



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 12:44 AM UTC

PDB ID : 1YPZ / pdb_00001ypz
Title : Immune receptor
Authors : Adams, E.J.; Garcia, K.C.
Deposited on : 2005-01-31
Resolution : 3.40 Å(reported)

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

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

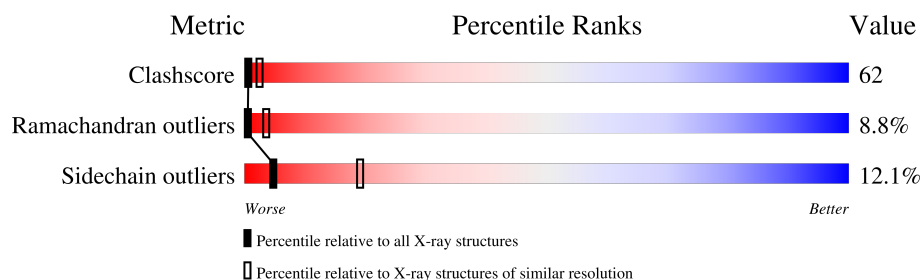
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.40 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	190562	1022 (3.44-3.36)
Ramachandran outliers	187476	1012 (3.44-3.36)
Sidechain outliers	187428	1012 (3.44-3.36)

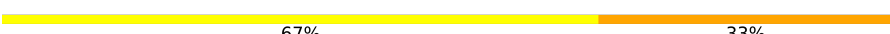
The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	260	23% 57% 18% .
1	C	260	20% 62% 17% .
2	B	102	32% 57% 10% .
2	D	102	22% 62% 16% .
3	E	207	28% 53% 17% .
3	G	207	35% 49% 14% .
4	F	230	27% 52% 18% .
4	H	230	34% 50% 14% .

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Mol	Chain	Length	Quality of chain
5	I	3	 33% 67%
5	L	3	 67% 33%
6	J	2	 100%
6	K	2	 50% 50%
7	M	3	 67% 33%

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
5	NAG	I	1	X	-	X	-
5	NAG	L	1	X	-	X	-
6	NAG	J	1	X	-	X	-
6	NAG	K	1	-	-	X	-
7	NAG	M	1	X	-	X	-

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 12984 atoms, of which 0 are hydrogens and 0 are deuteriums.

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

- Molecule 1 is a protein called H2-T22 protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	260	Total	C	N	O	S	0	0	0
			2106	1330	363	402	11			
1	C	260	Total	C	N	O	S	0	0	0
			2106	1330	363	402	11			

- Molecule 2 is a protein called Beta-2-microglobulin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	102	Total	C	N	O	S	0	0	0
			849	540	143	163	3			
2	D	102	Total	C	N	O	S	0	0	0
			849	540	143	163	3			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-1	ASP	-	cloning artifact	UNP P61769
B	0	PRO	-	cloning artifact	UNP P61769
D	-1	ASP	-	cloning artifact	UNP P61769
D	0	PRO	-	cloning artifact	UNP P61769

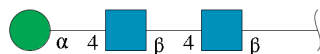
- Molecule 3 is a protein called T cell receptor delta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	E	207	Total	C	N	O	S	0	0	0
			1603	1000	276	321	6			
3	G	207	Total	C	N	O	S	281	0	0
			1592	992	274	321	5			

- Molecule 4 is a protein called T-cell receptor gamma chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	F	230	Total	C	N	O	S	0	0	0
			1852	1187	305	353	7			
4	H	230	Total	C	N	O	S	371	0	0
			1844	1181	303	353	7			

- Molecule 5 is an oligosaccharide called alpha-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



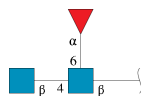
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
5	I	3	Total	C	N	O	0	0	0
			42	22	2	18			
5	L	3	Total	C	N	O	0	0	0
			42	22	2	18			

- Molecule 6 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
6	J	2	Total	C	N	O	0	0	0
			30	16	2	12			
6	K	2	Total	C	N	O	0	0	0
			30	16	2	12			

- Molecule 7 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose.



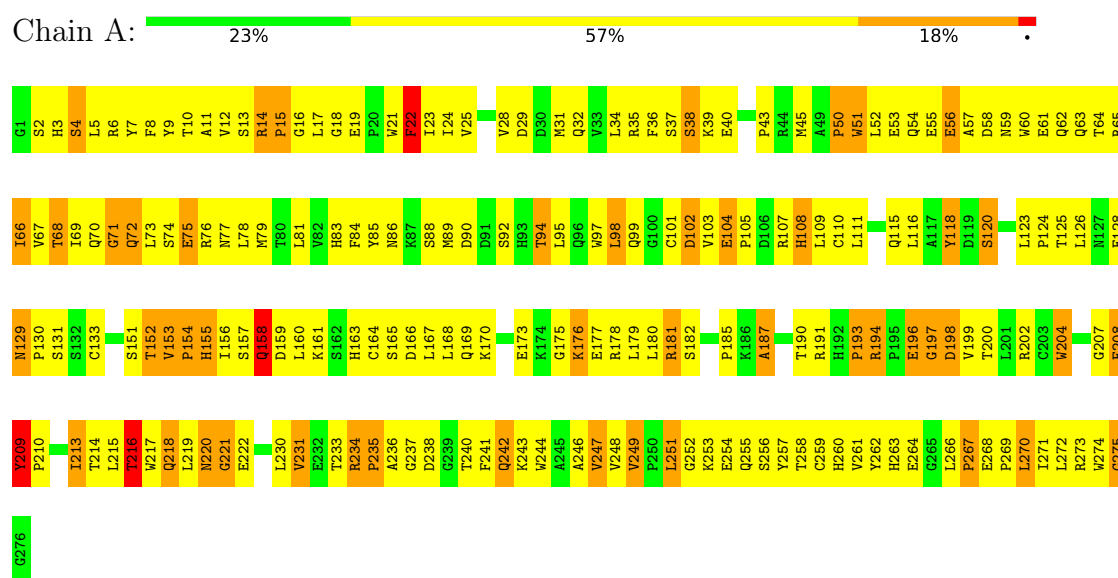
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
7	M	3	Total	C	N	O	0	0	0
			39	22	2	15			

3 Residue-property plots

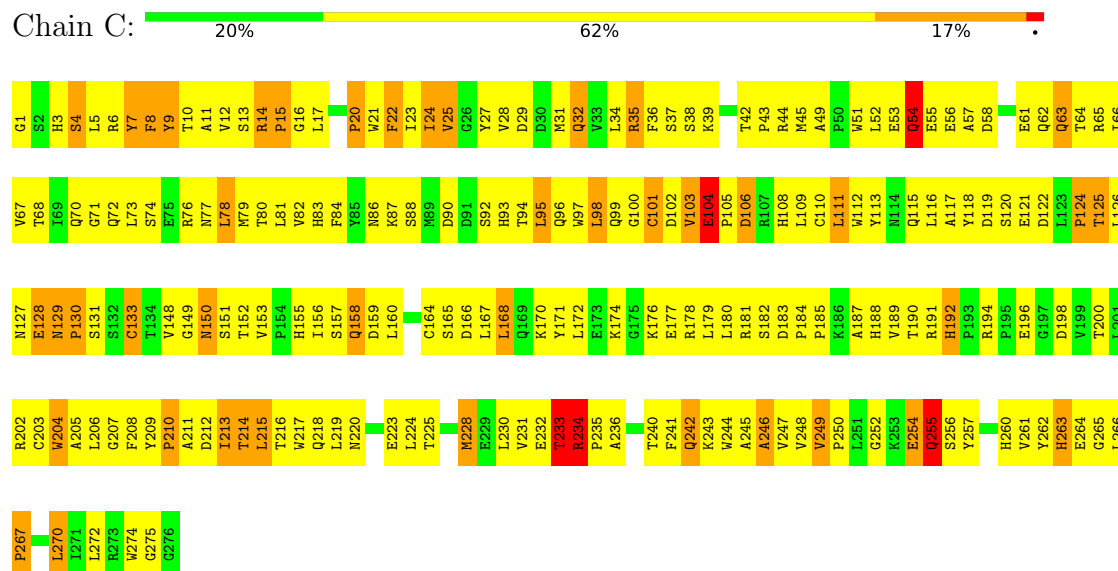
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded 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. 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.

Note EDS was not executed.

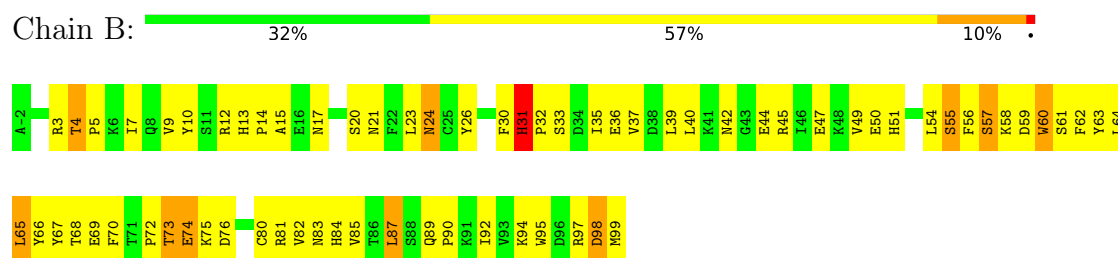
• Molecule 1: H2-T22 protein



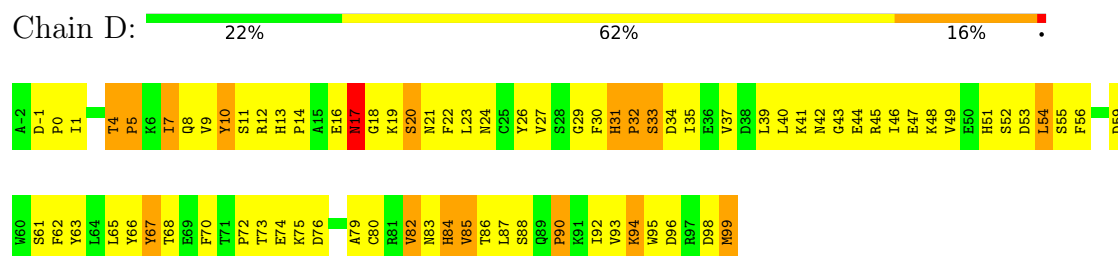
• Molecule 1: H2-T22 protein



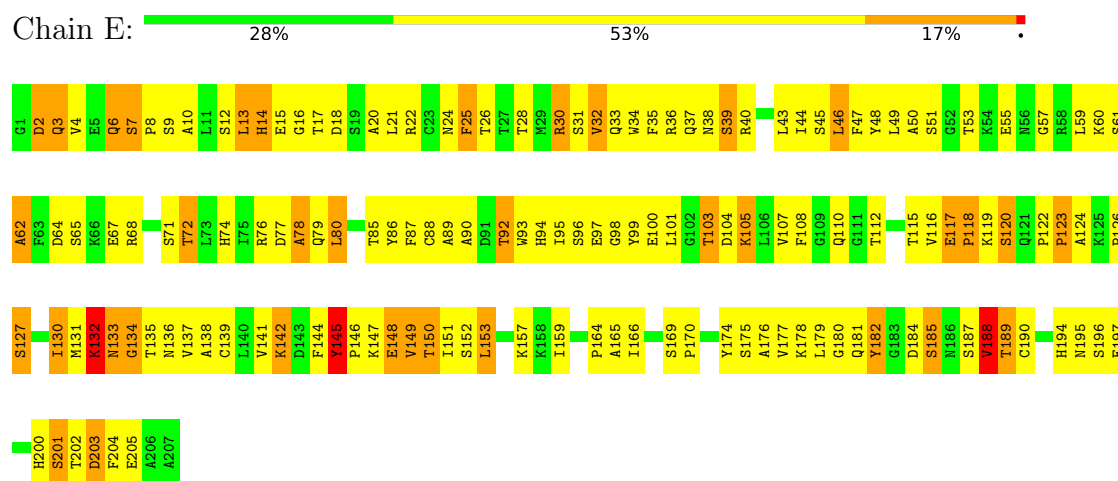
- Molecule 2: Beta-2-microglobulin



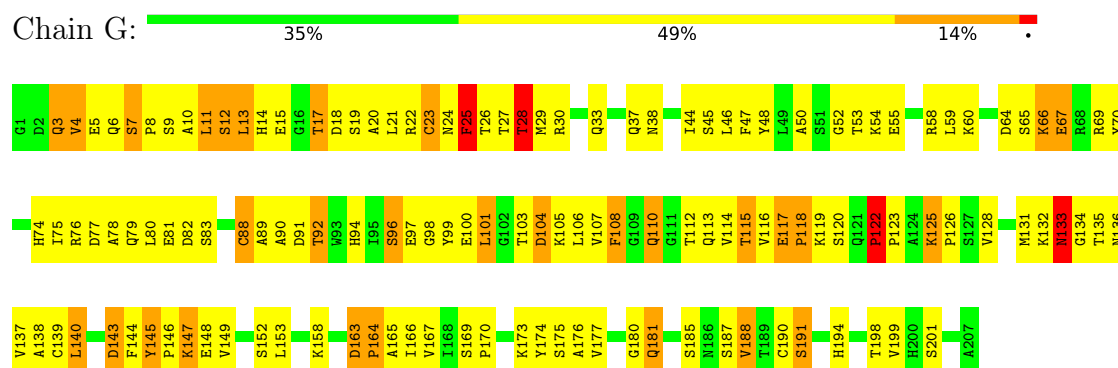
- Molecule 2: Beta-2-microglobulin



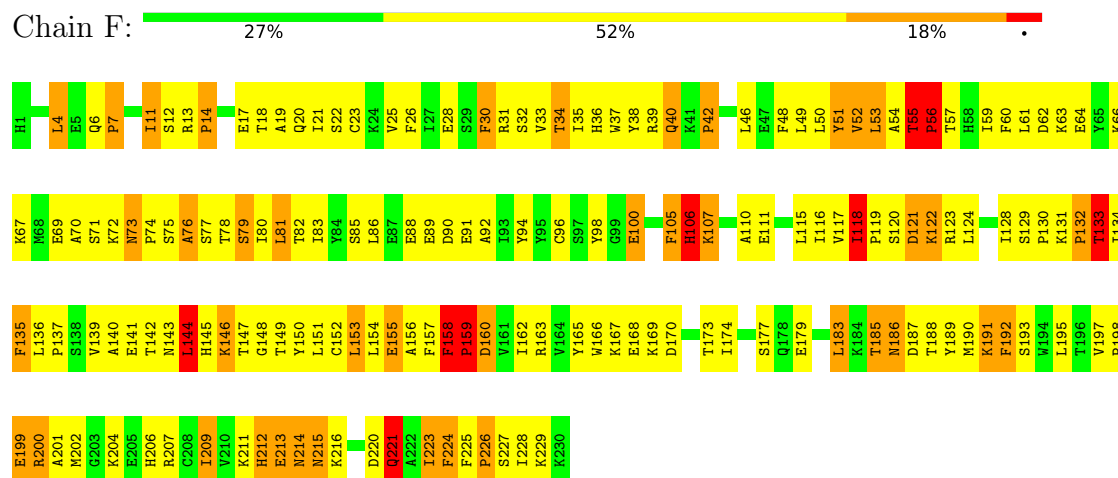
- Molecule 3: T cell receptor delta



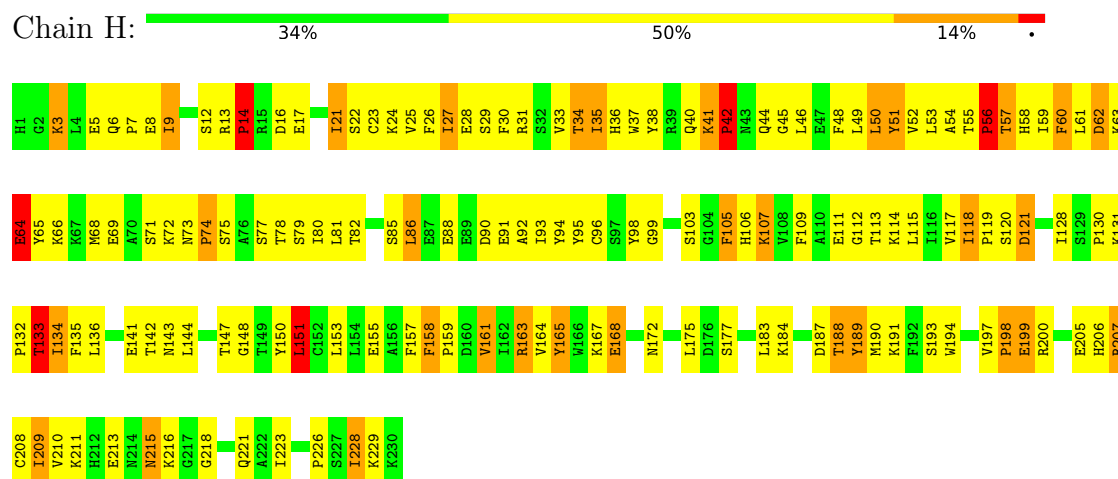
- Molecule 3: T cell receptor delta



- Molecule 4: T-cell receptor gamma chain



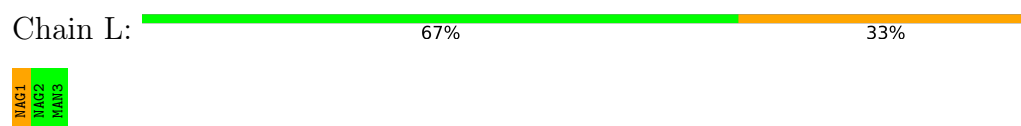
- Molecule 4: T-cell receptor gamma chain



- Molecule 5: alpha-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



- Molecule 5: alpha-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

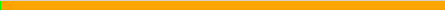


- Molecule 6: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain J:  100%

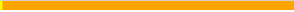
NAG1
NAG2

- Molecule 6: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain K:  50%  50%

NAG1
NAG2

- Molecule 7: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose

Chain M:  67%  33%

NAG1
NAG2
FUC3

4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	110.53Å 113.05Å 167.97Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 3.40	Depositor
% Data completeness (in resolution range)	97.4 (20.00-3.40)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.264 , 0.330	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	12984	wwPDB-VP
Average B, all atoms (Å ²)	103.0	wwPDB-VP

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: NAG, FUC, MAN

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	A	0.76	1/2166 (0.0%)	1.31	29/2951 (1.0%)
1	C	0.64	1/2166 (0.0%)	1.32	35/2951 (1.2%)
2	B	0.58	0/873	1.22	7/1182 (0.6%)
2	D	0.52	0/873	1.07	8/1182 (0.7%)
3	E	0.71	0/1636	1.29	22/2212 (1.0%)
3	G	0.67	0/1625	1.33	27/2201 (1.2%)
4	F	0.67	0/1898	1.29	21/2565 (0.8%)
4	H	0.58	0/1889	1.32	26/2553 (1.0%)
All	All	0.66	2/13126 (0.0%)	1.29	175/17797 (1.0%)

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
2	D	0	1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	155	HIS	CB-CG	5.58	1.57	1.50
1	C	228	MET	SD-CE	-5.25	1.66	1.79

