



Full wwPDB EM Validation Report ⓘ

Mar 8, 2026 – 04:15 AM UTC

PDB ID : 6M32 / pdb_00006m32
EMDB ID : EMD-30069
Title : Cryo-EM structure of FMO-RC complex from green sulfur bacteria
Authors : Chen, J.H.; Zhang, X.
Deposited on : 2020-03-02
Resolution : 2.70 Å(reported)

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A user guide is available at

<https://www.wwpdb.org/validation/2017/EMValidationReportHelp>

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:

EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

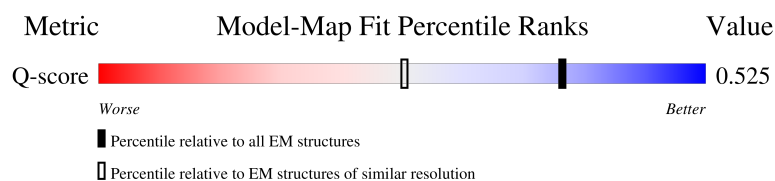
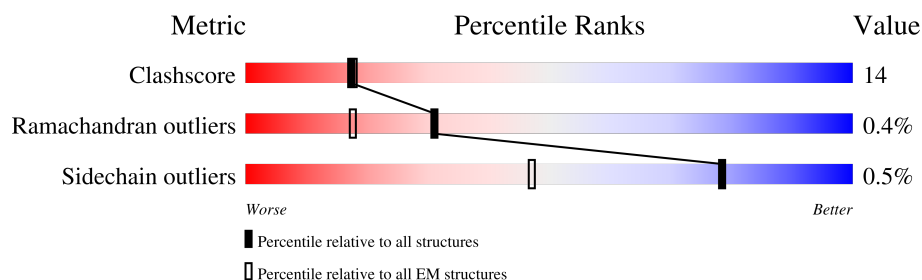
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 2.70 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.



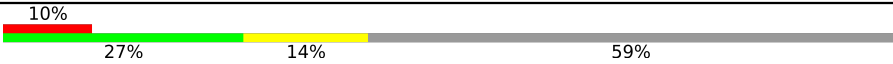
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	10327 (2.20 - 3.20)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the map. The red, orange, yellow and green segments of the 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 $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the EM map (all-atom inclusion $< 40\%$). The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	E	366	
1	F	366	
1	G	366	
2	D	143	

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Mol	Chain	Length	Quality of chain
3	B	231	
4	A	731	
4	a	731	

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
6	SF4	B	301	-	-	X	-
9	F39	A	817	-	X	-	-
9	F39	a	817	-	X	-	-

2 Entry composition

There are 13 unique types of molecules in this entry. The entry contains 23566 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, 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 Bacteriochlorophyll a protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	E	362	Total	C	N	O	S	0	0
			2815	1783	501	524	7		
1	F	358	Total	C	N	O	S	0	0
			2789	1770	496	516	7		
1	G	358	Total	C	N	O	S	0	0
			2789	1770	496	516	7		

- Molecule 2 is a protein called P840 reaction center 17 kDa protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	D	94	Total	C	N	O	S	0	0
			766	484	137	141	4		

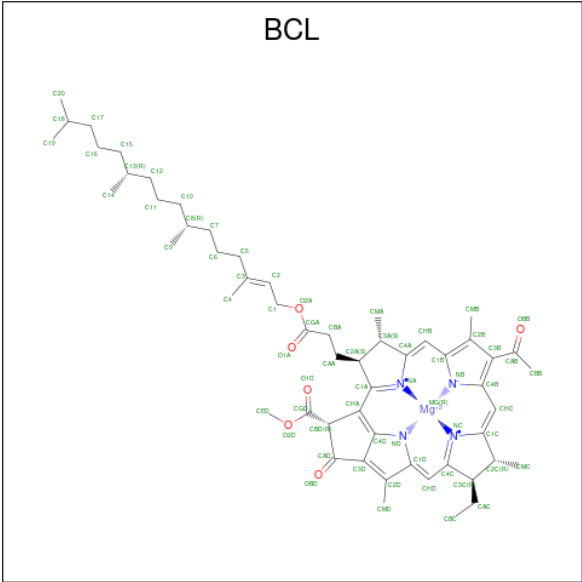
- Molecule 3 is a protein called Photosystem P840 reaction center iron-sulfur protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	B	95	Total	C	N	O	S	0	0
			731	469	115	138	9		

- Molecule 4 is a protein called Photosystem P840 reaction center, large subunit.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	A	628	Total	C	N	O	S	0	0
			5030	3360	798	846	26		
4	a	628	Total	C	N	O	S	0	0
			5030	3360	798	846	26		

- Molecule 5 is BACTERIOCHLOROPHYLL A (CCD ID: BCL) (formula: C₅₅H₇₄MgN₄O₆).



Mol	Chain	Residues	Atoms					AltConf
5	E	1	Total 66	C 55	Mg 1	N 4	O 6	0
5	E	1	Total 66	C 55	Mg 1	N 4	O 6	0
5	E	1	Total 66	C 55	Mg 1	N 4	O 6	0
5	E	1	Total 66	C 55	Mg 1	N 4	O 6	0
5	E	1	Total 66	C 55	Mg 1	N 4	O 6	0
5	E	1	Total 66	C 55	Mg 1	N 4	O 6	0
5	E	1	Total 46	C 35	Mg 1	N 4	O 6	1
5	E	1	Total 66	C 55	Mg 1	N 4	O 6	0
5	F	1	Total 66	C 55	Mg 1	N 4	O 6	0
5	F	1	Total 66	C 55	Mg 1	N 4	O 6	0
5	F	1	Total 66	C 55	Mg 1	N 4	O 6	0
5	F	1	Total 66	C 55	Mg 1	N 4	O 6	0
5	F	1	Total 66	C 55	Mg 1	N 4	O 6	0
5	F	1	Total 66	C 55	Mg 1	N 4	O 6	0

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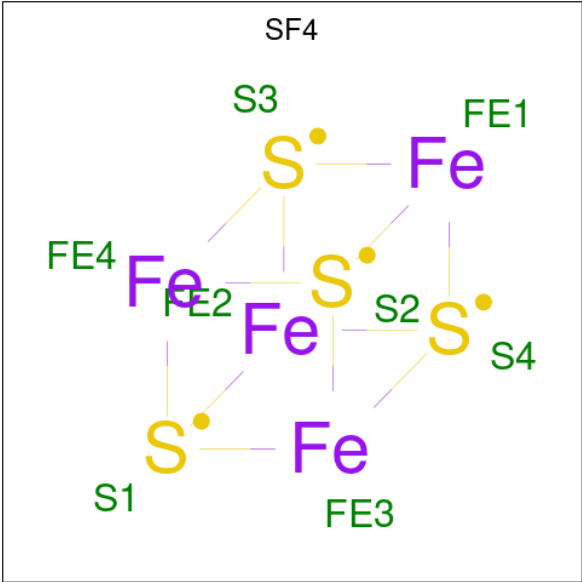
Mol	Chain	Residues	Atoms					AltConf
5	F	1	Total	C	Mg	N	O	0
			66	55	1	4	6	
5	F	1	Total	C	Mg	N	O	1
			46	35	1	4	6	
5	G	1	Total	C	Mg	N	O	1
			46	35	1	4	6	
5	G	1	Total	C	Mg	N	O	0
			66	55	1	4	6	
5	G	1	Total	C	Mg	N	O	0
			66	55	1	4	6	
5	G	1	Total	C	Mg	N	O	0
			66	55	1	4	6	
5	G	1	Total	C	Mg	N	O	0
			66	55	1	4	6	
5	G	1	Total	C	Mg	N	O	0
			66	55	1	4	6	
5	A	1	Total	C	Mg	N	O	0
			46	35	1	4	6	
5	A	1	Total	C	Mg	N	O	0
			56	45	1	4	6	
5	A	1	Total	C	Mg	N	O	0
			66	55	1	4	6	
5	A	1	Total	C	Mg	N	O	0
			46	35	1	4	6	
5	A	1	Total	C	Mg	N	O	0
			66	55	1	4	6	
5	A	1	Total	C	Mg	N	O	0
			46	35	1	4	6	
5	A	1	Total	C	Mg	N	O	0
			66	55	1	4	6	
5	A	1	Total	C	Mg	N	O	0
			66	55	1	4	6	
5	A	1	Total	C	Mg	N	O	0
			46	35	1	4	6	
5	A	1	Total	C	Mg	N	O	0
			46	35	1	4	6	

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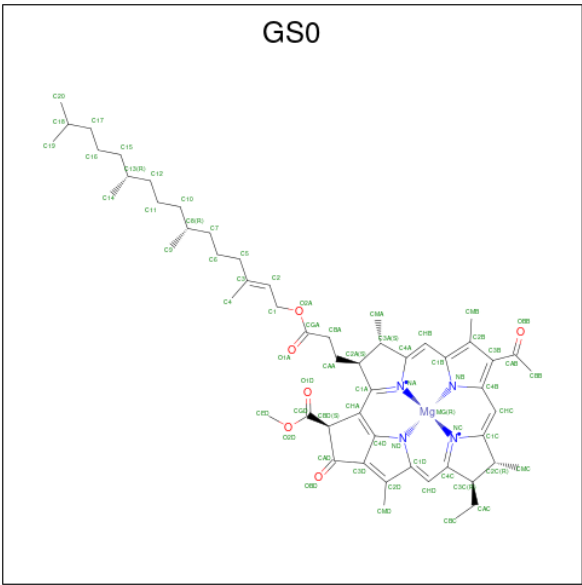
Mol	Chain	Residues	Atoms					AltConf
5	A	1	Total	C	Mg	N	O	0
			46	35	1	4	6	
5	a	1	Total	C	Mg	N	O	0
			46	35	1	4	6	
5	a	1	Total	C	Mg	N	O	0
			56	45	1	4	6	
5	a	1	Total	C	Mg	N	O	0
			66	55	1	4	6	
5	a	1	Total	C	Mg	N	O	0
			46	35	1	4	6	
5	a	1	Total	C	Mg	N	O	0
			66	55	1	4	6	
5	a	1	Total	C	Mg	N	O	0
			46	35	1	4	6	
5	a	1	Total	C	Mg	N	O	0
			66	55	1	4	6	
5	a	1	Total	C	Mg	N	O	0
			46	35	1	4	6	
5	a	1	Total	C	Mg	N	O	0
			46	35	1	4	6	
5	a	1	Total	C	Mg	N	O	0
			46	35	1	4	6	

- Molecule 6 is IRON/SULFUR CLUSTER (CCD ID: SF4) (formula: Fe₄S₄).



Mol	Chain	Residues	Atoms			AltConf
6	B	1	Total	Fe	S	0
			8	4	4	
6	B	1	Total	Fe	S	0
			8	4	4	
6	A	1	Total	Fe	S	0
			8	4	4	

- Molecule 7 is Bacteriochlorophyll A isomer (CCD ID: GS0) (formula: C₅₅H₇₄MgN₄O₆).



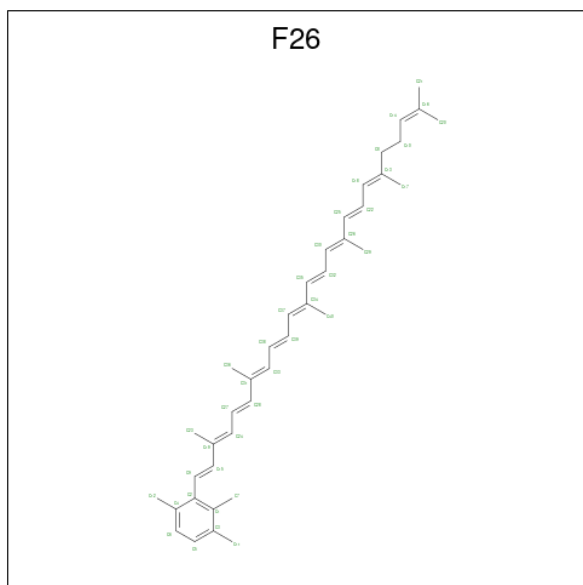
Mol	Chain	Residues	Atoms					AltConf
7	A	1	Total	C	Mg	N	O	0
			66	55	1	4	6	

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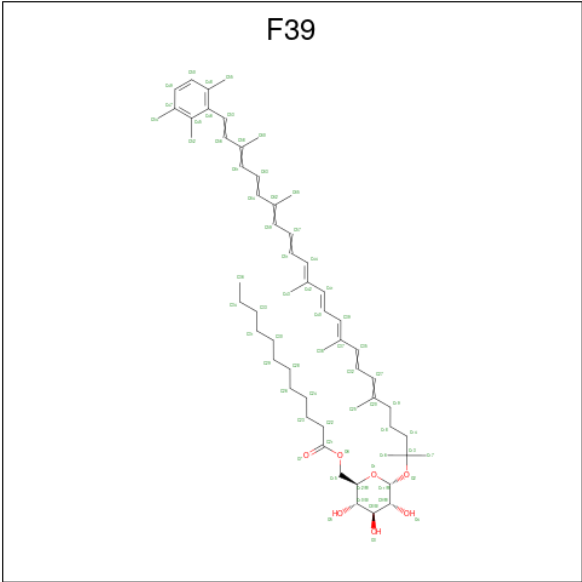
Mol	Chain	Residues	Atoms					AltConf
7	a	1	Total	C	Mg	N	O	0
			66	55	1	4	6	

- Molecule 8 is 2-[(1E,3E,5E,7E,9E,11E,13E,15E,17E,19E)-3,7,12,16,20,24-hexamethylpentacos-1,3,5,7,9,11,13,15,17,19,23-undecaenyl]-1,3,4-trimethyl-benzene (CCD ID: F26) (formula: C₄₀H₅₂).



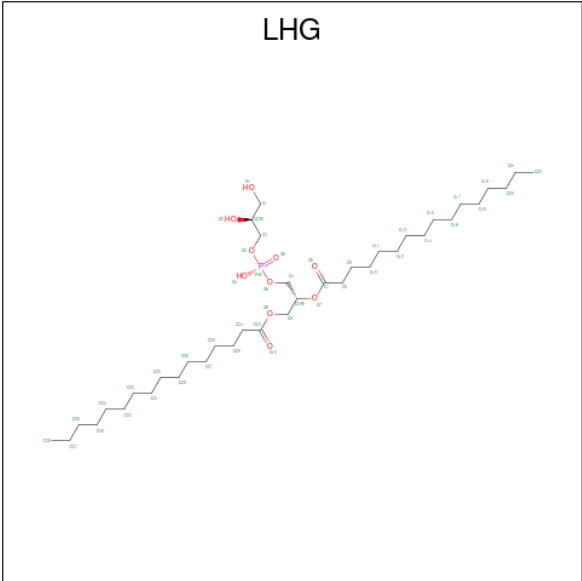
Mol	Chain	Residues	Atoms		AltConf
8	A	1	Total	C	0
			40	40	
8	a	1	Total	C	0
			40	40	

- Molecule 9 is [(2R,3S,4S,5R,6R)-6-[(10E,12E,14E)-2,6,10,14,19,23-hexamethyl-25-(2,3,6-trimethylphenyl)pentacos-6,8,10,12,14,16,18,20,22,24-decaen-2-yl]oxy-3,4,5-tris(oxidanyl)oxan-2-yl]methyl dodecanoate (CCD ID: F39) (formula: C₅₈H₈₆O₇).



Mol	Chain	Residues	Atoms			AltConf
9	A	1	Total	C	O	0
			65	58	7	
9	a	1	Total	C	O	0
			65	58	7	

- Molecule 10 is 1,2-DIPALMITOYL-PHOSPHATIDYL-GLYCEROLE (CCD ID: LHG) (formula: C₃₈H₇₅O₁₀P).



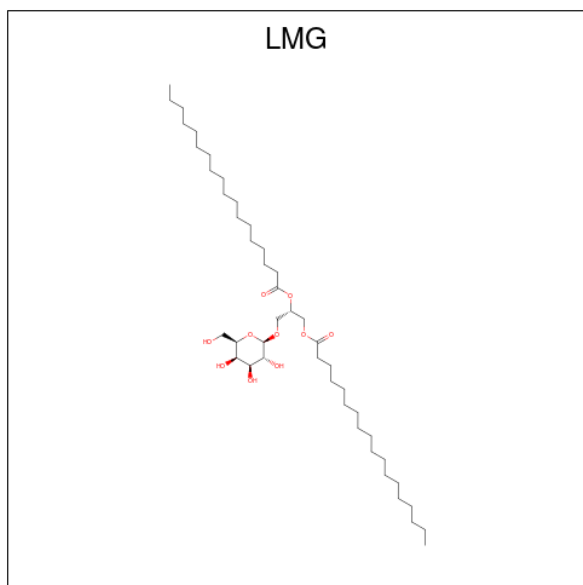
Mol	Chain	Residues	Atoms				AltConf
10	A	1	Total	C	O	P	0
			46	35	10	1	

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Mol	Chain	Residues	Atoms				AltConf
10	a	1	Total	C	O	P	0
			46	35	10	1	

- Molecule 11 is 1,2-DISTEAROYL-MONOGALACTOSYL-DIGLYCERIDE (CCD ID: LMG) (formula: $C_{45}H_{86}O_{10}$).

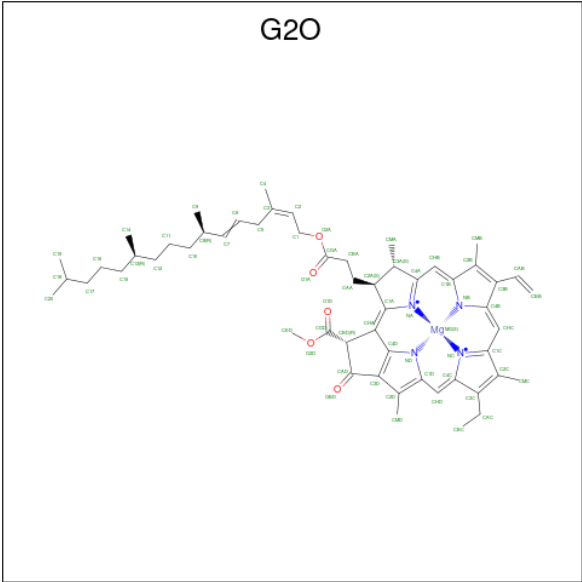


Mol	Chain	Residues	Atoms				AltConf
11	A	1	Total	C	O		0
			44	34	10		
11	a	1	Total	C	O		0
			44	34	10		

- Molecule 12 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		AltConf
12	A	1	Total	Ca	0
			1	1	
12	a	1	Total	Ca	0
			1	1	

- Molecule 13 is Chlorophyll A ester (CCD ID: G2O) (formula: $C_{55}H_{70}MgN_4O_5$).

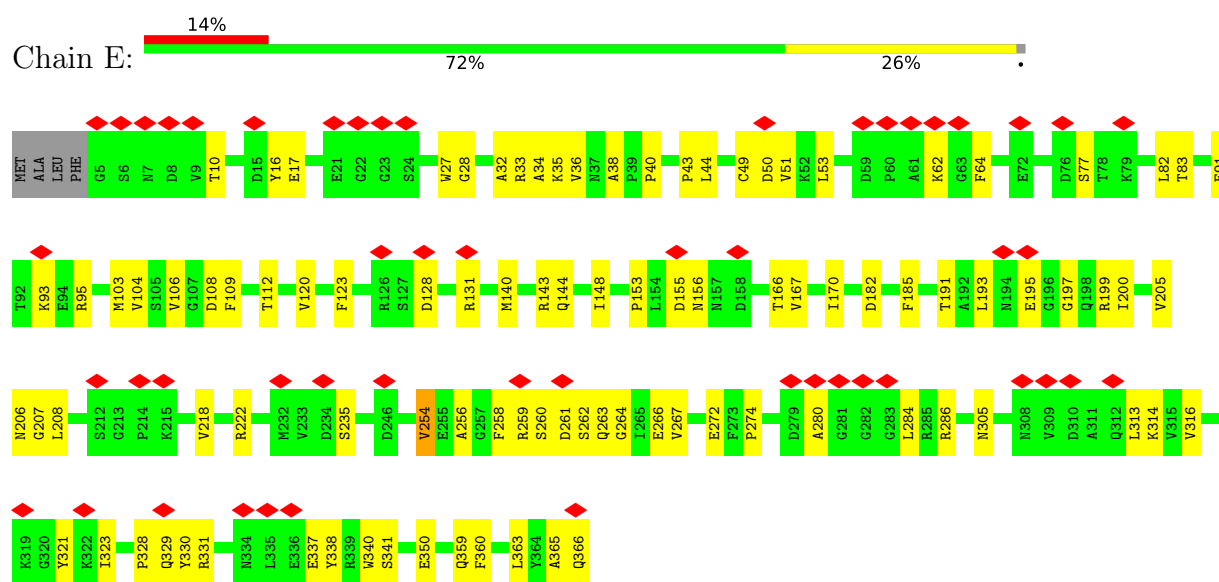


Mol	Chain	Residues	Atoms					AltConf
13	A	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
13	A	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
13	a	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
13	a	1	Total	C	Mg	N	O	0
			65	55	1	4	5	

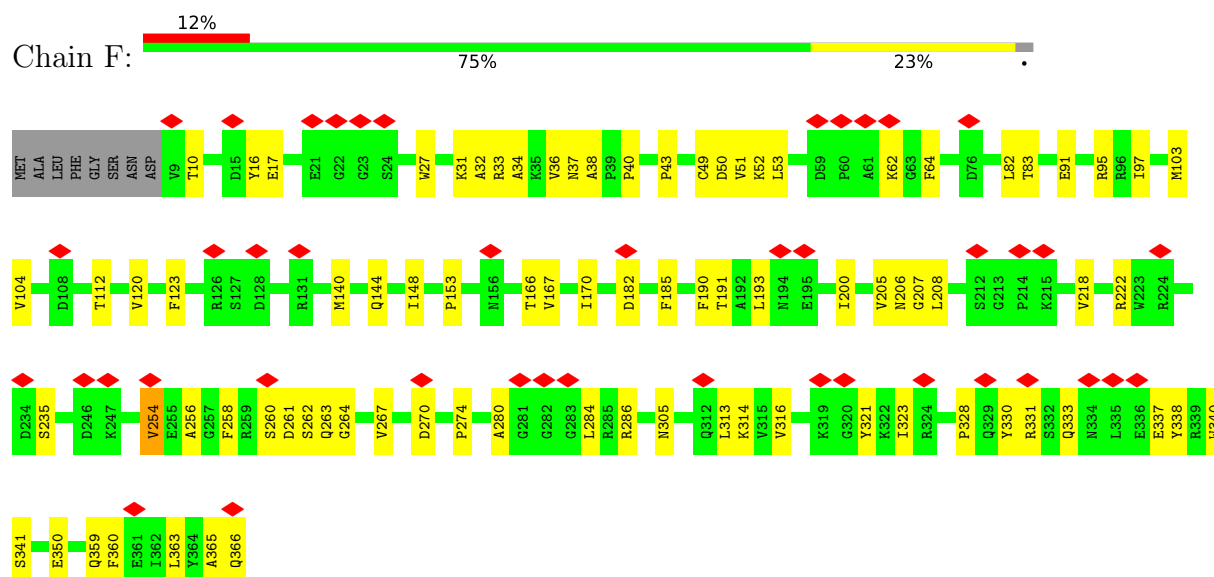
3 Residue-property plots

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 atom inclusion in map 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 diamond above a residue indicates a poor fit to the EM map for this residue (all-atom inclusion < 40%). 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.

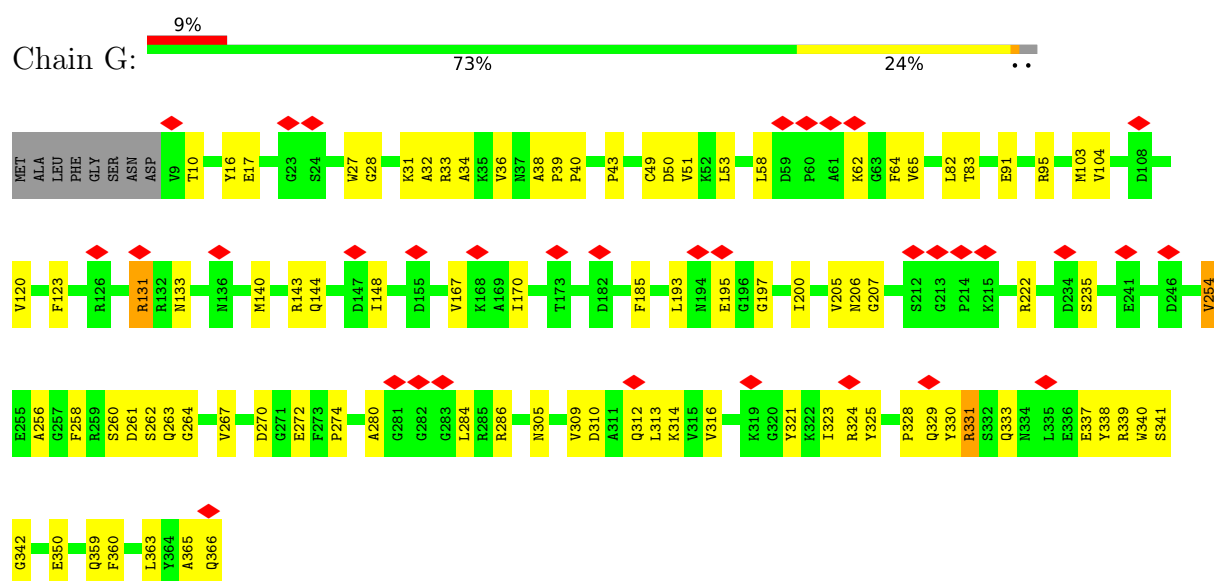
- Molecule 1: Bacteriochlorophyll a protein



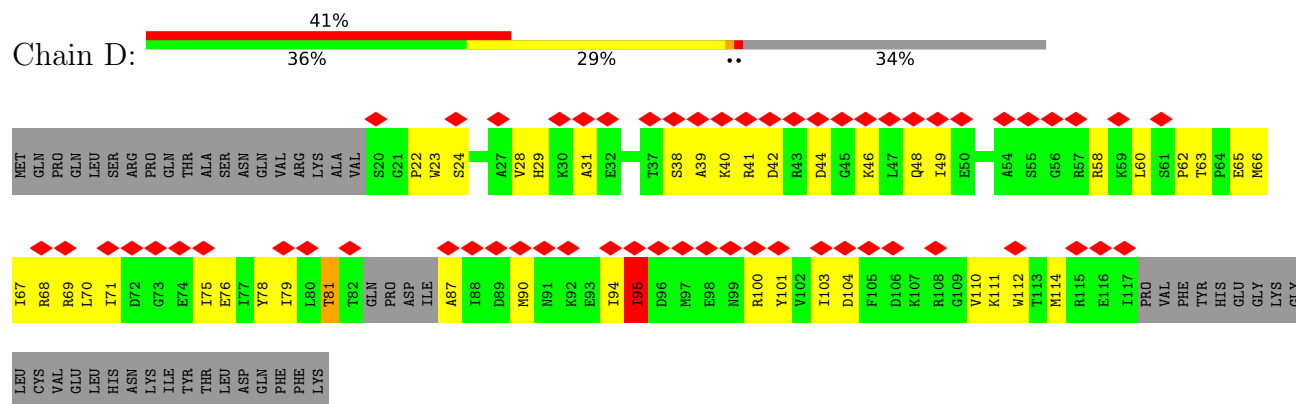
- Molecule 1: Bacteriochlorophyll a protein



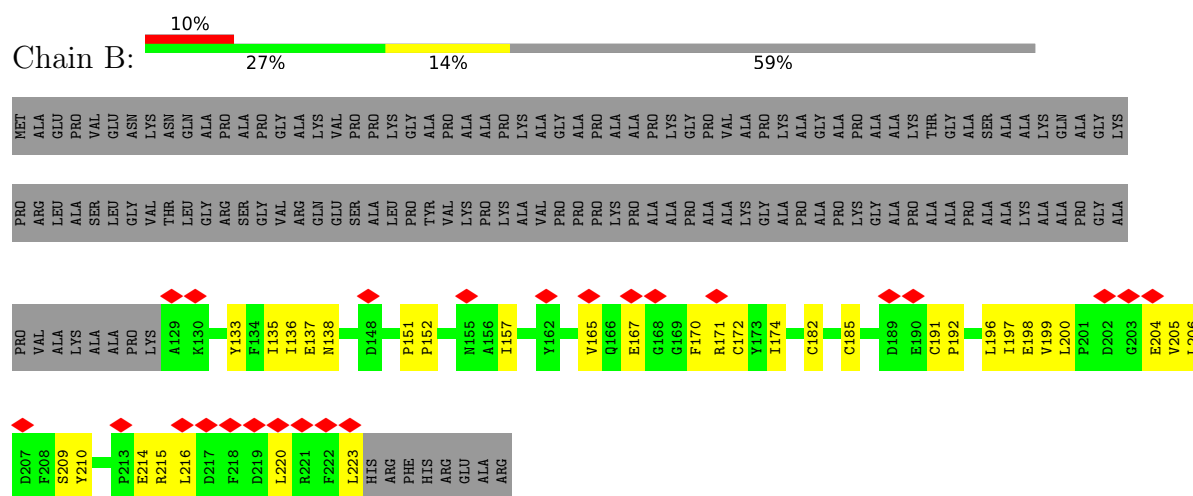
- Molecule 1: Bacteriochlorophyll a protein



- Molecule 2: P840 reaction center 17 kDa protein

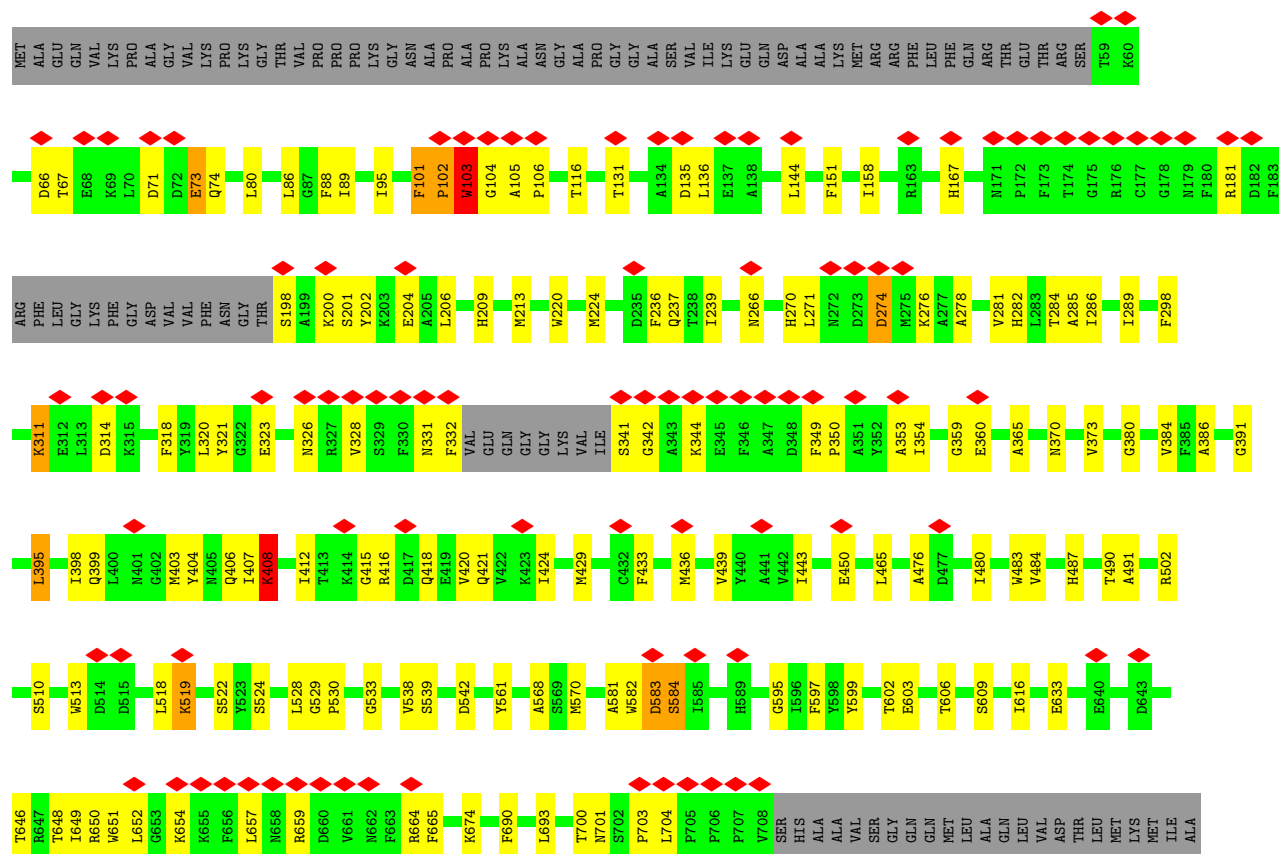


- Molecule 3: Photosystem P840 reaction center iron-sulfur protein

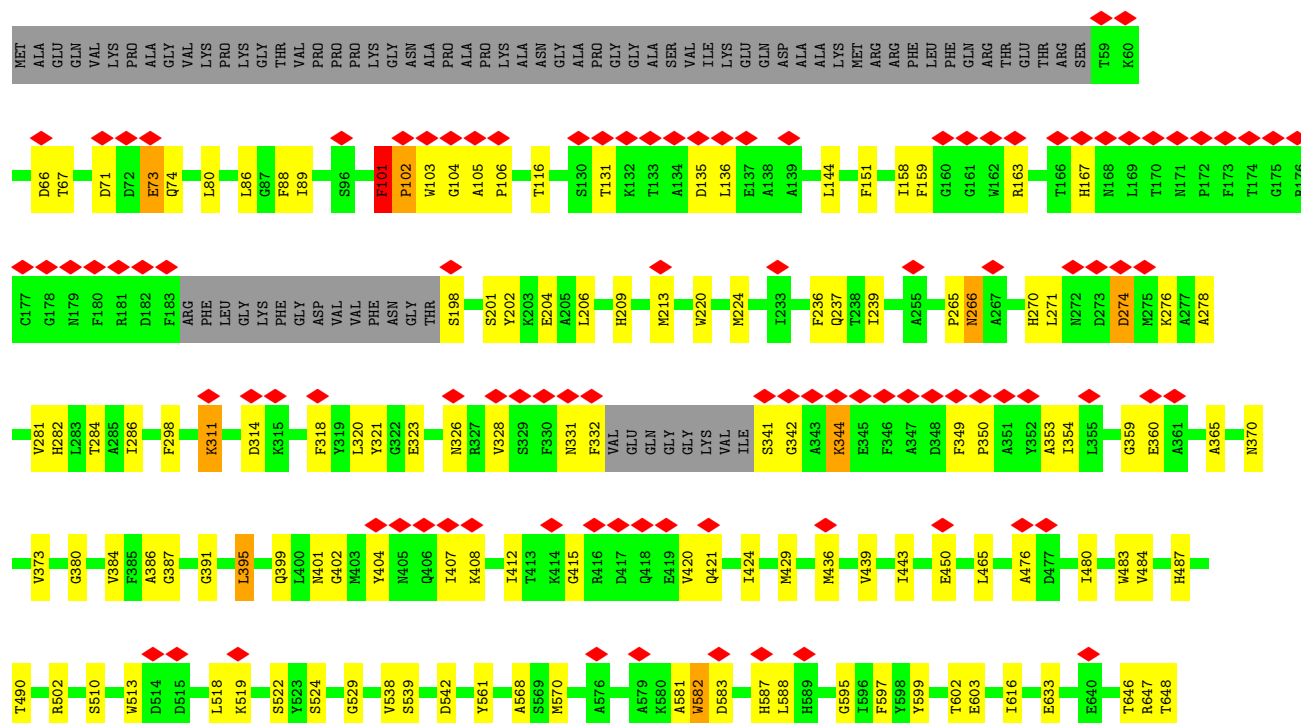


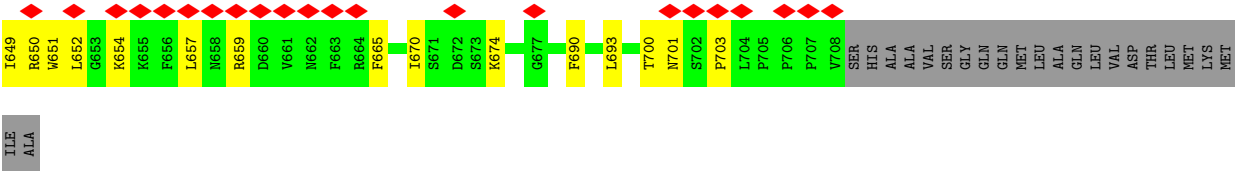
- Molecule 4: Photosystem P840 reaction center, large subunit





• Molecule 4: Photosystem P840 reaction center, large subunit





4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	268430	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	NONE	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	47	Depositor
Minimum defocus (nm)	1500	Depositor
Maximum defocus (nm)	2500	Depositor
Magnification	38244	Depositor
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	0.163	Depositor
Minimum map value	-0.048	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.009	Depositor
Recommended contour level	0.0341	Depositor
Map size ($\text{\AA}$)	209.12001, 209.12001, 209.12001	wwPDB
Map dimensions	160, 160, 160	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	1.307, 1.307, 1.307	Depositor

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: LMG, G2O, F26, GS0, LHG, SF4, F39, BCL, CA

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	E	0.24	0/2885	0.52	1/3910 (0.0%)
1	F	0.23	0/2859	0.53	1/3875 (0.0%)
1	G	0.24	0/2859	0.53	1/3875 (0.0%)
2	D	0.22	0/778	0.64	1/1043 (0.1%)
3	B	0.25	0/749	0.55	0/1013
4	A	0.22	0/5207	0.58	9/7097 (0.1%)
4	a	0.22	0/5207	0.55	5/7097 (0.1%)
All	All	0.23	0/20544	0.55	18/27910 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
3	B	0	1
4	A	0	3
4	a	0	3
All	All	0	7

There are no bond length outliers.

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	a	101	PHE	CA-C-N	10.50	132.97	119.84
4	a	101	PHE	C-N-CA	10.50	132.97	119.84
4	A	101	PHE	CA-C-N	9.56	131.79	119.84
4	A	101	PHE	C-N-CA	9.56	131.79	119.84
4	a	266	ASN	N-CA-C	9.27	126.56	107.37
4	A	266	ASN	N-CA-C	9.26	126.55	107.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	103	TRP	N-CA-C	-8.44	102.77	113.23
4	a	103	TRP	N-CA-C	-8.37	102.85	113.23
4	A	583	ASP	N-CA-C	-5.93	101.07	109.96
4	a	274	ASP	N-CA-C	5.78	119.95	113.01
4	A	274	ASP	N-CA-C	5.68	119.83	113.01
2	D	95	ILE	N-CA-C	5.62	121.03	109.34
4	A	584	SER	CA-C-N	-5.40	111.68	120.64
4	A	584	SER	C-N-CA	-5.40	111.68	120.64
1	E	254	VAL	N-CA-C	-5.34	100.37	108.17
1	F	254	VAL	N-CA-C	-5.34	100.37	108.17
1	G	254	VAL	N-CA-C	-5.33	100.38	108.17
4	A	408	LYS	N-CA-C	-5.04	107.31	113.50

There are no chirality outliers.

All (7) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
4	A	311	LYS	Peptide
4	A	344	LYS	Peptide
4	A	519	LYS	Peptide
3	B	216	LEU	Peptide
4	a	311	LYS	Peptide
4	a	344	LYS	Peptide
4	a	519	LYS	Peptide

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	E	2815	0	2749	83	0
1	F	2789	0	2731	67	0
1	G	2789	0	2731	88	0
2	D	766	0	783	54	0
3	B	731	0	692	35	0
4	A	5030	0	4916	146	0
4	a	5030	0	4916	121	0
5	A	642	0	584	31	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	E	508	0	551	25	0
5	F	508	0	551	21	0
5	G	508	0	551	25	0
5	a	642	0	584	27	0
6	A	8	0	0	0	0
6	B	16	0	0	4	0
7	A	66	0	0	0	0
7	a	66	0	0	0	0
8	A	40	0	0	0	0
8	a	40	0	0	0	0
9	A	65	0	0	0	0
9	a	65	0	0	0	0
10	A	46	0	65	0	0
10	a	46	0	65	0	0
11	A	44	0	57	3	0
11	a	44	0	57	3	0
12	A	1	0	0	0	0
12	a	1	0	0	0	0
13	A	130	0	0	1	0
13	a	130	0	0	2	0
All	All	23566	0	22583	627	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 14.

All (627) 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)
4:a:583:ASP:OD1	4:a:588:LEU:HB2	1.39	1.18
4:a:202:TYR:CE1	4:a:274:ASP:OD2	2.06	1.09
1:G:331:ARG:NH2	4:A:271:LEU:O	1.87	1.07
1:E:259:ARG:HH11	3:B:215:ARG:HH12	1.05	0.95
1:G:341:SER:OG	4:A:200:LYS:NZ	2.01	0.93
4:A:450:GLU:OE2	4:A:582:TRP:NE1	2.05	0.89
4:a:450:GLU:OE2	4:a:582:TRP:NE1	2.06	0.89
4:A:101:PHE:HB3	4:A:103:TRP:NE1	1.88	0.88
4:A:350:PRO:HB2	4:A:353:ALA:HB2	1.55	0.88
4:A:651:TRP:HD1	4:A:654:LYS:HD3	1.39	0.88
4:A:101:PHE:O	4:A:104:GLY:N	2.08	0.87
4:a:651:TRP:HD1	4:a:654:LYS:HD3	1.39	0.87
4:a:350:PRO:HB2	4:a:353:ALA:HB2	1.55	0.86

