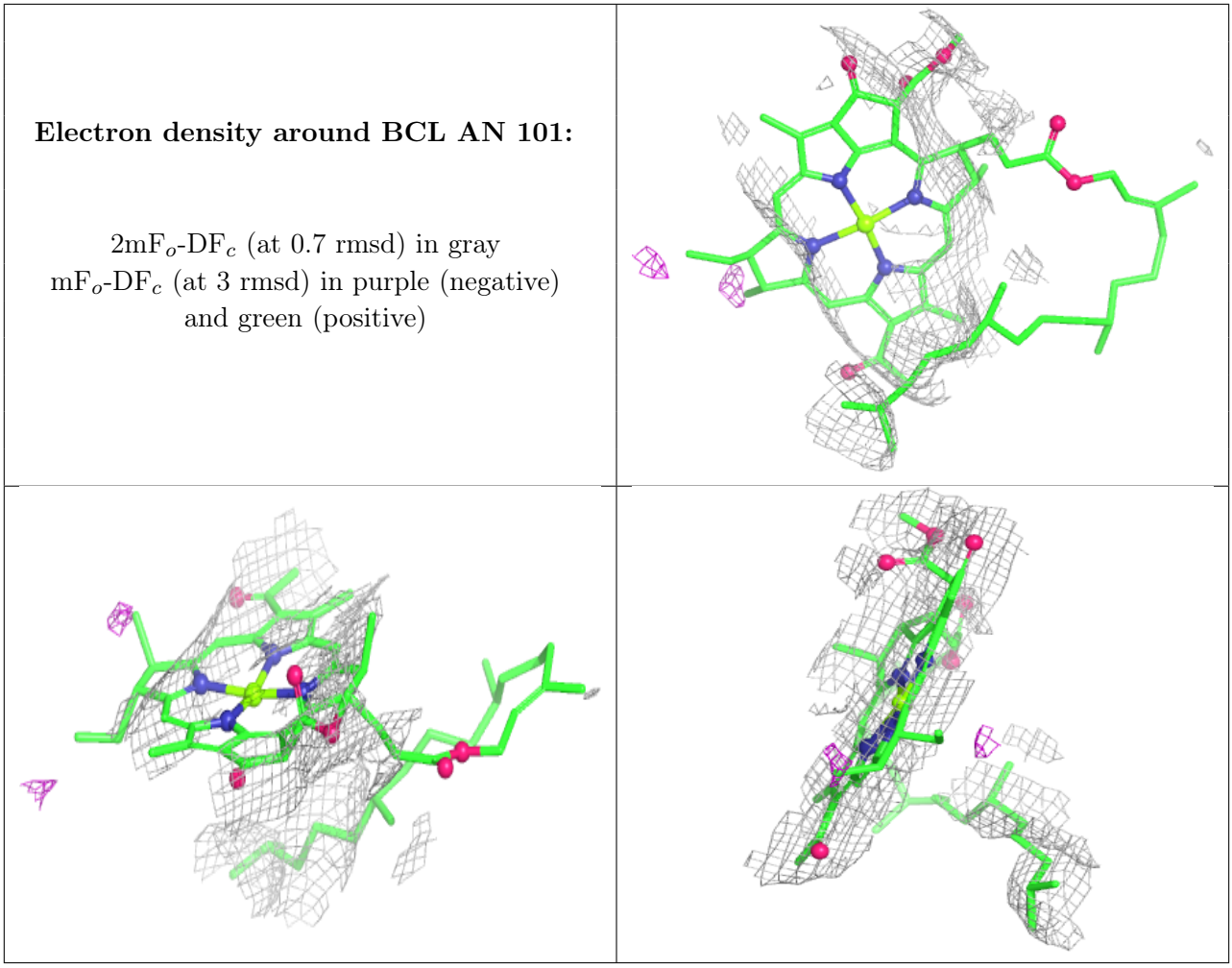






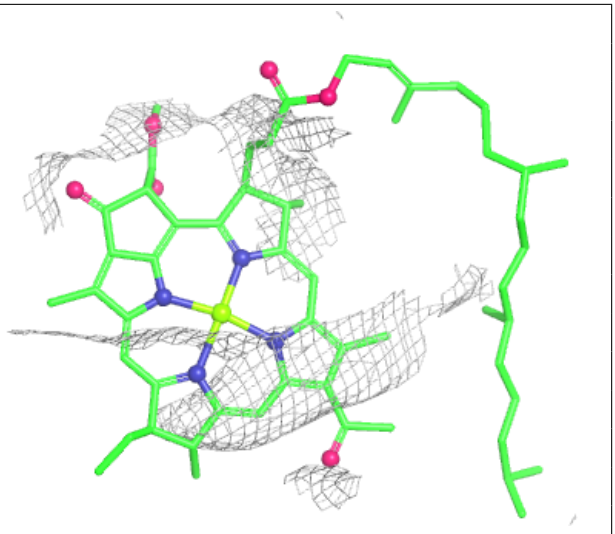
Electron density around BCL AN 101:

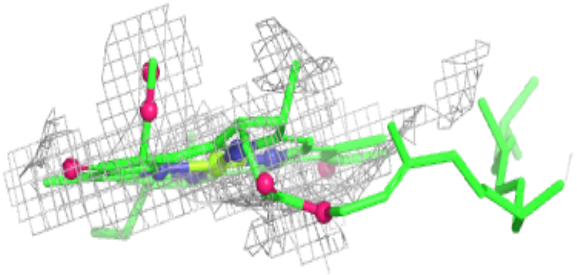
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

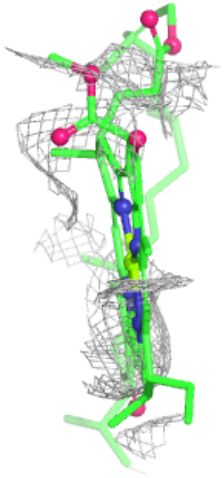


Electron density around BCL BT 101:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

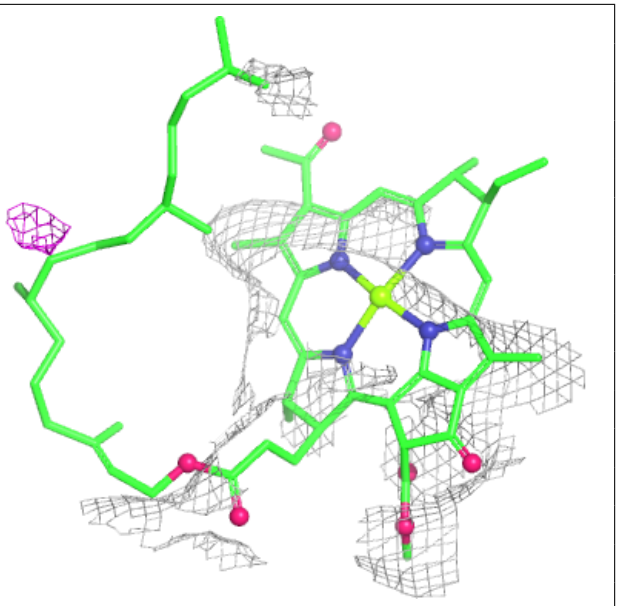


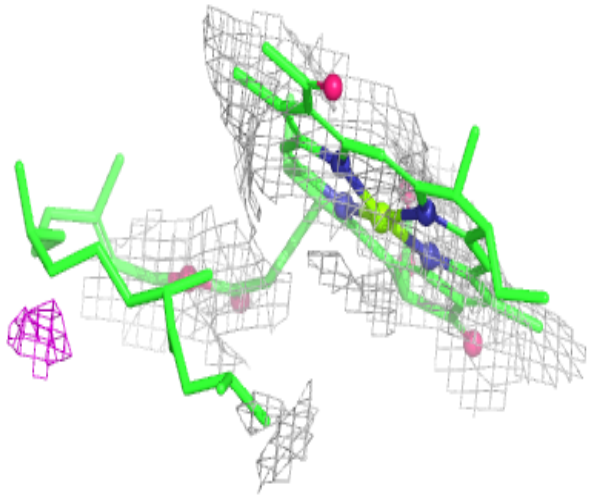




Electron density around BCL BV 101:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

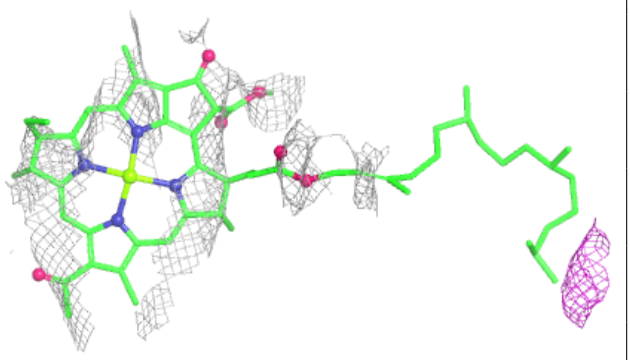


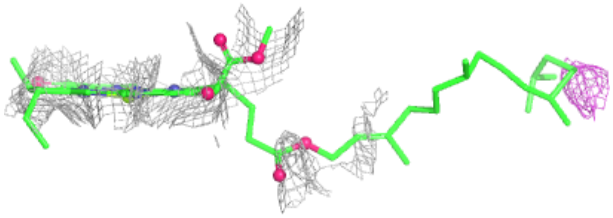


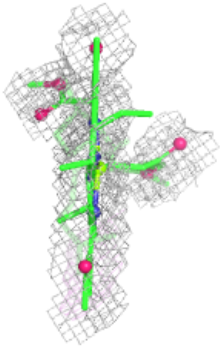


Electron density around BCL AF 102:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

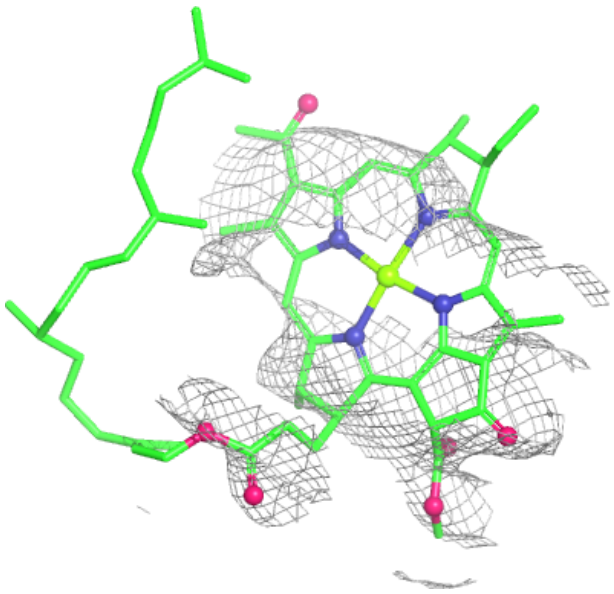


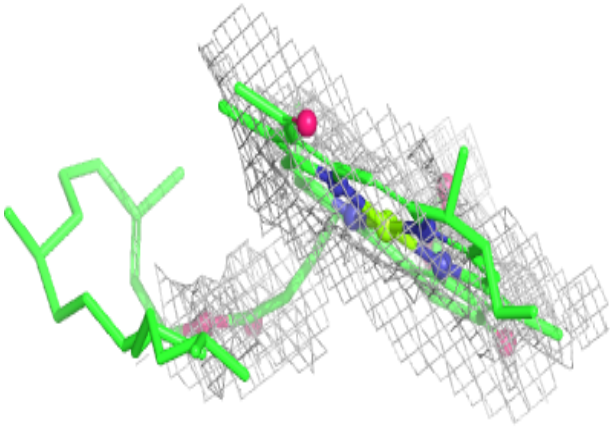


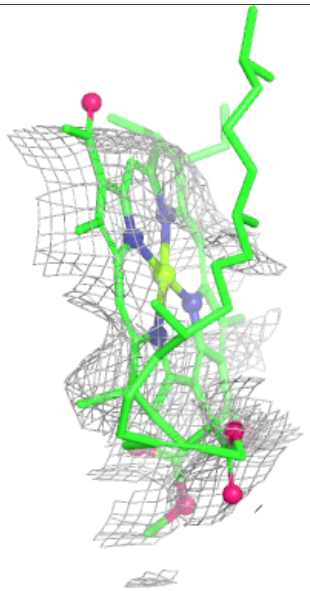


Electron density around BCL AG 101:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

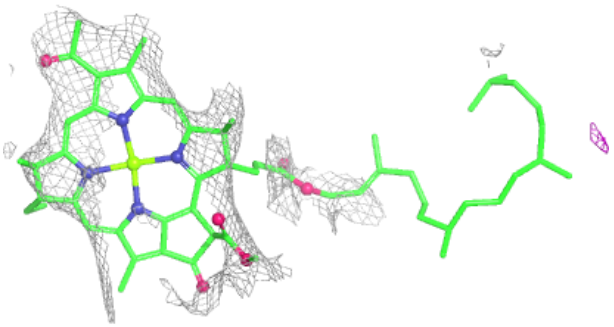


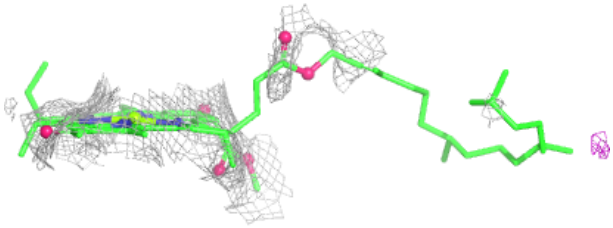


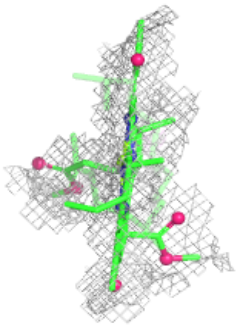


Electron density around BCL BY 102:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

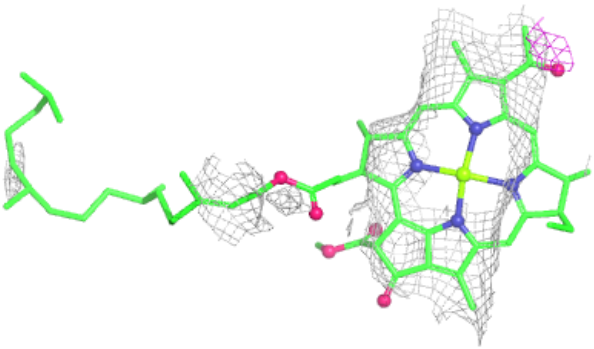


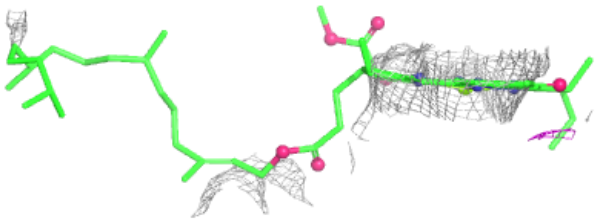




Electron density around BCL AI 102:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

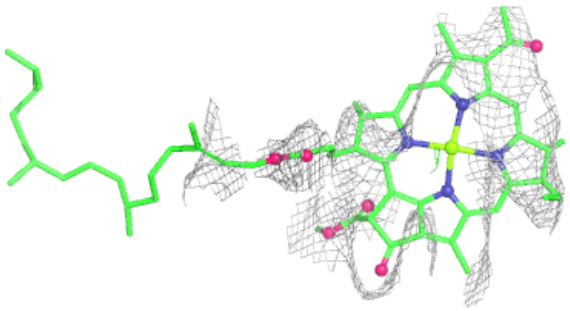


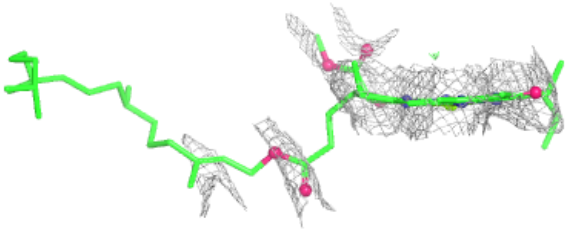




Electron density around BCL B1 102:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