All (175) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	H	29	SER	N-CA-C	-10.95	99.50	112.86
1	C	32	GLN	OE1-CD-NE2	-10.94	111.66	122.60
1	A	158	GLN	OE1-CD-NE2	-10.79	111.81	122.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	234	ARG	CA-C-N	10.77	133.47	120.23
1	C	234	ARG	C-N-CA	10.77	133.47	120.23
4	H	106	HIS	N-CA-C	-10.58	100.38	113.18
4	H	44	GLN	OE1-CD-NE2	-10.06	112.54	122.60
3	E	201	SER	N-CA-C	-9.87	98.98	111.02
4	F	13	ARG	CA-C-N	9.80	132.09	119.84
4	F	13	ARG	C-N-CA	9.80	132.09	119.84
1	C	255	GLN	OE1-CD-NE2	-9.80	112.80	122.60
1	A	72	GLN	OE1-CD-NE2	-9.79	112.81	122.60
1	C	54	GLN	OE1-CD-NE2	-9.72	112.88	122.60
1	A	54	GLN	OE1-CD-NE2	-9.70	112.90	122.60
1	A	32	GLN	OE1-CD-NE2	-9.69	112.91	122.60
2	B	89	GLN	OE1-CD-NE2	-9.66	112.94	122.60
1	C	72	GLN	OE1-CD-NE2	-9.56	113.04	122.60
4	F	221	GLN	OE1-CD-NE2	-9.55	113.05	122.60
3	G	125	LYS	CA-C-N	9.50	131.72	119.84
3	G	125	LYS	C-N-CA	9.50	131.72	119.84
4	F	106	HIS	N-CA-C	-9.41	99.83	112.26
3	G	37	GLN	OE1-CD-NE2	-9.35	113.25	122.60
4	F	40	GLN	OE1-CD-NE2	-9.30	113.30	122.60
1	A	218	GLN	OE1-CD-NE2	-9.30	113.30	122.60
3	G	3	GLN	OE1-CD-NE2	-9.20	113.40	122.60
3	E	6	GLN	OE1-CD-NE2	-8.99	113.61	122.60
1	C	242	GLN	OE1-CD-NE2	-8.96	113.64	122.60
3	E	3	GLN	N-CA-C	8.91	123.86	111.56
3	G	110	GLN	OE1-CD-NE2	-8.83	113.77	122.60
4	H	57	THR	N-CA-C	-8.77	102.75	113.19
4	H	206	HIS	N-CA-C	-8.33	97.89	110.14
3	G	147	LYS	N-CA-C	-8.30	102.26	112.38
4	F	209	ILE	N-CA-C	8.18	119.66	107.80
1	C	192	HIS	CA-C-N	8.11	128.11	119.76
1	C	192	HIS	C-N-CA	8.11	128.11	119.76
3	G	117	GLU	CA-C-N	8.10	128.05	120.03
3	G	117	GLU	C-N-CA	8.10	128.05	120.03
4	F	146	LYS	N-CA-C	-7.79	100.42	111.56
3	G	66	LYS	N-CA-C	-7.62	102.18	113.61
4	F	105	PHE	N-CA-C	7.55	121.41	107.60
2	B	20	SER	N-CA-C	7.53	120.30	110.43
4	H	207	ARG	N-CA-C	7.51	121.15	108.90
1	C	4	SER	N-CA-C	7.26	119.19	108.60
3	G	97	GLU	N-CA-C	-7.20	103.16	112.23
3	E	149	VAL	N-CA-C	7.17	118.67	108.42

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	158	GLN	CG-CD-NE2	7.14	127.11	116.40
2	B	97	ARG	N-CA-C	7.13	120.09	111.82
1	A	66	ILE	N-CA-C	7.06	117.57	110.23
1	A	208	PHE	N-CA-C	-7.06	98.43	109.72
2	D	52	SER	N-CA-C	7.01	119.57	110.53
1	C	125	THR	N-CA-C	-6.88	99.38	110.32
4	F	221	GLN	CG-CD-NE2	6.88	126.72	116.40
1	A	32	GLN	CG-CD-NE2	6.82	126.63	116.40
4	H	161	VAL	N-CA-C	6.82	117.30	106.32
3	G	37	GLN	CG-CD-NE2	6.82	126.63	116.40
1	A	72	GLN	N-CA-C	6.82	118.40	110.97
1	A	54	GLN	CG-CD-NE2	6.80	126.60	116.40
1	C	32	GLN	CG-CD-NE2	6.79	126.58	116.40
1	C	22	PHE	N-CA-C	6.77	118.44	108.14
2	B	89	GLN	CG-CD-NE2	6.75	126.52	116.40
4	H	106	HIS	CA-CB-CG	-6.74	107.06	113.80
3	E	6	GLN	CG-CD-NE2	6.73	126.50	116.40
4	H	44	GLN	CG-CD-NE2	6.72	126.48	116.40
4	H	3	LYS	N-CA-C	-6.69	104.69	114.39
4	H	88	GLU	N-CA-C	6.66	119.11	111.11
1	C	242	GLN	CG-CD-NE2	6.65	126.38	116.40
1	C	255	GLN	CG-CD-NE2	6.63	126.35	116.40
4	F	40	GLN	CG-CD-NE2	6.63	126.34	116.40
1	A	187	ALA	N-CA-C	6.62	119.94	108.76
1	A	218	GLN	CG-CD-NE2	6.58	126.27	116.40
1	C	54	GLN	CG-CD-NE2	6.52	126.18	116.40
1	C	72	GLN	CG-CD-NE2	6.42	126.03	116.40
1	C	270	LEU	N-CA-C	-6.35	99.66	109.76
1	A	72	GLN	CG-CD-NE2	6.30	125.85	116.40
3	G	110	GLN	CG-CD-NE2	6.27	125.80	116.40
4	H	228	ILE	N-CA-C	6.24	119.24	108.95
4	H	163	ARG	N-CA-C	6.21	118.41	108.41
3	G	3	GLN	CG-CD-NE2	6.19	125.69	116.40
3	G	50	ALA	N-CA-C	-6.17	105.44	114.39
2	D	5	PRO	N-CA-C	6.16	120.75	111.14
4	F	120	SER	N-CA-C	6.15	118.42	109.07
4	F	158	PHE	N-CA-C	6.14	123.39	109.81
3	G	110	GLN	CA-C-N	-6.10	116.40	122.20
3	G	110	GLN	C-N-CA	-6.10	116.40	122.20
3	E	180	GLY	N-CA-C	-6.10	101.79	110.63
4	F	118	ILE	CA-C-O	6.04	123.12	119.19
4	H	177	SER	N-CA-C	-6.03	100.44	109.96

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	10	TYR	N-CA-C	5.98	117.63	108.42
1	A	22	PHE	N-CA-C	5.97	118.20	108.76
4	H	151	LEU	N-CA-C	5.97	116.98	108.74
1	C	246	ALA	N-CA-C	5.96	117.26	108.86
1	A	209	TYR	CA-C-N	-5.95	112.40	119.84
1	A	209	TYR	C-N-CA	-5.95	112.40	119.84
1	C	158	GLN	N-CA-C	5.93	123.42	110.80
1	A	154	PRO	CA-N-CD	-5.92	103.72	112.00
3	E	62	ALA	N-CA-C	5.92	117.24	108.60
1	C	168	LEU	N-CA-C	-5.90	104.48	111.03
4	F	73	ASN	CA-C-N	5.86	127.16	119.84
4	F	73	ASN	C-N-CA	5.86	127.16	119.84
1	C	249	VAL	CA-C-N	5.84	127.15	119.84
1	C	249	VAL	C-N-CA	5.84	127.15	119.84
4	H	105	PHE	CB-CA-C	-5.84	102.67	110.79
3	E	105	LYS	N-CA-C	-5.81	98.42	110.80
3	E	142	LYS	N-CA-C	5.80	118.99	109.72
3	G	65	SER	N-CA-C	5.79	118.60	108.52
3	E	120	SER	N-CA-C	-5.78	101.48	110.10
3	E	130	ILE	N-CA-C	5.76	116.30	107.77
4	H	205	GLU	N-CA-C	5.75	118.83	108.48
3	G	145	TYR	N-CA-C	5.75	122.51	109.81
3	E	182	TYR	N-CA-C	5.74	117.69	109.14
1	A	155	HIS	CB-CA-C	-5.72	99.03	110.42
2	D	20	SER	N-CA-C	5.70	117.90	110.43
2	D	4	THR	CA-C-N	5.67	125.61	119.76
2	D	4	THR	C-N-CA	5.67	125.61	119.76
1	C	63	GLN	N-CA-C	-5.66	105.87	112.89
4	H	50	LEU	N-CA-C	5.64	116.53	108.74
4	H	136	LEU	N-CA-C	5.63	116.91	109.93
4	H	35	ILE	N-CA-C	5.60	115.92	107.80
1	C	159	ASP	N-CA-C	5.53	117.87	110.35
4	F	144	LEU	N-CA-C	5.51	120.12	112.90
3	G	158	LYS	N-CA-C	5.50	119.07	110.32
4	H	86	LEU	N-CA-C	5.47	117.63	109.59
3	E	132	LYS	N-CA-C	5.46	117.49	109.24
4	F	197	VAL	N-CA-C	5.45	113.27	107.76
1	A	216	THR	N-CA-C	5.45	118.33	108.69
3	G	118	PRO	N-CA-C	5.43	119.56	111.41
1	A	249	VAL	N-CA-C	5.43	113.81	107.89
1	A	4	SER	N-CA-C	5.42	117.45	108.13
2	B	24	ASN	N-CA-C	5.42	117.24	108.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	F	191	LYS	N-CA-C	5.42	116.75	108.52
2	B	31	HIS	CA-C-N	-5.38	113.12	119.84
2	B	31	HIS	C-N-CA	-5.38	113.12	119.84
1	C	14	ARG	CA-C-N	5.37	126.55	119.84
1	C	14	ARG	C-N-CA	5.37	126.55	119.84
4	F	56	PRO	N-CA-C	5.32	123.42	112.47
3	G	188	VAL	N-CA-C	5.30	120.37	109.34
3	G	116	VAL	N-CA-C	5.30	114.94	106.88
3	G	25	PHE	N-CA-C	-5.29	100.69	108.99
2	D	82	VAL	N-CA-C	5.28	115.65	107.78
3	E	127	SER	CA-C-N	-5.28	116.22	123.14
3	E	127	SER	C-N-CA	-5.28	116.22	123.14
1	C	65	ARG	N-CA-C	-5.28	105.42	111.07
3	E	85	THR	N-CA-C	5.27	118.21	109.46
3	E	14	HIS	N-CA-C	-5.23	103.00	110.59
3	E	9	SER	N-CA-C	-5.23	106.16	112.54
3	E	117	GLU	CA-C-O	-5.21	115.91	120.60
1	C	104	GLU	CA-C-N	5.21	126.35	119.84
1	C	104	GLU	C-N-CA	5.21	126.35	119.84
1	A	66	ILE	CB-CA-C	-5.20	105.04	111.70
4	H	68	MET	N-CA-C	5.18	119.60	113.12
1	A	197	GLY	N-CA-C	5.18	125.46	113.18
3	E	145	TYR	N-CA-C	5.18	121.25	109.81
4	F	133	THR	N-CA-C	-5.17	99.78	110.80
4	H	199	GLU	N-CA-C	5.17	121.81	110.80
3	E	166	ILE	N-CA-C	5.15	114.19	106.42
3	E	118	PRO	N-CA-C	5.14	119.38	111.21
2	D	45	ARG	N-CA-C	5.11	116.99	107.99
4	F	57	THR	N-CA-C	-5.11	104.84	111.74
1	C	103	VAL	N-CA-C	5.10	114.91	107.88
1	C	204	TRP	CA-C-N	5.09	129.45	121.76
1	C	204	TRP	C-N-CA	5.09	129.45	121.76
3	G	122	PRO	CA-C-N	5.08	126.19	119.84
3	G	122	PRO	C-N-CA	5.08	126.19	119.84
1	A	71	GLY	CA-C-N	5.07	127.34	120.65
1	A	71	GLY	C-N-CA	5.07	127.34	120.65
4	H	134	ILE	N-CA-C	5.06	115.92	110.21
4	H	188	THR	N-CA-C	-5.05	101.06	108.99
1	C	248	VAL	N-CA-C	-5.05	101.20	108.53
1	A	209	TYR	N-CA-C	5.03	120.92	109.81
3	G	88	CYS	N-CA-C	-5.03	102.35	110.14
1	C	35	ARG	N-CA-C	5.02	116.16	108.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	194	ARG	O-C-N	5.02	124.75	121.14
3	G	176	ALA	N-CA-C	5.01	116.04	108.07
1	A	204	TRP	N-CA-C	5.01	117.92	109.95
4	H	27	ILE	N-CA-C	5.01	115.06	107.80

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	D	63	TYR	Sidechain

5.2 Too-close contacts

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	A	2106	0	2006	291	1
1	C	2106	0	2009	289	4
2	B	849	0	810	85	0
2	D	849	0	810	123	0
3	E	1603	0	1554	203	1
3	G	1592	0	1524	168	0
4	F	1852	0	1832	246	4
4	H	1844	0	1816	171	0
5	I	42	0	38	10	0
5	L	42	0	36	7	0
6	J	30	0	27	9	0
6	K	30	0	27	22	0
7	M	39	0	36	12	0
All	All	12984	0	12525	1504	5

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 62.

All (1504) 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)
3:G:24:ASN:HD21	7:M:1:NAG:C1	1.10	1.58
4:F:186:ASN:HD22	6:K:1:NAG:C1	0.99	1.57
3:E:133:ASN:HD21	5:I:1:NAG:C1	0.88	1.53
1:A:155:HIS:CD2	1:A:157:SER:N	1.88	1.41
4:F:186:ASN:CB	6:K:1:NAG:O1	1.67	1.38
1:A:219:LEU:O	1:A:221:GLY:N	1.58	1.32
3:E:7:SER:OG	3:E:8:PRO:HD2	1.24	1.31
3:E:133:ASN:ND2	5:I:1:NAG:H1	0.97	1.28
1:C:84:PHE:CE1	3:G:30:ARG:NH2	2.07	1.23
3:E:22:ARG:NH2	6:J:2:NAG:O7	1.72	1.21
3:G:24:ASN:ND2	7:M:1:NAG:H1	0.88	1.19
4:F:186:ASN:ND2	6:K:1:NAG:H1	1.43	1.16
1:A:155:HIS:NE2	1:A:157:SER:CB	2.10	1.15
1:C:228:MET:HE2	1:C:245:ALA:HB1	1.24	1.14
1:C:194:ARG:HB2	1:C:198:ASP:O	1.46	1.11
4:F:224:PHE:CE1	5:I:3:MAN:O2	2.03	1.11
4:F:186:ASN:ND2	6:K:1:NAG:O5	1.71	1.10
4:H:158:PHE:H	4:H:159:PRO:HD3	1.17	1.09
3:G:58:ARG:HH22	3:G:81:GLU:HG3	0.95	1.09
4:F:40:GLN:O	4:F:92:ALA:HB1	1.52	1.08
4:H:134:ILE:N	5:L:1:NAG:O6	1.86	1.08
4:H:40:GLN:HB2	4:H:46:LEU:HD12	1.32	1.08
4:F:186:ASN:HB2	6:K:1:NAG:O1	1.46	1.07
4:F:186:ASN:HB3	6:K:1:NAG:O1	1.38	1.07
1:C:81:LEU:HD21	3:G:101:LEU:HD12	1.34	1.07
4:F:6:GLN:HE21	4:F:110:ALA:HB3	1.19	1.06
4:F:50:LEU:HD21	4:F:61:LEU:HG	1.38	1.06
1:A:155:HIS:HE2	1:A:157:SER:CB	1.67	1.05
3:E:7:SER:OG	3:E:8:PRO:CD	2.02	1.05
1:C:129:ASN:HB3	1:C:130:PRO:CD	1.86	1.05
4:F:186:ASN:CB	6:K:1:NAG:C1	2.33	1.05
4:F:186:ASN:HB2	6:K:1:NAG:C1	1.86	1.05
3:E:126:PRO:HG3	3:E:144:PHE:HB3	1.36	1.04
3:E:24:ASN:ND2	6:J:1:NAG:N2	2.06	1.04
3:G:11:LEU:HD21	3:G:13:LEU:HD21	1.35	1.04
4:F:128:ILE:HG21	4:F:158:PHE:HB2	1.36	1.02
3:E:132:LYS:HE2	3:E:132:LYS:H	1.23	1.02
4:H:52:VAL:HG22	4:H:71:SER:HB2	1.40	1.01
3:E:24:ASN:ND2	6:J:1:NAG:C2	2.24	1.01
1:A:155:HIS:CD2	1:A:157:SER:H	1.74	1.00
1:C:228:MET:CE	1:C:245:ALA:HB1	1.92	1.00
1:C:49:ALA:HB3	1:C:52:LEU:HD12	1.38	1.00

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:31:MET:HE3	1:C:31:MET:HA	1.39	1.00
1:A:155:HIS:CD2	1:A:157:SER:CB	2.44	0.99
1:C:80:THR:HB	4:H:103:SER:HB3	1.41	0.99
4:F:186:ASN:CG	6:K:1:NAG:C1	2.34	0.99
3:G:15:GLU:HB3	3:G:119:LYS:HG3	1.42	0.99
1:A:22:PHE:O	1:A:37:SER:HB2	1.62	0.98
1:A:155:HIS:CD2	1:A:157:SER:CA	2.47	0.98
1:A:31:MET:HG2	1:A:209:TYR:OH	1.63	0.97
4:F:224:PHE:HE1	5:I:3:MAN:O2	1.45	0.97
1:C:129:ASN:HB3	1:C:130:PRO:HD3	1.42	0.97
4:F:185:THR:HG23	6:K:1:NAG:H82	1.46	0.97
3:G:58:ARG:NH2	3:G:81:GLU:HG3	1.78	0.97
4:H:134:ILE:CA	5:L:1:NAG:O6	2.12	0.97
1:C:32:GLN:HE22	2:D:53:ASP:CG	1.72	0.97
1:A:129:ASN:HB2	1:A:130:PRO:HD3	1.47	0.94
1:A:266:LEU:HD22	1:A:270:LEU:HD13	1.48	0.94
1:C:96:GLN:NE2	2:D:31:HIS:NE2	2.15	0.94
1:C:194:ARG:CB	1:C:198:ASP:O	2.14	0.94
4:F:207:ARG:HG2	4:F:224:PHE:HA	1.49	0.94
1:A:155:HIS:HD2	1:A:157:SER:N	1.44	0.93
1:A:218:GLN:HB2	1:A:222:GLU:O	1.69	0.93
4:F:186:ASN:CG	6:K:1:NAG:H1	1.93	0.92
3:G:58:ARG:NH1	3:G:79:GLN:HB2	1.84	0.92
3:G:24:ASN:ND2	7:M:1:NAG:C2	2.32	0.92
4:H:17:GLU:O	4:H:86:LEU:HG	1.68	0.92
1:A:219:LEU:O	1:A:220:ASN:C	2.13	0.91
4:H:52:VAL:CG2	4:H:71:SER:HB2	2.01	0.91
3:G:3:GLN:HE21	3:G:26:THR:H	0.98	0.91
1:C:99:GLN:NE2	1:C:160:LEU:HD23	1.86	0.91
3:G:4:VAL:HG13	3:G:25:PHE:HD2	1.34	0.91
3:G:24:ASN:HD21	7:M:1:NAG:C2	1.83	0.91
3:G:24:ASN:HD22	7:M:1:NAG:H1	1.32	0.91
1:A:155:HIS:CD2	1:A:157:SER:HB3	2.06	0.91
1:A:157:SER:O	1:A:158:GLN:CD	2.15	0.90
4:F:185:THR:HG23	6:K:1:NAG:C8	2.02	0.89
3:E:37:GLN:HB3	3:E:43:LEU:HD23	1.51	0.89
3:G:166:ILE:HG23	3:G:175:SER:O	1.72	0.89
4:H:134:ILE:N	5:L:1:NAG:C6	2.36	0.89
4:H:34:THR:HA	4:H:52:VAL:O	1.72	0.88
1:C:98:LEU:HD12	1:C:99:GLN:N	1.86	0.88
1:C:202:ARG:HH11	1:C:202:ARG:HB2	1.38	0.88