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:E:371:BCL:HAA1	5:E:371:BCL:HBD	1.60	0.84
5:F:371:BCL:HAA1	5:F:371:BCL:HBD	1.60	0.84
2:D:79:ILE:HB	2:D:111:LYS:HG3	1.59	0.84
3:B:137:GLU:O	3:B:171:ARG:NH1	2.11	0.84
4:a:202:TYR:HE1	4:a:274:ASP:OD2	1.60	0.84
5:G:371:BCL:HBD	5:G:371:BCL:HAA1	1.60	0.83
4:a:599:TYR:CZ	4:a:603:GLU:OE1	2.31	0.83
1:E:259:ARG:NH1	3:B:215:ARG:HH12	1.77	0.82
4:a:101:PHE:O	4:a:104:GLY:N	2.13	0.82
4:A:599:TYR:CZ	4:A:603:GLU:OE1	2.32	0.82
1:G:83:THR:CG2	1:G:103:MET:HG2	2.10	0.81
1:F:95:ARG:NH2	1:F:97:ILE:HD11	1.96	0.80
2:D:79:ILE:HB	2:D:111:LYS:CG	2.11	0.80
4:A:510:SER:HB3	4:A:513:TRP:HD1	1.47	0.80
2:D:114:MET:HE2	3:B:136:ILE:HG12	1.64	0.79
2:D:95:ILE:HG12	2:D:95:ILE:O	1.81	0.79
4:a:265:PRO:O	4:a:266:ASN:OD1	2.00	0.79
4:a:201:SER:HB2	4:a:271:LEU:HD23	1.65	0.79
4:a:510:SER:HB3	4:a:513:TRP:HD1	1.47	0.78
4:a:649:ILE:HD13	4:a:665:PHE:CE2	2.19	0.78
2:D:79:ILE:HB	2:D:111:LYS:CD	2.13	0.78
2:D:40:LYS:HB2	2:D:48:GLN:HG3	1.65	0.78
4:A:583:ASP:OD1	4:A:704:LEU:HD23	1.84	0.77
4:A:101:PHE:HB3	4:A:103:TRP:HE1	1.48	0.77
4:a:602:THR:O	4:a:603:GLU:HB2	1.85	0.77
1:E:259:ARG:HH11	3:B:215:ARG:NH1	1.83	0.76
1:F:51:VAL:HG12	1:F:53:LEU:CD1	2.16	0.76
4:A:602:THR:O	4:A:603:GLU:HB2	1.86	0.76
1:E:259:ARG:CG	1:E:266:GLU:HB3	2.16	0.75
1:G:324:ARG:HH12	3:B:209:SER:HB3	1.50	0.74
1:E:260:SER:O	1:E:261:ASP:OD1	2.07	0.73
1:G:260:SER:O	1:G:261:ASP:OD1	2.07	0.73
1:F:37:ASN:OD1	1:F:263:GLN:O	2.06	0.73
3:B:199:VAL:HG12	3:B:205:VAL:HG22	1.70	0.73
2:D:46:LYS:HE3	2:D:49:ILE:HG12	1.70	0.72
4:A:201:SER:OG	4:A:204:GLU:OE1	2.06	0.72
4:a:646:THR:O	4:a:650:ARG:HB2	1.89	0.72
4:a:583:ASP:OD2	4:a:703:PRO:O	2.09	0.71
4:a:583:ASP:OD1	4:a:588:LEU:CB	2.29	0.70
4:a:284:THR:OG1	4:a:387:GLY:HA3	1.92	0.70
1:G:312:GLN:HE22	1:G:339:ARG:HH11	1.40	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:94:ILE:HD12	2:D:100:ARG:HG2	1.73	0.70
4:a:651:TRP:CD1	4:a:654:LYS:HD3	2.26	0.70
1:E:259:ARG:HG2	1:E:266:GLU:HB3	1.72	0.70
1:G:83:THR:HG22	1:G:103:MET:HG2	1.73	0.69
4:a:201:SER:OG	4:a:204:GLU:OE1	2.06	0.69
4:A:513:TRP:HZ2	4:A:538:VAL:HG11	1.58	0.69
4:a:513:TRP:HZ2	4:a:538:VAL:HG11	1.58	0.68
1:G:324:ARG:HH12	3:B:209:SER:CB	2.07	0.68
1:E:254:VAL:CG1	5:E:375:BCL:H72	2.23	0.68
5:G:373:BCL:H42	5:G:375:BCL:H151	1.76	0.68
5:E:373:BCL:H42	5:E:375:BCL:H151	1.76	0.67
4:A:651:TRP:CD1	4:A:654:LYS:HD3	2.26	0.67
1:G:312:GLN:NE2	1:G:339:ARG:HH11	1.93	0.67
1:G:254:VAL:CG1	5:G:375:BCL:H72	2.24	0.67
4:A:386:ALA:HB2	5:A:813:BCL:HMC3	1.77	0.67
4:A:649:ILE:HD12	4:A:665:PHE:CE2	2.29	0.67
4:A:66:ASP:OD1	4:A:67:THR:N	2.28	0.66
4:A:649:ILE:HD12	4:A:665:PHE:CZ	2.30	0.66
4:a:66:ASP:OD1	4:a:67:THR:N	2.28	0.66
4:a:386:ALA:HB2	5:a:813:BCL:HMC3	1.77	0.66
4:a:570:MET:HE2	4:a:597:PHE:HE2	1.61	0.66
5:F:373:BCL:H42	5:F:375:BCL:H151	1.76	0.66
1:G:329:GLN:NE2	4:A:398:ILE:HG22	2.11	0.66
1:G:254:VAL:HG11	5:G:375:BCL:H72	1.77	0.66
4:A:570:MET:HE2	4:A:597:PHE:HE2	1.61	0.65
1:E:254:VAL:HG11	5:E:375:BCL:H72	1.77	0.65
5:A:806:BCL:HBB3	5:A:808:BCL:H91	1.79	0.65
5:a:806:BCL:HBB3	5:a:808:BCL:H91	1.79	0.65
2:D:95:ILE:O	2:D:95:ILE:CG1	2.41	0.65
4:a:490:THR:HG22	11:a:819:LMG:H381	1.79	0.64
4:a:649:ILE:CD1	4:a:665:PHE:CE2	2.79	0.64
1:G:254:VAL:HG23	5:G:376:BCL:H18	1.80	0.64
1:F:261:ASP:OD1	1:F:262:SER:N	2.28	0.64
3:B:196:LEU:CD1	6:B:301:SF4:S1	2.85	0.64
1:E:77:SER:O	2:D:58:ARG:NH2	2.31	0.64
4:a:649:ILE:HD13	4:a:665:PHE:HE2	1.63	0.64
2:D:103:ILE:HG13	2:D:104:ASP:H	1.62	0.64
4:A:490:THR:HG22	11:A:819:LMG:H381	1.79	0.64
1:F:51:VAL:CG1	1:F:53:LEU:CD1	2.76	0.63
3:B:196:LEU:HD12	6:B:301:SF4:S1	2.38	0.63
1:E:16:TYR:HE1	5:E:374:BCL:H43	1.63	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:16:TYR:HE1	5:G:374:BCL:H43	1.63	0.63
1:F:16:TYR:HE1	5:F:374:BCL:H43	1.63	0.63
1:E:144:GLN:NE2	1:E:235:SER:O	2.32	0.63
2:D:79:ILE:HB	2:D:111:LYS:HD2	1.79	0.63
4:a:599:TYR:CE2	4:a:603:GLU:OE1	2.51	0.63
1:F:144:GLN:NE2	1:F:235:SER:O	2.32	0.63
1:G:312:GLN:HE22	1:G:339:ARG:NH1	1.97	0.62
4:A:599:TYR:CE2	4:A:603:GLU:OE1	2.52	0.62
1:G:144:GLN:NE2	1:G:235:SER:O	2.32	0.62
4:A:690:PHE:HB2	5:A:811:BCL:H42	1.82	0.62
1:F:262:SER:O	1:F:263:GLN:HG3	2.00	0.62
1:F:254:VAL:CG1	5:F:375:BCL:H72	2.30	0.62
1:E:10:THR:HG22	1:E:10:THR:O	2.00	0.62
4:a:649:ILE:CD1	4:a:665:PHE:HE2	2.13	0.62
2:D:112:TRP:CE3	3:B:138:ASN:HB2	2.34	0.61
4:a:690:PHE:HB2	5:a:811:BCL:H42	1.82	0.61
2:D:94:ILE:HG22	2:D:95:ILE:HG22	1.82	0.61
2:D:49:ILE:HD12	2:D:62:PRO:HG3	1.82	0.61
1:E:262:SER:O	1:E:263:GLN:HG3	2.00	0.61
4:A:135:ASP:OD1	4:A:136:LEU:N	2.34	0.61
1:G:36:VAL:O	1:G:264:GLY:HA3	2.01	0.61
1:G:262:SER:O	1:G:263:GLN:HG3	2.00	0.61
1:E:36:VAL:O	1:E:264:GLY:HA3	2.01	0.61
1:G:53:LEU:HB2	1:G:254:VAL:HB	1.83	0.61
4:a:135:ASP:OD1	4:a:136:LEU:N	2.34	0.61
1:G:10:THR:HG22	1:G:10:THR:O	2.00	0.61
1:F:10:THR:HG22	1:F:10:THR:O	2.00	0.61
4:a:201:SER:HB2	4:a:271:LEU:CD2	2.30	0.61
1:G:27:TRP:HA	1:G:274:PRO:HA	1.83	0.60
4:A:311:LYS:HG2	4:A:365:ALA:CB	2.31	0.60
4:a:239:ILE:HA	5:a:809:BCL:H3C	1.83	0.60
1:E:27:TRP:HA	1:E:274:PRO:HA	1.83	0.60
1:F:36:VAL:O	1:F:264:GLY:HA3	2.01	0.60
4:A:236:PHE:O	4:A:237:GLN:NE2	2.34	0.60
4:a:311:LYS:HG2	4:a:365:ALA:CB	2.32	0.60
1:F:27:TRP:HA	1:F:274:PRO:HA	1.83	0.60
1:E:259:ARG:HG3	1:E:266:GLU:HB3	1.84	0.60
1:E:254:VAL:HG23	5:E:376:BCL:H18	1.84	0.60
1:G:331:ARG:CG	1:G:331:ARG:HH11	2.15	0.59
1:E:106:VAL:O	1:G:133:ASN:ND2	2.35	0.59
4:a:236:PHE:O	4:a:237:GLN:NE2	2.34	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:51:VAL:HG12	1:F:53:LEU:HD13	1.83	0.59
1:G:34:ALA:HB3	1:G:267:VAL:HB	1.85	0.59
4:a:201:SER:CB	4:a:271:LEU:HD23	2.32	0.59
1:G:83:THR:HG22	1:G:103:MET:CG	2.32	0.59
4:A:570:MET:HE2	4:A:597:PHE:CE2	2.37	0.59
1:E:195:GLU:O	1:F:191:THR:HG21	2.03	0.59
4:A:239:ILE:HA	5:A:809:BCL:H3C	1.84	0.59
1:F:34:ALA:HB3	1:F:267:VAL:HB	1.85	0.59
4:A:311:LYS:HG2	4:A:365:ALA:HB1	1.84	0.59
1:F:254:VAL:HG11	5:F:375:BCL:H72	1.85	0.58
4:A:529:GLY:HA2	4:a:529:GLY:HA2	1.85	0.58
4:a:311:LYS:HG2	4:a:365:ALA:HB1	1.84	0.58
4:a:570:MET:HE2	4:a:597:PHE:CE2	2.37	0.58
4:a:700:THR:HG22	4:a:701:ASN:H	1.68	0.58
4:A:646:THR:O	4:A:650:ARG:HB2	2.03	0.58
1:E:34:ALA:HB3	1:E:267:VAL:HB	1.85	0.58
1:G:38:ALA:H	1:G:264:GLY:HA2	1.69	0.58
2:D:87:ALA:HA	2:D:90:MET:HE2	1.85	0.58
5:A:811:BCL:H202	5:A:813:BCL:H2C	1.86	0.58
4:a:483:TRP:O	4:a:487:HIS:ND1	2.37	0.58
1:F:38:ALA:H	1:F:264:GLY:HA2	1.69	0.57
4:A:700:THR:HG22	4:A:701:ASN:H	1.68	0.57
1:F:52:LYS:C	1:F:53:LEU:HD12	2.29	0.57
3:B:192:PRO:HA	4:A:519:LYS:HD2	1.85	0.57
2:D:103:ILE:HG13	2:D:104:ASP:N	2.20	0.57
2:D:63:THR:O	2:D:66:MET:N	2.35	0.57
4:A:483:TRP:O	4:A:487:HIS:ND1	2.37	0.57
1:E:350:GLU:OE1	1:F:305:ASN:ND2	2.37	0.57
1:E:38:ALA:H	1:E:264:GLY:HA2	1.69	0.57
1:G:331:ARG:HH12	4:A:274:ASP:HB3	1.70	0.57
2:D:28:VAL:HG13	2:D:29:HIS:CG	2.40	0.57
1:E:53:LEU:HB2	1:E:254:VAL:HB	1.86	0.56
4:a:616:ILE:HG12	5:a:815:BCL:HMC3	1.86	0.56
2:D:78:TYR:HA	2:D:110:VAL:O	2.05	0.56
4:A:510:SER:HB3	4:A:513:TRP:CD1	2.36	0.56
4:A:583:ASP:OD1	4:A:704:LEU:CD2	2.53	0.56
4:A:616:ILE:HG12	5:A:815:BCL:HMC3	1.86	0.56
4:a:570:MET:CE	4:a:597:PHE:HE2	2.18	0.56
2:D:81:THR:HG22	2:D:111:LYS:HE2	1.86	0.56
4:A:206:LEU:HB2	5:A:810:BCL:HHD	1.88	0.56
4:a:206:LEU:HB2	5:a:810:BCL:HHD	1.88	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:39:ALA:HA	2:D:46:LYS:HZ2	1.70	0.56
4:A:67:THR:HG21	4:A:276:LYS:HE2	1.87	0.56
4:a:67:THR:HG21	4:a:276:LYS:HE2	1.87	0.56
4:A:570:MET:CE	4:A:597:PHE:HE2	2.18	0.56
4:a:206:LEU:HB2	5:a:810:BCL:CHD	2.36	0.56
5:a:811:BCL:H202	5:a:813:BCL:H2C	1.86	0.56
4:A:206:LEU:HB2	5:A:810:BCL:CHD	2.36	0.55
1:F:316:VAL:HG22	1:F:337:GLU:HG2	1.88	0.55
4:a:328:VAL:HG13	4:a:332:PHE:CD2	2.42	0.55
4:a:583:ASP:HA	4:a:588:LEU:HG	1.89	0.55
1:E:316:VAL:HG22	1:E:337:GLU:HG2	1.88	0.55
4:A:328:VAL:HG13	4:A:332:PHE:CD2	2.42	0.55
4:a:401:ASN:OD1	4:a:402:GLY:N	2.38	0.55
2:D:70:LEU:HA	2:D:75:ILE:HB	1.89	0.55
4:a:510:SER:HB3	4:a:513:TRP:CD1	2.36	0.55
1:G:331:ARG:HH11	1:G:331:ARG:HG2	1.71	0.55
2:D:112:TRP:HE3	3:B:138:ASN:HB2	1.71	0.55
4:A:326:ASN:N	4:A:353:ALA:O	2.38	0.54
2:D:65:GLU:O	2:D:68:ARG:HG3	2.08	0.54
1:G:254:VAL:CG2	5:G:376:BCL:H18	2.37	0.54
1:G:316:VAL:HG22	1:G:337:GLU:HG2	1.88	0.54
4:A:116:THR:HG21	5:A:811:BCL:HMA2	1.89	0.54
4:a:326:ASN:N	4:a:353:ALA:O	2.38	0.54
1:F:321:TYR:HA	1:F:365:ALA:O	2.08	0.54
1:G:312:GLN:NE2	1:G:339:ARG:NH1	2.54	0.54
2:D:28:VAL:HG13	2:D:29:HIS:CD2	2.42	0.54
1:F:262:SER:O	1:F:263:GLN:CG	2.56	0.54
4:A:581:ALA:O	4:A:583:ASP:O	2.26	0.54
1:G:262:SER:O	1:G:263:GLN:CG	2.56	0.54
1:G:321:TYR:HA	1:G:365:ALA:O	2.08	0.54
2:D:49:ILE:HB	2:D:62:PRO:HD3	1.89	0.54
4:a:116:THR:HG21	5:a:811:BCL:HMA2	1.89	0.54
1:G:286:ARG:HB3	1:G:363:LEU:HD23	1.90	0.54
4:A:649:ILE:CD1	4:A:665:PHE:HZ	2.21	0.53
4:A:101:PHE:O	4:A:103:TRP:N	2.41	0.53
4:A:209:HIS:ND1	5:A:810:BCL:OBD	2.39	0.53
1:F:359:GLN:NE2	1:G:43:PRO:O	2.41	0.53
3:B:172:CYS:SG	6:B:301:SF4:S3	3.04	0.53
1:E:262:SER:O	1:E:263:GLN:CG	2.56	0.53
4:A:416:ARG:HG2	4:A:416:ARG:O	2.08	0.53
4:a:209:HIS:ND1	5:a:810:BCL:OBD	2.39	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:35:LYS:HE3	4:A:408:LYS:HB3	1.91	0.53
1:E:305:ASN:ND2	1:G:350:GLU:OE1	2.40	0.53
1:E:321:TYR:HA	1:E:365:ALA:O	2.08	0.53
4:a:80:LEU:HD11	5:a:807:BCL:HMC3	1.91	0.53
2:D:63:THR:O	2:D:65:GLU:N	2.43	0.52
4:A:80:LEU:HD11	5:A:807:BCL:HMC3	1.91	0.52
1:G:342:GLY:N	4:A:200:LYS:HE3	2.24	0.52
3:B:182:CYS:N	6:B:302:SF4:S2	2.83	0.52
1:F:286:ARG:HB3	1:F:363:LEU:HD23	1.90	0.52
4:a:399:GLN:HG2	4:a:404:TYR:CG	2.45	0.52
4:a:420:VAL:HG12	4:a:424:ILE:HD12	1.92	0.52
1:E:286:ARG:HB3	1:E:363:LEU:HD23	1.90	0.52
4:A:476:ALA:HB3	11:A:819:LMG:O3	2.09	0.52
4:a:476:ALA:HB3	11:a:819:LMG:O3	2.09	0.52
4:A:206:LEU:HB2	5:A:810:BCL:C1D	2.31	0.52
4:A:420:VAL:HG12	4:A:424:ILE:HD12	1.92	0.52
1:E:331:ARG:HD2	1:E:341:SER:O	2.09	0.52
1:F:53:LEU:HB2	1:F:254:VAL:HB	1.92	0.51
4:A:399:GLN:HG2	4:A:404:TYR:CG	2.45	0.51
4:a:349:PHE:HB3	4:a:350:PRO:HD3	1.92	0.51
1:E:93:LYS:HG2	1:E:93:LYS:O	2.11	0.51
1:F:331:ARG:HD2	1:F:341:SER:O	2.09	0.51
5:F:371:BCL:HAA1	5:F:371:BCL:CBD	2.35	0.51
4:A:95:ILE:HG22	4:A:95:ILE:O	2.10	0.51
1:E:328:PRO:HD2	1:E:360:PHE:CD1	2.45	0.51
1:F:328:PRO:HD2	1:F:360:PHE:CD1	2.45	0.51
1:G:328:PRO:HD2	1:G:360:PHE:CD1	2.45	0.51
4:A:202:TYR:H	4:A:271:LEU:HD22	1.75	0.51
1:E:32:ALA:O	1:E:33:ARG:HD3	2.11	0.51
4:A:67:THR:HG21	4:A:276:LYS:HB2	1.93	0.51
4:A:380:GLY:O	4:A:384:VAL:HG23	2.11	0.51
1:G:32:ALA:O	1:G:33:ARG:HD3	2.11	0.51
1:G:310:ASP:HA	4:A:200:LYS:HE2	1.93	0.50
4:A:281:VAL:O	4:A:284:THR:HG22	2.11	0.50
4:A:349:PHE:HB3	4:A:350:PRO:HD3	1.92	0.50
1:G:329:GLN:HE22	4:A:398:ILE:HG22	1.77	0.50
1:F:32:ALA:O	1:F:33:ARG:HD3	2.11	0.50
4:a:67:THR:HG21	4:a:276:LYS:HB2	1.93	0.50
4:A:332:PHE:CE2	4:A:342:GLY:HA3	2.47	0.50
4:a:380:GLY:O	4:a:384:VAL:HG23	2.11	0.50
1:E:254:VAL:CG2	5:E:376:BCL:H18	2.41	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:131:ARG:O	1:G:131:ARG:HG2	2.12	0.50
1:E:82:LEU:HD12	1:E:104:VAL:HB	1.94	0.50
2:D:67:ILE:O	2:D:71:ILE:HG23	2.12	0.50
4:A:403:MET:SD	4:A:406:GLN:OE1	2.70	0.50
5:F:372:BCL:H171	5:F:371:BCL:HBC3	1.94	0.50
3:B:165:VAL:CG1	3:B:170:PHE:HB3	2.42	0.50
5:E:371:BCL:HBC3	5:E:372:BCL:H171	1.94	0.49
3:B:185:CYS:O	3:B:196:LEU:CD2	2.60	0.49
1:G:342:GLY:H	4:A:200:LYS:HE3	1.76	0.49
4:a:651:TRP:O	4:a:654:LYS:HB3	2.13	0.49
1:E:191:THR:HG21	1:G:195:GLU:O	2.12	0.49
4:A:651:TRP:O	4:A:654:LYS:HB3	2.13	0.49
1:G:82:LEU:HD12	1:G:104:VAL:HB	1.94	0.49
4:A:181:ARG:O	4:A:198:SER:OG	2.17	0.49
3:B:165:VAL:HG12	3:B:170:PHE:HB3	1.94	0.49
3:B:200:LEU:HB2	3:B:204:GLU:HB2	1.93	0.49
2:D:94:ILE:HD12	2:D:100:ARG:CG	2.41	0.49
4:a:332:PHE:CE2	4:a:342:GLY:HA3	2.47	0.49
4:a:465:LEU:HB3	4:a:480:ILE:HD13	1.95	0.49
1:F:82:LEU:HD12	1:F:104:VAL:HB	1.94	0.49
4:A:370:ASN:HB3	4:A:693:LEU:HD22	1.94	0.49
4:a:436:MET:HA	4:a:439:VAL:HG22	1.94	0.49
1:G:254:VAL:HG23	5:G:376:BCL:C18	2.43	0.48
4:a:159:PHE:CZ	4:a:163:ARG:NE	2.70	0.48
1:F:206:ASN:HB2	1:F:222:ARG:HG3	1.96	0.48
1:G:331:ARG:CG	1:G:331:ARG:NH1	2.73	0.48
2:D:62:PRO:O	2:D:66:MET:HB2	2.14	0.48
4:a:102:PRO:HG2	4:a:144:LEU:CD1	2.43	0.48
5:G:378[B]:BCL:H2C	5:G:378[B]:BCL:HBC3	1.66	0.48
4:A:71:ASP:OD1	4:A:74:GLN:HG3	2.14	0.48
5:A:812:BCL:H141	5:A:812:BCL:H161	1.68	0.48
4:A:101:PHE:O	4:A:103:TRP:CD1	2.67	0.48
4:A:88:PHE:HE2	4:A:158:ILE:HD12	1.79	0.48
4:A:465:LEU:HB3	4:A:480:ILE:HD13	1.95	0.48
1:G:206:ASN:HB2	1:G:222:ARG:HG3	1.96	0.48
5:G:372:BCL:H171	5:G:371:BCL:HBC3	1.94	0.48
4:A:278:ALA:O	4:A:282:HIS:ND1	2.46	0.48
4:a:278:ALA:O	4:a:282:HIS:ND1	2.46	0.48
4:a:370:ASN:HB3	4:a:693:LEU:HD22	1.94	0.48
5:a:812:BCL:H143	5:a:812:BCL:H111	1.59	0.48
1:E:284:LEU:HG	1:E:365:ALA:HB2	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:284:LEU:HG	1:F:365:ALA:HB2	1.96	0.48
4:A:436:MET:HA	4:A:439:VAL:HG22	1.94	0.48
1:G:284:LEU:HG	1:G:365:ALA:HB2	1.96	0.47
2:D:79:ILE:CB	2:D:111:LYS:HG3	2.35	0.47
4:A:561:TYR:CE2	11:A:819:LMG:HC61	2.49	0.47
4:A:674:LYS:HE2	4:a:633:GLU:HB2	1.95	0.47
4:a:88:PHE:HD1	4:a:151:PHE:HD1	1.62	0.47
4:a:561:TYR:CE2	11:a:819:LMG:HC61	2.49	0.47
1:E:206:ASN:HB2	1:E:222:ARG:HG3	1.96	0.47
2:D:79:ILE:HG13	2:D:101:TYR:HD2	1.79	0.47
1:G:321:TYR:CD1	1:G:366:GLN:HB3	2.50	0.47
1:G:329:GLN:HE22	4:A:398:ILE:CG2	2.27	0.47
4:A:429:MET:HE1	13:a:803:G2O:CED	2.44	0.47
4:A:649:ILE:CD1	4:A:665:PHE:CZ	2.95	0.47
4:a:71:ASP:OD1	4:a:74:GLN:HG3	2.14	0.47
4:a:88:PHE:HE2	4:a:158:ILE:HD12	1.79	0.47
1:F:170:ILE:HD13	1:F:208:LEU:HD12	1.96	0.47
1:G:262:SER:C	1:G:263:GLN:HG3	2.39	0.47
2:D:81:THR:HG23	2:D:112:TRP:O	2.14	0.47
5:A:807:BCL:O1D	5:A:807:BCL:H2A	2.15	0.47
1:E:259:ARG:HD3	1:E:266:GLU:HG2	1.97	0.47
1:E:262:SER:C	1:E:263:GLN:HG3	2.39	0.47
1:F:123:PHE:HB3	1:F:140:MET:HE3	1.97	0.47
1:F:262:SER:C	1:F:263:GLN:HG3	2.39	0.47
1:F:321:TYR:CD1	1:F:366:GLN:HB3	2.50	0.47
5:F:372:BCL:H203	5:F:372:BCL:H162	1.77	0.47
5:F:371:BCL:H112	5:F:371:BCL:H152	1.59	0.47
4:A:331:ASN:O	4:A:341:SER:OG	2.27	0.47
4:A:408:LYS:N	4:A:408:LYS:HD3	2.28	0.47
4:A:415:GLY:O	4:A:421:GLN:NE2	2.41	0.47
1:E:323:ILE:HD12	1:E:338:TYR:CG	2.50	0.47
4:A:101:PHE:HA	4:A:102:PRO:HD2	1.70	0.47
4:A:320:LEU:O	4:A:323:GLU:HG3	2.15	0.47
1:E:17:GLU:OE2	1:E:314:LYS:HE2	2.15	0.47
1:E:321:TYR:CD1	1:E:366:GLN:HB3	2.50	0.47
1:F:17:GLU:OE2	1:F:314:LYS:HE2	2.15	0.47
1:F:167:VAL:HA	1:F:170:ILE:HG12	1.96	0.47
1:F:323:ILE:HD12	1:F:338:TYR:CG	2.50	0.47
2:D:94:ILE:HB	2:D:100:ARG:HD3	1.95	0.47
1:E:166:THR:O	1:E:170:ILE:HG23	2.16	0.46
2:D:42:ASP:OD2	2:D:42:ASP:N	2.47	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:a:807:BCL:H2A	5:a:807:BCL:O1D	2.15	0.46
1:E:83:THR:OG1	1:E:103:MET:HG2	2.15	0.46
1:E:170:ILE:HD13	1:E:208:LEU:HD12	1.96	0.46
5:F:373:BCL:H61	5:F:373:BCL:H41	1.81	0.46
1:G:167:VAL:O	1:G:170:ILE:HG22	2.15	0.46
4:A:599:TYR:CE1	4:A:603:GLU:OE1	2.69	0.46
4:a:67:THR:CG2	4:a:276:LYS:HE2	2.46	0.46
4:a:443:ILE:CG2	4:a:570:MET:HG3	2.44	0.46
1:E:167:VAL:HA	1:E:170:ILE:HG12	1.96	0.46
1:F:62:LYS:HB3	1:F:64:PHE:HD2	1.80	0.46
1:F:166:THR:O	1:F:170:ILE:HG23	2.16	0.46
1:G:17:GLU:OE2	1:G:314:LYS:HE2	2.15	0.46
4:A:67:THR:CG2	4:A:276:LYS:HE2	2.46	0.46
4:A:213:MET:HE1	5:A:810:BCL:O1D	2.15	0.46
4:A:403:MET:SD	4:A:406:GLN:CD	2.99	0.46
1:F:83:THR:OG1	1:F:103:MET:HG2	2.15	0.46
1:G:123:PHE:HB3	1:G:140:MET:HE3	1.97	0.46
4:a:318:PHE:HE1	4:a:359:GLY:HA3	1.81	0.46
4:a:320:LEU:O	4:a:323:GLU:HG3	2.15	0.46
1:G:331:ARG:CZ	4:A:271:LEU:O	2.61	0.46
4:A:443:ILE:CG2	4:A:570:MET:HG3	2.44	0.46
5:a:809:BCL:H2C	5:a:809:BCL:HBC3	1.60	0.46
1:E:329:GLN:HG2	1:E:330:TYR:O	2.15	0.46
5:E:371:BCL:H192	5:E:371:BCL:H161	1.71	0.46
1:G:36:VAL:O	1:G:264:GLY:CA	2.63	0.46
5:E:375:BCL:H2	5:E:375:BCL:H61	1.67	0.46
1:F:166:THR:O	1:F:170:ILE:HG12	2.16	0.46
5:a:812:BCL:H141	5:a:812:BCL:H161	1.68	0.46
1:E:62:LYS:HB3	1:E:64:PHE:HD2	1.81	0.46
4:A:633:GLU:HB2	4:a:674:LYS:HE2	1.97	0.46
4:a:602:THR:O	4:a:603:GLU:CB	2.60	0.46
5:E:372:BCL:H203	5:E:372:BCL:H162	1.77	0.46
1:F:36:VAL:O	1:F:264:GLY:CA	2.63	0.46
1:E:123:PHE:HB3	1:E:140:MET:HE3	1.97	0.46
5:E:373:BCL:H61	5:E:373:BCL:H41	1.81	0.46
5:A:806:BCL:H93	5:A:806:BCL:H61	1.80	0.46
4:a:213:MET:HE1	5:a:810:BCL:O1D	2.15	0.46
4:a:321:TYR:HB2	5:a:815:BCL:O1D	2.16	0.46
1:E:36:VAL:O	1:E:264:GLY:CA	2.63	0.45
1:E:166:THR:O	1:E:170:ILE:HG12	2.16	0.45
1:G:323:ILE:HD12	1:G:338:TYR:CG	2.50	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:A:321:TYR:HB2	5:A:815:BCL:O1D	2.16	0.45
4:a:513:TRP:CZ3	4:a:518:LEU:HD13	2.51	0.45
1:G:62:LYS:HB3	1:G:64:PHE:HD2	1.81	0.45
1:G:167:VAL:HA	1:G:170:ILE:HG22	1.98	0.45
4:A:88:PHE:HD1	4:A:151:PHE:HD1	1.62	0.45
4:a:599:TYR:CE1	4:a:603:GLU:OE1	2.69	0.45
1:E:50:ASP:OD1	1:E:51:VAL:N	2.49	0.45
1:E:109:PHE:H	1:G:133:ASN:HD21	1.64	0.45
1:F:50:ASP:OD1	1:F:51:VAL:N	2.49	0.45
1:G:50:ASP:OD1	1:G:51:VAL:N	2.49	0.45
3:B:167:GLU:OE2	3:B:167:GLU:N	2.49	0.45
3:B:185:CYS:O	3:B:196:LEU:HD22	2.17	0.45
5:F:372:BCL:H141	5:F:372:BCL:H161	1.72	0.45
2:D:65:GLU:HG2	2:D:68:ARG:HE	1.81	0.45
4:A:318:PHE:HE1	4:A:359:GLY:HA3	1.81	0.45
4:a:647:ARG:HG2	4:a:650:ARG:NH2	2.31	0.45
3:B:223:LEU:HD23	3:B:223:LEU:H	1.80	0.45
4:a:101:PHE:HA	4:a:102:PRO:HD2	1.59	0.45
4:a:202:TYR:CZ	4:a:274:ASP:OD2	2.64	0.45
1:E:44:LEU:HD23	1:G:359:GLN:HE22	1.82	0.45
5:G:372:BCL:H141	5:G:372:BCL:H161	1.72	0.45
4:A:513:TRP:CZ3	4:A:518:LEU:HD13	2.51	0.45
4:a:415:GLY:O	4:a:421:GLN:NE2	2.41	0.45
2:D:49:ILE:HB	2:D:62:PRO:CD	2.47	0.45
4:A:539:SER:OG	4:A:542:ASP:OD2	2.30	0.45
4:A:648:THR:O	4:A:652:LEU:N	2.45	0.45
5:A:809:BCL:HBC3	5:A:809:BCL:H2C	1.60	0.45
1:E:195:GLU:O	1:F:191:THR:CG2	2.65	0.45
1:F:185:PHE:CZ	1:F:200:ILE:HG13	2.52	0.45
1:G:185:PHE:CZ	1:G:200:ILE:HG13	2.52	0.45
4:a:73:GLU:OE1	4:a:167:HIS:HA	2.17	0.45
4:a:399:GLN:HG2	4:a:404:TYR:CB	2.47	0.45
5:a:806:BCL:H93	5:a:806:BCL:H61	1.80	0.45
1:E:91:GLU:OE1	1:E:95:ARG:NH2	2.50	0.45
5:E:378[B]:BCL:H2C	5:E:378[B]:BCL:HBC3	1.66	0.45
1:F:91:GLU:OE1	1:F:95:ARG:NH2	2.50	0.45
1:G:91:GLU:OE1	1:G:95:ARG:NH2	2.50	0.45
1:E:108:ASP:OD1	2:D:60:LEU:HD22	2.16	0.44
5:G:374:BCL:H162	5:G:374:BCL:H192	1.76	0.44
5:G:375:BCL:H2	5:G:375:BCL:H61	1.67	0.44
4:a:595:GLY:O	4:a:599:TYR:HB2	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:185:PHE:CZ	1:E:200:ILE:HG13	2.52	0.44
1:F:330:TYR:HB3	1:F:340:TRP:CE2	2.52	0.44
5:G:376:BCL:H41	5:G:376:BCL:H61	1.57	0.44
4:A:102:PRO:HG2	4:A:144:LEU:CD1	2.47	0.44
4:A:399:GLN:HG2	4:A:404:TYR:CB	2.47	0.44
4:A:490:THR:HG21	5:A:814:BCL:HMA1	2.00	0.44
4:a:86:LEU:HD12	4:a:89:ILE:HD11	1.99	0.44
4:a:539:SER:OG	4:a:542:ASP:OD2	2.30	0.44
4:a:648:THR:O	4:a:652:LEU:N	2.45	0.44
1:E:254:VAL:HG23	5:E:376:BCL:C18	2.45	0.44
1:E:330:TYR:HB3	1:E:340:TRP:CE2	2.52	0.44
1:F:40:PRO:HG3	1:F:260:SER:HB2	2.00	0.44
5:F:376:BCL:H61	5:F:376:BCL:H41	1.57	0.44
13:A:803:G2O:CED	4:a:429:MET:HE1	2.48	0.44
5:F:377:BCL:H162	5:F:377:BCL:H122	1.74	0.44
4:A:198:SER:HA	5:A:808:BCL:HBB1	1.99	0.44
1:G:330:TYR:HB3	1:G:340:TRP:CE2	2.52	0.44
4:A:648:THR:C	4:A:650:ARG:N	2.75	0.44
4:A:649:ILE:HD12	4:A:665:PHE:HE2	1.79	0.44
4:a:159:PHE:CE1	4:a:163:ARG:NE	2.85	0.44
4:A:86:LEU:HD12	4:A:89:ILE:HD11	1.99	0.44
4:a:648:THR:C	4:a:650:ARG:N	2.74	0.44
1:F:313:LEU:HD23	1:F:340:TRP:CE3	2.53	0.44
4:A:73:GLU:OE1	4:A:167:HIS:HA	2.17	0.44
4:A:105:ALA:O	4:A:106:PRO:C	2.61	0.44
5:a:806:BCL:H122	5:a:806:BCL:H161	1.69	0.44
5:E:376:BCL:H2C	5:E:376:BCL:HBC3	1.66	0.43
3:B:170:PHE:CZ	4:A:530:PRO:HD2	2.52	0.43
4:A:103:TRP:HD1	4:A:104:GLY:N	2.16	0.43
4:A:314:ASP:O	4:A:360:GLU:HB3	2.18	0.43
4:A:391:GLY:O	4:A:395:LEU:HB2	2.18	0.43
1:G:325:TYR:HB2	3:B:210:TYR:CE2	2.53	0.43
4:A:286:ILE:HD11	5:A:810:BCL:HBA1	1.99	0.43
4:a:198:SER:HA	5:a:808:BCL:HBB1	1.99	0.43
4:a:314:ASP:O	4:a:360:GLU:HB3	2.18	0.43
4:a:391:GLY:O	4:a:395:LEU:HB2	2.18	0.43
4:a:583:ASP:O	4:a:587:HIS:N	2.50	0.43
1:E:313:LEU:HD23	1:E:340:TRP:CE3	2.53	0.43
2:D:22:PRO:HD2	2:D:23:TRP:CZ3	2.53	0.43
2:D:76:GLU:OE2	2:D:76:GLU:N	2.51	0.43
4:a:331:ASN:O	4:a:341:SER:OG	2.27	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:G:377:BCL:H122	5:G:377:BCL:H162	1.74	0.43
2:D:38:SER:O	2:D:46:LYS:NZ	2.51	0.43
2:D:94:ILE:HD12	2:D:100:ARG:CD	2.48	0.43
4:A:583:ASP:HB3	4:A:703:PRO:O	2.18	0.43
4:A:595:GLY:O	4:A:599:TYR:HB2	2.18	0.43
5:A:809:BCL:HAA2	5:A:809:BCL:HBD	1.99	0.43
4:a:286:ILE:HD11	5:a:810:BCL:HBA1	1.99	0.43
1:E:40:PRO:HG3	1:E:260:SER:HB3	2.00	0.43
1:F:95:ARG:HG3	1:F:120:VAL:HG22	2.01	0.43
4:A:103:TRP:CD1	4:A:104:GLY:N	2.87	0.43
5:E:371:BCL:H112	5:E:371:BCL:H152	1.59	0.43
5:F:373:BCL:H111	5:F:373:BCL:H142	1.78	0.43
5:F:375:BCL:H61	5:F:375:BCL:H2	1.67	0.43
1:G:95:ARG:HG3	1:G:120:VAL:HG22	2.01	0.43
1:G:312:GLN:HE21	1:G:339:ARG:HD3	1.84	0.43
2:D:114:MET:HE1	3:B:197:ILE:HD13	1.99	0.43
4:a:490:THR:HG21	5:a:814:BCL:HMA1	2.00	0.43
5:E:373:BCL:H142	5:E:373:BCL:H111	1.78	0.43
1:G:40:PRO:HG3	1:G:260:SER:HB3	2.00	0.43
5:G:372:BCL:H203	5:G:372:BCL:H162	1.77	0.43
3:B:135:ILE:HD12	3:B:174:ILE:HG13	2.01	0.43
1:F:333:GLN:HE21	1:F:338:TYR:HE1	1.66	0.43
1:G:313:LEU:HD23	1:G:340:TRP:CE3	2.53	0.43
2:D:68:ARG:HD2	2:D:69:ARG:N	2.33	0.43
4:A:484:VAL:HG21	4:A:568:ALA:HA	2.01	0.43
5:a:809:BCL:HAA2	5:a:809:BCL:HBD	1.99	0.43
1:E:199:ARG:HD2	1:F:190:PHE:CZ	2.53	0.42
1:F:49:CYS:HB3	1:F:258:PHE:CE2	2.54	0.42
5:G:371:BCL:H192	5:G:371:BCL:H161	1.71	0.42
3:B:191:CYS:O	4:A:519:LYS:HD2	2.18	0.42
5:A:812:BCL:H143	5:A:812:BCL:H111	1.59	0.42
5:a:811:BCL:H62	5:a:811:BCL:H41	1.77	0.42
1:E:43:PRO:O	1:G:359:GLN:NE2	2.52	0.42
1:E:49:CYS:HB3	1:E:258:PHE:CE2	2.55	0.42
1:E:148:ILE:O	1:E:222:ARG:HA	2.19	0.42
5:E:377:BCL:H122	5:E:377:BCL:H162	1.74	0.42
1:G:148:ILE:O	1:G:222:ARG:HA	2.19	0.42
2:D:42:ASP:C	2:D:44:ASP:H	2.27	0.42
3:B:220:LEU:HD12	3:B:220:LEU:C	2.44	0.42
4:A:270:HIS:CE1	4:A:281:VAL:HG21	2.54	0.42
4:a:201:SER:CB	4:a:271:LEU:CD2	2.95	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:a:657:LEU:H	4:a:657:LEU:HD23	1.84	0.42
1:E:195:GLU:C	1:E:197:GLY:N	2.77	0.42
1:F:148:ILE:O	1:F:222:ARG:HA	2.19	0.42
2:D:31:ALA:HB3	2:D:76:GLU:OE1	2.19	0.42
4:A:202:TYR:HB2	4:A:271:LEU:CD2	2.49	0.42
4:A:657:LEU:H	4:A:657:LEU:HD23	1.84	0.42
2:D:24:SER:HB2	3:B:214:GLU:HA	2.01	0.42
5:E:374:BCL:H162	5:E:374:BCL:H192	1.76	0.42
4:A:533:GLY:O	4:a:670:ILE:HG13	2.19	0.42
4:a:270:HIS:CE1	4:a:281:VAL:HG21	2.54	0.42
4:a:323:GLU:HB2	4:a:354:ILE:HG23	2.02	0.42
5:G:377:BCL:H161	5:G:377:BCL:H192	1.75	0.42
1:G:49:CYS:HB3	1:G:258:PHE:CE2	2.55	0.42
2:D:39:ALA:HA	2:D:46:LYS:NZ	2.35	0.42
4:a:105:ALA:O	4:a:106:PRO:C	2.61	0.42
4:a:395:LEU:O	4:a:399:GLN:HG3	2.20	0.42
4:a:484:VAL:HG21	4:a:568:ALA:HA	2.01	0.42
5:E:373:BCL:H141	5:E:373:BCL:H162	1.84	0.42
5:F:374:BCL:H192	5:F:374:BCL:H162	1.76	0.42
4:A:101:PHE:HB3	4:A:103:TRP:CD1	2.55	0.42
4:A:220:TRP:O	4:A:224:MET:HG3	2.19	0.42
5:A:814:BCL:HBC3	5:A:814:BCL:H2C	1.79	0.42
1:E:95:ARG:HG3	1:E:120:VAL:HG22	2.01	0.42
1:G:333:GLN:HE21	1:G:338:TYR:HE1	1.66	0.42
2:D:23:TRP:CH2	4:A:418:GLN:HB2	2.55	0.42
2:D:114:MET:HA	3:B:137:GLU:HG2	2.00	0.42
5:A:812:BCL:H112	5:A:812:BCL:H72	1.73	0.42
4:a:522:SER:C	4:a:524:SER:H	2.27	0.42
1:G:256:ALA:HB3	5:G:375:BCL:H92	2.02	0.42
1:E:182:ASP:OD2	1:G:143:ARG:NH2	2.46	0.41
1:E:259:ARG:HG3	1:E:259:ARG:O	2.20	0.41
1:E:323:ILE:HD12	1:E:338:TYR:CD2	2.55	0.41
5:F:374:BCL:H92	5:F:374:BCL:H62	1.82	0.41
3:B:192:PRO:HA	4:A:519:LYS:HB2	2.02	0.41
3:B:206:LEU:HD23	3:B:206:LEU:HA	1.92	0.41
4:A:522:SER:C	4:A:524:SER:H	2.27	0.41
4:a:102:PRO:HG2	4:a:144:LEU:HD12	2.02	0.41
4:A:395:LEU:O	4:A:399:GLN:HG3	2.20	0.41
5:E:376:BCL:HMB2	5:E:372:BCL:H18	2.02	0.41
1:G:323:ILE:HD12	1:G:338:TYR:CD2	2.55	0.41
5:G:372:BCL:H18	5:G:376:BCL:HMB2	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:A:298:PHE:HB2	4:A:373:VAL:HG21	2.02	0.41
5:A:808:BCL:H93	5:A:808:BCL:H62	1.85	0.41
1:E:128:ASP:OD1	1:E:131:ARG:NH2	2.53	0.41
1:E:143:ARG:NH2	1:F:182:ASP:OD2	2.46	0.41
1:E:359:GLN:NE2	1:F:43:PRO:O	2.54	0.41
1:G:205:VAL:HG12	1:G:207:GLY:H	1.86	0.41
1:G:280:ALA:HB2	1:G:286:ARG:HG2	2.02	0.41
3:B:157:ILE:HG12	3:B:174:ILE:HG12	2.01	0.41
4:A:201:SER:HB2	4:A:271:LEU:HD22	2.02	0.41
5:a:807:BCL:H3A	5:a:807:BCL:HBA2	1.90	0.41
5:E:371:BCL:H112	5:E:371:BCL:H72	1.85	0.41
1:F:51:VAL:CG1	1:F:53:LEU:HD11	2.47	0.41
1:F:205:VAL:HG12	1:F:207:GLY:H	1.86	0.41
1:F:256:ALA:HB3	5:F:375:BCL:H92	2.02	0.41
1:F:323:ILE:HD12	1:F:338:TYR:CD2	2.55	0.41
5:G:371:BCL:HAA1	5:G:371:BCL:CB D	2.35	0.41
5:A:806:BCL:H161	5:A:806:BCL:H122	1.69	0.41
5:A:806:BCL:CAD	5:A:812:BCL:H192	2.50	0.41
1:E:280:ALA:HB2	1:E:286:ARG:HG2	2.02	0.41
1:F:280:ALA:HB2	1:F:286:ARG:HG2	2.02	0.41
4:A:202:TYR:HB2	4:A:271:LEU:HD23	2.03	0.41
4:A:323:GLU:HB2	4:A:354:ILE:HG23	2.02	0.41
5:a:806:BCL:CAD	5:a:812:BCL:H192	2.50	0.41
5:E:377:BCL:HED2	5:E:377:BCL:H43	2.02	0.41
5:G:372:BCL:H92	5:G:372:BCL:H61	1.83	0.41
5:G:376:BCL:HBC3	5:G:376:BCL:H2C	1.66	0.41
5:A:811:BCL:H92	5:A:811:BCL:H61	1.88	0.41
4:a:298:PHE:HB2	4:a:373:VAL:HG21	2.02	0.41
1:E:256:ALA:HB3	5:E:375:BCL:H92	2.02	0.41
1:F:350:GLU:OE1	1:G:305:ASN:ND2	2.52	0.41
4:A:407:ILE:HG21	4:A:412:ILE:HG13	2.03	0.41
4:a:220:TRP:O	4:a:224:MET:HG3	2.19	0.41
4:a:407:ILE:HG21	4:a:412:ILE:HG13	2.03	0.41
5:a:814:BCL:HBC3	5:a:814:BCL:H2C	1.79	0.41
1:E:193:LEU:HD11	1:E:200:ILE:HG12	2.03	0.41
1:F:103:MET:HB3	1:F:112:THR:HG22	2.03	0.41
1:F:193:LEU:HD11	1:F:200:ILE:HG12	2.03	0.41
1:G:195:GLU:C	1:G:197:GLY:N	2.78	0.41
5:G:377:BCL:HED2	5:G:377:BCL:H43	2.02	0.41
3:B:133:TYR:HA	3:B:198:GLU:HA	2.03	0.41
4:A:285:ALA:O	4:A:289:ILE:HG12	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:A:646:THR:HG21	4:A:664:ARG:HG2	2.03	0.41
4:a:332:PHE:CZ	4:a:344:LYS:HD3	2.56	0.41
4:a:399:GLN:HG2	4:a:404:TYR:HB2	2.03	0.41
1:F:153:PRO:HA	1:F:218:VAL:HG22	2.03	0.41
5:F:372:BCL:H18	5:F:376:BCL:HMB2	2.02	0.41
5:F:375:BCL:H193	5:F:377:BCL:H203	2.03	0.41
2:D:41:ARG:HH21	2:D:90:MET:HE1	1.85	0.41
2:D:58:ARG:HB3	2:D:60:LEU:HD21	2.02	0.41
4:A:102:PRO:HG2	4:A:144:LEU:HD12	2.03	0.41
4:A:131:THR:HG23	4:A:131:THR:O	2.21	0.41
4:a:320:LEU:O	4:a:354:ILE:HD13	2.21	0.41
4:a:648:THR:C	4:a:650:ARG:H	2.29	0.41
1:F:31:LYS:HD3	1:F:270:ASP:OD1	2.21	0.40
1:E:153:PRO:HA	1:E:218:VAL:HG22	2.02	0.40
1:E:155:ASP:OD1	1:E:156:ASN:N	2.55	0.40
5:F:377:BCL:H161	5:F:377:BCL:H192	1.75	0.40
5:A:812:BCL:HBA2	5:A:812:BCL:H3A	1.87	0.40
4:a:131:THR:HG23	4:a:131:THR:O	2.21	0.40
1:E:205:VAL:HG12	1:E:207:GLY:H	1.86	0.40
1:G:31:LYS:HD3	1:G:270:ASP:OD1	2.21	0.40
1:G:39:PRO:HA	1:G:40:PRO:HD2	2.00	0.40
1:G:58:LEU:H	1:G:65:VAL:HA	1.87	0.40
5:G:376:BCL:H61	5:G:376:BCL:H102	1.96	0.40
3:B:151:PRO:HA	3:B:152:PRO:HD3	1.95	0.40
4:A:274:ASP:O	4:A:274:ASP:CG	2.64	0.40
4:A:528:LEU:HD23	4:A:528:LEU:HA	1.96	0.40
1:E:103:MET:HB3	1:E:112:THR:HG22	2.03	0.40
1:G:28:GLY:O	1:G:272:GLU:HA	2.22	0.40
1:G:193:LEU:HD11	1:G:200:ILE:HG12	2.03	0.40
2:D:94:ILE:O	2:D:95:ILE:HB	2.21	0.40
4:A:95:ILE:O	4:A:95:ILE:CG2	2.70	0.40
4:A:433:PHE:HA	13:a:803:G2O:C4	2.52	0.40
4:A:606:THR:O	4:A:609:SER:OG	2.36	0.40
5:A:807:BCL:H3A	5:A:807:BCL:HBA2	1.90	0.40
1:E:28:GLY:O	1:E:272:GLU:HA	2.22	0.40
5:E:375:BCL:H193	5:E:377:BCL:H203	2.03	0.40
1:G:309:VAL:HG21	1:G:350:GLU:HB2	2.04	0.40
4:A:436:MET:HB3	4:A:491:ALA:HB1	2.04	0.40
4:A:700:THR:HG22	4:A:701:ASN:N	2.36	0.40
4:a:450:GLU:CD	4:a:582:TRP:HE1	2.15	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all EM entries.

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	E	360/366 (98%)	339 (94%)	21 (6%)	0	100	100
1	F	356/366 (97%)	338 (95%)	18 (5%)	0	100	100
1	G	356/366 (97%)	338 (95%)	18 (5%)	0	100	100
2	D	90/143 (63%)	71 (79%)	17 (19%)	2 (2%)	5	14
3	B	93/231 (40%)	76 (82%)	17 (18%)	0	100	100
4	A	622/731 (85%)	568 (91%)	51 (8%)	3 (0%)	24	48
4	a	622/731 (85%)	571 (92%)	46 (7%)	5 (1%)	16	37
All	All	2499/2934 (85%)	2301 (92%)	188 (8%)	10 (0%)	31	54

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	D	95	ILE
4	A	102	PRO
4	a	102	PRO
4	A	659	ARG
4	a	659	ARG
4	A	584	SER
4	a	581	ALA
4	a	582	TRP
2	D	81	THR
4	a	101	PHE

5.3.2 Protein sidechains

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all EM entries.

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	E	299/302 (99%)	299 (100%)	0	100	100
1	F	296/302 (98%)	296 (100%)	0	100	100
1	G	296/302 (98%)	294 (99%)	2 (1%)	76	90
2	D	83/128 (65%)	83 (100%)	0	100	100
3	B	80/162 (49%)	80 (100%)	0	100	100
4	A	521/599 (87%)	516 (99%)	5 (1%)	68	86
4	a	521/599 (87%)	517 (99%)	4 (1%)	73	88
All	All	2096/2394 (88%)	2085 (100%)	11 (0%)	78	92

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	G	131	ARG
1	G	331	ARG
4	A	73	GLU
4	A	103	TRP
4	A	395	LEU
4	A	408	LYS
4	A	502	ARG
4	a	73	GLU
4	a	395	LEU
4	a	408	LYS
4	a	502	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	E	7	ASN
1	E	297	HIS
1	F	37	ASN
1	F	297	HIS
1	G	297	HIS
1	G	312	GLN
1	G	329	GLN
1	G	359	GLN
2	D	48	GLN
3	B	166	GLN
4	A	150	HIS
4	A	237	GLN

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Mol	Chain	Res	Type
4	A	303	GLN
4	A	521	ASN
4	A	612	HIS
4	a	150	HIS
4	a	237	GLN
4	a	303	GLN
4	a	587	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 67 ligands modelled in this entry, 2 are monoatomic - leaving 65 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
5	BCL	E	378[B]	1	49,54,74	1.94	14 (28%)	55,91,115	2.22	22 (40%)
13	G2O	A	802	-	68,73,73	2.59	22 (32%)	81,113,113	1.34	3 (3%)
5	BCL	F	374	1	69,74,74	1.74	14 (20%)	79,115,115	2.04	28 (35%)
13	G2O	a	803	-	68,73,73	2.63	22 (32%)	81,113,113	1.41	6 (7%)
5	BCL	G	376	1	69,74,74	1.72	15 (21%)	79,115,115	2.10	26 (32%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	BCL	a	805	-	59,64,74	1.84	16 (27%)	67,103,115	2.16	23 (34%)
7	GS0	A	801	4	69,74,74	1.74	14 (20%)	79,115,115	1.98	25 (31%)
11	LMG	a	819	-	44,44,55	0.78	2 (4%)	52,52,63	1.33	4 (7%)
5	BCL	G	378[B]	1	49,54,74	1.95	14 (28%)	55,91,115	2.22	22 (40%)
5	BCL	G	372	-	69,74,74	1.73	14 (20%)	79,115,115	2.06	27 (34%)
5	BCL	a	811	4	69,74,74	1.72	13 (18%)	79,115,115	2.09	25 (31%)
5	BCL	F	372	-	69,74,74	1.73	14 (20%)	79,115,115	2.06	26 (32%)
5	BCL	a	804	4	49,54,74	1.98	15 (30%)	55,91,115	2.25	24 (43%)
9	F39	a	817	-	66,66,66	6.15	59 (89%)	83,85,85	4.32	32 (38%)
5	BCL	E	377	1	69,74,74	1.72	15 (21%)	79,115,115	2.10	27 (34%)
5	BCL	E	372	-	69,74,74	1.73	14 (20%)	79,115,115	2.07	27 (34%)
5	BCL	A	805	-	59,64,74	1.84	16 (27%)	67,103,115	2.15	23 (34%)
5	BCL	E	373	1	69,74,74	1.73	15 (21%)	79,115,115	2.14	27 (34%)
5	BCL	G	375	1	69,74,74	1.71	15 (21%)	79,115,115	2.08	29 (36%)
13	G2O	a	802	-	68,73,73	2.59	22 (32%)	81,113,113	1.33	3 (3%)
10	LHG	A	818	-	45,45,48	0.65	1 (2%)	48,51,54	1.29	6 (12%)
5	BCL	A	814	4	49,54,74	1.96	14 (28%)	55,91,115	2.26	21 (38%)
6	SF4	B	302	3	0,12,12	-	-	-	-	-
5	BCL	F	377	1	69,74,74	1.73	15 (21%)	79,115,115	2.09	26 (32%)
5	BCL	A	806	4	69,74,74	1.71	15 (21%)	79,115,115	2.09	27 (34%)
5	BCL	A	808	4	69,74,74	1.73	14 (20%)	79,115,115	2.04	25 (31%)
5	BCL	a	810	4	49,54,74	1.91	11 (22%)	55,91,115	2.41	20 (36%)
5	BCL	A	804	4	49,54,74	1.98	14 (28%)	55,91,115	2.25	24 (43%)
5	BCL	E	374	1	69,74,74	1.74	14 (20%)	79,115,115	2.03	29 (36%)
5	BCL	E	376	1	69,74,74	1.72	15 (21%)	79,115,115	2.10	26 (32%)
5	BCL	A	809	4	49,54,74	1.97	13 (26%)	55,91,115	2.30	22 (40%)
5	BCL	E	375	1	69,74,74	1.72	16 (23%)	79,115,115	2.08	29 (36%)
5	BCL	A	812	4	69,74,74	1.73	14 (20%)	79,115,115	2.03	25 (31%)
5	BCL	E	371	1	69,74,74	1.74	14 (20%)	79,115,115	2.18	28 (35%)
5	BCL	a	814	4	49,54,74	1.96	12 (24%)	55,91,115	2.26	21 (38%)
10	LHG	a	818	-	45,45,48	0.65	1 (2%)	48,51,54	1.29	6 (12%)
13	G2O	A	803	-	68,73,73	2.62	22 (32%)	81,113,113	1.41	6 (7%)
5	BCL	G	373	1	69,74,74	1.73	16 (23%)	79,115,115	2.14	27 (34%)
5	BCL	F	375	1	69,74,74	1.71	15 (21%)	79,115,115	2.08	29 (36%)
5	BCL	G	374	1	69,74,74	1.74	14 (20%)	79,115,115	2.04	28 (35%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	BCL	a	808	4	69,74,74	1.73	14 (20%)	79,115,115	2.03	24 (30%)
5	BCL	a	807	-	49,54,74	1.96	13 (26%)	55,91,115	2.21	21 (38%)
5	BCL	G	371	1	69,74,74	1.74	15 (21%)	79,115,115	2.19	28 (35%)
5	BCL	A	815	-	49,54,74	1.97	14 (28%)	55,91,115	2.25	22 (40%)
5	BCL	a	809	4	49,54,74	1.97	14 (28%)	55,91,115	2.30	22 (40%)
5	BCL	F	373	1	69,74,74	1.73	15 (21%)	79,115,115	2.14	27 (34%)
5	BCL	a	806	4	69,74,74	1.71	15 (21%)	79,115,115	2.09	27 (34%)
5	BCL	A	810	4	49,54,74	1.91	12 (24%)	55,91,115	2.40	20 (36%)
5	BCL	A	813	-	49,54,74	1.99	15 (30%)	55,91,115	2.24	23 (41%)
8	F26	a	816	-	40,40,40	1.76	12 (30%)	49,50,50	2.09	12 (24%)
5	BCL	a	813	-	49,54,74	1.98	14 (28%)	55,91,115	2.24	22 (40%)
6	SF4	A	821	4	0,12,12	-	-	-	-	-
5	BCL	a	815	-	49,54,74	1.97	14 (28%)	55,91,115	2.26	22 (40%)
5	BCL	G	377	1	69,74,74	1.72	15 (21%)	79,115,115	2.10	26 (32%)
5	BCL	a	812	4	69,74,74	1.72	14 (20%)	79,115,115	2.03	25 (31%)
8	F26	A	816	-	40,40,40	1.75	12 (30%)	49,50,50	2.09	12 (24%)
5	BCL	A	811	4	69,74,74	1.72	15 (21%)	79,115,115	2.09	25 (31%)
5	BCL	F	376	1	69,74,74	1.72	15 (21%)	79,115,115	2.10	26 (32%)
5	BCL	A	807	-	49,54,74	1.95	13 (26%)	55,91,115	2.20	21 (38%)
11	LMG	A	819	-	44,44,55	0.78	2 (4%)	52,52,63	1.33	3 (5%)
6	SF4	B	301	3	0,12,12	-	-	-	-	-
7	GS0	a	801	4	69,74,74	1.75	14 (20%)	79,115,115	1.98	25 (31%)
5	BCL	F	371	1	69,74,74	1.75	15 (21%)	79,115,115	2.19	28 (35%)
5	BCL	F	378[B]	1	49,54,74	1.95	14 (28%)	55,91,115	2.22	22 (40%)
9	F39	A	817	-	66,66,66	6.14	59 (89%)	83,85,85	4.31	32 (38%)

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
5	BCL	E	378[B]	1	-	2/17/113/137	-
13	G2O	A	802	-	-	12/39/115/115	-
5	BCL	F	374	1	-	14/41/137/137	-
13	G2O	a	803	-	-	18/39/115/115	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	BCL	G	376	1	-	8/41/137/137	-
5	BCL	a	805	-	-	11/29/125/137	-
7	GS0	A	801	4	-	13/41/137/137	-
11	LMG	a	819	-	-	20/39/59/70	0/1/1/1
5	BCL	G	378[B]	1	-	2/17/113/137	-
5	BCL	G	372	-	-	17/41/137/137	-
5	BCL	a	811	4	-	16/41/137/137	-
5	BCL	F	372	-	-	17/41/137/137	-
5	BCL	a	804	4	-	9/17/113/137	-
9	F39	a	817	-	-	41/58/78/78	0/2/2/2
5	BCL	E	377	1	-	12/41/137/137	-
5	BCL	E	372	-	-	17/41/137/137	-
5	BCL	A	805	-	-	11/29/125/137	-
5	BCL	E	373	1	-	11/41/137/137	-
5	BCL	G	375	1	-	10/41/137/137	-
13	G2O	a	802	-	-	12/39/115/115	-
10	LHG	A	818	-	-	21/50/50/53	-
5	BCL	A	814	4	-	3/17/113/137	-
6	SF4	B	302	3	-	-	0/6/5/5
5	BCL	F	377	1	-	12/41/137/137	-
5	BCL	A	806	4	-	10/41/137/137	-
5	BCL	A	808	4	-	17/41/137/137	-
5	BCL	a	810	4	-	8/17/113/137	-
5	BCL	A	804	4	-	9/17/113/137	-
5	BCL	E	374	1	-	14/41/137/137	-
5	BCL	E	376	1	-	8/41/137/137	-
5	BCL	A	809	4	-	5/17/113/137	-
5	BCL	E	375	1	-	10/41/137/137	-
5	BCL	A	812	4	-	13/41/137/137	-
5	BCL	E	371	1	-	12/41/137/137	-
5	BCL	a	814	4	-	3/17/113/137	-
10	LHG	a	818	-	-	21/50/50/53	-
13	G2O	A	803	-	-	18/39/115/115	-
5	BCL	G	373	1	-	11/41/137/137	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	BCL	F	375	1	-	10/41/137/137	-
5	BCL	G	374	1	-	14/41/137/137	-
5	BCL	a	808	4	-	17/41/137/137	-
5	BCL	a	807	-	-	4/17/113/137	-
5	BCL	G	371	1	-	12/41/137/137	-
5	BCL	A	815	-	-	5/17/113/137	-
5	BCL	a	809	4	-	5/17/113/137	-
5	BCL	F	373	1	-	11/41/137/137	-
5	BCL	a	806	4	-	10/41/137/137	-
5	BCL	A	810	4	-	8/17/113/137	-
5	BCL	A	813	-	-	5/17/113/137	-
8	F26	a	816	-	-	15/36/36/36	0/1/1/1
5	BCL	a	813	-	-	5/17/113/137	-
6	SF4	A	821	4	-	-	0/6/5/5
5	BCL	a	815	-	-	5/17/113/137	-
5	BCL	G	377	1	-	12/41/137/137	-
5	BCL	a	812	4	-	14/41/137/137	-
8	F26	A	816	-	-	15/36/36/36	0/1/1/1
5	BCL	A	811	4	-	16/41/137/137	-
5	BCL	F	376	1	-	8/41/137/137	-
5	BCL	A	807	-	-	4/17/113/137	-
11	LMG	A	819	-	-	20/39/59/70	0/1/1/1
6	SF4	B	301	3	-	-	0/6/5/5
7	GS0	a	801	4	-	13/41/137/137	-
5	BCL	F	371	1	-	11/41/137/137	-
5	BCL	F	378[B]	1	-	2/17/113/137	-
9	F39	A	817	-	-	41/58/78/78	0/2/2/2

All (950) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	a	817	F39	C39-C37	12.63	1.65	1.35
9	A	817	F39	C39-C37	12.61	1.65	1.35
9	A	817	F39	C56-C53	11.83	1.63	1.33
9	a	817	F39	C56-C53	11.81	1.63	1.33
9	A	817	F39	C40-C41	11.71	1.65	1.34

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	a	817	F39	C40-C41	11.71	1.65	1.34
9	A	817	F39	C32-C35	11.63	1.65	1.34
9	a	817	F39	C32-C35	11.62	1.65	1.34
9	A	817	F39	C44-C42	11.52	1.62	1.35
9	a	817	F39	C44-C42	11.50	1.62	1.35
9	a	817	F39	C61-C58	11.40	1.62	1.35
9	A	817	F39	C61-C58	11.34	1.62	1.35
9	A	817	F39	C59-C62	11.22	1.61	1.35
9	a	817	F39	C59-C62	11.21	1.61	1.35
9	A	817	F39	C63-C64	11.21	1.63	1.34
9	a	817	F39	C63-C64	11.20	1.63	1.34
13	a	803	G2O	CAA-C2A	-10.90	1.34	1.54
13	A	803	G2O	CAA-C2A	-10.85	1.34	1.54
13	A	802	G2O	CAA-C2A	-10.48	1.35	1.54
13	a	802	G2O	CAA-C2A	-10.42	1.35	1.54
9	a	817	F39	C57-C51	10.18	1.64	1.36
9	A	817	F39	C57-C51	10.15	1.64	1.36
9	a	817	F39	C14-C13	8.81	1.66	1.53
9	A	817	F39	C14-C13	8.78	1.66	1.53
9	A	817	F39	C35-C37	8.77	1.64	1.46
9	a	817	F39	C35-C37	8.77	1.64	1.46
9	A	817	F39	C27-C20	8.33	1.62	1.35
9	a	817	F39	C27-C20	8.31	1.62	1.35
13	a	802	G2O	CMA-C3A	-8.12	1.36	1.53
13	A	802	G2O	CMA-C3A	-8.12	1.36	1.53
13	A	803	G2O	CMA-C3A	-8.04	1.36	1.53
13	a	803	G2O	CMA-C3A	-8.00	1.36	1.53
9	A	817	F39	C41-C42	7.81	1.62	1.46
9	a	817	F39	C41-C42	7.77	1.62	1.46
9	a	817	F39	C64-C62	7.75	1.62	1.46
9	A	817	F39	C64-C62	7.73	1.62	1.46
9	A	817	F39	C56-C58	7.55	1.62	1.46
9	a	817	F39	C56-C58	7.55	1.62	1.46
9	a	817	F39	C46-C53	7.53	1.63	1.47
9	A	817	F39	C46-C53	7.53	1.63	1.47
9	a	817	F39	C51-C44	6.94	1.64	1.43
9	A	817	F39	C51-C44	6.93	1.64	1.43
9	a	817	F39	C57-C59	6.92	1.64	1.43
9	A	817	F39	C57-C59	6.90	1.64	1.43
9	a	817	F39	C40-C39	6.90	1.64	1.43
9	A	817	F39	C40-C39	6.87	1.64	1.43
9	a	817	F39	C63-C61	6.77	1.64	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	a	817	F39	C32-C27	6.73	1.64	1.43
9	A	817	F39	C63-C61	6.72	1.64	1.43
9	A	817	F39	C32-C27	6.72	1.64	1.43
13	a	802	G2O	CBD-CGD	-6.51	1.32	1.52
13	A	802	G2O	CBD-CGD	-6.49	1.32	1.52
13	A	803	G2O	CBD-CGD	-6.43	1.33	1.52
13	a	803	G2O	CBD-CGD	-6.39	1.33	1.52
5	a	808	BCL	MG-ND	-5.98	1.93	2.05
5	A	808	BCL	MG-ND	-5.98	1.93	2.05
7	a	801	GS0	MG-ND	-5.97	1.94	2.05
5	E	374	BCL	MG-ND	-5.96	1.94	2.05
5	F	374	BCL	MG-ND	-5.95	1.94	2.05
5	G	371	BCL	MG-ND	-5.94	1.94	2.05
5	G	374	BCL	MG-ND	-5.93	1.94	2.05
5	F	371	BCL	MG-ND	-5.93	1.94	2.05
7	A	801	GS0	MG-ND	-5.93	1.94	2.05
5	E	372	BCL	MG-ND	-5.93	1.94	2.05
5	G	372	BCL	MG-ND	-5.91	1.94	2.05
5	E	373	BCL	MG-ND	-5.88	1.94	2.05
5	A	813	BCL	MG-ND	-5.88	1.94	2.05
5	F	372	BCL	MG-ND	-5.88	1.94	2.05
5	a	804	BCL	MG-ND	-5.86	1.94	2.05
5	G	373	BCL	MG-ND	-5.85	1.94	2.05
5	A	804	BCL	MG-ND	-5.83	1.94	2.05
5	E	371	BCL	MG-ND	-5.83	1.94	2.05
5	F	373	BCL	MG-ND	-5.83	1.94	2.05
5	A	812	BCL	MG-ND	-5.81	1.94	2.05
5	A	809	BCL	MG-ND	-5.81	1.94	2.05
5	a	813	BCL	MG-ND	-5.81	1.94	2.05
5	E	377	BCL	MG-ND	-5.80	1.94	2.05
5	A	815	BCL	MG-ND	-5.80	1.94	2.05
5	a	809	BCL	MG-ND	-5.79	1.94	2.05
5	a	815	BCL	MG-ND	-5.79	1.94	2.05
5	a	812	BCL	MG-ND	-5.78	1.94	2.05
5	G	377	BCL	MG-ND	-5.78	1.94	2.05
5	F	377	BCL	MG-ND	-5.77	1.94	2.05
5	F	376	BCL	MG-ND	-5.77	1.94	2.05
5	A	814	BCL	MG-ND	-5.77	1.94	2.05
5	G	376	BCL	MG-ND	-5.75	1.94	2.05
5	a	814	BCL	MG-ND	-5.74	1.94	2.05
9	a	817	F39	C19-C20	5.74	1.63	1.51
9	A	817	F39	C19-C20	5.73	1.63	1.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	E	376	BCL	MG-ND	-5.72	1.94	2.05
9	a	817	F39	C54-C47	5.72	1.61	1.51
5	A	811	BCL	MG-ND	-5.71	1.94	2.05
5	a	811	BCL	MG-ND	-5.71	1.94	2.05
5	E	375	BCL	MG-ND	-5.70	1.94	2.05
9	A	817	F39	C54-C47	5.69	1.61	1.51
9	A	817	F39	O2-C13	5.69	1.54	1.46
9	a	817	F39	O2-C13	5.68	1.54	1.46
5	F	375	BCL	MG-ND	-5.67	1.94	2.05
5	a	806	BCL	MG-ND	-5.67	1.94	2.05
5	A	806	BCL	MG-ND	-5.64	1.94	2.05
5	F	378[B]	BCL	MG-ND	-5.63	1.94	2.05
5	G	378[B]	BCL	MG-ND	-5.62	1.94	2.05
9	A	817	F39	C55-C48	5.61	1.61	1.51
5	G	375	BCL	MG-ND	-5.60	1.94	2.05
5	A	805	BCL	MG-ND	-5.60	1.94	2.05
5	a	805	BCL	MG-ND	-5.58	1.94	2.05
5	E	378[B]	BCL	MG-ND	-5.58	1.94	2.05
5	a	807	BCL	MG-ND	-5.57	1.94	2.05
5	A	807	BCL	MG-ND	-5.55	1.94	2.05
9	a	817	F39	C55-C48	5.53	1.61	1.51
5	a	810	BCL	MG-ND	-5.48	1.94	2.05
5	A	810	BCL	MG-ND	-5.45	1.95	2.05
9	a	817	F39	O6-C21	5.44	1.49	1.33
9	A	817	F39	O6-C21	5.42	1.49	1.33
9	A	817	F39	O1-C11	5.26	1.55	1.41
9	a	817	F39	O1-C11	5.23	1.55	1.41
9	a	817	F39	C50-C49	5.17	1.47	1.38
13	a	803	G2O	C2A-C1A	-5.17	1.40	1.52
9	A	817	F39	C50-C49	5.15	1.47	1.38
13	A	803	G2O	C2A-C1A	-5.12	1.40	1.52
7	A	801	GS0	MG-NB	-5.06	1.95	2.05
7	a	801	GS0	MG-NB	-5.06	1.95	2.05
13	a	803	G2O	C3A-C2A	-5.03	1.40	1.54
13	A	803	G2O	C3A-C2A	-5.03	1.40	1.54
9	a	817	F39	C38-C37	5.01	1.61	1.50
5	G	374	BCL	MG-NB	-5.01	1.95	2.05
9	A	817	F39	C38-C37	4.99	1.60	1.50
5	a	813	BCL	MG-NB	-4.98	1.95	2.05
5	E	374	BCL	MG-NB	-4.96	1.96	2.05
5	A	813	BCL	MG-NB	-4.95	1.96	2.05
5	F	374	BCL	MG-NB	-4.95	1.96	2.05

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	a	815	BCL	MG-NB	-4.93	1.96	2.05
5	E	371	BCL	MG-NB	-4.93	1.96	2.05
13	a	802	G2O	C3A-C2A	-4.93	1.41	1.54
5	F	372	BCL	MG-NB	-4.93	1.96	2.05
5	G	372	BCL	MG-NB	-4.92	1.96	2.05
5	A	815	BCL	MG-NB	-4.91	1.96	2.05
13	a	802	G2O	C2A-C1A	-4.91	1.41	1.52
5	a	804	BCL	MG-NB	-4.90	1.96	2.05
5	A	808	BCL	MG-NB	-4.90	1.96	2.05
5	a	809	BCL	MG-NB	-4.89	1.96	2.05
5	G	371	BCL	MG-NB	-4.89	1.96	2.05
5	E	372	BCL	MG-NB	-4.88	1.96	2.05
5	F	373	BCL	MG-NB	-4.88	1.96	2.05
13	A	802	G2O	C3A-C2A	-4.88	1.41	1.54
5	F	371	BCL	MG-NB	-4.88	1.96	2.05
5	F	377	BCL	MG-NB	-4.88	1.96	2.05
5	A	804	BCL	MG-NB	-4.88	1.96	2.05
5	A	809	BCL	MG-NB	-4.87	1.96	2.05
5	a	808	BCL	MG-NB	-4.87	1.96	2.05
5	G	373	BCL	MG-NB	-4.87	1.96	2.05
5	G	377	BCL	MG-NB	-4.86	1.96	2.05
5	a	814	BCL	MG-NB	-4.86	1.96	2.05
13	A	802	G2O	C2A-C1A	-4.86	1.41	1.52
5	E	376	BCL	MG-NB	-4.85	1.96	2.05
5	E	377	BCL	MG-NB	-4.85	1.96	2.05
5	E	373	BCL	MG-NB	-4.84	1.96	2.05
5	a	812	BCL	MG-NB	-4.84	1.96	2.05
5	F	376	BCL	MG-NB	-4.84	1.96	2.05
5	G	376	BCL	MG-NB	-4.84	1.96	2.05
5	A	812	BCL	MG-NB	-4.84	1.96	2.05
5	A	814	BCL	MG-NB	-4.84	1.96	2.05
5	a	811	BCL	MG-NB	-4.83	1.96	2.05
5	A	811	BCL	MG-NB	-4.81	1.96	2.05
13	A	803	G2O	MG-NB	-4.78	1.96	2.05
13	a	803	G2O	MG-NB	-4.77	1.96	2.05
9	a	817	F39	O1-C12	4.77	1.56	1.44
9	A	817	F39	O1-C12	4.77	1.56	1.44
5	E	378[B]	BCL	MG-NB	-4.77	1.96	2.05
9	A	817	F39	C52-C45	4.75	1.61	1.51
5	F	378[B]	BCL	MG-NB	-4.74	1.96	2.05
9	a	817	F39	C52-C45	4.74	1.61	1.51
5	A	807	BCL	OBD-CAD	4.73	1.30	1.22

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
13	a	802	G2O	MG-NB	-4.73	1.96	2.05
5	a	807	BCL	OBD-CAD	4.73	1.30	1.22
13	A	802	G2O	MG-NB	-4.73	1.96	2.05
5	E	374	BCL	OBD-CAD	4.72	1.30	1.22
5	G	378[B]	BCL	MG-NB	-4.72	1.96	2.05
5	G	375	BCL	MG-NB	-4.70	1.96	2.05
5	G	374	BCL	OBD-CAD	4.69	1.30	1.22
5	F	375	BCL	MG-NB	-4.69	1.96	2.05
5	A	811	BCL	OBD-CAD	4.68	1.30	1.22
5	E	375	BCL	OBD-CAD	4.67	1.30	1.22
5	A	806	BCL	MG-NB	-4.67	1.96	2.05
5	F	373	BCL	OBD-CAD	4.67	1.30	1.22
5	a	810	BCL	MG-NB	-4.67	1.96	2.05
5	a	805	BCL	OBD-CAD	4.67	1.30	1.22
5	a	811	BCL	OBD-CAD	4.66	1.30	1.22
5	F	374	BCL	OBD-CAD	4.66	1.30	1.22
5	A	805	BCL	OBD-CAD	4.66	1.30	1.22
5	A	814	BCL	OBD-CAD	4.66	1.30	1.22
5	a	808	BCL	OBD-CAD	4.66	1.30	1.22
5	a	806	BCL	MG-NB	-4.66	1.96	2.05
5	E	371	BCL	OBD-CAD	4.65	1.30	1.22
5	E	375	BCL	MG-NB	-4.65	1.96	2.05
5	G	376	BCL	OBD-CAD	4.64	1.30	1.22
5	F	371	BCL	OBD-CAD	4.64	1.30	1.22
5	A	812	BCL	OBD-CAD	4.64	1.30	1.22
5	G	371	BCL	OBD-CAD	4.63	1.30	1.22
5	A	810	BCL	MG-NB	-4.63	1.96	2.05
5	G	375	BCL	OBD-CAD	4.63	1.30	1.22
5	G	373	BCL	OBD-CAD	4.63	1.30	1.22
5	a	814	BCL	OBD-CAD	4.63	1.30	1.22
5	F	376	BCL	OBD-CAD	4.62	1.30	1.22
5	F	377	BCL	OBD-CAD	4.62	1.30	1.22
5	a	805	BCL	MG-NB	-4.61	1.96	2.05
5	E	373	BCL	OBD-CAD	4.61	1.30	1.22
5	A	808	BCL	OBD-CAD	4.60	1.30	1.22
5	A	805	BCL	MG-NB	-4.60	1.96	2.05
5	a	807	BCL	MG-NB	-4.60	1.96	2.05
5	F	372	BCL	OBD-CAD	4.60	1.30	1.22
5	G	378[B]	BCL	OBD-CAD	4.59	1.30	1.22
5	F	375	BCL	OBD-CAD	4.59	1.30	1.22
5	A	806	BCL	OBD-CAD	4.59	1.30	1.22
5	A	815	BCL	OBD-CAD	4.59	1.30	1.22

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	E	378[B]	BCL	OBD-CAD	4.59	1.30	1.22
5	E	376	BCL	OBD-CAD	4.59	1.30	1.22
5	a	813	BCL	OBD-CAD	4.59	1.30	1.22
5	F	378[B]	BCL	OBD-CAD	4.59	1.30	1.22
5	a	804	BCL	OBD-CAD	4.58	1.30	1.22
5	G	377	BCL	OBD-CAD	4.58	1.30	1.22
5	E	372	BCL	OBD-CAD	4.58	1.30	1.22
5	A	804	BCL	OBD-CAD	4.58	1.30	1.22
5	a	806	BCL	OBD-CAD	4.58	1.30	1.22
5	a	815	BCL	OBD-CAD	4.57	1.30	1.22
5	G	372	BCL	OBD-CAD	4.57	1.30	1.22
5	a	809	BCL	OBD-CAD	4.57	1.30	1.22
5	A	807	BCL	MG-NB	-4.57	1.96	2.05
5	a	812	BCL	OBD-CAD	4.57	1.30	1.22
5	A	809	BCL	OBD-CAD	4.56	1.30	1.22
5	A	813	BCL	OBD-CAD	4.56	1.30	1.22
7	a	801	GS0	OBD-CAD	4.51	1.30	1.22
5	E	377	BCL	OBD-CAD	4.50	1.30	1.22
7	A	801	GS0	OBD-CAD	4.49	1.30	1.22
9	a	817	F39	C46-C48	4.36	1.47	1.41
9	A	817	F39	C46-C48	4.36	1.47	1.41
9	a	817	F39	O3-C8	4.29	1.53	1.43
9	A	817	F39	O3-C8	4.27	1.53	1.43
9	a	817	F39	C60-C58	4.24	1.59	1.50
9	A	817	F39	C60-C58	4.23	1.59	1.50
9	A	817	F39	C22-C21	4.18	1.62	1.50
9	a	817	F39	C22-C21	4.17	1.62	1.50
9	a	817	F39	C46-C45	4.16	1.47	1.41
5	A	815	BCL	O1D-CGD	-4.14	1.10	1.21
5	a	815	BCL	O1D-CGD	-4.14	1.10	1.21
5	a	810	BCL	O1D-CGD	-4.11	1.10	1.21
7	a	801	GS0	O1D-CGD	-4.11	1.10	1.21
5	F	376	BCL	O1D-CGD	-4.10	1.10	1.21
5	A	810	BCL	OBD-CAD	4.10	1.29	1.22
9	A	817	F39	C46-C45	4.09	1.47	1.41
5	E	371	BCL	O1D-CGD	-4.09	1.10	1.21
5	F	373	BCL	O1D-CGD	-4.09	1.11	1.21
5	A	809	BCL	O1D-CGD	-4.09	1.11	1.21
5	A	812	BCL	O1D-CGD	-4.09	1.11	1.21
5	G	373	BCL	O1D-CGD	-4.09	1.11	1.21
5	F	371	BCL	O1D-CGD	-4.09	1.11	1.21
5	A	810	BCL	O1D-CGD	-4.08	1.11	1.21