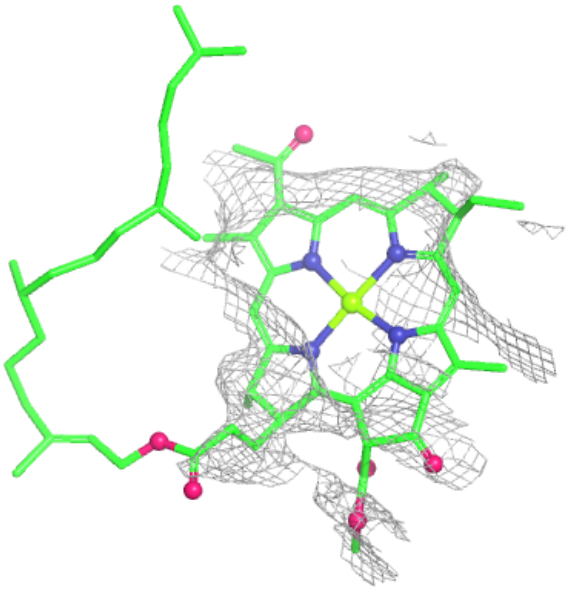


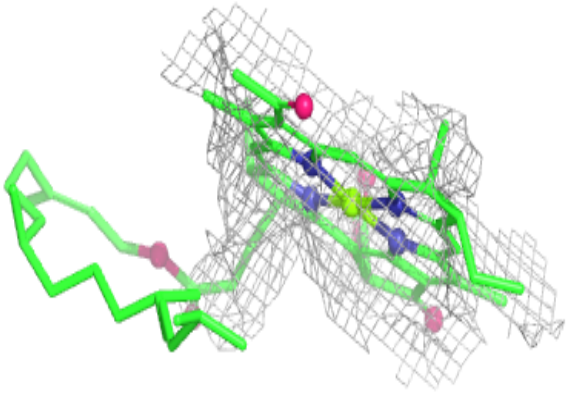


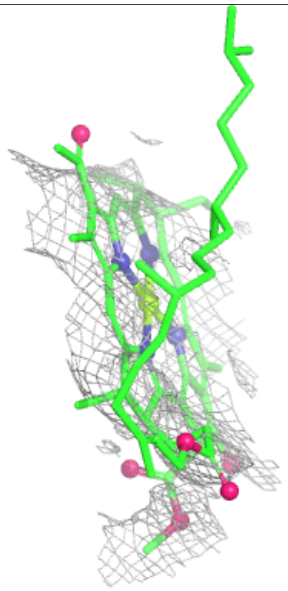


Electron density around BCL B2 101:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

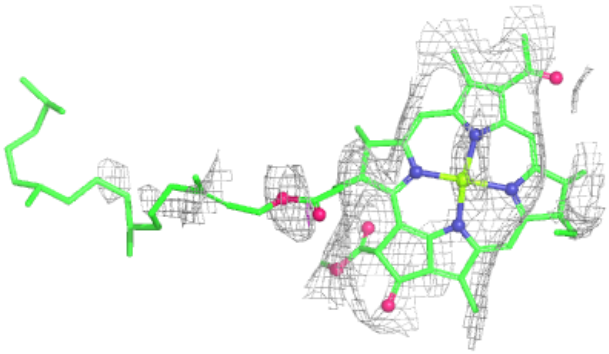


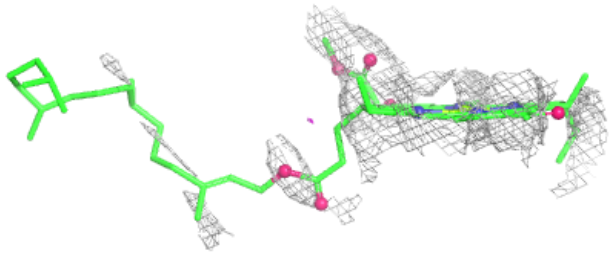




Electron density around BCL B3 102:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

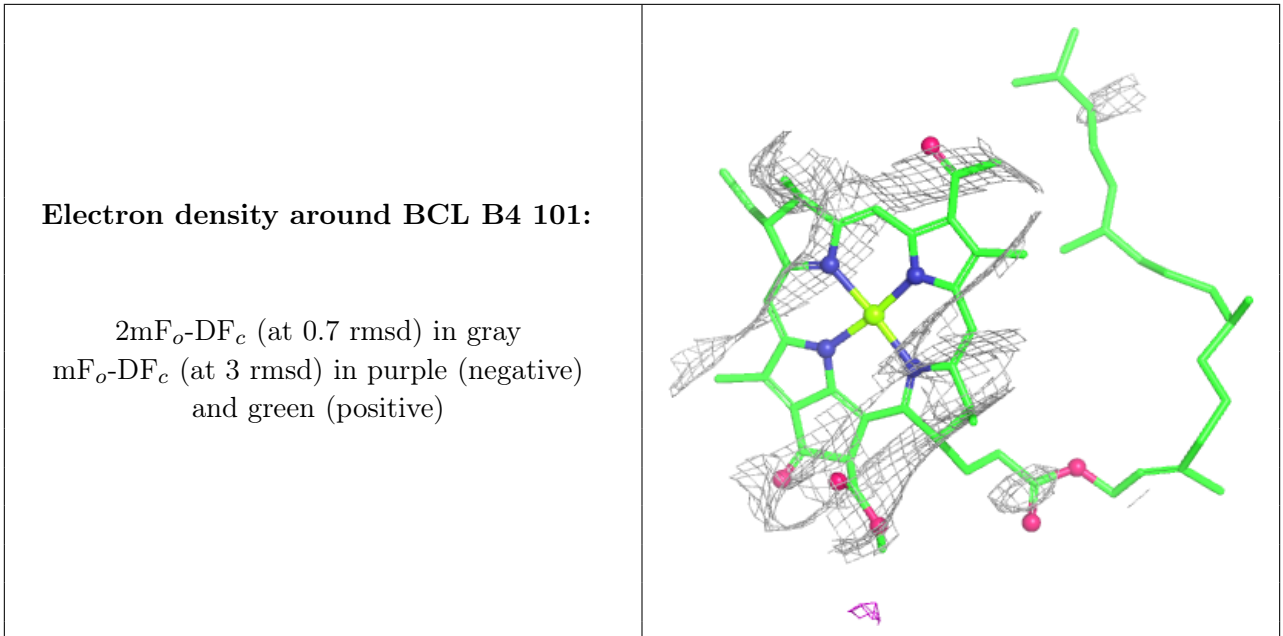


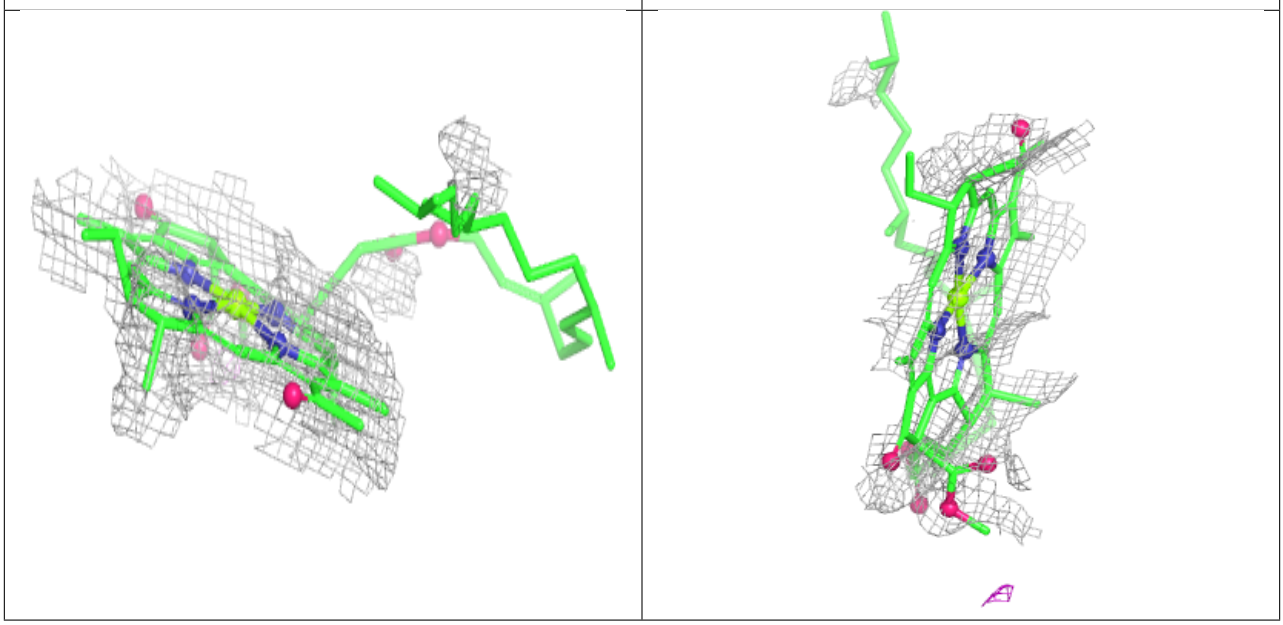




Electron density around BCL B4 101:

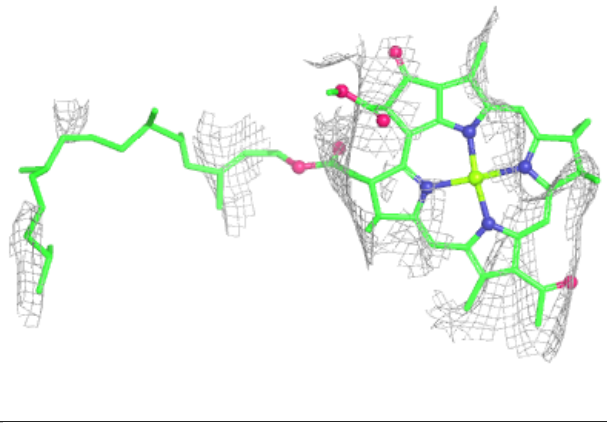
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

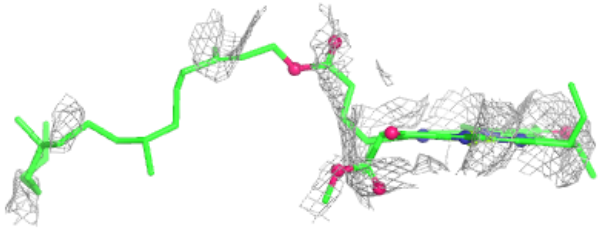


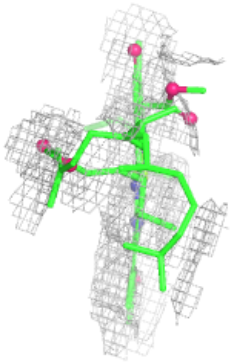


Electron density around BCL B5 102:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

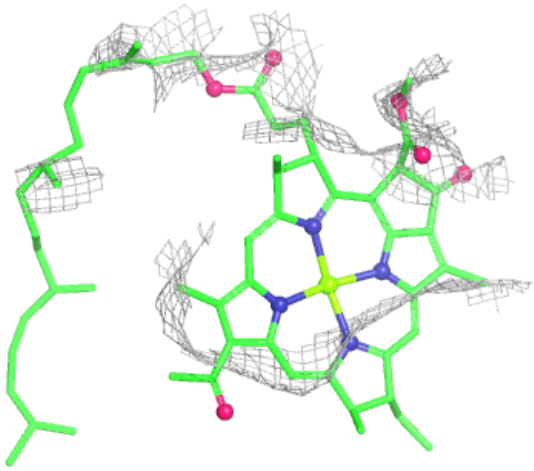


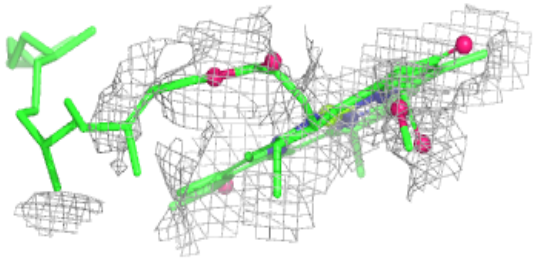


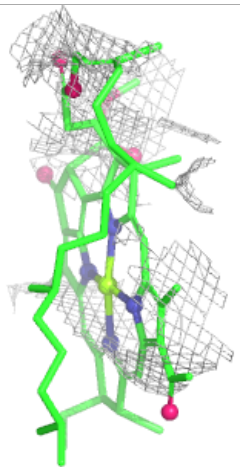


Electron density around BCL B6 101:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

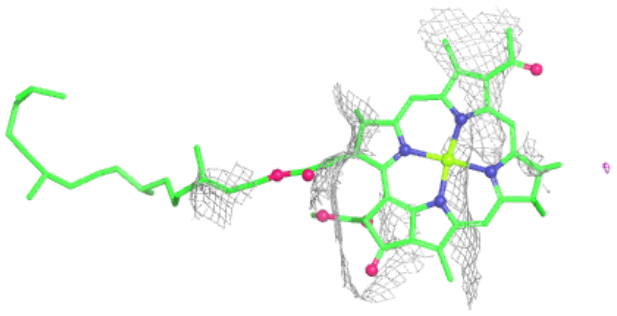


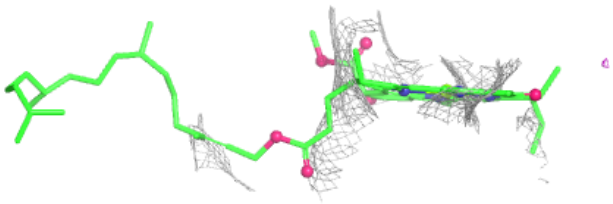




Electron density around BCL B7 103:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

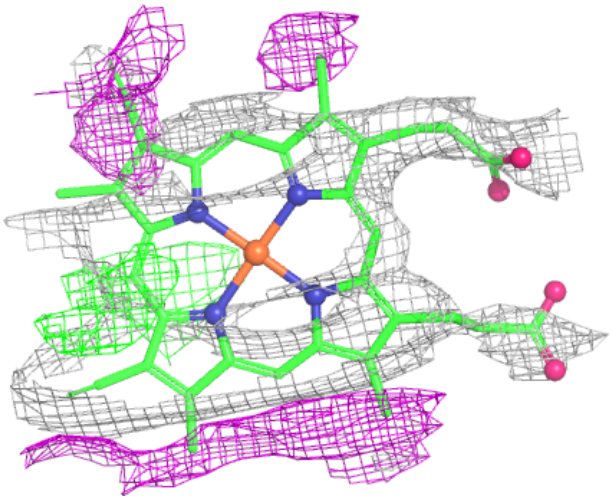


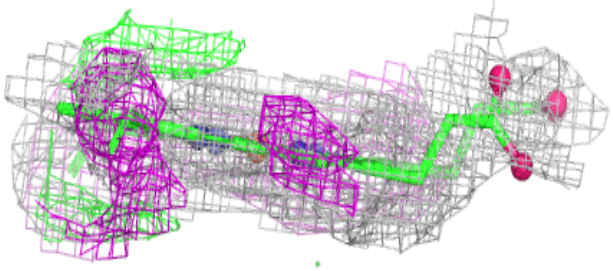


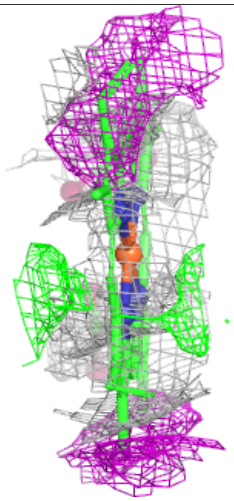


Electron density around HEM AC 502:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