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:H:12:SER:HB2	4:H:118:ILE:HD11	1.53	0.88
4:F:91:GLU:HG3	4:F:115:LEU:O	1.73	0.87
1:A:103:VAL:HG12	1:A:105:PRO:HD3	1.55	0.87
1:A:84:PHE:CZ	3:E:30:ARG:NE	2.43	0.87
1:C:84:PHE:CD1	3:G:30:ARG:NH2	2.43	0.87
1:C:194:ARG:HD2	1:C:198:ASP:HB2	1.57	0.87
2:B:40:LEU:HA	2:B:44:GLU:O	1.75	0.87
4:H:30:PHE:O	4:H:54:ALA:HB1	1.74	0.87
1:A:51:TRP:CE3	1:A:51:TRP:O	2.27	0.87
3:G:11:LEU:C	3:G:11:LEU:HD23	2.00	0.87
3:G:4:VAL:HG13	3:G:25:PHE:CD2	2.09	0.86
1:C:81:LEU:CD2	3:G:101:LEU:HD12	2.04	0.86
1:C:9:TYR:OH	1:C:74:SER:HB2	1.76	0.86
4:H:25:VAL:HG12	4:H:27:ILE:H	1.39	0.86
2:B:83:ASN:ND2	2:B:90:PRO:HG3	1.90	0.86
4:F:129:SER:HB2	4:F:215:ASN:HA	1.56	0.86
1:C:32:GLN:NE2	2:D:53:ASP:OD2	2.09	0.85
2:D:86:THR:HG22	2:D:87:LEU:H	1.38	0.85
3:G:3:GLN:HE21	3:G:26:THR:N	1.72	0.85
4:H:134:ILE:H	5:L:1:NAG:C6	1.88	0.85
1:C:31:MET:HB2	1:C:209:TYR:OH	1.76	0.85
1:C:189:VAL:HG12	1:C:190:THR:N	1.91	0.85
1:C:24:ILE:HD12	1:C:67:VAL:HG13	1.58	0.85
3:G:58:ARG:HH22	3:G:81:GLU:CG	1.86	0.85
4:F:156:ALA:HA	4:F:188:THR:HG21	1.57	0.84
1:A:124:PRO:HB3	3:E:28:THR:HG22	1.58	0.84
1:C:103:VAL:HG21	1:C:168:LEU:HD23	1.59	0.84
1:C:111:LEU:HD12	1:C:111:LEU:N	1.93	0.84
3:E:53:THR:HB	3:E:62:ALA:HB2	1.60	0.84
3:E:126:PRO:HG3	3:E:144:PHE:CB	2.07	0.84
4:F:186:ASN:ND2	6:K:1:NAG:C5	2.41	0.83
3:E:24:ASN:ND2	6:J:1:NAG:O1	2.09	0.83
1:C:234:ARG:HG2	2:D:8:GLN:OE1	1.79	0.83
2:D:59:ASP:OD1	2:D:61:SER:HB2	1.79	0.83
1:A:155:HIS:NE2	1:A:157:SER:OG	1.66	0.83
4:H:158:PHE:H	4:H:159:PRO:CD	1.92	0.83
2:D:37:VAL:HB	2:D:66:TYR:CZ	2.14	0.83
2:D:20:SER:O	2:D:21:ASN:CG	2.21	0.82
4:H:93:ILE:HG12	4:H:114:LYS:HB2	1.62	0.82
2:B:31:HIS:HB3	2:B:32:PRO:HD3	1.62	0.82
4:H:52:VAL:HG22	4:H:71:SER:CB	2.10	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:84:HIS:H	2:B:87:LEU:CD2	1.92	0.81
3:G:46:LEU:HD22	3:G:59:LEU:HD12	1.60	0.81
3:E:24:ASN:OD1	3:E:24:ASN:O	1.97	0.81
2:B:84:HIS:H	2:B:87:LEU:HD21	1.45	0.81
1:A:155:HIS:CE1	1:A:157:SER:HG	1.98	0.81
4:F:224:PHE:CZ	5:I:3:MAN:O2	2.33	0.81
3:G:11:LEU:HD21	3:G:13:LEU:CD2	2.10	0.81
2:D:17:ASN:N	2:D:17:ASN:HD22	1.79	0.81
3:E:78:ALA:O	3:E:79:GLN:HG2	1.81	0.81
3:E:164:PRO:HG3	3:E:178:LYS:HD2	1.62	0.81
4:F:159:PRO:HD2	4:F:212:HIS:NE2	1.94	0.81
1:C:78:LEU:O	1:C:82:VAL:HG23	1.81	0.80
2:D:31:HIS:HB3	2:D:32:PRO:HD3	1.61	0.80
1:C:23:ILE:HG22	1:C:24:ILE:H	1.46	0.80
1:C:178:ARG:O	1:C:179:LEU:HD23	1.81	0.80
4:F:152:CYS:HB2	4:F:166:TRP:CZ2	2.16	0.80
1:C:84:PHE:CE1	3:G:30:ARG:CZ	2.65	0.80
1:A:155:HIS:NE2	1:A:157:SER:N	2.29	0.80
4:H:134:ILE:O	5:L:1:NAG:O5	1.99	0.80
1:A:24:ILE:HD12	1:A:67:VAL:HG13	1.62	0.80
3:E:7:SER:CB	3:E:8:PRO:CD	2.60	0.80
3:G:7:SER:HB2	3:G:8:PRO:HD3	1.64	0.80
1:A:155:HIS:HD2	1:A:157:SER:CA	1.92	0.80
3:E:15:GLU:OE1	3:E:118:PRO:HA	1.82	0.79
1:A:234:ARG:HG3	2:B:10:TYR:CZ	2.17	0.79
4:F:153:LEU:HD12	4:F:154:LEU:N	1.96	0.79
1:A:155:HIS:CD2	1:A:156:ILE:C	2.60	0.79
3:E:169:SER:OG	3:E:175:SER:HB2	1.82	0.79
4:H:40:GLN:HB2	4:H:46:LEU:CD1	2.12	0.79
3:G:3:GLN:NE2	3:G:26:THR:H	1.78	0.79
1:A:77:ASN:OD1	3:E:101:LEU:HB2	1.83	0.79
1:A:153:VAL:HB	1:A:154:PRO:CD	2.13	0.79
3:E:145:TYR:HB3	3:E:146:PRO:HD3	1.65	0.79
4:F:186:ASN:HD22	6:K:1:NAG:C5	1.94	0.79
1:C:81:LEU:HD21	3:G:101:LEU:CD1	2.13	0.78
3:G:30:ARG:HG2	3:G:91:ASP:O	1.83	0.78
4:F:130:PRO:HD2	4:F:215:ASN:HB2	1.65	0.78
1:C:97:TRP:HD1	1:C:98:LEU:N	1.79	0.78
2:D:23:LEU:O	2:D:67:TYR:HA	1.84	0.78
3:E:26:THR:HG22	3:E:26:THR:O	1.81	0.78
3:G:144:PHE:O	3:G:173:LYS:HB3	1.84	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:H:93:ILE:HA	4:H:114:LYS:HA	1.66	0.78
4:H:158:PHE:N	4:H:159:PRO:HD3	1.95	0.78
2:D:5:PRO:HB3	2:D:30:PHE:HB3	1.66	0.78
1:C:87:LYS:HB3	1:C:93:HIS:HE1	1.49	0.78
2:D:20:SER:O	2:D:21:ASN:OD1	2.02	0.78
2:D:23:LEU:HB3	2:D:68:THR:HG22	1.63	0.78
1:C:23:ILE:HG22	1:C:24:ILE:N	1.99	0.77
3:G:3:GLN:HG3	3:G:3:GLN:O	1.83	0.77
1:A:259:CYS:HB3	1:A:272:LEU:HB2	1.64	0.77
1:C:115:GLN:C	1:C:116:LEU:HD12	2.10	0.77
4:H:36:HIS:HA	4:H:50:LEU:O	1.83	0.77
2:D:19:LYS:O	2:D:72:PRO:HD2	1.84	0.77
4:F:12:SER:HA	4:F:116:ILE:O	1.84	0.77
2:D:54:LEU:HD11	2:D:62:PHE:HD2	1.50	0.77
3:G:148:GLU:O	3:G:149:VAL:HG23	1.83	0.77
4:F:6:GLN:NE2	4:F:110:ALA:HB3	1.97	0.76
4:F:25:VAL:HG23	4:F:78:THR:HA	1.66	0.76
3:E:108:PHE:CD1	4:F:46:LEU:HB2	2.19	0.76
2:D:17:ASN:HA	2:D:72:PRO:HB2	1.68	0.76
1:C:98:LEU:HD12	1:C:98:LEU:C	2.10	0.76
1:C:209:TYR:O	1:C:211:ALA:N	2.19	0.76
4:H:92:ALA:O	4:H:94:TYR:CE1	2.39	0.76
4:F:25:VAL:HG21	4:F:35:ILE:HD11	1.66	0.76
2:D:80:CYS:O	2:D:92:ILE:HA	1.86	0.76
4:F:188:THR:HG22	4:F:189:TYR:H	1.50	0.76
1:C:176:LYS:HB3	1:C:180:LEU:HD12	1.67	0.75
1:A:5:LEU:HD12	1:A:6:ARG:N	2.02	0.75
4:F:62:ASP:HA	4:F:67:LYS:HA	1.68	0.75
2:B:70:PHE:CZ	2:B:72:PRO:HG3	2.22	0.75
3:E:7:SER:CB	3:E:8:PRO:HD2	2.15	0.75
4:F:152:CYS:HB2	4:F:166:TRP:HZ2	1.51	0.75
1:C:103:VAL:CG2	1:C:168:LEU:HD23	2.17	0.75
1:A:266:LEU:HD12	1:A:266:LEU:O	1.87	0.75
4:F:69:GLU:OE1	4:F:81:LEU:HD21	1.87	0.75
4:H:74:PRO:HD2	4:H:78:THR:O	1.87	0.75
4:H:107:LYS:HG3	4:H:109:PHE:CE1	2.22	0.75
2:D:23:LEU:HD23	2:D:39:LEU:HD13	1.67	0.74
1:A:238:ASP:OD1	1:A:240:THR:HG23	1.86	0.74
2:B:17:ASN:HA	2:B:72:PRO:O	1.87	0.74
1:A:202:ARG:HD3	1:A:204:TRP:CZ2	2.22	0.74
1:A:209:TYR:HB3	1:A:210:PRO:HD3	1.69	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:G:131:MET:CB	3:G:140:LEU:HD21	2.18	0.74
1:A:155:HIS:HD2	1:A:156:ILE:C	1.95	0.74
1:A:194:ARG:HG2	1:A:198:ASP:O	1.87	0.74
3:E:53:THR:HB	3:E:62:ALA:CB	2.18	0.74
4:H:210:VAL:HG12	4:H:211:LYS:H	1.53	0.74
4:F:128:ILE:CG2	4:F:158:PHE:HB2	2.15	0.74
4:F:168:GLU:HG2	4:F:170:ASP:OD1	1.87	0.74
1:A:208:PHE:CZ	1:A:213:ILE:HD12	2.23	0.74
1:A:209:TYR:O	1:A:210:PRO:C	2.31	0.74
2:B:23:LEU:HD23	2:B:39:LEU:HD13	1.69	0.74
3:E:33:GLN:HE22	4:F:107:LYS:HB2	1.52	0.74
1:C:182:SER:O	1:C:184:PRO:HD3	1.87	0.73
7:M:1:NAG:H83	7:M:1:NAG:O1	1.87	0.73
3:E:133:ASN:HD22	5:I:1:NAG:H1	1.47	0.73
3:E:185:SER:O	3:E:188:VAL:HG23	1.87	0.73
4:F:147:THR:HG22	4:F:148:GLY:N	2.02	0.73
4:F:165:TYR:HE1	4:F:167:LYS:HD2	1.52	0.73
3:G:7:SER:CB	3:G:8:PRO:CD	2.67	0.73
3:G:98:GLY:C	3:G:99:TYR:HD1	1.97	0.73
4:F:33:VAL:HG12	4:F:34:THR:H	1.52	0.73
1:C:17:LEU:O	1:C:17:LEU:HD23	1.88	0.73
3:E:133:ASN:ND2	5:I:1:NAG:O1	2.21	0.73
4:H:37:TRP:CE3	4:H:81:LEU:HD22	2.23	0.73
4:H:41:LYS:HB3	4:H:42:PRO:HD2	1.71	0.73
1:C:80:THR:CB	4:H:103:SER:HB3	2.17	0.73
1:A:153:VAL:HB	1:A:154:PRO:HD2	1.69	0.73
3:E:53:THR:CB	3:E:62:ALA:HB2	2.18	0.72
4:F:33:VAL:HG12	4:F:34:THR:N	2.04	0.72
3:G:5:GLU:O	3:G:23:CYS:HA	1.89	0.72
4:H:210:VAL:HG12	4:H:211:LYS:N	2.02	0.72
3:E:34:TRP:CZ3	3:E:88:CYS:HB3	2.24	0.72
1:C:183:ASP:O	1:C:208:PHE:HA	1.89	0.72
1:C:189:VAL:CG1	1:C:190:THR:N	2.52	0.72
4:F:162:ILE:O	4:F:163:ARG:HD3	1.90	0.72
3:G:38:ASN:ND2	3:G:44:ILE:HD11	2.04	0.72
3:E:96:SER:C	3:E:98:GLY:H	1.96	0.72
1:C:202:ARG:HB2	1:C:202:ARG:NH1	2.04	0.72
4:H:33:VAL:HB	4:H:54:ALA:HB2	1.70	0.72
1:A:155:HIS:NE2	1:A:157:SER:HB3	2.01	0.72
1:C:249:VAL:HG13	1:C:250:PRO:HD2	1.71	0.72
1:C:5:LEU:O	1:C:6:ARG:HD3	1.89	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:13:SER:HB3	1:C:93:HIS:H	1.55	0.72
4:H:33:VAL:HB	4:H:54:ALA:CB	2.20	0.72
1:C:185:PRO:HD2	1:C:266:LEU:HD11	1.70	0.72
3:E:124:ALA:HB3	3:E:145:TYR:H	1.55	0.72
2:B:55:SER:HB3	2:B:63:TYR:CZ	2.25	0.72
2:D:54:LEU:HD11	2:D:62:PHE:CD2	2.25	0.72
1:A:39:LYS:O	1:A:40:GLU:HG3	1.89	0.71
1:A:215:LEU:CD2	1:A:261:VAL:HG13	2.20	0.71
3:E:61:SER:HB2	3:E:72:THR:O	1.88	0.71
4:F:213:GLU:O	4:F:215:ASN:N	2.23	0.71
4:H:167:LYS:HE2	4:H:209:ILE:CD1	2.20	0.71
4:F:121:ASP:O	4:F:122:LYS:C	2.31	0.71
4:H:158:PHE:N	4:H:159:PRO:CD	2.49	0.71
3:E:7:SER:HB3	3:E:22:ARG:HB2	1.72	0.71
1:C:129:ASN:CB	1:C:130:PRO:HD3	2.18	0.71
3:G:147:LYS:HG3	3:G:174:TYR:HE2	1.55	0.71
4:H:134:ILE:N	5:L:1:NAG:H62	2.05	0.71
1:A:251:LEU:HD23	1:A:251:LEU:H	1.54	0.71
4:F:132:PRO:O	4:F:133:THR:OG1	2.09	0.71
4:H:22:SER:HB3	4:H:80:ILE:HG12	1.71	0.71
4:H:41:LYS:HB3	4:H:42:PRO:CD	2.19	0.71
1:C:52:LEU:HD23	1:C:174:LYS:C	2.15	0.71
3:E:138:ALA:HB2	3:E:179:LEU:HD23	1.71	0.71
4:F:4:LEU:HD22	4:F:98:TYR:HD2	1.53	0.71
4:F:6:GLN:HE21	4:F:110:ALA:CB	2.02	0.71
3:E:179:LEU:HD13	4:F:153:LEU:HD22	1.73	0.71
1:A:50:PRO:O	1:A:52:LEU:N	2.24	0.70
1:A:208:PHE:CE2	1:A:213:ILE:HD12	2.25	0.70
2:B:17:ASN:HB2	2:B:73:THR:HA	1.72	0.70
1:C:99:GLN:HA	1:C:113:TYR:O	1.92	0.70
1:C:189:VAL:HB	1:C:274:TRP:CB	2.22	0.70
3:E:48:TYR:O	3:E:49:LEU:HD23	1.91	0.70
3:G:24:ASN:HD22	7:M:1:NAG:C1	1.89	0.70
3:G:139:CYS:O	3:G:140:LEU:HD23	1.91	0.70
1:C:184:PRO:HB3	1:C:266:LEU:CD2	2.21	0.70
2:D:41:LYS:HB2	2:D:46:ILE:HD11	1.73	0.70
4:F:228:ILE:HG22	4:F:229:LYS:N	2.06	0.70
3:E:132:LYS:H	3:E:132:LYS:CE	2.02	0.70
3:G:7:SER:HB2	3:G:8:PRO:CD	2.22	0.70
1:C:124:PRO:O	3:G:92:THR:HG23	1.91	0.70
1:C:34:LEU:HD23	1:C:35:ARG:N	2.05	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:81:LEU:HB3	1:A:118:TYR:CE1	2.27	0.70
4:F:131:LYS:HD2	4:F:156:ALA:HB3	1.73	0.70
1:A:57:ALA:C	1:A:59:ASN:H	1.98	0.69
4:F:55:THR:HG22	4:F:56:PRO:HD3	1.72	0.69
3:G:48:TYR:CE1	4:H:105:PHE:O	2.44	0.69
1:C:203:CYS:HB2	1:C:217:TRP:CZ2	2.28	0.69
3:G:58:ARG:HH12	3:G:79:GLN:HB2	1.57	0.69
1:C:66:ILE:HD13	1:C:160:LEU:HB2	1.72	0.69
4:H:62:ASP:HA	4:H:66:LYS:O	1.92	0.69
1:A:129:ASN:CB	1:A:130:PRO:HD3	2.23	0.69
1:C:213:ILE:HG12	1:C:214:THR:N	2.07	0.69
1:A:61:GLU:HA	1:A:64:THR:OG1	1.92	0.69
1:A:160:LEU:HD12	1:A:160:LEU:O	1.93	0.69
2:B:42:ASN:HD21	2:B:76:ASP:HA	1.57	0.69
3:E:144:PHE:CE1	3:E:149:VAL:HG21	2.27	0.69
4:H:151:LEU:HD23	4:H:151:LEU:N	2.08	0.69
1:A:6:ARG:HB3	1:A:8:PHE:HE1	1.58	0.69
1:C:105:PRO:O	1:C:108:HIS:N	2.24	0.69
4:H:157:PHE:CD1	4:H:159:PRO:HD2	2.27	0.69
1:C:167:LEU:HD21	1:C:171:TYR:CE2	2.28	0.68
1:A:202:ARG:HA	1:A:246:ALA:HB2	1.75	0.68
4:H:131:LYS:H	4:H:155:GLU:H	1.39	0.68
1:C:211:ALA:HB2	1:C:241:PHE:CE1	2.29	0.68
3:E:133:ASN:CG	5:I:1:NAG:H1	2.07	0.68
1:C:49:ALA:HB1	1:C:51:TRP:NE1	2.07	0.68
1:C:184:PRO:HG3	1:C:265:GLY:O	1.94	0.68
3:E:126:PRO:CG	3:E:144:PHE:HB3	2.19	0.68
4:H:208:CYS:O	4:H:208:CYS:SG	2.51	0.68
1:A:51:TRP:O	1:A:51:TRP:HE3	1.74	0.68
1:C:97:TRP:CD1	1:C:98:LEU:N	2.62	0.68
1:A:81:LEU:HB3	1:A:118:TYR:HE1	1.59	0.68
3:E:179:LEU:CD1	4:F:153:LEU:HD22	2.24	0.68
3:G:25:PHE:O	3:G:69:ARG:HB2	1.94	0.68
2:B:9:VAL:HG21	2:B:80:CYS:HB2	1.76	0.67
4:F:31:ARG:O	4:F:54:ALA:HB1	1.94	0.67
4:F:53:LEU:HD13	4:F:59:ILE:HD11	1.76	0.67
1:A:5:LEU:HD12	1:A:6:ARG:H	1.59	0.67
1:C:76:ARG:O	1:C:79:MET:HB2	1.94	0.67
3:E:184:ASP:HB2	3:E:187:SER:OG	1.93	0.67
4:F:52:VAL:HG22	4:F:71:SER:HB2	1.75	0.67
2:B:73:THR:O	2:B:75:LYS:N	2.27	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:E:14:HIS:HA	3:E:117:GLU:O	1.95	0.67
4:F:28:GLU:C	4:F:30:PHE:H	2.01	0.67
4:H:153:LEU:HD23	4:H:153:LEU:C	2.20	0.67
3:G:12:SER:HB2	3:G:115:THR:HB	1.77	0.67
4:H:3:LYS:HE3	4:H:98:TYR:OH	1.95	0.67
4:H:60:PHE:N	4:H:60:PHE:CD2	2.62	0.67
4:F:23:CYS:CB	4:F:96:CYS:SG	2.83	0.67
1:C:63:GLN:O	1:C:67:VAL:HG23	1.95	0.67
4:F:88:GLU:C	4:F:90:ASP:H	2.02	0.67
4:F:130:PRO:CD	4:F:215:ASN:HB2	2.25	0.67
1:A:14:ARG:HB2	1:A:15:PRO:CD	2.25	0.66
3:E:24:ASN:HD21	6:J:1:NAG:HN2	1.42	0.66
4:F:30:PHE:CZ	4:F:77:SER:HA	2.31	0.66
1:A:219:LEU:O	1:A:221:GLY:CA	2.42	0.66
1:C:21:TRP:CH2	1:C:23:ILE:HD11	2.30	0.66
2:D:17:ASN:OD1	2:D:74:GLU:HB2	1.96	0.66
4:F:185:THR:CG2	6:K:1:NAG:H82	2.24	0.66
3:G:3:GLN:NE2	3:G:27:THR:H	1.92	0.66
3:G:19:SER:O	3:G:75:ILE:N	2.23	0.66
1:C:150:ASN:HB3	1:C:153:VAL:HG21	1.76	0.66
3:G:139:CYS:C	3:G:140:LEU:HD23	2.21	0.66
3:G:98:GLY:C	3:G:99:TYR:CD1	2.73	0.66
4:F:153:LEU:HD12	4:F:153:LEU:C	2.20	0.66
1:C:45:MET:CE	1:C:54:GLN:HE22	2.09	0.66
1:A:23:ILE:HG22	1:A:24:ILE:N	2.10	0.65
3:E:133:ASN:CG	5:I:1:NAG:C1	2.67	0.65
4:F:51:TYR:O	4:F:59:ILE:HD12	1.96	0.65
4:F:32:SER:HA	4:F:54:ALA:HB1	1.78	0.65
1:A:128:GLU:O	1:A:129:ASN:C	2.39	0.65
1:C:254:GLU:O	1:C:256:SER:N	2.29	0.65
4:F:179:GLU:HA	4:F:193:SER:HA	1.77	0.65
3:G:96:SER:OG	3:G:100:GLU:N	2.30	0.65
4:H:63:LYS:O	4:H:63:LYS:HG3	1.96	0.65
1:A:11:ALA:O	1:A:94:THR:HA	1.96	0.65
2:D:16:GLU:O	2:D:16:GLU:HG3	1.95	0.65
1:C:185:PRO:HD2	1:C:266:LEU:CG	2.27	0.65
1:A:6:ARG:NH1	1:A:102:ASP:OD1	2.30	0.65
1:C:170:LYS:HA	1:C:170:LYS:HE2	1.79	0.65
1:C:185:PRO:HD2	1:C:266:LEU:CD1	2.26	0.65
4:F:146:LYS:HG3	4:F:199:GLU:HG3	1.77	0.65
3:G:7:SER:OG	3:G:8:PRO:HD2	1.97	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:65:ARG:HH12	1:A:158:GLN:HG3	1.62	0.64
3:G:58:ARG:HB2	3:G:76:ARG:O	1.96	0.64
4:H:33:VAL:CG1	4:H:34:THR:N	2.59	0.64
1:A:219:LEU:C	1:A:221:GLY:N	2.49	0.64
1:C:191:ARG:HB3	1:C:274:TRP:HZ3	1.62	0.64
2:D:5:PRO:CB	2:D:30:PHE:HB3	2.27	0.64
4:H:107:LYS:HG3	4:H:109:PHE:HE1	1.59	0.64
3:E:94:HIS:C	3:E:95:ILE:HG13	2.21	0.64
1:C:8:PHE:CD1	1:C:8:PHE:N	2.65	0.64
1:C:191:ARG:H	1:C:274:TRP:HZ3	1.45	0.64
1:C:200:THR:HA	1:C:247:VAL:O	1.98	0.64
4:F:157:PHE:O	4:F:188:THR:HG22	1.96	0.64
4:F:158:PHE:O	4:F:159:PRO:C	2.40	0.64
3:G:29:MET:HE3	3:G:90:ALA:HB3	1.80	0.64
4:H:35:ILE:O	4:H:51:TYR:HA	1.96	0.64
2:B:30:PHE:HE1	2:B:33:SER:HA	1.62	0.64
1:C:189:VAL:CG1	1:C:190:THR:H	2.09	0.64
1:A:101:CYS:HA	1:A:111:LEU:O	1.97	0.64
1:A:213:ILE:HG13	1:A:263:HIS:HD2	1.61	0.64
1:C:119:ASP:O	1:C:120:SER:HB3	1.97	0.64
4:F:52:VAL:HG22	4:F:71:SER:CB	2.28	0.64
4:F:147:THR:HG22	4:F:148:GLY:H	1.60	0.64
4:H:93:ILE:HG12	4:H:114:LYS:CB	2.27	0.64
1:A:73:LEU:HD12	3:E:99:TYR:C	2.23	0.63
1:A:190:THR:HA	1:A:274:TRP:HZ3	1.63	0.63
4:F:37:TRP:CD2	4:F:81:LEU:HD13	2.32	0.63
4:F:111:GLU:OE2	4:F:111:GLU:N	2.31	0.63
3:G:152:SER:O	3:G:153:LEU:HD23	1.98	0.63
1:C:167:LEU:O	1:C:167:LEU:HD23	1.98	0.63
2:D:79:ALA:HA	2:D:93:VAL:O	1.98	0.63
4:F:30:PHE:CZ	4:F:73:ASN:OD1	2.52	0.63
4:F:157:PHE:CE1	4:F:189:TYR:HB2	2.34	0.63
1:A:110:CYS:O	1:A:111:LEU:HD12	1.98	0.63
1:C:67:VAL:O	1:C:70:GLN:HB3	1.98	0.63
1:C:189:VAL:HB	1:C:274:TRP:HB3	1.80	0.63
3:E:18:ASP:OD1	3:E:76:ARG:HA	1.98	0.63
1:C:8:PHE:HD2	2:D:56:PHE:CD1	2.17	0.63
1:A:45:MET:HE1	1:A:60:TRP:CG	2.33	0.63
2:B:70:PHE:CE2	2:B:72:PRO:HG3	2.34	0.63
2:D:31:HIS:O	2:D:32:PRO:C	2.39	0.63
3:G:17:THR:HG22	3:G:18:ASP:H	1.64	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:G:110:GLN:CD	3:G:110:GLN:H	2.07	0.63
1:A:24:ILE:HG22	1:A:24:ILE:O	1.97	0.63
1:C:208:PHE:CE2	1:C:241:PHE:O	2.52	0.63
4:F:134:ILE:HG23	4:F:225:PHE:CE2	2.34	0.63
1:A:14:ARG:CB	1:A:15:PRO:CD	2.76	0.62
1:A:66:ILE:HD13	1:A:160:LEU:HB2	1.81	0.62
1:C:111:LEU:N	1:C:111:LEU:CD1	2.61	0.62
3:E:30:ARG:HG3	3:E:31:SER:N	2.13	0.62
1:A:163:HIS:O	1:A:164:CYS:C	2.40	0.62
3:E:133:ASN:O	3:E:134:GLY:C	2.42	0.62
1:A:24:ILE:HD12	1:A:67:VAL:CG1	2.29	0.62
1:C:31:MET:HA	1:C:31:MET:CE	2.21	0.62
2:D:17:ASN:HB2	2:D:73:THR:HA	1.80	0.62
3:E:46:LEU:HD22	3:E:59:LEU:CD1	2.29	0.62
3:G:89:ALA:HA	3:G:107:VAL:O	1.98	0.62
4:H:51:TYR:O	4:H:59:ILE:HD12	1.99	0.62
1:C:261:VAL:HG12	1:C:261:VAL:O	1.99	0.62
4:F:186:ASN:HB2	6:K:1:NAG:H1	1.80	0.62
1:C:191:ARG:N	1:C:274:TRP:CZ3	2.68	0.62
3:G:6:GLN:HA	3:G:23:CYS:HB2	1.80	0.62
1:C:249:VAL:CG1	1:C:250:PRO:HD2	2.29	0.62
3:G:67:GLU:OE1	3:G:69:ARG:NE	2.32	0.62
4:H:56:PRO:HD3	4:H:73:ASN:HB3	1.80	0.62
2:B:57:SER:O	2:B:60:TRP:N	2.33	0.62