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	811	BCL	O1D-CGD	-4.08	1.11	1.21
5	a	810	BCL	OBD-CAD	4.08	1.29	1.22
5	G	372	BCL	O1D-CGD	-4.07	1.11	1.21
5	A	813	BCL	O1D-CGD	-4.07	1.11	1.21
5	a	811	BCL	O1D-CGD	-4.07	1.11	1.21
5	F	377	BCL	O1D-CGD	-4.07	1.11	1.21
5	G	374	BCL	O1D-CGD	-4.07	1.11	1.21
5	E	375	BCL	O1D-CGD	-4.07	1.11	1.21
5	E	372	BCL	O1D-CGD	-4.07	1.11	1.21
5	G	375	BCL	O1D-CGD	-4.06	1.11	1.21
7	A	801	GS0	O1D-CGD	-4.06	1.11	1.21
5	G	376	BCL	O1D-CGD	-4.06	1.11	1.21
5	a	813	BCL	O1D-CGD	-4.06	1.11	1.21
5	E	374	BCL	O1D-CGD	-4.05	1.11	1.21
5	F	372	BCL	O1D-CGD	-4.04	1.11	1.21
5	G	371	BCL	O1D-CGD	-4.04	1.11	1.21
5	a	805	BCL	O1D-CGD	-4.04	1.11	1.21
5	G	377	BCL	O1D-CGD	-4.04	1.11	1.21
5	a	812	BCL	O1D-CGD	-4.04	1.11	1.21
5	E	376	BCL	O1D-CGD	-4.04	1.11	1.21
5	E	373	BCL	O1D-CGD	-4.04	1.11	1.21
5	F	375	BCL	O1D-CGD	-4.04	1.11	1.21
5	a	808	BCL	O1D-CGD	-4.04	1.11	1.21
5	A	804	BCL	O1D-CGD	-4.03	1.11	1.21
5	F	374	BCL	O1D-CGD	-4.03	1.11	1.21
5	A	805	BCL	O1D-CGD	-4.03	1.11	1.21
5	E	377	BCL	O1D-CGD	-4.03	1.11	1.21
5	a	809	BCL	O1D-CGD	-4.03	1.11	1.21
5	a	804	BCL	O1D-CGD	-4.03	1.11	1.21
5	A	806	BCL	O1D-CGD	-4.02	1.11	1.21
5	a	806	BCL	O1D-CGD	-4.01	1.11	1.21
5	A	808	BCL	O1D-CGD	-4.01	1.11	1.21
5	F	378[B]	BCL	O1D-CGD	-4.00	1.11	1.21
5	G	378[B]	BCL	O1D-CGD	-4.00	1.11	1.21
5	A	814	BCL	O1D-CGD	-3.99	1.11	1.21
5	a	814	BCL	O1D-CGD	-3.99	1.11	1.21
9	A	817	F39	C17-C13	3.98	1.61	1.52
9	a	817	F39	C17-C13	3.97	1.61	1.52
5	E	378[B]	BCL	O1D-CGD	-3.97	1.11	1.21
5	A	807	BCL	O1D-CGD	-3.94	1.11	1.21
5	a	807	BCL	O1D-CGD	-3.93	1.11	1.21
13	a	803	G2O	C10-C8	-3.91	1.42	1.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
13	a	802	G2O	C10-C8	-3.89	1.42	1.54
13	A	803	G2O	C10-C8	-3.89	1.42	1.54
13	A	802	G2O	C10-C8	-3.89	1.42	1.54
5	A	807	BCL	O2D-CED	3.87	1.53	1.45
9	a	817	F39	C65-C62	3.86	1.58	1.50
9	A	817	F39	C18-C19	3.86	1.66	1.52
9	a	817	F39	C14-C18	3.85	1.65	1.52
5	a	807	BCL	O2D-CED	3.84	1.53	1.45
9	A	817	F39	C65-C62	3.84	1.58	1.50
9	A	817	F39	C14-C18	3.84	1.65	1.52
9	A	817	F39	C16-C13	3.83	1.61	1.52
9	a	817	F39	C43-C42	3.83	1.58	1.50
9	a	817	F39	C16-C13	3.83	1.61	1.52
9	a	817	F39	C18-C19	3.83	1.66	1.52
13	a	802	G2O	C4B-NB	3.82	1.42	1.37
9	A	817	F39	C47-C45	3.81	1.47	1.40
9	A	817	F39	C43-C42	3.80	1.58	1.50
13	A	802	G2O	C4B-NB	3.80	1.42	1.37
9	a	817	F39	C47-C45	3.78	1.47	1.40
13	a	803	G2O	CHC-C4B	-3.76	1.31	1.38
13	A	803	G2O	CHC-C4B	-3.73	1.31	1.38
5	a	810	BCL	O2D-CED	3.69	1.53	1.45
13	A	802	G2O	CHC-C4B	-3.67	1.31	1.38
5	A	810	BCL	O2D-CED	3.67	1.53	1.45
13	a	802	G2O	CHC-C4B	-3.65	1.31	1.38
7	A	801	GS0	O2D-CED	3.64	1.53	1.45
5	G	374	BCL	O2D-CED	3.64	1.53	1.45
5	F	374	BCL	O2D-CED	3.64	1.53	1.45
5	E	373	BCL	O2D-CED	3.63	1.53	1.45
5	E	377	BCL	O2D-CED	3.63	1.53	1.45
5	A	813	BCL	O2D-CED	3.62	1.53	1.45
5	G	375	BCL	O2D-CED	3.62	1.53	1.45
9	A	817	F39	C10-C12	3.62	1.60	1.53
7	a	801	GS0	O2D-CED	3.62	1.53	1.45
5	F	377	BCL	O2D-CED	3.62	1.53	1.45
5	A	814	BCL	O2D-CED	3.61	1.53	1.45
5	E	375	BCL	O2D-CED	3.61	1.53	1.45
5	a	804	BCL	O2D-CED	3.61	1.53	1.45
5	E	374	BCL	O2D-CED	3.61	1.53	1.45
5	A	811	BCL	O2D-CED	3.60	1.53	1.45
5	F	375	BCL	O2D-CED	3.60	1.53	1.45
5	G	373	BCL	O2D-CED	3.60	1.53	1.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	a	817	F39	C10-C12	3.60	1.60	1.53
5	a	809	BCL	O2D-CED	3.60	1.53	1.45
5	a	805	BCL	O2D-CED	3.60	1.53	1.45
5	A	804	BCL	O2D-CED	3.60	1.53	1.45
5	A	805	BCL	O2D-CED	3.60	1.53	1.45
5	F	378[B]	BCL	O2D-CED	3.59	1.53	1.45
5	a	812	BCL	O2D-CED	3.59	1.53	1.45
5	A	806	BCL	O2D-CED	3.59	1.53	1.45
5	A	809	BCL	O2D-CED	3.59	1.53	1.45
5	a	806	BCL	O2D-CED	3.59	1.53	1.45
5	A	812	BCL	O2D-CED	3.59	1.53	1.45
5	a	813	BCL	O2D-CED	3.59	1.53	1.45
5	a	814	BCL	O2D-CED	3.58	1.53	1.45
5	A	815	BCL	O2D-CED	3.58	1.53	1.45
13	a	803	G2O	C14-C13	-3.58	1.41	1.52
5	a	811	BCL	O2D-CED	3.58	1.53	1.45
5	G	377	BCL	O2D-CED	3.57	1.53	1.45
5	F	371	BCL	O2D-CED	3.57	1.53	1.45
5	F	373	BCL	O2D-CED	3.57	1.53	1.45
13	A	803	G2O	C14-C13	-3.57	1.41	1.52
5	F	376	BCL	O2D-CED	3.57	1.53	1.45
5	a	815	BCL	O2D-CED	3.56	1.53	1.45
9	a	817	F39	C49-C47	3.55	1.46	1.39
5	F	372	BCL	O2D-CED	3.55	1.53	1.45
5	G	371	BCL	O2D-CED	3.55	1.53	1.45
5	G	378[B]	BCL	O2D-CED	3.55	1.53	1.45
5	E	378[B]	BCL	O2D-CED	3.55	1.53	1.45
5	G	372	BCL	O2D-CED	3.55	1.53	1.45
5	a	808	BCL	O2D-CED	3.55	1.53	1.45
5	E	371	BCL	O2D-CED	3.54	1.53	1.45
5	A	808	BCL	O2D-CED	3.54	1.53	1.45
5	G	376	BCL	O2D-CED	3.53	1.53	1.45
9	a	817	F39	O5-C10	3.53	1.51	1.43
9	A	817	F39	C49-C47	3.52	1.46	1.39
5	E	372	BCL	O2D-CED	3.52	1.53	1.45
13	A	803	G2O	C4B-NB	3.51	1.42	1.37
9	A	817	F39	O5-C10	3.50	1.51	1.43
13	a	803	G2O	C4B-NB	3.50	1.42	1.37
9	A	817	F39	C50-C48	3.50	1.46	1.39
9	a	817	F39	C50-C48	3.49	1.46	1.39
5	E	376	BCL	O2D-CED	3.49	1.53	1.45
13	a	802	G2O	C14-C13	-3.46	1.42	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
13	A	802	G2O	C14-C13	-3.45	1.42	1.52
13	a	803	G2O	C3A-C4A	-3.36	1.40	1.51
13	a	802	G2O	C9-C8	-3.34	1.41	1.53
13	A	803	G2O	C3A-C4A	-3.33	1.40	1.51
13	A	802	G2O	C9-C8	-3.33	1.41	1.53
13	A	803	G2O	C9-C8	-3.29	1.41	1.53
13	a	803	G2O	C9-C8	-3.28	1.41	1.53
13	a	802	G2O	C3A-C4A	-3.28	1.41	1.51
13	A	802	G2O	C3A-C4A	-3.27	1.41	1.51
9	a	817	F39	C25-C20	3.20	1.58	1.50
8	A	816	F26	C15-C19	3.19	1.52	1.46
8	a	816	F26	C15-C19	3.19	1.52	1.46
9	A	817	F39	C25-C20	3.18	1.58	1.50
5	A	811	BCL	O2A-CGA	-3.12	1.24	1.33
5	F	374	BCL	O2A-CGA	-3.11	1.24	1.33
5	G	374	BCL	O2A-CGA	-3.10	1.24	1.33
5	a	811	BCL	O2A-CGA	-3.09	1.24	1.33
5	E	372	BCL	O2A-CGA	-3.09	1.24	1.33
5	E	371	BCL	O2A-CGA	-3.08	1.24	1.33
5	E	374	BCL	O2A-CGA	-3.08	1.24	1.33
5	A	806	BCL	O2A-CGA	-3.08	1.24	1.33
5	F	371	BCL	O2A-CGA	-3.08	1.24	1.33
5	A	812	BCL	O2A-CGA	-3.08	1.24	1.33
5	G	372	BCL	O2A-CGA	-3.07	1.24	1.33
5	F	372	BCL	O2A-CGA	-3.07	1.24	1.33
5	E	373	BCL	O2A-CGA	-3.06	1.24	1.33
5	F	373	BCL	O2A-CGA	-3.06	1.24	1.33
5	G	371	BCL	O2A-CGA	-3.05	1.24	1.33
5	a	806	BCL	O2A-CGA	-3.05	1.24	1.33
13	a	803	G2O	C8-C7	-3.05	1.42	1.51
5	E	375	BCL	O2A-CGA	-3.04	1.24	1.33
5	a	812	BCL	O2A-CGA	-3.04	1.24	1.33
5	G	373	BCL	O2A-CGA	-3.03	1.24	1.33
5	G	375	BCL	O2A-CGA	-3.03	1.24	1.33
7	a	801	GS0	O2A-CGA	-3.03	1.24	1.33
5	F	376	BCL	O2A-CGA	-3.03	1.24	1.33
5	G	376	BCL	O2A-CGA	-3.03	1.24	1.33
7	A	801	GS0	O2A-CGA	-3.02	1.24	1.33
13	A	803	G2O	C8-C7	-3.02	1.42	1.51
5	E	376	BCL	O2A-CGA	-3.02	1.24	1.33
7	a	801	GS0	O2D-CGD	-3.01	1.25	1.33
5	F	377	BCL	O2A-CGA	-3.01	1.24	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	F	372	BCL	O2D-CGD	-3.01	1.25	1.33
5	G	377	BCL	O2A-CGA	-3.00	1.24	1.33
5	G	373	BCL	O2D-CGD	-3.00	1.25	1.33
5	G	372	BCL	O2D-CGD	-3.00	1.25	1.33
7	A	801	GS0	O2D-CGD	-3.00	1.25	1.33
5	A	808	BCL	O2D-CGD	-3.00	1.25	1.33
5	F	375	BCL	O2A-CGA	-2.99	1.24	1.33
5	E	372	BCL	O2D-CGD	-2.99	1.25	1.33
5	E	371	BCL	O2D-CGD	-2.99	1.25	1.33
5	E	377	BCL	O2A-CGA	-2.98	1.24	1.33
5	a	808	BCL	O2D-CGD	-2.98	1.25	1.33
5	a	812	BCL	O2D-CGD	-2.98	1.25	1.33
8	a	816	F26	C39-C37	2.98	1.52	1.43
5	G	371	BCL	O2D-CGD	-2.98	1.25	1.33
5	G	377	BCL	O2D-CGD	-2.97	1.25	1.33
8	A	816	F26	C39-C37	2.97	1.52	1.43
5	E	373	BCL	O2D-CGD	-2.97	1.25	1.33
5	a	813	BCL	O2D-CGD	-2.97	1.25	1.33
9	a	817	F39	O4-C9	2.97	1.50	1.43
5	F	373	BCL	O2D-CGD	-2.97	1.25	1.33
5	F	377	BCL	O2D-CGD	-2.97	1.25	1.33
9	A	817	F39	O4-C9	2.96	1.50	1.43
5	G	376	BCL	O2D-CGD	-2.96	1.25	1.33
5	F	376	BCL	O2D-CGD	-2.96	1.25	1.33
5	A	810	BCL	O2D-CGD	-2.96	1.26	1.33
5	A	812	BCL	O2D-CGD	-2.96	1.26	1.33
5	E	377	BCL	O2D-CGD	-2.95	1.26	1.33
5	a	805	BCL	O2A-CGA	-2.95	1.24	1.33
5	A	805	BCL	O2A-CGA	-2.95	1.24	1.33
5	a	804	BCL	O2D-CGD	-2.95	1.26	1.33
5	A	814	BCL	O2D-CGD	-2.95	1.26	1.33
5	A	815	BCL	O2D-CGD	-2.94	1.26	1.33
5	a	810	BCL	O2D-CGD	-2.94	1.26	1.33
5	F	375	BCL	O2D-CGD	-2.94	1.26	1.33
5	A	804	BCL	O2D-CGD	-2.94	1.26	1.33
5	a	811	BCL	O2D-CGD	-2.94	1.26	1.33
8	a	816	F26	C28-C31	2.94	1.52	1.46
5	E	375	BCL	O2D-CGD	-2.94	1.26	1.33
5	A	811	BCL	O2D-CGD	-2.94	1.26	1.33
5	A	813	BCL	O2D-CGD	-2.93	1.26	1.33
5	G	375	BCL	O2D-CGD	-2.93	1.26	1.33
5	a	814	BCL	O2D-CGD	-2.93	1.26	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	F	371	BCL	O2D-CGD	-2.93	1.26	1.33
5	F	374	BCL	O2D-CGD	-2.92	1.26	1.33
5	E	376	BCL	O2D-CGD	-2.92	1.26	1.33
5	G	374	BCL	O2D-CGD	-2.92	1.26	1.33
5	F	378[B]	BCL	O2D-CGD	-2.92	1.26	1.33
5	G	378[B]	BCL	O2D-CGD	-2.91	1.26	1.33
5	a	815	BCL	O2D-CGD	-2.91	1.26	1.33
5	a	809	BCL	O2D-CGD	-2.91	1.26	1.33
8	a	816	F26	C22-C18	2.91	1.52	1.43
5	A	809	BCL	O2D-CGD	-2.90	1.26	1.33
5	a	808	BCL	O2A-CGA	-2.90	1.24	1.33
8	A	816	F26	C22-C18	2.89	1.52	1.43
8	a	816	F26	C38-C33	2.89	1.52	1.43
8	A	816	F26	C28-C31	2.88	1.52	1.46
5	A	808	BCL	O2A-CGA	-2.88	1.25	1.33
5	A	805	BCL	O2D-CGD	-2.88	1.26	1.33
5	E	374	BCL	O2D-CGD	-2.88	1.26	1.33
5	a	805	BCL	O2D-CGD	-2.88	1.26	1.33
8	A	816	F26	C38-C33	2.87	1.52	1.43
5	A	806	BCL	O2D-CGD	-2.86	1.26	1.33
5	a	806	BCL	O2D-CGD	-2.86	1.26	1.33
8	a	816	F26	C32-C30	2.86	1.52	1.43
5	A	812	BCL	O1A-CGA	-2.86	1.14	1.22
5	E	378[B]	BCL	O2D-CGD	-2.85	1.26	1.33
8	A	816	F26	C32-C30	2.85	1.52	1.43
9	a	817	F39	C23-C22	2.84	1.62	1.52
5	a	806	BCL	O1A-CGA	-2.84	1.14	1.22
5	E	372	BCL	O1A-CGA	-2.84	1.14	1.22
5	a	811	BCL	O1A-CGA	-2.84	1.14	1.22
5	A	811	BCL	O1A-CGA	-2.84	1.14	1.22
8	a	816	F26	C27-C24	2.83	1.52	1.43
7	A	801	GS0	O1A-CGA	-2.83	1.14	1.22
5	G	373	BCL	O1A-CGA	-2.83	1.14	1.22
5	a	812	BCL	O1A-CGA	-2.83	1.14	1.22
5	E	374	BCL	O1A-CGA	-2.82	1.14	1.22
5	A	806	BCL	O1A-CGA	-2.82	1.14	1.22
7	a	801	GS0	O1A-CGA	-2.82	1.14	1.22
9	A	817	F39	C23-C22	2.82	1.62	1.52
5	G	376	BCL	O1A-CGA	-2.82	1.14	1.22
8	A	816	F26	C27-C24	2.82	1.51	1.43
5	E	373	BCL	O1A-CGA	-2.82	1.14	1.22
5	F	373	BCL	O1A-CGA	-2.81	1.14	1.22

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	F	376	BCL	O1A-CGA	-2.81	1.14	1.22
5	E	377	BCL	O1A-CGA	-2.81	1.14	1.22
5	F	377	BCL	O1A-CGA	-2.80	1.14	1.22
5	G	372	BCL	O1A-CGA	-2.80	1.14	1.22
5	F	372	BCL	O1A-CGA	-2.80	1.14	1.22
5	a	808	BCL	O1A-CGA	-2.80	1.14	1.22
5	G	371	BCL	O1A-CGA	-2.79	1.14	1.22
5	F	371	BCL	O1A-CGA	-2.79	1.14	1.22
5	E	376	BCL	O1A-CGA	-2.79	1.14	1.22
5	G	377	BCL	O1A-CGA	-2.78	1.14	1.22
5	F	374	BCL	O1A-CGA	-2.78	1.14	1.22
5	G	374	BCL	O1A-CGA	-2.78	1.14	1.22
5	A	808	BCL	O1A-CGA	-2.78	1.14	1.22
5	E	371	BCL	O1A-CGA	-2.77	1.14	1.22
5	a	805	BCL	O1A-CGA	-2.77	1.14	1.22
5	A	805	BCL	O1A-CGA	-2.77	1.14	1.22
5	F	375	BCL	O1A-CGA	-2.75	1.14	1.22
13	A	802	G2O	MG-NA	2.74	2.12	2.06
5	G	375	BCL	O1A-CGA	-2.74	1.14	1.22
5	E	375	BCL	O1A-CGA	-2.72	1.14	1.22
9	A	817	F39	C11-C9	2.70	1.60	1.52
13	a	802	G2O	MG-NA	2.70	2.12	2.06
13	A	802	G2O	C8-C7	-2.69	1.43	1.51
13	A	803	G2O	MG-ND	-2.69	2.00	2.05
9	a	817	F39	C11-C9	2.69	1.60	1.52
13	a	803	G2O	MG-ND	-2.69	2.00	2.05
13	a	802	G2O	C8-C7	-2.69	1.43	1.51
5	E	371	BCL	C1D-C2D	-2.68	1.40	1.45
5	G	371	BCL	C1D-C2D	-2.66	1.40	1.45
5	F	371	BCL	C1D-C2D	-2.64	1.40	1.45
5	a	807	BCL	O2D-CGD	-2.64	1.26	1.33
8	A	816	F26	C35-C34	2.64	1.51	1.46
13	a	802	G2O	MG-ND	-2.64	2.00	2.05
13	A	803	G2O	MG-NA	2.64	2.12	2.06
8	a	816	F26	C35-C34	2.63	1.51	1.46
13	a	802	G2O	CMB-C2B	-2.63	1.45	1.50
13	a	803	G2O	MG-NA	2.62	2.12	2.06
13	A	802	G2O	MG-ND	-2.62	2.00	2.05
5	A	807	BCL	O2D-CGD	-2.60	1.26	1.33
7	a	801	GS0	OBB-CAB	-2.59	1.17	1.23
13	A	802	G2O	CMB-C2B	-2.58	1.45	1.50
7	a	801	GS0	C1D-C2D	-2.56	1.40	1.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
8	a	816	F26	C2-C9	2.56	1.53	1.47
13	a	803	G2O	CMB-C2B	-2.56	1.45	1.50
5	E	376	BCL	C1D-C2D	-2.56	1.40	1.45
7	A	801	GS0	OBB-CAB	-2.56	1.17	1.23
5	E	375	BCL	OBB-CAB	-2.54	1.17	1.23
5	A	807	BCL	C3B-CAB	2.54	1.54	1.46
5	A	808	BCL	C1B-C2B	-2.53	1.37	1.43
5	a	807	BCL	C3B-CAB	2.53	1.54	1.46
7	A	801	GS0	C1D-C2D	-2.53	1.40	1.45
8	A	816	F26	C2-C9	2.53	1.53	1.47
13	A	803	G2O	CMB-C2B	-2.51	1.45	1.50
5	G	376	BCL	C1D-C2D	-2.51	1.40	1.45
5	a	808	BCL	C1B-C2B	-2.50	1.37	1.43
5	F	376	BCL	C1D-C2D	-2.49	1.40	1.45
5	F	375	BCL	OBB-CAB	-2.49	1.17	1.23
5	G	375	BCL	OBB-CAB	-2.49	1.17	1.23
5	A	813	BCL	C1D-C2D	-2.49	1.40	1.45
5	G	374	BCL	OBB-CAB	-2.49	1.17	1.23
8	A	816	F26	C25-C26	2.49	1.51	1.46
5	a	812	BCL	OBB-CAB	-2.48	1.17	1.23
5	G	377	BCL	OBB-CAB	-2.48	1.17	1.23
5	a	811	BCL	OBB-CAB	-2.48	1.17	1.23
5	G	378[B]	BCL	C3B-CAB	2.48	1.54	1.46
5	F	374	BCL	OBB-CAB	-2.48	1.17	1.23
5	E	372	BCL	C3B-CAB	2.48	1.54	1.46
5	E	376	BCL	OBB-CAB	-2.47	1.17	1.23
5	A	811	BCL	OBB-CAB	-2.47	1.17	1.23
5	E	378[B]	BCL	C3B-CAB	2.47	1.54	1.46
5	A	812	BCL	OBB-CAB	-2.47	1.17	1.23
5	F	376	BCL	OBB-CAB	-2.47	1.17	1.23
5	a	813	BCL	OBB-CAB	-2.47	1.17	1.23
5	G	373	BCL	C1D-C2D	-2.46	1.40	1.45
5	G	377	BCL	C1D-C2D	-2.46	1.40	1.45
5	a	806	BCL	C1D-C2D	-2.46	1.40	1.45
5	a	813	BCL	C1D-C2D	-2.46	1.40	1.45
8	a	816	F26	C25-C26	2.46	1.51	1.46
5	E	374	BCL	OBB-CAB	-2.46	1.17	1.23
5	G	371	BCL	OBB-CAB	-2.46	1.17	1.23
5	F	378[B]	BCL	C3B-CAB	2.46	1.54	1.46
5	F	371	BCL	OBB-CAB	-2.46	1.17	1.23
5	a	805	BCL	C1D-C2D	-2.46	1.40	1.45
5	G	376	BCL	OBB-CAB	-2.46	1.17	1.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	F	377	BCL	OBB-CAB	-2.46	1.17	1.23
5	A	813	BCL	C1B-C2B	-2.46	1.37	1.43
5	A	805	BCL	OBB-CAB	-2.45	1.17	1.23
5	E	377	BCL	OBB-CAB	-2.45	1.17	1.23
13	A	802	G2O	C1B-C2B	2.45	1.50	1.45
5	A	804	BCL	OBB-CAB	-2.45	1.17	1.23
5	a	809	BCL	OBB-CAB	-2.45	1.17	1.23
5	E	373	BCL	OBB-CAB	-2.45	1.17	1.23
5	A	813	BCL	OBB-CAB	-2.45	1.17	1.23
5	F	373	BCL	OBB-CAB	-2.45	1.17	1.23
5	a	813	BCL	C1B-C2B	-2.45	1.37	1.43
5	F	377	BCL	C1D-C2D	-2.45	1.40	1.45
5	A	815	BCL	OBB-CAB	-2.45	1.17	1.23
5	E	371	BCL	OBB-CAB	-2.45	1.17	1.23
5	G	372	BCL	C3B-CAB	2.45	1.54	1.46
5	a	815	BCL	OBB-CAB	-2.44	1.17	1.23
5	G	373	BCL	OBB-CAB	-2.44	1.17	1.23
5	a	810	BCL	OBB-CAB	-2.44	1.17	1.23
5	a	806	BCL	C1B-C2B	-2.44	1.37	1.43
5	A	810	BCL	C3B-CAB	2.44	1.54	1.46
5	a	810	BCL	C3B-CAB	2.44	1.54	1.46
5	E	373	BCL	C1D-C2D	-2.44	1.40	1.45
5	a	812	BCL	C3B-CAB	2.44	1.54	1.46
5	A	809	BCL	OBB-CAB	-2.43	1.17	1.23
5	a	805	BCL	OBB-CAB	-2.43	1.17	1.23
5	F	372	BCL	C3B-CAB	2.43	1.54	1.46
5	A	806	BCL	C3B-CAB	2.43	1.54	1.46
5	a	813	BCL	O1A-CGA	-2.43	1.14	1.22
5	A	806	BCL	OBB-CAB	-2.43	1.17	1.23
5	F	372	BCL	OBB-CAB	-2.43	1.17	1.23
5	A	814	BCL	C3B-CAB	2.43	1.54	1.46
5	A	810	BCL	O1A-CGA	-2.42	1.14	1.22
5	A	809	BCL	C1D-C2D	-2.42	1.40	1.45
5	A	805	BCL	C3B-CAB	2.42	1.54	1.46
5	F	378[B]	BCL	O1A-CGA	-2.42	1.14	1.22
5	F	373	BCL	C1D-C2D	-2.42	1.40	1.45
5	E	378[B]	BCL	O1A-CGA	-2.42	1.14	1.22
5	E	378[B]	BCL	OBB-CAB	-2.42	1.17	1.23
5	E	376	BCL	C3B-CAB	2.42	1.54	1.46
5	G	376	BCL	C3B-CAB	2.42	1.54	1.46
5	a	806	BCL	C3B-CAB	2.42	1.54	1.46
5	F	378[B]	BCL	OBB-CAB	-2.42	1.17	1.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
13	a	802	G2O	C1B-C2B	2.42	1.50	1.45
5	A	806	BCL	C1B-C2B	-2.42	1.37	1.43
5	G	378[B]	BCL	O1A-CGA	-2.42	1.14	1.22
5	A	813	BCL	O1A-CGA	-2.42	1.14	1.22
5	A	808	BCL	OBB-CAB	-2.42	1.17	1.23
5	F	375	BCL	C1D-C2D	-2.42	1.40	1.45
13	A	803	G2O	C1A-CHA	2.42	1.42	1.37
5	A	804	BCL	C1B-C2B	-2.42	1.37	1.43
5	E	377	BCL	C3B-CAB	2.42	1.54	1.46
5	a	809	BCL	C1D-C2D	-2.41	1.40	1.45
5	E	372	BCL	OBB-CAB	-2.41	1.17	1.23
5	a	811	BCL	C3B-CAB	2.41	1.54	1.46
5	a	810	BCL	O1A-CGA	-2.41	1.14	1.22
5	E	377	BCL	C1D-C2D	-2.41	1.40	1.45
5	a	810	BCL	C1B-C2B	-2.41	1.37	1.43
5	A	812	BCL	C3B-CAB	2.41	1.54	1.46
5	A	806	BCL	C1D-C2D	-2.41	1.40	1.45
5	a	805	BCL	C3B-CAB	2.41	1.54	1.46
5	E	375	BCL	C1D-C2D	-2.41	1.40	1.45
5	a	808	BCL	OBB-CAB	-2.40	1.17	1.23
5	F	373	BCL	C3B-CAB	2.40	1.54	1.46
5	A	812	BCL	C1D-C2D	-2.40	1.40	1.45
5	A	810	BCL	C1B-C2B	-2.40	1.37	1.43
5	a	804	BCL	C1B-C2B	-2.40	1.37	1.43
8	A	816	F26	C30-C26	-2.40	1.30	1.35
5	G	378[B]	BCL	OBB-CAB	-2.40	1.17	1.23
5	a	804	BCL	OBB-CAB	-2.40	1.17	1.23
5	A	804	BCL	C1D-C2D	-2.40	1.40	1.45
5	a	807	BCL	O1A-CGA	-2.40	1.14	1.22
5	F	376	BCL	C3B-CAB	2.40	1.54	1.46
5	A	811	BCL	C3B-CAB	2.40	1.54	1.46
5	G	373	BCL	C3B-CAB	2.40	1.54	1.46
13	A	803	G2O	C1B-C2B	2.40	1.50	1.45
5	a	815	BCL	C1D-C2D	-2.40	1.40	1.45
5	a	814	BCL	C3B-CAB	2.40	1.54	1.46
13	a	803	G2O	C1A-CHA	2.39	1.42	1.37
5	A	805	BCL	C1D-C2D	-2.39	1.40	1.45
5	a	811	BCL	C1D-C2D	-2.39	1.40	1.45
5	A	815	BCL	O1A-CGA	-2.39	1.14	1.22
5	a	809	BCL	O1A-CGA	-2.39	1.14	1.22
5	A	807	BCL	O1A-CGA	-2.39	1.14	1.22
5	A	807	BCL	OBB-CAB	-2.39	1.17	1.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	a	815	BCL	C3B-CAB	2.39	1.54	1.46
5	a	806	BCL	OBB-CAB	-2.39	1.17	1.23
5	E	373	BCL	C3B-CAB	2.39	1.54	1.46
5	G	374	BCL	C1D-C2D	-2.39	1.40	1.45
5	E	377	BCL	C1B-C2B	-2.39	1.37	1.43
5	A	810	BCL	OBB-CAB	-2.39	1.17	1.23
5	a	804	BCL	O1A-CGA	-2.39	1.14	1.22
5	F	378[B]	BCL	C1D-C2D	-2.39	1.40	1.45
5	E	371	BCL	C3B-CAB	2.39	1.54	1.46
5	G	377	BCL	C3B-CAB	2.39	1.54	1.46
5	A	815	BCL	C3B-CAB	2.38	1.54	1.46
5	a	814	BCL	C1D-C2D	-2.38	1.40	1.45
5	A	804	BCL	O1A-CGA	-2.38	1.14	1.22
5	F	371	BCL	C3B-CAB	2.38	1.54	1.46
5	A	814	BCL	OBB-CAB	-2.38	1.17	1.23
5	a	807	BCL	OBB-CAB	-2.38	1.17	1.23
5	A	811	BCL	C1D-C2D	-2.38	1.40	1.45
5	G	372	BCL	OBB-CAB	-2.38	1.17	1.23
5	a	804	BCL	C1D-C2D	-2.38	1.40	1.45
5	G	371	BCL	C3B-CAB	2.38	1.54	1.46
5	E	374	BCL	C3B-CAB	2.38	1.54	1.46
5	F	377	BCL	C3B-CAB	2.38	1.54	1.46
5	G	374	BCL	C3B-CAB	2.38	1.54	1.46
5	a	804	BCL	C3B-CAB	2.37	1.54	1.46
5	G	378[B]	BCL	C1D-C2D	-2.37	1.40	1.45
7	a	801	GS0	C1B-C2B	-2.37	1.37	1.43
5	G	375	BCL	C1D-C2D	-2.37	1.40	1.45
5	a	815	BCL	O1A-CGA	-2.37	1.14	1.22
5	E	378[B]	BCL	C1D-C2D	-2.37	1.40	1.45
5	a	812	BCL	C1D-C2D	-2.37	1.40	1.45
5	A	815	BCL	C1D-C2D	-2.36	1.40	1.45
5	G	373	BCL	C1B-C2B	-2.36	1.37	1.43
5	F	374	BCL	C1D-C2D	-2.36	1.40	1.45
5	A	809	BCL	O1A-CGA	-2.36	1.14	1.22
5	F	374	BCL	C3B-CAB	2.36	1.54	1.46
5	E	374	BCL	C1D-C2D	-2.36	1.40	1.45
5	A	814	BCL	C1D-C2D	-2.36	1.40	1.45
5	a	814	BCL	O1A-CGA	-2.36	1.14	1.22
8	a	816	F26	C30-C26	-2.36	1.30	1.35
5	A	814	BCL	O1A-CGA	-2.36	1.14	1.22
5	A	811	BCL	C1B-C2B	-2.36	1.37	1.43
5	F	377	BCL	C1B-C2B	-2.35	1.37	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	a	811	BCL	C1B-C2B	-2.35	1.37	1.43
5	A	805	BCL	C1B-C2B	-2.35	1.37	1.43
5	E	375	BCL	C3B-CAB	2.35	1.54	1.46
5	A	809	BCL	C3B-CAB	2.35	1.54	1.46
5	a	814	BCL	OBB-CAB	-2.35	1.17	1.23
5	A	804	BCL	C3D-C4D	-2.35	1.38	1.44
5	A	812	BCL	C1B-C2B	-2.35	1.37	1.43
5	A	804	BCL	C3B-CAB	2.35	1.54	1.46
5	A	813	BCL	C3B-CAB	2.35	1.54	1.46
5	A	808	BCL	C3B-CAB	2.35	1.54	1.46
13	a	803	G2O	C1B-C2B	2.34	1.50	1.45
5	G	375	BCL	C3B-CAB	2.34	1.54	1.46
13	a	802	G2O	C1A-CHA	2.34	1.42	1.37
7	A	801	GS0	C1B-C2B	-2.34	1.37	1.43
5	F	375	BCL	C3B-CAB	2.34	1.54	1.46
5	E	375	BCL	C1B-C2B	-2.34	1.37	1.43
5	F	373	BCL	C1B-C2B	-2.34	1.37	1.43
5	a	808	BCL	C3B-CAB	2.34	1.54	1.46
5	a	809	BCL	C3B-CAB	2.34	1.54	1.46
5	A	808	BCL	C1D-C2D	-2.33	1.40	1.45
5	E	374	BCL	C1B-C2B	-2.33	1.37	1.43
5	a	813	BCL	C3B-CAB	2.33	1.54	1.46
5	F	375	BCL	C1B-C2B	-2.32	1.37	1.43
5	a	804	BCL	C3D-C4D	-2.32	1.39	1.44
7	a	801	GS0	C3D-C4D	-2.32	1.39	1.44
5	F	376	BCL	C1B-C2B	-2.32	1.37	1.43
5	a	809	BCL	C1B-C2B	-2.32	1.37	1.43
5	E	378[B]	BCL	C3D-C4D	-2.32	1.39	1.44
5	a	808	BCL	C1D-C2D	-2.32	1.40	1.45
5	a	805	BCL	C1B-C2B	-2.31	1.37	1.43
5	E	376	BCL	C1B-C2B	-2.31	1.37	1.43
5	A	809	BCL	C1B-C2B	-2.31	1.37	1.43
7	a	801	GS0	C3B-CAB	2.31	1.54	1.46
5	G	377	BCL	C1B-C2B	-2.31	1.37	1.43
5	a	807	BCL	C1B-C2B	-2.31	1.37	1.43
5	E	373	BCL	C1B-C2B	-2.31	1.37	1.43
7	A	801	GS0	C3B-CAB	2.31	1.54	1.46
5	A	807	BCL	C1B-C2B	-2.30	1.37	1.43
5	E	378[B]	BCL	C1B-C2B	-2.30	1.37	1.43
5	G	371	BCL	C1B-C2B	-2.30	1.37	1.43
5	F	374	BCL	C1B-C2B	-2.30	1.37	1.43
5	G	375	BCL	C1B-C2B	-2.30	1.37	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	a	812	BCL	C1B-C2B	-2.30	1.37	1.43
5	F	371	BCL	C1B-C2B	-2.30	1.37	1.43
5	a	815	BCL	C1B-C2B	-2.30	1.37	1.43
5	F	378[B]	BCL	C3D-C4D	-2.30	1.39	1.44
5	G	378[B]	BCL	C1B-C2B	-2.30	1.37	1.43
5	A	808	BCL	C3D-C4D	-2.30	1.39	1.44
5	A	815	BCL	C1B-C2B	-2.29	1.37	1.43
8	a	816	F26	C37-C34	-2.29	1.30	1.35
5	G	374	BCL	C1B-C2B	-2.29	1.37	1.43
7	A	801	GS0	C3D-C4D	-2.29	1.39	1.44
5	G	376	BCL	C1B-C2B	-2.29	1.37	1.43
5	a	815	BCL	C3D-C4D	-2.29	1.39	1.44
5	G	378[B]	BCL	C3D-C4D	-2.29	1.39	1.44
5	a	814	BCL	C1B-C2B	-2.28	1.37	1.43
13	A	802	G2O	C1A-CHA	2.28	1.42	1.37
5	a	808	BCL	C3D-C4D	-2.28	1.39	1.44
5	E	371	BCL	C1B-C2B	-2.28	1.37	1.43
5	G	377	BCL	C3D-C4D	-2.28	1.39	1.44
5	E	372	BCL	C3D-C4D	-2.27	1.39	1.44
5	E	374	BCL	C3D-C4D	-2.27	1.39	1.44
8	A	816	F26	C37-C34	-2.27	1.30	1.35
10	a	818	LHG	O7-C5	-2.27	1.41	1.46
5	G	372	BCL	C3D-C4D	-2.27	1.39	1.44
5	E	372	BCL	C1D-C2D	-2.27	1.40	1.45
5	A	807	BCL	C1D-C2D	-2.26	1.40	1.45
5	F	372	BCL	C3D-C4D	-2.26	1.39	1.44
10	A	818	LHG	O7-C5	-2.26	1.41	1.46
5	F	374	BCL	C3D-C4D	-2.25	1.39	1.44
5	a	807	BCL	C1D-C2D	-2.25	1.40	1.45
5	a	813	BCL	C3D-C4D	-2.25	1.39	1.44
5	G	374	BCL	C3D-C4D	-2.25	1.39	1.44
5	A	814	BCL	C1B-C2B	-2.24	1.37	1.43
5	A	813	BCL	C3D-C4D	-2.24	1.39	1.44
5	A	815	BCL	C3D-C4D	-2.24	1.39	1.44
5	a	811	BCL	C3D-C4D	-2.24	1.39	1.44
5	F	378[B]	BCL	C1B-C2B	-2.24	1.37	1.43
5	F	377	BCL	C3D-C4D	-2.23	1.39	1.44
5	A	811	BCL	C3D-C4D	-2.23	1.39	1.44
5	F	372	BCL	C1B-C2B	-2.23	1.37	1.43
5	E	377	BCL	C3D-C4D	-2.23	1.39	1.44
9	a	817	F39	C30-C29	2.23	1.62	1.51
9	A	817	F39	C30-C29	2.23	1.62	1.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	804	BCL	C3B-C2B	-2.23	1.35	1.42
5	G	372	BCL	C1D-C2D	-2.22	1.40	1.45
5	G	372	BCL	C1B-C2B	-2.22	1.37	1.43
5	a	805	BCL	C3D-C4D	-2.22	1.39	1.44
5	E	373	BCL	C3D-C4D	-2.22	1.39	1.44
5	a	804	BCL	C3B-C2B	-2.21	1.35	1.42
5	A	806	BCL	C3D-C4D	-2.21	1.39	1.44
5	a	809	BCL	C3D-C4D	-2.21	1.39	1.44
5	F	372	BCL	C1D-C2D	-2.21	1.41	1.45
5	A	809	BCL	C3D-C4D	-2.21	1.39	1.44
5	F	375	BCL	C3D-C4D	-2.21	1.39	1.44
5	E	372	BCL	C1B-C2B	-2.21	1.37	1.43
5	a	814	BCL	C3D-C4D	-2.21	1.39	1.44
5	E	375	BCL	C3D-C4D	-2.21	1.39	1.44
13	A	803	G2O	CMD-C2D	-2.20	1.46	1.50
5	A	805	BCL	C3D-C4D	-2.20	1.39	1.44
5	a	806	BCL	C3D-C4D	-2.20	1.39	1.44
5	a	808	BCL	C3B-C2B	-2.20	1.35	1.42
13	A	803	G2O	C1B-NB	-2.20	1.35	1.37
5	E	376	BCL	C3D-C4D	-2.20	1.39	1.44
13	a	803	G2O	CMD-C2D	-2.20	1.46	1.50
9	A	817	F39	C29-C28	2.19	1.62	1.51
5	A	808	BCL	C3B-C2B	-2.19	1.35	1.42
9	a	817	F39	C29-C28	2.19	1.62	1.51
5	G	375	BCL	C3D-C4D	-2.19	1.39	1.44
9	a	817	F39	C26-C24	2.18	1.62	1.51
9	a	817	F39	C33-C31	2.18	1.62	1.51
5	a	812	BCL	C3D-C4D	-2.18	1.39	1.44
9	A	817	F39	C33-C31	2.18	1.62	1.51
9	a	817	F39	C28-C26	2.18	1.62	1.51
9	a	817	F39	C31-C30	2.18	1.62	1.51
9	A	817	F39	C26-C24	2.17	1.62	1.51
5	A	812	BCL	C3D-C4D	-2.17	1.39	1.44
9	A	817	F39	C28-C26	2.17	1.62	1.51
9	A	817	F39	C31-C30	2.17	1.62	1.51
5	G	376	BCL	C3D-C4D	-2.17	1.39	1.44
5	a	807	BCL	C3D-C4D	-2.17	1.39	1.44
5	G	373	BCL	C3D-C4D	-2.17	1.39	1.44
5	a	806	BCL	C3B-C2B	-2.16	1.35	1.42
13	a	802	G2O	CMC-C2C	-2.16	1.46	1.50
13	a	803	G2O	C1B-NB	-2.16	1.35	1.37
13	A	802	G2O	CMC-C2C	-2.15	1.46	1.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	814	BCL	C3D-C4D	-2.15	1.39	1.44
5	A	813	BCL	C3B-C2B	-2.15	1.35	1.42
13	A	803	G2O	CMC-C2C	-2.15	1.46	1.50
5	A	806	BCL	C3B-C2B	-2.15	1.35	1.42
5	F	373	BCL	C3D-C4D	-2.14	1.39	1.44
5	a	813	BCL	C3B-C2B	-2.14	1.35	1.42
13	a	803	G2O	CMC-C2C	-2.14	1.46	1.50
9	A	817	F39	C24-C23	2.14	1.62	1.51
5	a	810	BCL	C3B-C2B	-2.14	1.35	1.42
5	A	810	BCL	C3B-C2B	-2.13	1.35	1.42
5	F	376	BCL	C3D-C4D	-2.13	1.39	1.44
5	A	807	BCL	C3D-C4D	-2.13	1.39	1.44
9	a	817	F39	C24-C23	2.12	1.62	1.51
9	A	817	F39	O6-C15	2.12	1.49	1.45
5	a	806	BCL	C3D-C2D	-2.11	1.33	1.39
9	a	817	F39	O6-C15	2.10	1.49	1.45
5	E	376	BCL	C2C-C3C	-2.10	1.48	1.54
5	a	813	BCL	C2C-C3C	-2.10	1.48	1.54
5	F	373	BCL	C3B-C2B	-2.10	1.35	1.42
5	F	371	BCL	C3D-C4D	-2.10	1.39	1.44
5	A	806	BCL	C3D-C2D	-2.10	1.33	1.39
5	A	815	BCL	C2C-C3C	-2.10	1.48	1.54
13	a	802	G2O	C3B-C4B	2.10	1.49	1.46
5	A	805	BCL	C3D-C2D	-2.09	1.33	1.39
5	A	813	BCL	C2C-C3C	-2.09	1.48	1.54
5	E	373	BCL	C3B-C2B	-2.09	1.35	1.42
5	a	805	BCL	C3D-C2D	-2.09	1.33	1.39
5	a	812	BCL	C3B-C2B	-2.08	1.35	1.42
13	A	803	G2O	C15-C13	-2.08	1.42	1.52
5	G	373	BCL	C3B-C2B	-2.08	1.35	1.42
13	a	802	G2O	CMD-C2D	-2.08	1.46	1.50
5	E	375	BCL	C3D-C2D	-2.08	1.33	1.39
5	G	371	BCL	C3D-C4D	-2.08	1.39	1.44
5	A	805	BCL	C2C-C3C	-2.08	1.48	1.54
13	A	802	G2O	C3B-C4B	2.08	1.49	1.46
5	a	815	BCL	C2C-C3C	-2.07	1.48	1.54
13	a	803	G2O	C15-C13	-2.07	1.42	1.52
5	A	812	BCL	C3B-C2B	-2.07	1.35	1.42
13	A	802	G2O	C1B-NB	-2.07	1.35	1.37
5	F	376	BCL	C2C-C3C	-2.07	1.48	1.54
5	E	371	BCL	C3D-C4D	-2.07	1.39	1.44
5	A	807	BCL	C3B-C2B	-2.07	1.35	1.42

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	a	807	BCL	C3B-C2B	-2.07	1.35	1.42
5	a	809	BCL	C2C-C3C	-2.07	1.48	1.54
5	a	805	BCL	C2C-C3C	-2.06	1.48	1.54
5	a	804	BCL	C3D-C2D	-2.06	1.33	1.39
5	F	378[B]	BCL	C3D-C2D	-2.06	1.33	1.39
5	G	375	BCL	C3D-C2D	-2.06	1.33	1.39
11	a	819	LMG	O7-C8	-2.05	1.41	1.46
5	A	809	BCL	C2C-C3C	-2.05	1.48	1.54
5	F	375	BCL	C3D-C2D	-2.05	1.33	1.39
5	E	372	BCL	C2C-C3C	-2.05	1.48	1.54
5	G	373	BCL	C2C-C3C	-2.04	1.48	1.54
13	a	802	G2O	C1B-NB	-2.04	1.35	1.37
5	G	376	BCL	C2C-C3C	-2.04	1.48	1.54
5	G	378[B]	BCL	C3D-C2D	-2.04	1.33	1.39
5	G	371	BCL	C3D-C2D	-2.04	1.33	1.39
5	F	373	BCL	C2C-C3C	-2.04	1.49	1.54
5	A	810	BCL	C2C-C3C	-2.04	1.49	1.54
5	G	376	BCL	C3D-C2D	-2.03	1.33	1.39
5	G	374	BCL	C2C-C3C	-2.03	1.49	1.54
5	F	376	BCL	C3D-C2D	-2.03	1.33	1.39
5	F	377	BCL	C3D-C2D	-2.03	1.33	1.39
5	F	375	BCL	C2C-C3C	-2.03	1.49	1.54
5	E	378[B]	BCL	C3D-C2D	-2.03	1.33	1.39
5	E	378[B]	BCL	C2C-C3C	-2.03	1.49	1.54
5	E	373	BCL	C2C-C3C	-2.03	1.49	1.54
5	A	815	BCL	C3D-C2D	-2.03	1.33	1.39
11	A	819	LMG	O7-C8	-2.02	1.41	1.46
5	a	815	BCL	C3D-C2D	-2.02	1.33	1.39
5	A	814	BCL	C3D-C2D	-2.02	1.33	1.39
5	E	375	BCL	C2C-C3C	-2.02	1.49	1.54
5	F	371	BCL	C2C-C3C	-2.02	1.49	1.54
13	A	802	G2O	CMD-C2D	-2.02	1.46	1.50
5	F	371	BCL	C3D-C2D	-2.02	1.33	1.39
5	E	376	BCL	C3D-C2D	-2.02	1.33	1.39
5	E	371	BCL	C3D-C2D	-2.02	1.33	1.39
5	F	372	BCL	C2C-C3C	-2.02	1.49	1.54
5	G	372	BCL	C2C-C3C	-2.02	1.49	1.54
11	a	819	LMG	O6-C5	-2.02	1.39	1.44
5	F	378[B]	BCL	C2C-C3C	-2.02	1.49	1.54
5	A	804	BCL	C3D-C2D	-2.02	1.33	1.39
5	G	375	BCL	C2C-C3C	-2.01	1.49	1.54
5	G	373	BCL	C3D-C2D	-2.01	1.33	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	E	377	BCL	C3B-C2B	-2.01	1.35	1.42
5	G	371	BCL	C2C-C3C	-2.01	1.49	1.54
5	a	809	BCL	C3D-C2D	-2.01	1.33	1.39
11	A	819	LMG	O6-C5	-2.01	1.39	1.44
5	E	374	BCL	C2C-C3C	-2.01	1.49	1.54
5	G	377	BCL	C3D-C2D	-2.01	1.33	1.39
5	F	377	BCL	C2C-C3C	-2.01	1.49	1.54
5	E	377	BCL	C2C-C3C	-2.01	1.49	1.54
5	F	374	BCL	C2C-C3C	-2.01	1.49	1.54
7	A	801	GS0	C3D-C2D	-2.01	1.33	1.39
5	G	378[B]	BCL	C2C-C3C	-2.01	1.49	1.54
5	A	811	BCL	C2C-C3C	-2.01	1.49	1.54
5	G	377	BCL	C2C-C3C	-2.00	1.49	1.54
5	A	813	BCL	C3D-C2D	-2.00	1.33	1.39
5	A	814	BCL	C2C-C3C	-2.00	1.49	1.54
5	A	811	BCL	C3D-C2D	-2.00	1.33	1.39
5	a	805	BCL	C3B-C2B	-2.00	1.35	1.42
5	a	804	BCL	C2C-C3C	-2.00	1.49	1.54
5	A	805	BCL	C3B-C2B	-2.00	1.35	1.42
7	a	801	GS0	C3D-C2D	-2.00	1.33	1.39
5	E	375	BCL	C3B-C2B	-2.00	1.35	1.42

All (1369) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
9	a	817	F39	C57-C59-C62	-11.93	110.55	127.28
9	A	817	F39	C57-C59-C62	-11.92	110.56	127.28
9	A	817	F39	C51-C44-C42	-11.38	111.31	127.28
9	a	817	F39	C51-C44-C42	-11.37	111.33	127.28
9	A	817	F39	C40-C39-C37	-11.34	111.37	127.28
9	a	817	F39	C40-C39-C37	-11.34	111.38	127.28
9	a	817	F39	C32-C27-C20	-11.19	111.95	127.69
9	A	817	F39	C32-C27-C20	-11.18	111.96	127.69
9	a	817	F39	C63-C61-C58	-10.90	112.00	127.28
9	A	817	F39	C63-C61-C58	-10.85	112.06	127.28
9	A	817	F39	C65-C62-C59	-8.60	108.88	122.82
9	a	817	F39	C65-C62-C59	-8.57	108.94	122.82
9	a	817	F39	C60-C58-C61	-8.43	109.16	122.82
9	A	817	F39	C60-C58-C61	-8.40	109.21	122.82
9	a	817	F39	C38-C37-C39	-8.31	109.35	122.82
9	A	817	F39	C38-C37-C39	-8.29	109.38	122.82
9	a	817	F39	C43-C42-C44	-8.14	109.63	122.82

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
9	A	817	F39	C43-C42-C44	-8.11	109.67	122.82
13	A	802	G2O	C5-C6-C7	-7.13	108.86	125.05
13	a	802	G2O	C5-C6-C7	-7.12	108.87	125.05
9	A	817	F39	C46-C53-C56	-6.96	112.73	128.50
9	a	817	F39	C46-C53-C56	-6.94	112.76	128.50
9	A	817	F39	C41-C42-C44	-6.76	108.37	119.01
9	a	817	F39	C41-C42-C44	-6.75	108.39	119.01
5	a	805	BCL	C1D-ND-C4D	-6.65	101.65	106.31
5	A	805	BCL	C1D-ND-C4D	-6.58	101.70	106.31
9	a	817	F39	C64-C62-C59	-6.58	108.67	119.01
13	A	803	G2O	C5-C6-C7	-6.57	110.14	125.05
9	A	817	F39	C64-C62-C59	-6.54	108.72	119.01
13	a	803	G2O	C5-C6-C7	-6.54	110.20	125.05
5	E	375	BCL	C1D-ND-C4D	-6.47	101.78	106.31
5	F	375	BCL	C1D-ND-C4D	-6.43	101.80	106.31
5	a	806	BCL	C1D-ND-C4D	-6.36	101.85	106.31
5	A	806	BCL	C1D-ND-C4D	-6.34	101.87	106.31
5	G	375	BCL	C1D-ND-C4D	-6.31	101.89	106.31
5	a	807	BCL	C1D-ND-C4D	-6.22	101.94	106.31
5	F	378[B]	BCL	C1D-ND-C4D	-6.18	101.98	106.31
5	A	807	BCL	C1D-ND-C4D	-6.13	102.01	106.31
5	a	811	BCL	C1D-ND-C4D	-6.12	102.02	106.31
5	E	376	BCL	C1D-ND-C4D	-6.11	102.02	106.31
5	G	377	BCL	C1D-ND-C4D	-6.11	102.02	106.31
5	G	376	BCL	C1D-ND-C4D	-6.10	102.03	106.31
5	A	811	BCL	C1D-ND-C4D	-6.10	102.03	106.31
5	E	378[B]	BCL	C1D-ND-C4D	-6.10	102.03	106.31
5	G	378[B]	BCL	C1D-ND-C4D	-6.09	102.04	106.31
5	F	377	BCL	C1D-ND-C4D	-6.09	102.04	106.31
5	E	377	BCL	C1D-ND-C4D	-6.08	102.05	106.31
5	F	376	BCL	C1D-ND-C4D	-6.06	102.06	106.31
5	a	804	BCL	C1D-ND-C4D	-6.05	102.07	106.31
5	A	804	BCL	C1D-ND-C4D	-6.04	102.07	106.31
5	a	814	BCL	C1D-ND-C4D	-6.04	102.08	106.31
5	A	814	BCL	C1D-ND-C4D	-6.02	102.09	106.31
9	a	817	F39	C65-C62-C64	-5.95	109.00	118.09
5	a	810	BCL	C1D-ND-C4D	-5.95	102.14	106.31
9	a	817	F39	C56-C58-C61	-5.94	109.67	119.01
9	A	817	F39	C56-C58-C61	-5.92	109.69	119.01
5	A	813	BCL	C1D-ND-C4D	-5.90	102.17	106.31
9	A	817	F39	C65-C62-C64	-5.89	109.09	118.09
5	A	810	BCL	C1D-ND-C4D	-5.89	102.18	106.31

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
9	a	817	F39	C35-C37-C39	-5.89	109.75	119.01
9	A	817	F39	C35-C37-C39	-5.88	109.76	119.01
9	A	817	F39	C60-C58-C56	-5.86	109.13	118.09
9	a	817	F39	C60-C58-C56	-5.86	109.13	118.09
9	A	817	F39	C53-C56-C58	-5.86	112.33	126.25
5	G	373	BCL	C1C-NC-C4C	-5.86	104.00	106.68
9	a	817	F39	C53-C56-C58	-5.85	112.36	126.25
5	A	809	BCL	C1D-ND-C4D	-5.85	102.20	106.31
5	A	812	BCL	C1D-ND-C4D	-5.84	102.22	106.31
5	a	809	BCL	C1D-ND-C4D	-5.83	102.22	106.31
5	a	815	BCL	C1D-ND-C4D	-5.81	102.23	106.31
5	F	373	BCL	C1C-NC-C4C	-5.81	104.03	106.68
8	a	816	F26	C27-C24-C19	-5.80	119.14	127.28
5	E	372	BCL	C1D-ND-C4D	-5.80	102.24	106.31
8	A	816	F26	C27-C24-C19	-5.80	119.14	127.28
5	E	373	BCL	C1C-NC-C4C	-5.80	104.03	106.68
5	a	813	BCL	C1D-ND-C4D	-5.79	102.25	106.31
5	G	372	BCL	C1D-ND-C4D	-5.78	102.25	106.31
5	A	808	BCL	C1D-ND-C4D	-5.77	102.27	106.31
5	a	812	BCL	C1D-ND-C4D	-5.76	102.27	106.31
8	a	816	F26	C38-C33-C31	-5.76	119.20	127.28
5	F	372	BCL	C1D-ND-C4D	-5.76	102.27	106.31
5	A	815	BCL	C1D-ND-C4D	-5.75	102.28	106.31
8	A	816	F26	C38-C33-C31	-5.74	119.23	127.28
5	G	371	BCL	O2D-CGD-CBD	5.73	121.24	111.23
5	a	808	BCL	C1D-ND-C4D	-5.72	102.30	106.31
5	E	371	BCL	O2D-CGD-CBD	5.72	121.23	111.23
5	F	371	BCL	O2D-CGD-CBD	5.71	121.21	111.23
5	E	373	BCL	C1D-ND-C4D	-5.67	102.33	106.31
5	G	371	BCL	C1D-ND-C4D	-5.64	102.35	106.31
5	F	371	BCL	C1D-ND-C4D	-5.63	102.36	106.31
5	G	373	BCL	C1D-ND-C4D	-5.63	102.36	106.31
13	A	802	G2O	C1A-NA-C4A	5.61	109.24	106.68
9	A	817	F39	C57-C51-C44	-5.58	112.10	123.52
9	a	817	F39	C57-C51-C44	-5.58	112.10	123.52
5	E	371	BCL	C1D-ND-C4D	-5.57	102.40	106.31
5	F	373	BCL	C1D-ND-C4D	-5.57	102.41	106.31
13	A	803	G2O	C1A-NA-C4A	5.56	109.21	106.68
5	G	374	BCL	C1D-ND-C4D	-5.54	102.43	106.31
5	E	374	BCL	C1D-ND-C4D	-5.52	102.44	106.31
5	F	374	BCL	C1D-ND-C4D	-5.51	102.44	106.31
13	a	803	G2O	C1A-NA-C4A	5.49	109.18	106.68

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
13	a	802	G2O	C1A-NA-C4A	5.45	109.16	106.68
9	A	817	F39	C32-C35-C37	-5.36	111.67	126.36
9	a	817	F39	C32-C35-C37	-5.35	111.68	126.36
7	a	801	GS0	C1D-ND-C4D	-5.34	102.56	106.31
7	A	801	GS0	O2D-CGD-CBD	5.33	120.54	111.23
7	a	801	GS0	O2D-CGD-CBD	5.30	120.49	111.23
5	G	373	BCL	O2D-CGD-CBD	5.28	120.47	111.23
5	E	373	BCL	O2D-CGD-CBD	5.28	120.45	111.23
5	F	373	BCL	O2D-CGD-CBD	5.27	120.44	111.23
9	a	817	F39	C63-C64-C62	-5.25	111.95	126.36
9	A	817	F39	C63-C64-C62	-5.25	111.97	126.36
9	a	817	F39	C51-C57-C59	-5.25	112.79	123.52
9	A	817	F39	C51-C57-C59	-5.21	112.86	123.52
7	A	801	GS0	C1D-ND-C4D	-5.21	102.66	106.31
5	E	374	BCL	O2D-CGD-CBD	5.11	120.17	111.23
5	F	374	BCL	O2D-CGD-CBD	5.11	120.16	111.23
5	G	374	BCL	O2D-CGD-CBD	5.08	120.12	111.23
5	E	376	BCL	O2D-CGD-CBD	5.06	120.08	111.23
5	G	376	BCL	O2D-CGD-CBD	5.05	120.06	111.23
5	A	810	BCL	O2D-CGD-CBD	5.05	120.05	111.23
5	A	808	BCL	O2D-CGD-CBD	5.03	120.03	111.23
5	a	810	BCL	O2D-CGD-CBD	5.03	120.02	111.23
5	a	811	BCL	O2D-CGD-CBD	5.03	120.02	111.23
5	a	812	BCL	O2D-CGD-CBD	5.01	119.99	111.23
5	F	376	BCL	O2D-CGD-CBD	5.01	119.99	111.23
5	A	811	BCL	O2D-CGD-CBD	5.00	119.98	111.23
5	a	808	BCL	O2D-CGD-CBD	5.00	119.97	111.23
5	a	813	BCL	O2D-CGD-CBD	4.98	119.93	111.23
5	A	812	BCL	O2D-CGD-CBD	4.97	119.93	111.23
5	F	375	BCL	O2D-CGD-CBD	4.96	119.90	111.23
5	E	375	BCL	O2D-CGD-CBD	4.96	119.89	111.23
5	A	814	BCL	O2D-CGD-CBD	4.95	119.88	111.23
5	A	813	BCL	O2D-CGD-CBD	4.95	119.88	111.23
5	G	375	BCL	O2D-CGD-CBD	4.95	119.88	111.23
5	a	814	BCL	O2D-CGD-CBD	4.93	119.86	111.23
5	G	371	BCL	O2D-CGD-O1D	-4.92	114.27	123.85
5	E	371	BCL	O2D-CGD-O1D	-4.92	114.28	123.85
5	F	371	BCL	O2D-CGD-O1D	-4.91	114.29	123.85
5	a	810	BCL	C2D-C1D-ND	4.87	114.94	110.13
5	A	810	BCL	C2D-C1D-ND	4.87	114.94	110.13
5	A	815	BCL	C1C-NC-C4C	-4.84	104.47	106.68
9	A	817	F39	C43-C42-C41	-4.84	110.70	118.09

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	805	BCL	O2D-CGD-CBD	4.84	119.69	111.23
5	a	804	BCL	O2D-CGD-CBD	4.84	119.68	111.23
9	a	817	F39	C43-C42-C41	-4.83	110.70	118.09
5	a	805	BCL	O2D-CGD-CBD	4.81	119.64	111.23
5	A	804	BCL	O2D-CGD-CBD	4.80	119.62	111.23
5	F	372	BCL	O2D-CGD-CBD	4.79	119.60	111.23
5	G	372	BCL	O2D-CGD-CBD	4.78	119.59	111.23
9	A	817	F39	C39-C40-C41	-4.76	109.42	123.20
5	A	815	BCL	O2D-CGD-CBD	4.76	119.54	111.23
9	a	817	F39	C39-C40-C41	-4.75	109.42	123.20
5	E	372	BCL	O2D-CGD-CBD	4.75	119.53	111.23
5	G	377	BCL	O2D-CGD-CBD	4.75	119.53	111.23
5	a	807	BCL	O2D-CGD-CBD	4.74	119.52	111.23
5	a	815	BCL	O2D-CGD-CBD	4.74	119.52	111.23
5	a	809	BCL	O2D-CGD-CBD	4.74	119.51	111.23
5	a	815	BCL	C1C-NC-C4C	-4.73	104.52	106.68
5	E	377	BCL	O2D-CGD-CBD	4.73	119.49	111.23
5	A	807	BCL	O2D-CGD-CBD	4.72	119.47	111.23
5	A	809	BCL	O2D-CGD-CBD	4.71	119.46	111.23
5	A	806	BCL	O2D-CGD-CBD	4.70	119.45	111.23
5	a	806	BCL	O2D-CGD-CBD	4.70	119.45	111.23
5	a	809	BCL	C1C-NC-C4C	-4.70	104.54	106.68
5	F	377	BCL	O2D-CGD-CBD	4.68	119.42	111.23
5	A	809	BCL	C1C-NC-C4C	-4.65	104.56	106.68
9	A	817	F39	C40-C41-C42	-4.63	113.67	126.36
9	a	817	F39	C40-C41-C42	-4.62	113.69	126.36
5	F	371	BCL	C2A-C3A-C4A	-4.62	94.41	101.87
5	G	371	BCL	C2A-C3A-C4A	-4.62	94.41	101.87
5	E	371	BCL	C2A-C3A-C4A	-4.60	94.43	101.87
9	a	817	F39	C38-C37-C35	-4.58	111.09	118.09
9	A	817	F39	C38-C37-C35	-4.56	111.13	118.09
8	A	816	F26	C22-C18-C13	-4.55	121.29	127.69
5	A	813	BCL	C1C-NC-C4C	-4.54	104.61	106.68
8	a	816	F26	C22-C18-C13	-4.50	121.36	127.69
5	G	378[B]	BCL	O2D-CGD-CBD	4.47	119.05	111.23
5	E	378[B]	BCL	O2D-CGD-CBD	4.47	119.04	111.23
5	F	378[B]	BCL	O2D-CGD-CBD	4.47	119.04	111.23
5	a	813	BCL	C1C-NC-C4C	-4.46	104.64	106.68
5	E	378[B]	BCL	CMB-C2B-C1B	-4.44	118.66	125.42
5	G	378[B]	BCL	CMB-C2B-C1B	-4.44	118.66	125.42
5	a	807	BCL	O2D-CGD-O1D	-4.43	115.22	123.85
5	E	373	BCL	O2D-CGD-O1D	-4.43	115.22	123.85

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	807	BCL	O2D-CGD-O1D	-4.43	115.23	123.85
5	A	811	BCL	CMB-C2B-C1B	-4.42	118.68	125.42
5	F	378[B]	BCL	CMB-C2B-C1B	-4.42	118.68	125.42
5	a	811	BCL	CMB-C2B-C1B	-4.42	118.69	125.42
5	a	809	BCL	CMB-C2B-C1B	-4.42	118.70	125.42
5	F	373	BCL	O2D-CGD-O1D	-4.41	115.26	123.85
5	G	373	BCL	O2D-CGD-O1D	-4.41	115.26	123.85
5	A	809	BCL	CMB-C2B-C1B	-4.41	118.71	125.42
5	A	814	BCL	CMB-C2B-C1B	-4.41	118.71	125.42
10	A	818	LHG	O4-P-O5	4.40	132.92	112.44
10	a	818	LHG	O4-P-O5	4.39	132.88	112.44
5	G	377	BCL	CMB-C2B-C1B	-4.39	118.73	125.42
5	E	377	BCL	CMB-C2B-C1B	-4.39	118.74	125.42
5	a	814	BCL	CMB-C2B-C1B	-4.39	118.74	125.42
7	A	801	GS0	O2D-CGD-O1D	-4.37	115.34	123.85
5	G	375	BCL	CMB-C2B-C1B	-4.36	118.77	125.42
5	F	377	BCL	CMB-C2B-C1B	-4.36	118.78	125.42
5	a	815	BCL	CMB-C2B-C1B	-4.35	118.80	125.42
5	E	377	BCL	C1C-NC-C4C	-4.35	104.70	106.68
5	F	375	BCL	CMB-C2B-C1B	-4.34	118.81	125.42
5	E	375	BCL	CMB-C2B-C1B	-4.34	118.81	125.42
5	E	372	BCL	CMB-C2B-C1B	-4.34	118.81	125.42
5	A	815	BCL	CMB-C2B-C1B	-4.33	118.82	125.42
7	a	801	GS0	O2D-CGD-O1D	-4.32	115.44	123.85
5	F	376	BCL	C1C-NC-C4C	-4.32	104.71	106.68
5	F	374	BCL	CMB-C2B-C1B	-4.31	118.85	125.42
5	E	374	BCL	CMB-C2B-C1B	-4.31	118.85	125.42
5	G	374	BCL	CMB-C2B-C1B	-4.31	118.86	125.42
5	a	810	BCL	C1C-NC-C4C	-4.31	104.71	106.68
5	A	810	BCL	C1C-NC-C4C	-4.29	104.72	106.68
5	F	372	BCL	CMB-C2B-C1B	-4.28	118.90	125.42
5	a	805	BCL	CMB-C2B-C1B	-4.28	118.91	125.42
5	G	376	BCL	CMB-C2B-C1B	-4.28	118.91	125.42
8	A	816	F26	C32-C30-C26	-4.27	121.29	127.28
5	G	372	BCL	CMB-C2B-C1B	-4.27	118.92	125.42
5	A	810	BCL	C3D-C2D-C1D	-4.27	100.00	105.83
5	G	371	BCL	CMB-C2B-C1B	-4.27	118.92	125.42
7	A	801	GS0	CMB-C2B-C1B	-4.27	118.92	125.42
5	A	805	BCL	CMB-C2B-C1B	-4.27	118.92	125.42
5	G	377	BCL	C1C-NC-C4C	-4.26	104.73	106.68
5	F	371	BCL	CMB-C2B-C1B	-4.26	118.93	125.42
5	E	376	BCL	CMB-C2B-C1B	-4.26	118.94	125.42

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	a	801	GS0	CMB-C2B-C1B	-4.26	118.94	125.42
8	a	816	F26	C32-C30-C26	-4.25	121.32	127.28
5	a	810	BCL	C3D-C2D-C1D	-4.25	100.03	105.83
5	a	810	BCL	O2D-CGD-O1D	-4.24	115.59	123.85
5	E	371	BCL	CMB-C2B-C1B	-4.24	118.97	125.42
5	F	374	BCL	O2D-CGD-O1D	-4.24	115.60	123.85
5	E	374	BCL	O2D-CGD-O1D	-4.23	115.61	123.85
5	F	376	BCL	CMB-C2B-C1B	-4.22	118.99	125.42
5	E	376	BCL	C1C-NC-C4C	-4.22	104.75	106.68
5	A	808	BCL	O2D-CGD-O1D	-4.21	115.64	123.85
5	A	810	BCL	O2D-CGD-O1D	-4.21	115.66	123.85
5	G	374	BCL	O2D-CGD-O1D	-4.21	115.66	123.85
5	F	377	BCL	C1C-NC-C4C	-4.20	104.77	106.68
5	A	813	BCL	O2D-CGD-O1D	-4.19	115.69	123.85
5	a	808	BCL	O2D-CGD-O1D	-4.19	115.69	123.85
5	G	375	BCL	O2D-CGD-O1D	-4.17	115.73	123.85
5	E	375	BCL	O2D-CGD-O1D	-4.17	115.74	123.85
5	a	813	BCL	O2D-CGD-O1D	-4.17	115.74	123.85
5	F	375	BCL	O2D-CGD-O1D	-4.16	115.74	123.85
5	G	376	BCL	C1C-NC-C4C	-4.16	104.78	106.68
5	A	806	BCL	C2D-C1D-ND	4.15	114.23	110.13
5	A	811	BCL	O2D-CGD-O1D	-4.14	115.79	123.85
5	E	376	BCL	O2D-CGD-O1D	-4.14	115.80	123.85
5	a	811	BCL	O2D-CGD-O1D	-4.14	115.80	123.85
5	G	376	BCL	O2D-CGD-O1D	-4.13	115.81	123.85
5	E	376	BCL	C2D-C1D-ND	4.13	114.21	110.13
5	F	376	BCL	C2D-C1D-ND	4.13	114.21	110.13
5	G	376	BCL	C2D-C1D-ND	4.13	114.21	110.13
5	a	812	BCL	O2D-CGD-O1D	-4.12	115.83	123.85
5	a	806	BCL	C2D-C1D-ND	4.12	114.20	110.13
5	E	377	BCL	O2D-CGD-O1D	-4.12	115.83	123.85
5	a	805	BCL	C2D-C1D-ND	4.12	114.20	110.13
5	a	815	BCL	O2D-CGD-O1D	-4.12	115.84	123.85
5	F	375	BCL	C2D-C1D-ND	4.12	114.20	110.13
9	A	817	F39	C27-C32-C35	-4.11	111.28	123.20
5	F	376	BCL	O2D-CGD-O1D	-4.11	115.84	123.85
5	A	814	BCL	O2D-CGD-O1D	-4.11	115.85	123.85
5	G	377	BCL	O2D-CGD-O1D	-4.11	115.85	123.85
5	A	812	BCL	O2D-CGD-O1D	-4.11	115.86	123.85
9	a	817	F39	C27-C32-C35	-4.10	111.31	123.20
5	A	815	BCL	O2D-CGD-O1D	-4.10	115.86	123.85
5	a	814	BCL	O2D-CGD-O1D	-4.09	115.89	123.85