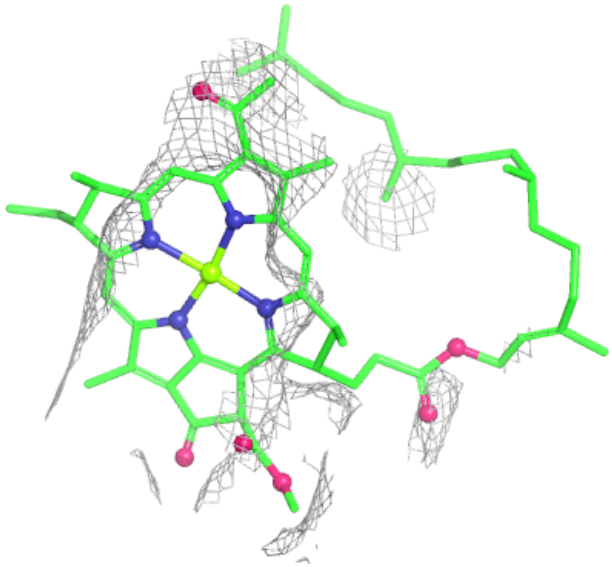


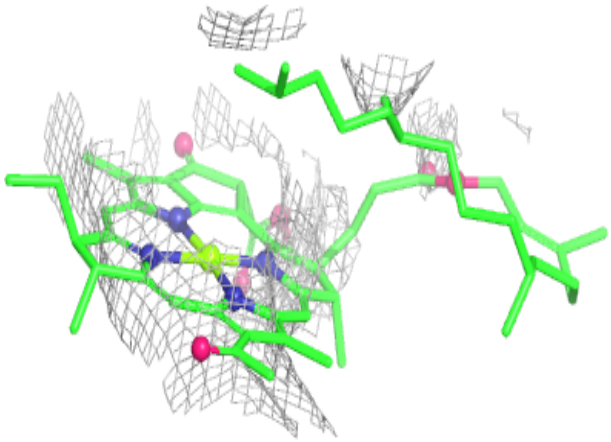


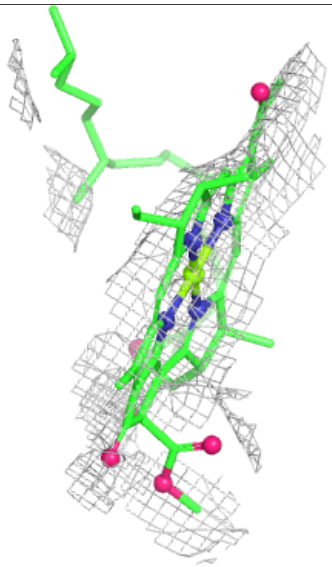


Electron density around BCL B0 102:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

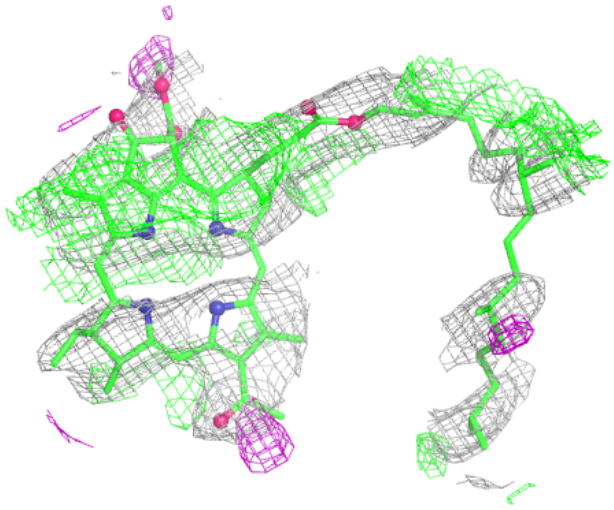


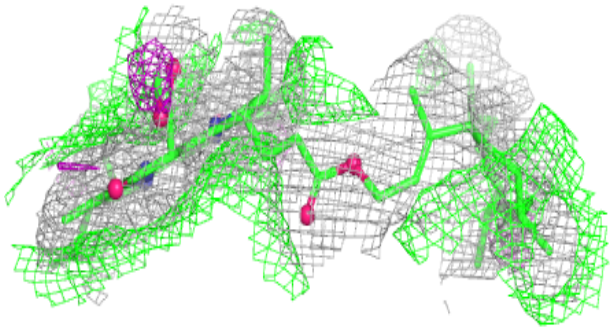




Electron density around BPH AL 302:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

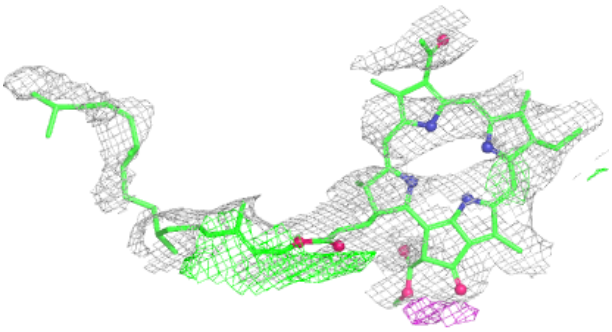


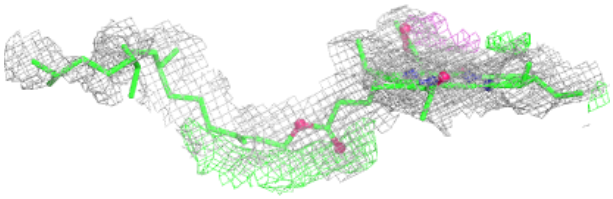


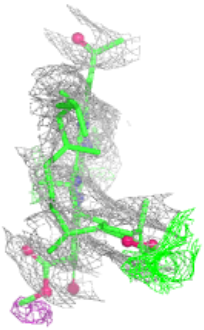


Electron density around BPH BM 403:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

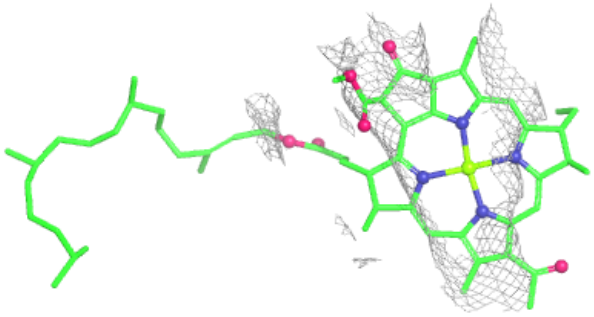


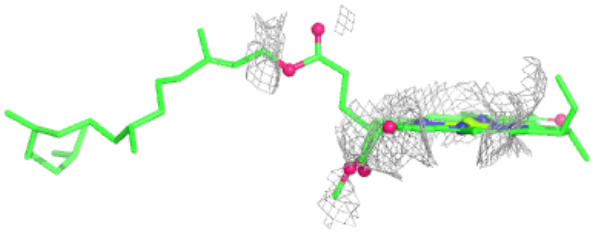


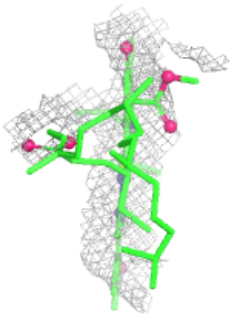


Electron density around BCL AS 103:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

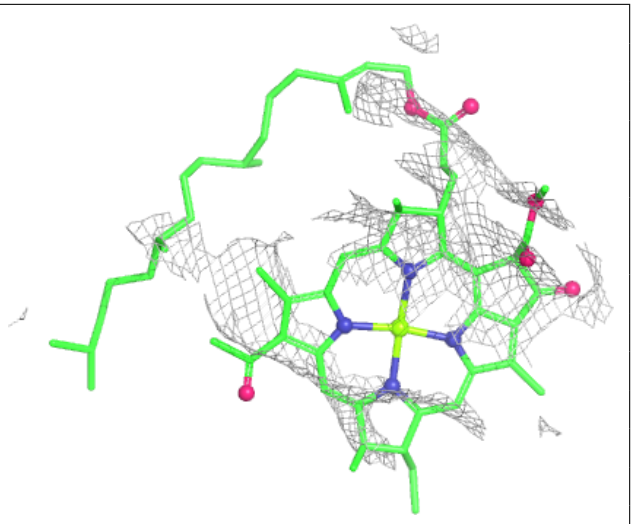


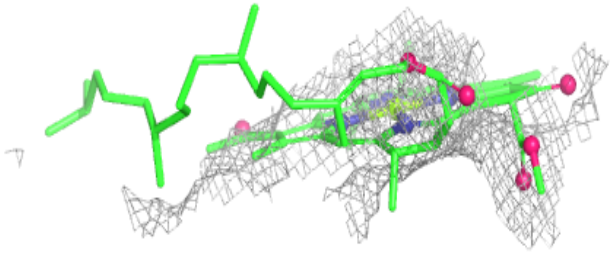




Electron density around BCL AT 101:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

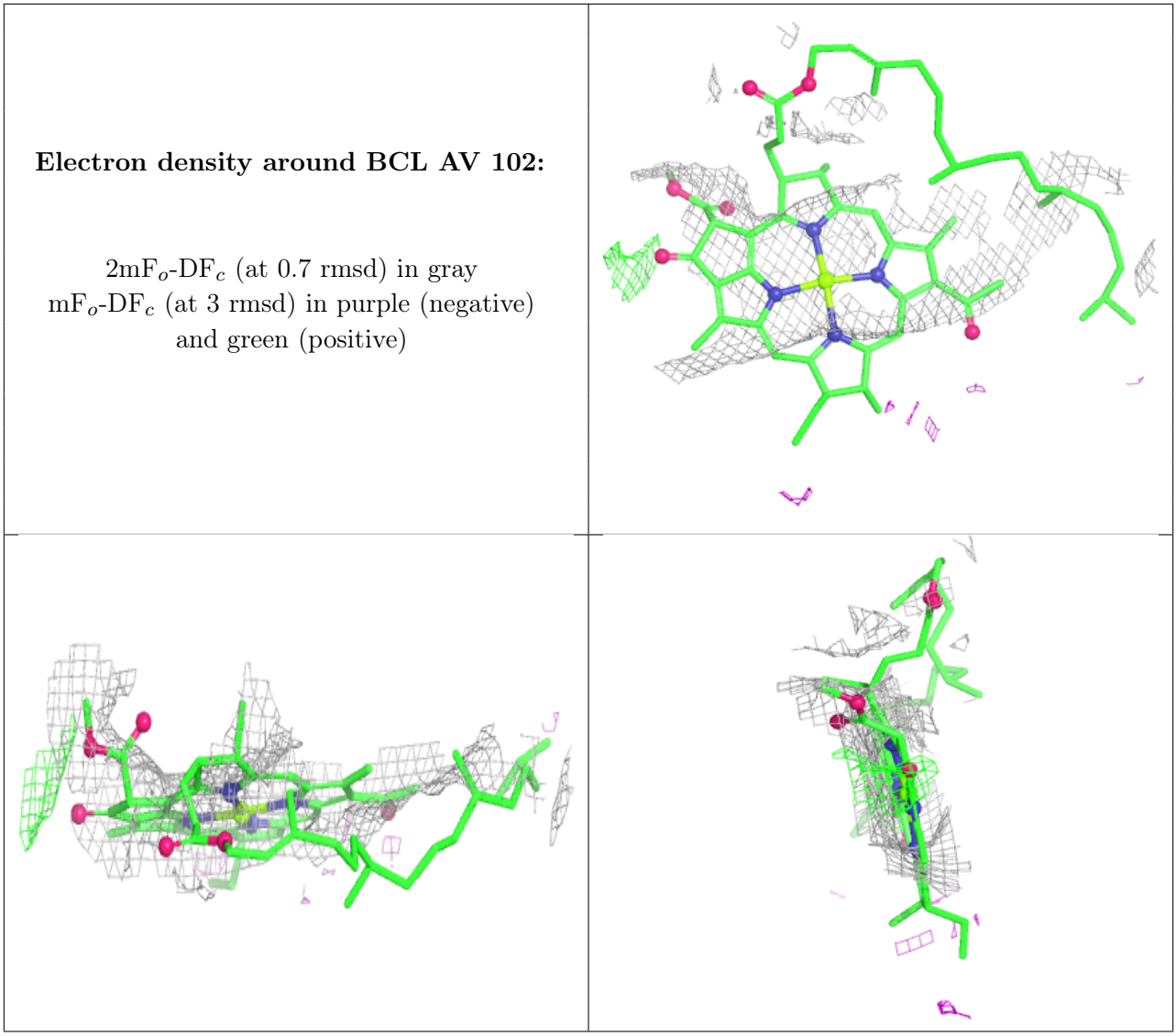






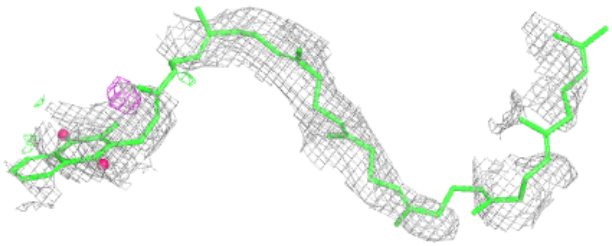
Electron density around BCL AV 102:

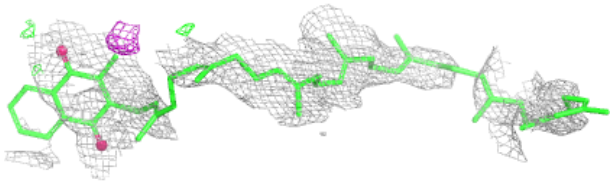
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

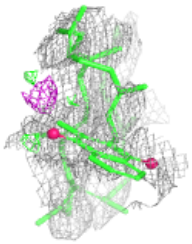


Electron density around MQ8 BM 405:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

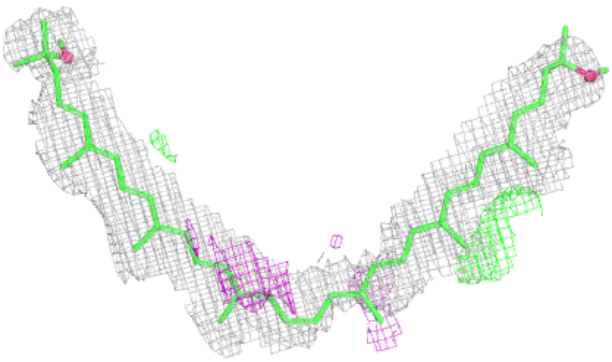


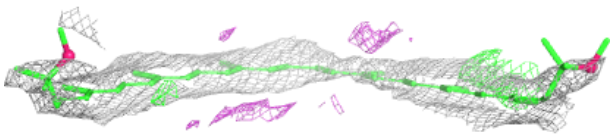


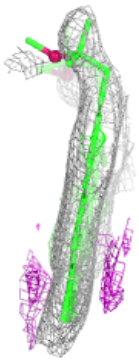


Electron density around CRT AM 406:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

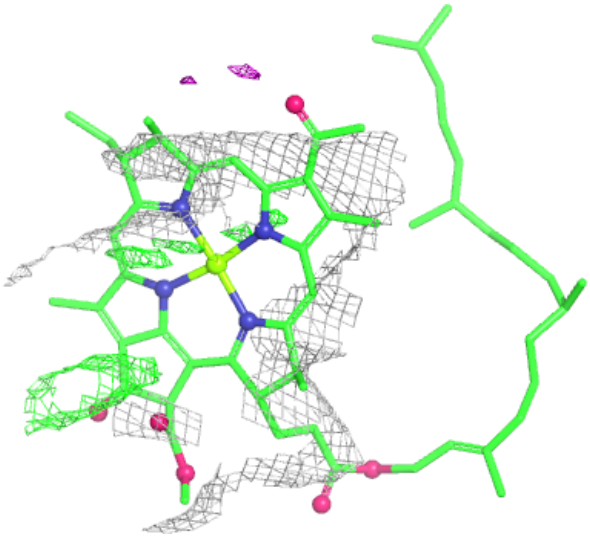


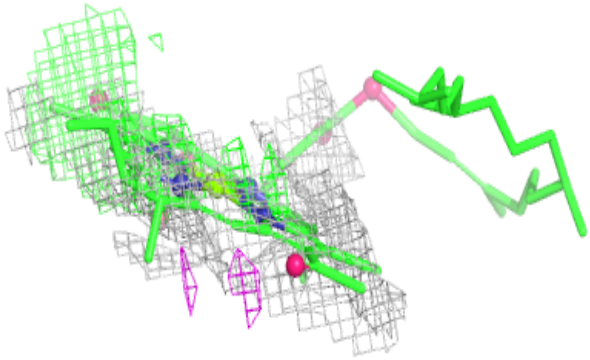


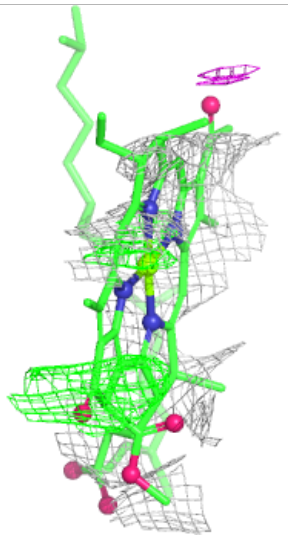


Electron density around BCL AX 101:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