1:A:234:ARG:NH1	1:A:242:GLN:OE1	2.33	0.62
2:B:17:ASN:CB	2:B:73:THR:HA	2.30	0.62
2:D:70:PHE:CZ	2:D:72:PRO:HG3	2.34	0.62
3:E:30:ARG:NH2	3:E:103:THR:HG21	2.15	0.61
3:G:26:THR:HA	3:G:69:ARG:HB2	1.82	0.61
1:C:230:LEU:HD21	1:C:245:ALA:HB2	1.82	0.61
3:E:17:THR:HG22	3:E:18:ASP:N	2.15	0.61
2:D:5:PRO:CA	2:D:30:PHE:HB3	2.31	0.61
2:B:3:ARG:O	2:B:30:PHE:HA	1.99	0.61
2:D:59:ASP:OD1	2:D:61:SER:CB	2.49	0.61
3:E:24:ASN:CG	6:J:1:NAG:C1	2.65	0.61
4:F:88:GLU:O	4:F:90:ASP:N	2.29	0.61
1:A:103:VAL:HG12	1:A:104:GLU:N	2.14	0.61
1:A:234:ARG:NH1	2:B:10:TYR:CG	2.69	0.61
1:C:249:VAL:HG22	1:C:257:TYR:CE2	2.36	0.61
3:E:134:GLY:O	3:E:135:THR:HB	2.00	0.61
4:F:14:PRO:HB2	4:F:17:GLU:HG3	1.82	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:F:158:PHE:O	4:F:160:ASP:N	2.34	0.61
3:G:8:PRO:O	3:G:112:THR:OG1	2.11	0.61
4:F:25:VAL:HG21	4:F:35:ILE:CD1	2.29	0.61
3:G:106:LEU:HD22	4:H:107:LYS:HE2	1.83	0.61
2:D:84:HIS:O	2:D:86:THR:N	2.34	0.61
3:G:25:PHE:O	3:G:69:ARG:O	2.18	0.61
4:H:75:SER:C	4:H:77:SER:H	2.08	0.61
2:D:41:LYS:HG2	2:D:42:ASN:ND2	2.16	0.61
1:C:73:LEU:HD12	3:G:99:TYR:HB2	1.81	0.60
4:H:132:PRO:HG2	4:H:221:GLN:HE22	1.65	0.60
1:A:115:GLN:C	1:A:116:LEU:HD12	2.26	0.60
1:C:24:ILE:HG22	1:C:24:ILE:O	2.00	0.60
1:C:84:PHE:HE1	3:G:30:ARG:NH2	1.92	0.60
4:F:63:LYS:HG2	4:F:66:LYS:O	2.01	0.60
3:G:128:VAL:HG13	3:G:140:LEU:O	2.01	0.60
4:H:56:PRO:C	4:H:58:HIS:H	2.07	0.60
1:A:71:GLY:O	1:A:74:SER:HB3	2.01	0.60
2:B:9:VAL:CG2	2:B:80:CYS:HB2	2.30	0.60
4:F:192:PHE:HD1	4:F:192:PHE:H	1.49	0.60
3:G:143:ASP:HB3	3:G:173:LYS:HD2	1.83	0.60
1:A:219:LEU:HD21	1:A:257:TYR:CE2	2.35	0.60
1:C:189:VAL:HG12	1:C:190:THR:H	1.64	0.60
3:E:132:LYS:HE2	3:E:132:LYS:N	2.05	0.60
3:G:133:ASN:C	3:G:133:ASN:OD1	2.44	0.60
1:A:9:TYR:OH	1:A:74:SER:CB	2.49	0.60
1:A:34:LEU:CD1	1:A:63:GLN:OE1	2.49	0.60
1:A:152:THR:HG22	1:A:153:VAL:H	1.66	0.60
1:C:105:PRO:O	1:C:106:ASP:C	2.44	0.60
3:E:12:SER:C	3:E:13:LEU:HD23	2.26	0.60
4:F:211:LYS:HE2	4:F:220:ASP:OD1	2.02	0.60
3:G:29:MET:HE3	3:G:90:ALA:CB	2.30	0.60
1:C:217:TRP:HD1	1:C:228:MET:HE3	1.66	0.60
3:G:137:VAL:N	3:G:180:GLY:O	2.30	0.60
4:H:105:PHE:CZ	4:H:107:LYS:NZ	2.68	0.60
1:A:215:LEU:HD22	1:A:261:VAL:HG22	1.83	0.60
4:F:139:VAL:O	4:F:143:ASN:HB2	2.01	0.60
4:F:6:GLN:HG2	4:F:23:CYS:HB2	1.83	0.60
1:A:235:PRO:HG2	2:B:65:LEU:HD13	1.83	0.60
1:C:80:THR:HB	4:H:103:SER:CB	2.24	0.60
4:F:149:THR:HA	4:F:195:LEU:O	2.02	0.60
4:F:169:LYS:CG	4:F:207:ARG:HB2	2.32	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:M:1:NAG:H83	7:M:1:NAG:H3	1.84	0.60
1:C:32:GLN:NE2	2:D:53:ASP:CG	2.54	0.60
3:E:133:ASN:O	3:E:136:ASN:HB2	2.02	0.60
1:A:22:PHE:HZ	1:A:36:PHE:HE2	1.48	0.59
1:A:58:ASP:O	1:A:61:GLU:HG3	2.02	0.59
1:C:254:GLU:C	1:C:256:SER:H	2.10	0.59
4:F:51:TYR:H	4:F:51:TYR:HD2	1.48	0.59
1:A:6:ARG:HB3	1:A:8:PHE:CE1	2.36	0.59
1:A:9:TYR:HB3	1:A:97:TRP:HB3	1.84	0.59
1:A:263:HIS:ND1	1:A:264:GLU:N	2.50	0.59
1:C:4:SER:O	1:C:28:VAL:HG13	2.01	0.59
3:G:94:HIS:HB2	3:G:104:ASP:N	2.17	0.59
1:A:69:ILE:O	1:A:72:GLN:N	2.35	0.59
1:A:129:ASN:HB2	1:A:130:PRO:CD	2.27	0.59
3:E:35:PHE:CD1	3:E:108:PHE:HE2	2.20	0.59
1:A:9:TYR:OH	1:A:74:SER:HB2	2.02	0.59
1:A:187:ALA:HB1	1:A:272:LEU:HD11	1.83	0.59
1:C:34:LEU:HD23	1:C:34:LEU:C	2.27	0.59
1:C:129:ASN:CB	1:C:130:PRO:CD	2.68	0.59
1:C:160:LEU:O	1:C:160:LEU:HD12	2.02	0.59
4:F:156:ALA:CA	4:F:188:THR:HG21	2.31	0.59
3:G:3:GLN:NE2	3:G:27:THR:N	2.51	0.59
3:G:6:GLN:NE2	3:G:88:CYS:H	1.99	0.59
3:G:15:GLU:HB2	3:G:118:PRO:HA	1.84	0.59
1:A:24:ILE:CD1	1:A:67:VAL:HG13	2.32	0.59
1:A:28:VAL:HG11	1:A:179:LEU:HD13	1.84	0.59
1:A:230:LEU:HD22	1:A:243:LYS:HE3	1.84	0.59
2:D:54:LEU:HD12	2:D:54:LEU:C	2.26	0.59
3:E:195:ASN:O	3:E:196:SER:OG	2.16	0.59
4:F:150:TYR:HB3	4:F:225:PHE:CE2	2.37	0.59
3:G:3:GLN:O	3:G:3:GLN:CG	2.48	0.59
4:H:55:THR:O	4:H:57:THR:N	2.36	0.59
4:H:168:GLU:HG3	4:H:175:LEU:CD1	2.32	0.59
1:A:22:PHE:CD1	1:A:22:PHE:C	2.81	0.59
3:E:26:THR:O	3:E:26:THR:CG2	2.51	0.59
3:E:164:PRO:HG3	3:E:178:LYS:CD	2.32	0.59
4:F:192:PHE:HD1	4:F:192:PHE:N	2.00	0.59
7:M:1:NAG:O1	7:M:1:NAG:C8	2.51	0.59
3:E:46:LEU:HD22	3:E:59:LEU:HD12	1.84	0.59
4:H:165:TYR:CD1	4:H:165:TYR:C	2.80	0.59
1:A:175:GLY:O	1:A:177:GLU:N	2.36	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:202:ARG:HD2	2:B:98:ASP:OD1	2.01	0.59
4:H:168:GLU:HG3	4:H:175:LEU:HD11	1.84	0.59
1:A:55:GLU:O	1:A:56:GLU:C	2.46	0.59
1:A:266:LEU:CD2	1:A:270:LEU:HD13	2.26	0.59
1:C:228:MET:CE	1:C:245:ALA:CB	2.76	0.59
3:E:148:GLU:H	3:E:148:GLU:CD	2.11	0.59
1:A:234:ARG:HG3	2:B:10:TYR:CE2	2.38	0.59
3:E:137:VAL:HG21	3:E:188:VAL:HG21	1.84	0.59
4:F:137:PRO:HD3	4:F:150:TYR:CE2	2.37	0.59
4:F:192:PHE:N	4:F:192:PHE:CD1	2.70	0.59
1:A:62:GLN:O	1:A:66:ILE:HG13	2.03	0.58
1:A:36:PHE:CD2	1:A:36:PHE:C	2.81	0.58
3:E:12:SER:O	3:E:13:LEU:HD23	2.02	0.58
4:H:41:LYS:HD3	4:H:92:ALA:HB2	1.84	0.58
1:A:202:ARG:HG2	1:A:204:TRP:NE1	2.18	0.58
1:C:9:TYR:C	1:C:9:TYR:CD2	2.81	0.58
1:C:264:GLU:C	1:C:266:LEU:H	2.10	0.58
4:F:186:ASN:CB	6:K:1:NAG:H1	2.23	0.58
1:A:178:ARG:O	1:A:181:ARG:HD2	2.03	0.58
2:B:65:LEU:O	2:B:65:LEU:HD23	2.04	0.58
1:A:153:VAL:CB	1:A:154:PRO:CD	2.80	0.58
1:C:189:VAL:HB	1:C:274:TRP:HB2	1.85	0.58
4:H:37:TRP:CD2	4:H:81:LEU:HD13	2.38	0.58
3:G:26:THR:HG23	3:G:69:ARG:HH11	1.69	0.58
2:D:96:ASP:OD2	2:D:98:ASP:HB2	2.04	0.58
4:H:120:SER:O	4:H:121:ASP:HB2	2.04	0.58
1:A:3:HIS:CE1	1:A:180:LEU:HD21	2.39	0.58
3:G:11:LEU:CD2	3:G:13:LEU:HD21	2.23	0.58
4:H:215:ASN:OD1	4:H:215:ASN:O	2.22	0.58
4:F:204:LYS:O	4:F:227:SER:HB3	2.04	0.58
1:A:63:GLN:NE2	1:A:167:LEU:HD22	2.18	0.57
4:H:6:GLN:HG2	4:H:23:CYS:CB	2.34	0.57
7:M:2:NAG:H82	7:M:2:NAG:H3	1.85	0.57
4:F:28:GLU:C	4:F:30:PHE:N	2.61	0.57
4:H:167:LYS:HE2	4:H:209:ILE:HD11	1.85	0.57
4:H:210:VAL:CG1	4:H:211:LYS:H	2.17	0.57
1:A:21:TRP:CZ2	1:A:23:ILE:HG13	2.38	0.57
1:A:110:CYS:C	1:A:111:LEU:HD12	2.30	0.57
2:D:26:TYR:HB2	2:D:65:LEU:HD13	1.86	0.57
3:E:34:TRP:CE3	3:E:88:CYS:HB3	2.39	0.57
1:A:71:GLY:O	1:A:74:SER:CB	2.52	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:F:165:TYR:CE1	4:F:167:LYS:HD2	2.37	0.57
4:F:228:ILE:HG22	4:F:229:LYS:H	1.69	0.57
4:H:22:SER:CB	4:H:80:ILE:HG12	2.33	0.57
4:H:33:VAL:CB	4:H:54:ALA:HB2	2.32	0.57
1:C:110:CYS:C	1:C:111:LEU:HD12	2.30	0.57
3:G:6:GLN:NE2	3:G:88:CYS:N	2.52	0.57
3:G:24:ASN:ND2	7:M:1:NAG:N2	2.51	0.57
4:H:53:LEU:O	4:H:53:LEU:HG	2.05	0.57
1:A:12:VAL:HG12	1:A:13:SER:N	2.20	0.57
1:C:81:LEU:HB3	1:C:118:TYR:CE1	2.40	0.57
1:C:233:THR:HB	1:C:243:LYS:HD2	1.87	0.57
2:D:22:PHE:HA	2:D:68:THR:O	2.04	0.57
1:A:253:LYS:O	1:A:255:GLN:N	2.38	0.57
4:H:99:GLY:HA3	4:H:105:PHE:CD2	2.40	0.57
3:G:33:GLN:NE2	3:G:47:PHE:O	2.38	0.57
3:G:66:LYS:O	3:G:67:GLU:HG3	2.05	0.57
1:A:234:ARG:HD3	1:A:242:GLN:OE1	2.05	0.56
1:C:5:LEU:HD13	1:C:28:VAL:HG22	1.87	0.56
3:E:96:SER:O	3:E:98:GLY:N	2.38	0.56
3:E:36:ARG:HB2	3:E:46:LEU:HD11	1.86	0.56
4:H:46:LEU:HD11	4:H:95:TYR:CE1	2.40	0.56
1:C:249:VAL:HG13	1:C:257:TYR:CZ	2.39	0.56
2:D:-1:ASP:HB2	2:D:0:PRO:HD3	1.87	0.56
2:D:31:HIS:HB3	2:D:32:PRO:CD	2.32	0.56
3:E:153:LEU:HD12	3:E:182:TYR:OH	2.04	0.56
4:F:206:HIS:O	4:F:225:PHE:HB2	2.06	0.56
3:G:132:LYS:HA	4:H:135:PHE:HE2	1.69	0.56
2:D:59:ASP:OD1	2:D:59:ASP:C	2.48	0.56
3:E:146:PRO:HD2	3:E:194:HIS:CE1	2.39	0.56
4:F:185:THR:O	4:F:186:ASN:C	2.49	0.56
3:G:33:GLN:OE1	3:G:48:TYR:CD1	2.59	0.56
1:C:209:TYR:HB3	1:C:210:PRO:HD3	1.88	0.56
4:H:165:TYR:CE1	4:H:209:ILE:HD12	2.40	0.56
3:E:6:GLN:HE22	3:E:87:PHE:HA	1.69	0.56
3:E:80:LEU:HD23	3:E:116:VAL:O	2.05	0.56
3:E:201:SER:OG	3:E:202:THR:N	2.38	0.56
4:F:128:ILE:HG22	4:F:158:PHE:N	2.21	0.56
4:H:37:TRP:CG	4:H:81:LEU:HD13	2.40	0.56
1:A:16:GLY:C	1:A:18:GLY:H	2.14	0.56
1:C:12:VAL:HG12	1:C:13:SER:N	2.19	0.56
2:D:11:SER:HB3	2:D:95:TRP:CZ2	2.40	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:E:134:GLY:O	3:E:135:THR:CB	2.54	0.56
3:E:189:THR:HG23	3:E:201:SER:OG	2.06	0.56
4:F:158:PHE:HB3	4:F:159:PRO:HD3	1.88	0.56
1:C:194:ARG:HB2	1:C:198:ASP:C	2.29	0.56
3:E:14:HIS:NE2	3:G:9:SER:OG	2.39	0.56
4:F:23:CYS:HB3	4:F:96:CYS:SG	2.46	0.56
4:H:57:THR:HB	4:H:59:ILE:HG13	1.87	0.56
1:A:74:SER:O	1:A:78:LEU:HB2	2.06	0.56
2:B:23:LEU:CD2	2:B:39:LEU:HD22	2.36	0.56
1:C:191:ARG:HD2	1:C:275:GLY:O	2.06	0.56
1:C:203:CYS:SG	1:C:272:LEU:HD12	2.46	0.56
3:E:3:GLN:CB	3:E:26:THR:HB	2.36	0.56
4:H:60:PHE:N	4:H:60:PHE:HD2	2.03	0.56
1:A:236:ALA:HB2	1:A:242:GLN:HE21	1.70	0.56
2:D:84:HIS:CG	2:D:85:VAL:N	2.74	0.56
4:H:28:GLU:C	4:H:30:PHE:N	2.61	0.56
2:D:24:ASN:ND2	2:D:67:TYR:CB	2.69	0.55
4:H:56:PRO:C	4:H:58:HIS:N	2.60	0.55
4:H:134:ILE:HD11	4:H:223:ILE:CG2	2.36	0.55
2:D:16:GLU:O	2:D:17:ASN:C	2.47	0.55
3:E:38:ASN:O	3:E:39:SER:C	2.49	0.55
3:E:130:ILE:HD11	3:E:201:SER:OG	2.06	0.55
3:G:3:GLN:O	3:G:25:PHE:HB3	2.06	0.55
2:D:74:GLU:HG2	2:D:74:GLU:O	2.06	0.55
3:E:108:PHE:CE1	4:F:46:LEU:HB2	2.41	0.55
1:A:105:PRO:C	1:A:108:HIS:H	2.15	0.55
1:A:202:ARG:HB2	1:A:202:ARG:CZ	2.37	0.55
1:C:184:PRO:HB3	1:C:266:LEU:HD23	1.86	0.55
2:D:79:ALA:HB2	2:D:94:LYS:HA	1.88	0.55
3:E:55:GLU:HB2	3:E:60:LYS:HG2	1.88	0.55
3:E:185:SER:C	3:E:188:VAL:HG23	2.30	0.55
1:C:194:ARG:CD	1:C:198:ASP:HB2	2.33	0.55
4:F:146:LYS:CG	4:F:199:GLU:HG3	2.36	0.55
4:F:121:ASP:O	4:F:123:ARG:N	2.40	0.55
4:F:212:HIS:O	4:F:215:ASN:HB3	2.07	0.55
4:F:213:GLU:HG2	4:F:214:ASN:N	2.21	0.55
1:A:116:LEU:HD12	1:A:116:LEU:N	2.22	0.55
1:A:163:HIS:O	1:A:166:ASP:N	2.40	0.55
2:B:50:GLU:HB2	2:B:67:TYR:CZ	2.42	0.55
1:C:49:ALA:HB1	1:C:51:TRP:CD1	2.42	0.55
1:C:266:LEU:CD1	1:C:270:LEU:HD22	2.36	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:G:30:ARG:NH1	3:G:91:ASP:OD2	2.40	0.55
3:G:53:THR:HG22	3:G:54:LYS:N	2.21	0.55
1:A:97:TRP:HD1	1:A:98:LEU:N	2.05	0.55
2:B:36:GLU:O	2:B:82:VAL:HA	2.07	0.55
4:H:51:TYR:O	4:H:51:TYR:CD2	2.60	0.55
1:A:215:LEU:HD23	1:A:261:VAL:HG13	1.88	0.55
1:C:6:ARG:HD3	1:C:100:GLY:HA3	1.89	0.55
2:D:17:ASN:N	2:D:17:ASN:ND2	2.49	0.55
4:F:55:THR:HB	4:F:56:PRO:HD2	1.89	0.55
1:C:13:SER:HB3	1:C:93:HIS:N	2.22	0.55
2:D:54:LEU:HD12	2:D:55:SER:N	2.21	0.55
3:G:27:THR:O	3:G:28:THR:C	2.50	0.55
2:B:84:HIS:H	2:B:87:LEU:HD22	1.71	0.54
3:G:11:LEU:C	3:G:11:LEU:CD2	2.74	0.54
1:A:249:VAL:HG13	1:A:257:TYR:CE1	2.43	0.54
4:F:157:PHE:CE1	4:F:189:TYR:CB	2.90	0.54
2:D:9:VAL:HG21	2:D:93:VAL:HB	1.88	0.54
4:F:159:PRO:HD3	4:F:214:ASN:HD22	1.71	0.54
2:D:73:THR:HB	2:D:76:ASP:OD2	2.07	0.54
1:C:1:GLY:O	1:C:3:HIS:CE1	2.60	0.54
1:C:13:SER:N	1:C:93:HIS:O	2.34	0.54
3:E:144:PHE:O	3:E:174:TYR:HB2	2.08	0.54
4:H:63:LYS:C	4:H:64:GLU:HG3	2.31	0.54
1:A:103:VAL:HG12	1:A:104:GLU:H	1.71	0.54
1:C:88:SER:C	1:C:90:ASP:H	2.15	0.54
2:D:7:ILE:HG23	2:D:27:VAL:HG22	1.88	0.54
3:E:7:SER:HB3	3:E:22:ARG:O	2.07	0.54
3:E:20:ALA:HB2	3:E:74:HIS:CD2	2.42	0.54
4:F:23:CYS:SG	4:F:96:CYS:CB	2.96	0.54
1:A:9:TYR:CD1	1:A:70:GLN:HG2	2.42	0.54
1:A:269:PRO:O	1:A:271:ILE:HD12	2.08	0.54
2:B:40:LEU:HD23	2:B:45:ARG:HA	1.90	0.54
1:C:99:GLN:HE21	1:C:160:LEU:HD23	1.69	0.54
1:C:204:TRP:CH2	2:D:99:MET:HA	2.43	0.54
4:F:134:ILE:HG23	4:F:225:PHE:CZ	2.43	0.54
3:E:96:SER:C	3:E:98:GLY:N	2.60	0.54
4:H:111:GLU:N	4:H:111:GLU:OE1	2.41	0.54
1:A:83:HIS:C	1:A:85:TYR:N	2.64	0.54
2:B:26:TYR:HD1	2:B:65:LEU:HB2	1.73	0.54
1:C:104:GLU:OE1	1:C:109:LEU:HD12	2.07	0.54
1:C:266:LEU:HD11	1:C:270:LEU:HD22	1.90	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:23:LEU:HB3	2:D:68:THR:CG2	2.35	0.54
2:D:84:HIS:ND1	2:D:84:HIS:C	2.66	0.54
3:E:124:ALA:HB3	3:E:145:TYR:N	2.21	0.54
4:F:25:VAL:C	4:F:26:PHE:HD1	2.15	0.54
3:G:94:HIS:HB2	3:G:104:ASP:CA	2.37	0.54
1:A:88:SER:C	1:A:90:ASP:H	2.16	0.54
3:E:76:ARG:O	3:E:77:ASP:C	2.51	0.54
3:E:204:PHE:O	3:E:205:GLU:C	2.51	0.54
4:F:186:ASN:CG	6:K:1:NAG:O1	2.37	0.54
1:C:8:PHE:N	1:C:8:PHE:HD1	2.05	0.53
2:D:12:ARG:HG2	2:D:13:HIS:CE1	2.43	0.53
4:F:169:LYS:HG3	4:F:207:ARG:HD2	1.89	0.53
4:H:132:PRO:CG	4:H:221:GLN:HE22	2.20	0.53
4:H:210:VAL:CG1	4:H:211:LYS:N	2.69	0.53
4:H:27:ILE:CD1	4:H:98:TYR:OH	2.55	0.53
4:H:51:TYR:CE2	4:H:59:ILE:HD13	2.43	0.53
1:C:52:LEU:HD23	1:C:174:LYS:O	2.06	0.53
3:E:89:ALA:HB2	3:E:108:PHE:CD2	2.43	0.53
1:A:187:ALA:CB	1:A:272:LEU:HD11	2.38	0.53
1:C:9:TYR:CD1	1:C:70:GLN:NE2	2.74	0.53
4:F:4:LEU:HD22	4:F:98:TYR:CD2	2.39	0.53
4:F:61:LEU:HD12	4:F:69:GLU:HG3	1.91	0.53
1:C:7:TYR:N	1:C:7:TYR:CD1	2.76	0.53
1:C:216:THR:HG22	1:C:260:HIS:HB2	1.89	0.53
1:C:92:SER:O	1:C:93:HIS:HD2	1.91	0.53
1:C:95:LEU:C	1:C:95:LEU:HD12	2.33	0.53
1:A:23:ILE:CG2	1:A:24:ILE:N	2.72	0.53
2:B:5:PRO:CA	2:B:30:PHE:HB3	2.39	0.53
1:C:68:THR:O	1:C:71:GLY:N	2.42	0.53
1:C:148:VAL:HG12	1:C:149:GLY:N	2.24	0.53
2:D:35:ILE:HG12	2:D:37:VAL:HG23	1.90	0.53
3:E:15:GLU:C	3:E:17:THR:H	2.17	0.53
3:E:24:ASN:OD1	3:E:24:ASN:C	2.48	0.53
4:F:33:VAL:CG1	4:F:34:THR:H	2.21	0.53
4:H:64:GLU:O	4:H:65:TYR:HB2	2.08	0.53
3:E:13:LEU:HB3	3:E:17:THR:HG21	1.90	0.53
4:H:134:ILE:HG22	4:H:135:PHE:N	2.23	0.53
1:C:61:GLU:O	1:C:64:THR:N	2.40	0.53
1:C:185:PRO:HD2	1:C:266:LEU:HG	1.90	0.53
1:C:252:GLY:N	1:C:254:GLU:OE2	2.42	0.53
4:F:50:LEU:CD2	4:F:61:LEU:HG	2.24	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:H:133:THR:HG22	4:H:153:LEU:O	2.09	0.53
1:A:157:SER:O	1:A:158:GLN:CG	2.57	0.53
1:A:208:PHE:CE2	1:A:242:GLN:HA	2.44	0.53
1:C:216:THR:CG2	1:C:260:HIS:HB2	2.39	0.53
4:F:38:TYR:CD2	4:F:48:PHE:HA	2.43	0.53
4:F:169:LYS:HG2	4:F:207:ARG:HB2	1.91	0.53
4:H:92:ALA:O	4:H:94:TYR:HE1	1.92	0.53
1:A:217:TRP:CD1	1:A:247:VAL:HG22	2.44	0.52
1:C:190:THR:HB	1:C:192:HIS:HE1	1.74	0.52
1:C:260:HIS:HB3	1:C:262:TYR:CE1	2.44	0.52
2:D:98:ASP:CB	2:D:99:MET:SD	2.98	0.52
1:C:242:GLN:OE1	2:D:10:TYR:CE1	2.62	0.52
3:E:2:ASP:CG	3:E:2:ASP:O	2.52	0.52
3:E:15:GLU:O	3:E:17:THR:N	2.41	0.52
4:F:153:LEU:HD11	4:F:155:GLU:HG2	1.90	0.52
2:B:21:ASN:O	2:B:70:PHE:O	2.28	0.52
1:C:12:VAL:O	1:C:20:PRO:HB3	2.10	0.52
4:F:55:THR:HG22	4:F:56:PRO:CD	2.38	0.52
4:F:185:THR:O	4:F:188:THR:N	2.42	0.52
1:A:166:ASP:O	1:A:169:GLN:HB3	2.10	0.52
1:A:216:THR:HG22	1:A:260:HIS:HB2	1.92	0.52
3:G:6:GLN:HE22	3:G:88:CYS:N	2.07	0.52
1:A:83:HIS:C	1:A:85:TYR:H	2.16	0.52
1:A:105:PRO:C	1:A:107:ARG:N	2.68	0.52
1:A:207:GLY:HA2	1:A:240:THR:HB	1.90	0.52
3:E:200:HIS:HB2	3:E:203:ASP:OD2	2.10	0.52
1:A:8:PHE:HB2	1:A:25:VAL:CG2	2.40	0.52
1:A:31:MET:HG3	1:A:179:LEU:HD23	1.90	0.52
1:A:57:ALA:C	1:A:59:ASN:N	2.66	0.52
1:A:255:GLN:C	1:A:257:TYR:H	2.18	0.52
1:C:204:TRP:HH2	2:D:99:MET:HA	1.75	0.52
3:E:145:TYR:CD2	3:E:146:PRO:N	2.77	0.52
4:F:30:PHE:CE2	4:F:77:SER:HA	2.45	0.52
4:F:190:MET:HG2	4:F:191:LYS:N	2.23	0.52
3:G:11:LEU:HD23	3:G:12:SER:N	2.24	0.52
3:G:106:LEU:HB3	4:H:38:TYR:OH	2.09	0.52
1:A:61:GLU:O	1:A:64:THR:N	2.43	0.52
2:D:9:VAL:CG2	2:D:93:VAL:HB	2.39	0.52
3:E:3:GLN:HB2	3:E:26:THR:HB	1.92	0.52
3:G:20:ALA:HA	3:G:74:HIS:HA	1.91	0.52
3:E:133:ASN:HD21	5:I:1:NAG:H1	0.36	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:F:49:LEU:HD13	4:F:69:GLU:CD	2.35	0.52
3:G:24:ASN:HB3	3:G:70:TYR:HD1	1.73	0.52
4:H:8:GLU:OE2	4:H:213:GLU:HG3	2.09	0.52
1:A:45:MET:HE1	1:A:60:TRP:CD2	2.44	0.52
3:E:20:ALA:CB	3:E:74:HIS:CD2	2.93	0.52
4:F:185:THR:HG23	6:K:1:NAG:C7	2.40	0.52
4:H:14:PRO:O	4:H:17:GLU:HB2	2.10	0.52
4:H:54:ALA:HA	4:H:73:ASN:HD22	1.75	0.52
1:A:204:TRP:CH2	2:B:98:ASP:O	2.63	0.51
1:C:187:ALA:CB	1:C:272:LEU:HD11	2.39	0.51