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	805	BCL	C2D-C1D-ND	4.09	114.17	110.13
5	F	377	BCL	O2D-CGD-O1D	-4.08	115.90	123.85
5	E	375	BCL	C2D-C1D-ND	4.08	114.16	110.13
5	a	804	BCL	O2D-CGD-O1D	-4.08	115.91	123.85
5	A	804	BCL	O2D-CGD-O1D	-4.08	115.91	123.85
5	G	375	BCL	C2D-C1D-ND	4.08	114.16	110.13
5	A	809	BCL	O2D-CGD-O1D	-4.04	115.99	123.85
5	G	371	BCL	C2D-C1D-ND	4.04	114.12	110.13
5	a	805	BCL	C1C-NC-C4C	-4.03	104.84	106.68
5	A	806	BCL	O2D-CGD-O1D	-4.02	116.03	123.85
5	a	809	BCL	O2D-CGD-O1D	-4.02	116.03	123.85
5	F	372	BCL	O2D-CGD-O1D	-4.01	116.04	123.85
5	a	806	BCL	O2D-CGD-O1D	-4.01	116.04	123.85
5	F	371	BCL	C2D-C1D-ND	4.01	114.10	110.13
5	E	371	BCL	C2D-C1D-ND	4.01	114.09	110.13
5	A	805	BCL	C1C-NC-C4C	-4.00	104.86	106.68
5	A	805	BCL	O2D-CGD-O1D	-3.99	116.08	123.85
5	G	372	BCL	O2D-CGD-O1D	-3.99	116.09	123.85
5	a	805	BCL	O2D-CGD-O1D	-3.98	116.10	123.85
5	E	372	BCL	O2D-CGD-O1D	-3.98	116.11	123.85
5	a	810	BCL	CHD-C1D-ND	-3.97	119.22	124.80
5	A	810	BCL	CHD-C1D-ND	-3.96	119.23	124.80
5	F	377	BCL	C2D-C1D-ND	3.96	114.04	110.13
5	E	377	BCL	C2D-C1D-ND	3.95	114.04	110.13
5	a	811	BCL	C2D-C1D-ND	3.93	114.02	110.13
5	E	378[B]	BCL	O2D-CGD-O1D	-3.93	116.20	123.85
5	E	372	BCL	C1C-NC-C4C	-3.93	104.89	106.68
9	a	817	F39	C61-C63-C64	-3.92	111.83	123.20
5	G	377	BCL	C2D-C1D-ND	3.92	114.00	110.13
5	a	807	BCL	CMB-C2B-C1B	-3.92	119.45	125.42
5	a	814	BCL	C2D-C1D-ND	3.91	114.00	110.13
5	G	378[B]	BCL	O2D-CGD-O1D	-3.91	116.23	123.85
9	A	817	F39	C61-C63-C64	-3.91	111.88	123.20
5	A	811	BCL	C2D-C1D-ND	3.89	113.98	110.13
5	G	372	BCL	C1C-NC-C4C	-3.89	104.90	106.68
5	F	378[B]	BCL	O2D-CGD-O1D	-3.89	116.28	123.85
5	A	813	BCL	C2D-C1D-ND	3.88	113.97	110.13
5	F	378[B]	BCL	C2D-C1D-ND	3.87	113.96	110.13
5	A	814	BCL	C2D-C1D-ND	3.87	113.96	110.13
5	E	378[B]	BCL	C2D-C1D-ND	3.87	113.96	110.13
5	A	807	BCL	CMB-C2B-C1B	-3.87	119.53	125.42
5	A	812	BCL	C2D-C1D-ND	3.87	113.95	110.13

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	807	BCL	C1C-NC-C4C	-3.86	104.92	106.68
5	a	812	BCL	C2D-C1D-ND	3.86	113.94	110.13
5	a	813	BCL	C2D-C1D-ND	3.86	113.94	110.13
5	F	373	BCL	C2C-C3C-C4C	-3.84	95.58	101.34
5	A	804	BCL	C2D-C1D-ND	3.83	113.92	110.13
5	G	373	BCL	C2C-C3C-C4C	-3.83	95.60	101.34
5	a	807	BCL	C2D-C1D-ND	3.83	113.92	110.13
5	G	378[B]	BCL	C2D-C1D-ND	3.83	113.92	110.13
5	E	373	BCL	C2C-C3C-C4C	-3.82	95.62	101.34
5	F	372	BCL	C1C-NC-C4C	-3.81	104.94	106.68
5	A	809	BCL	C2D-C1D-ND	3.81	113.89	110.13
5	G	373	BCL	C2D-C1D-ND	3.81	113.89	110.13
5	A	807	BCL	C2D-C1D-ND	3.80	113.89	110.13
5	a	804	BCL	C2D-C1D-ND	3.80	113.88	110.13
5	a	812	BCL	C1C-NC-C4C	-3.80	104.95	106.68
11	A	819	LMG	O6-C1-O1	-3.79	101.08	110.04
11	a	819	LMG	O6-C1-O1	-3.79	101.09	110.04
5	a	807	BCL	OBB-CAB-CBB	-3.79	111.28	119.77
5	A	807	BCL	OBB-CAB-CBB	-3.78	111.30	119.77
5	E	373	BCL	C2D-C1D-ND	3.77	113.85	110.13
5	F	373	BCL	C2D-C1D-ND	3.77	113.85	110.13
5	F	378[B]	BCL	OBB-CAB-CBB	-3.76	111.34	119.77
5	a	809	BCL	C2D-C1D-ND	3.76	113.85	110.13
5	G	378[B]	BCL	OBB-CAB-CBB	-3.76	111.35	119.77
5	a	815	BCL	C2D-C1D-ND	3.76	113.84	110.13
5	E	378[B]	BCL	OBB-CAB-CBB	-3.75	111.37	119.77
5	A	815	BCL	C2D-C1D-ND	3.75	113.83	110.13
5	A	805	BCL	OBB-CAB-CBB	-3.74	111.39	119.77
5	a	805	BCL	OBB-CAB-CBB	-3.74	111.40	119.77
5	A	808	BCL	C2D-C1D-ND	3.74	113.82	110.13
5	G	371	BCL	OBB-CAB-CBB	-3.73	111.42	119.77
5	F	371	BCL	OBB-CAB-CBB	-3.72	111.44	119.77
5	G	376	BCL	OBB-CAB-CBB	-3.71	111.46	119.77
5	F	376	BCL	OBB-CAB-CBB	-3.71	111.47	119.77
5	F	372	BCL	C2D-C1D-ND	3.71	113.79	110.13
5	a	808	BCL	C2D-C1D-ND	3.69	113.78	110.13
5	E	371	BCL	OBB-CAB-CBB	-3.69	111.50	119.77
5	E	376	BCL	OBB-CAB-CBB	-3.69	111.51	119.77
5	A	808	BCL	C4D-CHA-C1A	3.68	125.64	121.24
5	A	812	BCL	C1C-NC-C4C	-3.68	105.00	106.68
5	F	372	BCL	OBB-CAB-CBB	-3.68	111.53	119.77
5	G	372	BCL	OBB-CAB-CBB	-3.68	111.53	119.77

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	a	808	BCL	C4D-CHA-C1A	3.68	125.63	121.24
5	E	378[B]	BCL	C1C-NC-C4C	-3.68	105.00	106.68
5	G	372	BCL	C2D-C1D-ND	3.68	113.77	110.13
5	A	812	BCL	OBB-CAB-CBB	-3.68	111.53	119.77
5	a	815	BCL	OBB-CAB-CBB	-3.67	111.56	119.77
5	E	372	BCL	C2D-C1D-ND	3.66	113.75	110.13
5	G	375	BCL	OBB-CAB-CBB	-3.66	111.57	119.77
5	A	815	BCL	OBB-CAB-CBB	-3.66	111.57	119.77
5	a	810	BCL	OBB-CAB-CBB	-3.66	111.57	119.77
5	A	811	BCL	OBB-CAB-CBB	-3.66	111.57	119.77
5	E	373	BCL	CMB-C2B-C1B	-3.66	119.85	125.42
5	G	373	BCL	OBB-CAB-CBB	-3.66	111.58	119.77
5	a	804	BCL	OBB-CAB-CBB	-3.66	111.58	119.77
5	a	811	BCL	C1C-NC-C4C	-3.65	105.01	106.68
5	A	810	BCL	OBB-CAB-CBB	-3.65	111.59	119.77
5	E	372	BCL	OBB-CAB-CBB	-3.65	111.59	119.77
5	F	373	BCL	OBB-CAB-CBB	-3.65	111.59	119.77
5	a	806	BCL	OBB-CAB-CBB	-3.65	111.59	119.77
5	A	806	BCL	OBB-CAB-CBB	-3.65	111.59	119.77
5	F	375	BCL	OBB-CAB-CBB	-3.65	111.60	119.77
5	A	810	BCL	CMD-C2D-C1D	3.65	131.15	124.73
7	a	801	GS0	C2D-C1D-ND	3.65	113.73	110.13
5	E	375	BCL	OBB-CAB-CBB	-3.65	111.61	119.77
5	a	812	BCL	OBB-CAB-CBB	-3.64	111.62	119.77
5	E	377	BCL	OBB-CAB-CBB	-3.64	111.62	119.77
5	a	811	BCL	OBB-CAB-CBB	-3.64	111.62	119.77
5	A	804	BCL	OBB-CAB-CBB	-3.63	111.63	119.77
5	G	374	BCL	C2D-C1D-ND	3.63	113.72	110.13
5	a	812	BCL	CMB-C2B-C1B	-3.63	119.89	125.42
5	E	373	BCL	OBB-CAB-CBB	-3.63	111.64	119.77
5	a	810	BCL	CMD-C2D-C1D	3.63	131.12	124.73
5	a	814	BCL	OBB-CAB-CBB	-3.63	111.64	119.77
5	G	378[B]	BCL	C1C-NC-C4C	-3.63	105.02	106.68
5	a	807	BCL	C1C-NC-C4C	-3.63	105.02	106.68
5	G	377	BCL	OBB-CAB-CBB	-3.63	111.65	119.77
5	F	373	BCL	CMB-C2B-C1B	-3.63	119.90	125.42
5	G	373	BCL	CMB-C2B-C1B	-3.63	119.90	125.42
5	A	814	BCL	OBB-CAB-CBB	-3.62	111.66	119.77
5	F	377	BCL	OBB-CAB-CBB	-3.62	111.66	119.77
7	A	801	GS0	C2D-C1D-ND	3.62	113.70	110.13
5	F	374	BCL	C2D-C1D-ND	3.61	113.70	110.13
5	E	374	BCL	OBB-CAB-CBB	-3.61	111.68	119.77

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	F	374	BCL	OBB-CAB-CBB	-3.61	111.68	119.77
5	A	812	BCL	CMB-C2B-C1B	-3.61	119.92	125.42
5	E	374	BCL	C2D-C1D-ND	3.61	113.70	110.13
5	G	374	BCL	OBB-CAB-CBB	-3.61	111.69	119.77
5	F	372	BCL	C4D-CHA-C1A	3.60	125.54	121.24
5	G	372	BCL	C4D-CHA-C1A	3.60	125.54	121.24
5	A	809	BCL	OBB-CAB-CBB	-3.60	111.70	119.77
5	a	809	BCL	OBB-CAB-CBB	-3.60	111.72	119.77
5	A	808	BCL	CMB-C2B-C1B	-3.59	119.95	125.42
5	E	372	BCL	C4D-CHA-C1A	3.59	125.53	121.24
7	A	801	GS0	OBB-CAB-CBB	-3.59	111.73	119.77
5	a	808	BCL	CMB-C2B-C1B	-3.59	119.95	125.42
5	a	813	BCL	OBB-CAB-CBB	-3.59	111.73	119.77
5	G	371	BCL	C4A-NA-C1A	-3.58	105.05	106.68
5	a	806	BCL	CMB-C2B-C1B	-3.58	119.97	125.42
5	A	806	BCL	CMB-C2B-C1B	-3.58	119.97	125.42
5	F	378[B]	BCL	C1C-NC-C4C	-3.57	105.05	106.68
7	a	801	GS0	OBB-CAB-CBB	-3.57	111.77	119.77
5	A	813	BCL	OBB-CAB-CBB	-3.57	111.77	119.77
5	A	811	BCL	C1C-NC-C4C	-3.56	105.06	106.68
5	A	810	BCL	CMB-C2B-C1B	-3.55	120.01	125.42
5	a	808	BCL	OBB-CAB-CBB	-3.54	111.84	119.77
5	a	810	BCL	CMB-C2B-C1B	-3.53	120.04	125.42
5	a	809	BCL	C4D-CHA-C1A	3.53	125.45	121.24
5	A	809	BCL	C4D-CHA-C1A	3.52	125.45	121.24
5	A	810	BCL	CHA-C1A-NA	-3.52	118.42	126.39
5	A	808	BCL	OBB-CAB-CBB	-3.52	111.89	119.77
5	a	810	BCL	CHA-C1A-NA	-3.51	118.44	126.39
5	A	804	BCL	CMB-C2B-C1B	-3.51	120.08	125.42
5	a	804	BCL	CMB-C2B-C1B	-3.51	120.08	125.42
5	F	371	BCL	C4A-NA-C1A	-3.50	105.08	106.68
5	a	813	BCL	CMB-C2B-C1B	-3.49	120.10	125.42
5	A	813	BCL	CMB-C2B-C1B	-3.47	120.13	125.42
5	E	371	BCL	C4A-NA-C1A	-3.44	105.11	106.68
5	A	808	BCL	CHA-C1A-NA	-3.40	118.69	126.39
5	a	808	BCL	CHA-C1A-NA	-3.40	118.70	126.39
5	G	373	BCL	C4D-CHA-C1A	3.39	125.28	121.24
5	E	373	BCL	C4D-CHA-C1A	3.39	125.28	121.24
9	a	817	F39	C25-C20-C19	-3.37	109.39	115.23
9	A	817	F39	C25-C20-C19	-3.36	109.39	115.23
5	a	811	BCL	C16-C15-C13	-3.36	104.79	115.97
5	A	811	BCL	C16-C15-C13	-3.36	104.81	115.97

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	F	373	BCL	C4D-CHA-C1A	3.35	125.24	121.24
5	E	376	BCL	C16-C15-C13	-3.34	104.87	115.97
5	F	376	BCL	C16-C15-C13	-3.34	104.88	115.97
5	G	376	BCL	C16-C15-C13	-3.33	104.88	115.97
5	E	371	BCL	C16-C15-C13	-3.33	104.89	115.97
5	A	811	BCL	O2A-CGA-O1A	-3.33	115.31	123.63
5	F	371	BCL	CHA-C1A-NA	-3.32	118.86	126.39
5	G	371	BCL	C16-C15-C13	-3.31	104.95	115.97
8	A	816	F26	C9-C15-C19	-3.31	118.39	126.25
5	a	811	BCL	O2A-CGA-O1A	-3.31	115.34	123.63
5	G	371	BCL	CHA-C1A-NA	-3.31	118.89	126.39
5	E	371	BCL	C1C-NC-C4C	-3.31	105.17	106.68
5	F	371	BCL	C16-C15-C13	-3.31	104.96	115.97
8	a	816	F26	C9-C15-C19	-3.31	118.40	126.25
5	F	372	BCL	CHA-C1A-NA	-3.30	118.92	126.39
5	F	371	BCL	C1C-NC-C4C	-3.30	105.17	106.68
5	G	372	BCL	CHA-C1A-NA	-3.30	118.92	126.39
7	A	801	GS0	CBB-CAB-C3B	3.29	127.34	120.40
5	G	371	BCL	C1C-NC-C4C	-3.29	105.18	106.68
5	E	372	BCL	CHA-C1A-NA	-3.29	118.94	126.39
7	a	801	GS0	CBB-CAB-C3B	3.29	127.33	120.40
5	A	804	BCL	CHA-C1A-NA	-3.29	118.95	126.39
5	E	371	BCL	CHA-C1A-NA	-3.29	118.95	126.39
8	A	816	F26	C39-C37-C34	-3.28	122.68	127.28
5	A	814	BCL	C4D-CHA-C1A	3.28	125.16	121.24
8	a	816	F26	C39-C37-C34	-3.28	122.68	127.28
5	a	804	BCL	CHA-C1A-NA	-3.27	118.98	126.39
13	a	803	G2O	C1-C2-C3	-3.27	120.83	126.20
5	A	812	BCL	C16-C15-C13	-3.27	105.09	115.97
5	a	812	BCL	C16-C15-C13	-3.27	105.09	115.97
5	G	377	BCL	C4D-CHA-C1A	3.27	125.14	121.24
5	a	814	BCL	C4D-CHA-C1A	3.26	125.13	121.24
5	G	377	BCL	C11-C10-C8	-3.26	105.14	115.97
5	F	377	BCL	C11-C10-C8	-3.25	105.15	115.97
5	G	374	BCL	C1C-NC-C4C	-3.24	105.20	106.68
5	F	371	BCL	C4D-CHA-C1A	3.24	125.11	121.24
13	A	803	G2O	C1-C2-C3	-3.24	120.89	126.20
5	A	812	BCL	C4D-CHA-C1A	3.24	125.11	121.24
5	E	377	BCL	C11-C10-C8	-3.23	105.24	115.97
5	F	374	BCL	CHA-C1A-NA	-3.22	119.09	126.39
5	G	374	BCL	CHA-C1A-NA	-3.22	119.09	126.39
5	E	374	BCL	CHA-C1A-NA	-3.22	119.10	126.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	a	801	GS0	CHA-C1A-NA	-3.22	119.11	126.39
5	A	810	BCL	CGD-CBD-CAD	-3.21	100.46	110.85
5	G	371	BCL	C4D-CHA-C1A	3.21	125.07	121.24
5	a	810	BCL	CGD-CBD-CAD	-3.20	100.47	110.85
5	G	376	BCL	CBB-CAB-C3B	3.20	127.15	120.40
7	A	801	GS0	CHA-C1A-NA	-3.20	119.16	126.39
7	a	801	GS0	C16-C15-C13	-3.19	105.35	115.97
5	E	376	BCL	CBB-CAB-C3B	3.19	127.13	120.40
5	E	377	BCL	C4D-CHA-C1A	3.19	125.05	121.24
5	F	377	BCL	C4D-CHA-C1A	3.19	125.05	121.24
5	A	815	BCL	CBB-CAB-C3B	3.19	127.13	120.40
5	a	815	BCL	CBB-CAB-C3B	3.19	127.12	120.40
5	E	375	BCL	C16-C15-C13	-3.19	105.37	115.97
7	A	801	GS0	C16-C15-C13	-3.19	105.37	115.97
5	F	376	BCL	CBB-CAB-C3B	3.18	127.11	120.40
5	A	804	BCL	C4D-CHA-C1A	3.18	125.03	121.24
5	G	377	BCL	CBB-CAB-C3B	3.18	127.10	120.40
5	F	375	BCL	C16-C15-C13	-3.18	105.41	115.97
5	A	805	BCL	CBB-CAB-C3B	3.17	127.09	120.40
5	F	377	BCL	CBB-CAB-C3B	3.17	127.09	120.40
5	G	375	BCL	C16-C15-C13	-3.17	105.44	115.97
5	G	378[B]	BCL	CBB-CAB-C3B	3.17	127.08	120.40
5	a	804	BCL	C4D-CHA-C1A	3.17	125.02	121.24
5	F	374	BCL	C1C-NC-C4C	-3.17	105.23	106.68
5	F	373	BCL	CHA-C1A-NA	-3.17	119.22	126.39
5	E	373	BCL	CHA-C1A-NA	-3.17	119.22	126.39
5	G	373	BCL	CHA-C1A-NA	-3.16	119.23	126.39
5	G	372	BCL	CBB-CAB-C3B	3.16	127.07	120.40
5	a	812	BCL	C4D-CHA-C1A	3.16	125.01	121.24
5	F	378[B]	BCL	CBB-CAB-C3B	3.16	127.06	120.40
5	a	805	BCL	CBB-CAB-C3B	3.16	127.06	120.40
5	E	378[B]	BCL	CBB-CAB-C3B	3.16	127.06	120.40
5	F	374	BCL	O2A-CGA-O1A	-3.16	115.73	123.63
5	E	372	BCL	CBB-CAB-C3B	3.16	127.06	120.40
5	A	804	BCL	CHB-C1B-NB	3.16	128.78	124.05
5	E	377	BCL	CBB-CAB-C3B	3.15	127.05	120.40
13	A	803	G2O	O2D-CGD-O1D	-3.15	117.72	123.85
5	a	812	BCL	C7-C6-C5	-3.14	104.88	113.26
5	E	371	BCL	C4D-CHA-C1A	3.14	124.99	121.24
5	E	374	BCL	CBB-CAB-C3B	3.14	127.03	120.40
5	a	811	BCL	CBB-CAB-C3B	3.14	127.03	120.40
5	A	811	BCL	CBB-CAB-C3B	3.14	127.03	120.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	F	374	BCL	CBB-CAB-C3B	3.14	127.02	120.40
5	F	373	BCL	C7-C6-C5	-3.14	104.89	113.26
5	E	373	BCL	C7-C6-C5	-3.14	104.90	113.26
5	a	806	BCL	C16-C15-C13	-3.14	105.54	115.97
5	F	372	BCL	CBB-CAB-C3B	3.13	127.01	120.40
5	E	374	BCL	O2A-CGA-O1A	-3.13	115.79	123.63
5	A	806	BCL	C16-C15-C13	-3.13	105.55	115.97
13	a	803	G2O	O2D-CGD-O1D	-3.13	117.75	123.85
5	a	804	BCL	CHB-C1B-NB	3.13	128.75	124.05
5	A	806	BCL	O2A-CGA-O1A	-3.13	115.80	123.63
5	G	371	BCL	CBB-CAB-C3B	3.13	126.99	120.40
5	E	374	BCL	C16-C15-C13	-3.13	105.58	115.97
5	a	806	BCL	C1C-NC-C4C	-3.13	105.25	106.68
5	G	374	BCL	CBB-CAB-C3B	3.13	126.99	120.40
5	G	373	BCL	C7-C6-C5	-3.13	104.93	113.26
5	a	806	BCL	O2A-CGA-O1A	-3.12	115.81	123.63
5	G	374	BCL	O2A-CGA-O1A	-3.12	115.82	123.63
5	A	809	BCL	C2C-C3C-C4C	-3.12	96.66	101.34
5	G	374	BCL	C16-C15-C13	-3.12	105.59	115.97
5	A	811	BCL	C4D-CHA-C1A	3.12	124.97	121.24
5	A	812	BCL	C7-C6-C5	-3.12	104.95	113.26
5	A	808	BCL	CHB-C1B-NB	3.12	128.72	124.05
5	F	371	BCL	CBB-CAB-C3B	3.11	126.96	120.40
5	A	814	BCL	CBB-CAB-C3B	3.11	126.96	120.40
5	a	814	BCL	CBB-CAB-C3B	3.11	126.95	120.40
5	A	806	BCL	CMC-C2C-C3C	-3.11	101.97	113.98
5	a	806	BCL	CMC-C2C-C3C	-3.11	101.97	113.98
5	F	374	BCL	C16-C15-C13	-3.11	105.64	115.97
5	a	811	BCL	C4D-CHA-C1A	3.10	124.95	121.24
5	F	375	BCL	CBB-CAB-C3B	3.10	126.94	120.40
5	G	375	BCL	CBB-CAB-C3B	3.10	126.94	120.40
5	F	374	BCL	C7-C6-C5	-3.10	105.00	113.26
5	E	371	BCL	CBB-CAB-C3B	3.10	126.93	120.40
5	a	813	BCL	C4D-CHA-C1A	3.09	124.94	121.24
5	G	374	BCL	C7-C6-C5	-3.09	105.02	113.26
5	E	374	BCL	C7-C6-C5	-3.09	105.02	113.26
5	F	371	BCL	C2A-C1A-CHA	3.09	129.23	123.87
5	A	806	BCL	CHB-C1B-NB	3.09	128.68	124.05
5	A	812	BCL	O2A-CGA-O1A	-3.08	115.91	123.63
5	a	812	BCL	O2A-CGA-O1A	-3.08	115.91	123.63
5	A	806	BCL	C1C-NC-C4C	-3.08	105.27	106.68
5	a	808	BCL	CHB-C1B-NB	3.08	128.67	124.05

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	a	809	BCL	C2C-C3C-C4C	-3.08	96.73	101.34
5	A	814	BCL	CHA-C1A-NA	-3.08	119.42	126.39
5	A	813	BCL	C4D-CHA-C1A	3.07	124.91	121.24
5	E	375	BCL	CBB-CAB-C3B	3.07	126.88	120.40
8	a	816	F26	C38-C39-C37	-3.07	117.23	123.52
8	A	816	F26	C38-C39-C37	-3.07	117.24	123.52
5	a	814	BCL	CHA-C1A-NA	-3.06	119.45	126.39
5	E	373	BCL	C11-C10-C8	-3.06	105.78	115.97
7	A	801	GS0	C7-C6-C5	-3.06	105.12	113.26
5	a	806	BCL	CHB-C1B-NB	3.06	128.63	124.05
5	E	376	BCL	C11-C10-C8	-3.05	105.81	115.97
5	G	371	BCL	C2A-C1A-CHA	3.05	129.17	123.87
5	E	376	BCL	O2A-CGA-O1A	-3.05	115.99	123.63
7	a	801	GS0	C7-C6-C5	-3.05	105.13	113.26
5	F	376	BCL	C11-C10-C8	-3.05	105.82	115.97
5	F	373	BCL	O2A-CGA-O1A	-3.04	116.02	123.63
5	G	372	BCL	C7-C6-C5	-3.04	105.16	113.26
5	G	376	BCL	C11-C10-C8	-3.04	105.86	115.97
5	A	809	BCL	CHA-C1A-NA	-3.04	119.52	126.39
5	F	376	BCL	O2A-CGA-O1A	-3.04	116.04	123.63
5	A	806	BCL	C11-C10-C8	-3.04	105.88	115.97
5	F	373	BCL	C11-C10-C8	-3.03	105.89	115.97
5	G	373	BCL	C11-C10-C8	-3.03	105.89	115.97
5	G	376	BCL	C7-C6-C5	-3.03	105.19	113.26
5	E	373	BCL	C16-C15-C13	-3.03	105.90	115.97
5	a	806	BCL	C11-C10-C8	-3.03	105.90	115.97
5	a	806	BCL	C4D-CHA-C1A	3.03	124.86	121.24
5	E	373	BCL	O2A-CGA-O1A	-3.03	116.06	123.63
5	A	806	BCL	C4D-CHA-C1A	3.03	124.85	121.24
5	G	376	BCL	O2A-CGA-O1A	-3.03	116.06	123.63
5	a	809	BCL	CHA-C1A-NA	-3.02	119.54	126.39
5	A	806	BCL	C7-C6-C5	-3.02	105.22	113.26
5	G	372	BCL	C16-C15-C13	-3.02	105.93	115.97
5	F	373	BCL	C16-C15-C13	-3.02	105.93	115.97
5	G	371	BCL	O2A-CGA-O1A	-3.02	116.08	123.63
5	A	814	BCL	C1C-NC-C4C	-3.02	105.30	106.68
5	E	372	BCL	C7-C6-C5	-3.02	105.22	113.26
5	E	371	BCL	C2A-C1A-CHA	3.01	129.10	123.87
5	F	372	BCL	C7-C6-C5	-3.01	105.23	113.26
5	a	806	BCL	C7-C6-C5	-3.01	105.23	113.26
5	F	376	BCL	C7-C6-C5	-3.01	105.23	113.26
5	G	373	BCL	O2A-CGA-O1A	-3.01	116.09	123.63

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	F	376	BCL	C4D-CHA-C1A	3.01	124.84	121.24
5	E	377	BCL	C16-C15-C13	-3.01	105.96	115.97
5	E	374	BCL	C1C-NC-C4C	-3.01	105.31	106.68
5	A	807	BCL	CHB-C1B-NB	3.01	128.56	124.05
5	G	376	BCL	C4D-CHA-C1A	3.01	124.83	121.24
5	E	371	BCL	O2A-CGA-O1A	-3.01	116.10	123.63
5	E	372	BCL	C16-C15-C13	-3.01	105.97	115.97
5	E	376	BCL	C4D-CHA-C1A	3.01	124.83	121.24
5	G	373	BCL	C16-C15-C13	-3.00	105.98	115.97
5	E	376	BCL	C7-C6-C5	-3.00	105.27	113.26
5	a	807	BCL	CHB-C1B-NB	3.00	128.55	124.05
5	F	372	BCL	C16-C15-C13	-3.00	106.00	115.97
5	A	812	BCL	CHA-C1A-NA	-3.00	119.60	126.39
5	G	377	BCL	C16-C15-C13	-3.00	106.00	115.97
5	F	371	BCL	O2A-CGA-O1A	-3.00	116.13	123.63
5	F	377	BCL	C16-C15-C13	-3.00	106.01	115.97
5	A	808	BCL	CHD-C1D-ND	-2.99	120.59	124.80
5	a	806	BCL	CHA-C1A-NA	-2.99	119.62	126.39
5	a	808	BCL	CHD-C1D-ND	-2.98	120.61	124.80
5	a	814	BCL	C1C-NC-C4C	-2.98	105.32	106.68
5	A	806	BCL	CHA-C1A-NA	-2.97	119.66	126.39
5	a	808	BCL	CMA-C3A-C4A	-2.97	103.79	111.77
5	a	812	BCL	CHA-C1A-NA	-2.97	119.67	126.39
5	a	809	BCL	CBB-CAB-C3B	2.97	126.66	120.40
5	A	809	BCL	CBB-CAB-C3B	2.96	126.65	120.40
5	A	805	BCL	CHA-C1A-NA	-2.96	119.68	126.39
5	A	808	BCL	C16-C15-C13	-2.96	106.12	115.97
5	A	808	BCL	CMA-C3A-C4A	-2.96	103.82	111.77
9	a	817	F39	C25-C20-C27	-2.96	109.72	122.50
5	E	372	BCL	C11-C10-C8	-2.95	106.14	115.97
9	A	817	F39	C25-C20-C27	-2.95	109.73	122.50
5	G	372	BCL	CHD-C1D-ND	-2.95	120.65	124.80
5	E	372	BCL	CHD-C1D-ND	-2.95	120.65	124.80
5	a	808	BCL	C16-C15-C13	-2.95	106.17	115.97
5	F	372	BCL	CHD-C1D-ND	-2.95	120.66	124.80
5	a	805	BCL	CHA-C1A-NA	-2.95	119.72	126.39
5	A	805	BCL	C7-C6-C5	-2.94	105.42	113.26
5	E	371	BCL	C7-C6-C5	-2.94	105.43	113.26
5	F	372	BCL	C11-C10-C8	-2.94	106.20	115.97
5	G	372	BCL	C11-C10-C8	-2.94	106.20	115.97
5	a	808	BCL	CMC-C2C-C3C	-2.94	102.63	113.98
5	F	371	BCL	C7-C6-C5	-2.94	105.44	113.26

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	a	805	BCL	C7-C6-C5	-2.93	105.44	113.26
5	A	808	BCL	CMC-C2C-C3C	-2.93	102.65	113.98
5	F	377	BCL	CMC-C2C-C3C	-2.93	102.65	113.98
5	G	378[B]	BCL	CHA-C1A-NA	-2.93	119.76	126.39
5	E	374	BCL	CHD-C1D-ND	-2.93	120.68	124.80
5	a	815	BCL	C4D-CHA-C1A	2.93	124.74	121.24
5	E	377	BCL	CMC-C2C-C3C	-2.93	102.67	113.98
5	G	371	BCL	C7-C6-C5	-2.92	105.48	113.26
5	F	378[B]	BCL	CHA-C1A-NA	-2.92	119.78	126.39
5	F	373	BCL	CHB-C1B-NB	2.92	128.43	124.05
5	E	373	BCL	CHB-C1B-NB	2.92	128.42	124.05
5	G	373	BCL	CHB-C1B-NB	2.91	128.42	124.05
5	G	377	BCL	CMC-C2C-C3C	-2.91	102.72	113.98
5	A	811	BCL	CHA-C1A-NA	-2.91	119.80	126.39
5	E	378[B]	BCL	C2C-C3C-C4C	-2.91	96.98	101.34
5	E	375	BCL	CMC-C2C-C3C	-2.91	102.73	113.98
5	G	374	BCL	CHD-C1D-ND	-2.91	120.71	124.80
5	a	812	BCL	CHB-C1B-NB	2.91	128.41	124.05
5	A	809	BCL	CHB-C1B-NB	2.91	128.41	124.05
5	F	376	BCL	CMC-C2C-C3C	-2.91	102.74	113.98
5	A	809	BCL	CHD-C1D-ND	-2.91	120.71	124.80
5	a	811	BCL	CHA-C1A-NA	-2.91	119.81	126.39
5	G	376	BCL	CMC-C2C-C3C	-2.91	102.75	113.98
5	a	809	BCL	CHB-C1B-NB	2.90	128.41	124.05
5	G	375	BCL	CMC-C2C-C3C	-2.90	102.78	113.98
5	E	376	BCL	CMC-C2C-C3C	-2.90	102.78	113.98
5	F	375	BCL	CMC-C2C-C3C	-2.90	102.79	113.98
5	a	809	BCL	CHD-C1D-ND	-2.89	120.73	124.80
5	F	374	BCL	CHD-C1D-ND	-2.89	120.73	124.80
5	E	378[B]	BCL	CHA-C1A-NA	-2.89	119.84	126.39
5	F	371	BCL	CMC-C2C-C3C	-2.89	102.81	113.98
5	E	371	BCL	CMC-C2C-C3C	-2.89	102.82	113.98
5	a	811	BCL	C7-C6-C5	-2.89	105.57	113.26
5	E	375	BCL	C7-C6-C5	-2.89	105.57	113.26
5	a	810	BCL	O2A-CGA-O1A	-2.88	115.92	123.33
5	G	378[B]	BCL	C2C-C3C-C4C	-2.88	97.02	101.34
5	A	815	BCL	C4D-CHA-C1A	2.88	124.68	121.24
5	F	375	BCL	C7-C6-C5	-2.88	105.59	113.26
5	G	371	BCL	CMC-C2C-C3C	-2.88	102.86	113.98
5	G	375	BCL	C7-C6-C5	-2.88	105.59	113.26
5	F	375	BCL	CHA-C1A-NA	-2.88	119.88	126.39
5	A	814	BCL	CHB-C1B-NB	2.88	128.36	124.05

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	812	BCL	CHB-C1B-NB	2.88	128.36	124.05
10	a	818	LHG	O8-C23-C24	2.87	120.60	111.83
7	a	801	GS0	C11-C10-C8	-2.87	106.42	115.97
5	F	378[B]	BCL	C2C-C3C-C4C	-2.87	97.04	101.34
5	A	806	BCL	C3D-C2D-C1D	-2.87	101.91	105.83
5	A	812	BCL	CHD-C1D-ND	-2.87	120.76	124.80
5	A	811	BCL	C7-C6-C5	-2.87	105.62	113.26
5	E	375	BCL	O2A-CGA-O1A	-2.86	116.46	123.63
10	A	818	LHG	O8-C23-C24	2.86	120.57	111.83
7	A	801	GS0	CHB-C1B-NB	2.86	128.34	124.05
5	G	371	BCL	C3D-C2D-C1D	-2.86	101.92	105.83
5	G	376	BCL	C3D-C2D-C1D	-2.86	101.92	105.83
5	a	815	BCL	CHA-C1A-NA	-2.86	119.91	126.39
5	A	811	BCL	C11-C10-C8	-2.86	106.45	115.97
5	F	375	BCL	C4D-CHA-C1A	2.86	124.66	121.24
5	E	376	BCL	C3D-C2D-C1D	-2.86	101.93	105.83
7	A	801	GS0	C11-C10-C8	-2.86	106.47	115.97
5	a	814	BCL	CMC-C2C-C3C	-2.86	102.93	113.98
5	G	377	BCL	CHA-C1A-NA	-2.86	119.92	126.39
5	a	811	BCL	C11-C10-C8	-2.86	106.47	115.97
5	A	814	BCL	CMC-C2C-C3C	-2.86	102.94	113.98
5	A	810	BCL	O2A-CGA-O1A	-2.86	115.99	123.33
5	A	815	BCL	CHA-C1A-NA	-2.85	119.93	126.39
5	F	371	BCL	C3D-C2D-C1D	-2.85	101.94	105.83
5	F	376	BCL	C3D-C2D-C1D	-2.85	101.94	105.83
7	A	801	GS0	CMC-C2C-C3C	-2.85	102.97	113.98
5	A	804	BCL	CMC-C2C-C3C	-2.85	102.97	113.98
5	F	375	BCL	O2A-CGA-O1A	-2.85	116.50	123.63
5	G	375	BCL	CHA-C1A-NA	-2.85	119.94	126.39
5	a	804	BCL	CMC-C2C-C3C	-2.85	102.97	113.98
5	a	812	BCL	CHD-C1D-ND	-2.85	120.80	124.80
7	a	801	GS0	CMC-C2C-C3C	-2.85	102.98	113.98
7	a	801	GS0	CHB-C1B-NB	2.85	128.32	124.05
5	a	809	BCL	O2A-CGA-O1A	-2.84	116.02	123.33
5	E	375	BCL	C4D-CHA-C1A	2.84	124.64	121.24
5	A	807	BCL	CMC-C2C-C3C	-2.84	102.99	113.98
5	A	815	BCL	CMA-C3A-C4A	-2.84	104.14	111.77
5	E	371	BCL	C3D-C2D-C1D	-2.84	101.96	105.83
5	F	374	BCL	CMC-C2C-C3C	-2.84	103.01	113.98
5	F	375	BCL	C3D-C2D-C1D	-2.84	101.96	105.83
5	a	814	BCL	O2A-CGA-O1A	-2.84	116.03	123.33
5	A	809	BCL	O2A-CGA-O1A	-2.84	116.04	123.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	a	807	BCL	CMC-C2C-C3C	-2.84	103.02	113.98
5	G	378[B]	BCL	O2A-CGA-O1A	-2.83	116.04	123.33
5	a	814	BCL	CHB-C1B-NB	2.83	128.30	124.05
5	F	378[B]	BCL	O2A-CGA-O1A	-2.83	116.05	123.33
5	a	805	BCL	CHB-C1B-NB	2.83	128.30	124.05
5	F	378[B]	BCL	CHB-C1B-NB	2.83	128.30	124.05
5	E	377	BCL	C7-C6-C5	-2.83	105.72	113.26
5	E	374	BCL	CMC-C2C-C3C	-2.83	103.03	113.98
5	G	374	BCL	CMC-C2C-C3C	-2.83	103.03	113.98
5	E	377	BCL	CHA-C1A-NA	-2.83	119.99	126.39
5	E	375	BCL	CHA-C1A-NA	-2.83	119.99	126.39
5	a	815	BCL	CMA-C3A-C4A	-2.83	104.17	111.77
5	a	815	BCL	CMC-C2C-C3C	-2.83	103.05	113.98
5	F	377	BCL	CHA-C1A-NA	-2.83	119.99	126.39
5	F	372	BCL	C3D-C2D-C1D	-2.83	101.97	105.83
5	a	806	BCL	CHD-C1D-ND	-2.82	120.83	124.80
5	A	804	BCL	CHD-C1D-ND	-2.82	120.83	124.80
5	a	806	BCL	C3D-C2D-C1D	-2.82	101.98	105.83
5	G	375	BCL	C3D-C2D-C1D	-2.82	101.98	105.83
5	G	375	BCL	O2A-CGA-O1A	-2.82	116.57	123.63
5	a	808	BCL	C11-C10-C8	-2.82	106.59	115.97
5	A	807	BCL	CHA-C1A-NA	-2.82	120.01	126.39
5	a	807	BCL	CHA-C1A-NA	-2.82	120.01	126.39
5	a	812	BCL	C3D-C2D-C1D	-2.82	101.98	105.83
5	A	815	BCL	CMC-C2C-C3C	-2.82	103.09	113.98
5	F	378[B]	BCL	CGD-CBD-CAD	-2.82	101.72	110.85
5	a	804	BCL	CHD-C1D-ND	-2.82	120.84	124.80
5	A	815	BCL	CGD-CBD-CAD	-2.81	101.73	110.85
5	E	378[B]	BCL	O2A-CGA-O1A	-2.81	116.09	123.33
5	E	375	BCL	C11-C10-C8	-2.81	106.61	115.97
5	F	375	BCL	C11-C10-C8	-2.81	106.62	115.97
5	A	808	BCL	C11-C10-C8	-2.81	106.62	115.97
5	G	375	BCL	C11-C10-C8	-2.81	106.62	115.97
5	a	815	BCL	CGD-CBD-CAD	-2.81	101.75	110.85
5	A	806	BCL	CHD-C1D-ND	-2.81	120.85	124.80
5	G	373	BCL	C3D-C2D-C1D	-2.81	102.00	105.83
5	E	375	BCL	CHD-C1D-ND	-2.81	120.85	124.80
5	E	378[B]	BCL	C3D-C2D-C1D	-2.81	102.00	105.83
5	G	377	BCL	C7-C6-C5	-2.81	105.78	113.26
5	a	811	BCL	CMC-C2C-C3C	-2.80	103.14	113.98
5	A	812	BCL	C3D-C2D-C1D	-2.80	102.00	105.83
13	a	802	G2O	O2D-CGD-O1D	-2.80	118.39	123.85

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	814	BCL	O2A-CGA-O1A	-2.80	116.12	123.33
5	G	375	BCL	C4D-CHA-C1A	2.80	124.59	121.24
5	A	811	BCL	CMC-C2C-C3C	-2.80	103.15	113.98
5	F	373	BCL	C3D-C2D-C1D	-2.80	102.01	105.83
5	F	377	BCL	C7-C6-C5	-2.80	105.80	113.26
5	G	378[B]	BCL	CHB-C1B-NB	2.80	128.25	124.05
5	F	376	BCL	CHA-C1A-NA	-2.80	120.05	126.39
5	G	378[B]	BCL	CGD-CBD-CAD	-2.80	101.79	110.85
5	a	813	BCL	C3B-C4B-NB	-2.80	106.07	109.76
5	A	805	BCL	CHB-C1B-NB	2.80	128.24	124.05
5	F	374	BCL	CHB-C1B-NB	2.79	128.24	124.05
5	a	805	BCL	CMC-C2C-C3C	-2.79	103.18	113.98
5	G	372	BCL	C3D-C2D-C1D	-2.79	102.02	105.83
5	a	813	BCL	C3D-C2D-C1D	-2.79	102.02	105.83
7	a	801	GS0	O2A-CGA-O1A	-2.79	116.65	123.63
5	E	378[B]	BCL	CGD-CBD-CAD	-2.79	101.81	110.85
7	A	801	GS0	O2A-CGA-O1A	-2.79	116.65	123.63
5	G	378[B]	BCL	CMC-C2C-C3C	-2.79	103.20	113.98
5	a	810	BCL	CMC-C2C-C3C	-2.79	103.20	113.98
5	A	805	BCL	CMC-C2C-C3C	-2.79	103.21	113.98
5	G	376	BCL	CHA-C1A-NA	-2.78	120.09	126.39
5	A	810	BCL	CMC-C2C-C3C	-2.78	103.23	113.98
5	E	373	BCL	C3D-C2D-C1D	-2.78	102.04	105.83
5	A	813	BCL	C3D-C2D-C1D	-2.78	102.04	105.83
5	F	378[B]	BCL	CMC-C2C-C3C	-2.78	103.24	113.98
5	G	375	BCL	CHB-C1B-NB	2.78	128.22	124.05
5	F	375	BCL	CHD-C1D-ND	-2.78	120.89	124.80
5	A	813	BCL	C3B-C4B-NB	-2.78	106.09	109.76
5	A	808	BCL	C7-C6-C5	-2.78	105.86	113.26
5	E	376	BCL	CHA-C1A-NA	-2.78	120.11	126.39
5	E	377	BCL	C3D-C2D-C1D	-2.78	102.04	105.83
5	E	378[B]	BCL	CHB-C1B-NB	2.77	128.21	124.05
5	a	815	BCL	CHB-C1B-NB	2.77	128.21	124.05
5	a	814	BCL	C3D-C2D-C1D	-2.77	102.05	105.83
7	A	801	GS0	C3D-C2D-C1D	-2.77	102.05	105.83
5	a	814	BCL	CHD-C1D-ND	-2.77	120.90	124.80
5	A	808	BCL	O2A-CGA-O1A	-2.77	116.69	123.63
5	a	808	BCL	O2A-CGA-O1A	-2.77	116.69	123.63
5	a	805	BCL	C4D-CHA-C1A	2.77	124.55	121.24
5	A	815	BCL	CHB-C1B-NB	2.77	128.21	124.05
5	a	808	BCL	C7-C6-C5	-2.77	105.88	113.26
5	E	372	BCL	C3D-C2D-C1D	-2.77	102.05	105.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	812	BCL	CMC-C2C-C3C	-2.77	103.28	113.98
5	E	378[B]	BCL	CMC-C2C-C3C	-2.77	103.28	113.98
5	E	375	BCL	C3D-C2D-C1D	-2.77	102.05	105.83
5	F	377	BCL	C3D-C2D-C1D	-2.77	102.05	105.83
5	G	374	BCL	C11-C10-C8	-2.77	106.77	115.97
5	a	812	BCL	CMC-C2C-C3C	-2.77	103.29	113.98
5	a	811	BCL	CHD-C1D-ND	-2.77	120.91	124.80
5	F	377	BCL	O2A-CGA-O1A	-2.76	116.72	123.63
7	a	801	GS0	C3D-C2D-C1D	-2.76	102.06	105.83
5	E	374	BCL	C11-C10-C8	-2.76	106.78	115.97
8	A	816	F26	C27-C28-C31	-2.76	118.79	126.36
5	a	815	BCL	C3D-C2D-C1D	-2.76	102.06	105.83
5	G	375	BCL	CHD-C1D-ND	-2.76	120.92	124.80
5	a	813	BCL	CHA-C1A-NA	-2.76	120.15	126.39
13	A	802	G2O	O2D-CGD-O1D	-2.76	118.48	123.85
5	F	378[B]	BCL	C3D-C2D-C1D	-2.76	102.07	105.83
5	G	374	BCL	CHB-C1B-NB	2.76	128.18	124.05
5	F	375	BCL	CHB-C1B-NB	2.75	128.18	124.05
5	A	814	BCL	CMA-C3A-C4A	-2.75	104.37	111.77
5	a	814	BCL	CMA-C3A-C4A	-2.75	104.37	111.77
5	G	378[B]	BCL	C3D-C2D-C1D	-2.75	102.07	105.83
5	A	815	BCL	C3D-C2D-C1D	-2.75	102.08	105.83
5	E	375	BCL	CHB-C1B-NB	2.75	128.17	124.05
5	A	814	BCL	CHD-C1D-ND	-2.75	120.93	124.80
5	A	804	BCL	C3D-C2D-C1D	-2.75	102.08	105.83
5	A	811	BCL	CHD-C1D-ND	-2.75	120.94	124.80
5	A	811	BCL	C3D-C2D-C1D	-2.75	102.08	105.83
5	a	811	BCL	C3D-C2D-C1D	-2.75	102.08	105.83
5	G	377	BCL	C3D-C2D-C1D	-2.75	102.08	105.83
5	F	374	BCL	C11-C10-C8	-2.75	106.84	115.97
8	a	816	F26	C27-C28-C31	-2.74	118.84	126.36
5	A	808	BCL	C3D-C2D-C1D	-2.74	102.09	105.83
5	E	377	BCL	O2A-CGA-O1A	-2.74	116.77	123.63
5	G	376	BCL	CHD-C1D-ND	-2.74	120.94	124.80
5	F	376	BCL	CHD-C1D-ND	-2.74	120.94	124.80
5	a	808	BCL	C3D-C2D-C1D	-2.74	102.09	105.83
5	a	815	BCL	C2C-C3C-C4C	-2.74	97.24	101.34
5	G	377	BCL	O2A-CGA-O1A	-2.74	116.78	123.63
5	A	813	BCL	CHA-C1A-NA	-2.74	120.19	126.39
5	A	805	BCL	C4D-CHA-C1A	2.74	124.51	121.24
5	A	805	BCL	O2A-CGA-O1A	-2.73	116.79	123.63
5	E	376	BCL	CHD-C1D-ND	-2.73	120.95	124.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	a	805	BCL	O2A-CGA-O1A	-2.73	116.79	123.63
5	A	809	BCL	C3D-C2D-C1D	-2.73	102.10	105.83
5	a	813	BCL	CHB-C1B-NB	2.73	128.15	124.05
5	A	814	BCL	C3D-C2D-C1D	-2.73	102.11	105.83
5	A	805	BCL	C3D-C2D-C1D	-2.72	102.11	105.83
5	G	377	BCL	CHB-C1B-NB	2.72	128.14	124.05
9	A	817	F39	C13-O2-C11	-2.72	114.20	119.78
5	a	804	BCL	C3D-C2D-C1D	-2.72	102.12	105.83
5	A	810	BCL	CMA-C3A-C4A	-2.72	104.46	111.77
5	E	374	BCL	CHB-C1B-NB	2.72	128.13	124.05
5	A	809	BCL	CMC-C2C-C3C	-2.72	103.47	113.98
5	A	815	BCL	C2C-C3C-C4C	-2.72	97.27	101.34
5	F	372	BCL	CMC-C2C-C3C	-2.72	103.47	113.98
5	a	804	BCL	CMA-C3A-C4A	-2.72	104.47	111.77
9	A	817	F39	O6-C21-C22	2.72	120.12	111.83
5	A	813	BCL	CHB-C1B-NB	2.72	128.12	124.05
5	a	810	BCL	CMA-C3A-C4A	-2.72	104.47	111.77
9	a	817	F39	O6-C21-C22	2.71	120.11	111.83
5	G	372	BCL	CMC-C2C-C3C	-2.71	103.50	113.98
5	A	804	BCL	CMA-C3A-C4A	-2.71	104.49	111.77
5	F	372	BCL	O2A-CGA-O1A	-2.71	116.85	123.63
5	E	372	BCL	CHB-C1B-NB	2.71	128.11	124.05
5	a	809	BCL	CMC-C2C-C3C	-2.70	103.53	113.98
5	E	372	BCL	CMC-C2C-C3C	-2.70	103.53	113.98
9	a	817	F39	C13-O2-C11	-2.70	114.25	119.78
5	a	809	BCL	C3D-C2D-C1D	-2.70	102.15	105.83
5	F	377	BCL	CHB-C1B-NB	2.70	128.10	124.05
5	a	805	BCL	C3D-C2D-C1D	-2.69	102.16	105.83
5	a	805	BCL	CMA-C3A-C4A	-2.69	104.53	111.77
5	G	377	BCL	CHD-C1D-ND	-2.69	121.02	124.80
5	a	806	BCL	CMA-C3A-C4A	-2.69	104.54	111.77
5	F	377	BCL	CHD-C1D-ND	-2.69	121.02	124.80
5	A	806	BCL	CMA-C3A-C4A	-2.69	104.55	111.77
5	G	372	BCL	CHB-C1B-NB	2.69	128.08	124.05
5	G	372	BCL	O2A-CGA-O1A	-2.68	116.91	123.63
5	E	377	BCL	CHD-C1D-ND	-2.68	121.03	124.80
5	A	815	BCL	O2A-CGA-O1A	-2.68	116.44	123.33
5	E	372	BCL	O2A-CGA-O1A	-2.68	116.93	123.63
5	F	374	BCL	C3D-C2D-C1D	-2.68	102.18	105.83
5	E	377	BCL	CHB-C1B-NB	2.68	128.06	124.05
5	A	805	BCL	CMA-C3A-C4A	-2.67	104.59	111.77
5	a	815	BCL	CHD-C1D-ND	-2.67	121.05	124.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	813	BCL	CMC-C2C-C3C	-2.67	103.67	113.98
5	a	807	BCL	C3D-C2D-C1D	-2.67	102.19	105.83
5	E	374	BCL	C3D-C2D-C1D	-2.67	102.19	105.83
5	A	804	BCL	O2A-CGA-O1A	-2.67	116.47	123.33
5	A	815	BCL	CHD-C1D-ND	-2.66	121.05	124.80
5	a	815	BCL	O2A-CGA-O1A	-2.66	116.48	123.33
5	F	372	BCL	CHB-C1B-NB	2.66	128.04	124.05
5	G	374	BCL	C3D-C2D-C1D	-2.66	102.20	105.83
5	G	375	BCL	CMA-C3A-C4A	-2.66	104.63	111.77
5	a	808	BCL	C3B-C4B-NB	-2.66	106.25	109.76
5	F	373	BCL	CMA-C3A-C4A	-2.66	104.63	111.77
5	E	373	BCL	CMA-C3A-C4A	-2.66	104.63	111.77
5	a	813	BCL	CMC-C2C-C3C	-2.66	103.71	113.98
5	a	804	BCL	O2A-CGA-O1A	-2.65	116.51	123.33
5	A	811	BCL	CMA-C3A-C4A	-2.65	104.65	111.77
5	a	810	BCL	C2C-C3C-C4C	-2.65	97.37	101.34
5	E	376	BCL	CMA-C3A-C4A	-2.65	104.65	111.77
5	G	376	BCL	CMA-C3A-C4A	-2.65	104.65	111.77
5	G	372	BCL	C11-C12-C13	-2.65	107.16	115.97
5	E	375	BCL	CMA-C3A-C4A	-2.65	104.65	111.77
5	a	811	BCL	CMA-C3A-C4A	-2.65	104.66	111.77
5	A	808	BCL	C3B-C4B-NB	-2.65	106.27	109.76
5	F	376	BCL	CMA-C3A-C4A	-2.65	104.66	111.77
5	F	375	BCL	CMA-C3A-C4A	-2.64	104.67	111.77
5	E	372	BCL	C11-C12-C13	-2.64	107.17	115.97
5	F	371	BCL	C11-C10-C8	-2.64	107.19	115.97
5	a	807	BCL	O2A-CGA-O1A	-2.64	116.55	123.33
5	A	807	BCL	C3D-C2D-C1D	-2.64	102.23	105.83
5	E	371	BCL	C11-C10-C8	-2.64	107.20	115.97
7	a	801	GS0	CHD-C1D-ND	-2.64	121.09	124.80
5	G	373	BCL	CMA-C3A-C4A	-2.64	104.69	111.77
5	a	808	BCL	C11-C12-C13	-2.63	107.21	115.97
5	G	371	BCL	C11-C10-C8	-2.63	107.22	115.97
5	F	372	BCL	C11-C12-C13	-2.63	107.22	115.97
5	A	807	BCL	O2A-CGA-O1A	-2.63	116.57	123.33
5	a	812	BCL	CBB-CAB-C3B	2.63	125.94	120.40
5	A	812	BCL	CBB-CAB-C3B	2.63	125.94	120.40
5	E	377	BCL	CMA-C3A-C4A	-2.62	104.72	111.77
5	F	378[B]	BCL	CHD-C1D-ND	-2.61	121.12	124.80
5	a	805	BCL	CHD-C1D-ND	-2.61	121.12	124.80
5	A	808	BCL	C11-C12-C13	-2.61	107.28	115.97
5	F	377	BCL	CMA-C3A-C4A	-2.61	104.76	111.77

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	813	BCL	CHD-C1D-ND	-2.61	121.13	124.80
5	G	377	BCL	CMA-C3A-C4A	-2.61	104.77	111.77
5	A	810	BCL	C2C-C3C-C4C	-2.60	97.44	101.34
5	A	805	BCL	CHD-C1D-ND	-2.60	121.14	124.80
5	E	372	BCL	CMA-C3A-C4A	-2.60	104.79	111.77
5	A	813	BCL	O2A-CGA-O1A	-2.60	116.66	123.33
5	a	806	BCL	C11-C12-C13	-2.59	107.34	115.97
7	A	801	GS0	CHD-C1D-ND	-2.59	121.15	124.80
5	A	809	BCL	C4A-NA-C1A	-2.59	105.50	106.68
5	G	378[B]	BCL	CHD-C1D-ND	-2.59	121.16	124.80
5	A	806	BCL	C11-C12-C13	-2.59	107.36	115.97
5	A	804	BCL	C3B-C4B-NB	-2.59	106.34	109.76
5	G	372	BCL	CMA-C3A-C4A	-2.58	104.83	111.77
5	A	805	BCL	C4A-NA-C1A	-2.58	105.50	106.68
5	E	378[B]	BCL	CHD-C1D-ND	-2.58	121.17	124.80
5	E	375	BCL	C1C-NC-C4C	-2.58	105.50	106.68
5	F	372	BCL	CMA-C3A-C4A	-2.58	104.85	111.77
5	a	807	BCL	CBB-CAB-C3B	2.57	125.83	120.40
5	F	373	BCL	C3B-C4B-NB	-2.57	106.37	109.76
5	F	373	BCL	C2A-C1A-CHA	2.57	128.33	123.87
5	a	813	BCL	CHD-C1D-ND	-2.57	121.19	124.80
5	G	376	BCL	CHB-C1B-NB	2.57	127.91	124.05
5	G	373	BCL	C3B-C4B-NB	-2.57	106.37	109.76
5	a	811	BCL	CHB-C1B-NB	2.57	127.90	124.05
5	F	378[B]	BCL	CMA-C3A-C4A	-2.57	104.88	111.77
5	a	813	BCL	CMA-C3A-C4A	-2.57	104.88	111.77
5	A	811	BCL	C11-C12-C13	-2.56	107.44	115.97
5	a	813	BCL	O2A-CGA-O1A	-2.56	116.74	123.33
5	A	807	BCL	CBB-CAB-C3B	2.56	125.81	120.40
5	E	372	BCL	CMD-C2D-C1D	2.56	129.24	124.73
5	G	378[B]	BCL	CMA-C3A-C4A	-2.56	104.89	111.77
5	F	372	BCL	CMD-C2D-C1D	2.56	129.24	124.73
5	a	804	BCL	C3B-C4B-NB	-2.56	106.38	109.76
5	a	811	BCL	C11-C12-C13	-2.56	107.46	115.97
5	G	373	BCL	C2A-C1A-CHA	2.56	128.31	123.87
5	E	376	BCL	CHB-C1B-NB	2.56	127.88	124.05
5	E	373	BCL	C3B-C4B-NB	-2.55	106.39	109.76
5	G	372	BCL	CMD-C2D-C1D	2.55	129.22	124.73
5	A	806	BCL	CMD-C2D-C1D	2.55	129.22	124.73
5	A	811	BCL	CHB-C1B-NB	2.55	127.88	124.05
5	E	378[B]	BCL	CMA-C3A-C4A	-2.55	104.92	111.77
5	A	813	BCL	CMA-C3A-C4A	-2.55	104.93	111.77

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	G	375	BCL	C1C-NC-C4C	-2.54	105.52	106.68
5	a	810	BCL	CHB-C1B-NB	2.54	127.86	124.05
5	a	809	BCL	CMA-C3A-C4A	-2.54	104.94	111.77
5	F	376	BCL	CHB-C1B-NB	2.54	127.86	124.05
5	A	814	BCL	C4A-NA-C1A	-2.54	105.52	106.68
5	A	809	BCL	CMA-C3A-C4A	-2.54	104.95	111.77
5	E	373	BCL	CGD-CBD-CAD	-2.54	102.63	110.85
5	G	373	BCL	CGD-CBD-CAD	-2.53	102.64	110.85
5	F	373	BCL	CGD-CBD-CAD	-2.53	102.64	110.85
5	E	373	BCL	C2A-C1A-CHA	2.53	128.26	123.87
5	a	806	BCL	CMD-C2D-C1D	2.53	129.19	124.73
5	a	808	BCL	C3C-C2C-C1C	2.53	105.96	101.87
5	G	371	BCL	CHD-C1D-ND	-2.53	121.25	124.80
5	F	371	BCL	CHB-C1B-NB	2.53	127.84	124.05
5	A	810	BCL	CHB-C1B-NB	2.53	127.84	124.05
5	A	808	BCL	C3C-C2C-C1C	2.52	105.95	101.87
5	E	371	BCL	CGD-CBD-CAD	-2.52	102.67	110.85
5	a	812	BCL	C3B-C4B-NB	-2.52	106.43	109.76
5	a	804	BCL	C2A-C3A-C4A	-2.52	97.80	101.87
5	F	371	BCL	CGD-CBD-CAD	-2.52	102.69	110.85
5	E	371	BCL	CHB-C1B-NB	2.52	127.83	124.05
5	E	373	BCL	CHD-C1D-ND	-2.51	121.27	124.80
5	A	812	BCL	C3B-C4B-NB	-2.51	106.44	109.76
5	G	373	BCL	CHD-C1D-ND	-2.51	121.27	124.80
5	A	804	BCL	C2A-C3A-C4A	-2.51	97.81	101.87
5	a	809	BCL	C4A-NA-C1A	-2.51	105.53	106.68
5	F	375	BCL	C1C-NC-C4C	-2.51	105.53	106.68
5	G	374	BCL	C4D-CHA-C1A	2.51	124.23	121.24
5	G	371	BCL	CGD-CBD-CAD	-2.50	102.74	110.85
10	a	818	LHG	C11-C10-C9	-2.50	101.73	114.37
5	F	373	BCL	CHD-C1D-ND	-2.50	121.28	124.80
10	A	818	LHG	C11-C10-C9	-2.50	101.75	114.37
5	F	375	BCL	CMD-C2D-C1D	2.50	129.12	124.73
5	F	374	BCL	C4D-CHA-C1A	2.49	124.22	121.24
5	E	371	BCL	CHD-C1D-ND	-2.49	121.30	124.80
5	a	812	BCL	CMA-C3A-C4A	-2.49	105.08	111.77
5	A	810	BCL	CBB-CAB-C3B	2.49	125.65	120.40
5	F	373	BCL	CMC-C2C-C3C	-2.49	104.36	113.98
5	F	371	BCL	CHD-C1D-ND	-2.49	121.30	124.80
5	A	812	BCL	CMA-C3A-C4A	-2.49	105.09	111.77
5	a	814	BCL	C4A-NA-C1A	-2.49	105.55	106.68
5	a	814	BCL	CMD-C2D-C1D	2.48	129.10	124.73

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	G	375	BCL	C11-C12-C13	-2.48	107.71	115.97
5	G	373	BCL	CMC-C2C-C3C	-2.48	104.38	113.98
5	E	375	BCL	C11-C12-C13	-2.48	107.71	115.97
5	E	373	BCL	CMC-C2C-C3C	-2.48	104.38	113.98
5	G	371	BCL	CHB-C1B-NB	2.48	127.77	124.05
5	E	374	BCL	CMA-C3A-C4A	-2.48	105.11	111.77
5	E	378[B]	BCL	CMD-C2D-C1D	2.48	129.09	124.73
5	a	810	BCL	CBB-CAB-C3B	2.48	125.63	120.40
5	F	375	BCL	C11-C12-C13	-2.48	107.73	115.97
5	A	806	BCL	CGD-CBD-CAD	-2.48	102.83	110.85
7	a	801	GS0	CMA-C3A-C4A	-2.48	105.11	111.77
5	E	374	BCL	C4D-CHA-C1A	2.48	124.20	121.24
9	A	817	F39	C19-C20-C27	-2.48	108.65	122.53
5	G	375	BCL	C3C-C2C-C1C	2.48	105.87	101.87
5	E	372	BCL	CGD-CBD-CAD	-2.47	102.83	110.85
5	A	814	BCL	CMD-C2D-C1D	2.47	129.08	124.73
7	A	801	GS0	C11-C12-C13	-2.47	107.75	115.97
9	a	817	F39	C19-C20-C27	-2.47	108.68	122.53
5	F	374	BCL	CMA-C3A-C4A	-2.47	105.14	111.77
5	G	374	BCL	CMA-C3A-C4A	-2.47	105.14	111.77
5	G	375	BCL	CMD-C2D-C1D	2.47	129.07	124.73
7	a	801	GS0	C11-C12-C13	-2.47	107.76	115.97
5	a	806	BCL	CGD-CBD-CAD	-2.46	102.87	110.85
7	A	801	GS0	CMA-C3A-C4A	-2.46	105.15	111.77
5	a	806	BCL	C3C-C2C-C1C	2.46	105.85	101.87
5	G	376	BCL	C11-C12-C13	-2.46	107.78	115.97
5	a	806	BCL	CBB-CAB-C3B	2.46	125.58	120.40
5	G	372	BCL	CGD-CBD-CAD	-2.46	102.89	110.85
5	G	378[B]	BCL	CMD-C2D-C1D	2.46	129.05	124.73
5	A	813	BCL	C2C-C3C-C4C	-2.45	97.67	101.34
5	F	372	BCL	CGD-CBD-CAD	-2.45	102.91	110.85
5	F	373	BCL	CBB-CAB-C3B	2.45	125.57	120.40
5	F	376	BCL	C11-C12-C13	-2.45	107.82	115.97
5	E	375	BCL	C3C-C2C-C1C	2.45	105.83	101.87
5	E	373	BCL	CBB-CAB-C3B	2.45	125.57	120.40
5	G	373	BCL	CBB-CAB-C3B	2.45	125.56	120.40
5	A	806	BCL	CBB-CAB-C3B	2.45	125.56	120.40
5	E	377	BCL	C11-C12-C13	-2.45	107.83	115.97
7	A	801	GS0	C3B-C4B-NB	-2.45	106.53	109.76
5	G	377	BCL	C11-C12-C13	-2.45	107.83	115.97
5	F	375	BCL	C3C-C2C-C1C	2.44	105.82	101.87
5	E	375	BCL	CMD-C2D-C1D	2.44	129.03	124.73