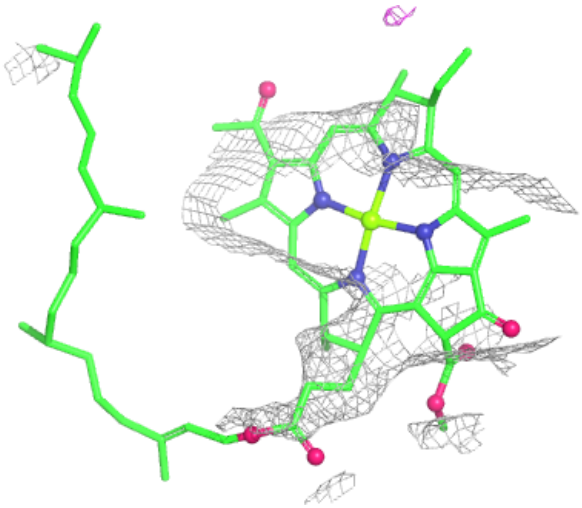


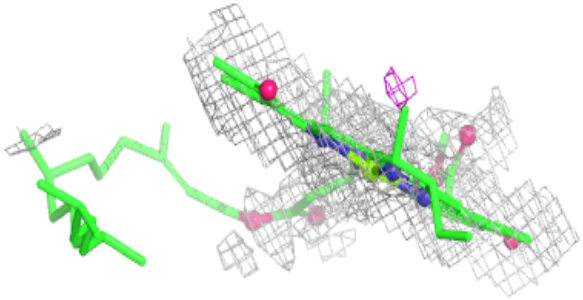


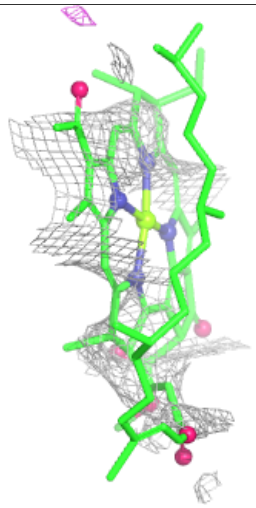


Electron density around BCL AZ 101:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

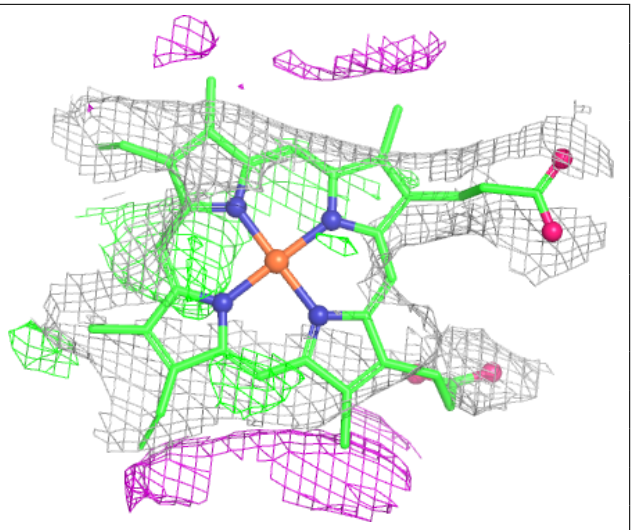


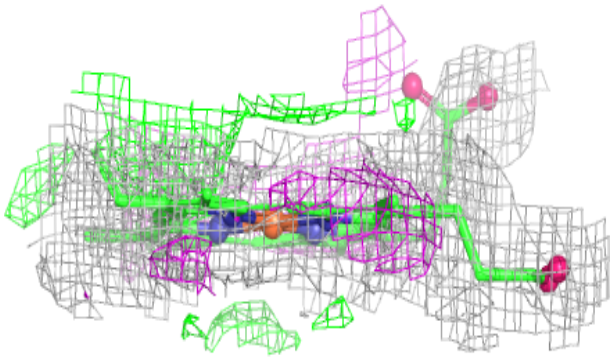


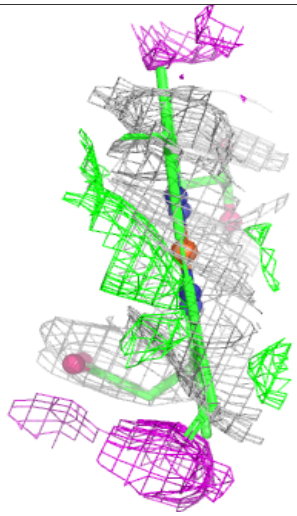


Electron density around HEM AC 503:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

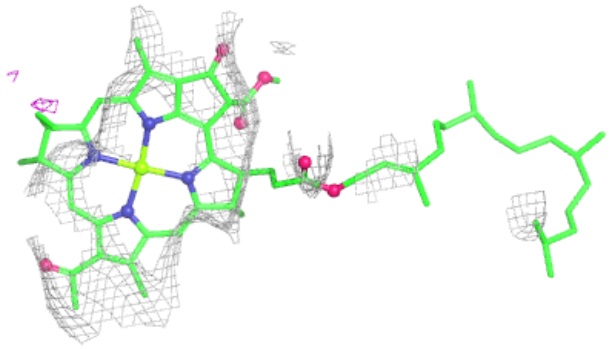


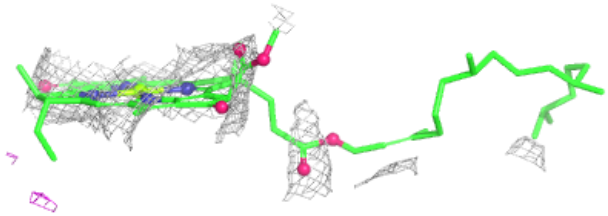


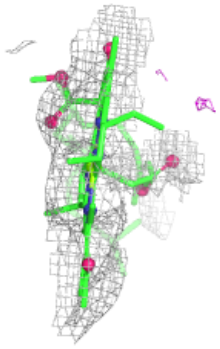


Electron density around BCL A5 102:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

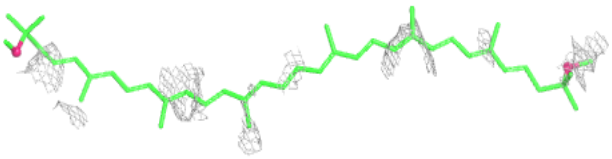


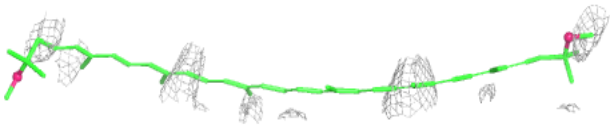


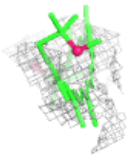


Electron density around CRT AR 102:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

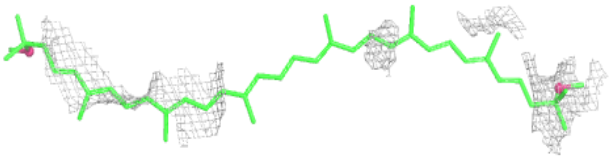


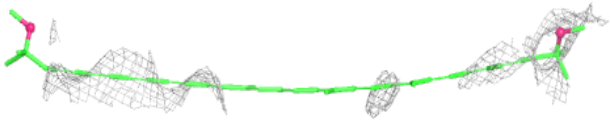




Electron density around CRT AT 102:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

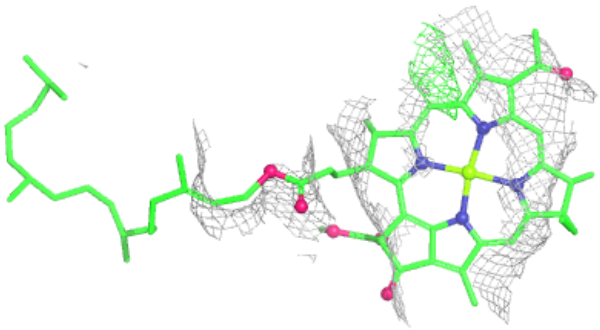


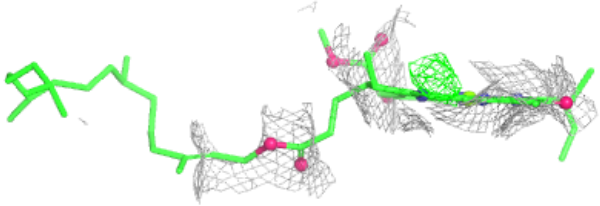




Electron density around BCL A7 103:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

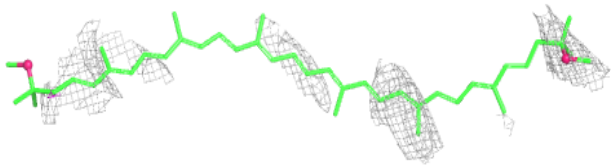


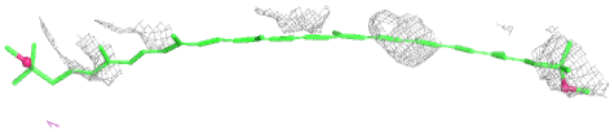


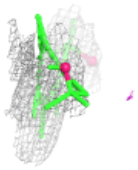


Electron density around CRT A1 103:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

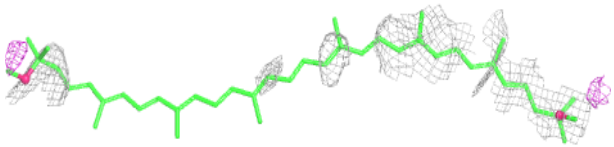


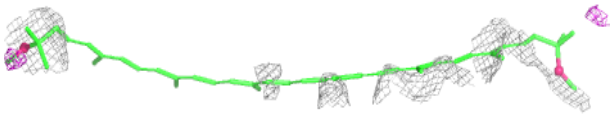


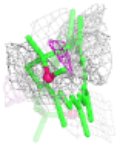


Electron density around CRT A2 102:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

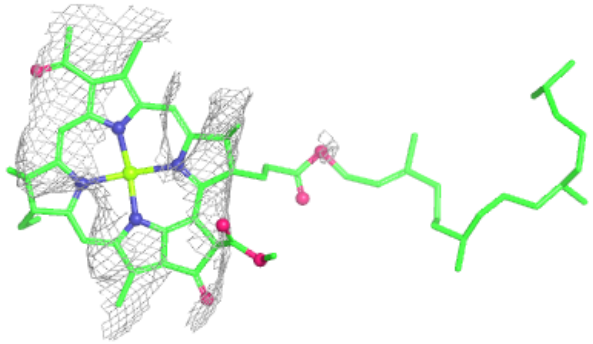


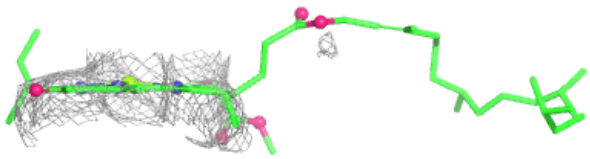


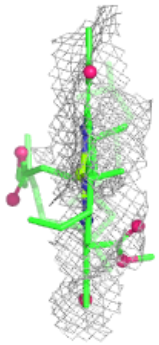


Electron density around BCL A9 102:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

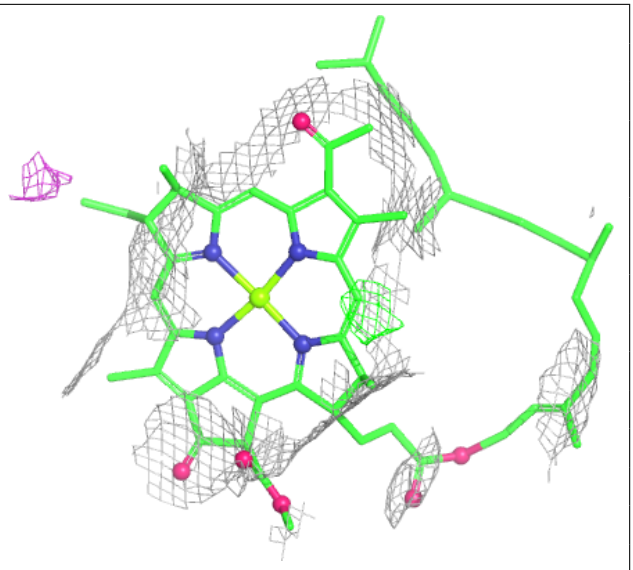


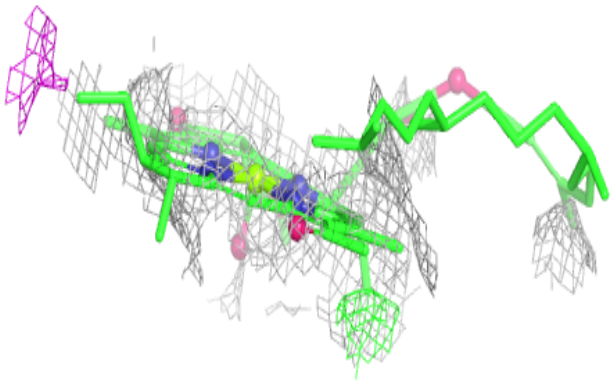




Electron density around BCL A0 102:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

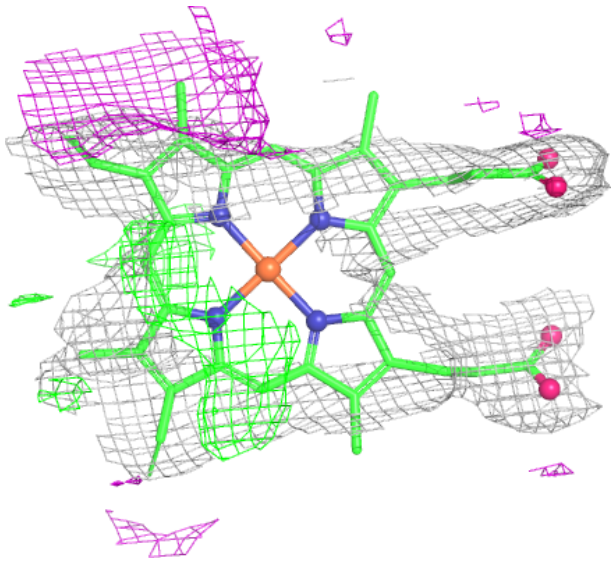


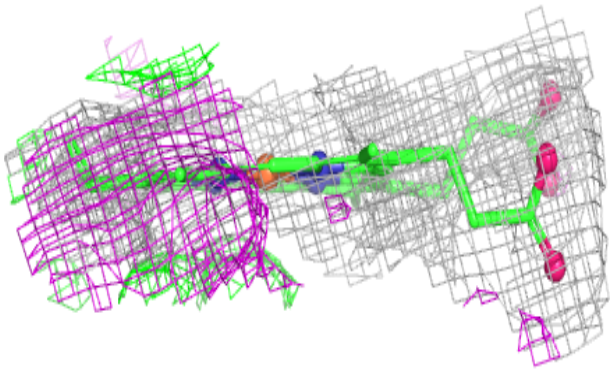


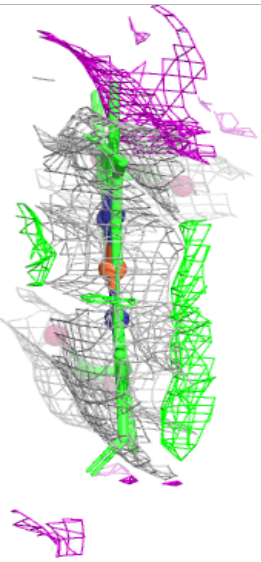


Electron density around HEM AC 504:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