3:E:145:TYR:CB	3:E:146:PRO:HD3	2.39	0.51
1:A:65:ARG:NH1	1:A:158:GLN:HG3	2.24	0.51
1:A:79:MET:C	1:A:81:LEU:H	2.19	0.51
4:F:141:GLU:O	4:F:142:THR:C	2.53	0.51
3:G:137:VAL:O	3:G:180:GLY:N	2.43	0.51
1:C:28:VAL:O	1:C:29:ASP:C	2.53	0.51
2:D:27:VAL:HG11	2:D:35:ILE:CD1	2.41	0.51
4:F:33:VAL:CG1	4:F:34:THR:N	2.73	0.51
4:F:79:SER:O	4:F:80:ILE:HG13	2.09	0.51
1:C:184:PRO:HB3	1:C:266:LEU:HD21	1.92	0.51
4:F:228:ILE:CG2	4:F:229:LYS:N	2.73	0.51
1:A:169:GLN:O	1:A:170:LYS:C	2.52	0.51
1:C:73:LEU:CD1	3:G:99:TYR:H	2.24	0.51
4:F:173:THR:HG22	4:F:174:ILE:N	2.25	0.51
3:G:14:HIS:HA	3:G:117:GLU:O	2.10	0.51
1:C:152:THR:HG22	1:C:152:THR:O	2.10	0.51
2:D:98:ASP:HB3	2:D:99:MET:SD	2.49	0.51
1:A:248:VAL:O	1:A:248:VAL:HG23	2.10	0.51
2:B:5:PRO:HB3	2:B:30:PHE:HB3	1.91	0.51
3:E:44:ILE:HG22	3:E:45:SER:N	2.26	0.51
2:B:7:ILE:HD13	2:B:82:VAL:HG21	1.92	0.51
2:B:23:LEU:HD23	2:B:39:LEU:HD22	1.92	0.51
3:E:30:ARG:HG2	3:E:30:ARG:HH11	1.76	0.51
4:H:63:LYS:O	4:H:63:LYS:CG	2.59	0.51
2:D:86:THR:HG22	2:D:87:LEU:N	2.16	0.51
3:E:133:ASN:HB3	3:E:136:ASN:HB2	1.93	0.51
3:E:200:HIS:O	3:E:201:SER:C	2.54	0.51
4:H:12:SER:HB2	4:H:118:ILE:CD1	2.34	0.51
4:H:208:CYS:HB3	4:H:223:ILE:HB	1.91	0.51
1:C:63:GLN:O	1:C:66:ILE:HB	2.11	0.51
1:C:208:PHE:O	1:C:240:THR:HB	2.11	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:F:212:HIS:ND1	4:F:212:HIS:C	2.69	0.51
4:H:36:HIS:ND1	4:H:51:TYR:HB3	2.25	0.51
1:A:175:GLY:O	1:A:176:LYS:C	2.54	0.50
2:B:5:PRO:HA	2:B:30:PHE:HB3	1.93	0.50
3:E:30:ARG:HH22	3:E:103:THR:HG21	1.76	0.50
4:F:38:TYR:HB2	4:F:46:LEU:HD12	1.92	0.50
4:F:75:SER:C	4:F:77:SER:H	2.19	0.50
3:G:3:GLN:O	3:G:4:VAL:HG22	2.12	0.50
1:A:199:VAL:O	1:A:249:VAL:N	2.38	0.50
1:C:215:LEU:HD12	1:C:245:ALA:HB2	1.93	0.50
3:G:92:THR:O	3:G:104:ASP:HB2	2.12	0.50
1:A:208:PHE:CE1	1:A:213:ILE:HD12	2.46	0.50
1:A:219:LEU:HD21	1:A:257:TYR:CD2	2.47	0.50
1:C:104:GLU:HB3	1:C:109:LEU:HD12	1.93	0.50
4:F:25:VAL:CG2	4:F:35:ILE:HD11	2.39	0.50
3:G:7:SER:OG	3:G:8:PRO:CD	2.57	0.50
4:H:133:THR:C	5:L:1:NAG:H62	2.36	0.50
1:C:126:LEU:HD21	3:G:105:LYS:HE2	1.94	0.50
4:F:86:LEU:HD22	4:F:117:VAL:CG2	2.41	0.50
4:H:63:LYS:O	4:H:64:GLU:HG3	2.11	0.50
1:A:22:PHE:O	1:A:37:SER:CB	2.48	0.50
4:F:52:VAL:HG22	4:F:71:SER:OG	2.11	0.50
3:G:21:LEU:CD2	3:G:112:THR:HG21	2.42	0.50
1:A:209:TYR:O	1:A:241:PHE:HD1	1.95	0.50
1:A:213:ILE:O	1:A:213:ILE:HG22	2.12	0.50
2:D:31:HIS:O	2:D:33:SER:N	2.45	0.50
2:D:35:ILE:HG23	2:D:35:ILE:O	2.12	0.50
3:E:185:SER:CA	3:E:188:VAL:HG23	2.42	0.50
4:F:162:ILE:HG22	4:F:163:ARG:N	2.27	0.50
1:C:133:CYS:O	1:C:133:CYS:SG	2.70	0.50
3:E:197:GLU:OE1	3:E:197:GLU:HA	2.12	0.50
1:A:7:TYR:OH	1:A:63:GLN:NE2	2.45	0.50
1:A:14:ARG:O	1:A:15:PRO:O	2.29	0.50
1:A:202:ARG:HA	1:A:246:ALA:CB	2.41	0.50
1:A:231:VAL:HG23	1:A:244:TRP:O	2.11	0.50
1:A:258:THR:HG22	1:A:273:ARG:CG	2.42	0.50
1:C:12:VAL:CG1	1:C:13:SER:N	2.75	0.50
1:C:262:TYR:O	1:C:263:HIS:HB2	2.11	0.50
1:C:9:TYR:CZ	1:C:11:ALA:HB2	2.47	0.49
4:F:52:VAL:HG11	4:F:73:ASN:HB2	1.93	0.49
4:H:209:ILE:HG22	4:H:209:ILE:O	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:3:HIS:NE2	1:A:180:LEU:HD21	2.27	0.49
1:A:157:SER:O	1:A:158:GLN:NE2	2.44	0.49
1:C:23:ILE:CG2	1:C:24:ILE:H	2.21	0.49
1:C:113:TYR:CE2	1:C:130:PRO:HG3	2.47	0.49
3:E:50:ALA:O	3:E:68:ARG:NH1	2.45	0.49
3:E:104:ASP:OD2	4:F:36:HIS:CE1	2.65	0.49
3:G:105:LYS:HD3	4:H:48:PHE:CD2	2.47	0.49
1:A:19:GLU:O	1:A:19:GLU:HG3	2.13	0.49
1:C:187:ALA:HB1	1:C:272:LEU:HD11	1.93	0.49
1:C:208:PHE:CE1	1:C:241:PHE:HB2	2.47	0.49
2:D:24:ASN:HD21	2:D:67:TYR:HB2	1.78	0.49
4:F:55:THR:O	4:F:56:PRO:O	2.30	0.49
4:F:136:LEU:HD21	4:F:226:PRO:HD2	1.94	0.49
2:B:51:HIS:HB3	2:B:66:TYR:CD2	2.47	0.49
1:C:71:GLY:O	1:C:74:SER:OG	2.24	0.49
2:D:11:SER:HB3	2:D:95:TRP:HZ2	1.77	0.49
3:E:24:ASN:HD21	6:J:1:NAG:C7	2.19	0.49
3:E:24:ASN:CB	6:J:1:NAG:C1	2.90	0.49
4:H:24:LYS:HA	4:H:78:THR:HB	1.93	0.49
1:A:268:GLU:O	1:A:269:PRO:C	2.55	0.49
2:B:70:PHE:HE2	2:B:95:TRP:CZ3	2.30	0.49
1:C:116:LEU:HD12	1:C:116:LEU:N	2.28	0.49
1:C:164:CYS:C	1:C:166:ASP:N	2.70	0.49
1:C:9:TYR:OH	1:C:11:ALA:HB2	2.12	0.49
1:C:96:GLN:HB3	2:D:56:PHE:CE2	2.48	0.49
2:D:67:TYR:N	2:D:67:TYR:CD1	2.79	0.49
3:E:37:GLN:O	3:E:37:GLN:HG3	2.11	0.49
3:G:138:ALA:HB1	3:G:177:VAL:CG1	2.43	0.49
4:H:33:VAL:HG12	4:H:34:THR:N	2.27	0.49
4:H:50:LEU:HD11	4:H:61:LEU:HB2	1.94	0.49
2:B:31:HIS:O	2:B:32:PRO:C	2.55	0.49
3:E:184:ASP:O	3:E:185:SER:C	2.56	0.49
4:F:37:TRP:CG	4:F:81:LEU:HD13	2.48	0.49
4:H:8:GLU:HA	4:H:8:GLU:OE1	2.13	0.49
4:H:27:ILE:HG22	4:H:28:GLU:O	2.13	0.49
1:A:63:GLN:O	1:A:67:VAL:HG23	2.13	0.49
3:E:185:SER:HA	3:E:188:VAL:CG2	2.43	0.49
4:F:11:ILE:O	4:F:116:ILE:N	2.40	0.49
4:F:159:PRO:O	4:F:160:ASP:OD1	2.31	0.49
4:F:185:THR:CG2	6:K:1:NAG:C8	2.83	0.49
4:F:186:ASN:ND2	6:K:1:NAG:H5	2.28	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:9:TYR:CB	1:A:70:GLN:NE2	2.75	0.48
1:A:16:GLY:C	1:A:18:GLY:N	2.71	0.48
4:F:53:LEU:HD13	4:F:59:ILE:CD1	2.43	0.48
4:H:12:SER:CB	4:H:118:ILE:HD11	2.35	0.48
1:A:2:SER:O	1:A:3:HIS:ND1	2.46	0.48
1:A:35:ARG:CZ	1:A:40:GLU:OE2	2.62	0.48
1:A:73:LEU:CD1	3:E:98:GLY:O	2.60	0.48
1:A:128:GLU:O	1:A:130:PRO:N	2.46	0.48
1:A:202:ARG:HB2	1:A:202:ARG:NH1	2.27	0.48
1:A:262:TYR:CD1	1:A:262:TYR:N	2.81	0.48
1:C:128:GLU:HG3	1:C:129:ASN:N	2.27	0.48
4:F:86:LEU:HD22	4:F:117:VAL:HG21	1.96	0.48
4:F:139:VAL:O	4:F:140:ALA:C	2.55	0.48
4:H:130:PRO:O	4:H:132:PRO:HD3	2.13	0.48
3:E:17:THR:HG22	3:E:18:ASP:H	1.76	0.48
3:E:32:VAL:HG22	3:E:90:ALA:HB2	1.95	0.48
4:H:27:ILE:HD13	4:H:98:TYR:OH	2.13	0.48
1:C:191:ARG:N	1:C:274:TRP:HZ3	2.10	0.48
3:E:7:SER:OG	3:E:8:PRO:N	2.44	0.48
1:A:71:GLY:O	1:A:74:SER:OG	2.30	0.48
1:A:208:PHE:CE1	1:A:213:ILE:HB	2.49	0.48
1:A:209:TYR:CG	1:A:210:PRO:N	2.80	0.48
3:E:169:SER:HB3	3:E:170:PRO:CD	2.44	0.48
3:G:98:GLY:O	3:G:99:TYR:CD1	2.66	0.48
3:G:126:PRO:HG3	3:G:144:PHE:CB	2.43	0.48
4:H:207:ARG:HA	4:H:223:ILE:O	2.13	0.48
1:A:14:ARG:HB2	1:A:15:PRO:HD2	1.95	0.48
1:A:218:GLN:OE1	1:A:260:HIS:CG	2.66	0.48
1:C:98:LEU:C	1:C:98:LEU:CD1	2.77	0.48
1:C:204:TRP:HZ2	2:D:98:ASP:O	1.97	0.48
4:F:157:PHE:CG	4:F:162:ILE:HD11	2.49	0.48
4:H:107:LYS:CG	4:H:109:PHE:HE1	2.25	0.48
1:A:69:ILE:O	1:A:72:GLN:HB2	2.13	0.48
1:A:238:ASP:OD1	1:A:238:ASP:C	2.57	0.48
2:B:73:THR:OG1	2:B:76:ASP:OD2	2.25	0.48
1:C:213:ILE:HG12	1:C:214:THR:H	1.78	0.48
3:E:149:VAL:CG1	3:E:150:THR:N	2.75	0.48
4:F:128:ILE:HG21	4:F:158:PHE:CB	2.26	0.48
4:H:134:ILE:CG2	4:H:135:PHE:N	2.76	0.48
1:A:61:GLU:C	1:A:64:THR:HG1	2.18	0.48
1:A:209:TYR:HB3	1:A:210:PRO:CD	2.40	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:7:ILE:HD13	2:B:82:VAL:CG2	2.44	0.48
1:C:76:ARG:C	1:C:79:MET:HB2	2.38	0.48
2:D:79:ALA:CA	2:D:93:VAL:O	2.62	0.48
4:H:75:SER:C	4:H:77:SER:N	2.72	0.48
1:A:69:ILE:O	1:A:70:GLN:C	2.55	0.48
1:A:124:PRO:HB3	3:E:28:THR:CG2	2.39	0.48
4:F:37:TRP:CZ2	4:F:96:CYS:SG	3.07	0.48
1:A:258:THR:HG22	1:A:273:ARG:HG2	1.95	0.48
3:G:15:GLU:HG2	3:G:119:LYS:HE3	1.95	0.48
4:H:73:ASN:OD1	4:H:73:ASN:O	2.32	0.48
4:H:99:GLY:HA3	4:H:105:PHE:HD2	1.79	0.48
2:B:24:ASN:HD21	2:B:67:TYR:HB2	1.78	0.47
4:F:146:LYS:HG2	4:F:199:GLU:OE1	2.14	0.47
4:F:199:GLU:O	4:F:201:ALA:N	2.47	0.47
4:H:209:ILE:HA	4:H:221:GLN:O	2.13	0.47
1:A:9:TYR:HB2	1:A:70:GLN:NE2	2.29	0.47
1:C:5:LEU:HD21	1:C:167:LEU:HD22	1.96	0.47
1:C:217:TRP:HE1	1:C:246:ALA:HA	1.79	0.47
3:E:55:GLU:O	3:E:55:GLU:HG3	2.14	0.47
4:F:145:HIS:O	4:F:146:LYS:HB3	2.14	0.47
4:F:147:THR:CG2	4:F:148:GLY:N	2.71	0.47
4:F:160:ASP:HB2	4:F:189:TYR:CG	2.48	0.47
3:G:144:PHE:CE2	3:G:175:SER:HA	2.49	0.47
1:A:34:LEU:HD12	1:A:63:GLN:OE1	2.13	0.47
1:C:92:SER:O	1:C:93:HIS:CD2	2.66	0.47
2:D:23:LEU:HD23	2:D:39:LEU:HD22	1.96	0.47
3:E:61:SER:CB	3:E:72:THR:O	2.59	0.47
4:F:23:CYS:SG	4:F:96:CYS:SG	3.12	0.47
4:F:51:TYR:CE2	4:F:59:ILE:HD13	2.49	0.47
4:F:147:THR:CG2	4:F:148:GLY:H	2.27	0.47
4:F:157:PHE:CD1	4:F:157:PHE:C	2.93	0.47
4:F:212:HIS:O	4:F:213:GLU:C	2.57	0.47
1:A:97:TRP:CE3	1:A:116:LEU:HD21	2.49	0.47
1:A:263:HIS:CG	1:A:264:GLU:N	2.82	0.47
2:B:4:THR:OG1	2:B:5:PRO:HD2	2.14	0.47
2:D:18:GLY:N	2:D:72:PRO:O	2.47	0.47
3:E:135:THR:O	3:E:181:GLN:HA	2.14	0.47
4:F:167:LYS:HB2	4:F:209:ILE:HD12	1.96	0.47
1:A:73:LEU:C	1:A:75:GLU:N	2.72	0.47
1:A:118:TYR:HB2	1:A:123:LEU:HD21	1.95	0.47
3:E:14:HIS:C	3:E:15:GLU:O	2.57	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:H:151:LEU:N	4:H:151:LEU:CD2	2.78	0.47
1:A:36:PHE:HD1	1:A:43:PRO:HA	1.78	0.47
1:A:199:VAL:CG1	1:A:200:THR:N	2.77	0.47
2:B:40:LEU:CA	2:B:44:GLU:O	2.56	0.47
2:B:59:ASP:C	2:B:61:SER:H	2.23	0.47
2:B:73:THR:O	2:B:74:GLU:C	2.58	0.47
2:D:37:VAL:HB	2:D:66:TYR:CE2	2.49	0.47
2:D:98:ASP:C	2:D:99:MET:SD	2.97	0.47
3:E:133:ASN:O	3:E:136:ASN:N	2.46	0.47
3:E:147:LYS:HD2	3:E:174:TYR:CE2	2.49	0.47
4:F:115:LEU:O	4:F:115:LEU:HG	2.15	0.47
1:A:126:LEU:N	3:E:92:THR:HG21	2.30	0.47
1:A:217:TRP:HE1	1:A:246:ALA:HA	1.79	0.47
1:A:274:TRP:CG	1:A:275:GLY:N	2.83	0.47
2:B:31:HIS:HB3	2:B:32:PRO:CD	2.39	0.47
1:C:9:TYR:OH	1:C:74:SER:CB	2.58	0.47
2:D:47:GLU:O	2:D:48:LYS:HG2	2.15	0.47
3:E:92:THR:CG2	3:E:93:TRP:N	2.78	0.47
3:E:148:GLU:N	3:E:148:GLU:OE1	2.44	0.47
4:F:21:ILE:CD1	4:F:94:TYR:HB2	2.45	0.47
4:F:55:THR:CB	4:F:56:PRO:CD	2.92	0.47
4:H:164:VAL:HG12	4:H:165:TYR:N	2.29	0.47
1:A:79:MET:C	1:A:81:LEU:N	2.70	0.47
1:A:209:TYR:CB	1:A:210:PRO:HD3	2.42	0.47
1:A:234:ARG:NH1	2:B:10:TYR:CD1	2.82	0.47
1:C:129:ASN:HB3	1:C:130:PRO:HD2	1.87	0.47
1:C:205:ALA:O	1:C:206:LEU:HD23	2.14	0.47
2:D:23:LEU:C	2:D:24:ASN:HD22	2.23	0.47
3:E:6:GLN:NE2	3:E:86:TYR:O	2.48	0.47
3:E:127:SER:OG	3:E:142:LYS:HB3	2.15	0.47
4:F:21:ILE:HG22	4:F:22:SER:N	2.30	0.47
4:F:72:LYS:O	4:F:79:SER:HA	2.13	0.47
1:A:153:VAL:HG12	1:A:154:PRO:HD3	1.97	0.47
1:C:126:LEU:HD21	3:G:105:LYS:CE	2.44	0.47
2:D:39:LEU:HD12	2:D:49:VAL:HG21	1.97	0.47
3:G:47:PHE:CZ	3:G:60:LYS:HA	2.50	0.47
3:G:152:SER:N	3:G:191:SER:O	2.48	0.47
4:H:105:PHE:C	4:H:107:LYS:N	2.64	0.47
1:A:97:TRP:CG	1:A:116:LEU:HG	2.50	0.47
2:B:24:ASN:ND2	2:B:67:TYR:CB	2.78	0.47
2:B:84:HIS:N	2:B:87:LEU:HD21	2.23	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:77:ASN:O	1:C:78:LEU:C	2.58	0.47
1:C:196:GLU:OE2	1:C:196:GLU:N	2.48	0.47
2:D:24:ASN:ND2	2:D:67:TYR:HB2	2.30	0.47
4:H:31:ARG:C	4:H:33:VAL:H	2.22	0.47
4:H:69:GLU:HA	4:H:82:THR:O	2.15	0.47
1:A:83:HIS:C	1:A:86:ASN:H	2.23	0.46
1:C:25:VAL:HG21	2:D:54:LEU:O	2.14	0.46
1:C:45:MET:SD	1:C:54:GLN:NE2	2.86	0.46
1:C:129:ASN:O	1:C:131:SER:N	2.49	0.46
4:F:49:LEU:HD13	4:F:69:GLU:OE2	2.15	0.46
4:F:72:LYS:O	4:F:79:SER:HB2	2.15	0.46
3:G:6:GLN:O	3:G:7:SER:O	2.32	0.46
1:C:3:HIS:ND1	1:C:3:HIS:N	2.63	0.46
4:F:198:PRO:O	4:F:202:MET:HG3	2.14	0.46
3:G:126:PRO:HG3	3:G:144:PHE:HB2	1.96	0.46
2:B:17:ASN:ND2	2:B:74:GLU:HB2	2.30	0.46
1:C:129:ASN:O	1:C:130:PRO:C	2.58	0.46
1:C:210:PRO:O	1:C:211:ALA:C	2.59	0.46
2:D:94:LYS:H	2:D:94:LYS:HG3	1.41	0.46
3:E:32:VAL:CG2	3:E:90:ALA:HB2	2.45	0.46
4:H:16:ASP:N	4:H:86:LEU:O	2.45	0.46
3:E:24:ASN:ND2	6:J:1:NAG:HN2	2.02	0.46
1:A:165:SER:OG	1:A:166:ASP:N	2.47	0.46
1:A:181:ARG:HG2	1:A:182:SER:N	2.30	0.46
1:C:51:TRP:HZ3	1:C:179:LEU:HG	1.81	0.46
3:E:145:TYR:CD2	3:E:146:PRO:CD	2.99	0.46
4:F:183:LEU:HB3	4:F:190:MET:HE2	1.97	0.46
4:H:105:PHE:C	4:H:107:LYS:H	2.13	0.46
2:B:39:LEU:O	2:B:45:ARG:HA	2.15	0.46
2:B:42:ASN:ND2	2:B:76:ASP:OD1	2.49	0.46
2:B:69:GLU:N	2:B:69:GLU:OE1	2.47	0.46
1:C:77:ASN:C	1:C:79:MET:N	2.73	0.46
2:D:31:HIS:HD1	2:D:32:PRO:N	2.13	0.46
2:D:67:TYR:N	2:D:67:TYR:HD1	2.13	0.46
3:E:139:CYS:HB2	3:E:153:LEU:HD21	1.97	0.46
4:F:48:PHE:O	4:F:49:LEU:HD23	2.15	0.46
2:B:33:SER:O	2:B:35:ILE:HG22	2.15	0.46
2:B:67:TYR:CD1	2:B:67:TYR:N	2.84	0.46
1:C:219:LEU:O	1:C:220:ASN:HB3	2.15	0.46
2:D:33:SER:O	2:D:34:ASP:C	2.58	0.46
2:D:41:LYS:HE2	2:D:42:ASN:ND2	2.30	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:E:7:SER:CB	3:E:22:ARG:HB2	2.44	0.46
4:F:185:THR:O	4:F:187:ASP:N	2.49	0.46
3:G:13:LEU:O	3:G:117:GLU:N	2.47	0.46
3:G:47:PHE:HZ	3:G:59:LEU:O	1.98	0.46
3:G:53:THR:CG2	3:G:54:LYS:N	2.79	0.46
1:C:176:LYS:HD3	1:C:180:LEU:HD11	1.97	0.46
3:E:21:LEU:HD22	3:E:112:THR:HG21	1.98	0.46
1:A:13:SER:O	1:A:92:SER:OG	2.27	0.46
1:A:266:LEU:HD22	1:A:270:LEU:CD1	2.34	0.46
1:C:181:ARG:HG2	1:C:182:SER:H	1.81	0.46
1:C:218:GLN:HG2	1:C:223:GLU:HA	1.97	0.46
3:E:132:LYS:CE	3:E:132:LYS:N	2.74	0.46
4:F:162:ILE:C	4:F:163:ARG:HD3	2.40	0.46
4:H:92:ALA:O	4:H:94:TYR:CD1	2.69	0.46
4:H:189:TYR:N	4:H:189:TYR:CD1	2.83	0.46
1:A:155:HIS:NE2	1:A:157:SER:CA	2.65	0.46
1:C:4:SER:O	1:C:28:VAL:HA	2.15	0.46
1:C:97:TRP:CG	1:C:116:LEU:HG	2.51	0.46
2:D:13:HIS:HB3	2:D:14:PRO:CD	2.46	0.46
4:F:151:LEU:HD12	4:F:151:LEU:HA	1.79	0.46
3:G:59:LEU:C	3:G:60:LYS:HG2	2.40	0.46
3:G:103:THR:HG23	3:G:103:THR:O	2.16	0.46
1:C:25:VAL:HG13	2:D:53:ASP:HB3	1.98	0.45
3:E:43:LEU:HD11	4:F:46:LEU:HD21	1.97	0.45
4:H:9:ILE:O	4:H:113:THR:HG23	2.16	0.45
1:A:84:PHE:HE1	3:E:30:ARG:HH21	1.65	0.45
1:A:180:LEU:HD23	1:A:180:LEU:HA	1.62	0.45
2:D:37:VAL:HG13	2:D:82:VAL:HG22	1.97	0.45
4:F:60:PHE:HB3	4:F:70:ALA:HB2	1.98	0.45
1:A:9:TYR:C	1:A:9:TYR:CD2	2.94	0.45
1:A:191:ARG:N	1:A:274:TRP:CZ3	2.84	0.45
1:A:196:GLU:H	1:A:196:GLU:HG3	1.53	0.45
2:B:59:ASP:O	2:B:61:SER:N	2.49	0.45
2:B:80:CYS:O	2:B:92:ILE:HA	2.17	0.45
1:C:55:GLU:HA	1:C:55:GLU:OE1	2.16	0.45
1:C:121:GLU:O	1:C:122:ASP:C	2.60	0.45
1:C:205:ALA:C	1:C:206:LEU:HD23	2.42	0.45
3:E:132:LYS:N	3:E:132:LYS:CD	2.79	0.45
3:E:185:SER:HA	3:E:188:VAL:HG23	1.98	0.45
1:A:73:LEU:HD23	1:A:73:LEU:N	2.31	0.45
1:A:103:VAL:HG23	1:A:168:LEU:HD22	1.97	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:97:TRP:CE2	1:C:116:LEU:HD11	2.52	0.45
1:C:234:ARG:HA	1:C:235:PRO:HD3	1.70	0.45
3:E:30:ARG:HG2	3:E:30:ARG:NH1	2.30	0.45
3:E:153:LEU:CD2	3:E:190:CYS:SG	3.05	0.45
1:C:51:TRP:CD1	1:C:51:TRP:H	2.34	0.45
3:G:147:LYS:HG3	3:G:174:TYR:CE2	2.44	0.45
1:C:9:TYR:CE2	1:C:11:ALA:HB2	2.52	0.45
4:F:54:ALA:C	4:F:55:THR:OG1	2.60	0.45
4:F:134:ILE:HG22	4:F:135:PHE:N	2.32	0.45
4:F:156:ALA:C	4:F:188:THR:HG21	2.41	0.45
4:F:215:ASN:OD1	4:F:215:ASN:O	2.34	0.45
3:G:67:GLU:OE1	3:G:69:ARG:CD	2.65	0.45
4:H:5:GLU:HG3	4:H:7:PRO:HD3	1.98	0.45
4:H:51:TYR:H	4:H:51:TYR:HD2	1.64	0.45
4:H:72:LYS:O	4:H:79:SER:HA	2.16	0.45
1:A:6:ARG:HG3	1:A:98:LEU:HD11	1.98	0.45
1:A:97:TRP:CZ3	1:A:116:LEU:HD21	2.52	0.45
2:B:23:LEU:HB3	2:B:68:THR:HG22	1.99	0.45
1:C:164:CYS:O	1:C:165:SER:C	2.60	0.45
2:D:5:PRO:HB3	2:D:30:PHE:CB	2.43	0.45
4:F:146:LYS:HG3	4:F:199:GLU:CG	2.46	0.45
1:A:202:ARG:CD	1:A:204:TRP:CZ2	2.97	0.45
1:C:25:VAL:HG12	1:C:27:TYR:CE1	2.52	0.45
1:C:55:GLU:O	1:C:57:ALA:N	2.48	0.45
4:F:199:GLU:C	4:F:201:ALA:H	2.25	0.45
3:G:26:THR:O	3:G:26:THR:HG22	2.16	0.45
3:G:27:THR:O	3:G:29:MET:SD	2.75	0.45
4:H:6:GLN:HG2	4:H:23:CYS:HB3	1.98	0.45
1:A:187:ALA:HB3	1:A:270:LEU:CD2	2.47	0.45
1:A:213:ILE:O	1:A:213:ILE:CG2	2.59	0.45
2:D:37:VAL:HG22	2:D:82:VAL:HG13	1.99	0.45
1:A:60:TRP:O	1:A:63:GLN:N	2.46	0.45
1:C:95:LEU:HD12	1:C:96:GLN:N	2.32	0.45
3:E:40:ARG:HB3	4:F:163:ARG:NH2	2.32	0.45
3:E:96:SER:OG	3:E:100:GLU:HB2	2.17	0.45
3:G:3:GLN:C	3:G:4:VAL:CG2	2.90	0.45
3:G:113:GLN:HG3	3:G:113:GLN:O	2.17	0.45
1:A:64:THR:HG1	1:A:65:ARG:H	1.64	0.44