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	F	378[B]	BCL	CMD-C2D-C1D	2.44	129.03	124.73
5	A	806	BCL	C3C-C2C-C1C	2.44	105.81	101.87
5	E	376	BCL	C11-C12-C13	-2.44	107.86	115.97
5	F	377	BCL	C11-C12-C13	-2.43	107.88	115.97
7	a	801	GS0	C3B-C4B-NB	-2.43	106.55	109.76
5	F	371	BCL	C3C-C2C-C1C	2.43	105.80	101.87
5	a	805	BCL	CMD-C2D-C1D	2.43	129.00	124.73
5	G	371	BCL	C3C-C2C-C1C	2.42	105.78	101.87
5	A	809	BCL	CMD-C2D-C1D	2.42	128.99	124.73
5	A	808	BCL	CMD-C2D-C1D	2.41	128.98	124.73
5	a	804	BCL	C1C-NC-C4C	-2.41	105.58	106.68
7	a	801	GS0	C3C-C2C-C1C	2.40	105.75	101.87
5	A	805	BCL	CMD-C2D-C1D	2.40	128.96	124.73
10	A	818	LHG	C20-C19-C18	-2.40	102.24	114.37
5	a	808	BCL	CMD-C2D-C1D	2.40	128.95	124.73
10	a	818	LHG	C20-C19-C18	-2.40	102.26	114.37
5	E	371	BCL	C3C-C2C-C1C	2.39	105.73	101.87
7	A	801	GS0	C3C-C2C-C1C	2.39	105.73	101.87
11	A	819	LMG	O3-C3-C2	-2.39	104.75	110.38
5	a	807	BCL	CHD-C1D-ND	-2.39	121.44	124.80
5	a	806	BCL	C3B-C4B-NB	-2.38	106.61	109.76
5	a	809	BCL	CMD-C2D-C1D	2.38	128.92	124.73
5	A	807	BCL	CHD-C1D-ND	-2.38	121.45	124.80
5	a	813	BCL	C2C-C3C-C4C	-2.38	97.77	101.34
5	a	813	BCL	CMD-C2D-C1D	2.38	128.92	124.73
5	a	805	BCL	C4A-NA-C1A	-2.38	105.59	106.68
5	E	376	BCL	CMD-C2D-C1D	2.38	128.91	124.73
5	G	374	BCL	C11-C12-C13	-2.38	108.06	115.97
5	A	806	BCL	C3B-C4B-NB	-2.38	106.62	109.76
5	A	812	BCL	C12-C11-C10	-2.38	102.63	113.28
11	a	819	LMG	O3-C3-C2	-2.38	104.78	110.38
5	a	812	BCL	C12-C11-C10	-2.38	102.64	113.28
5	a	815	BCL	CMD-C2D-C1D	2.37	128.91	124.73
5	A	804	BCL	CMD-C2D-C1D	2.37	128.90	124.73
5	A	813	BCL	CMD-C2D-C1D	2.37	128.90	124.73
5	E	374	BCL	C11-C12-C13	-2.37	108.09	115.97
5	F	374	BCL	C11-C12-C13	-2.37	108.10	115.97
5	G	376	BCL	C12-C11-C10	-2.36	102.70	113.28
5	G	374	BCL	C3B-C4B-NB	-2.36	106.65	109.76
5	E	376	BCL	C12-C11-C10	-2.35	102.73	113.28
5	A	804	BCL	C1C-NC-C4C	-2.35	105.61	106.68
5	A	811	BCL	CHC-C1C-NC	2.35	127.80	124.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	E	374	BCL	C3B-C4B-NB	-2.35	106.66	109.76
5	F	376	BCL	C12-C11-C10	-2.35	102.76	113.28
5	G	377	BCL	C3C-C2C-C1C	2.35	105.66	101.87
5	A	815	BCL	CMD-C2D-C1D	2.35	128.86	124.73
5	a	812	BCL	C11-C10-C8	-2.34	108.17	115.97
5	A	812	BCL	C11-C10-C8	-2.34	108.17	115.97
5	F	376	BCL	CMD-C2D-C1D	2.34	128.85	124.73
5	G	376	BCL	CMD-C2D-C1D	2.34	128.85	124.73
5	a	804	BCL	CMD-C2D-C1D	2.34	128.85	124.73
5	F	375	BCL	CGD-CBD-CAD	-2.34	103.28	110.85
5	a	811	BCL	CHC-C1C-NC	2.34	127.77	124.40
5	A	804	BCL	C2B-C1B-NB	-2.34	107.90	110.33
5	F	374	BCL	C3B-C4B-NB	-2.34	106.68	109.76
5	F	377	BCL	C3C-C2C-C1C	2.33	105.64	101.87
5	A	807	BCL	C4D-CHA-C1A	2.33	124.03	121.24
5	G	375	BCL	CGD-CBD-CAD	-2.33	103.31	110.85
5	E	375	BCL	CGD-CBD-CAD	-2.33	103.31	110.85
5	E	377	BCL	C3C-C2C-C1C	2.33	105.63	101.87
5	a	811	BCL	C12-C11-C10	-2.32	102.87	113.28
5	A	811	BCL	C12-C11-C10	-2.32	102.88	113.28
5	F	377	BCL	CMD-C2D-C1D	2.32	128.82	124.73
5	E	377	BCL	CMD-C2D-C1D	2.32	128.81	124.73
5	A	810	BCL	C3B-C4B-NB	-2.31	106.70	109.76
5	A	807	BCL	CMD-C2D-C1D	2.31	128.80	124.73
5	a	807	BCL	C4D-CHA-C1A	2.31	124.00	121.24
5	a	811	BCL	CMD-C2D-C1D	2.31	128.80	124.73
5	G	377	BCL	C12-C11-C10	-2.31	102.93	113.28
5	E	373	BCL	C2B-C1B-NB	-2.31	107.93	110.33
5	A	812	BCL	CMD-C2D-C1D	2.31	128.79	124.73
5	a	807	BCL	CMD-C2D-C1D	2.31	128.79	124.73
5	F	377	BCL	C12-C11-C10	-2.31	102.94	113.28
5	A	811	BCL	CMD-C2D-C1D	2.31	128.79	124.73
5	a	812	BCL	CMD-C2D-C1D	2.31	128.79	124.73
5	E	371	BCL	C11-C12-C13	-2.30	108.31	115.97
5	a	804	BCL	C2B-C1B-NB	-2.30	107.94	110.33
5	G	371	BCL	C11-C12-C13	-2.30	108.31	115.97
5	E	377	BCL	C12-C11-C10	-2.30	102.97	113.28
5	F	371	BCL	C11-C12-C13	-2.30	108.32	115.97
5	a	810	BCL	C3B-C4B-NB	-2.30	106.72	109.76
5	G	377	BCL	CMD-C2D-C1D	2.30	128.77	124.73
5	a	812	BCL	C2B-C1B-NB	-2.30	107.95	110.33
5	E	374	BCL	C12-C11-C10	-2.29	103.01	113.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	G	373	BCL	C2B-C1B-NB	-2.29	107.95	110.33
5	G	374	BCL	C12-C11-C10	-2.29	103.03	113.28
5	F	373	BCL	C2B-C1B-NB	-2.29	107.96	110.33
11	A	819	LMG	C6-C5-C4	-2.29	107.41	113.02
5	F	374	BCL	C12-C11-C10	-2.28	103.04	113.28
5	A	807	BCL	C3B-C4B-NB	-2.28	106.75	109.76
11	a	819	LMG	C6-C5-C4	-2.28	107.43	113.02
5	G	376	BCL	CHC-C1C-NC	2.27	127.68	124.40
5	G	373	BCL	CMD-C2D-C1D	2.27	128.73	124.73
5	F	373	BCL	C11-C12-C13	-2.27	108.42	115.97
5	G	377	BCL	C4A-NA-C1A	-2.27	105.64	106.68
5	a	804	BCL	CBB-CAB-C3B	2.27	125.19	120.40
5	A	813	BCL	CBB-CAB-C3B	2.27	125.18	120.40
5	a	807	BCL	CHC-C1C-NC	2.27	127.67	124.40
5	a	807	BCL	C3B-C4B-NB	-2.26	106.77	109.76
5	G	373	BCL	C11-C12-C13	-2.26	108.44	115.97
5	A	812	BCL	C2B-C1B-NB	-2.26	107.98	110.33
7	a	801	GS0	CMD-C2D-C1D	2.26	128.71	124.73
5	E	373	BCL	C11-C12-C13	-2.26	108.45	115.97
5	E	376	BCL	CHC-C1C-NC	2.26	127.66	124.40
5	F	373	BCL	CMD-C2D-C1D	2.26	128.70	124.73
5	A	804	BCL	C4A-NA-C1A	-2.26	105.65	106.68
5	a	813	BCL	C2B-C1B-NB	-2.26	107.98	110.33
5	E	377	BCL	C3B-C4B-NB	-2.26	106.78	109.76
5	A	804	BCL	CBB-CAB-C3B	2.26	125.16	120.40
5	a	813	BCL	CBB-CAB-C3B	2.26	125.16	120.40
7	A	801	GS0	CMD-C2D-C1D	2.26	128.70	124.73
5	G	374	BCL	CMD-C2D-C1D	2.25	128.70	124.73
5	A	804	BCL	C3C-C2C-C1C	2.25	105.51	101.87
13	a	803	G2O	O2A-CGA-O1A	-2.25	117.99	123.63
5	a	804	BCL	C3C-C2C-C1C	2.25	105.51	101.87
13	A	803	G2O	O2A-CGA-O1A	-2.25	117.99	123.63
5	E	374	BCL	CMD-C2D-C1D	2.25	128.69	124.73
13	A	803	G2O	CAA-CBA-CGA	-2.25	106.82	113.21
5	a	804	BCL	CGD-CBD-CAD	-2.25	103.57	110.85
5	E	373	BCL	CMD-C2D-C1D	2.25	128.68	124.73
13	a	803	G2O	CAA-CBA-CGA	-2.24	106.83	113.21
5	F	374	BCL	CGD-CBD-CAD	-2.24	103.58	110.85
8	A	816	F26	C23-C19-C15	2.24	121.51	118.09
5	A	804	BCL	CGD-CBD-CAD	-2.24	103.59	110.85
5	A	813	BCL	CGD-CBD-CAD	-2.24	103.60	110.85
5	a	813	BCL	CGD-CBD-CAD	-2.24	103.60	110.85

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	F	376	BCL	CHC-C1C-NC	2.24	127.63	124.40
5	F	374	BCL	CMD-C2D-C1D	2.24	128.66	124.73
10	A	818	LHG	C18-C17-C16	-2.23	103.08	114.37
10	a	818	LHG	C18-C17-C16	-2.23	103.08	114.37
5	G	372	BCL	C12-C11-C10	-2.23	103.28	113.28
5	G	374	BCL	CGD-CBD-CAD	-2.23	103.63	110.85
5	G	378[B]	BCL	C4D-CHA-C1A	2.23	123.90	121.24
5	A	805	BCL	CHC-C1C-NC	2.23	127.62	124.40
5	F	372	BCL	C12-C11-C10	-2.23	103.30	113.28
5	E	374	BCL	CGD-CBD-CAD	-2.23	103.64	110.85
8	a	816	F26	C2-C9-C15	-2.22	123.48	128.50
5	E	372	BCL	C12-C11-C10	-2.22	103.34	113.28
5	G	377	BCL	C3B-C4B-NB	-2.21	106.84	109.76
8	a	816	F26	C23-C19-C15	2.21	121.47	118.09
5	F	376	BCL	CGD-CBD-CAD	-2.21	103.69	110.85
5	a	815	BCL	C3B-C4B-NB	-2.21	106.84	109.76
5	F	378[B]	BCL	C4D-CHA-C1A	2.21	123.88	121.24
5	A	815	BCL	C3B-C4B-NB	-2.21	106.85	109.76
8	A	816	F26	C2-C9-C15	-2.21	123.50	128.50
5	a	808	BCL	CHC-C1C-NC	2.20	127.58	124.40
5	G	376	BCL	CGD-CBD-CAD	-2.20	103.71	110.85
5	E	376	BCL	CGD-CBD-CAD	-2.20	103.72	110.85
5	A	812	BCL	C2A-C3A-C4A	-2.19	98.33	101.87
5	a	804	BCL	C4A-NA-C1A	-2.19	105.68	106.68
5	E	372	BCL	CHC-C1C-NC	2.19	127.56	124.40
5	A	807	BCL	CHC-C1C-NC	2.19	127.56	124.40
5	a	807	BCL	C3C-C2C-C1C	2.19	105.40	101.87
5	A	811	BCL	C3B-C4B-NB	-2.19	106.87	109.76
5	E	378[B]	BCL	C4D-CHA-C1A	2.19	123.85	121.24
5	a	807	BCL	C2B-C1B-NB	-2.19	108.06	110.33
5	A	807	BCL	C3C-C2C-C1C	2.18	105.40	101.87
5	A	813	BCL	C2B-C1B-NB	-2.18	108.06	110.33
5	G	377	BCL	CMB-C2B-C3B	2.18	132.55	125.67
5	G	372	BCL	CHC-C1C-NC	2.18	127.54	124.40
5	F	377	BCL	C4A-NA-C1A	-2.18	105.69	106.68
5	F	372	BCL	CHC-C1C-NC	2.17	127.54	124.40
5	A	809	BCL	CMB-C2B-C3B	2.17	132.52	125.67
5	a	805	BCL	CHC-C1C-NC	2.17	127.53	124.40
5	A	806	BCL	C12-C11-C10	-2.17	103.55	113.28
7	A	801	GS0	C2B-C1B-NB	-2.17	108.07	110.33
5	G	375	BCL	CMB-C2B-C3B	2.17	132.52	125.67
5	F	377	BCL	C3B-C4B-NB	-2.17	106.90	109.76

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	a	811	BCL	C3B-C4B-NB	-2.17	106.90	109.76
5	E	371	BCL	CMD-C2D-C1D	2.17	128.55	124.73
5	F	371	BCL	CMD-C2D-C1D	2.17	128.55	124.73
5	a	806	BCL	C12-C11-C10	-2.17	103.56	113.28
5	A	806	BCL	C4A-NA-C1A	-2.17	105.69	106.68
5	a	806	BCL	C4A-NA-C1A	-2.17	105.69	106.68
5	F	377	BCL	CMB-C2B-C3B	2.17	132.50	125.67
5	G	371	BCL	CMD-C2D-C1D	2.17	128.54	124.73
5	a	809	BCL	CMB-C2B-C3B	2.16	132.50	125.67
5	A	814	BCL	CMB-C2B-C3B	2.16	132.49	125.67
5	A	814	BCL	C3C-C2C-C1C	2.16	105.36	101.87
5	E	378[B]	BCL	C3B-C4B-NB	-2.16	106.91	109.76
5	E	375	BCL	CMB-C2B-C3B	2.16	132.47	125.67
5	a	811	BCL	CMB-C2B-C3B	2.15	132.47	125.67
5	a	809	BCL	C3B-C4B-NB	-2.15	106.92	109.76
5	F	375	BCL	C2A-C3A-C4A	-2.15	98.39	101.87
5	A	811	BCL	CMB-C2B-C3B	2.15	132.46	125.67
5	a	809	BCL	C2A-C3A-C4A	-2.15	98.39	101.87
5	F	375	BCL	CMB-C2B-C3B	2.15	132.46	125.67
5	E	375	BCL	C2A-C3A-C4A	-2.15	98.39	101.87
5	a	812	BCL	C2A-C3A-C4A	-2.15	98.39	101.87
5	a	814	BCL	C3C-C2C-C1C	2.15	105.34	101.87
5	E	377	BCL	CMB-C2B-C3B	2.15	132.45	125.67
5	A	812	BCL	CHC-C1C-NC	2.15	127.50	124.40
5	A	807	BCL	C2B-C1B-NB	-2.15	108.10	110.33
5	G	375	BCL	C2A-C3A-C4A	-2.15	98.40	101.87
5	A	809	BCL	C3B-C4B-NB	-2.14	106.93	109.76
5	F	372	BCL	C3B-C4B-NB	-2.14	106.93	109.76
5	G	371	BCL	CMB-C2B-C3B	2.14	132.43	125.67
5	A	809	BCL	C2A-C3A-C4A	-2.14	98.41	101.87
5	F	378[B]	BCL	C3B-C4B-NB	-2.14	106.93	109.76
7	a	801	GS0	C2B-C1B-NB	-2.14	108.11	110.33
5	E	372	BCL	C3B-C4B-NB	-2.14	106.94	109.76
5	a	814	BCL	C3B-C4B-NB	-2.14	106.94	109.76
5	F	374	BCL	C2A-C1A-CHA	2.14	127.57	123.87
5	A	808	BCL	CHC-C1C-NC	2.13	127.48	124.40
5	a	814	BCL	CMB-C2B-C3B	2.13	132.39	125.67
5	E	374	BCL	C2B-C1B-NB	-2.13	108.12	110.33
5	F	371	BCL	CMB-C2B-C3B	2.13	132.38	125.67
5	F	374	BCL	C2B-C1B-NB	-2.13	108.12	110.33
5	G	374	BCL	C2A-C1A-CHA	2.13	127.56	123.87
5	a	812	BCL	CHC-C1C-NC	2.12	127.47	124.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	G	372	BCL	C3B-C4B-NB	-2.12	106.96	109.76
5	G	374	BCL	CMB-C2B-C3B	2.12	132.36	125.67
7	a	801	GS0	C4D-CHA-C1A	2.12	123.77	121.24
5	F	378[B]	BCL	CMB-C2B-C3B	2.12	132.36	125.67
5	A	805	BCL	C3C-C2C-C1C	2.12	105.29	101.87
5	A	804	BCL	CHC-C1C-NC	2.12	127.46	124.40
5	F	371	BCL	C3B-C4B-NB	-2.12	106.96	109.76
7	A	801	GS0	C12-C11-C10	-2.12	103.78	113.28
5	E	371	BCL	CMB-C2B-C3B	2.12	132.35	125.67
5	A	806	BCL	C2A-C3A-C4A	-2.12	98.45	101.87
5	G	375	BCL	CHC-C1C-NC	2.12	127.45	124.40
5	F	375	BCL	CHC-C1C-NC	2.12	127.45	124.40
5	G	378[B]	BCL	C3B-C4B-NB	-2.11	106.97	109.76
7	a	801	GS0	C12-C11-C10	-2.11	103.80	113.28
5	G	378[B]	BCL	CMB-C2B-C3B	2.11	132.34	125.67
5	a	808	BCL	CBB-CAB-C3B	2.11	124.86	120.40
5	a	806	BCL	C2A-C3A-C4A	-2.11	98.46	101.87
5	F	374	BCL	CMB-C2B-C3B	2.11	132.33	125.67
5	F	375	BCL	C3B-C4B-NB	-2.11	106.97	109.76
5	F	376	BCL	C3B-C4B-NB	-2.11	106.97	109.76
5	A	808	BCL	CBB-CAB-C3B	2.11	124.85	120.40
5	G	376	BCL	C3B-C4B-NB	-2.11	106.98	109.76
5	E	374	BCL	C2A-C1A-CHA	2.11	127.52	123.87
5	E	378[B]	BCL	CMB-C2B-C3B	2.10	132.31	125.67
5	E	372	BCL	CMB-C2B-C3B	2.10	132.31	125.67
5	E	375	BCL	C3B-C4B-NB	-2.10	106.98	109.76
5	A	815	BCL	CMB-C2B-C3B	2.10	132.31	125.67
5	G	371	BCL	C3B-C4B-NB	-2.10	106.98	109.76
5	E	377	BCL	C4A-NA-C1A	-2.10	105.72	106.68
5	a	815	BCL	CMB-C2B-C3B	2.10	132.30	125.67
5	E	372	BCL	C2B-C1B-NB	-2.10	108.15	110.33
5	E	374	BCL	CMB-C2B-C3B	2.10	132.29	125.67
5	a	805	BCL	CGD-CBD-CAD	-2.10	104.05	110.85
7	A	801	GS0	CMB-C2B-C3B	2.10	132.29	125.67
5	A	805	BCL	CGD-CBD-CAD	-2.10	104.06	110.85
5	F	372	BCL	CMB-C2B-C3B	2.10	132.28	125.67
5	A	814	BCL	C3B-C4B-NB	-2.09	107.00	109.76
5	G	375	BCL	C3B-C4B-NB	-2.09	107.00	109.76
5	G	372	BCL	CMB-C2B-C3B	2.09	132.27	125.67
5	E	375	BCL	CHC-C1C-NC	2.09	127.42	124.40
5	a	805	BCL	C3C-C2C-C1C	2.09	105.24	101.87
5	a	805	BCL	CMB-C2B-C3B	2.09	132.25	125.67

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	a	801	GS0	CMB-C2B-C3B	2.09	132.25	125.67
5	G	376	BCL	CMB-C2B-C3B	2.08	132.24	125.67
5	a	810	BCL	C4D-CHA-C1A	2.08	123.73	121.24
5	A	807	BCL	CMA-C3A-C4A	-2.08	106.18	111.77
5	A	805	BCL	CMB-C2B-C3B	2.08	132.23	125.67
5	G	374	BCL	C2B-C1B-NB	-2.08	108.17	110.33
5	E	373	BCL	C12-C11-C10	-2.08	103.97	113.28
5	A	810	BCL	C4D-CHA-C1A	2.08	123.72	121.24
10	A	818	LHG	C27-C26-C25	-2.08	103.88	114.37
10	a	818	LHG	C27-C26-C25	-2.08	103.88	114.37
5	E	375	BCL	C12-C11-C10	-2.08	103.98	113.28
5	E	376	BCL	CMB-C2B-C3B	2.07	132.21	125.67
5	a	804	BCL	CHC-C1C-NC	2.07	127.39	124.40
5	G	375	BCL	C12-C11-C10	-2.07	103.98	113.28
5	a	808	BCL	C12-C11-C10	-2.07	104.00	113.28
5	F	375	BCL	C12-C11-C10	-2.07	104.00	113.28
7	A	801	GS0	C4D-CHA-C1A	2.07	123.71	121.24
5	E	376	BCL	C3B-C4B-NB	-2.07	107.03	109.76
8	A	816	F26	C40-C34-C35	2.06	121.24	118.09
5	E	371	BCL	C12-C11-C10	-2.06	104.04	113.28
5	F	376	BCL	CMB-C2B-C3B	2.06	132.17	125.67
5	F	373	BCL	C12-C11-C10	-2.06	104.04	113.28
5	G	373	BCL	C12-C11-C10	-2.06	104.05	113.28
5	a	807	BCL	CMA-C3A-C4A	-2.06	106.24	111.77
5	G	371	BCL	C12-C11-C10	-2.06	104.06	113.28
5	F	371	BCL	C12-C11-C10	-2.06	104.07	113.28
5	E	371	BCL	C3B-C4B-NB	-2.06	107.05	109.76
5	A	808	BCL	C12-C11-C10	-2.05	104.08	113.28
5	a	815	BCL	C2B-C1B-NB	-2.05	108.20	110.33
5	A	815	BCL	C2B-C1B-NB	-2.05	108.20	110.33
5	E	374	BCL	CHC-C1C-NC	2.05	127.35	124.40
5	F	378[B]	BCL	C2A-C1A-CHA	2.04	127.41	123.87
5	G	378[B]	BCL	C2A-C1A-CHA	2.04	127.41	123.87
8	a	816	F26	C40-C34-C35	2.04	121.20	118.09
5	F	374	BCL	C2C-C3C-C4C	-2.03	98.29	101.34
8	a	816	F26	C22-C25-C26	-2.03	120.80	126.36
5	E	374	BCL	C2C-C3C-C4C	-2.02	98.31	101.34
5	F	375	BCL	C2A-C1A-CHA	2.02	127.37	123.87
5	G	372	BCL	C2B-C1B-NB	-2.02	108.24	110.33
5	A	813	BCL	CHC-C1C-NC	2.01	127.30	124.40
8	A	816	F26	C22-C25-C26	-2.01	120.85	126.36
5	G	374	BCL	C2C-C3C-C4C	-2.01	98.33	101.34

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	G	375	BCL	C2A-C1A-CHA	2.01	127.35	123.87
5	E	375	BCL	C2A-C1A-CHA	2.00	127.35	123.87
5	E	377	BCL	C6-C5-C3	-2.00	108.59	113.47
11	a	819	LMG	O3-C3-C4	-2.00	105.66	110.38
5	A	813	BCL	C2A-C3A-C4A	-2.00	98.63	101.87
5	a	813	BCL	C2A-C3A-C4A	-2.00	98.64	101.87
5	E	378[B]	BCL	C2A-C1A-CHA	2.00	127.34	123.87
5	A	808	BCL	C2B-C1B-NB	-2.00	108.25	110.33

There are no chirality outliers.

All (750) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	E	371	BCL	C1A-C2A-CAA-CBA
5	E	371	BCL	C3A-C2A-CAA-CBA
5	E	371	BCL	CBD-CGD-O2D-CED
5	E	371	BCL	O1D-CGD-O2D-CED
5	E	373	BCL	C2C-C3C-CAC-CBC
5	E	373	BCL	C4C-C3C-CAC-CBC
5	E	373	BCL	CAD-CBD-CGD-O2D
5	E	373	BCL	C2-C3-C5-C6
5	E	374	BCL	C2C-C3C-CAC-CBC
5	E	374	BCL	C4C-C3C-CAC-CBC
5	E	377	BCL	CHA-CBD-CGD-O1D
5	E	377	BCL	CHA-CBD-CGD-O2D
5	E	378[B]	BCL	C3A-C2A-CAA-CBA
5	E	372	BCL	C2C-C3C-CAC-CBC
5	E	372	BCL	C4C-C3C-CAC-CBC
5	E	372	BCL	C1-C2-C3-C5
5	F	372	BCL	C2C-C3C-CAC-CBC
5	F	372	BCL	C4C-C3C-CAC-CBC
5	F	372	BCL	C1-C2-C3-C5
5	F	371	BCL	C1A-C2A-CAA-CBA
5	F	371	BCL	C3A-C2A-CAA-CBA
5	F	371	BCL	CBD-CGD-O2D-CED
5	F	371	BCL	O1D-CGD-O2D-CED
5	F	373	BCL	C2C-C3C-CAC-CBC
5	F	373	BCL	C4C-C3C-CAC-CBC
5	F	373	BCL	CAD-CBD-CGD-O2D
5	F	373	BCL	C2-C3-C5-C6
5	F	374	BCL	C2C-C3C-CAC-CBC
5	F	374	BCL	C4C-C3C-CAC-CBC

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Mol	Chain	Res	Type	Atoms
5	F	377	BCL	CHA-CBD-CGD-O1D
5	F	377	BCL	CHA-CBD-CGD-O2D
5	F	378[B]	BCL	C3A-C2A-CAA-CBA
5	G	378[B]	BCL	C3A-C2A-CAA-CBA
5	G	372	BCL	C2C-C3C-CAC-CBC
5	G	372	BCL	C4C-C3C-CAC-CBC
5	G	372	BCL	C1-C2-C3-C5
5	G	371	BCL	C1A-C2A-CAA-CBA
5	G	371	BCL	C3A-C2A-CAA-CBA
5	G	371	BCL	CBD-CGD-O2D-CED
5	G	371	BCL	O1D-CGD-O2D-CED
5	G	373	BCL	C2C-C3C-CAC-CBC
5	G	373	BCL	C4C-C3C-CAC-CBC
5	G	373	BCL	CAD-CBD-CGD-O2D
5	G	373	BCL	C2-C3-C5-C6
5	G	374	BCL	C2C-C3C-CAC-CBC
5	G	374	BCL	C4C-C3C-CAC-CBC
5	G	377	BCL	CHA-CBD-CGD-O1D
5	G	377	BCL	CHA-CBD-CGD-O2D
5	A	804	BCL	C2C-C3C-CAC-CBC
5	A	804	BCL	C4C-C3C-CAC-CBC
5	A	804	BCL	CAD-CBD-CGD-O1D
5	A	804	BCL	CAD-CBD-CGD-O2D
5	A	805	BCL	CHA-CBD-CGD-O1D
5	A	805	BCL	CHA-CBD-CGD-O2D
5	A	805	BCL	C1-C2-C3-C5
5	A	806	BCL	C4C-C3C-CAC-CBC
5	A	807	BCL	C1A-C2A-CAA-CBA
5	A	807	BCL	C3A-C2A-CAA-CBA
5	A	807	BCL	C2C-C3C-CAC-CBC
5	A	808	BCL	C1-C2-C3-C4
5	A	808	BCL	C1-C2-C3-C5
5	A	810	BCL	C1A-C2A-CAA-CBA
5	A	810	BCL	C3A-C2A-CAA-CBA
5	A	810	BCL	C4C-C3C-CAC-CBC
5	A	811	BCL	C4C-C3C-CAC-CBC
5	A	811	BCL	C1-C2-C3-C5
5	A	812	BCL	C3A-C2A-CAA-CBA
5	A	812	BCL	C4C-C3C-CAC-CBC
5	A	813	BCL	C2C-C3C-CAC-CBC
5	A	813	BCL	C4C-C3C-CAC-CBC
5	A	813	BCL	CHA-CBD-CGD-O1D

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Mol	Chain	Res	Type	Atoms
5	A	813	BCL	CHA-CBD-CGD-O2D
5	A	814	BCL	CAD-CBD-CGD-O1D
5	A	814	BCL	CAD-CBD-CGD-O2D
5	A	815	BCL	C4C-C3C-CAC-CBC
5	A	815	BCL	CAD-CBD-CGD-O1D
5	A	815	BCL	CAD-CBD-CGD-O2D
5	a	804	BCL	C2C-C3C-CAC-CBC
5	a	804	BCL	C4C-C3C-CAC-CBC
5	a	804	BCL	CAD-CBD-CGD-O1D
5	a	804	BCL	CAD-CBD-CGD-O2D
5	a	805	BCL	CHA-CBD-CGD-O1D
5	a	805	BCL	CHA-CBD-CGD-O2D
5	a	805	BCL	C1-C2-C3-C5
5	a	806	BCL	C4C-C3C-CAC-CBC
5	a	807	BCL	C1A-C2A-CAA-CBA
5	a	807	BCL	C3A-C2A-CAA-CBA
5	a	807	BCL	C2C-C3C-CAC-CBC
5	a	808	BCL	C1-C2-C3-C4
5	a	808	BCL	C1-C2-C3-C5
5	a	810	BCL	C1A-C2A-CAA-CBA
5	a	810	BCL	C3A-C2A-CAA-CBA
5	a	810	BCL	C4C-C3C-CAC-CBC
5	a	811	BCL	C4C-C3C-CAC-CBC
5	a	811	BCL	C1-C2-C3-C5
5	a	812	BCL	C3A-C2A-CAA-CBA
5	a	812	BCL	C4C-C3C-CAC-CBC
5	a	813	BCL	C2C-C3C-CAC-CBC
5	a	813	BCL	C4C-C3C-CAC-CBC
5	a	813	BCL	CHA-CBD-CGD-O1D
5	a	813	BCL	CHA-CBD-CGD-O2D
5	a	814	BCL	CAD-CBD-CGD-O1D
5	a	814	BCL	CAD-CBD-CGD-O2D
5	a	815	BCL	C4C-C3C-CAC-CBC
5	a	815	BCL	CAD-CBD-CGD-O1D
5	a	815	BCL	CAD-CBD-CGD-O2D
7	A	801	GS0	C1A-C2A-CAA-CBA
7	A	801	GS0	C3A-C2A-CAA-CBA
7	A	801	GS0	CHA-CBD-CGD-O2D
7	a	801	GS0	C1A-C2A-CAA-CBA
7	a	801	GS0	C3A-C2A-CAA-CBA
7	a	801	GS0	CHA-CBD-CGD-O2D
8	A	816	F26	C19-C15-C9-C2

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Mol	Chain	Res	Type	Atoms
8	A	816	F26	C22-C25-C26-C29
8	A	816	F26	C27-C28-C31-C36
8	A	816	F26	C37-C34-C35-C32
8	a	816	F26	C19-C15-C9-C2
8	a	816	F26	C22-C25-C26-C29
8	a	816	F26	C27-C28-C31-C36
8	a	816	F26	C37-C34-C35-C32
9	A	817	F39	O2-C13-C14-C18
9	A	817	F39	C19-C20-C27-C32
9	A	817	F39	C25-C20-C27-C32
9	A	817	F39	C22-C21-O6-C15
9	A	817	F39	O7-C21-O6-C15
9	A	817	F39	C27-C32-C35-C37
9	A	817	F39	C32-C35-C37-C39
9	A	817	F39	C38-C37-C39-C40
9	A	817	F39	C37-C39-C40-C41
9	A	817	F39	C43-C42-C44-C51
9	A	817	F39	C44-C51-C57-C59
9	A	817	F39	C46-C53-C56-C58
9	A	817	F39	C53-C56-C58-C60
9	A	817	F39	C56-C58-C61-C63
9	A	817	F39	C60-C58-C61-C63
9	A	817	F39	C57-C59-C62-C64
9	A	817	F39	C57-C59-C62-C65
9	A	817	F39	C61-C63-C64-C62
9	a	817	F39	O2-C13-C14-C18
9	a	817	F39	C19-C20-C27-C32
9	a	817	F39	C25-C20-C27-C32
9	a	817	F39	C22-C21-O6-C15
9	a	817	F39	O7-C21-O6-C15
9	a	817	F39	C27-C32-C35-C37
9	a	817	F39	C32-C35-C37-C39
9	a	817	F39	C38-C37-C39-C40
9	a	817	F39	C37-C39-C40-C41
9	a	817	F39	C43-C42-C44-C51
9	a	817	F39	C44-C51-C57-C59
9	a	817	F39	C46-C53-C56-C58
9	a	817	F39	C53-C56-C58-C60
9	a	817	F39	C56-C58-C61-C63
9	a	817	F39	C60-C58-C61-C63
9	a	817	F39	C57-C59-C62-C64
9	a	817	F39	C57-C59-C62-C65

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Mol	Chain	Res	Type	Atoms
9	a	817	F39	C61-C63-C64-C62
10	A	818	LHG	C1-C2-C3-O3
10	A	818	LHG	C3-O3-P-O5
10	A	818	LHG	C3-O3-P-O6
10	A	818	LHG	C4-O6-P-O3
10	A	818	LHG	C4-O6-P-O4
10	A	818	LHG	C4-O6-P-O5
10	a	818	LHG	C1-C2-C3-O3
10	a	818	LHG	C3-O3-P-O5
10	a	818	LHG	C3-O3-P-O6
10	a	818	LHG	C4-O6-P-O3
10	a	818	LHG	C4-O6-P-O4
10	a	818	LHG	C4-O6-P-O5
11	A	819	LMG	O6-C1-O1-C7
11	A	819	LMG	C11-C10-O7-C8
11	a	819	LMG	O6-C1-O1-C7
11	a	819	LMG	C11-C10-O7-C8
13	A	802	G2O	C5-C6-C7-C8
13	A	802	G2O	C11-C10-C8-C7
13	A	802	G2O	C1A-C2A-CAA-CBA
13	A	802	G2O	C3A-C2A-CAA-CBA
13	A	803	G2O	C11-C10-C8-C7
13	A	803	G2O	C11-C10-C8-C9
13	A	803	G2O	CBD-CGD-O2D-CED
13	A	803	G2O	O1D-CGD-O2D-CED
13	a	802	G2O	C5-C6-C7-C8
13	a	802	G2O	C11-C10-C8-C7
13	a	802	G2O	C1A-C2A-CAA-CBA
13	a	802	G2O	C3A-C2A-CAA-CBA
13	a	803	G2O	C11-C10-C8-C7
13	a	803	G2O	C11-C10-C8-C9
13	a	803	G2O	CBD-CGD-O2D-CED
13	a	803	G2O	O1D-CGD-O2D-CED
9	A	817	F39	O1-C12-C15-O6
9	a	817	F39	O1-C12-C15-O6
11	A	819	LMG	O9-C10-O7-C8
11	a	819	LMG	O9-C10-O7-C8
5	E	375	BCL	C3-C5-C6-C7
5	F	375	BCL	C3-C5-C6-C7
5	G	375	BCL	C3-C5-C6-C7
13	A	803	G2O	C4-C3-C5-C6
13	a	803	G2O	C4-C3-C5-C6

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Mol	Chain	Res	Type	Atoms
5	E	376	BCL	C2-C3-C5-C6
5	F	376	BCL	C2-C3-C5-C6
5	G	376	BCL	C2-C3-C5-C6
5	A	811	BCL	C3-C5-C6-C7
5	a	811	BCL	C3-C5-C6-C7
5	A	812	BCL	C3-C5-C6-C7
5	a	812	BCL	C3-C5-C6-C7
10	A	818	LHG	O2-C2-C3-O3
10	a	818	LHG	O2-C2-C3-O3
9	A	817	F39	C10-C12-C15-O6
9	a	817	F39	C10-C12-C15-O6
13	A	803	G2O	C2-C3-C5-C6
13	a	803	G2O	C2-C3-C5-C6
8	A	816	F26	C14-C10-C8-C13
8	a	816	F26	C14-C10-C8-C13
9	A	817	F39	C13-C14-C18-C19
9	a	817	F39	C13-C14-C18-C19
5	A	811	BCL	C2-C3-C5-C6
5	a	811	BCL	C2-C3-C5-C6
5	A	808	BCL	C6-C7-C8-C9
5	A	812	BCL	C11-C12-C13-C14
5	a	808	BCL	C6-C7-C8-C9
5	a	812	BCL	C11-C12-C13-C14
9	A	817	F39	C65-C62-C64-C63
9	a	817	F39	C65-C62-C64-C63
8	A	816	F26	C22-C25-C26-C30
8	a	816	F26	C22-C25-C26-C30
9	A	817	F39	C40-C41-C42-C44
9	a	817	F39	C40-C41-C42-C44
7	A	801	GS0	C15-C16-C17-C18
7	a	801	GS0	C15-C16-C17-C18
7	A	801	GS0	C6-C7-C8-C10
7	a	801	GS0	C6-C7-C8-C10
8	A	816	F26	C19-C24-C27-C28
8	a	816	F26	C19-C24-C27-C28
9	A	817	F39	C20-C27-C32-C35
9	A	817	F39	C42-C44-C51-C57
9	a	817	F39	C20-C27-C32-C35
9	a	817	F39	C42-C44-C51-C57
11	A	819	LMG	C10-C11-C12-C13
11	a	819	LMG	C10-C11-C12-C13
5	E	372	BCL	C10-C11-C12-C13

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Mol	Chain	Res	Type	Atoms
5	F	372	BCL	C10-C11-C12-C13
5	G	372	BCL	C10-C11-C12-C13
5	E	373	BCL	C2A-CAA-CBA-CGA
5	E	372	BCL	C2A-CAA-CBA-CGA
5	F	372	BCL	C2A-CAA-CBA-CGA
5	F	373	BCL	C2A-CAA-CBA-CGA
5	G	372	BCL	C2A-CAA-CBA-CGA
5	G	373	BCL	C2A-CAA-CBA-CGA
5	A	810	BCL	C2A-CAA-CBA-CGA
5	a	810	BCL	C2A-CAA-CBA-CGA
13	A	802	G2O	C2A-CAA-CBA-CGA
13	A	803	G2O	C2A-CAA-CBA-CGA
13	a	802	G2O	C2A-CAA-CBA-CGA
13	a	803	G2O	C2A-CAA-CBA-CGA
10	A	818	LHG	C23-C24-C25-C26
10	a	818	LHG	C23-C24-C25-C26
5	E	372	BCL	C5-C6-C7-C8
5	F	372	BCL	C5-C6-C7-C8
5	G	372	BCL	C5-C6-C7-C8
11	A	819	LMG	C28-C29-C30-C31
11	a	819	LMG	C28-C29-C30-C31
5	a	806	BCL	C5-C6-C7-C8
5	A	806	BCL	C5-C6-C7-C8
5	A	806	BCL	C8-C10-C11-C12
5	E	376	BCL	O1A-CGA-O2A-C1
5	F	376	BCL	O1A-CGA-O2A-C1
5	G	376	BCL	O1A-CGA-O2A-C1
5	a	806	BCL	C8-C10-C11-C12
13	A	803	G2O	C13-C15-C16-C17
13	a	803	G2O	C13-C15-C16-C17
9	A	817	F39	C23-C24-C26-C28
9	a	817	F39	C23-C24-C26-C28
5	E	371	BCL	C15-C16-C17-C18
5	E	372	BCL	C15-C16-C17-C18
5	F	372	BCL	C15-C16-C17-C18
5	F	371	BCL	C15-C16-C17-C18
5	G	372	BCL	C15-C16-C17-C18
5	G	371	BCL	C15-C16-C17-C18
5	A	808	BCL	C5-C6-C7-C8
5	A	812	BCL	C10-C11-C12-C13
5	a	808	BCL	C5-C6-C7-C8
5	a	812	BCL	C10-C11-C12-C13

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Mol	Chain	Res	Type	Atoms
13	A	802	G2O	C10-C11-C12-C13
13	a	802	G2O	C10-C11-C12-C13
5	E	374	BCL	C10-C11-C12-C13
5	F	374	BCL	C10-C11-C12-C13
5	G	374	BCL	C10-C11-C12-C13
11	A	819	LMG	C2-C1-O1-C7
11	a	819	LMG	C2-C1-O1-C7
5	A	805	BCL	C3-C5-C6-C7
5	a	805	BCL	C3-C5-C6-C7
9	A	817	F39	C53-C56-C58-C61
9	a	817	F39	C53-C56-C58-C61
5	A	811	BCL	O1A-CGA-O2A-C1
5	A	812	BCL	O1A-CGA-O2A-C1
5	a	811	BCL	O1A-CGA-O2A-C1
5	A	815	BCL	C2A-CAA-CBA-CGA
5	a	815	BCL	C2A-CAA-CBA-CGA
13	A	803	G2O	C8-C10-C11-C12
13	a	803	G2O	C8-C10-C11-C12
5	A	805	BCL	O2A-C1-C2-C3
5	a	805	BCL	O2A-C1-C2-C3
5	A	808	BCL	C16-C17-C18-C19
5	a	808	BCL	C16-C17-C18-C19
5	a	812	BCL	O1A-CGA-O2A-C1
13	A	803	G2O	CBA-CGA-O2A-C1
13	a	803	G2O	CBA-CGA-O2A-C1
5	A	806	BCL	O1A-CGA-O2A-C1
5	a	806	BCL	O1A-CGA-O2A-C1
10	A	818	LHG	C13-C14-C15-C16
10	a	818	LHG	C13-C14-C15-C16
5	A	808	BCL	C16-C17-C18-C20
5	a	808	BCL	C16-C17-C18-C20
10	a	818	LHG	C24-C25-C26-C27
5	E	375	BCL	C8-C10-C11-C12
5	F	375	BCL	C8-C10-C11-C12
5	G	375	BCL	C8-C10-C11-C12
10	A	818	LHG	C24-C25-C26-C27
5	E	372	BCL	C3A-C2A-CAA-CBA
5	F	372	BCL	C3A-C2A-CAA-CBA
5	G	372	BCL	C3A-C2A-CAA-CBA
5	A	804	BCL	C3A-C2A-CAA-CBA
5	A	808	BCL	C3A-C2A-CAA-CBA
5	a	804	BCL	C3A-C2A-CAA-CBA

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Mol	Chain	Res	Type	Atoms
5	a	808	BCL	C3A-C2A-CAA-CBA
5	E	372	BCL	C8-C10-C11-C12
5	F	372	BCL	C8-C10-C11-C12
5	G	372	BCL	C8-C10-C11-C12
9	A	817	F39	C58-C61-C63-C64
9	a	817	F39	C58-C61-C63-C64
10	A	818	LHG	C11-C10-C9-C8
10	a	818	LHG	C11-C10-C9-C8
11	A	819	LMG	C29-C28-O8-C9
11	a	819	LMG	C29-C28-O8-C9
5	A	807	BCL	C2A-CAA-CBA-CGA
5	a	807	BCL	C2A-CAA-CBA-CGA
13	A	803	G2O	O1A-CGA-O2A-C1
13	a	803	G2O	O1A-CGA-O2A-C1
9	A	817	F39	C28-C29-C30-C31
9	a	817	F39	C28-C29-C30-C31
8	A	816	F26	C18-C22-C25-C26
8	a	816	F26	C18-C22-C25-C26
5	E	377	BCL	C10-C11-C12-C13
5	F	377	BCL	C10-C11-C12-C13
5	G	377	BCL	C10-C11-C12-C13
10	A	818	LHG	C25-C26-C27-C28
10	a	818	LHG	C25-C26-C27-C28
5	E	374	BCL	C6-C7-C8-C9
5	F	374	BCL	C6-C7-C8-C9
5	G	374	BCL	C6-C7-C8-C9
11	A	819	LMG	C32-C33-C34-C35
11	a	819	LMG	C32-C33-C34-C35
5	E	371	BCL	C5-C6-C7-C8
5	F	371	BCL	C5-C6-C7-C8
5	G	371	BCL	C5-C6-C7-C8
11	A	819	LMG	C15-C16-C17-C18
11	a	819	LMG	C15-C16-C17-C18
8	A	816	F26	C26-C30-C32-C35
8	a	816	F26	C26-C30-C32-C35
5	F	371	BCL	C2-C3-C5-C6
5	G	371	BCL	C2-C3-C5-C6
13	A	802	G2O	C2B-C3B-CAB-CBB
13	a	802	G2O	C2B-C3B-CAB-CBB
5	A	808	BCL	C3-C5-C6-C7
5	A	808	BCL	C10-C11-C12-C13
5	a	808	BCL	C10-C11-C12-C13

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Mol	Chain	Res	Type	Atoms
5	a	808	BCL	C3-C5-C6-C7
11	A	819	LMG	C18-C19-C20-C21
11	a	819	LMG	C18-C19-C20-C21
9	A	817	F39	C24-C26-C28-C29
9	a	817	F39	C24-C26-C28-C29
5	E	371	BCL	C2-C3-C5-C6
10	A	818	LHG	C26-C27-C28-C29
10	a	818	LHG	C26-C27-C28-C29
11	A	819	LMG	O10-C28-O8-C9
11	a	819	LMG	O10-C28-O8-C9
5	E	378[B]	BCL	C1A-C2A-CAA-CBA
5	E	372	BCL	C1A-C2A-CAA-CBA
5	F	372	BCL	C1A-C2A-CAA-CBA
5	F	378[B]	BCL	C1A-C2A-CAA-CBA
5	G	378[B]	BCL	C1A-C2A-CAA-CBA
5	G	372	BCL	C1A-C2A-CAA-CBA
5	A	808	BCL	C1A-C2A-CAA-CBA
5	A	811	BCL	C1A-C2A-CAA-CBA
5	A	812	BCL	C1A-C2A-CAA-CBA
5	A	813	BCL	C1A-C2A-CAA-CBA
5	a	808	BCL	C1A-C2A-CAA-CBA
5	a	811	BCL	C1A-C2A-CAA-CBA
5	a	812	BCL	C1A-C2A-CAA-CBA
5	a	813	BCL	C1A-C2A-CAA-CBA
5	E	375	BCL	C2C-C3C-CAC-CBC
5	F	375	BCL	C2C-C3C-CAC-CBC
5	G	375	BCL	C2C-C3C-CAC-CBC
5	A	805	BCL	C2C-C3C-CAC-CBC
5	A	810	BCL	C2C-C3C-CAC-CBC
5	A	811	BCL	C2C-C3C-CAC-CBC
5	A	812	BCL	C2C-C3C-CAC-CBC
5	A	815	BCL	C2C-C3C-CAC-CBC
5	a	805	BCL	C2C-C3C-CAC-CBC
5	a	810	BCL	C2C-C3C-CAC-CBC
5	a	811	BCL	C2C-C3C-CAC-CBC
5	a	812	BCL	C2C-C3C-CAC-CBC
5	a	815	BCL	C2C-C3C-CAC-CBC
13	A	803	G2O	C14-C13-C15-C16
13	a	803	G2O	C14-C13-C15-C16
11	a	819	LMG	C33-C34-C35-C36
11	A	819	LMG	C33-C34-C35-C36
5	F	375	BCL	C5-C6-C7-C8

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Mol	Chain	Res	Type	Atoms
11	A	819	LMG	C34-C35-C36-C37
11	a	819	LMG	C34-C35-C36-C37
13	A	803	G2O	C6-C7-C8-C10
13	a	803	G2O	C6-C7-C8-C10
5	G	375	BCL	C5-C6-C7-C8
5	E	375	BCL	C5-C6-C7-C8
8	A	816	F26	C36-C31-C33-C38
8	a	816	F26	C36-C31-C33-C38
11	A	819	LMG	O6-C5-C6-O5
11	a	819	LMG	O6-C5-C6-O5
5	a	808	BCL	C8-C10-C11-C12
5	E	376	BCL	CBA-CGA-O2A-C1
5	F	376	BCL	CBA-CGA-O2A-C1
5	G	376	BCL	CBA-CGA-O2A-C1
5	A	808	BCL	C8-C10-C11-C12
5	F	377	BCL	C2-C3-C5-C6
5	G	377	BCL	C2-C3-C5-C6
11	A	819	LMG	C16-C17-C18-C19
11	a	819	LMG	C16-C17-C18-C19
10	a	818	LHG	C29-C30-C31-C32
10	A	818	LHG	C29-C30-C31-C32
7	A	801	GS0	C3-C5-C6-C7
7	a	801	GS0	C3-C5-C6-C7
5	E	374	BCL	C5-C6-C7-C8
5	F	374	BCL	C5-C6-C7-C8
5	G	374	BCL	C5-C6-C7-C8
5	A	806	BCL	CBA-CGA-O2A-C1
5	a	806	BCL	CBA-CGA-O2A-C1
9	A	817	F39	C29-C30-C31-C33
5	E	376	BCL	C4-C3-C5-C6
5	F	376	BCL	C4-C3-C5-C6
5	G	376	BCL	C4-C3-C5-C6
5	E	377	BCL	C2-C3-C5-C6
9	a	817	F39	C29-C30-C31-C33
5	E	371	BCL	C14-C13-C15-C16
5	E	373	BCL	C11-C10-C8-C9
5	E	375	BCL	C11-C10-C8-C9
5	E	372	BCL	C11-C10-C8-C9
5	F	372	BCL	C11-C10-C8-C9
5	F	371	BCL	C14-C13-C15-C16
5	F	373	BCL	C11-C10-C8-C9
5	F	375	BCL	C11-C10-C8-C9

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Mol	Chain	Res	Type	Atoms
5	G	372	BCL	C11-C10-C8-C9
5	G	371	BCL	C14-C13-C15-C16
5	G	373	BCL	C11-C10-C8-C9
5	G	375	BCL	C11-C10-C8-C9
5	A	811	BCL	C11-C12-C13-C14
5	a	811	BCL	C11-C12-C13-C14
7	A	801	GS0	C6-C7-C8-C9
7	a	801	GS0	C6-C7-C8-C9
9	A	817	F39	C16-C13-C14-C18
9	A	817	F39	C17-C13-C14-C18
9	a	817	F39	C16-C13-C14-C18
9	a	817	F39	C17-C13-C14-C18
5	E	371	BCL	C12-C13-C15-C16
5	E	373	BCL	C11-C10-C8-C7
5	E	375	BCL	C11-C10-C8-C7
5	E	377	BCL	C11-C10-C8-C7
5	E	372	BCL	C11-C10-C8-C7
5	F	372	BCL	C11-C10-C8-C7
5	F	371	BCL	C12-C13-C15-C16
5	F	373	BCL	C11-C10-C8-C7
5	F	375	BCL	C11-C10-C8-C7
5	F	377	BCL	C11-C10-C8-C7
5	G	372	BCL	C11-C10-C8-C7
5	G	371	BCL	C12-C13-C15-C16
5	G	373	BCL	C11-C10-C8-C7
5	G	375	BCL	C11-C10-C8-C7
5	G	377	BCL	C11-C10-C8-C7
5	A	811	BCL	C11-C12-C13-C15
5	a	811	BCL	C11-C12-C13-C15
13	A	803	G2O	C12-C13-C15-C16
13	a	803	G2O	C12-C13-C15-C16
5	E	377	BCL	C4-C3-C5-C6
5	F	377	BCL	C4-C3-C5-C6
5	G	377	BCL	C4-C3-C5-C6
5	F	375	BCL	C16-C17-C18-C20
8	A	816	F26	C27-C28-C31-C33
8	a	816	F26	C27-C28-C31-C33
5	E	375	BCL	C16-C17-C18-C20
5	G	375	BCL	C16-C17-C18-C20
5	A	811	BCL	C5-C6-C7-C8
5	a	811	BCL	C5-C6-C7-C8
7	A	801	GS0	C10-C11-C12-C13

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Mol	Chain	Res	Type	Atoms
7	a	801	GS0	C10-C11-C12-C13
8	A	816	F26	C18-C13-C8-C10
8	a	816	F26	C18-C13-C8-C10
5	E	373	BCL	C1-C2-C3-C5
5	E	375	BCL	C1-C2-C3-C5
5	E	377	BCL	C1-C2-C3-C5
5	F	373	BCL	C1-C2-C3-C5
5	F	375	BCL	C1-C2-C3-C5
5	F	377	BCL	C1-C2-C3-C5
5	G	373	BCL	C1-C2-C3-C5
5	G	375	BCL	C1-C2-C3-C5
5	G	377	BCL	C1-C2-C3-C5
9	A	817	F39	C17-C13-O2-C11
9	a	817	F39	C17-C13-O2-C11
11	A	819	LMG	C17-C18-C19-C20
11	a	819	LMG	C17-C18-C19-C20
10	A	818	LHG	C15-C16-C17-C18
10	a	818	LHG	C15-C16-C17-C18
8	A	816	F26	C29-C26-C30-C32
8	a	816	F26	C29-C26-C30-C32
5	A	812	BCL	CBA-CGA-O2A-C1
5	a	812	BCL	CBA-CGA-O2A-C1
5	E	371	BCL	C11-C12-C13-C15
5	F	371	BCL	C11-C12-C13-C15
5	G	371	BCL	C11-C12-C13-C15
7	A	801	GS0	O1A-CGA-O2A-C1
7	a	801	GS0	O1A-CGA-O2A-C1
13	A	802	G2O	C3-C5-C6-C7
13	a	802	G2O	C3-C5-C6-C7
5	A	806	BCL	C3-C5-C6-C7
5	a	806	BCL	C3-C5-C6-C7
5	A	812	BCL	C8-C10-C11-C12
5	a	812	BCL	C8-C10-C11-C12
11	A	819	LMG	O1-C7-C8-O7
11	a	819	LMG	O1-C7-C8-O7
5	E	376	BCL	C3-C5-C6-C7
5	F	376	BCL	C3-C5-C6-C7
5	G	376	BCL	C3-C5-C6-C7
5	E	374	BCL	C8-C10-C11-C12
13	a	803	G2O	C15-C16-C17-C18
5	A	814	BCL	C2A-CAA-CBA-CGA
5	a	814	BCL	C2A-CAA-CBA-CGA

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Mol	Chain	Res	Type	Atoms
5	A	806	BCL	C16-C17-C18-C20
5	a	806	BCL	C16-C17-C18-C20
5	F	374	BCL	C8-C10-C11-C12
5	G	374	BCL	C8-C10-C11-C12
13	A	803	G2O	C15-C16-C17-C18
5	A	804	BCL	C1A-C2A-CAA-CBA
5	a	804	BCL	C1A-C2A-CAA-CBA
5	E	373	BCL	CBA-CGA-O2A-C1
5	F	373	BCL	CBA-CGA-O2A-C1
5	G	373	BCL	CBA-CGA-O2A-C1
5	E	371	BCL	C4-C3-C5-C6
5	F	371	BCL	C4-C3-C5-C6
5	G	371	BCL	C4-C3-C5-C6
13	A	803	G2O	C10-C11-C12-C13
13	a	803	G2O	C10-C11-C12-C13
13	A	802	G2O	C6-C7-C8-C9
13	a	802	G2O	C6-C7-C8-C9
5	E	374	BCL	C11-C12-C13-C15
5	E	372	BCL	C11-C12-C13-C15
5	F	372	BCL	C11-C12-C13-C15
5	F	374	BCL	C11-C12-C13-C15
5	G	372	BCL	C11-C12-C13-C15
5	G	374	BCL	C11-C12-C13-C15
5	A	806	BCL	C11-C12-C13-C15
5	a	806	BCL	C11-C12-C13-C15
5	A	808	BCL	C2C-C3C-CAC-CBC
5	a	808	BCL	C2C-C3C-CAC-CBC
7	A	801	GS0	C16-C17-C18-C20
7	a	801	GS0	C16-C17-C18-C20
5	E	377	BCL	C11-C10-C8-C9
5	F	377	BCL	C11-C10-C8-C9
5	G	377	BCL	C11-C10-C8-C9
5	A	811	BCL	C6-C7-C8-C9
5	a	811	BCL	C6-C7-C8-C9
5	A	808	BCL	C15-C16-C17-C18
5	a	808	BCL	C15-C16-C17-C18
5	A	808	BCL	CAD-CBD-CGD-O2D
5	a	808	BCL	CAD-CBD-CGD-O2D
10	A	818	LHG	C27-C28-C29-C30
10	a	818	LHG	C27-C28-C29-C30
5	E	373	BCL	CAD-CBD-CGD-O1D
5	F	373	BCL	CAD-CBD-CGD-O1D

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Mol	Chain	Res	Type	Atoms
5	G	373	BCL	CAD-CBD-CGD-O1D
5	A	808	BCL	CAD-CBD-CGD-O1D
5	A	809	BCL	CHA-CBD-CGD-O1D
5	A	809	BCL	CHA-CBD-CGD-O2D
5	A	810	BCL	CAD-CBD-CGD-O1D
5	a	808	BCL	CAD-CBD-CGD-O1D
5	a	809	BCL	CHA-CBD-CGD-O1D
5	a	809	BCL	CHA-CBD-CGD-O2D
7	A	801	GS0	CHA-CBD-CGD-O1D
7	a	801	GS0	CHA-CBD-CGD-O1D
10	A	818	LHG	C3-O3-P-O4
10	a	818	LHG	C3-O3-P-O4
8	A	816	F26	C17-C13-C8-C10
8	a	816	F26	C17-C13-C8-C10
9	A	817	F39	C18-C19-C20-C27
9	a	817	F39	C18-C19-C20-C27
10	A	818	LHG	C2-C3-O3-P
10	a	818	LHG	C2-C3-O3-P
5	E	376	BCL	C10-C11-C12-C13
5	F	376	BCL	C10-C11-C12-C13
5	G	376	BCL	C10-C11-C12-C13
5	F	374	BCL	C3-C5-C6-C7
5	G	374	BCL	C3-C5-C6-C7
5	E	374	BCL	C3-C5-C6-C7
5	a	805	BCL	C5-C6-C7-C8
5	A	805	BCL	C5-C6-C7-C8
5	E	374	BCL	C11-C12-C13-C14
5	F	374	BCL	C11-C12-C13-C14
5	G	374	BCL	C11-C12-C13-C14
5	A	806	BCL	C11-C12-C13-C14
5	a	806	BCL	C11-C12-C13-C14
9	A	817	F39	C31-C33-C34-C36
9	a	817	F39	C31-C33-C34-C36
11	A	819	LMG	O1-C7-C8-C9
11	a	819	LMG	O1-C7-C8-C9
5	E	374	BCL	C2A-CAA-CBA-CGA
5	F	374	BCL	C2A-CAA-CBA-CGA
5	G	374	BCL	C2A-CAA-CBA-CGA
5	E	375	BCL	C16-C17-C18-C19
5	F	375	BCL	C16-C17-C18-C19
5	G	375	BCL	C16-C17-C18-C19
5	E	372	BCL	C11-C12-C13-C14

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Mol	Chain	Res	Type	Atoms
5	F	372	BCL	C11-C12-C13-C14
5	G	372	BCL	C11-C12-C13-C14
5	A	811	BCL	CBA-CGA-O2A-C1
5	a	811	BCL	CBA-CGA-O2A-C1
13	A	802	G2O	C4B-C3B-CAB-CBB
13	a	802	G2O	C4B-C3B-CAB-CBB
5	A	812	BCL	C11-C12-C13-C15
5	a	812	BCL	C11-C12-C13-C15
5	a	804	BCL	C2A-CAA-CBA-CGA
5	A	804	BCL	C2A-CAA-CBA-CGA
10	A	818	LHG	O1-C1-C2-C3
10	a	818	LHG	O1-C1-C2-C3
5	A	805	BCL	C1-C2-C3-C4
5	a	805	BCL	C1-C2-C3-C4
5	A	805	BCL	C1A-C2A-CAA-CBA
5	a	805	BCL	C1A-C2A-CAA-CBA
5	A	809	BCL	CAA-CBA-CGA-O1A
5	a	809	BCL	CAA-CBA-CGA-O1A
13	A	803	G2O	C6-C7-C8-C9
13	a	803	G2O	C6-C7-C8-C9
5	a	808	BCL	C4-C3-C5-C6
5	A	808	BCL	C6-C7-C8-C10
5	a	808	BCL	C6-C7-C8-C10
5	a	810	BCL	CAA-CBA-CGA-O1A
5	A	810	BCL	CAA-CBA-CGA-O1A
5	A	808	BCL	C4-C3-C5-C6
7	A	801	GS0	C4-C3-C5-C6
7	a	801	GS0	C4-C3-C5-C6
5	A	812	BCL	C13-C15-C16-C17
5	a	812	BCL	C13-C15-C16-C17
9	A	817	F39	C16-C13-O2-C11
9	a	817	F39	C16-C13-O2-C11
5	E	376	BCL	C15-C16-C17-C18
9	a	817	F39	C22-C23-C24-C26
13	A	802	G2O	C6-C7-C8-C10
13	a	802	G2O	C6-C7-C8-C10
9	A	817	F39	C22-C23-C24-C26
5	F	376	BCL	C15-C16-C17-C18
5	G	376	BCL	C15-C16-C17-C18
10	a	818	LHG	C32-C33-C34-C35
7	A	801	GS0	CAA-CBA-CGA-O2A
10	A	818	LHG	C32-C33-C34-C35

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Mol	Chain	Res	Type	Atoms
5	A	805	BCL	C4-C3-C5-C6
5	a	805	BCL	C4-C3-C5-C6
13	A	803	G2O	C5-C6-C7-C8
13	a	803	G2O	C5-C6-C7-C8
7	a	801	GS0	CAA-CBA-CGA-O2A
5	A	812	BCL	C2A-CAA-CBA-CGA
5	a	812	BCL	C2A-CAA-CBA-CGA
5	E	374	BCL	C2-C1-O2A-CGA
5	E	377	BCL	C2-C1-O2A-CGA
5	F	374	BCL	C2-C1-O2A-CGA
5	F	377	BCL	C2-C1-O2A-CGA
5	G	374	BCL	C2-C1-O2A-CGA
5	G	377	BCL	C2-C1-O2A-CGA
5	A	805	BCL	C2-C1-O2A-CGA
5	a	805	BCL	C2-C1-O2A-CGA
5	F	374	BCL	O1A-CGA-O2A-C1
5	E	374	BCL	O1A-CGA-O2A-C1
9	a	817	F39	C18-C19-C20-C25
5	G	374	BCL	O1A-CGA-O2A-C1
9	A	817	F39	C18-C19-C20-C25
8	A	816	F26	C34-C37-C39-C38
8	a	816	F26	C34-C37-C39-C38
5	A	806	BCL	C6-C7-C8-C9
5	A	811	BCL	C14-C13-C15-C16
5	a	806	BCL	C6-C7-C8-C9
5	a	811	BCL	C14-C13-C15-C16
5	A	809	BCL	CAA-CBA-CGA-O2A
5	a	809	BCL	CAA-CBA-CGA-O2A
5	E	374	BCL	C6-C7-C8-C10
5	F	374	BCL	C6-C7-C8-C10
5	G	374	BCL	C6-C7-C8-C10
5	A	811	BCL	C12-C13-C15-C16
5	a	811	BCL	C12-C13-C15-C16
5	E	375	BCL	C2-C3-C5-C6
5	F	375	BCL	C2-C3-C5-C6
5	G	375	BCL	C2-C3-C5-C6
5	E	372	BCL	C2-C3-C5-C6
5	F	372	BCL	C2-C3-C5-C6
5	G	372	BCL	C2-C3-C5-C6
5	A	804	BCL	CAA-CBA-CGA-O2A
5	a	804	BCL	CAA-CBA-CGA-O2A
5	E	372	BCL	C6-C7-C8-C9

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Mol	Chain	Res	Type	Atoms
5	F	372	BCL	C6-C7-C8-C9
5	G	372	BCL	C6-C7-C8-C9
5	E	377	BCL	C15-C16-C17-C18
5	F	377	BCL	C15-C16-C17-C18
5	G	377	BCL	C15-C16-C17-C18
5	E	376	BCL	C2-C1-O2A-CGA
5	F	376	BCL	C2-C1-O2A-CGA
5	G	376	BCL	C2-C1-O2A-CGA
5	A	804	BCL	CAA-CBA-CGA-O1A
5	E	377	BCL	C2C-C3C-CAC-CBC
5	F	377	BCL	C2C-C3C-CAC-CBC
5	G	377	BCL	C2C-C3C-CAC-CBC
5	a	804	BCL	CAA-CBA-CGA-O1A
9	A	817	F39	C48-C46-C53-C56
9	a	817	F39	C48-C46-C53-C56
5	E	374	BCL	CAA-CBA-CGA-O2A
5	F	374	BCL	CAA-CBA-CGA-O2A
5	A	811	BCL	C3A-C2A-CAA-CBA
5	a	811	BCL	C3A-C2A-CAA-CBA
9	A	817	F39	C14-C13-O2-C11
9	a	817	F39	C14-C13-O2-C11
5	G	374	BCL	CAA-CBA-CGA-O2A
5	E	372	BCL	C13-C15-C16-C17
10	a	818	LHG	C17-C18-C19-C20
5	F	372	BCL	C13-C15-C16-C17
5	G	372	BCL	C13-C15-C16-C17
10	A	818	LHG	C17-C18-C19-C20
11	A	819	LMG	C8-C7-O1-C1
11	a	819	LMG	C8-C7-O1-C1
5	A	809	BCL	CAD-CBD-CGD-O2D
5	A	810	BCL	CAD-CBD-CGD-O2D
5	a	809	BCL	CAD-CBD-CGD-O2D
5	a	810	BCL	CAD-CBD-CGD-O2D
13	A	802	G2O	CAD-CBD-CGD-O2D
13	a	802	G2O	CAD-CBD-CGD-O2D
5	E	373	BCL	C2-C1-O2A-CGA
5	F	373	BCL	C2-C1-O2A-CGA
5	G	373	BCL	C2-C1-O2A-CGA
5	A	811	BCL	C2-C1-O2A-CGA
5	a	811	BCL	C2-C1-O2A-CGA
5	E	377	BCL	C1-C2-C3-C4
5	F	377	BCL	C1-C2-C3-C4

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Mol	Chain	Res	Type	Atoms
5	G	377	BCL	C1-C2-C3-C4
11	a	819	LMG	C31-C32-C33-C34
5	E	371	BCL	CAA-CBA-CGA-O2A
5	a	810	BCL	CAA-CBA-CGA-O2A
11	A	819	LMG	C31-C32-C33-C34
5	G	371	BCL	CAA-CBA-CGA-O2A
5	a	812	BCL	C15-C16-C17-C18

There are no ring outliers.