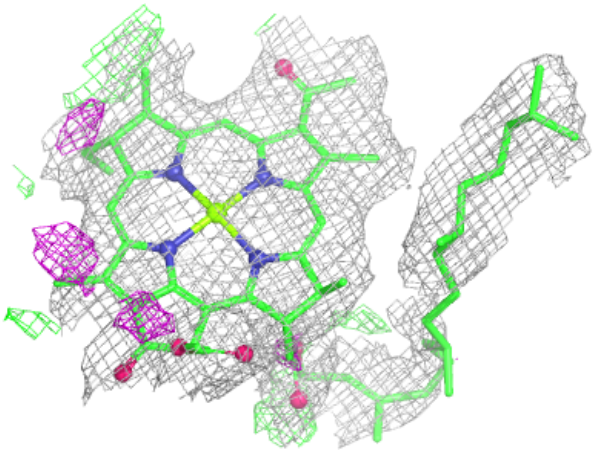


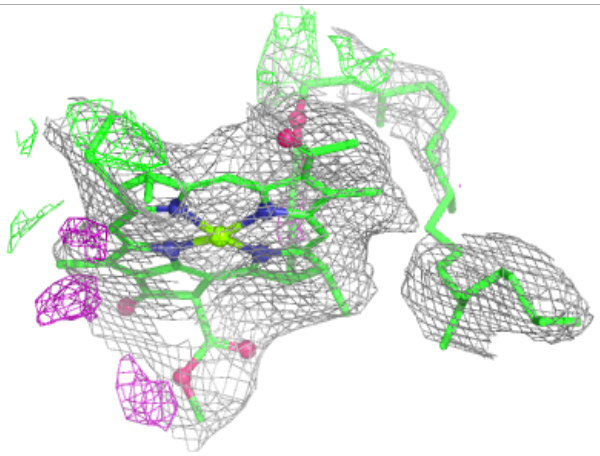


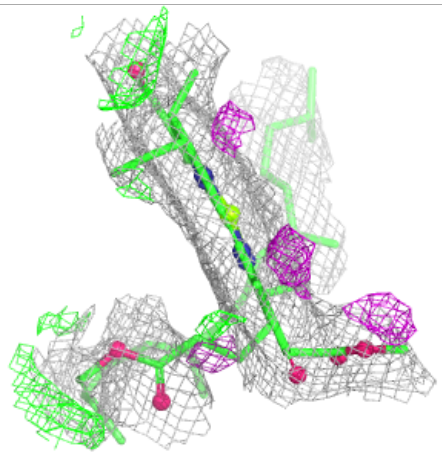


Electron density around BCL BL 303:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

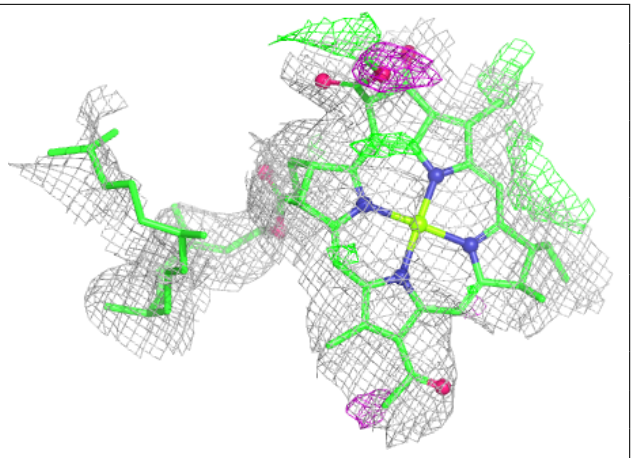


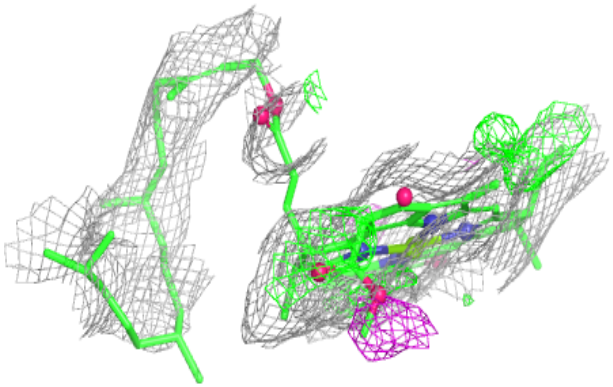


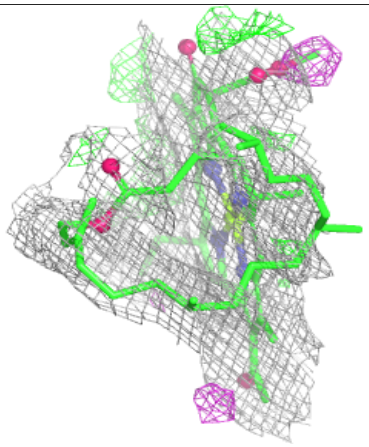


Electron density around BCL BM 401:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

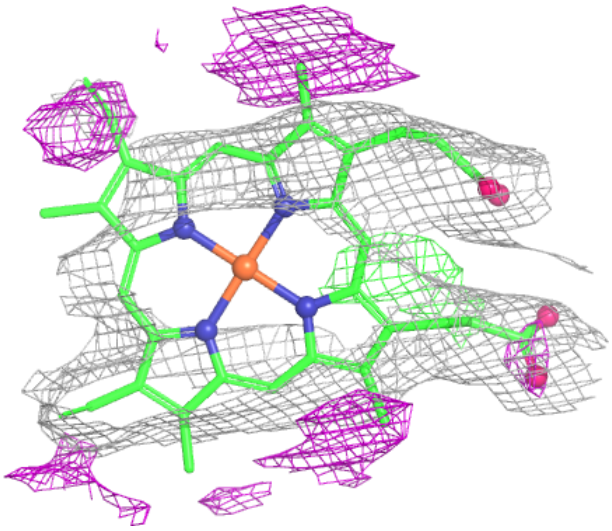


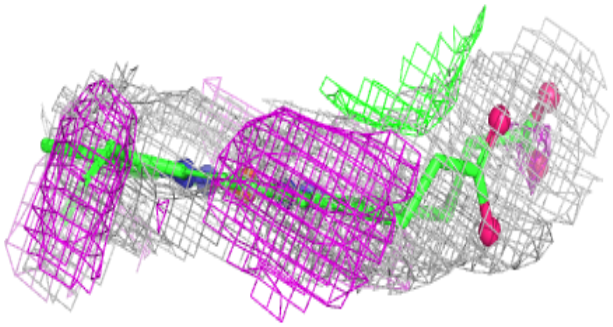


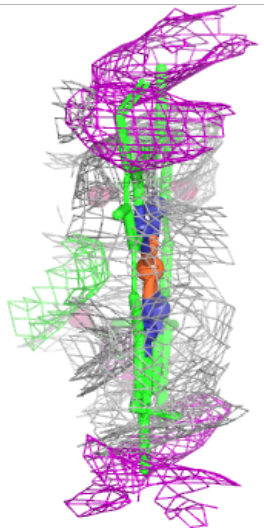


Electron density around HEM BC 502:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

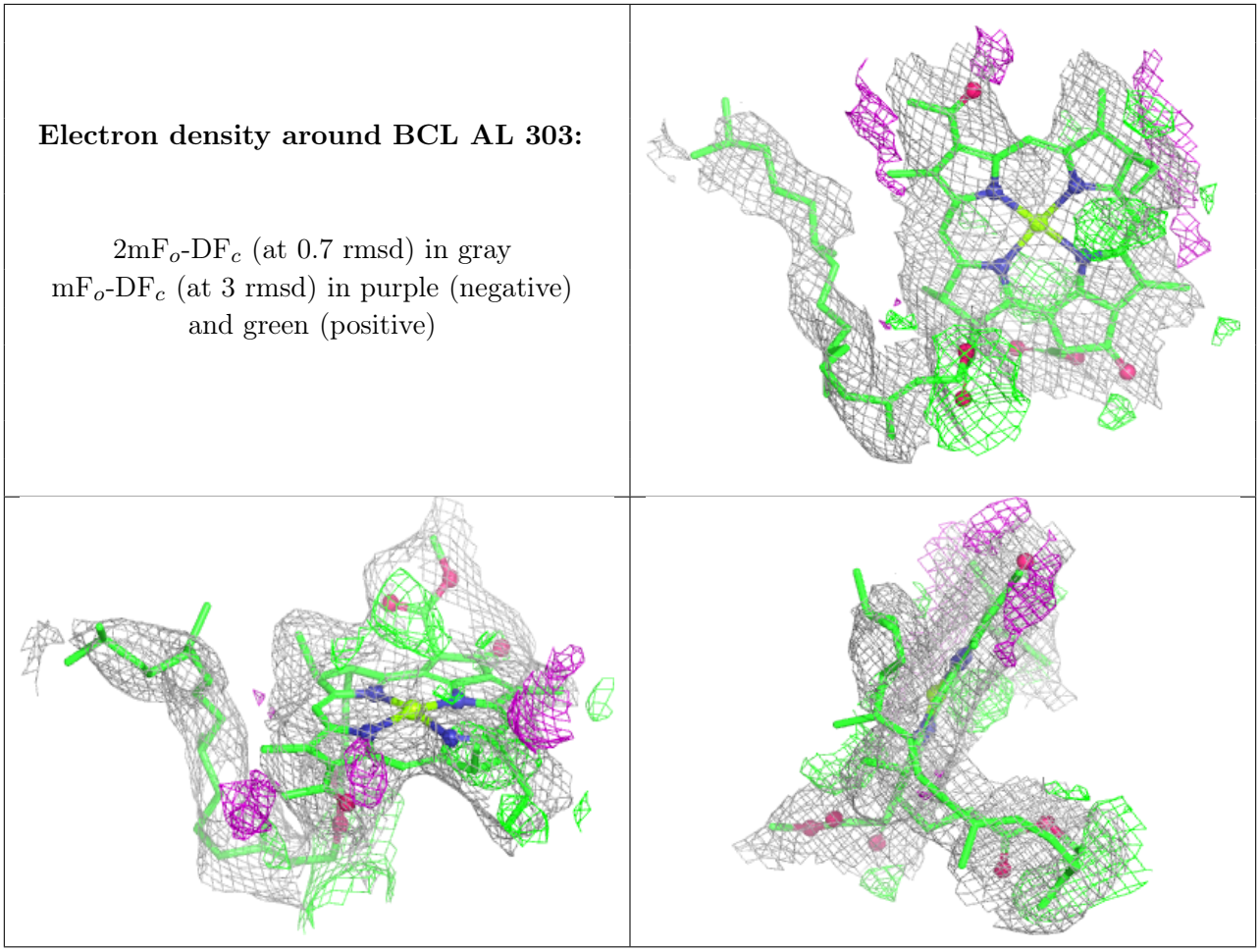






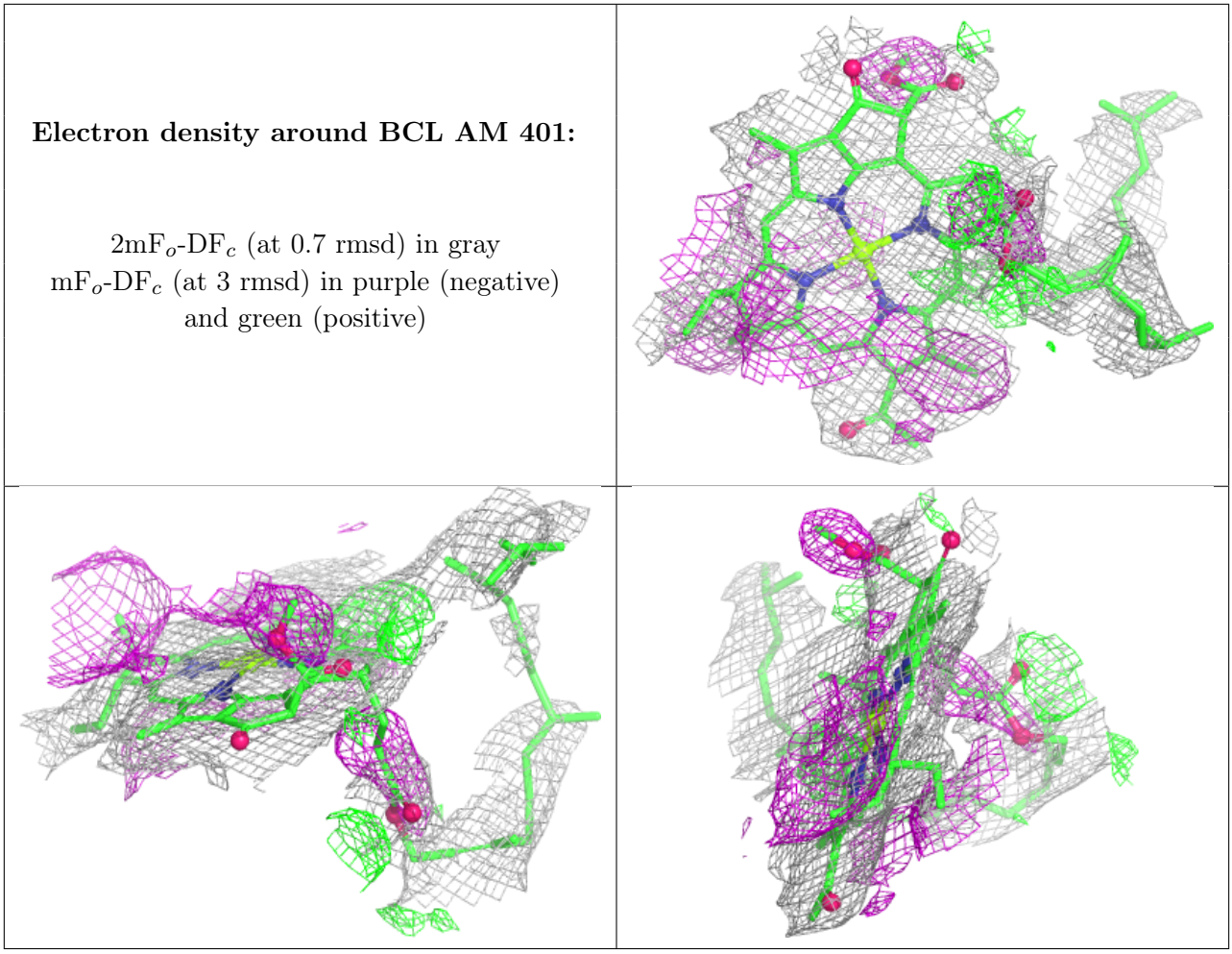
Electron density around BCL AL 303:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)



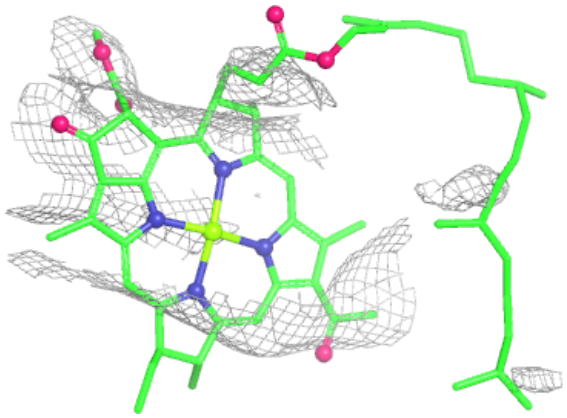
Electron density around BCL AM 401:

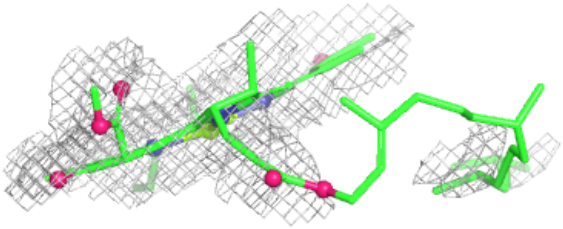
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