1:C:235:PRO:HG2	2:D:26:TYR:CE1	2.52	0.44
1:C:252:GLY:C	1:C:254:GLU:H	2.26	0.44
3:E:35:PHE:CE1	3:E:108:PHE:HE2	2.34	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:H:21:ILE:HD12	4:H:81:LEU:HD23	1.99	0.44
4:H:33:VAL:HG13	4:H:34:THR:H	1.81	0.44
1:A:105:PRO:C	1:A:107:ARG:H	2.23	0.44
2:B:83:ASN:CG	2:B:90:PRO:HG3	2.43	0.44
1:C:126:LEU:CA	3:G:92:THR:HG21	2.47	0.44
2:D:41:LYS:C	2:D:43:GLY:N	2.73	0.44
3:E:4:VAL:HG13	3:E:25:PHE:CD2	2.52	0.44
3:E:30:ARG:HH22	3:E:103:THR:CG2	2.30	0.44
1:A:103:VAL:CG1	1:A:104:GLU:N	2.80	0.44
1:C:117:ALA:HB1	1:C:121:GLU:H	1.83	0.44
1:C:232:GLU:HB3	2:D:8:GLN:NE2	2.32	0.44
2:D:13:HIS:HB3	2:D:14:PRO:HD2	1.99	0.44
4:F:7:PRO:HG2	4:F:22:SER:OG	2.17	0.44
4:F:23:CYS:SG	4:F:96:CYS:HB3	2.58	0.44
1:A:73:LEU:CB	1:A:77:ASN:ND2	2.81	0.44
1:C:230:LEU:CD2	1:C:245:ALA:HB2	2.47	0.44
1:C:254:GLU:C	1:C:256:SER:N	2.71	0.44
2:D:39:LEU:O	2:D:40:LEU:HD23	2.16	0.44
2:D:84:HIS:CE1	2:D:85:VAL:O	2.71	0.44
4:F:63:LYS:O	4:F:63:LYS:HG3	2.18	0.44
4:H:93:ILE:HG22	4:H:95:TYR:CE1	2.53	0.44
1:A:39:LYS:O	1:A:40:GLU:CG	2.64	0.44
1:A:191:ARG:HB3	1:A:274:TRP:CH2	2.53	0.44
1:A:199:VAL:N	1:A:249:VAL:O	2.51	0.44
2:B:49:VAL:HG22	2:B:68:THR:HB	2.00	0.44
1:C:29:ASP:CG	1:C:29:ASP:O	2.61	0.44
1:C:100:GLY:C	1:C:101:CYS:SG	2.98	0.44
1:C:191:ARG:CB	1:C:274:TRP:HZ3	2.28	0.44
3:E:89:ALA:HB1	3:E:107:VAL:O	2.17	0.44
3:E:165:ALA:O	3:E:176:ALA:HA	2.17	0.44
4:F:150:TYR:CD2	4:F:225:PHE:HD2	2.35	0.44
4:H:131:LYS:N	4:H:155:GLU:O	2.51	0.44
1:A:22:PHE:CZ	1:A:71:GLY:HA3	2.52	0.44
1:A:76:ARG:O	1:A:77:ASN:C	2.60	0.44
2:B:56:PHE:HB2	2:B:61:SER:O	2.18	0.44
1:C:188:HIS:CE1	1:C:204:TRP:HB2	2.52	0.44
4:F:159:PRO:O	4:F:160:ASP:CB	2.66	0.44
3:G:89:ALA:HB1	3:G:106:LEU:HG	1.99	0.44
3:G:198:THR:CG2	3:G:199:VAL:N	2.81	0.44
4:H:198:PRO:O	4:H:200:ARG:N	2.51	0.44
1:A:4:SER:HB2	1:A:102:ASP:OD1	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:76:ARG:O	1:A:79:MET:N	2.48	0.44
1:A:84:PHE:CE1	3:E:30:ARG:NE	2.77	0.44
1:A:234:ARG:HD2	1:A:242:GLN:O	2.17	0.44
2:B:84:HIS:O	2:B:87:LEU:CD2	2.66	0.44
1:C:87:LYS:HD3	1:C:93:HIS:CE1	2.52	0.44
3:E:145:TYR:CD2	3:E:146:PRO:HG3	2.53	0.44
4:F:132:PRO:CG	4:F:221:GLN:HE22	2.30	0.44
4:F:136:LEU:HB3	4:F:137:PRO:CD	2.48	0.44
4:H:193:SER:O	4:H:194:TRP:HB3	2.17	0.44
3:E:131:MET:HB2	3:E:138:ALA:HB3	2.00	0.44
3:E:179:LEU:HD11	4:F:153:LEU:HD22	1.99	0.44
4:H:33:VAL:O	4:H:53:LEU:C	2.61	0.44
4:H:94:TYR:O	4:H:112:GLY:HA2	2.17	0.44
1:A:12:VAL:CG1	1:A:13:SER:N	2.81	0.44
1:A:34:LEU:C	1:A:34:LEU:HD23	2.43	0.44
1:A:60:TRP:O	1:A:61:GLU:C	2.61	0.44
1:A:81:LEU:HD13	1:A:118:TYR:HD1	1.82	0.44
1:A:209:TYR:CB	1:A:210:PRO:CD	2.96	0.44
1:C:35:ARG:HG2	1:C:36:PHE:N	2.33	0.44
1:C:81:LEU:HG	1:C:118:TYR:CD1	2.53	0.44
1:C:208:PHE:CZ	1:C:241:PHE:O	2.70	0.44
2:D:29:GLY:HA2	2:D:61:SER:OG	2.18	0.44
3:E:35:PHE:CE1	3:E:108:PHE:CE2	3.06	0.44
3:G:24:ASN:N	3:G:24:ASN:OD1	2.50	0.44
4:H:60:PHE:HD2	4:H:60:PHE:H	1.55	0.44
1:C:97:TRP:HD1	1:C:98:LEU:H	1.61	0.43
2:D:56:PHE:HB2	2:D:61:SER:O	2.18	0.43
3:E:7:SER:HB3	3:E:22:ARG:C	2.44	0.43
1:A:10:THR:HG21	2:B:62:PHE:HE2	1.82	0.43
1:A:208:PHE:CD2	1:A:213:ILE:HD12	2.54	0.43
2:B:13:HIS:O	2:B:15:ALA:N	2.51	0.43
1:C:167:LEU:HD23	1:C:167:LEU:C	2.43	0.43
2:D:41:LYS:C	2:D:43:GLY:H	2.24	0.43
4:F:20:GLN:HA	4:F:81:LEU:O	2.19	0.43
4:F:30:PHE:O	4:F:31:ARG:C	2.61	0.43
3:G:77:ASP:O	3:G:78:ALA:C	2.61	0.43
4:H:148:GLY:O	4:H:197:VAL:HG23	2.18	0.43
7:M:2:NAG:H82	7:M:2:NAG:C3	2.47	0.43
1:A:231:VAL:CG2	1:A:244:TRP:O	2.67	0.43
1:A:274:TRP:CG	1:A:275:GLY:H	2.37	0.43
2:B:37:VAL:HA	2:B:81:ARG:O	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:87:LYS:HB3	1:C:93:HIS:CE1	2.41	0.43
1:C:166:ASP:O	1:C:170:LYS:HG2	2.18	0.43
3:E:6:GLN:OE1	3:E:88:CYS:SG	2.77	0.43
3:G:8:PRO:O	3:G:112:THR:HG23	2.18	0.43
3:G:133:ASN:O	3:G:133:ASN:CG	2.61	0.43
4:H:25:VAL:HG12	4:H:26:PHE:N	2.33	0.43
4:H:157:PHE:CG	4:H:159:PRO:HD2	2.54	0.43
1:A:28:VAL:O	1:A:29:ASP:HB2	2.17	0.43
1:A:38:SER:O	1:A:39:LYS:HB3	2.18	0.43
1:C:34:LEU:CD2	1:C:35:ARG:N	2.80	0.43
3:E:10:ALA:HB3	3:G:12:SER:HB3	2.00	0.43
4:F:55:THR:HB	4:F:56:PRO:CD	2.48	0.43
4:H:98:TYR:CD1	4:H:98:TYR:C	2.97	0.43
1:A:97:TRP:CE2	1:A:116:LEU:HD11	2.54	0.43
2:B:7:ILE:CD1	2:B:82:VAL:HG21	2.49	0.43
1:C:61:GLU:C	1:C:63:GLN:N	2.77	0.43
2:D:20:SER:C	2:D:21:ASN:CG	2.84	0.43
2:D:23:LEU:CD2	2:D:39:LEU:HD22	2.48	0.43
3:E:37:GLN:HG2	3:E:87:PHE:HE1	1.84	0.43
4:F:82:THR:HG22	4:F:83:ILE:N	2.33	0.43
4:F:105:PHE:HD2	4:F:107:LYS:HG2	1.82	0.43
4:F:158:PHE:C	4:F:160:ASP:N	2.77	0.43
1:A:152:THR:HG22	1:A:153:VAL:N	2.33	0.43
2:B:54:LEU:HD11	2:B:62:PHE:HD2	1.84	0.43
1:C:231:VAL:HG11	1:C:244:TRP:CZ2	2.54	0.43
1:C:234:ARG:HB2	2:D:10:TYR:OH	2.19	0.43
3:E:141:VAL:HG21	3:E:151:ILE:HD11	1.99	0.43
4:F:190:MET:CG	4:F:191:LYS:N	2.82	0.43
3:G:7:SER:CB	3:G:22:ARG:H	2.31	0.43
3:G:143:ASP:CB	3:G:173:LYS:HD2	2.47	0.43
1:A:43:PRO:HG2	1:A:64:THR:HA	2.01	0.43
2:B:42:ASN:ND2	2:B:76:ASP:HA	2.29	0.43
1:C:5:LEU:HD12	1:C:27:TYR:O	2.19	0.43
1:C:23:ILE:O	1:C:24:ILE:HG12	2.19	0.43
1:C:176:LYS:O	1:C:177:GLU:C	2.61	0.43
2:D:56:PHE:CB	2:D:61:SER:O	2.66	0.43
3:E:77:ASP:O	3:E:78:ALA:C	2.62	0.43
3:E:123:PRO:HA	3:E:146:PRO:HD3	1.99	0.43
4:F:106:HIS:O	4:F:107:LYS:C	2.61	0.43
4:F:124:LEU:O	4:F:124:LEU:HG	2.17	0.43
3:G:146:PRO:C	3:G:148:GLU:N	2.71	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:H:3:LYS:HE3	4:H:27:ILE:HD11	2.01	0.43
4:H:90:ASP:O	4:H:94:TYR:OH	2.37	0.43
4:H:147:THR:HG22	4:H:148:GLY:N	2.33	0.43
1:A:8:PHE:HB2	1:A:25:VAL:HG23	2.01	0.43
2:B:54:LEU:HA	2:B:64:LEU:CD2	2.49	0.43
1:C:97:TRP:CD2	1:C:116:LEU:HG	2.53	0.43
1:C:202:ARG:HA	1:C:246:ALA:HB2	2.01	0.43
3:E:31:SER:HA	3:E:50:ALA:HA	2.01	0.43
3:E:46:LEU:O	3:E:47:PHE:CD2	2.71	0.43
3:E:164:PRO:HG3	3:E:178:LYS:CE	2.49	0.43
4:F:75:SER:O	4:F:77:SER:N	2.46	0.43
3:G:125:LYS:HA	3:G:126:PRO:HD3	1.53	0.43
4:H:51:TYR:O	4:H:51:TYR:HD2	2.02	0.43
4:H:150:TYR:C	4:H:151:LEU:HD23	2.43	0.43
1:A:173:GLU:C	1:A:175:GLY:H	2.26	0.43
1:C:8:PHE:CD2	2:D:56:PHE:CE1	3.07	0.43
2:D:41:LYS:HE2	2:D:42:ASN:HD21	1.83	0.43
3:E:4:VAL:HG13	3:E:25:PHE:HD2	1.84	0.43
3:E:122:PRO:O	3:E:123:PRO:O	2.36	0.43
4:F:228:ILE:CG2	4:F:229:LYS:H	2.29	0.43
1:A:157:SER:C	1:A:158:GLN:HG2	2.43	0.43
1:A:185:PRO:CB	1:A:208:PHE:HB3	2.49	0.43
1:C:219:LEU:HB2	1:C:224:LEU:HD11	1.99	0.43
1:C:264:GLU:C	1:C:266:LEU:N	2.77	0.43
4:F:30:PHE:O	4:F:33:VAL:HG23	2.19	0.43
4:F:79:SER:C	4:F:80:ILE:HG13	2.44	0.43
4:F:132:PRO:HG3	4:F:223:ILE:HD11	2.00	0.43
3:G:7:SER:HB3	3:G:22:ARG:H	1.83	0.43
3:G:106:LEU:CB	4:H:38:TYR:OH	2.67	0.43
3:G:108:PHE:N	3:G:108:PHE:CD1	2.85	0.43
3:G:134:GLY:O	3:G:136:ASN:ND2	2.51	0.43
1:A:81:LEU:HD22	1:A:118:TYR:CD1	2.54	0.42
2:B:17:ASN:OD1	2:B:17:ASN:N	2.49	0.42
1:C:262:TYR:N	1:C:262:TYR:CD1	2.87	0.42
2:D:24:ASN:ND2	2:D:67:TYR:HB3	2.33	0.42
4:F:90:ASP:O	4:F:115:LEU:HD23	2.19	0.42
1:A:73:LEU:HB2	1:A:77:ASN:ND2	2.34	0.42
1:A:107:ARG:C	1:A:109:LEU:N	2.77	0.42
1:C:14:ARG:O	1:C:16:GLY:N	2.52	0.42
1:C:51:TRP:CZ3	1:C:179:LEU:HG	2.54	0.42
2:D:95:TRP:CD1	2:D:96:ASP:N	2.87	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:F:63:LYS:O	4:F:64:GLU:HB2	2.18	0.42
3:G:82:ASP:O	3:G:83:SER:C	2.62	0.42
3:G:148:GLU:O	3:G:149:VAL:CG2	2.61	0.42
1:A:252:GLY:O	1:A:253:LYS:C	2.61	0.42
1:C:73:LEU:CD1	3:G:99:TYR:HB2	2.46	0.42
1:C:102:ASP:O	1:C:103:VAL:HG23	2.19	0.42
4:F:160:ASP:HB2	4:F:189:TYR:CD2	2.54	0.42
4:F:169:LYS:HB2	4:F:207:ARG:HG3	2.00	0.42
3:G:18:ASP:HA	3:G:75:ILE:O	2.19	0.42
3:G:26:THR:O	3:G:26:THR:CG2	2.67	0.42
3:G:46:LEU:HA	3:G:46:LEU:HD23	1.80	0.42
3:G:47:PHE:CZ	3:G:59:LEU:O	2.72	0.42
1:A:157:SER:C	1:A:158:GLN:CG	2.92	0.42
2:B:57:SER:O	2:B:58:LYS:C	2.62	0.42
2:B:87:LEU:HD23	2:B:87:LEU:O	2.20	0.42
2:D:80:CYS:O	2:D:93:VAL:N	2.49	0.42
3:E:80:LEU:CD2	3:E:116:VAL:O	2.66	0.42
4:F:28:GLU:O	4:F:30:PHE:N	2.52	0.42
4:F:144:LEU:HA	4:F:144:LEU:HD22	1.64	0.42
3:G:166:ILE:CG2	3:G:175:SER:O	2.56	0.42
3:G:181:GLN:H	3:G:181:GLN:HG3	1.66	0.42
1:A:83:HIS:O	1:A:85:TYR:N	2.53	0.42
1:A:97:TRP:CD1	1:A:98:LEU:N	2.88	0.42
1:A:190:THR:OG1	1:A:202:ARG:HB3	2.20	0.42
1:A:204:TRP:CH2	1:A:244:TRP:CD1	3.08	0.42
1:C:236:ALA:CB	1:C:240:THR:O	2.67	0.42
3:E:12:SER:HA	3:E:115:THR:O	2.20	0.42
3:E:135:THR:O	3:E:135:THR:HG22	2.19	0.42
3:E:145:TYR:CG	3:E:146:PRO:N	2.86	0.42
1:A:61:GLU:O	1:A:65:ARG:N	2.38	0.42
2:B:5:PRO:CB	2:B:30:PHE:HB3	2.49	0.42
1:C:22:PHE:CE2	1:C:38:SER:HB3	2.54	0.42
1:C:191:ARG:HB3	1:C:274:TRP:CZ3	2.49	0.42
1:C:194:ARG:C	1:C:196:GLU:N	2.77	0.42
3:E:6:GLN:O	3:E:7:SER:O	2.37	0.42
3:E:61:SER:CA	3:E:72:THR:O	2.68	0.42
3:E:119:LYS:O	3:E:120:SER:C	2.61	0.42
4:F:19:ALA:O	4:F:83:ILE:N	2.53	0.42
3:G:99:TYR:CD1	3:G:99:TYR:N	2.88	0.42
4:H:221:GLN:NE2	4:H:223:ILE:HD11	2.35	0.42
1:A:153:VAL:CG1	1:A:154:PRO:HD3	2.49	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:49:VAL:HG13	2:B:67:TYR:O	2.19	0.42
2:B:75:LYS:HA	2:B:75:LYS:HD3	1.61	0.42
1:C:88:SER:C	1:C:90:ASP:N	2.77	0.42
3:E:117:GLU:OE1	3:E:117:GLU:HA	2.20	0.42
3:E:122:PRO:O	3:E:123:PRO:C	2.63	0.42
1:A:103:VAL:O	1:A:104:GLU:HB2	2.18	0.42
1:A:125:THR:C	3:E:92:THR:HG21	2.44	0.42
1:C:219:LEU:HB3	1:C:224:LEU:HD21	2.02	0.42
1:C:234:ARG:H	1:C:234:ARG:HG3	1.40	0.42
1:C:249:VAL:HG21	1:C:257:TYR:CD2	2.55	0.42
3:E:7:SER:HB3	3:E:22:ARG:CB	2.46	0.42
3:G:114:VAL:HG12	3:G:115:THR:N	2.35	0.42
3:G:128:VAL:HG11	3:G:201:SER:HB3	2.01	0.42
1:A:55:GLU:O	1:A:57:ALA:N	2.53	0.42
1:A:84:PHE:CD2	1:A:84:PHE:O	2.73	0.42
3:E:90:ALA:O	3:E:107:VAL:HG22	2.20	0.42
3:E:169:SER:HB3	3:E:170:PRO:HD2	2.01	0.42
4:F:39:ARG:HG3	4:F:94:TYR:CE2	2.55	0.42
1:A:65:ARG:NH1	1:A:158:GLN:CG	2.83	0.42
1:A:68:THR:O	1:A:69:ILE:C	2.62	0.42
1:A:185:PRO:HD2	1:A:266:LEU:HD23	2.00	0.42
1:C:97:TRP:CD1	1:C:97:TRP:C	2.98	0.42
1:C:233:THR:H	1:C:233:THR:HG22	1.49	0.42
3:E:40:ARG:HA	3:E:40:ARG:HD3	1.67	0.42
3:E:55:GLU:CB	3:E:60:LYS:HG2	2.49	0.42
4:F:132:PRO:C	4:F:133:THR:OG1	2.62	0.42
3:G:26:THR:CG2	3:G:69:ARG:HH11	2.33	0.42
1:A:193:PRO:HG3	1:A:199:VAL:HG22	2.01	0.41
1:C:151:SER:O	1:C:153:VAL:HG23	2.20	0.41
1:C:266:LEU:O	1:C:267:PRO:C	2.63	0.41
2:D:65:LEU:HD12	2:D:66:TYR:H	1.84	0.41
4:F:20:GLN:HA	4:F:82:THR:HA	2.02	0.41
4:F:88:GLU:C	4:F:90:ASP:N	2.70	0.41
4:F:186:ASN:HD21	6:K:1:NAG:C5	2.31	0.41
4:H:188:THR:C	4:H:189:TYR:CD1	2.98	0.41
1:A:64:THR:O	1:A:68:THR:OG1	2.38	0.41
1:A:263:HIS:CG	1:A:264:GLU:H	2.37	0.41
2:B:23:LEU:HD23	2:B:39:LEU:CD1	2.43	0.41
1:C:10:THR:HG21	2:D:62:PHE:CE2	2.55	0.41
1:C:234:ARG:HD3	2:D:10:TYR:CZ	2.55	0.41
2:D:79:ALA:CB	2:D:94:LYS:HA	2.49	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:G:190:CYS:H	3:G:201:SER:HG	1.67	0.41
4:H:28:GLU:C	4:H:30:PHE:H	2.21	0.41
1:A:8:PHE:CD1	1:A:8:PHE:N	2.88	0.41
1:C:39:LYS:O	1:C:39:LYS:HG2	2.20	0.41
1:C:249:VAL:CG2	1:C:257:TYR:CE2	3.02	0.41
3:E:92:THR:HG22	3:E:93:TRP:O	2.20	0.41
3:E:139:CYS:O	3:E:177:VAL:HA	2.21	0.41
3:G:25:PHE:CD1	3:G:25:PHE:N	2.87	0.41
3:G:101:LEU:HD23	3:G:101:LEU:HA	1.79	0.41
1:A:61:GLU:O	1:A:62:GLN:C	2.63	0.41
1:A:73:LEU:HD11	3:E:98:GLY:O	2.21	0.41
1:C:23:ILE:CG2	1:C:24:ILE:N	2.70	0.41
1:C:204:TRP:HE3	1:C:206:LEU:HD21	1.83	0.41
2:D:41:LYS:O	2:D:43:GLY:N	2.53	0.41
3:G:20:ALA:CB	3:G:74:HIS:HA	2.50	0.41
2:B:23:LEU:HG	2:B:24:ASN:N	2.36	0.41
1:C:126:LEU:N	3:G:92:THR:HG21	2.35	0.41
1:C:184:PRO:CG	1:C:265:GLY:O	2.64	0.41
2:D:40:LEU:HA	2:D:44:GLU:O	2.19	0.41
3:E:35:PHE:N	3:E:87:PHE:O	2.47	0.41
3:E:64:ASP:HB3	3:E:67:GLU:OE1	2.20	0.41
4:F:6:GLN:HA	4:F:7:PRO:HD3	1.84	0.41
4:F:86:LEU:HD23	4:F:86:LEU:HA	1.73	0.41
4:F:137:PRO:HD3	4:F:150:TYR:CD2	2.56	0.41
4:F:199:GLU:HA	4:F:202:MET:HE2	2.02	0.41
1:C:231:VAL:O	1:C:232:GLU:C	2.63	0.41
2:D:19:LYS:HD3	2:D:19:LYS:HA	1.84	0.41
3:E:13:LEU:HB3	3:E:17:THR:CG2	2.49	0.41
4:F:73:ASN:N	4:F:74:PRO:HD3	2.34	0.41
4:F:136:LEU:HB3	4:F:137:PRO:HD2	2.02	0.41
3:G:5:GLU:HA	3:G:5:GLU:OE1	2.20	0.41
3:G:112:THR:O	3:G:112:THR:HG22	2.20	0.41
3:G:163:ASP:HA	3:G:164:PRO:HD3	1.89	0.41
4:H:3:LYS:HE3	4:H:27:ILE:CD1	2.50	0.41
4:H:91:GLU:HG3	4:H:115:LEU:O	2.20	0.41
1:A:85:TYR:O	1:A:86:ASN:C	2.63	0.41
2:B:31:HIS:O	2:B:33:SER:N	2.54	0.41
1:C:219:LEU:O	1:C:220:ASN:CB	2.69	0.41
3:E:62:ALA:O	3:E:71:SER:HA	2.21	0.41
3:E:99:TYR:N	3:E:99:TYR:CD1	2.88	0.41
4:F:51:TYR:CD2	4:F:59:ILE:HD13	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:H:141:GLU:O	4:H:142:THR:C	2.62	0.41
2:B:42:ASN:HD21	2:B:76:ASP:CA	2.30	0.41
1:C:209:TYR:O	1:C:210:PRO:C	2.60	0.41
2:D:31:HIS:CB	2:D:32:PRO:CD	2.99	0.41
2:D:75:LYS:HD3	2:D:75:LYS:HA	1.93	0.41
2:D:99:MET:SD	2:D:99:MET:N	2.94	0.41
4:F:69:GLU:CB	4:F:82:THR:O	2.68	0.41
4:F:118:ILE:HD11	4:F:124:LEU:HD13	2.02	0.41
4:H:6:GLN:HG2	4:H:23:CYS:HB2	2.03	0.41
4:H:49:LEU:O	4:H:50:LEU:HG	2.21	0.41
2:B:23:LEU:HD21	2:B:39:LEU:HD22	2.03	0.41
2:B:26:TYR:CD1	2:B:65:LEU:HB2	2.53	0.41
1:C:6:ARG:CD	1:C:100:GLY:HA3	2.50	0.41
1:C:42:THR:HA	1:C:43:PRO:HD3	1.85	0.41
1:C:51:TRP:CH2	1:C:179:LEU:HD21	2.55	0.41
1:C:62:GLN:O	1:C:66:ILE:HG13	2.21	0.41
1:C:84:PHE:HD1	4:H:103:SER:O	2.03	0.41
1:C:127:ASN:C	1:C:128:GLU:O	2.62	0.41
1:C:202:ARG:HD3	2:D:98:ASP:O	2.21	0.41
2:D:59:ASP:OD1	2:D:59:ASP:O	2.39	0.41
3:E:14:HIS:O	3:E:15:GLU:C	2.64	0.41
3:E:61:SER:HA	3:E:72:THR:O	2.21	0.41
4:F:76:ALA:O	4:F:77:SER:C	2.64	0.41
4:F:199:GLU:C	4:F:201:ALA:N	2.78	0.41
3:G:15:GLU:N	3:G:117:GLU:O	2.54	0.41
4:H:73:ASN:OD1	4:H:73:ASN:C	2.64	0.41
4:H:99:GLY:HA3	4:H:105:PHE:CE2	2.55	0.41
1:A:17:LEU:HD23	1:A:17:LEU:HA	1.79	0.41
1:A:207:GLY:HA2	1:A:240:THR:CB	2.51	0.41
1:C:10:THR:HG21	2:D:62:PHE:HE2	1.86	0.41
2:D:83:ASN:HB2	2:D:90:PRO:HB3	2.03	0.41
3:E:17:THR:CG2	3:E:18:ASP:N	2.83	0.41
4:F:145:HIS:O	4:F:146:LYS:CB	2.66	0.41
4:H:13:ARG:HA	4:H:14:PRO:HD3	1.92	0.41
4:H:53:LEU:O	4:H:53:LEU:CG	2.69	0.41
4:H:60:PHE:CD1	4:H:62:ASP:OD1	2.74	0.41
4:H:157:PHE:CE1	4:H:189:TYR:HB2	2.56	0.41
1:A:14:ARG:CB	1:A:15:PRO:HD3	2.50	0.40
1:A:78:LEU:HD13	1:A:95:LEU:HB3	2.03	0.40
1:A:219:LEU:CD2	1:A:256:SER:O	2.69	0.40
2:B:55:SER:HB3	2:B:63:TYR:OH	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:156:ILE:O	1:C:156:ILE:HG22	2.21	0.40
1:C:164:CYS:O	1:C:166:ASP:N	2.54	0.40
4:F:128:ILE:CG2	4:F:158:PHE:CB	2.93	0.40
3:G:3:GLN:NE2	3:G:26:THR:N	2.52	0.40
3:G:55:GLU:HB2	3:G:60:LYS:HE2	2.03	0.40
1:C:184:PRO:HA	1:C:185:PRO:HD3	1.81	0.40
3:E:145:TYR:CB	3:E:146:PRO:CD	2.99	0.40
4:F:14:PRO:HB2	4:F:17:GLU:CG	2.51	0.40
4:F:91:GLU:OE2	4:F:116:ILE:HA	2.21	0.40
4:F:134:ILE:C	4:F:135:PHE:HD1	2.29	0.40
1:A:73:LEU:HD12	3:E:99:TYR:O	2.22	0.40
1:A:159:ASP:C	1:A:161:LYS:N	2.79	0.40
1:A:261:VAL:HB	1:A:270:LEU:HB3	2.03	0.40
1:A:266:LEU:HA	1:A:267:PRO:HD3	1.95	0.40
1:C:155:HIS:CD2	1:C:157:SER:OG	2.74	0.40
1:C:215:LEU:HD12	1:C:245:ALA:CB	2.52	0.40
4:H:81:LEU:HA	4:H:81:LEU:HD12	1.90	0.40
4:H:96:CYS:O	4:H:96:CYS:SG	2.79	0.40
1:A:36:PHE:HA	1:A:40:GLU:OE1	2.21	0.40
1:C:14:ARG:HA	1:C:15:PRO:HD3	1.93	0.40
4:F:52:VAL:CG2	4:F:71:SER:HB2	2.47	0.40
4:F:200:ARG:HD3	4:F:200:ARG:HA	1.57	0.40
4:F:202:MET:O	4:F:227:SER:HB2	2.22	0.40
3:G:81:GLU:H	3:G:81:GLU:HG2	1.70	0.40
1:C:28:VAL:HG11	1:C:172:LEU:HD21	2.04	0.40
3:G:8:PRO:C	3:G:10:ALA:N	2.78	0.40
3:G:58:ARG:CB	3:G:76:ARG:O	2.67	0.40