49 monomers are involved in 142 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	E	378[B]	BCL	1	0
5	F	374	BCL	3	0
13	a	803	G2O	2	0
5	G	376	BCL	7	0
11	a	819	LMG	3	0
5	G	378[B]	BCL	1	0
5	G	372	BCL	5	0
5	a	811	BCL	4	0
5	F	372	BCL	4	0
5	E	377	BCL	3	0
5	E	372	BCL	3	0
5	E	373	BCL	4	0
5	G	375	BCL	5	0
5	A	814	BCL	2	0
6	B	302	SF4	1	0
5	F	377	BCL	3	0
5	A	806	BCL	4	0
5	A	808	BCL	3	0
5	a	810	BCL	5	0
5	E	374	BCL	2	0
5	E	376	BCL	5	0
5	A	809	BCL	3	0
5	E	375	BCL	6	0
5	A	812	BCL	5	0
5	E	371	BCL	5	0
5	a	814	BCL	2	0
13	A	803	G2O	1	0
5	G	373	BCL	1	0
5	F	375	BCL	6	0
5	G	374	BCL	2	0

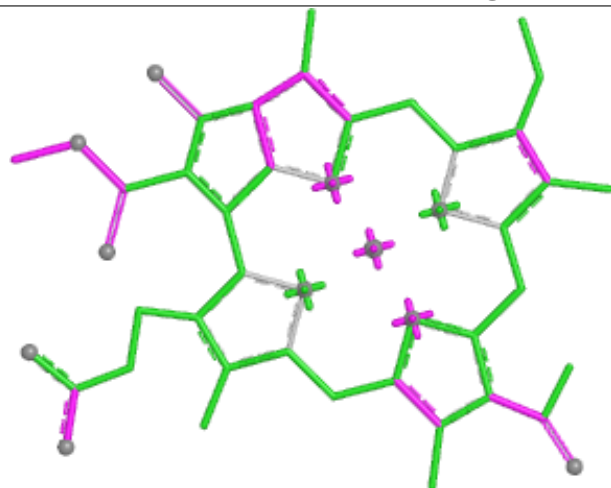
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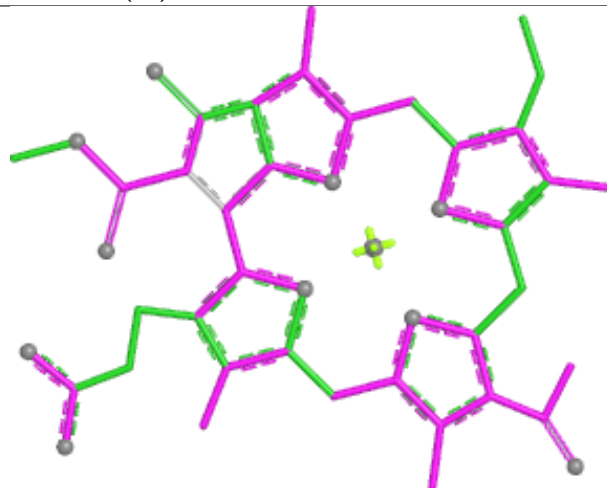
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	a	808	BCL	2	0
5	a	807	BCL	3	0
5	G	371	BCL	4	0
5	A	815	BCL	2	0
5	a	809	BCL	3	0
5	F	373	BCL	3	0
5	a	806	BCL	4	0
5	A	810	BCL	6	0
5	A	813	BCL	2	0
5	a	813	BCL	2	0
5	a	815	BCL	2	0
5	G	377	BCL	3	0
5	a	812	BCL	3	0
5	A	811	BCL	4	0
5	F	376	BCL	2	0
5	A	807	BCL	3	0
11	A	819	LMG	3	0
6	B	301	SF4	3	0
5	F	371	BCL	4	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 than 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.

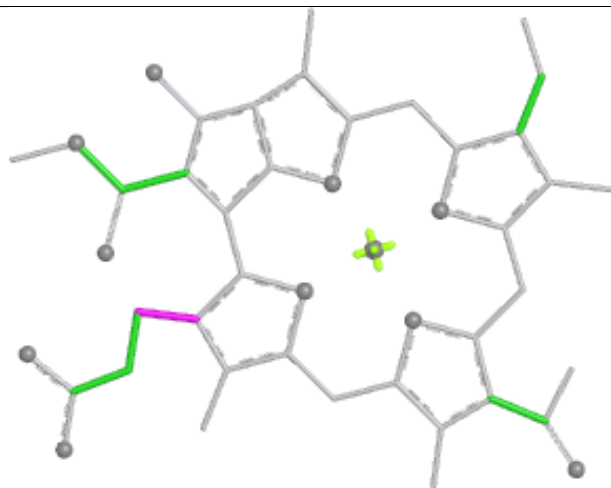
Ligand BCL E 378 (B)



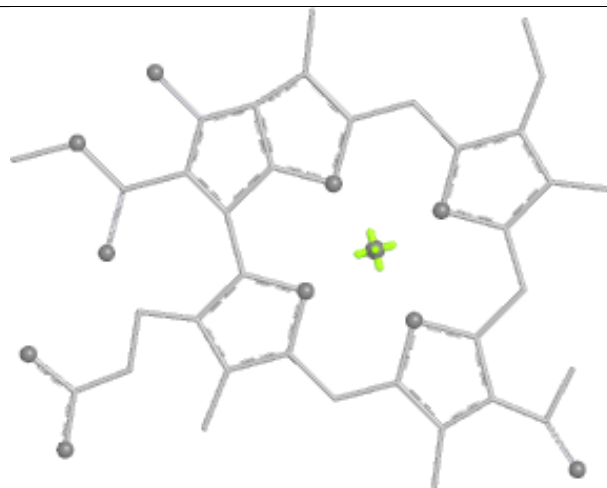
Bond lengths



Bond angles

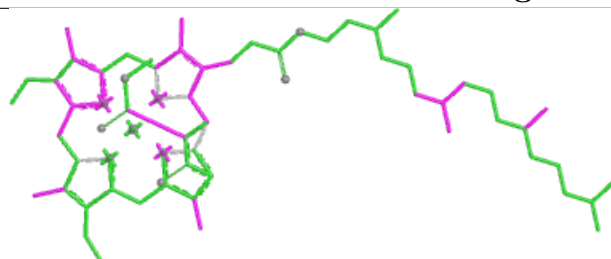


Torsions

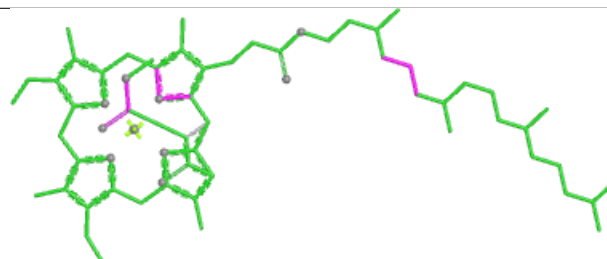


Rings

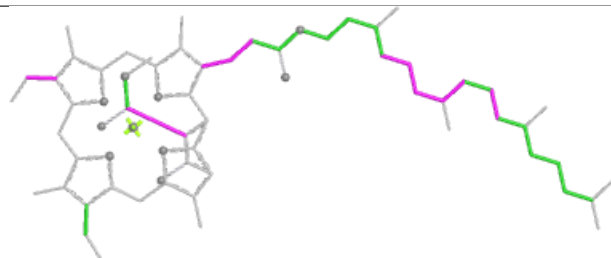
Ligand G2O A 802



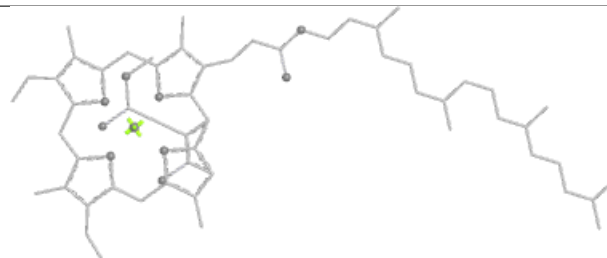
Bond lengths



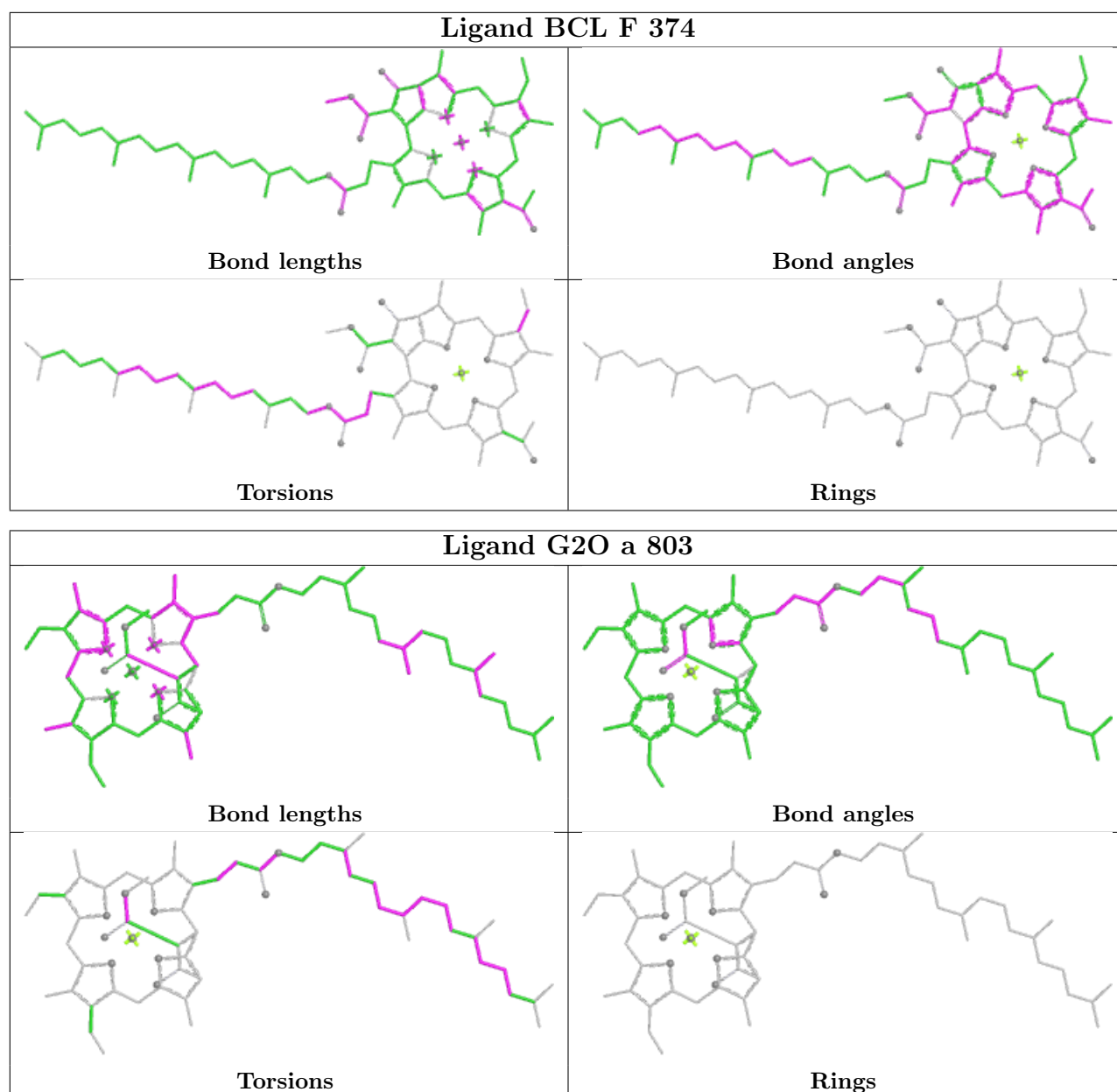
Bond angles

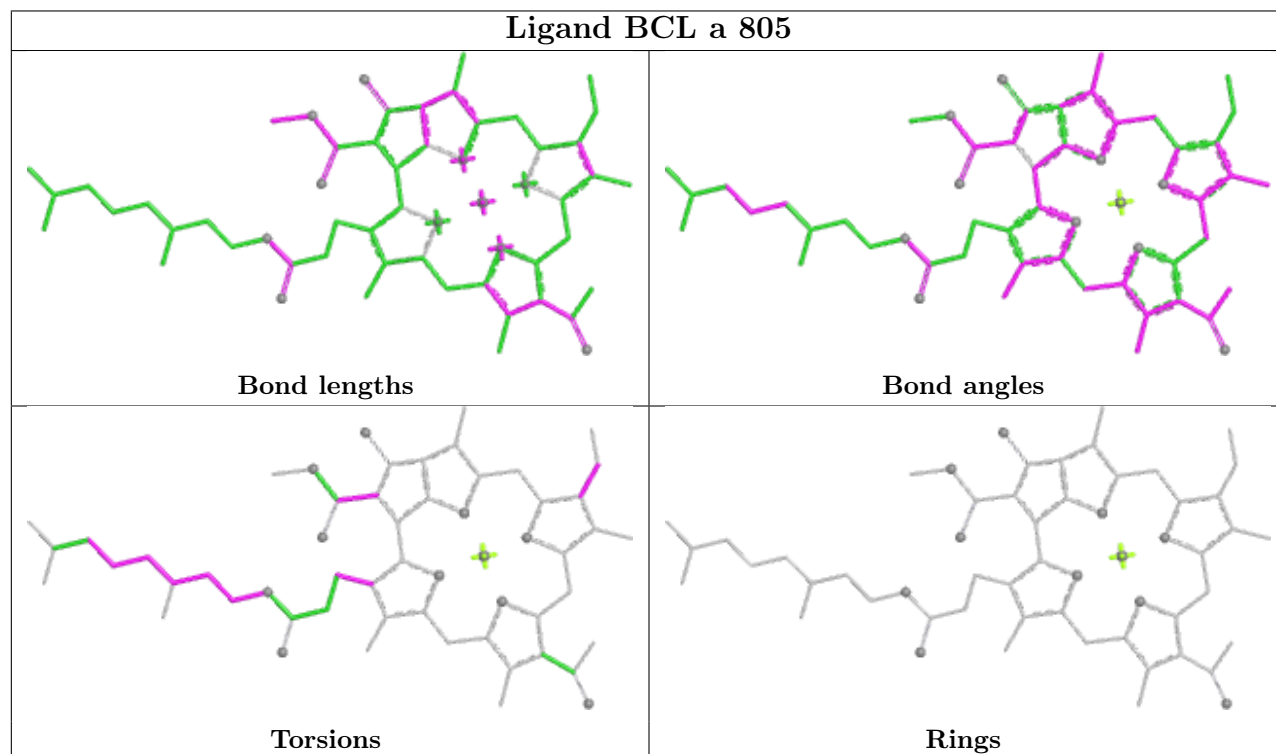
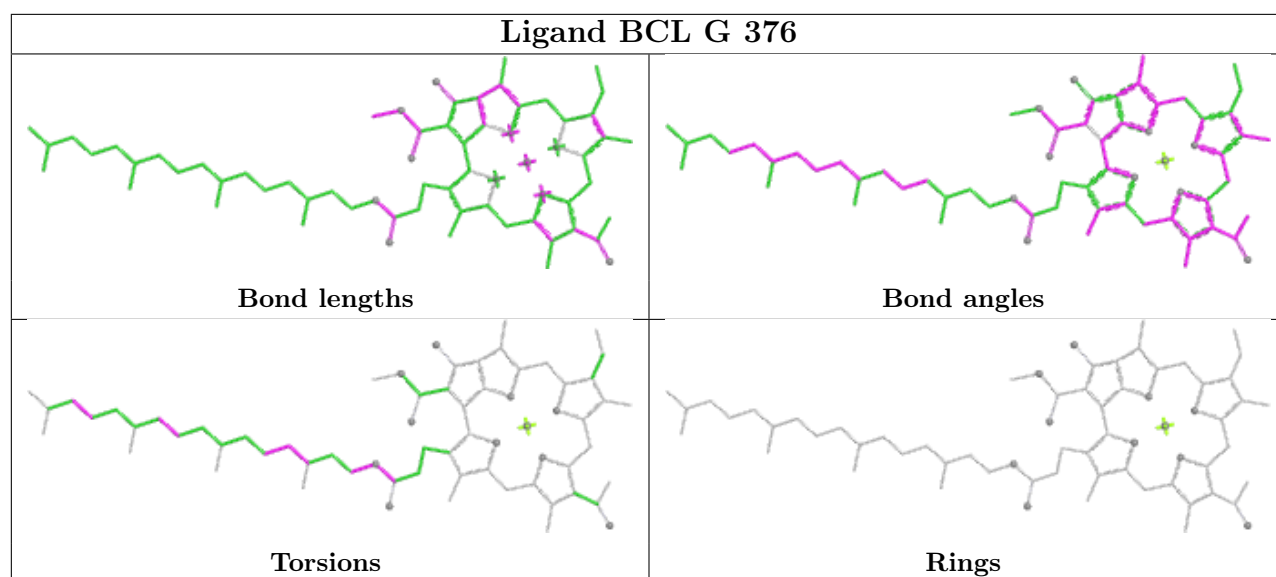


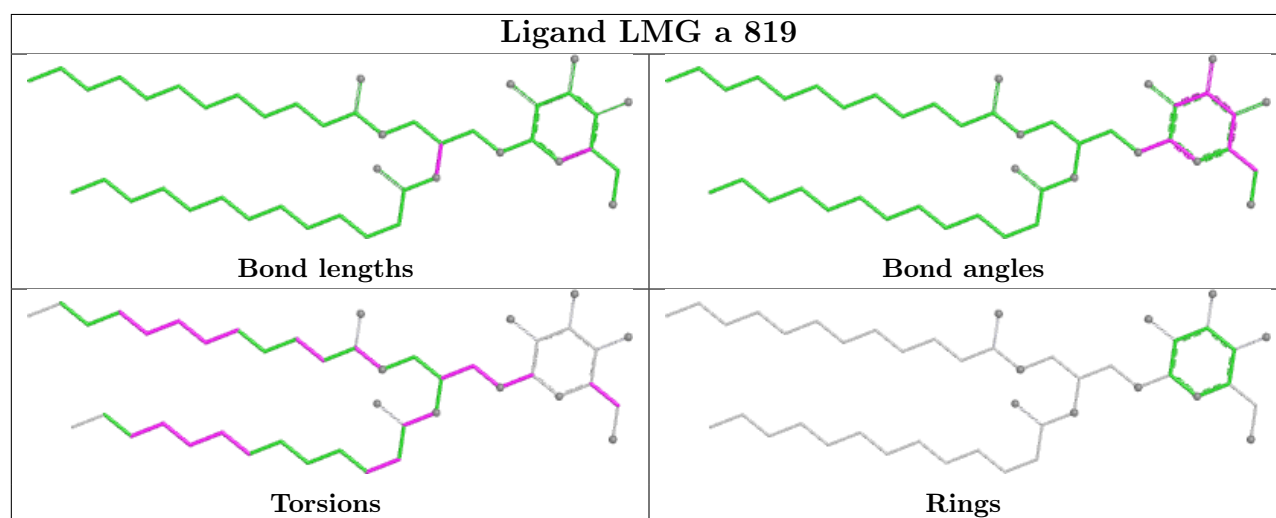
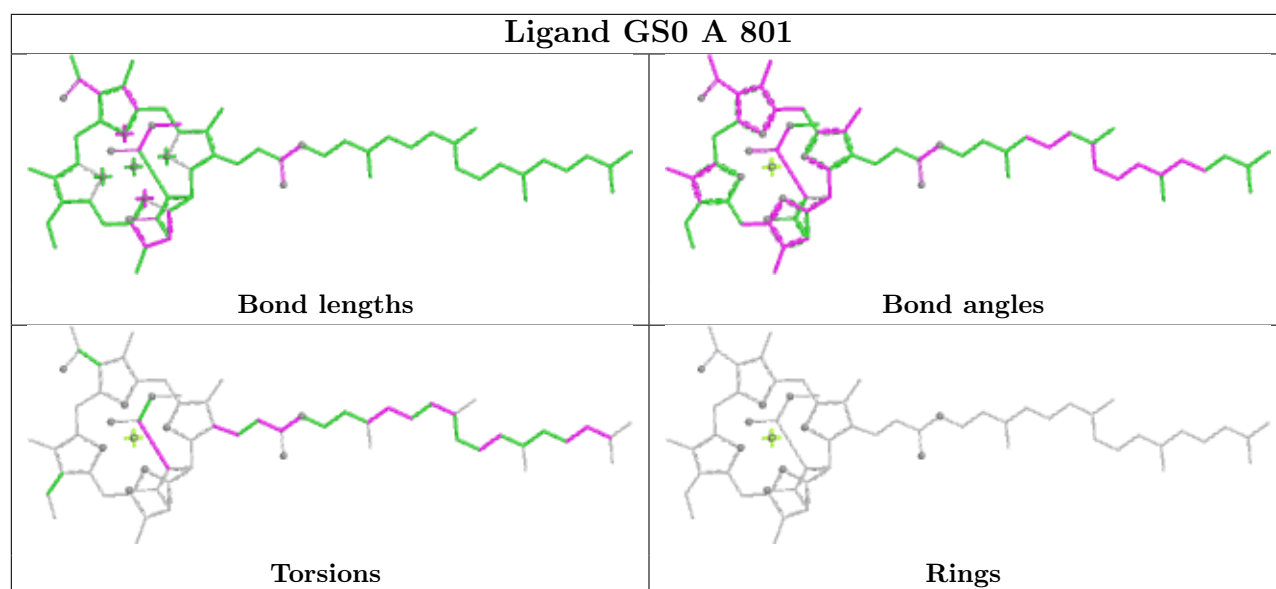
Torsions



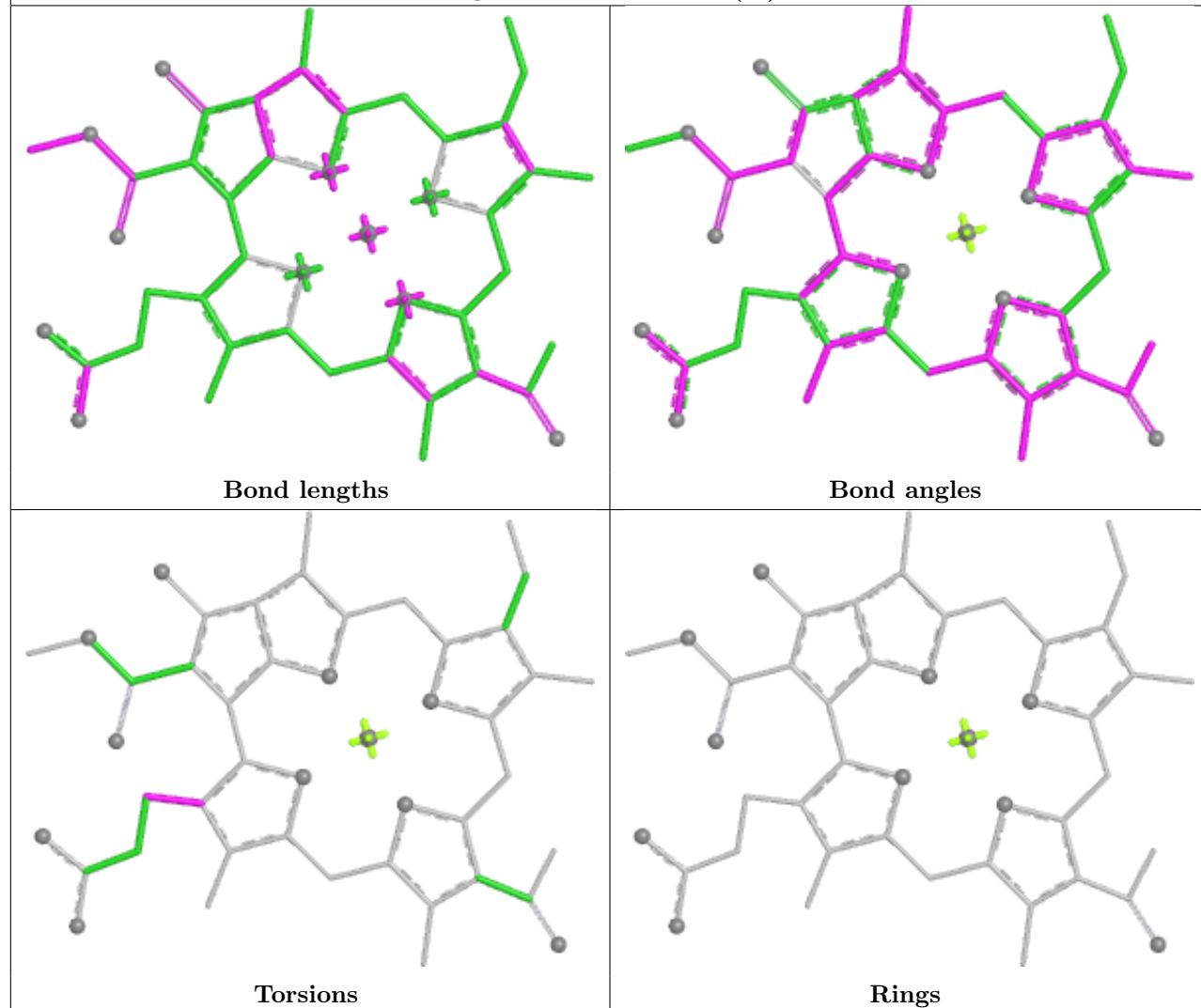
Rings



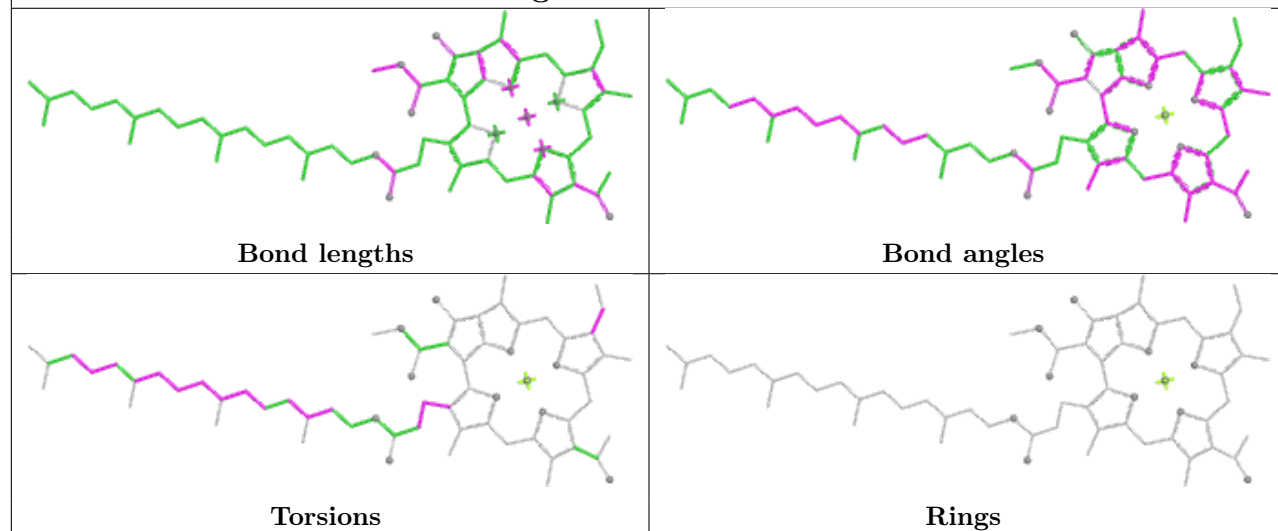


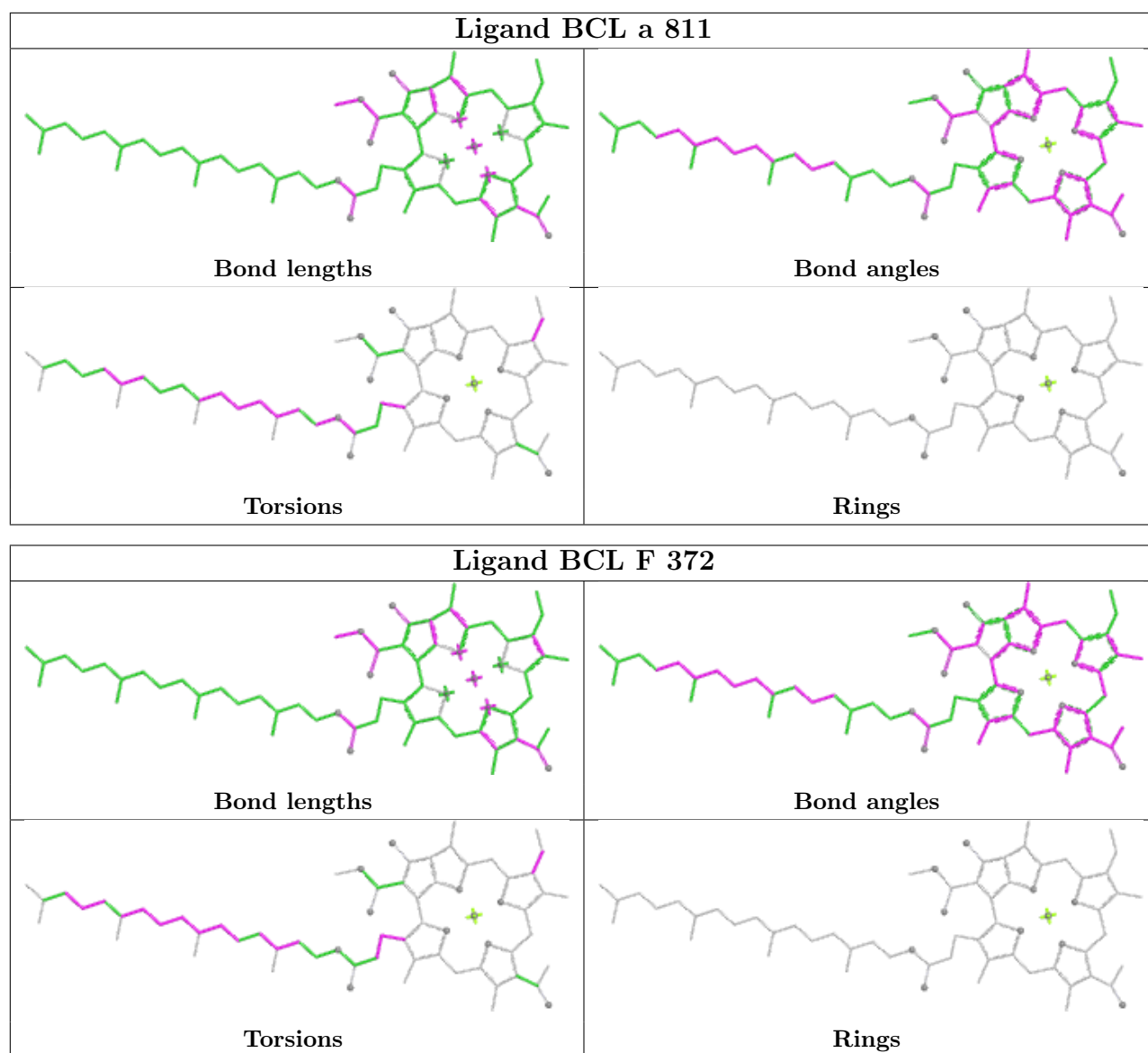


Ligand BCL G 378 (B)

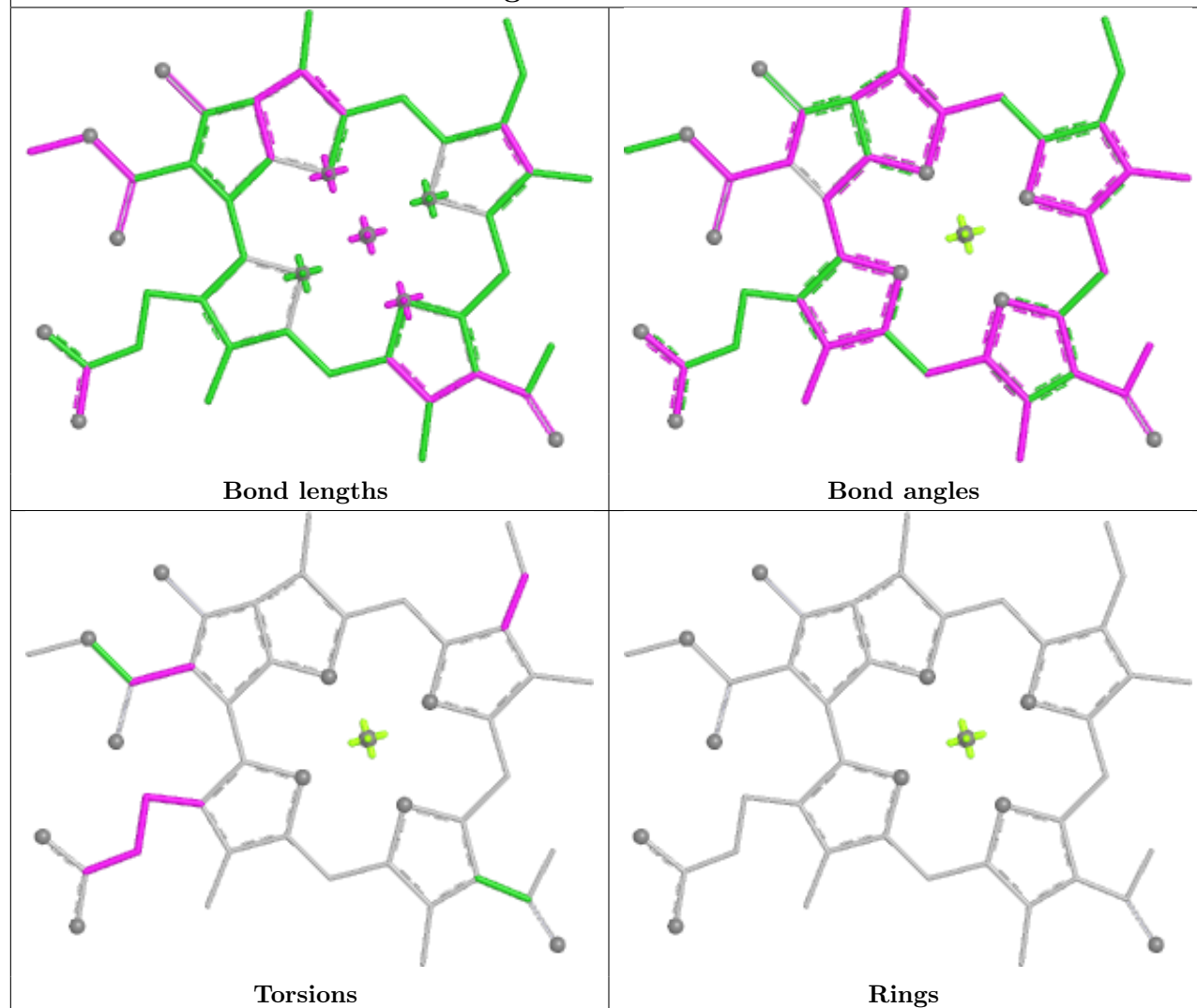


Ligand BCL G 372

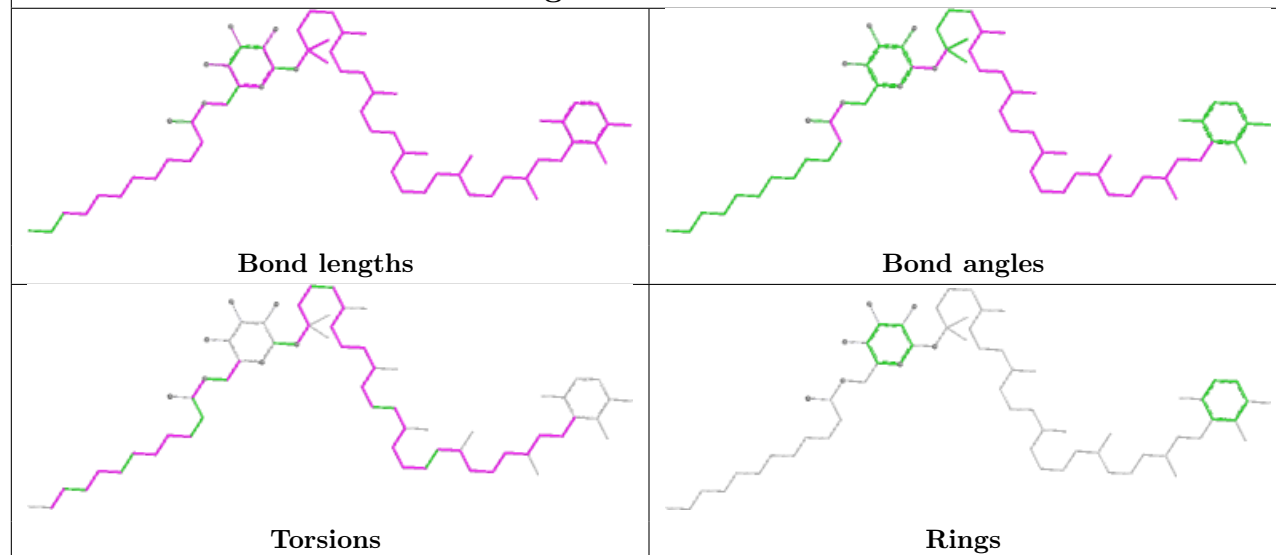


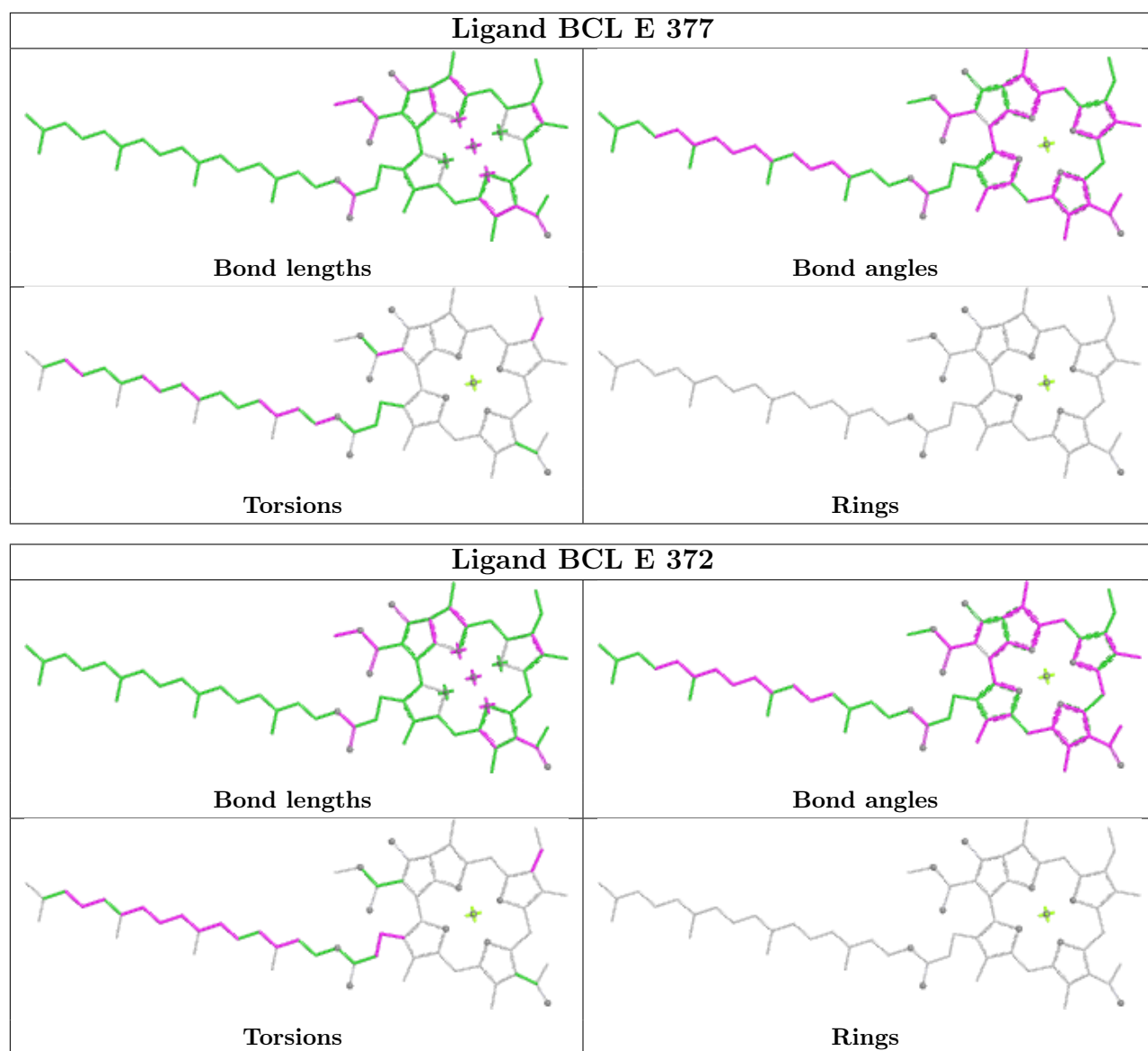


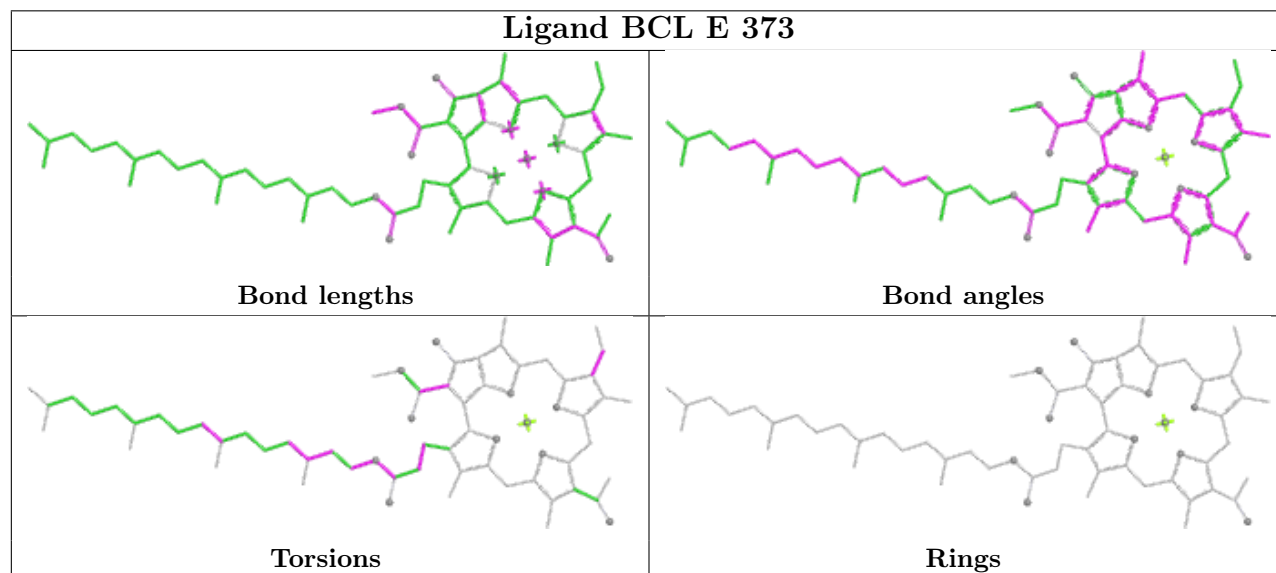
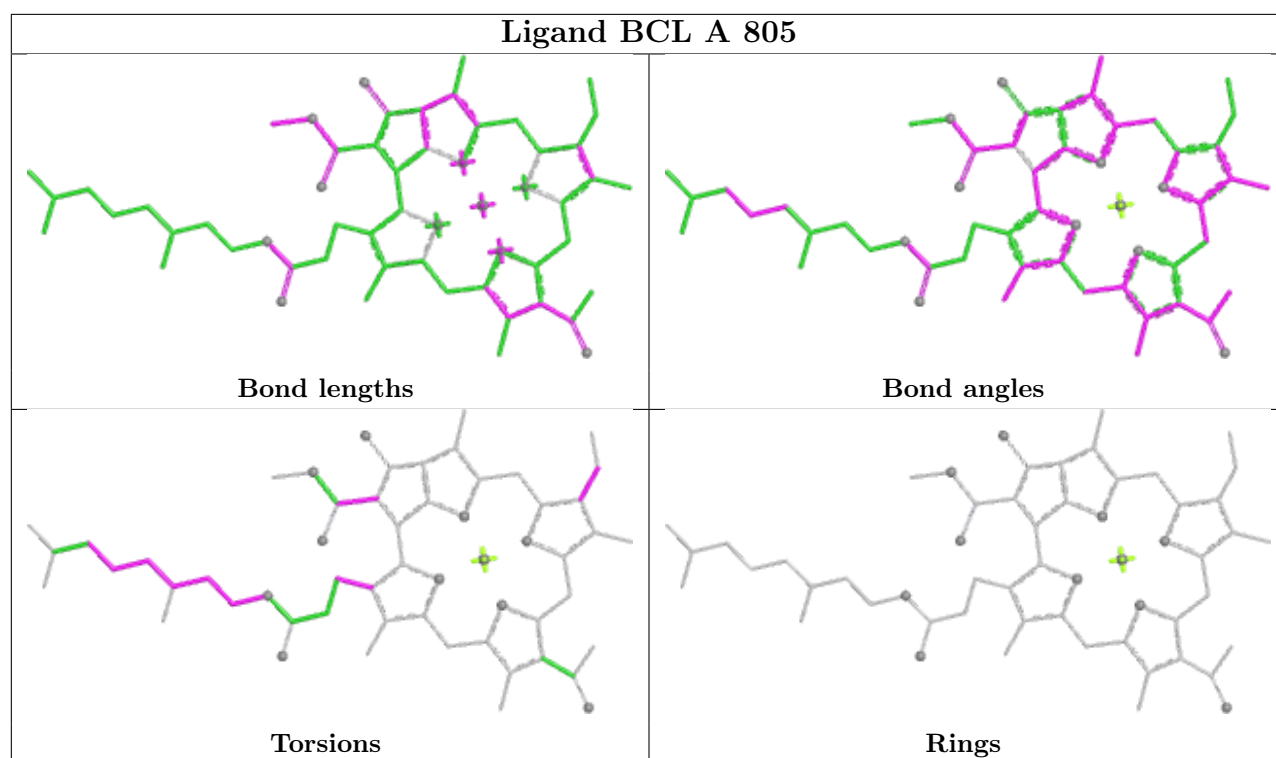
Ligand BCL a 804

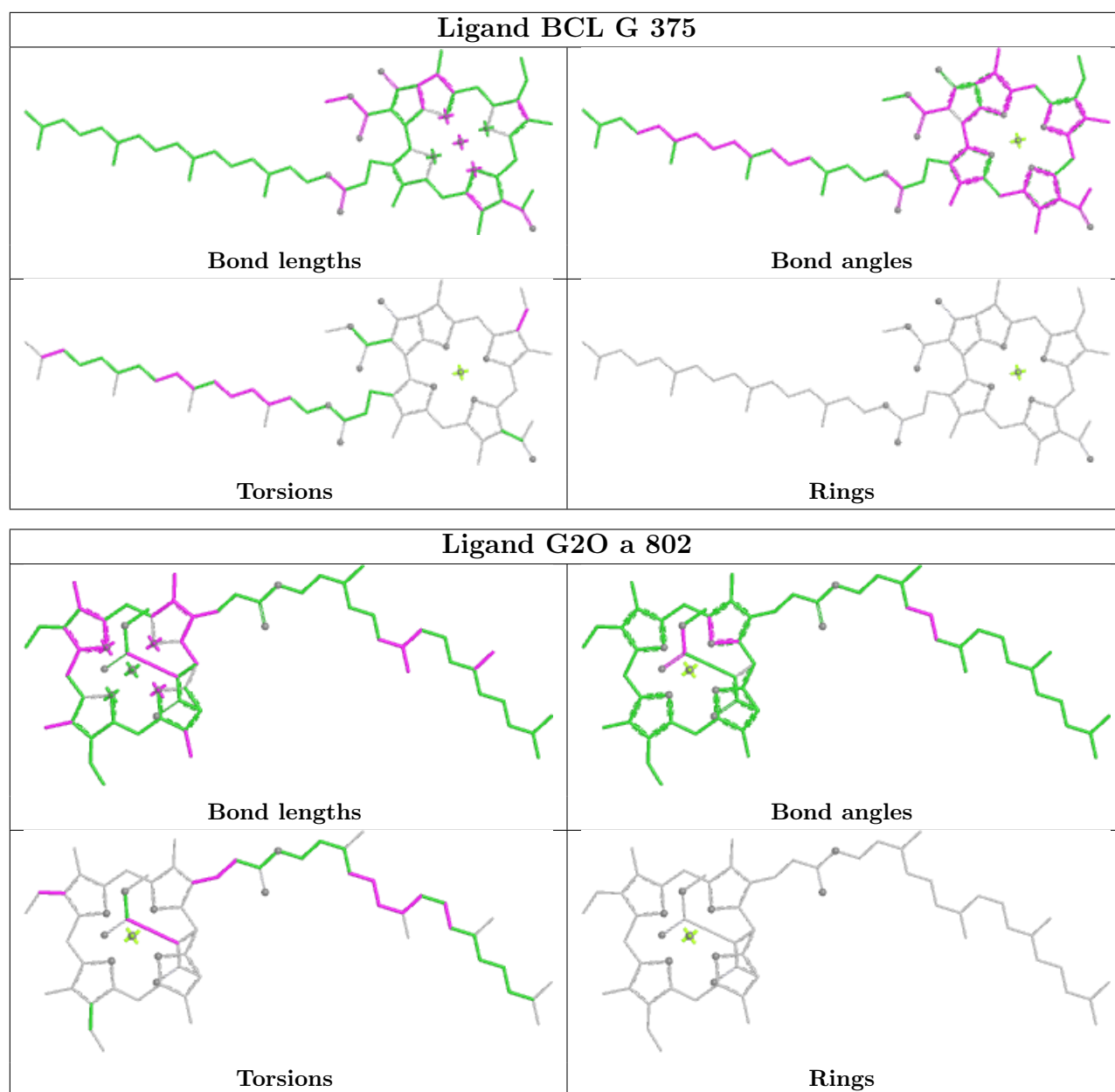


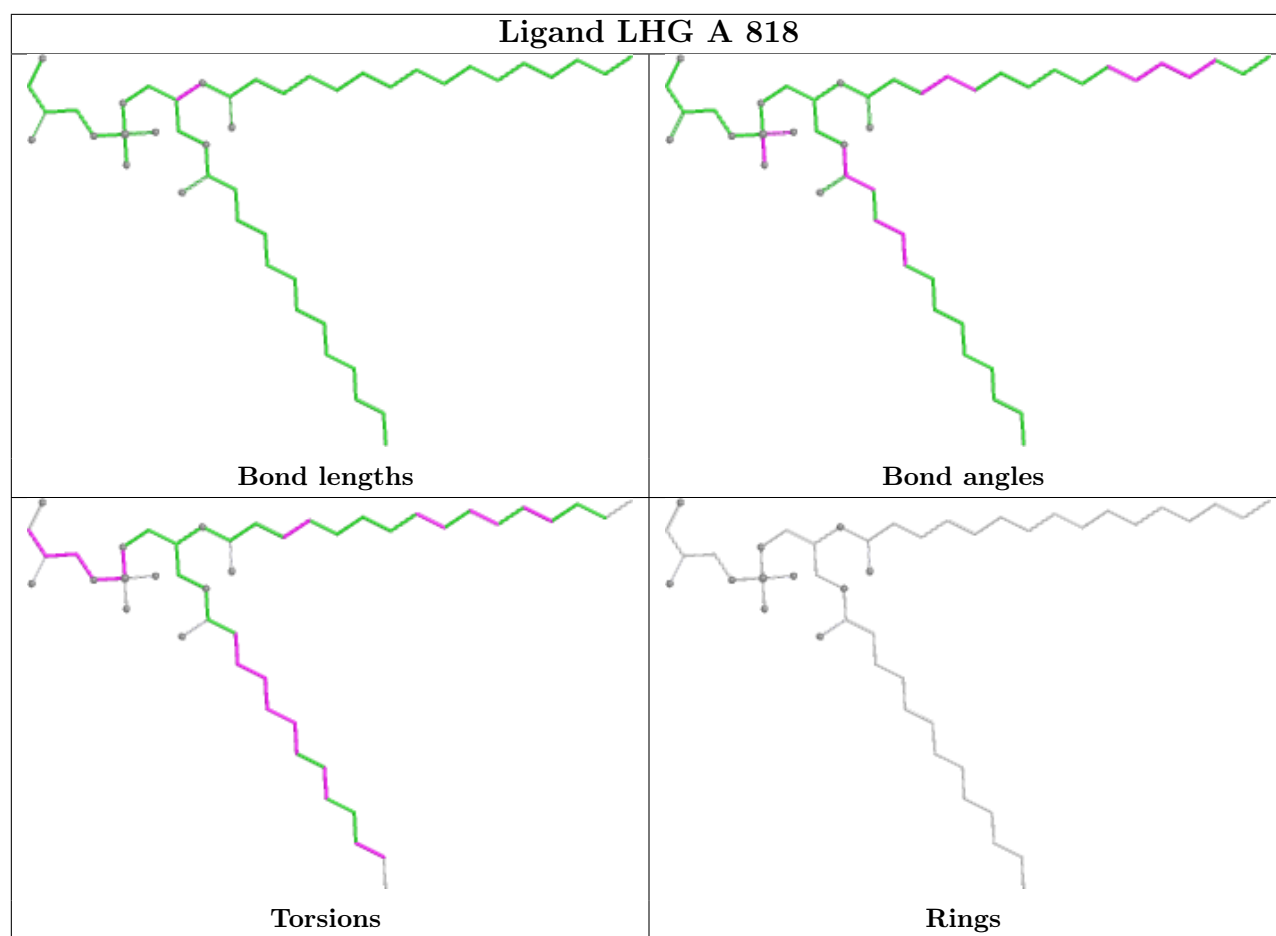
Ligand F39 a 817



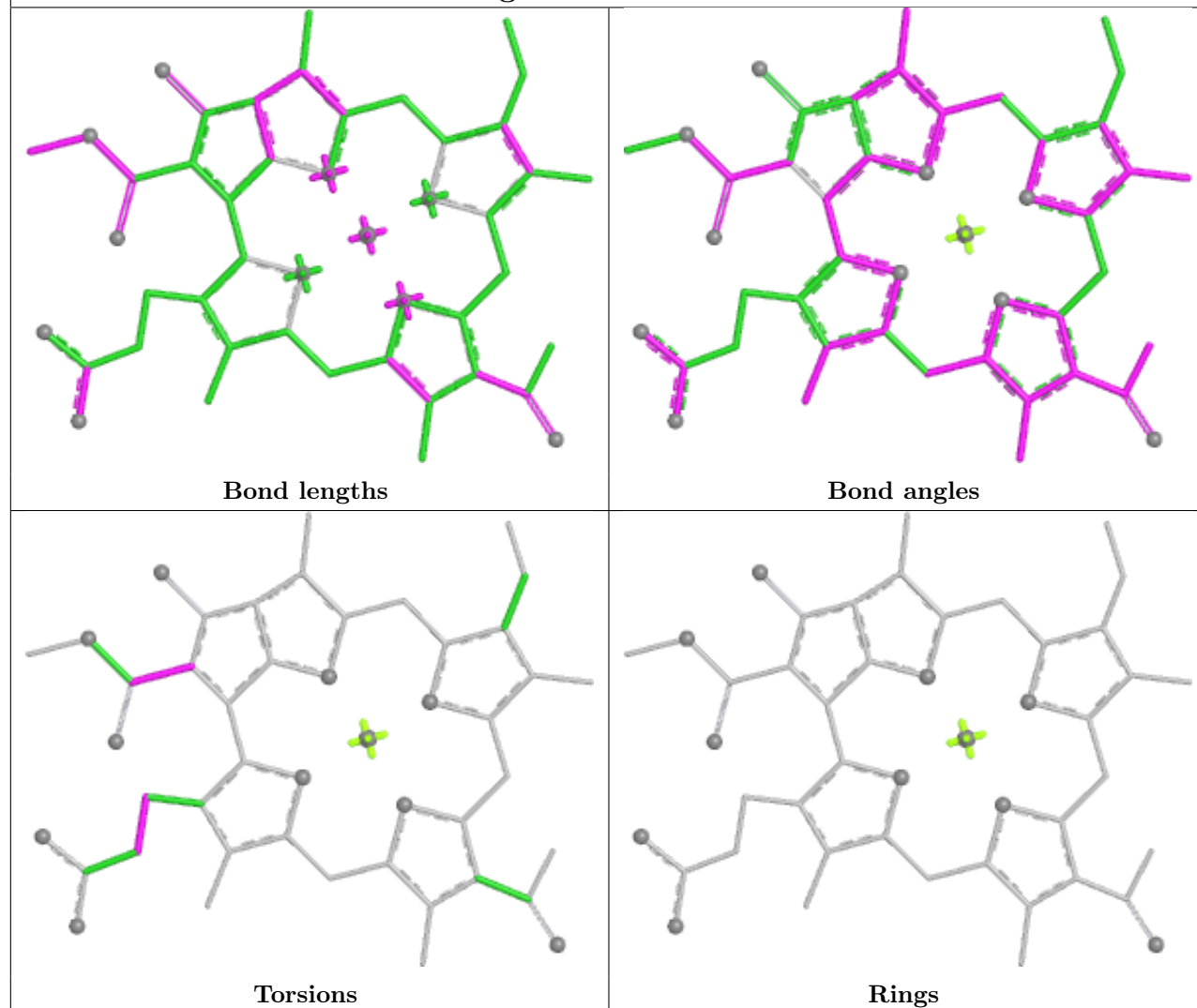




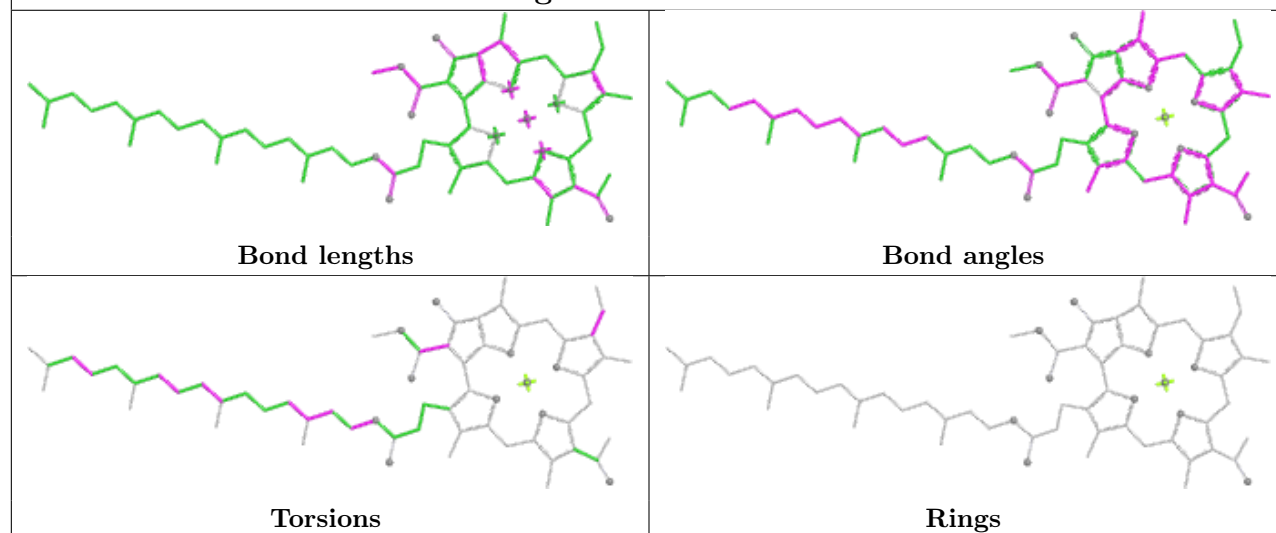


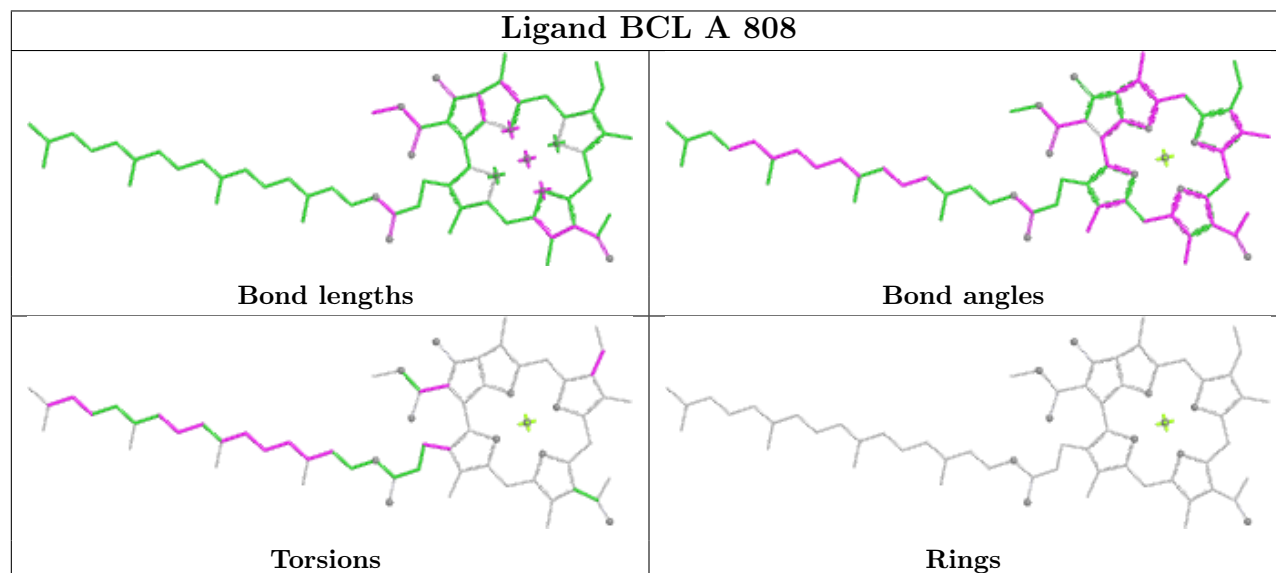
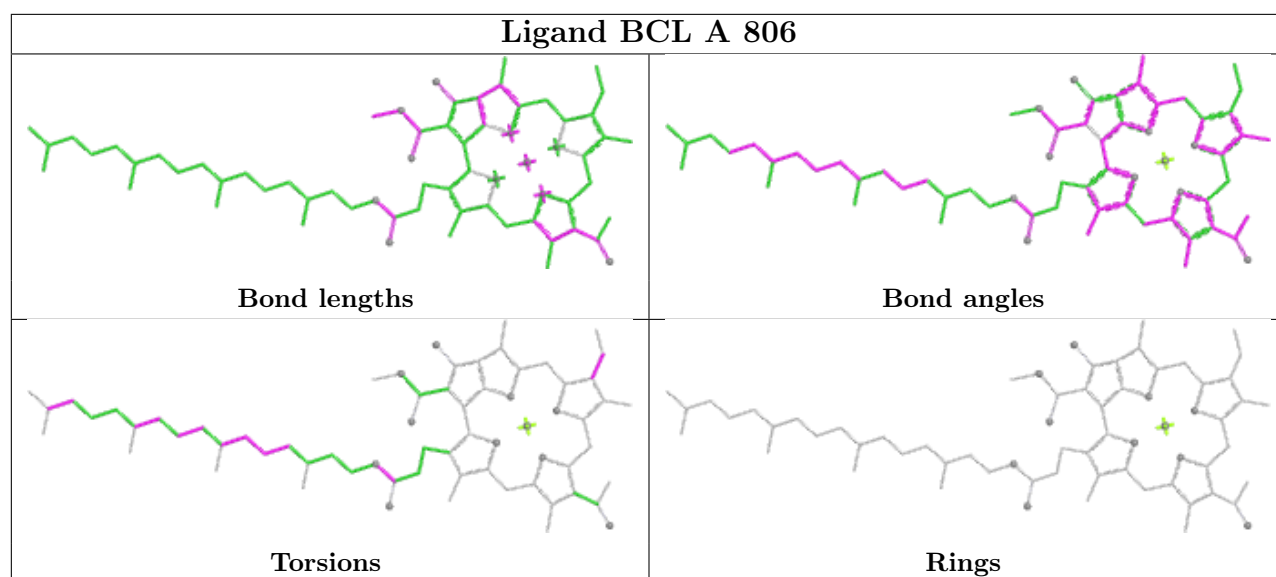