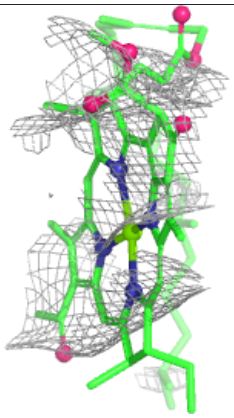


Electron density around BCL BG 101:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

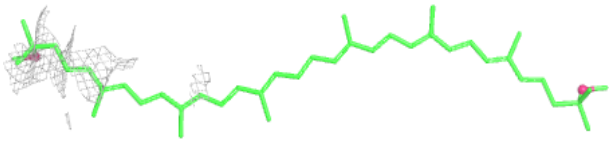


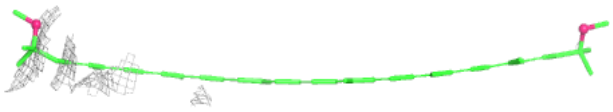


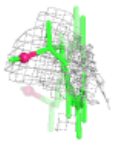


Electron density around CRT BS 103:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

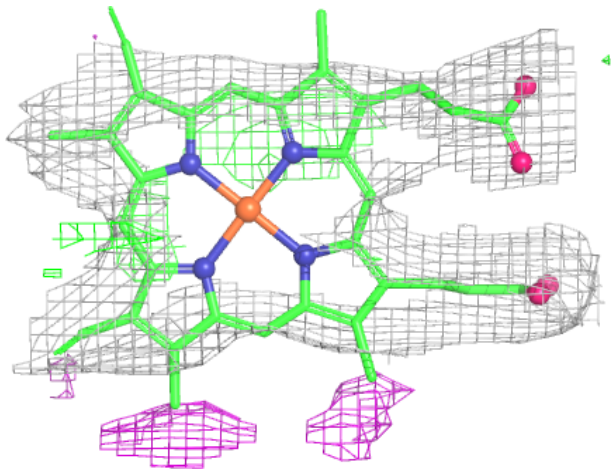


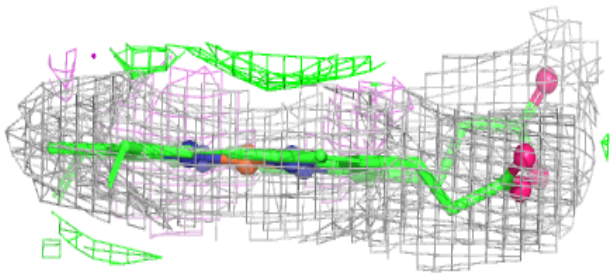


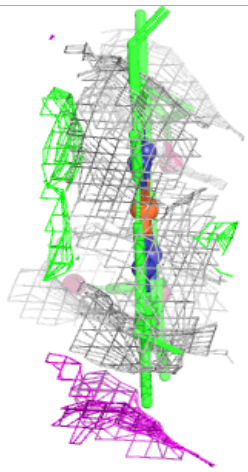


Electron density around HEM BC 504:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

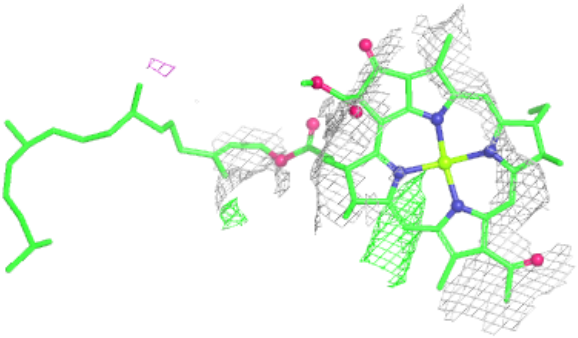


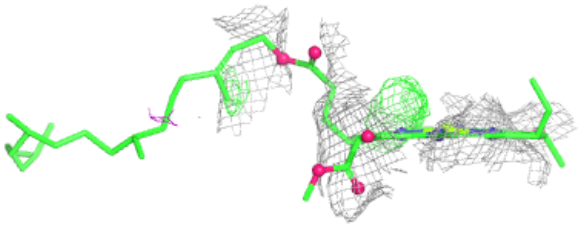


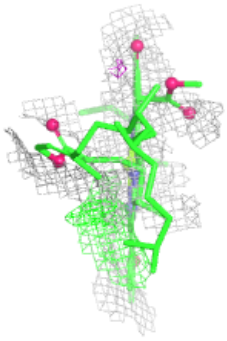


Electron density around BCL AA 101:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

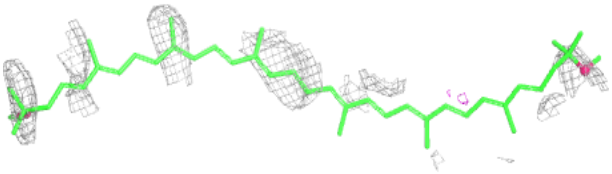


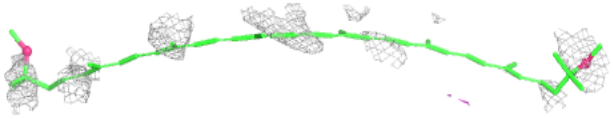


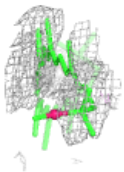


Electron density around CRT BW 103:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

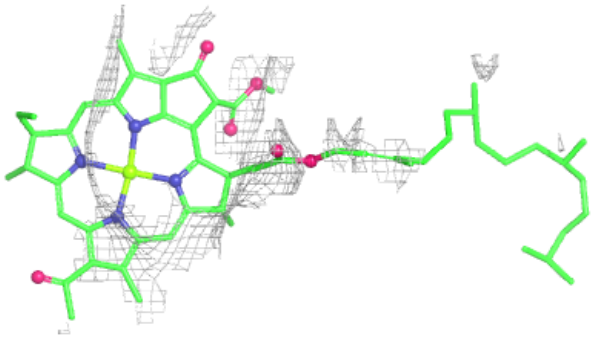


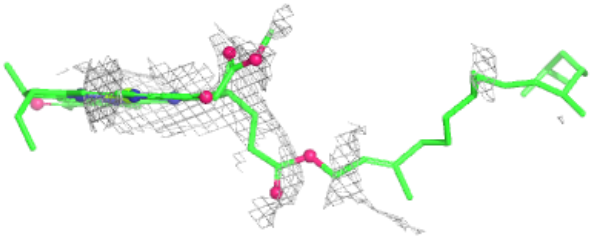


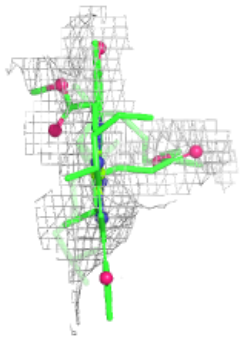


Electron density around BCL BK 102:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

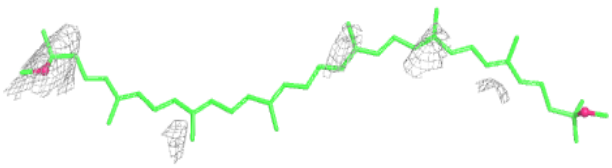


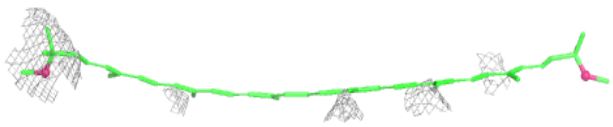


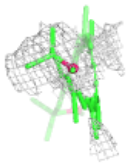


Electron density around CRT B2 102:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

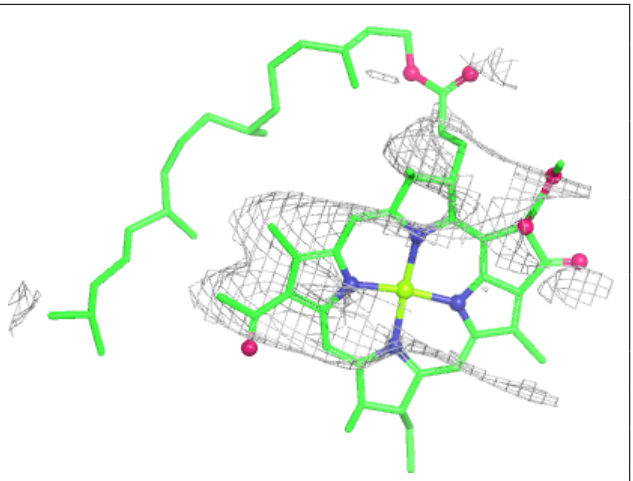


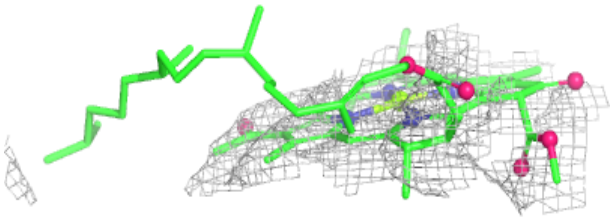


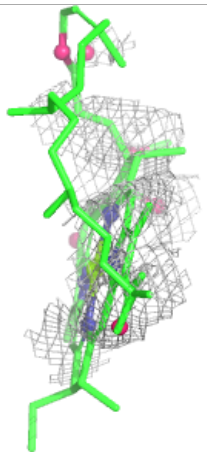


Electron density around BCL BN 101:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

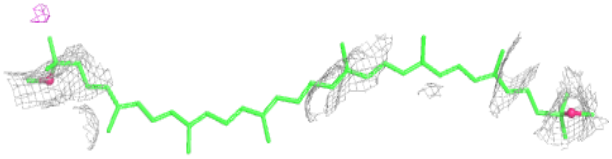


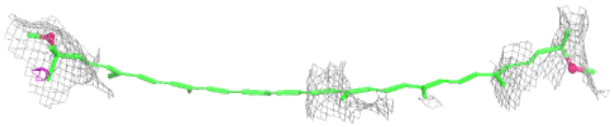


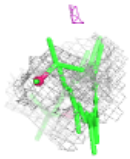


Electron density around CRT B7 102:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

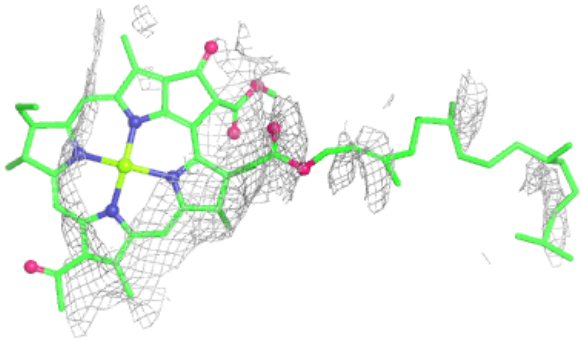


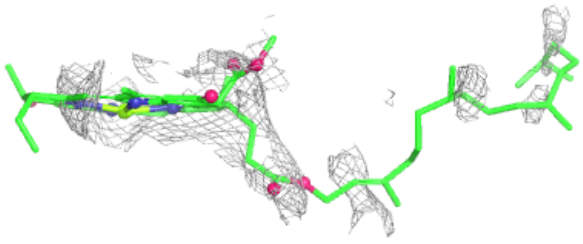


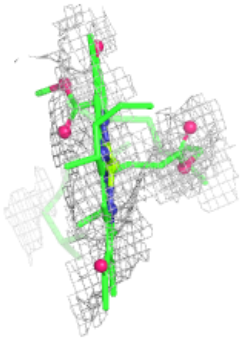


Electron density around BCL BO 102:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

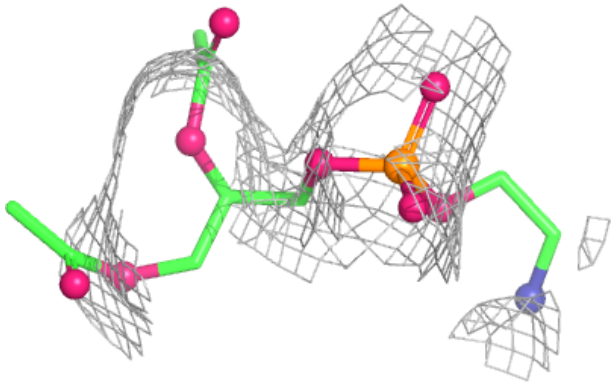


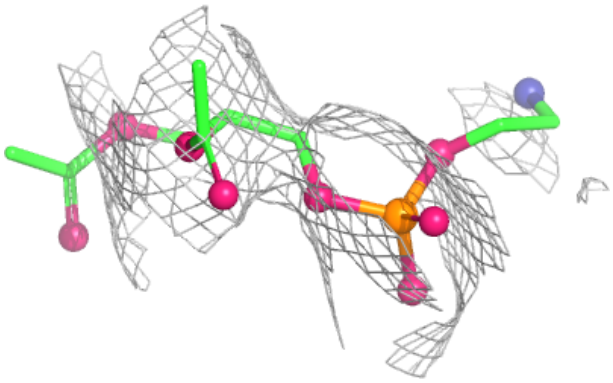


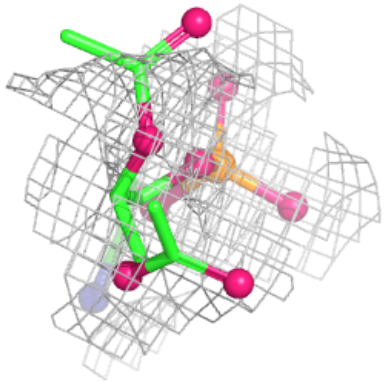


Electron density around PEF AM 407:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

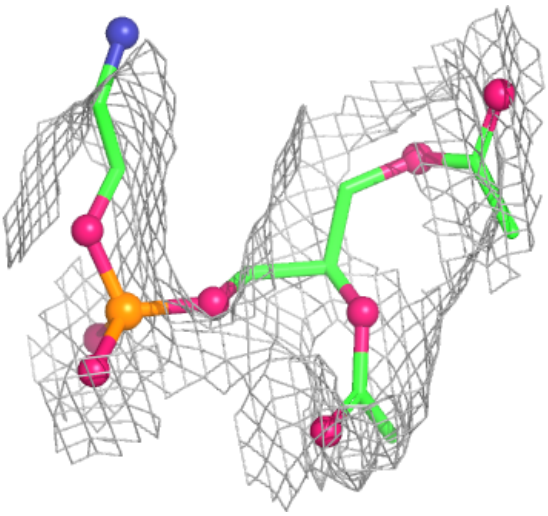


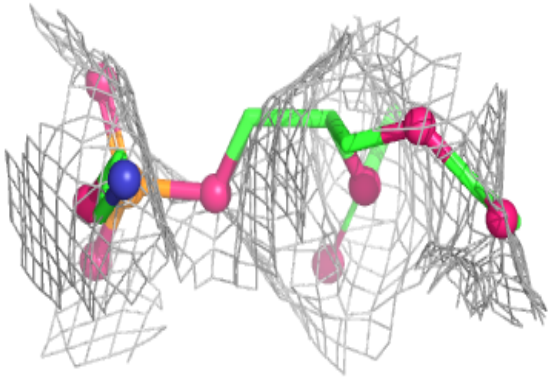


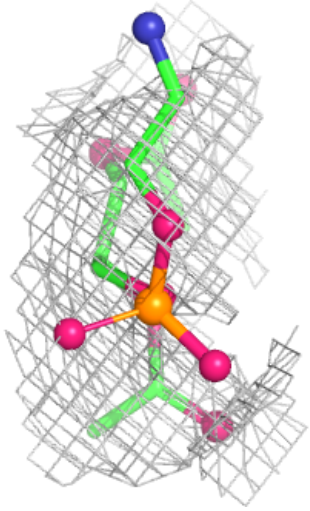


Electron density around PEF AH 301:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