All (5) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:255:GLN:NE2	4:F:66:LYS:CE[1_545]	1.35	0.85
1:A:53:GLU:OE2	3:E:157:LYS:NZ[1_455]	1.36	0.84
1:C:255:GLN:CD	4:F:66:LYS:NZ[1_545]	1.96	0.24
1:C:255:GLN:CD	4:F:66:LYS:CE[1_545]	2.13	0.07
1:C:255:GLN:NE2	4:F:66:LYS:NZ[1_545]	2.17	0.03

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 X-ray entries followed by that with respect to entries of similar resolution.

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	258/260 (99%)	181 (70%)	55 (21%)	22 (8%)	0	4
1	C	258/260 (99%)	192 (74%)	45 (17%)	21 (8%)	1	4
2	B	100/102 (98%)	86 (86%)	9 (9%)	5 (5%)	1	10
2	D	100/102 (98%)	79 (79%)	12 (12%)	9 (9%)	0	3
3	E	205/207 (99%)	157 (77%)	30 (15%)	18 (9%)	0	3
3	G	205/207 (99%)	158 (77%)	29 (14%)	18 (9%)	0	3
4	F	228/230 (99%)	167 (73%)	37 (16%)	24 (10%)	0	2
4	H	228/230 (99%)	172 (75%)	34 (15%)	22 (10%)	0	3
All	All	1582/1598 (99%)	1192 (75%)	251 (16%)	139 (9%)	0	3