Ligand BCL A 814

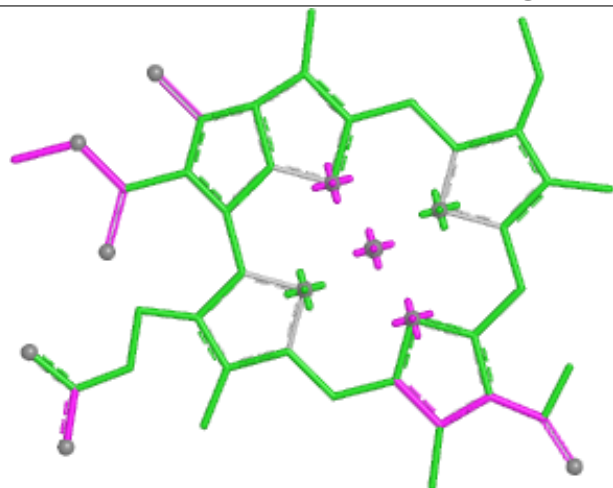


Ligand BCL F 377

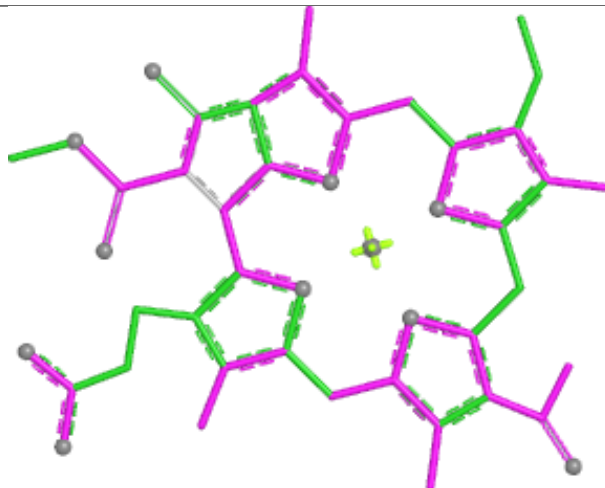




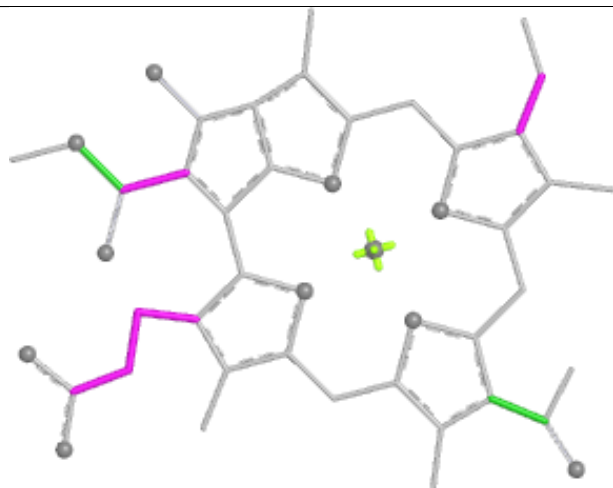
Ligand BCL a 810



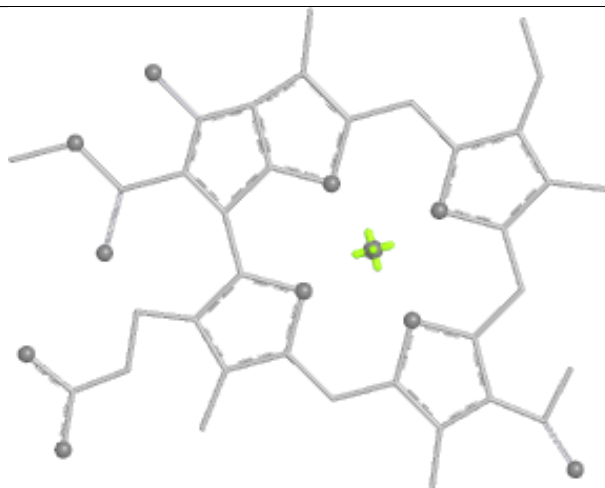
Bond lengths



Bond angles

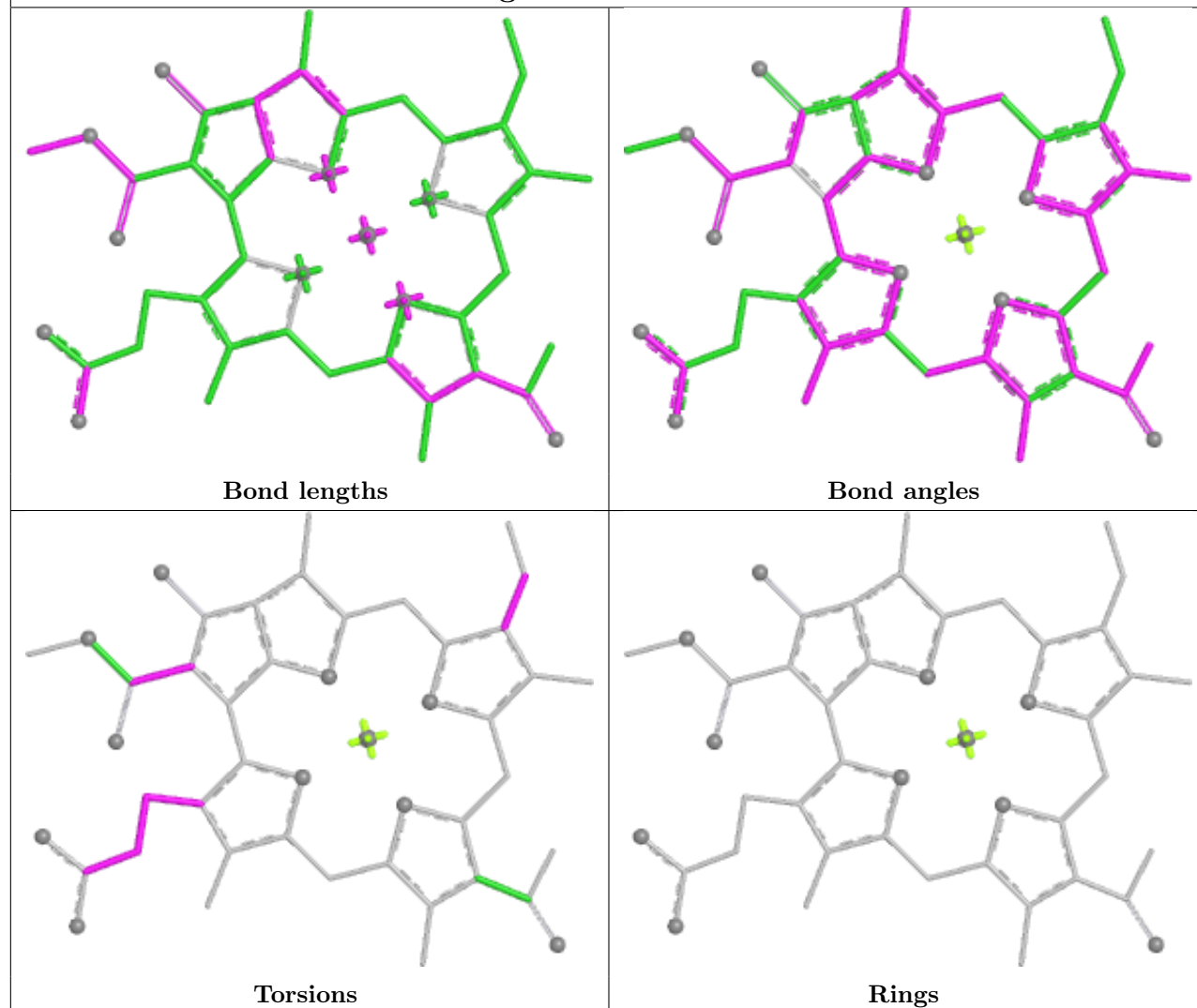


Torsions

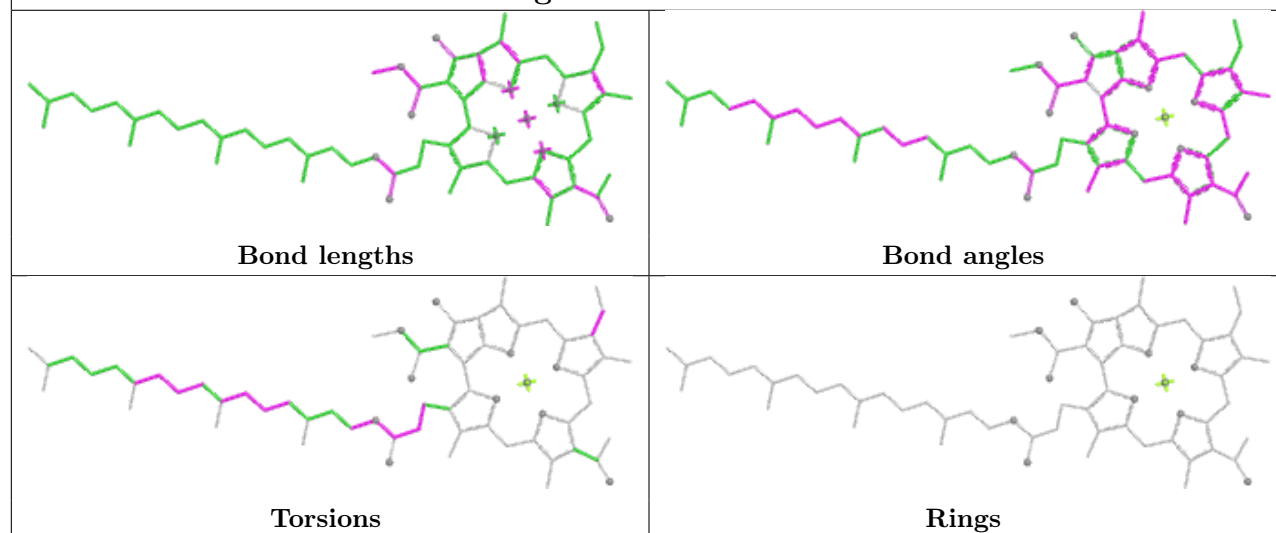


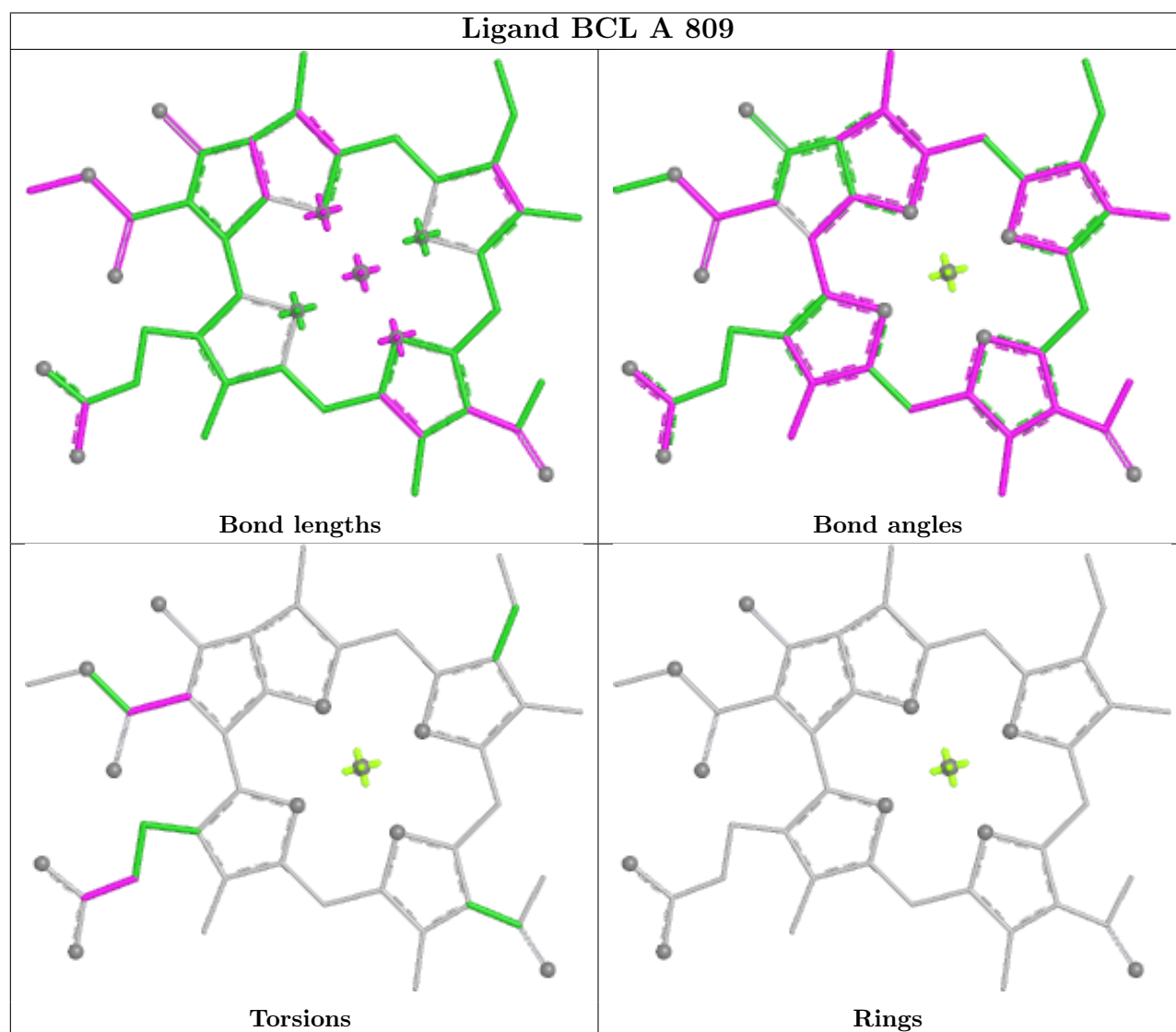
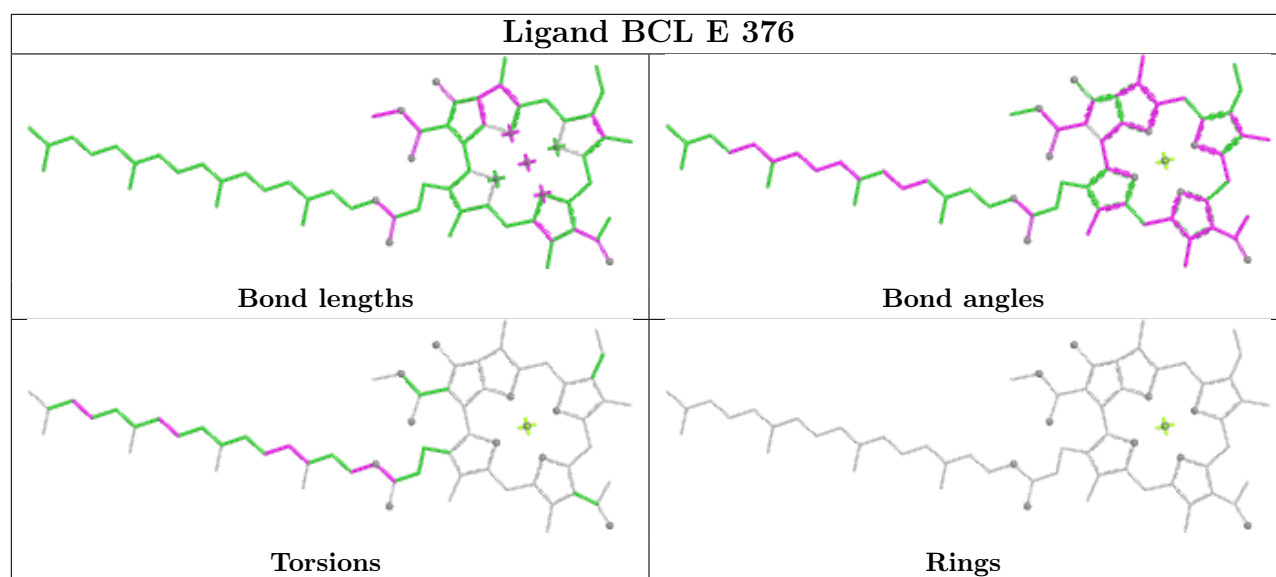
Rings

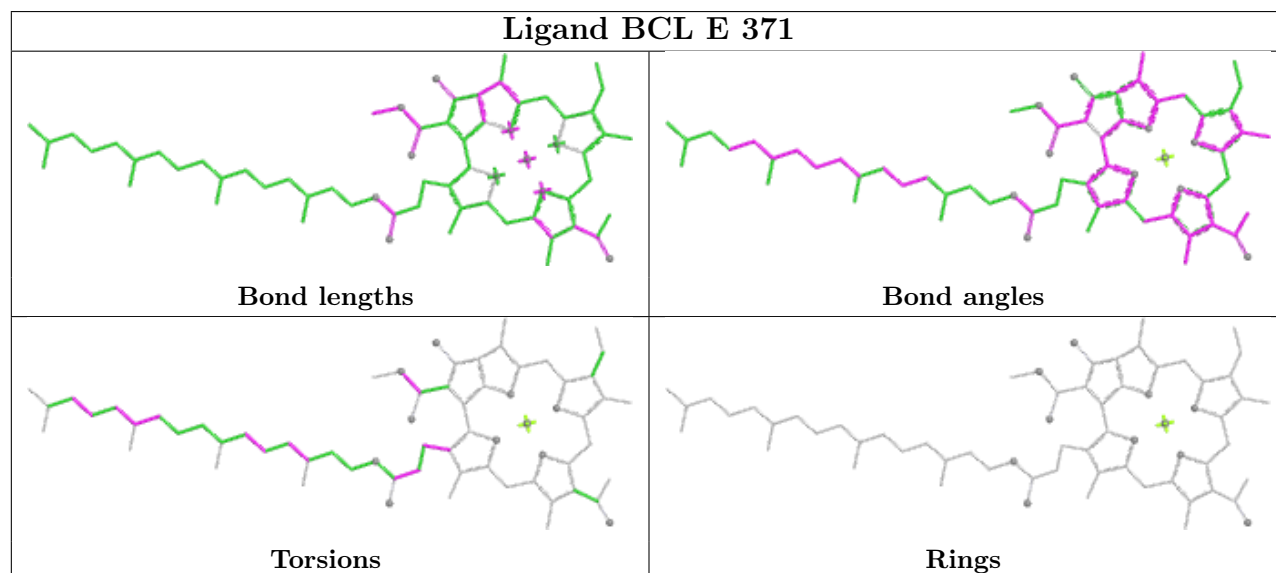
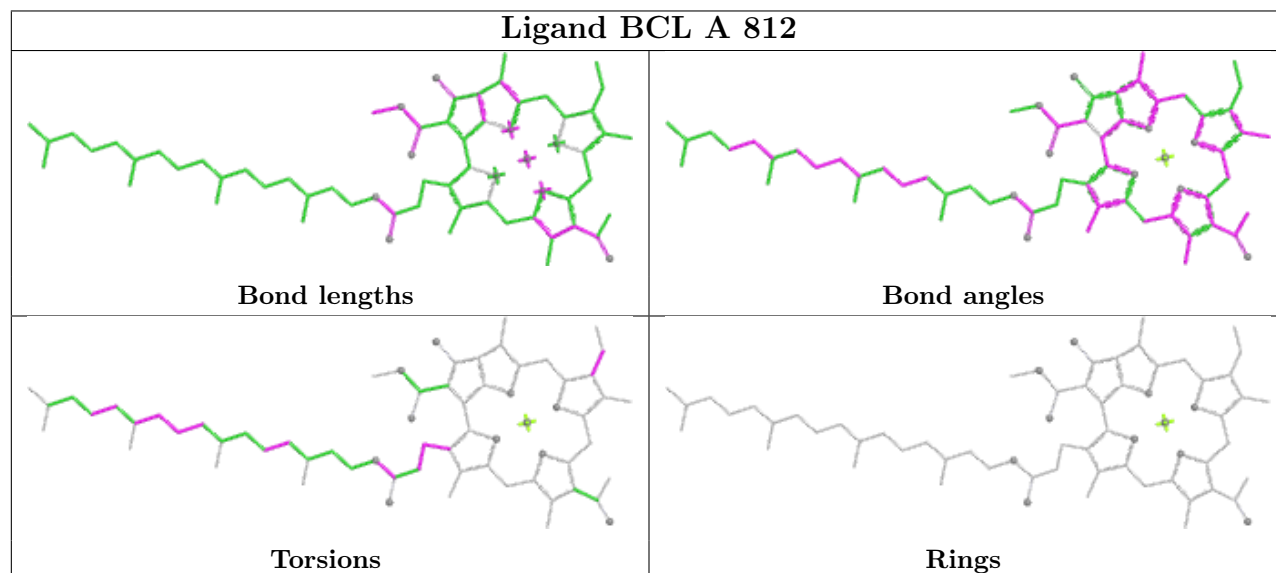
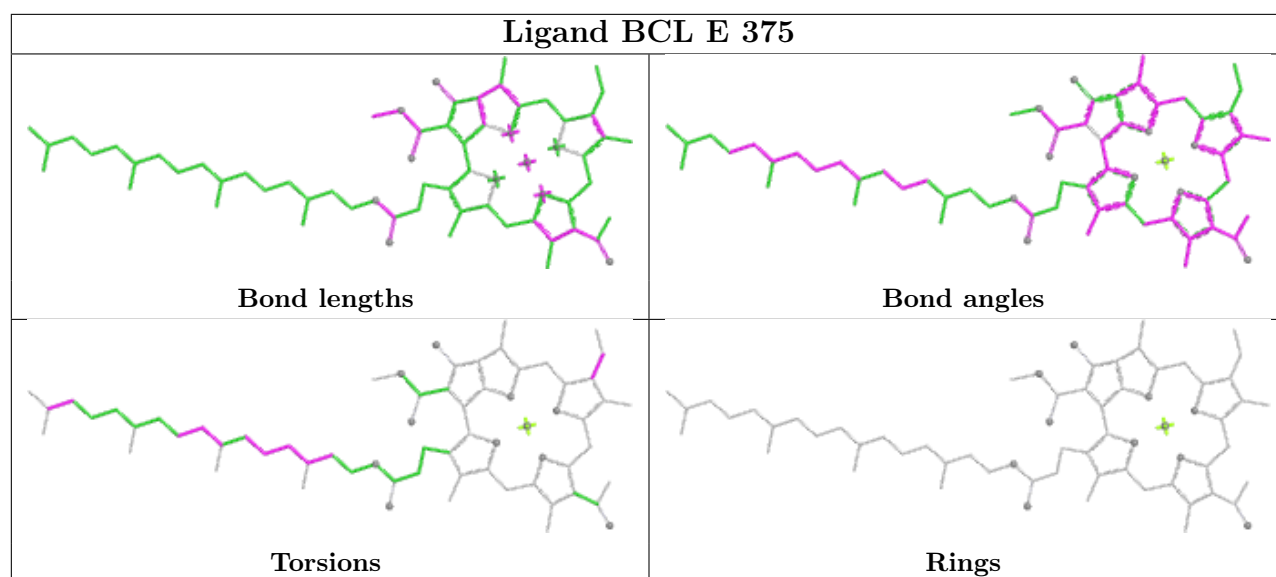
Ligand BCL A 804



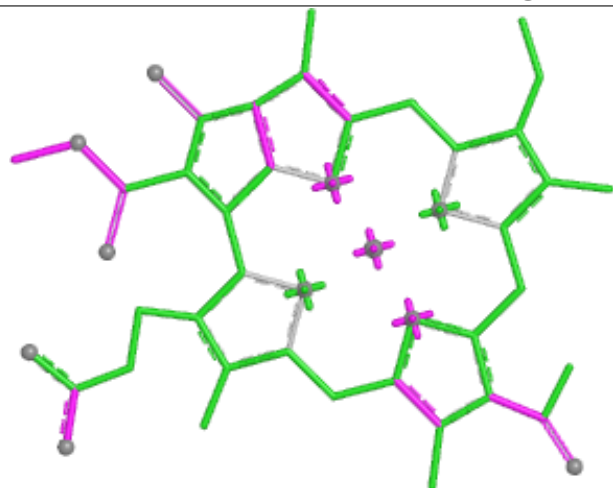
Ligand BCL E 374



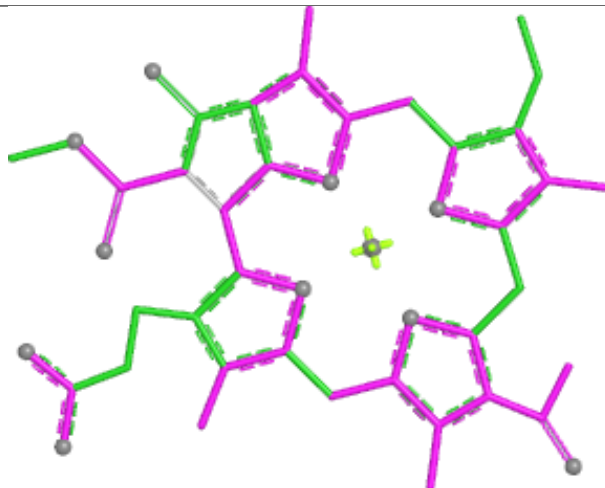




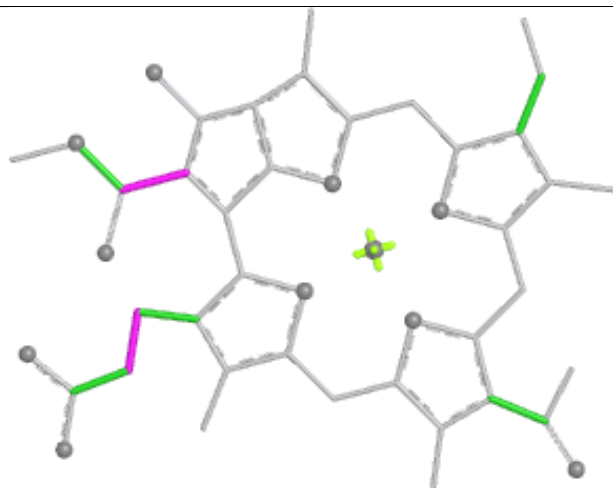
Ligand BCL a 814



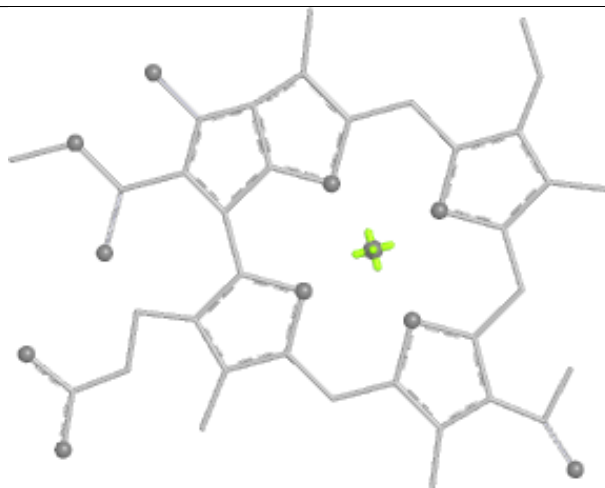
Bond lengths



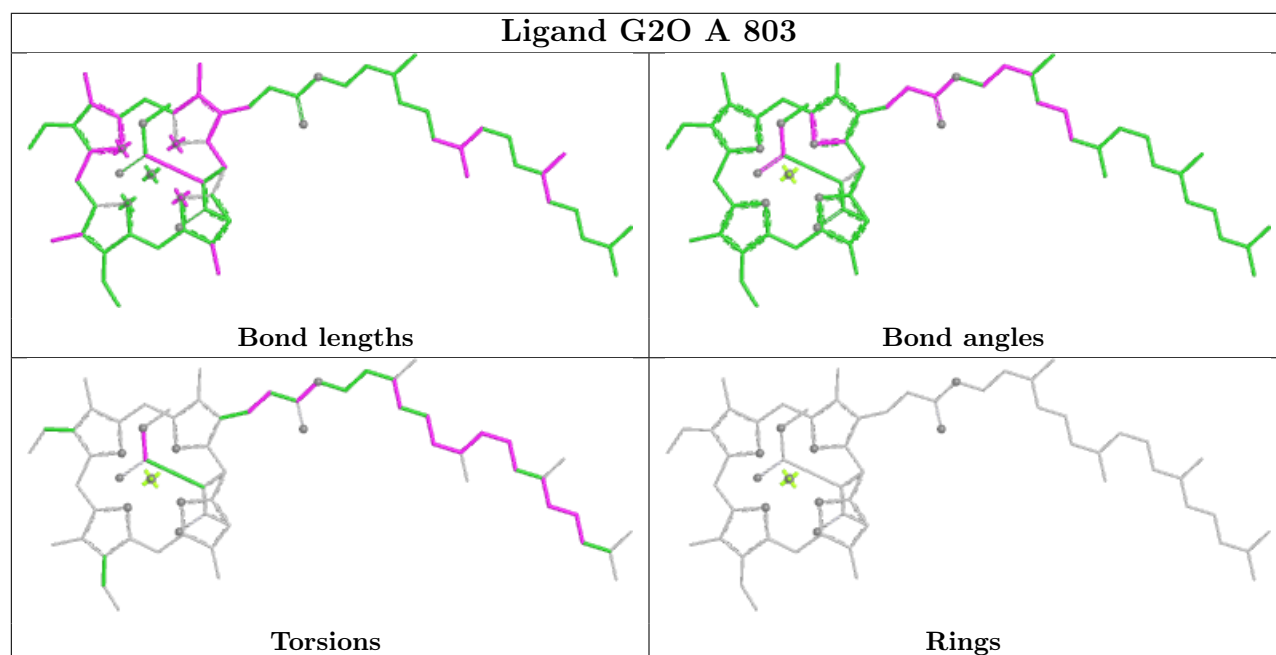
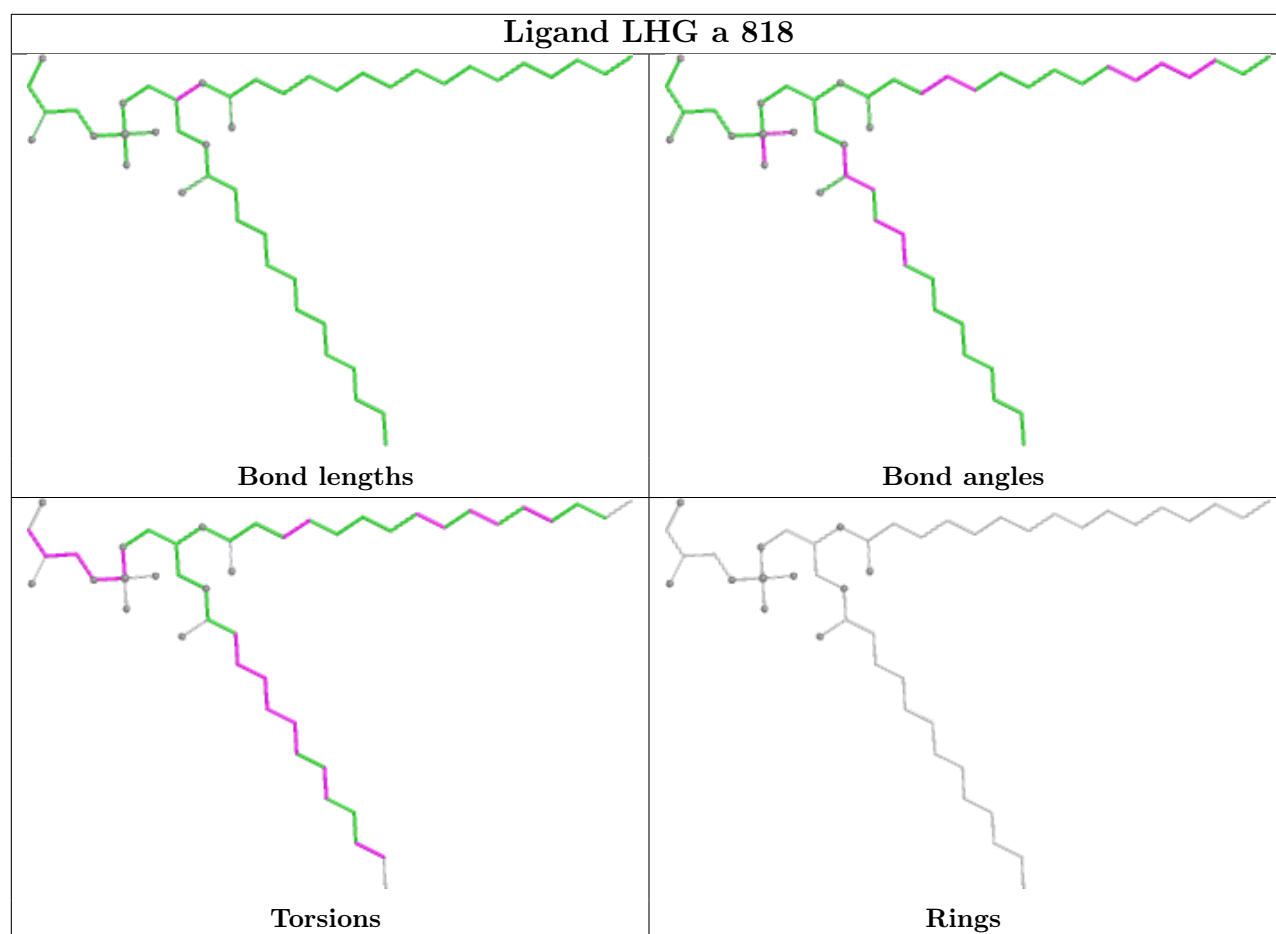
Bond angles

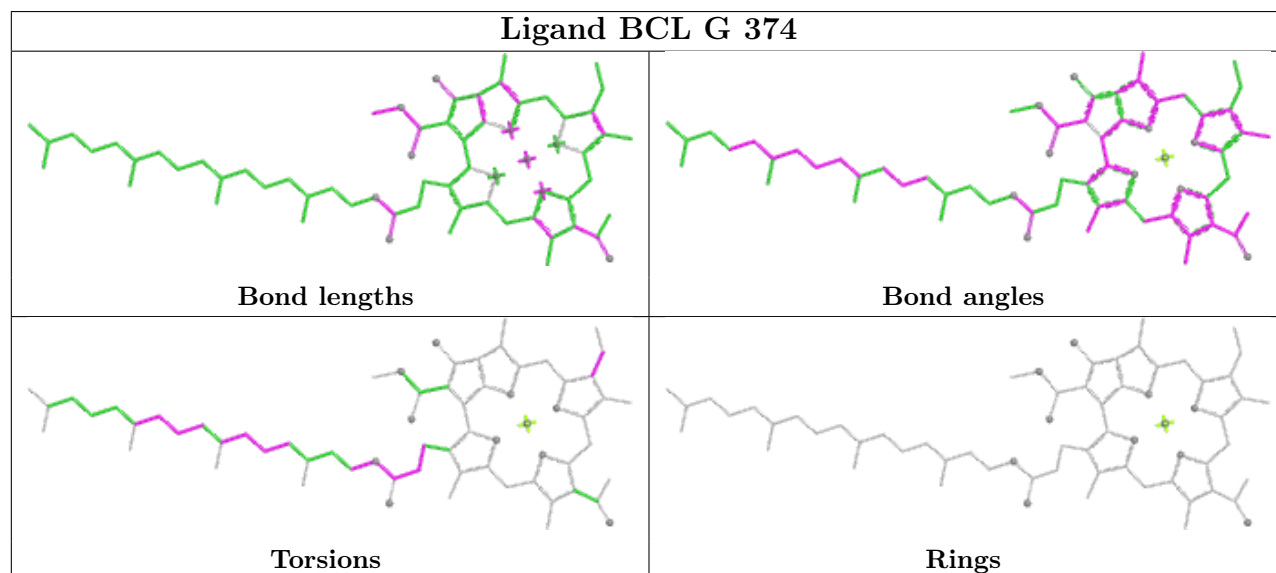
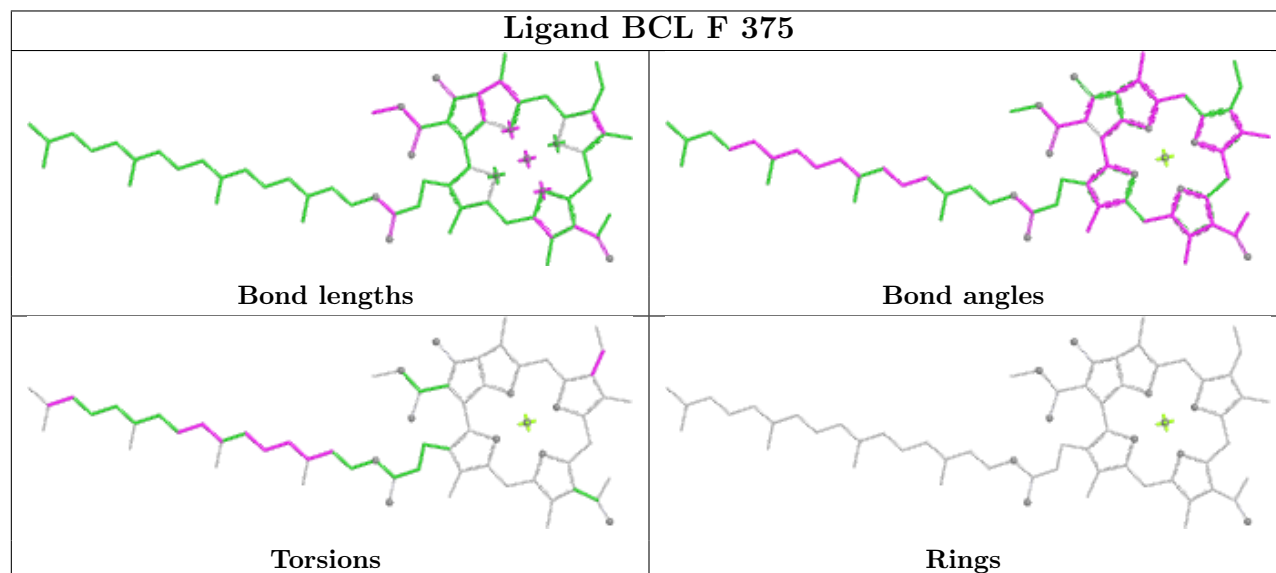
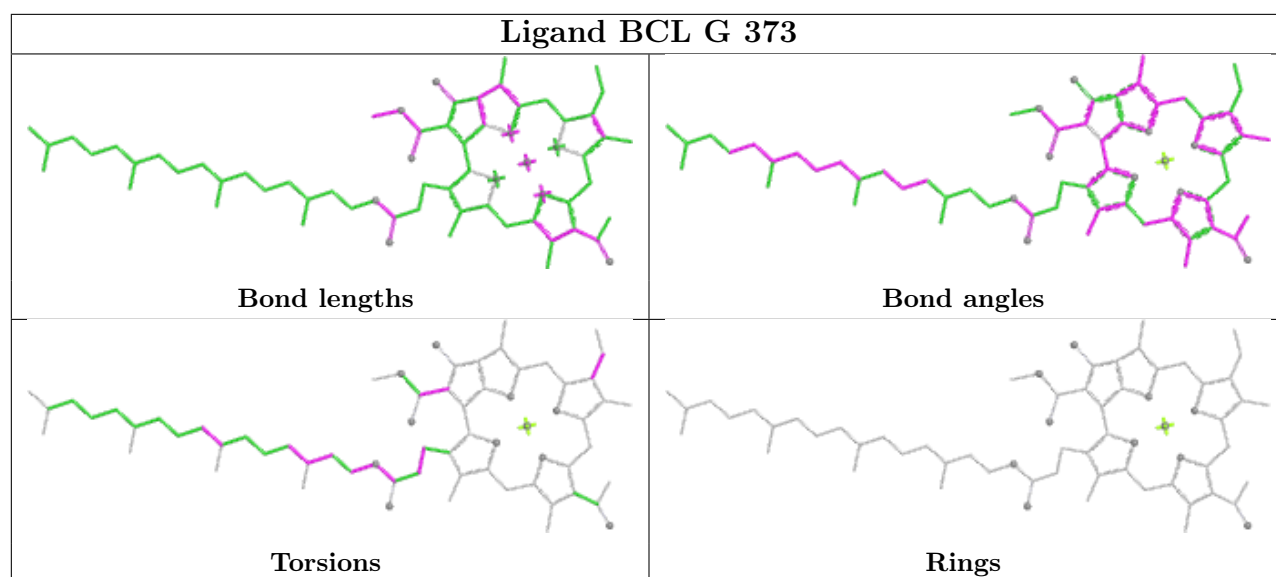


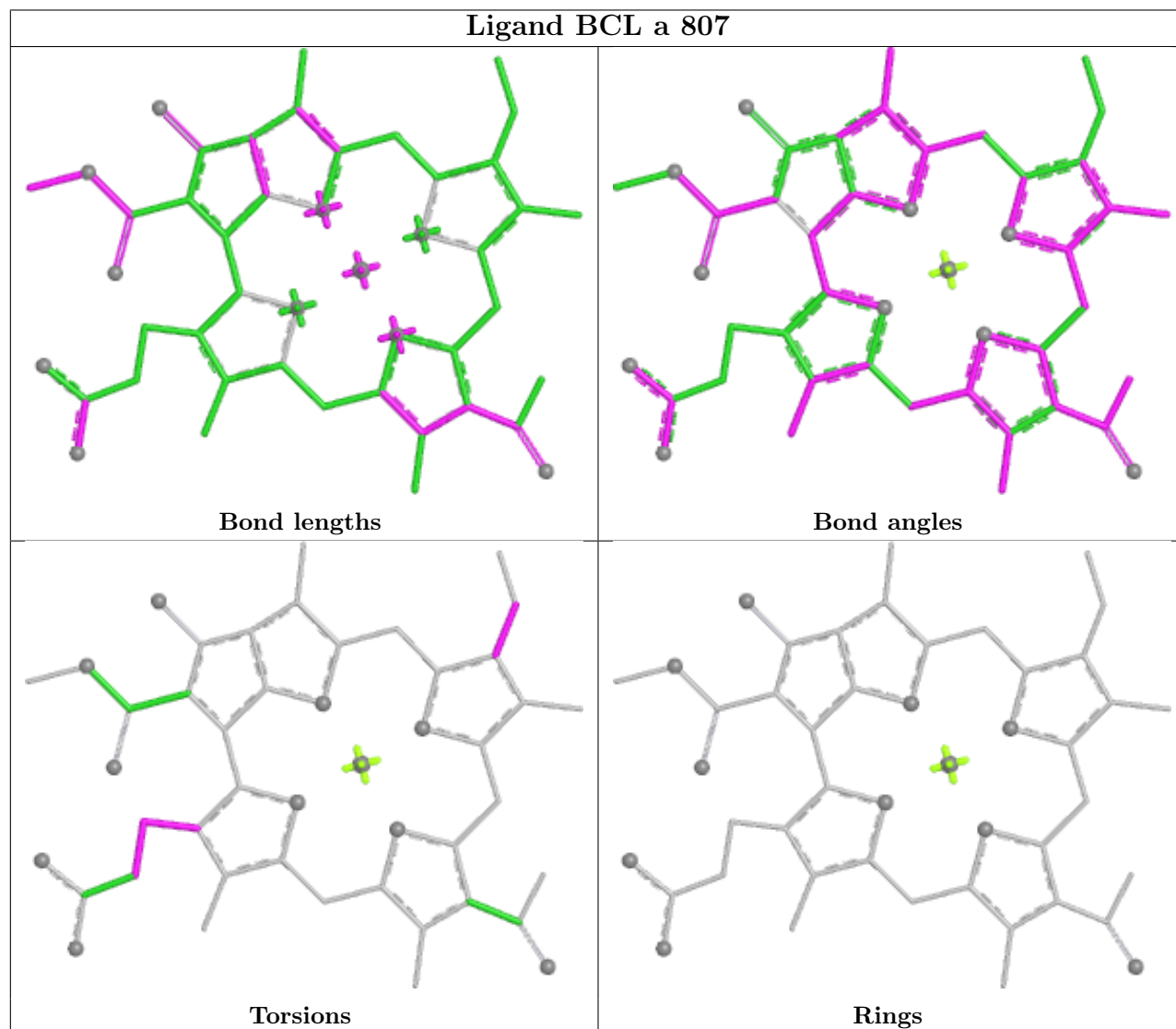
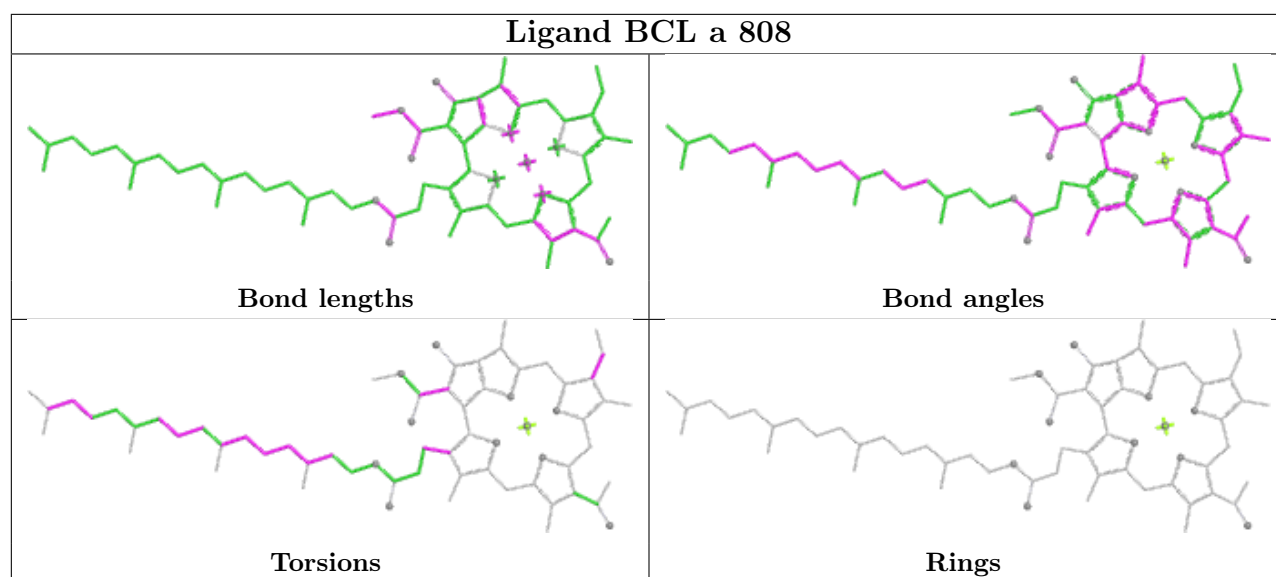
Torsions

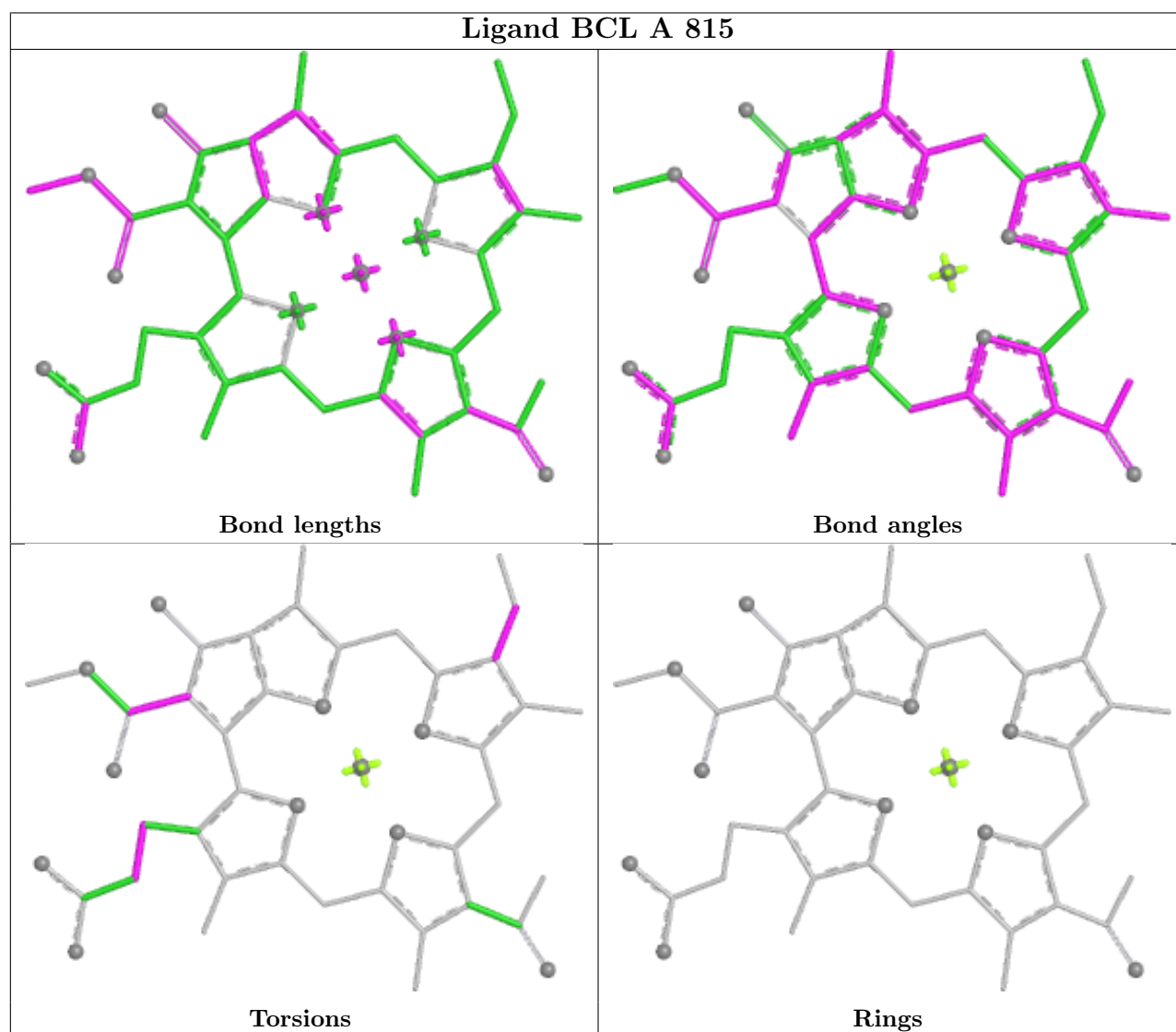
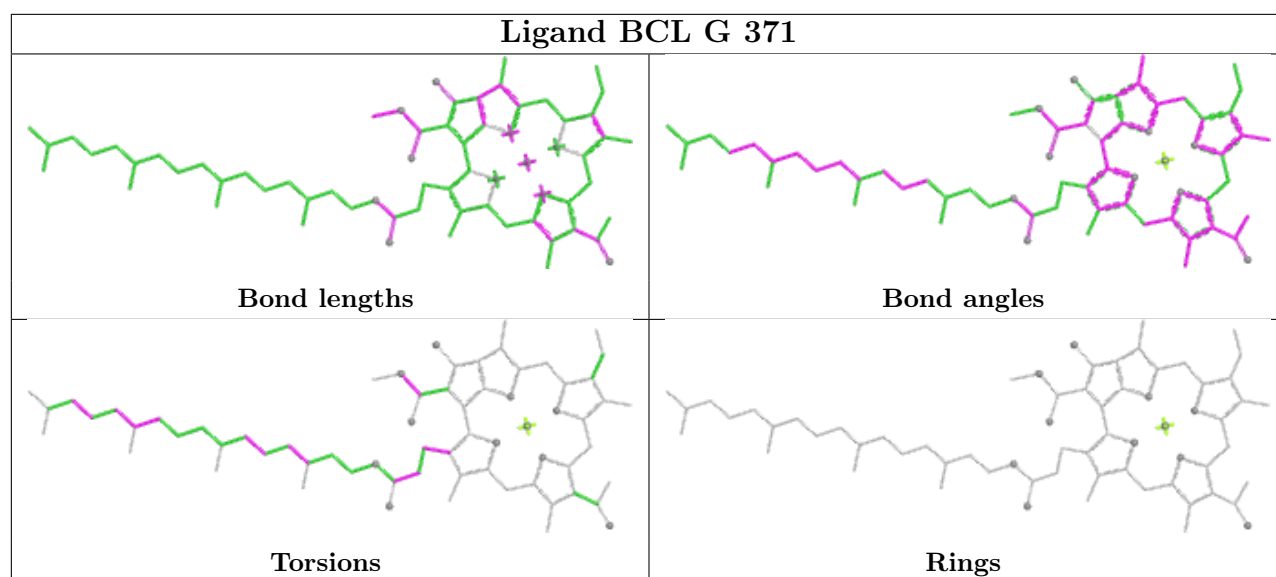


Rings

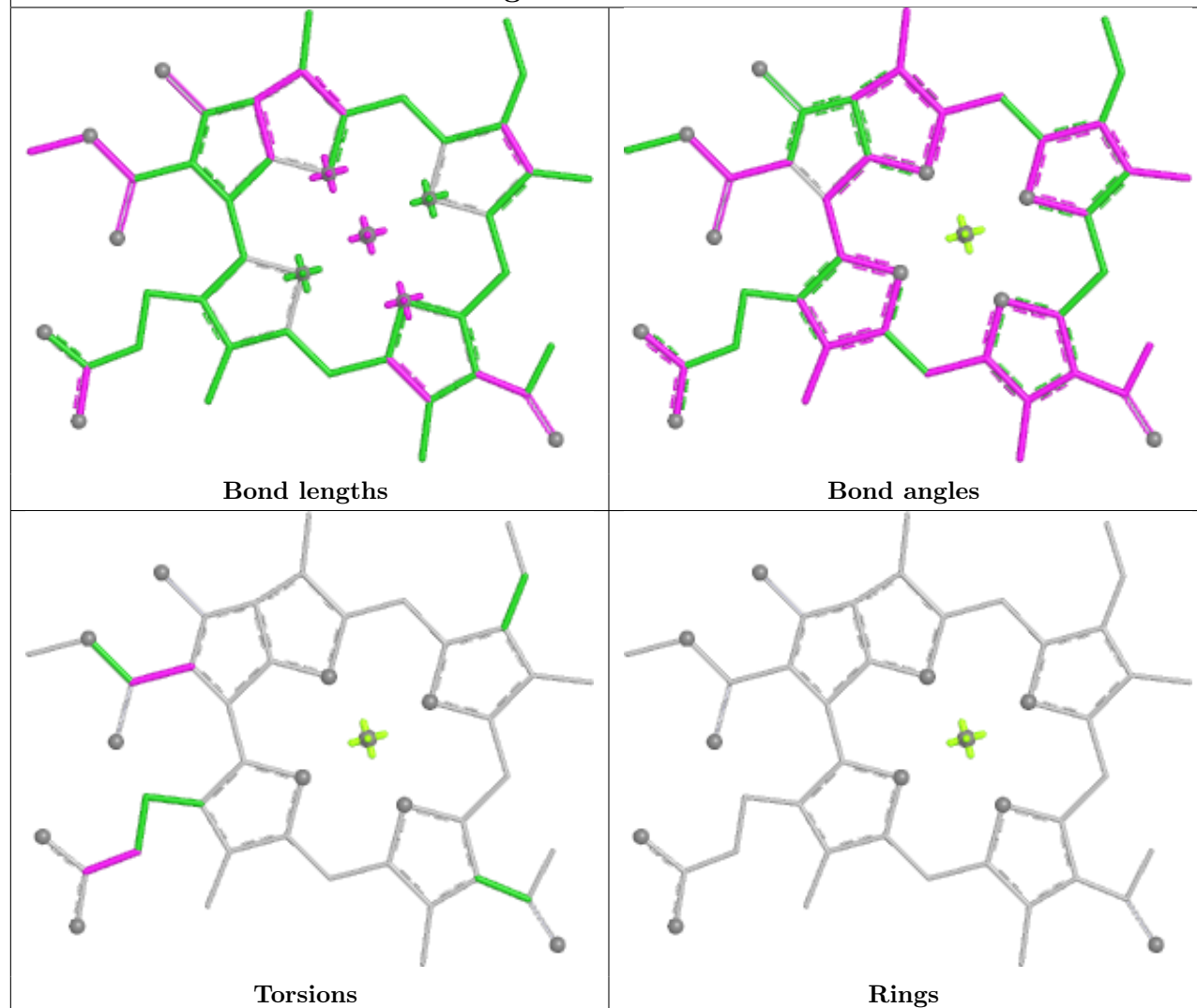




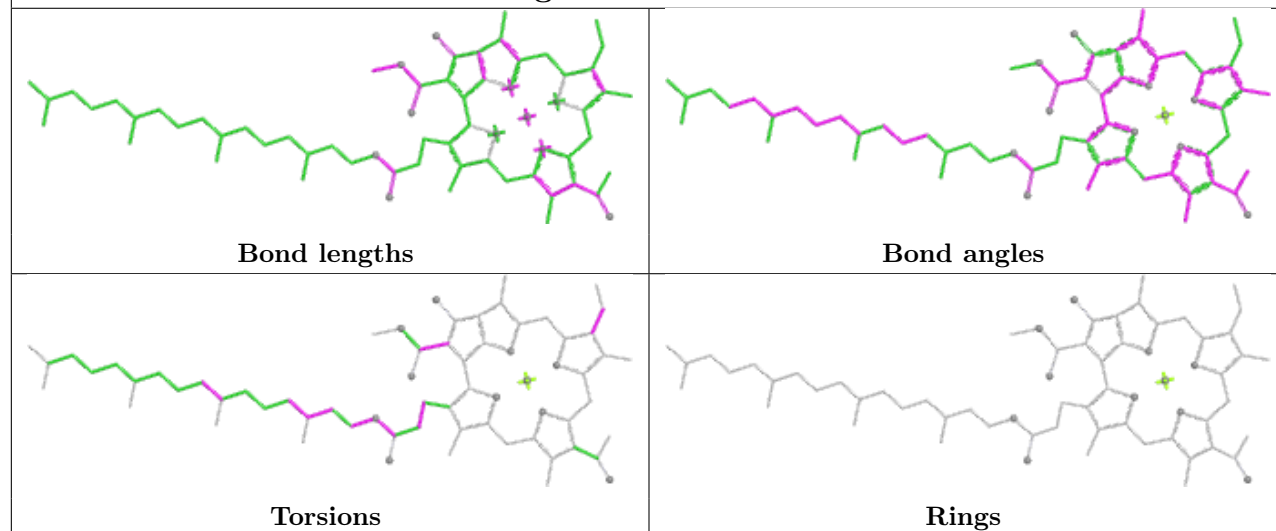


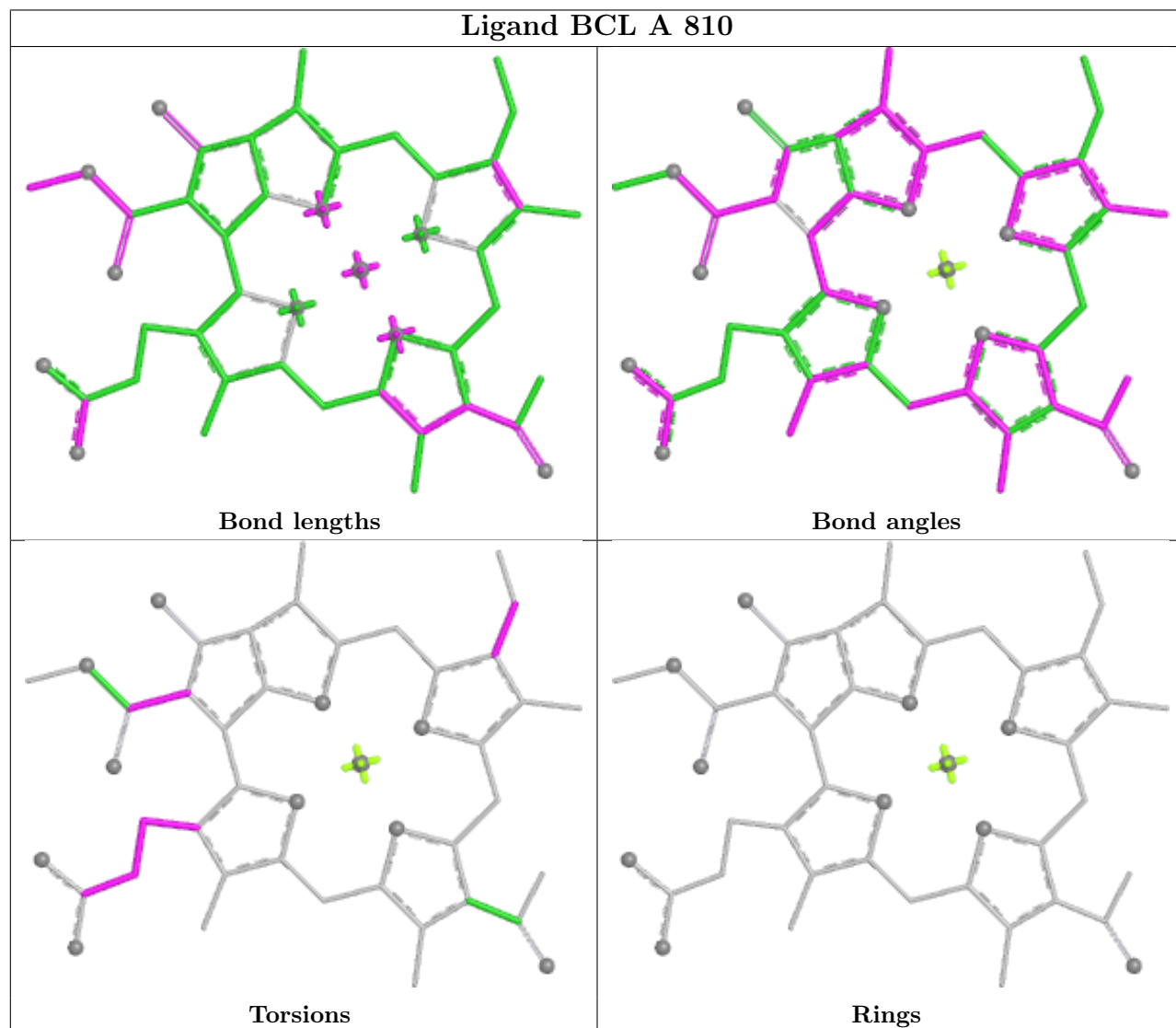
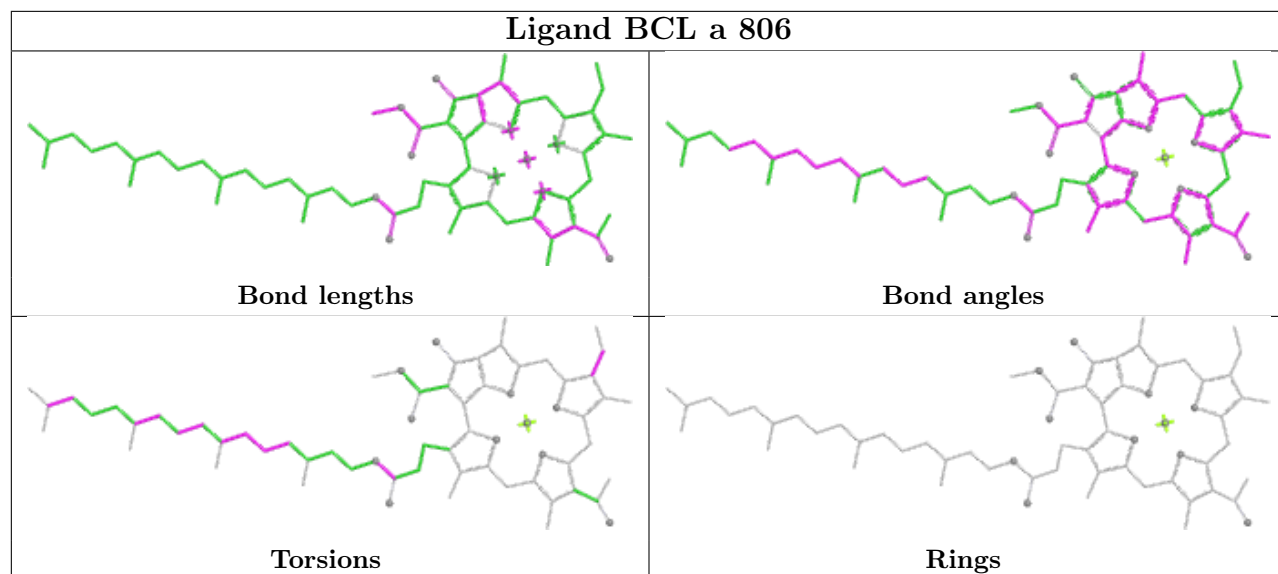