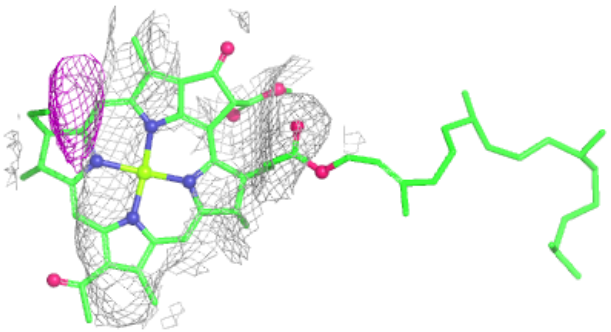


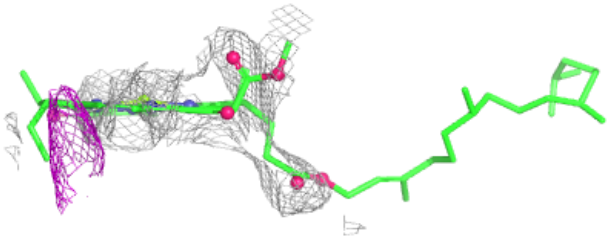


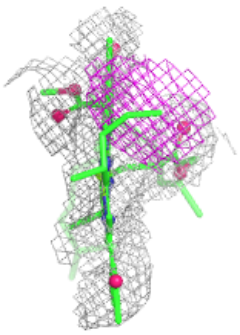


Electron density around BCL BQ 103:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)



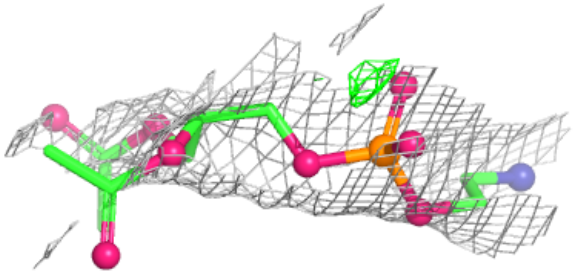


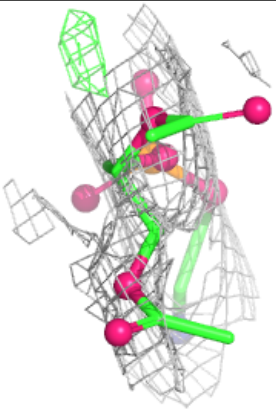


Electron density around PEF BM 407:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

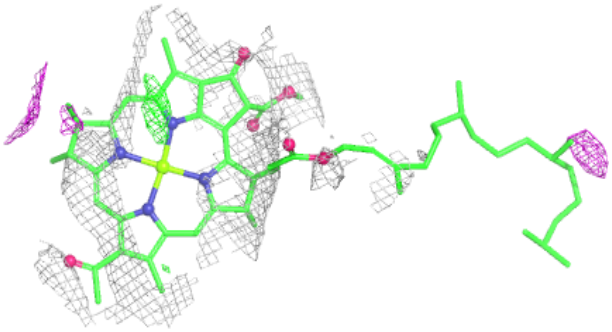


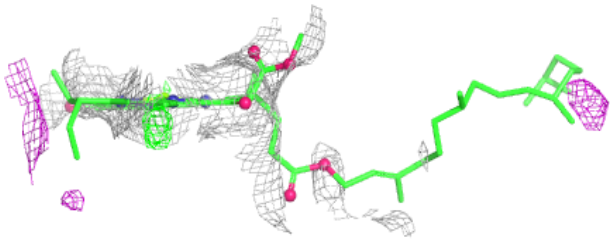




Electron density around BCL AD 102:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

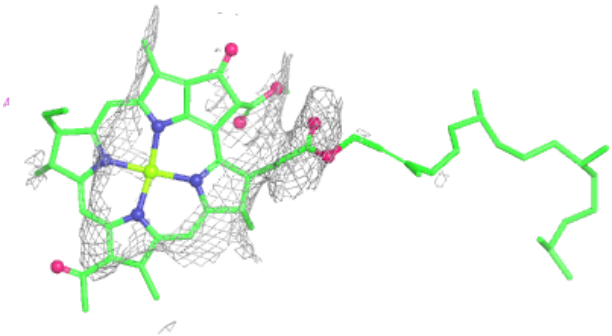


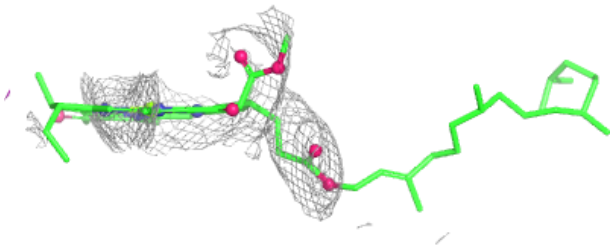


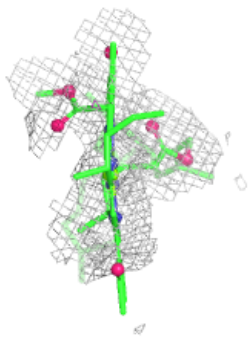


Electron density around BCL BS 102:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

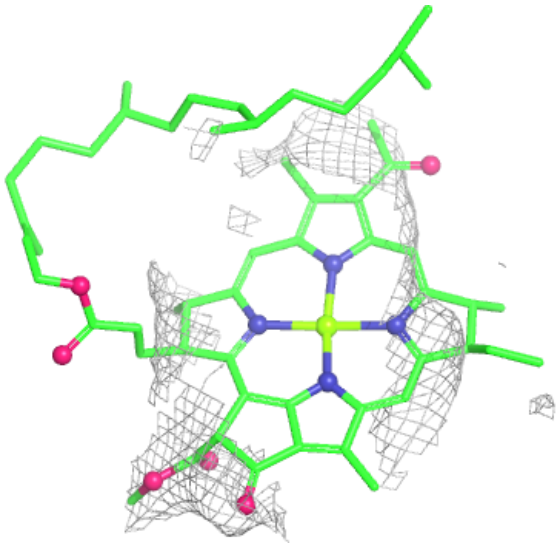


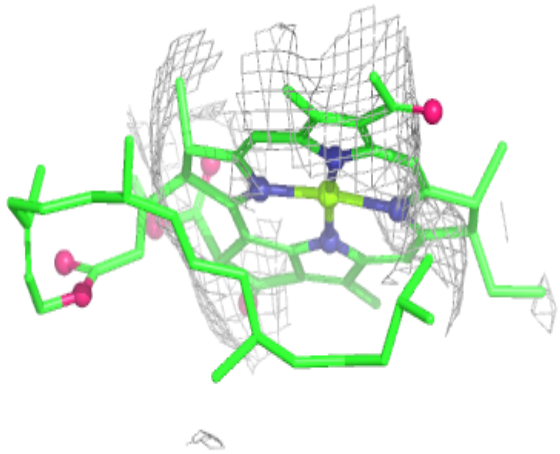




Electron density around BCL BE 101:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

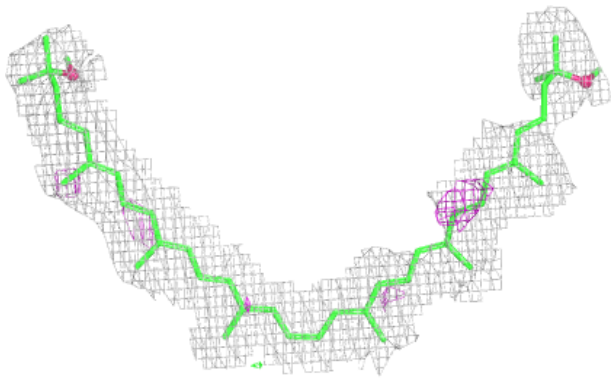


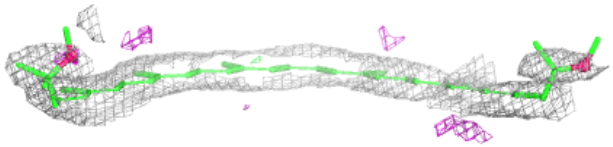


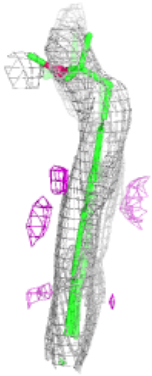


Electron density around CRT BM 406:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

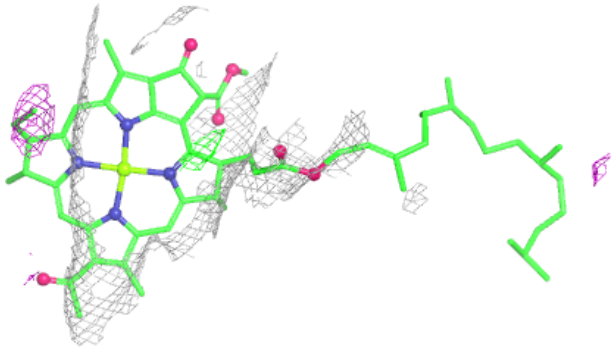


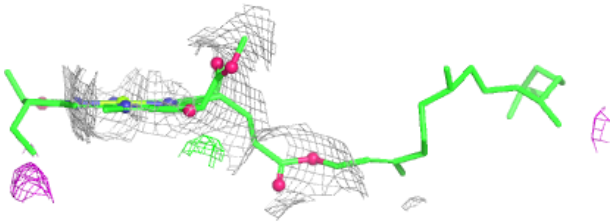


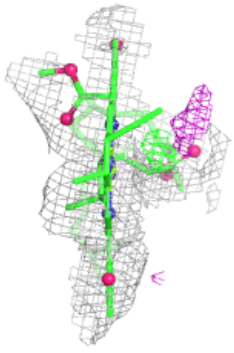


Electron density around BCL AU 102:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

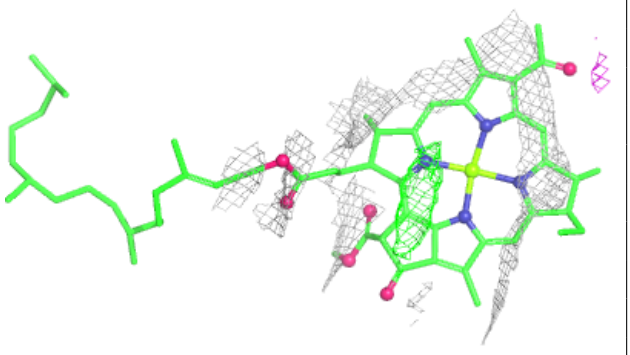


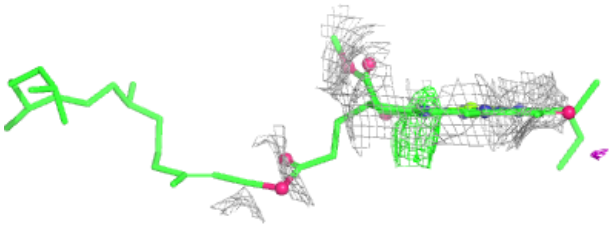




Electron density around BCL AW 101:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

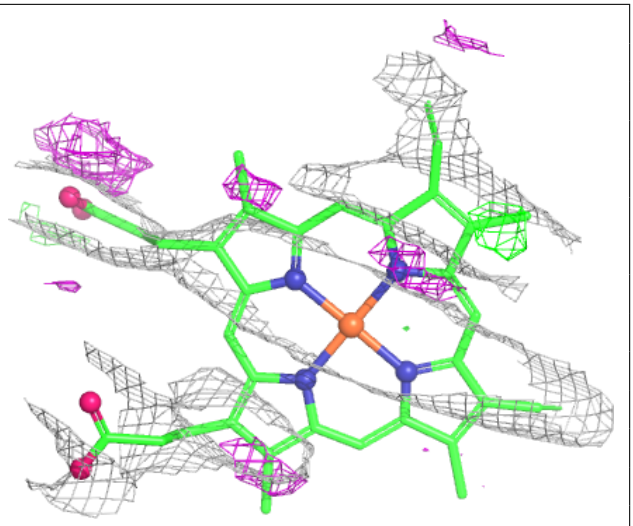


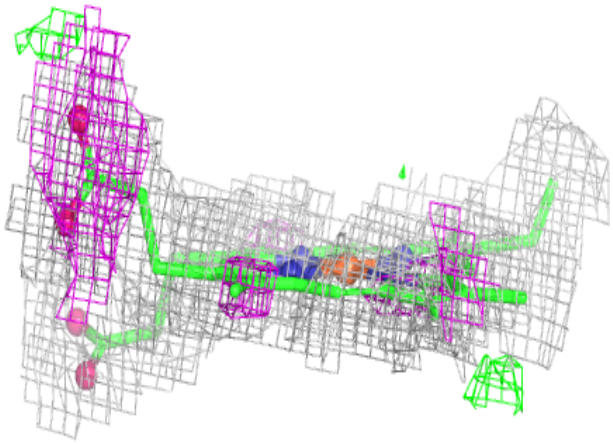


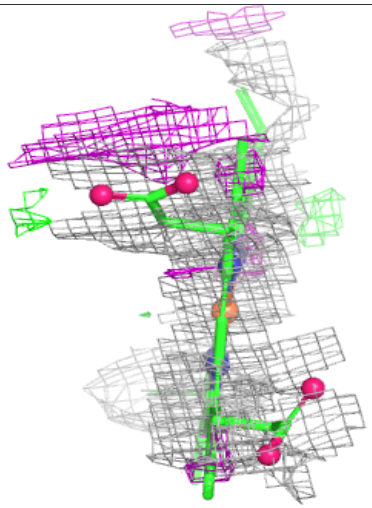


Electron density around HEM BC 501:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

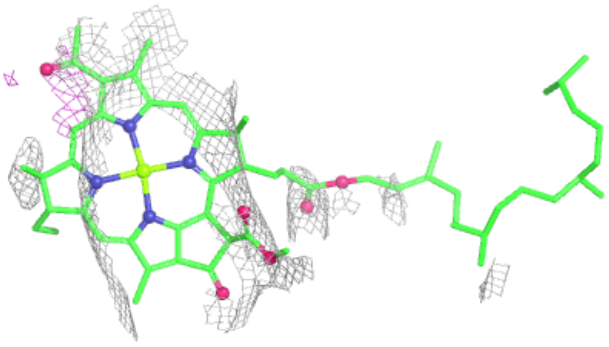


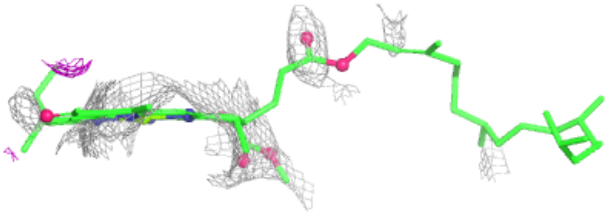


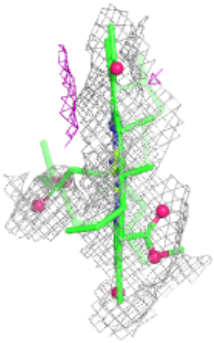


Electron density around BCL AY 102:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

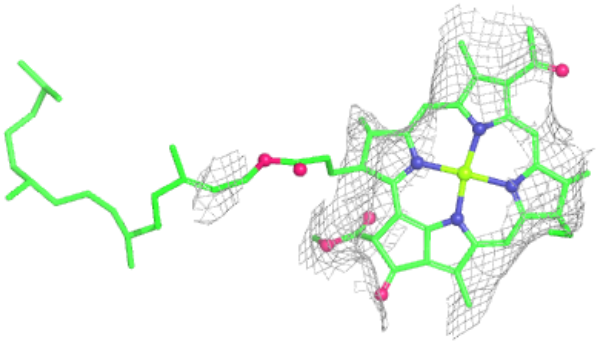


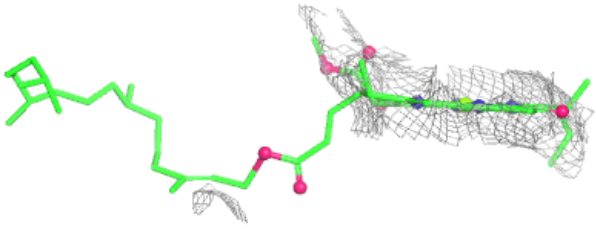


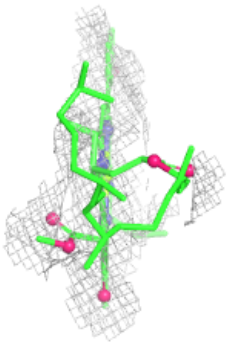


Electron density around BCL A1 102:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