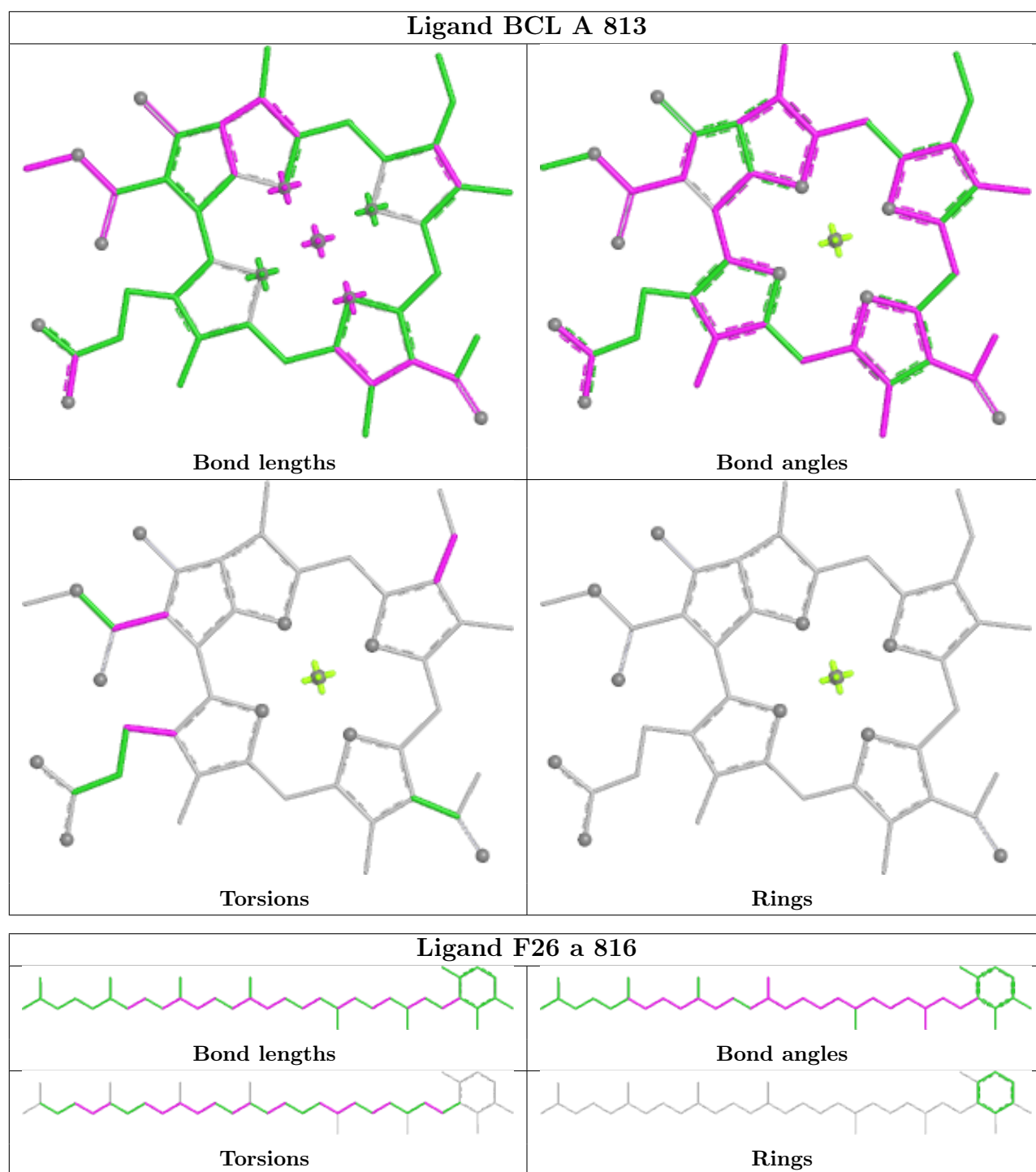
Ligand BCL a 809



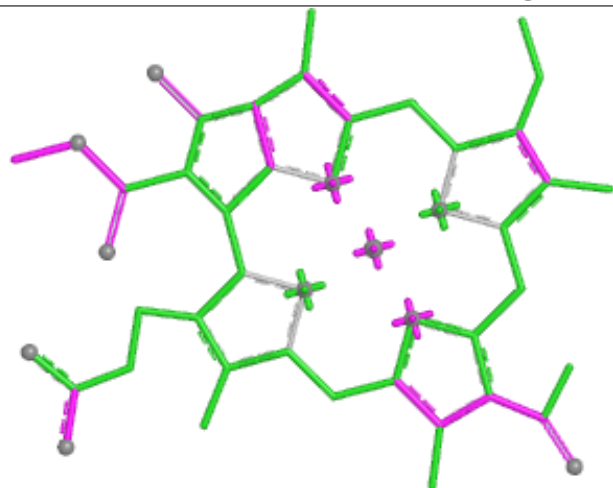
Ligand BCL F 373



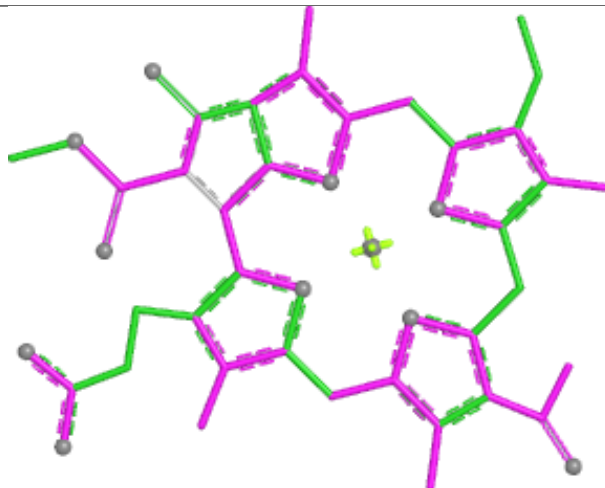




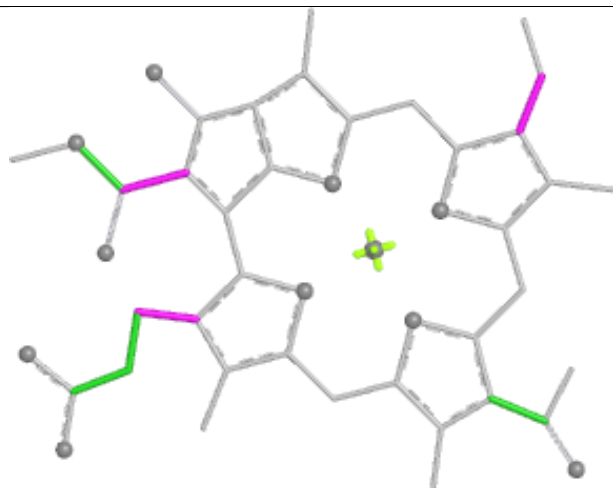
Ligand BCL a 813



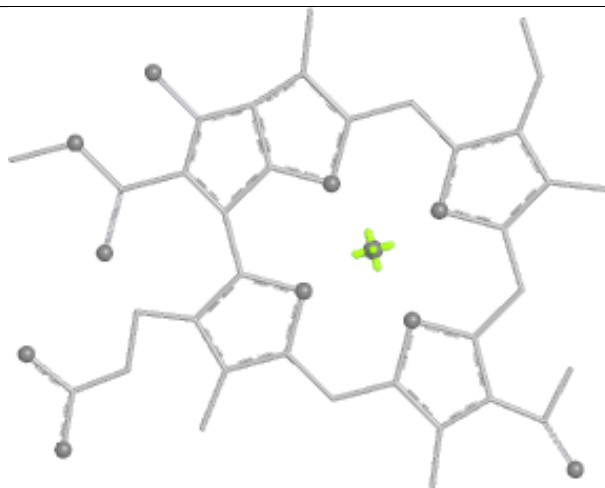
Bond lengths



Bond angles

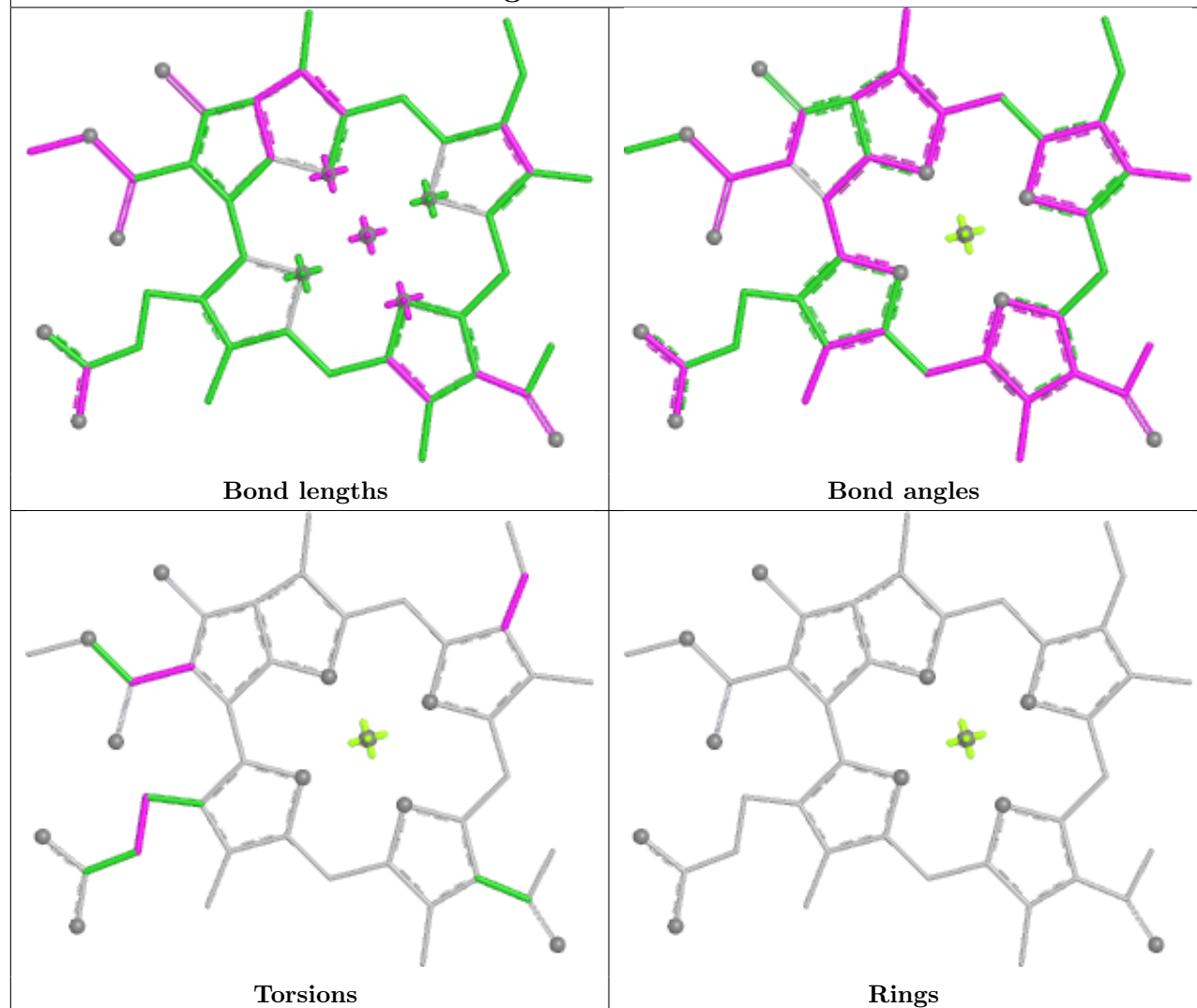


Torsions

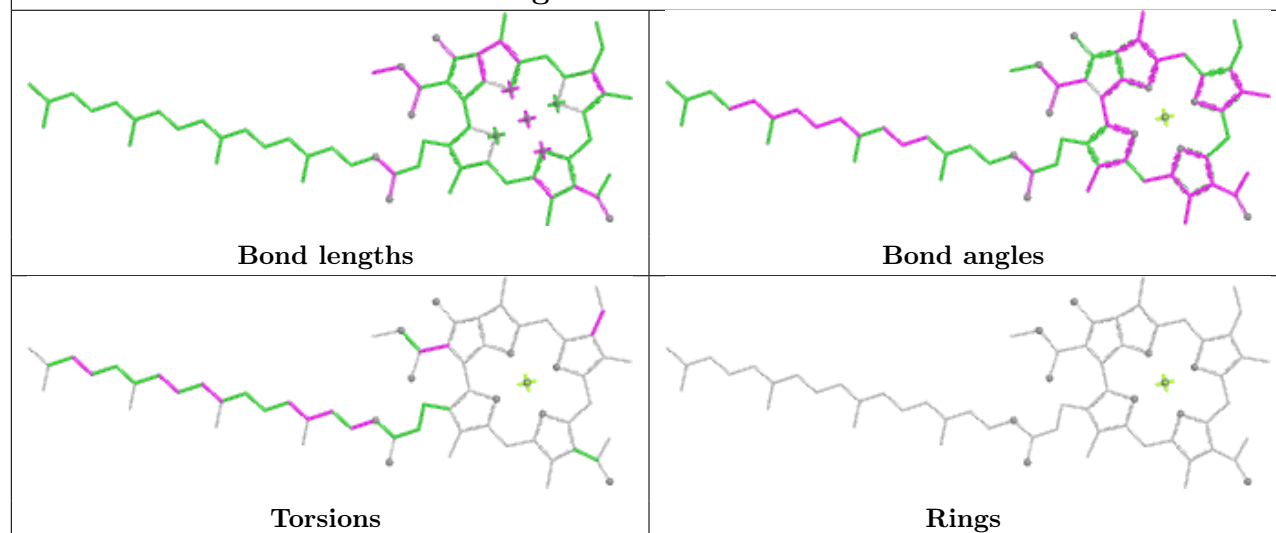


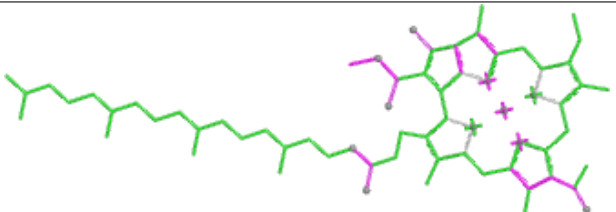
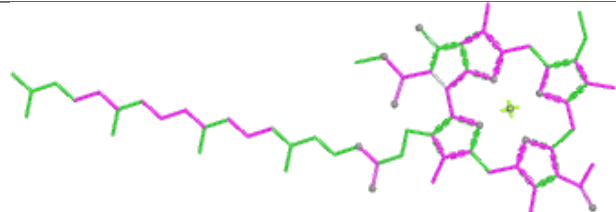
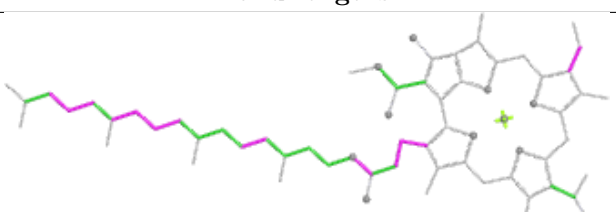
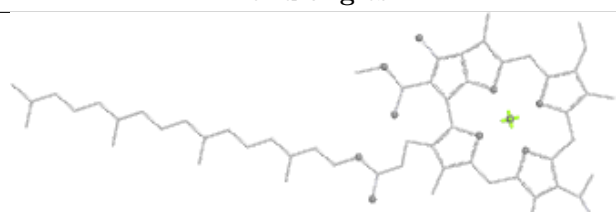
Rings

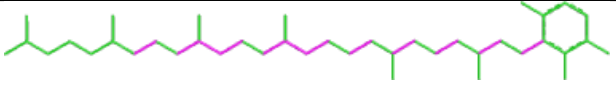
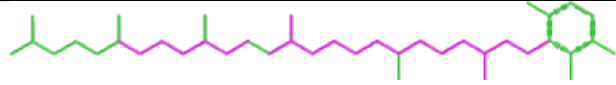
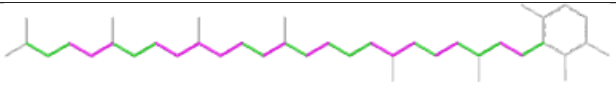
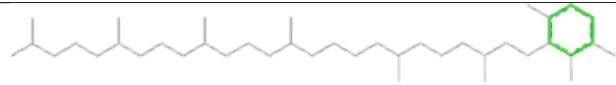
Ligand BCL a 815

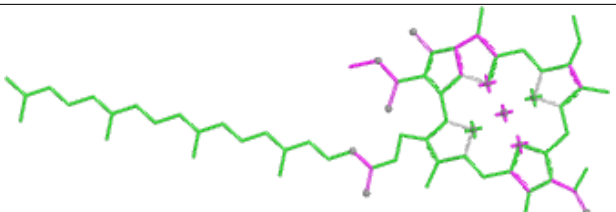
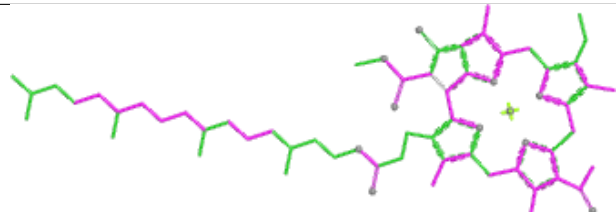
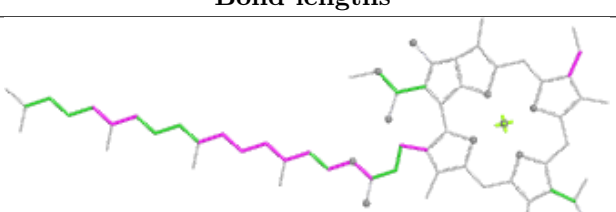
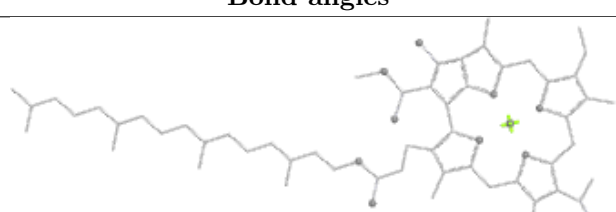


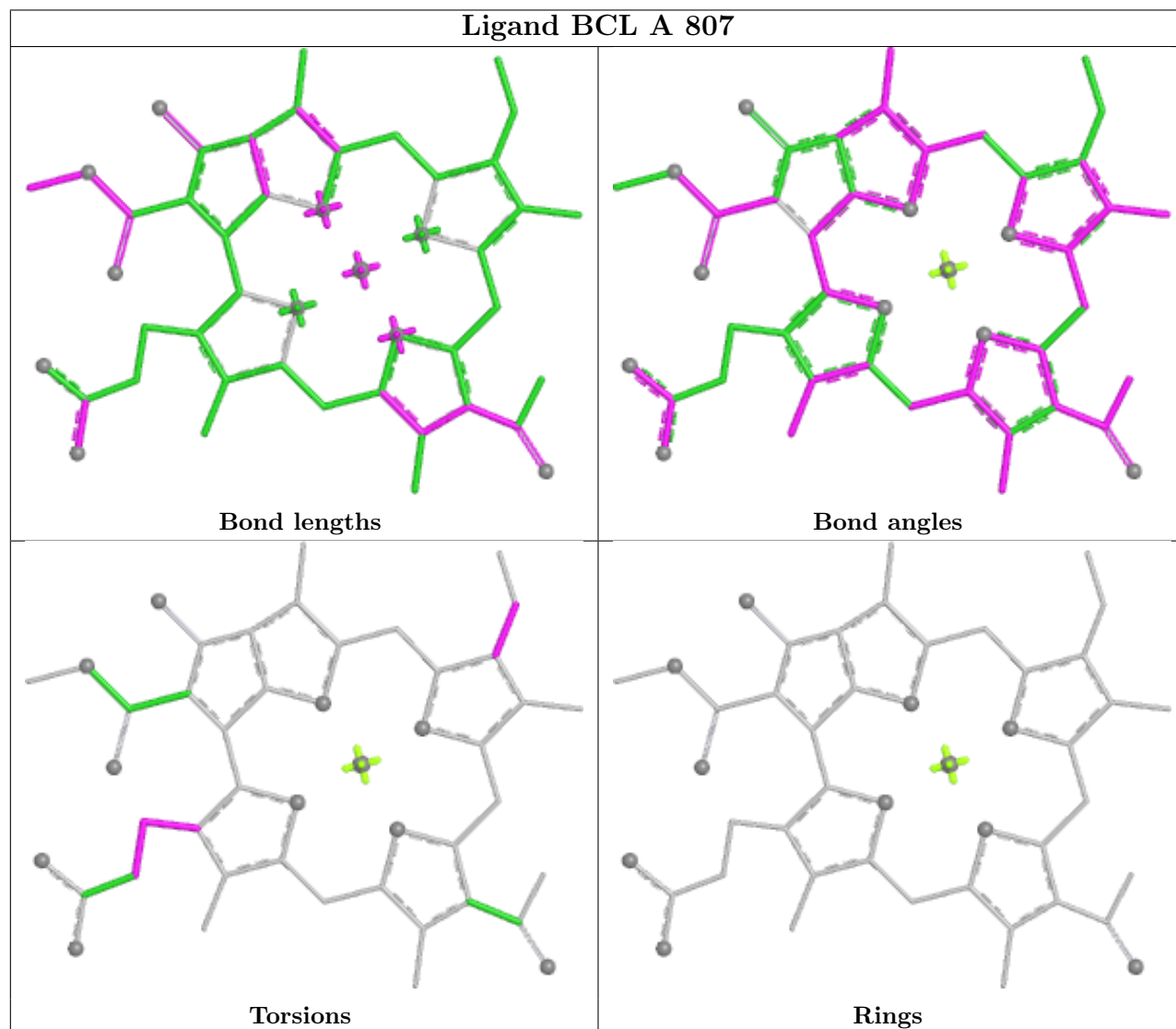
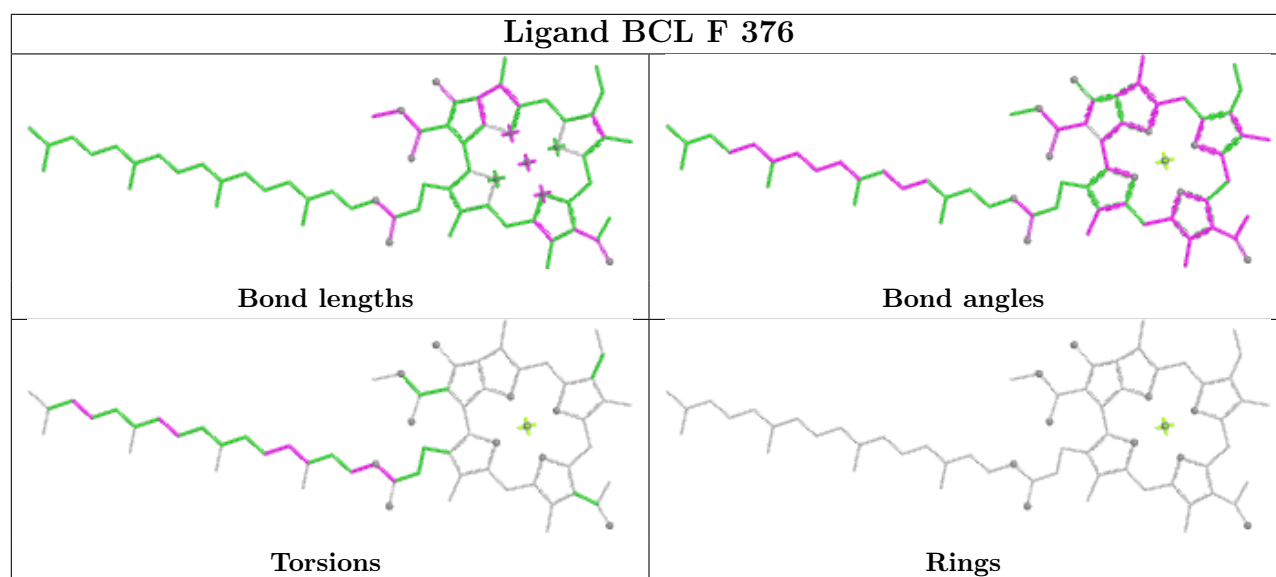
Ligand BCL G 377

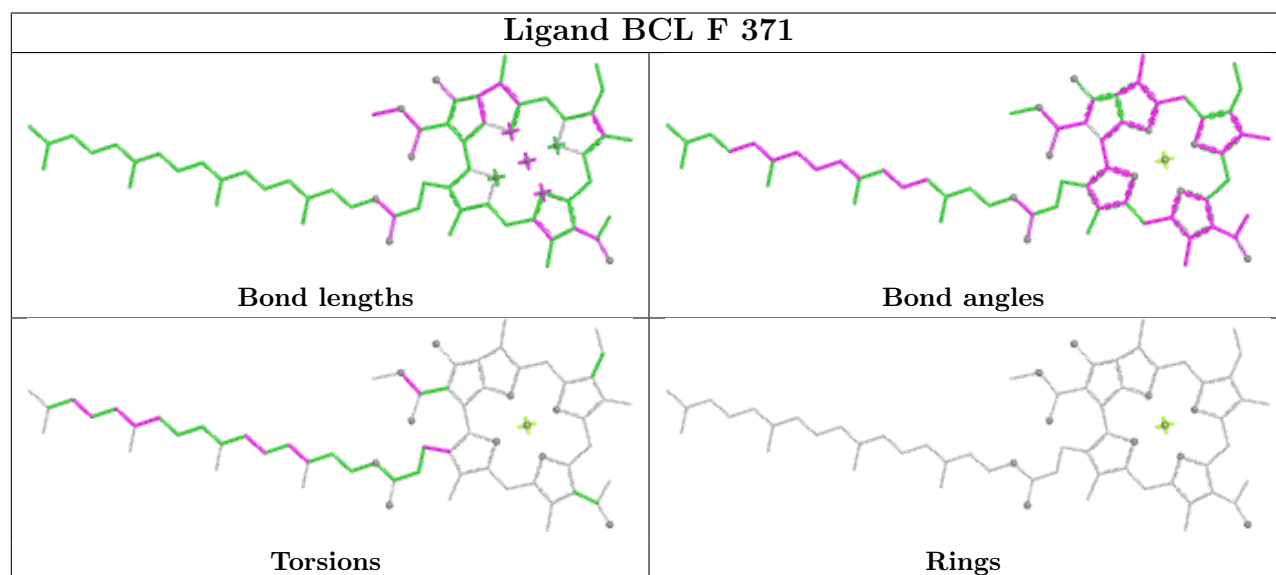
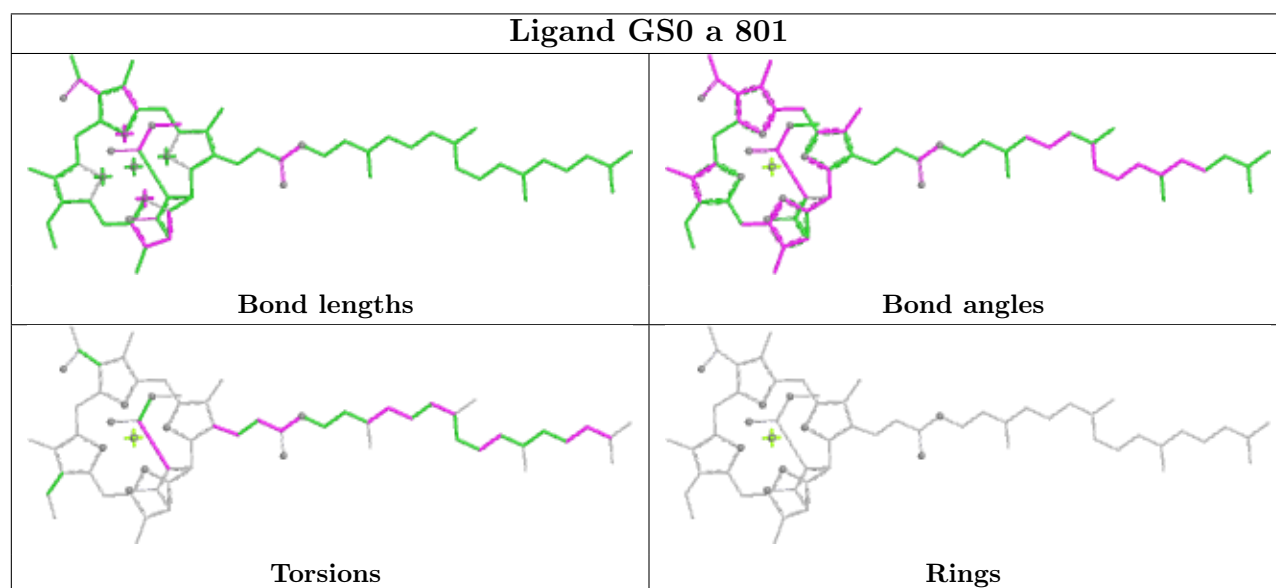
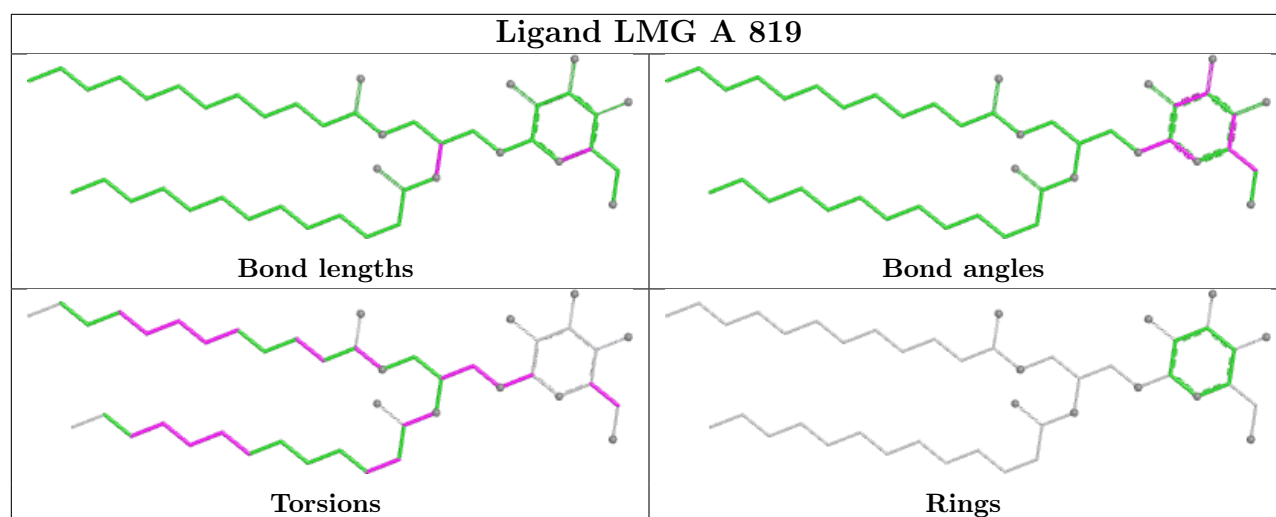


Ligand BCL a 812	
	
Bond lengths	Bond angles
	
Torsions	Rings

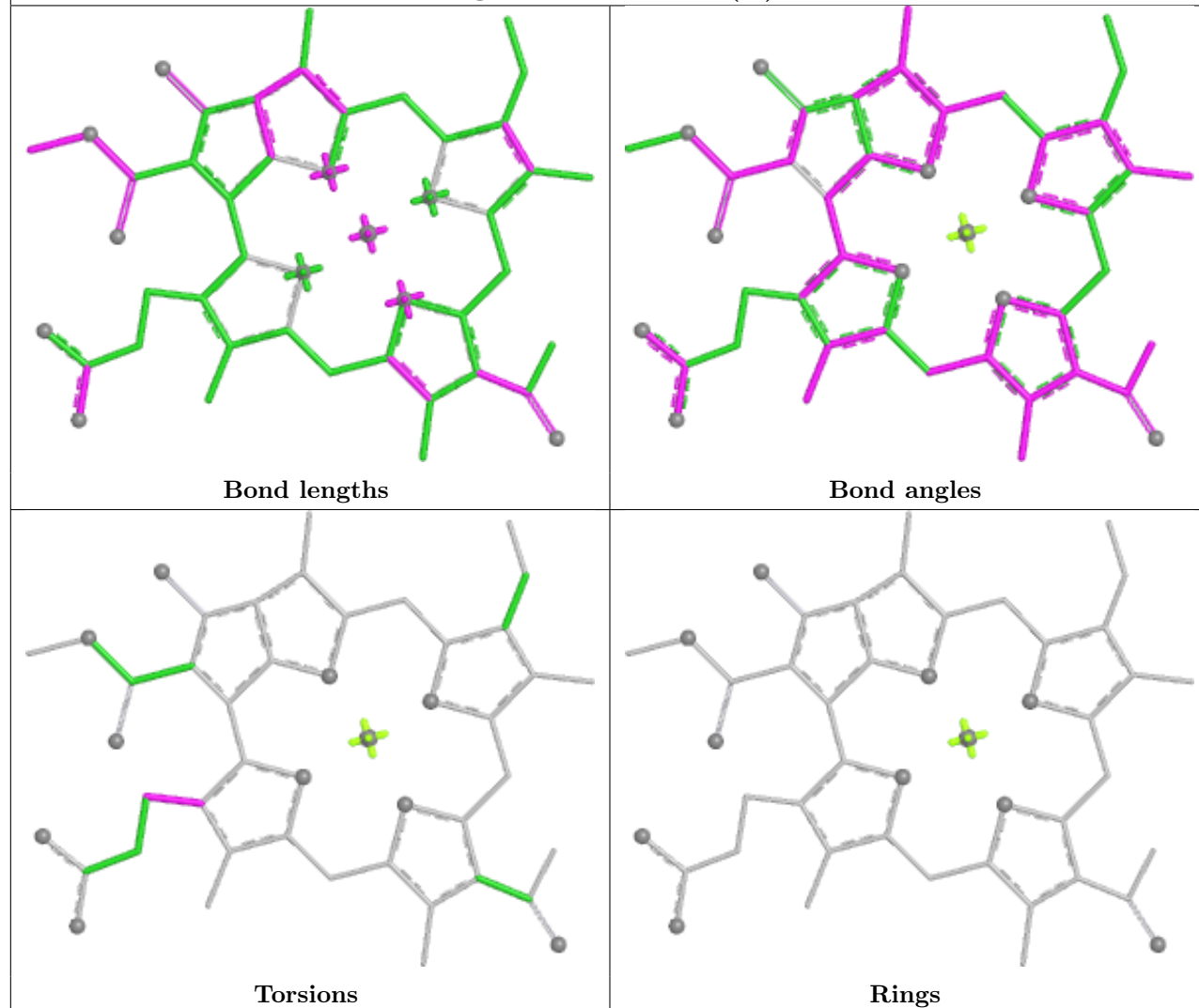
Ligand F26 A 816	
	
Bond lengths	Bond angles
	
Torsions	Rings

Ligand BCL A 811	
	
Bond lengths	Bond angles
	
Torsions	Rings

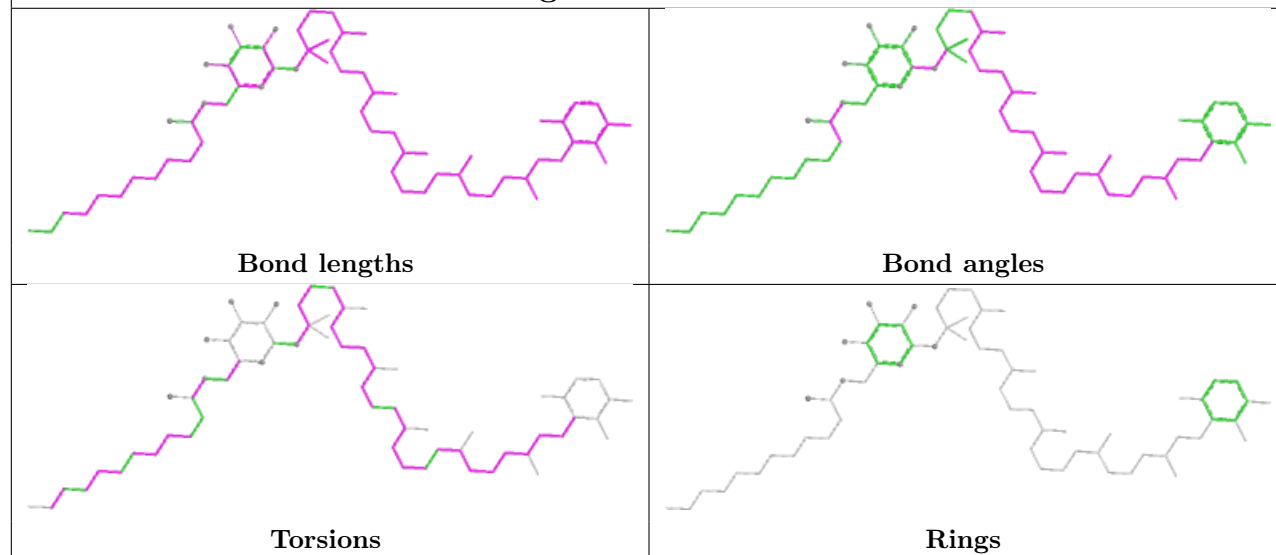




Ligand BCL F 378 (B)



Ligand F39 A 817



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.

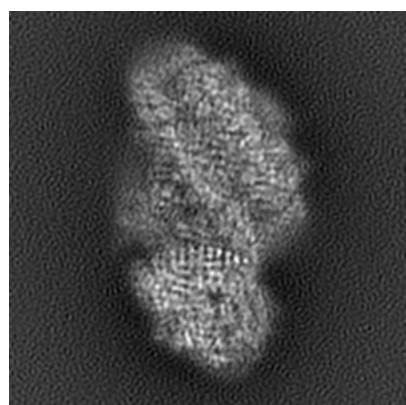
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-30069. These allow visual inspection of the internal detail of the map and identification of artifacts.

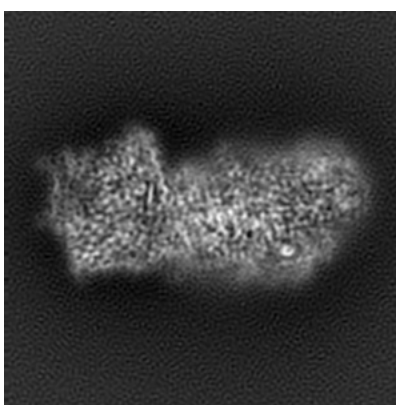
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

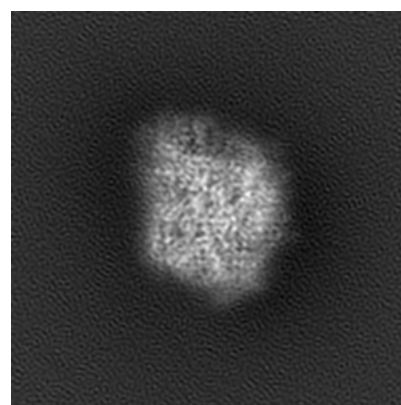
6.1.1 Primary map



X



Y

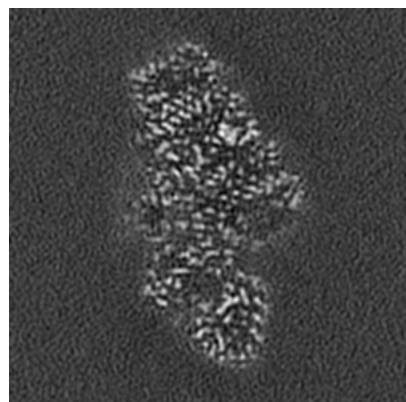


Z

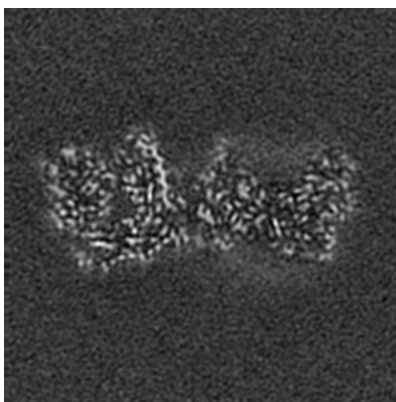
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

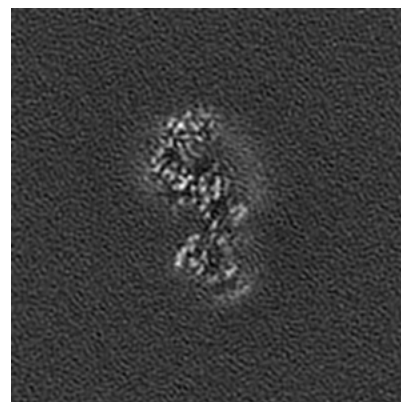
6.2.1 Primary map



X Index: 80



Y Index: 80

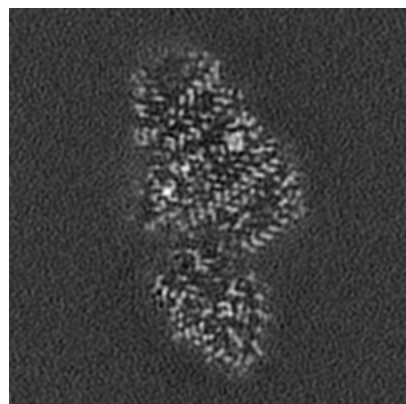


Z Index: 80

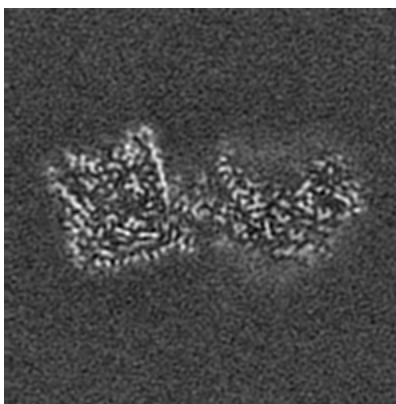
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

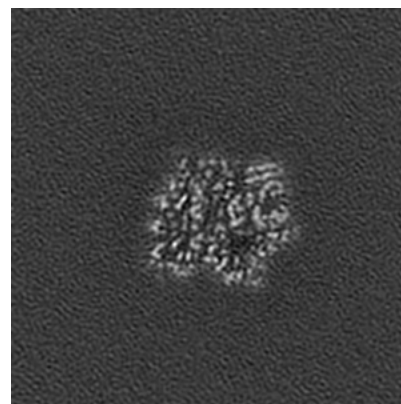
6.3.1 Primary map



X Index: 77



Y Index: 76

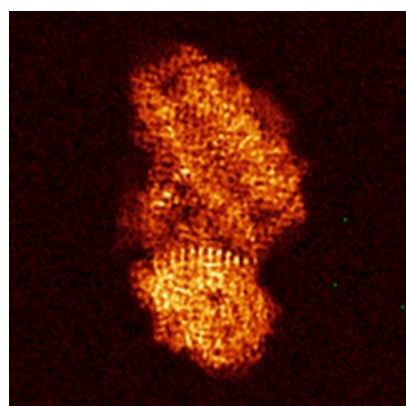


Z Index: 50

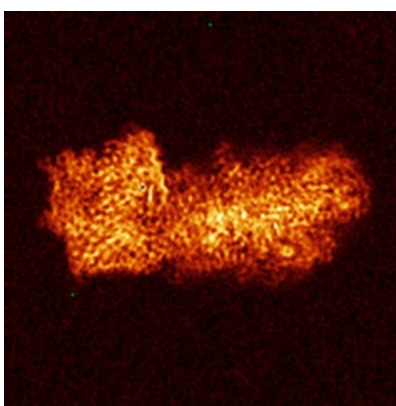
The images above show the largest variance slices of the map in three orthogonal directions.

6.4 Orthogonal standard-deviation projections (False-color) [i](#)

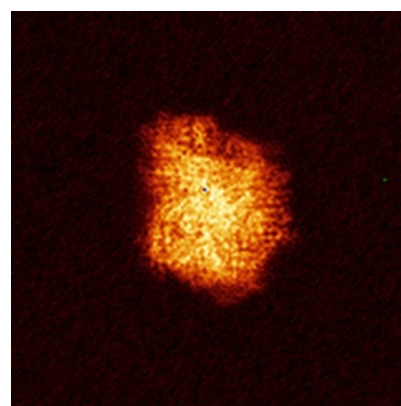
6.4.1 Primary map



X



Y



Z

The images above show the map standard deviation projections with false color in three orthogonal directions. Minimum values are shown in green, max in blue, and dark to light orange shades represent small to large values respectively.

6.5 Orthogonal surface views

This section was not generated.

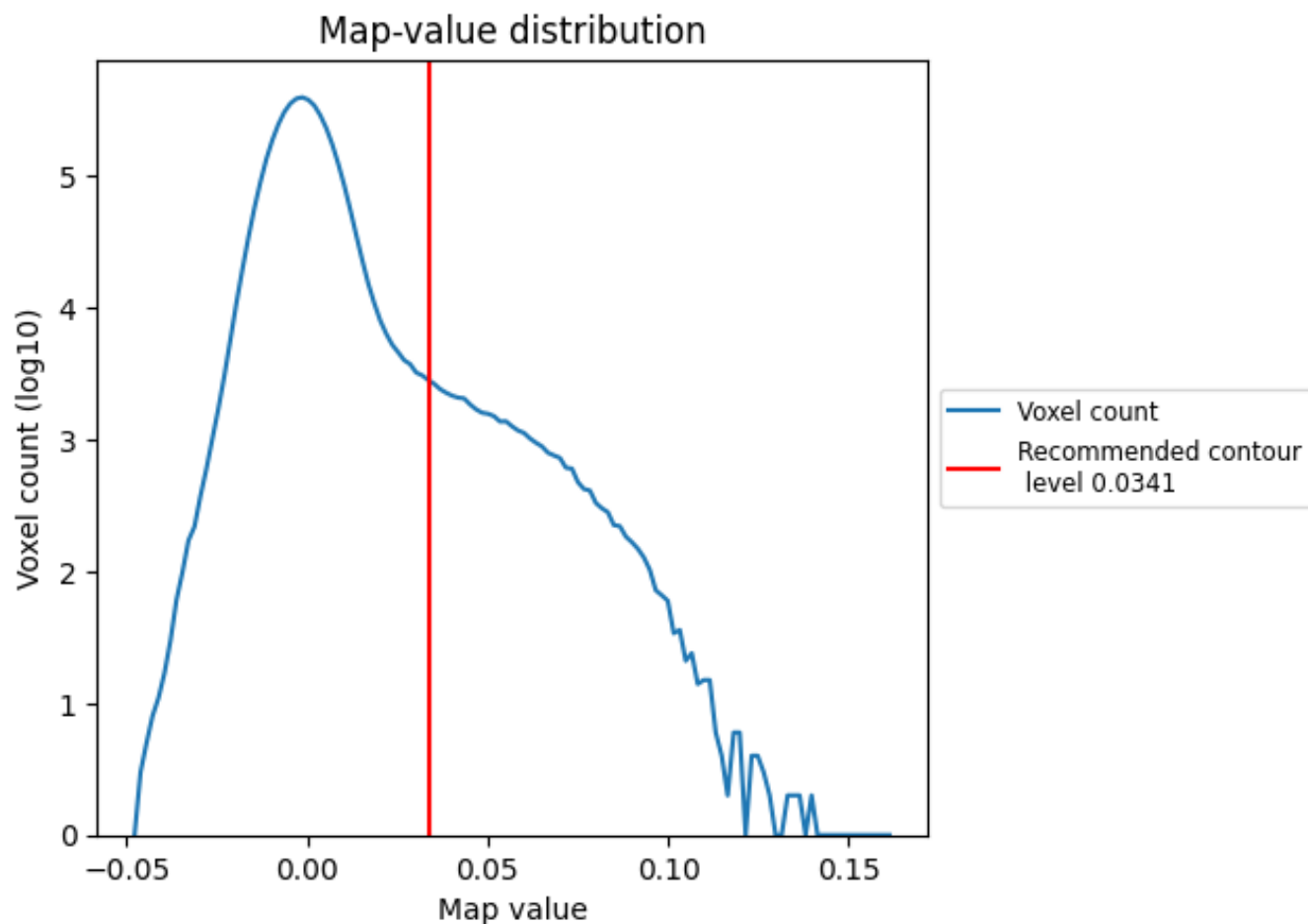
6.6 Mask visualisation

This section was not generated. No masks/segmentation were deposited.

7 Map analysis [i](#)

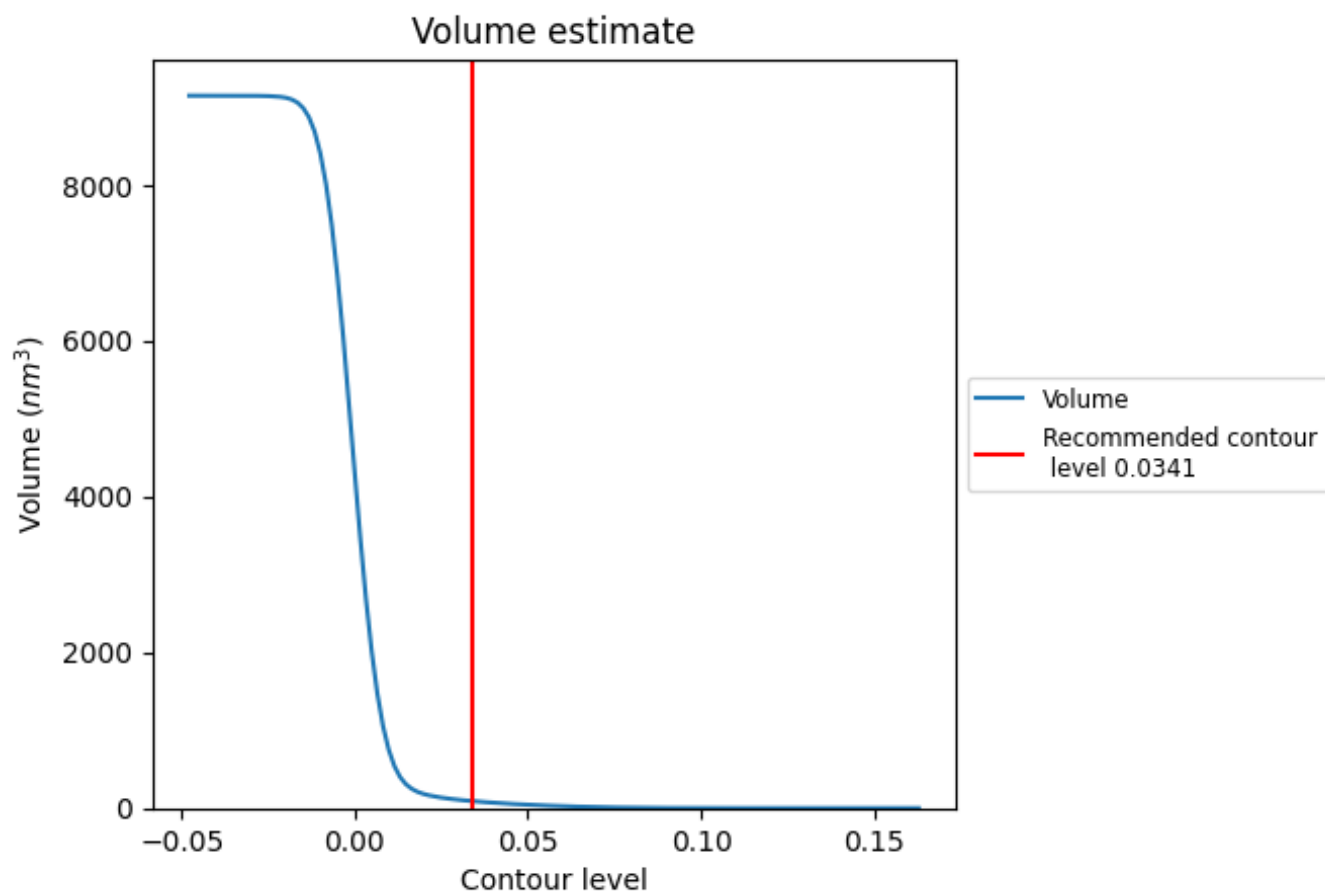
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



The map-value distribution is plotted in 128 intervals along the x-axis. The y-axis is logarithmic. A spike in this graph at zero usually indicates that the volume has been masked.

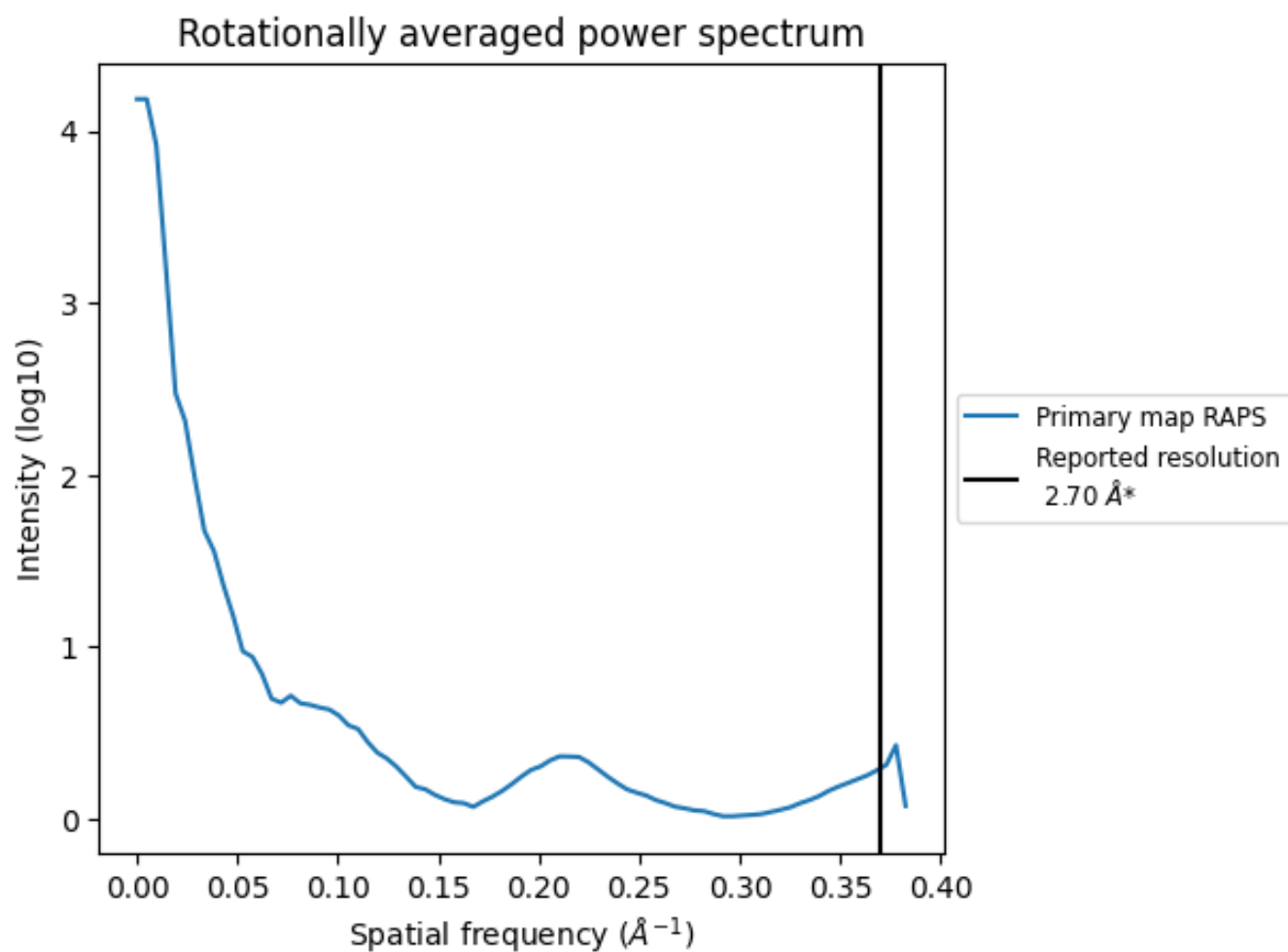
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 90 nm³; this corresponds to an approximate mass of 81 kDa.

The volume estimate graph shows how the enclosed volume varies with the contour level. The recommended contour level is shown as a vertical line and the intersection between the line and the curve gives the volume of the enclosed surface at the given level.

7.3 Rotationally averaged power spectrum ⓘ



*Reported resolution corresponds to spatial frequency of 0.370 Å⁻¹

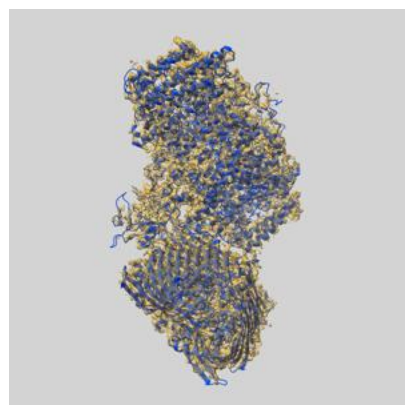
8 Fourier-Shell correlation

This section was not generated. No FSC curve or half-maps provided.

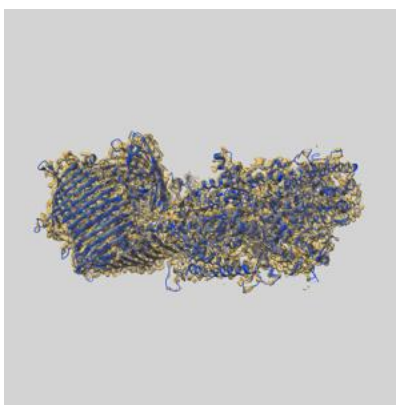
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-30069 and PDB model 6M32. Per-residue inclusion information can be found in section 3 on page 13.

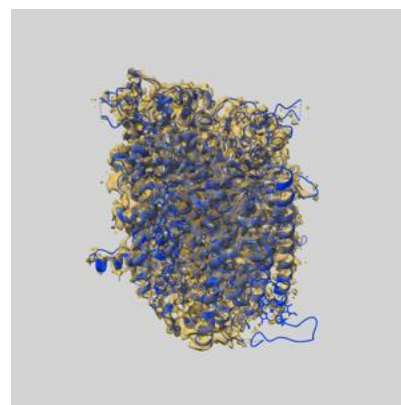
9.1 Map-model overlay [i](#)



X



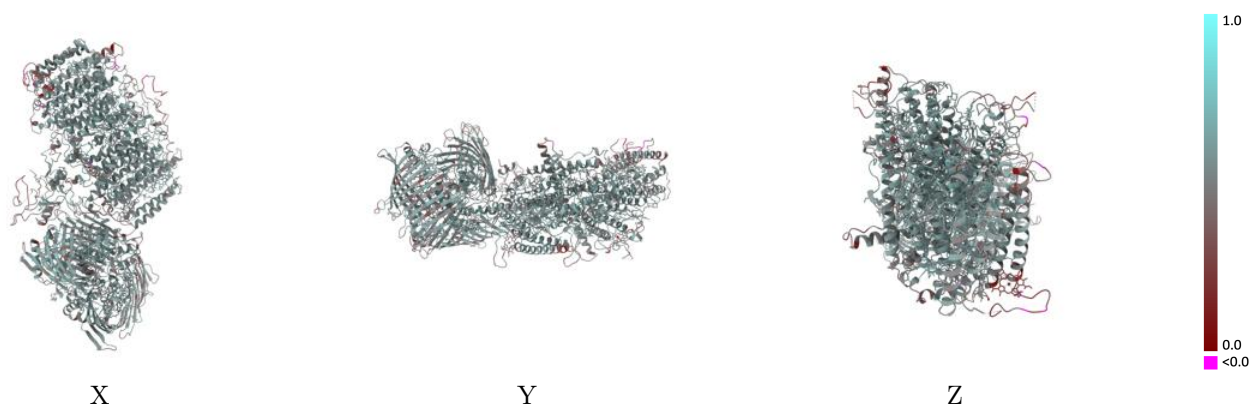
Y



Z

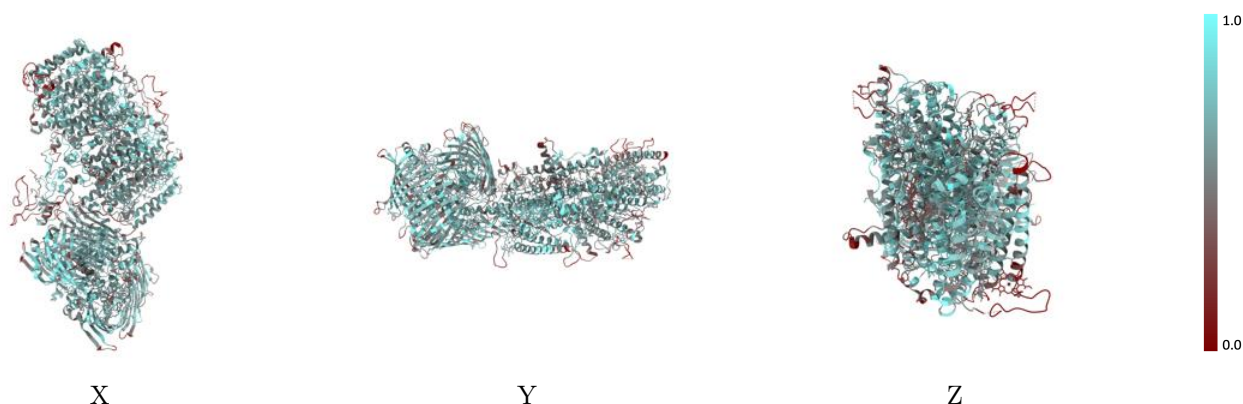
The images above show the 3D surface view of the map at the recommended contour level 0.0341 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

9.2 Q-score mapped to coordinate model [i](#)



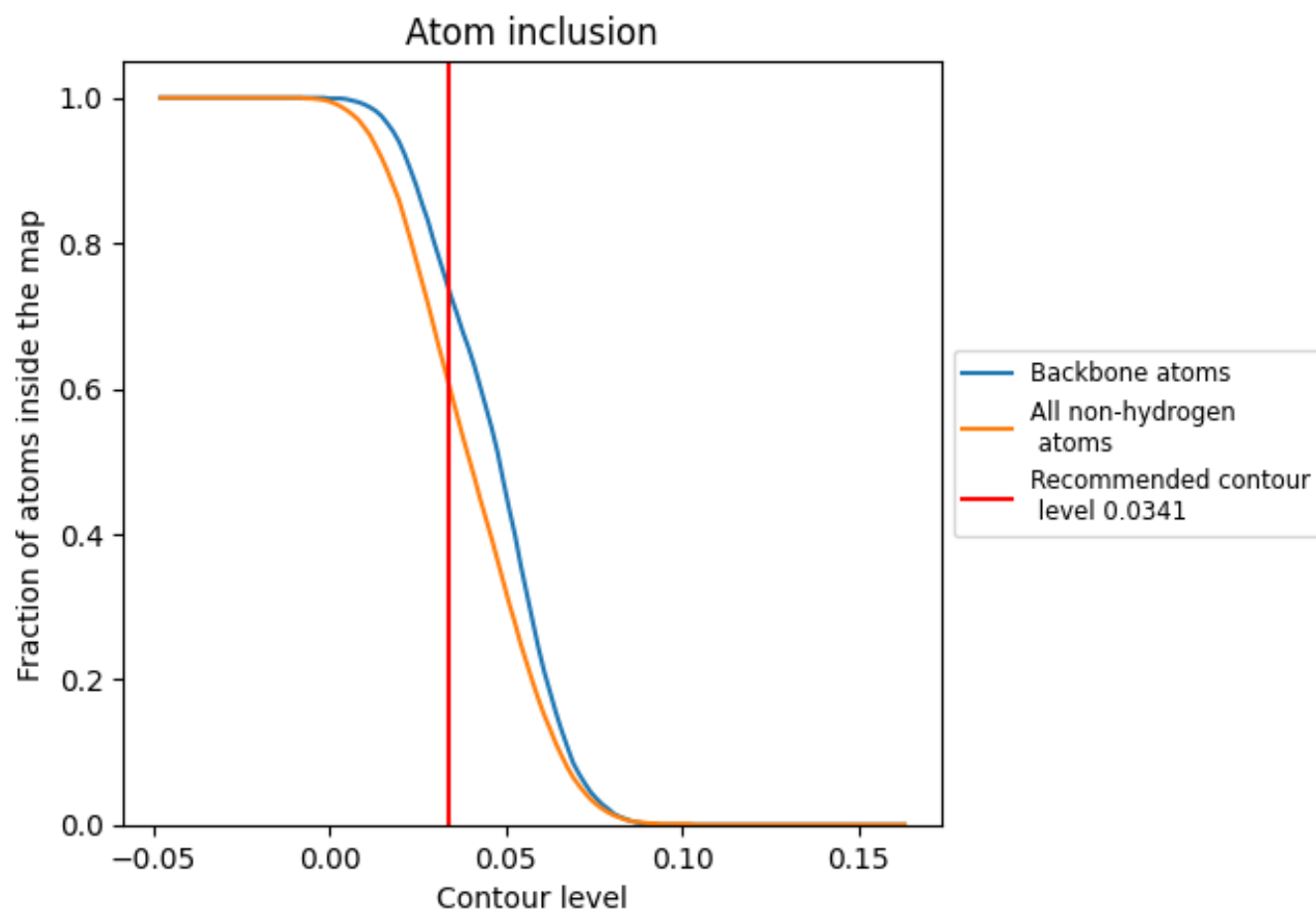
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions at which atoms can be resolved. Low Q-score values may therefore be expected for many entries.

9.3 Atom inclusion mapped to coordinate model [i](#)



The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (0.0341).

9.4 Atom inclusion ⓘ



At the recommended contour level, 73% of all backbone atoms, 60% of all non-hydrogen atoms, are inside the map.

9.5 Map-model fit summary ⓘ

The table lists the average atom inclusion at the recommended contour level (0.0341) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.6030	0.5250
A	0.6030	0.5350
B	0.5750	0.4900
D	0.3100	0.4350
E	0.6450	0.5350
F	0.6390	0.5260
G	0.6610	0.5420
a	0.5710	0.5150

1.0

0.0

<0.0