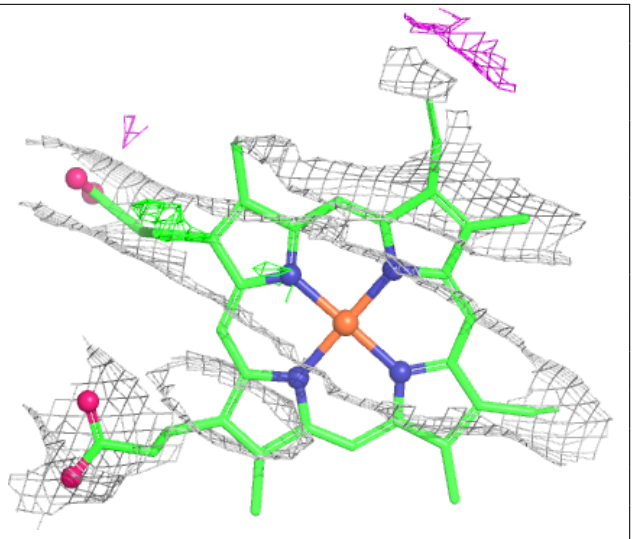


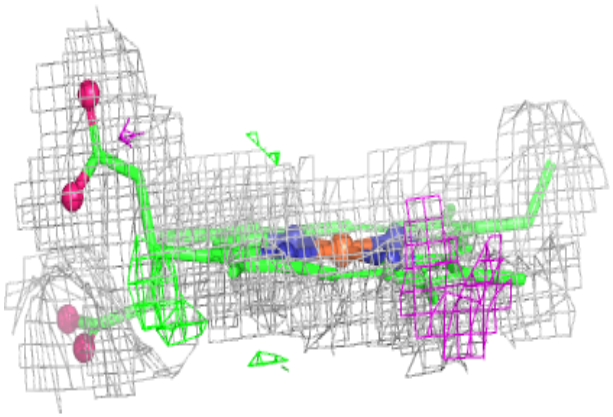




Electron density around HEM AC 501:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

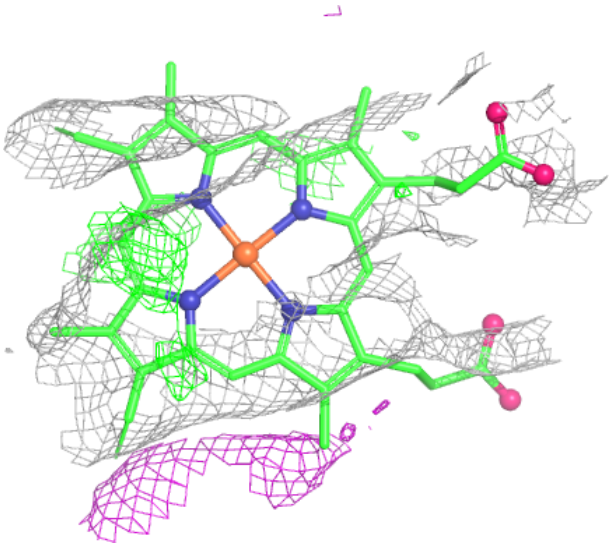


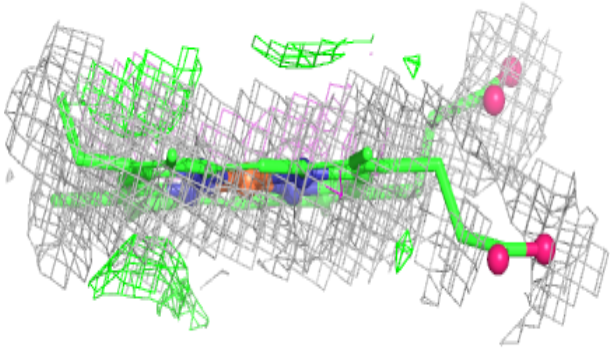


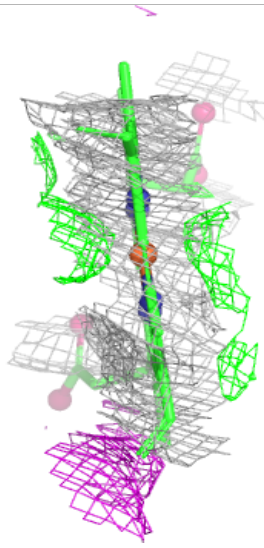


Electron density around HEM BC 503:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

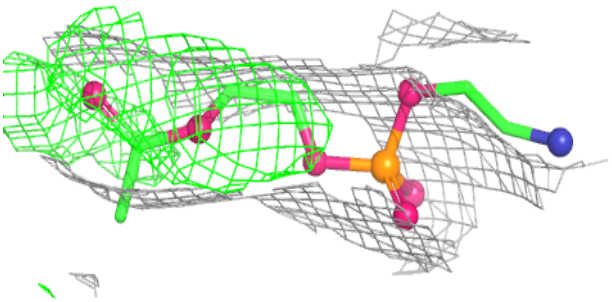


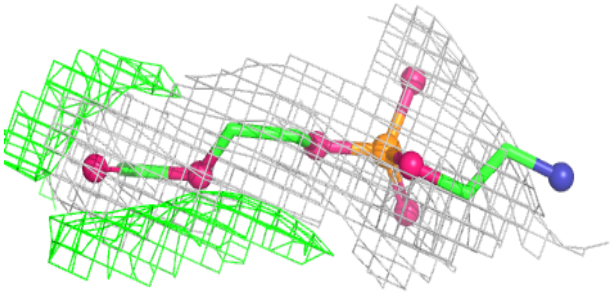


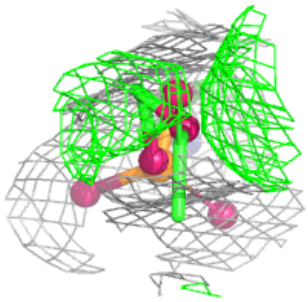


Electron density around PEF AM 408:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

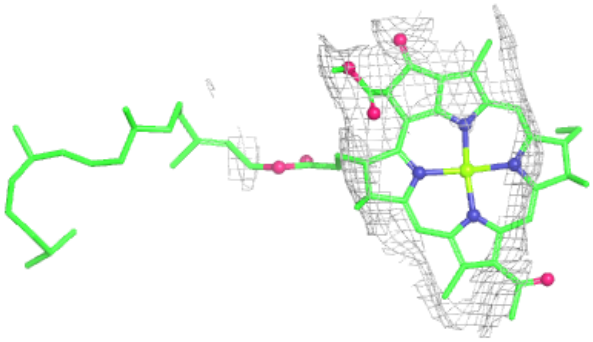


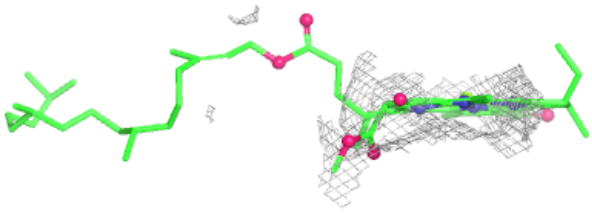




Electron density around BCL AO 102:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

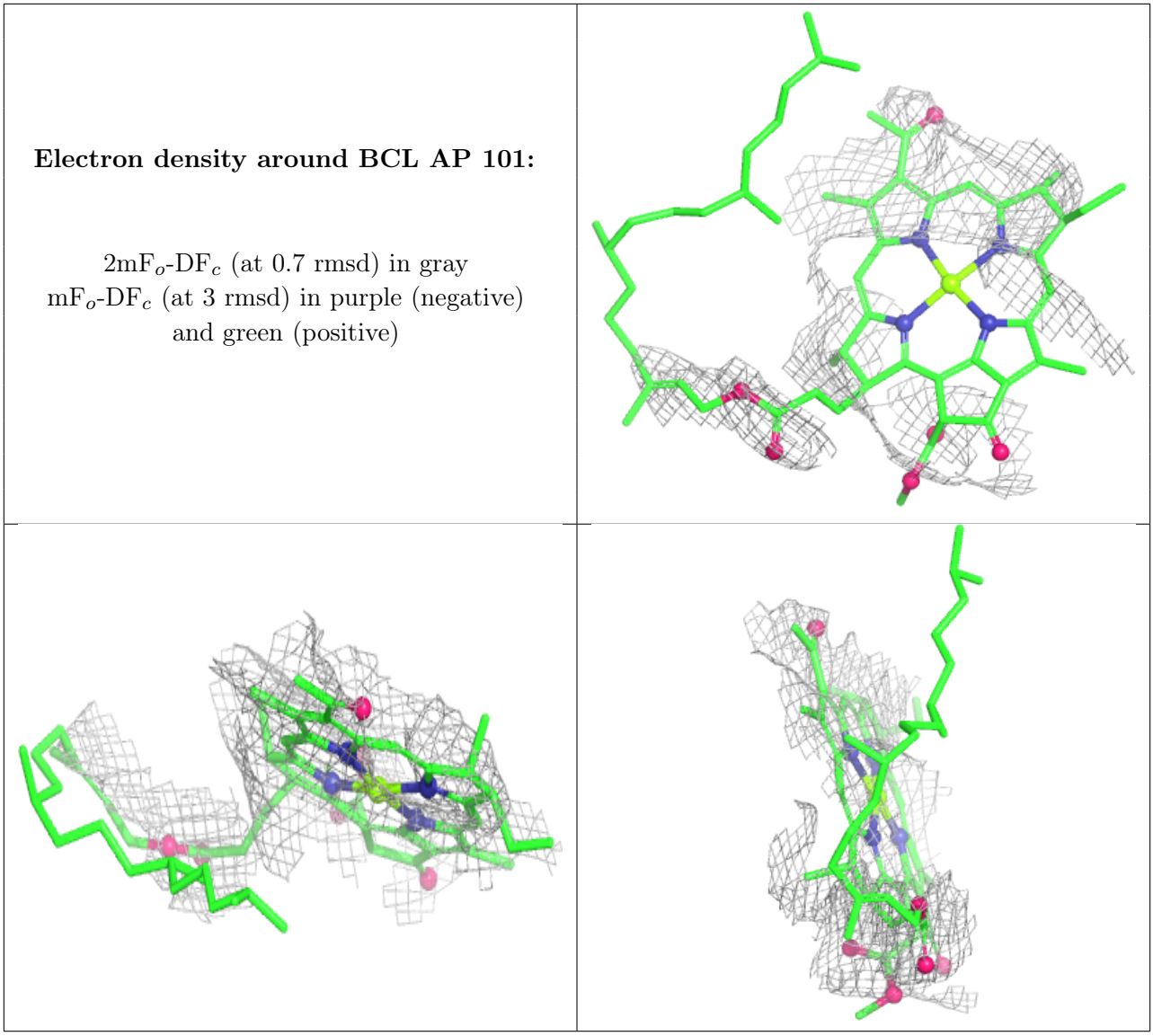






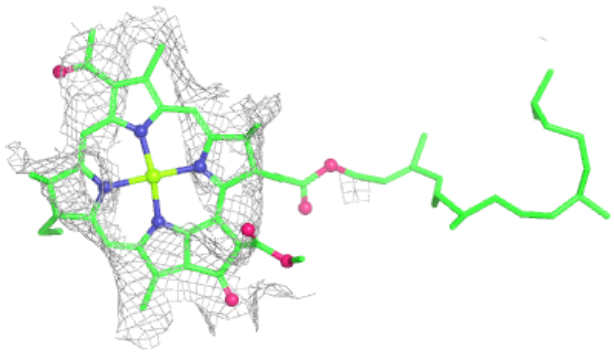
Electron density around BCL AP 101:

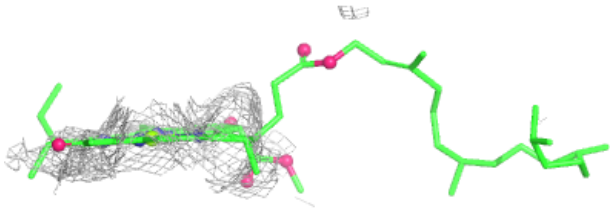
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

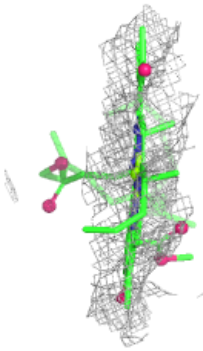


Electron density around BCL AQ 102:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

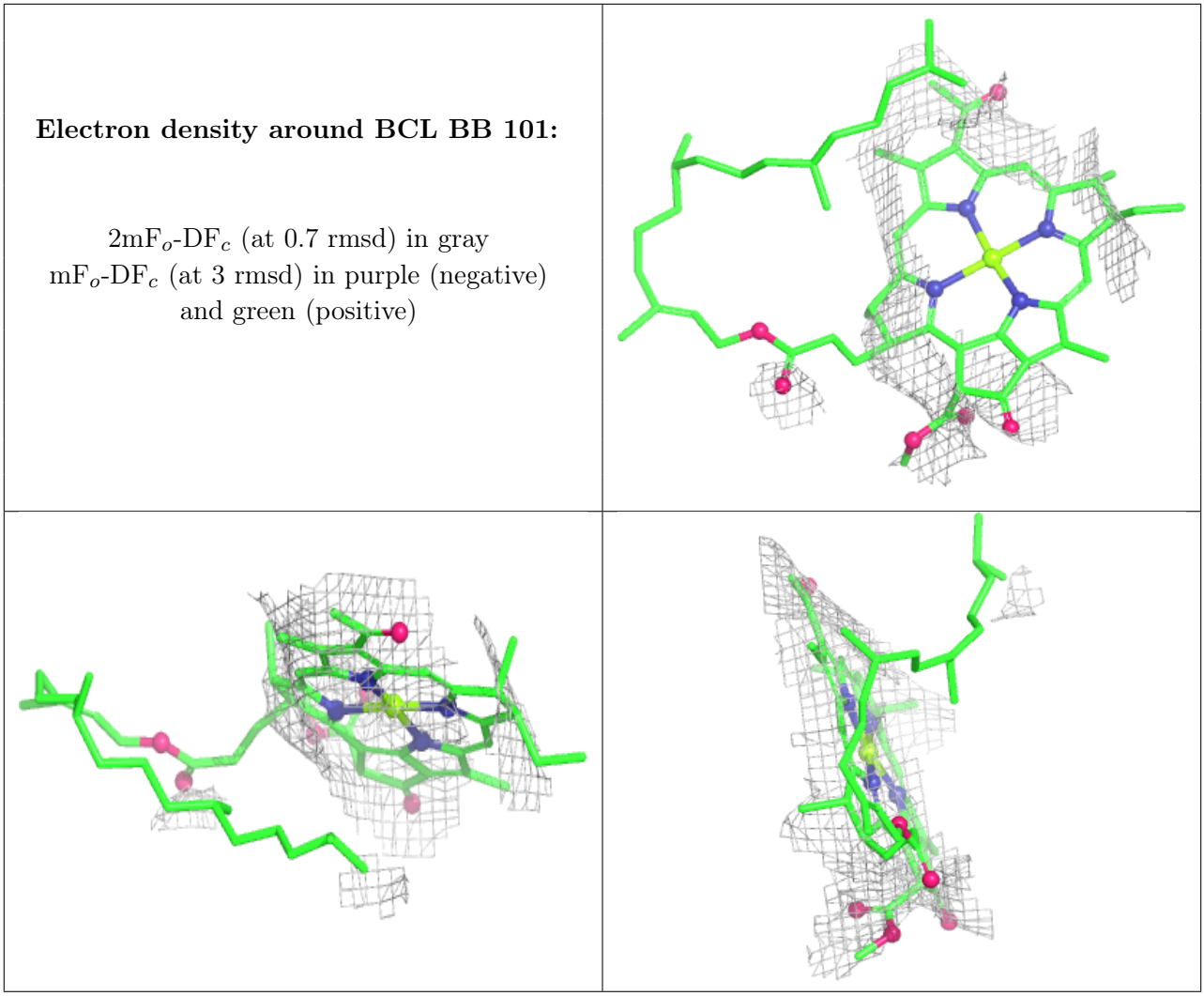






Electron density around BCL BB 101:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)



6.5 Other polymers

There are no such residues in this entry.