All (139) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	14	ARG
1	A	15	PRO
1	A	51	TRP
1	A	56	GLU
1	A	129	ASN
1	A	197	GLY
1	A	209	TYR
1	A	220	ASN
1	A	254	GLU
2	B	47	GLU
2	B	60	TRP
2	B	74	GLU
1	C	15	PRO
1	C	56	GLU
1	C	106	ASP
1	C	129	ASN
1	C	150	ASN

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Mol	Chain	Res	Type
1	C	158	GLN
1	C	255	GLN
2	D	31	HIS
2	D	33	SER
2	D	85	VAL
3	E	7	SER
3	E	39	SER
3	E	185	SER
4	F	7	PRO
4	F	56	PRO
4	F	107	LYS
4	F	119	PRO
4	F	122	LYS
4	F	133	THR
4	F	186	ASN
4	F	213	GLU
4	F	214	ASN
4	F	216	LYS
3	G	7	SER
3	G	122	PRO
3	G	188	VAL
4	H	56	PRO
4	H	64	GLU
4	H	121	ASP
4	H	133	THR
4	H	172	ASN
4	H	199	GLU
4	H	209	ILE
1	A	120	SER
1	A	176	LYS
1	A	221	GLY
1	C	212	ASP
1	C	233	THR
1	C	263	HIS
3	E	65	SER
3	E	78	ALA
3	E	97	GLU
3	E	105	LYS
3	E	123	PRO
3	E	134	GLY
3	E	188	VAL
3	E	203	ASP

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Mol	Chain	Res	Type
4	F	4	LEU
4	F	30	PHE
4	F	76	ALA
4	F	89	GLU
4	F	100	GLU
4	F	199	GLU
3	G	28	THR
3	G	52	GLY
3	G	67	GLU
3	G	133	ASN
3	G	145	TYR
3	G	165	ALA
3	G	185	SER
3	G	187	SER
4	H	45	GLY
4	H	85	SER
4	H	107	LYS
4	H	168	GLU
1	A	89	MET
1	A	118	TYR
1	A	158	GLN
1	A	193	PRO
1	A	275	GLY
1	C	24	ILE
1	C	124	PRO
1	C	130	PRO
1	C	133	CYS
1	C	207	GLY
1	C	267	PRO
2	D	1	ILE
2	D	17	ASN
2	D	32	PRO
2	D	90	PRO
3	E	103	THR
4	F	42	PRO
4	F	55	THR
4	F	132	PRO
4	F	159	PRO
4	F	200	ARG
3	G	80	LEU
3	G	96	SER
3	G	123	PRO

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Mol	Chain	Res	Type
3	G	164	PRO
4	H	128	ILE
4	H	226	PRO
1	A	153	VAL
1	C	44	ARG
1	C	210	PRO
1	C	254	GLU
2	D	54	LEU
3	E	51	SER
3	E	145	TYR
4	H	74	PRO
4	H	198	PRO
4	H	216	LYS
4	H	218	GLY
1	A	50	PRO
2	D	88	SER
3	E	16	GLY
3	E	80	LEU
4	F	14	PRO
4	F	158	PHE
3	G	169	SER
4	H	14	PRO
4	H	42	PRO
4	H	158	PHE
1	A	237	GLY
3	E	159	ILE
1	A	104	GLU
2	B	31	HIS
1	C	20	PRO
1	C	104	GLU
4	F	226	PRO
1	A	267	PRO
3	E	57	GLY
3	G	167	VAL
2	B	14	PRO
3	G	170	PRO
4	H	41	LYS
4	H	119	PRO

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 X-ray entries followed by that with respect to entries of similar resolution.

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	233/233 (100%)	205 (88%)	28 (12%)	5	19
1	C	233/233 (100%)	208 (89%)	25 (11%)	6	23
2	B	96/96 (100%)	85 (88%)	11 (12%)	5	20
2	D	96/96 (100%)	88 (92%)	8 (8%)	10	35
3	E	180/180 (100%)	163 (91%)	17 (9%)	8	29
3	G	177/180 (98%)	152 (86%)	25 (14%)	3	13
4	F	206/206 (100%)	175 (85%)	31 (15%)	3	11
4	H	204/206 (99%)	176 (86%)	28 (14%)	3	14
All	All	1425/1430 (100%)	1252 (88%)	173 (12%)	5	19

All (173) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	PHE
1	A	38	SER
1	A	68	THR
1	A	75	GLU
1	A	94	THR
1	A	98	LEU
1	A	99	GLN
1	A	102	ASP
1	A	108	HIS
1	A	120	SER
1	A	131	SER
1	A	133	CYS
1	A	151	SER
1	A	152	THR
1	A	181	ARG
1	A	196	GLU
1	A	198	ASP
1	A	213	ILE
1	A	214	THR

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Mol	Chain	Res	Type
1	A	216	THR
1	A	231	VAL
1	A	233	THR
1	A	234	ARG
1	A	235	PRO
1	A	242	GLN
1	A	247	VAL
1	A	251	LEU
1	A	270	LEU
2	B	4	THR
2	B	12	ARG
2	B	55	SER
2	B	57	SER
2	B	65	LEU
2	B	73	THR
2	B	85	VAL
2	B	87	LEU
2	B	94	LYS
2	B	98	ASP
2	B	99	MET
1	C	7	TYR
1	C	8	PHE
1	C	9	TYR
1	C	25	VAL
1	C	37	SER
1	C	53	GLU
1	C	54	GLN
1	C	58	ASP
1	C	78	LEU
1	C	83	HIS
1	C	86	ASN
1	C	94	THR
1	C	95	LEU
1	C	98	LEU
1	C	101	CYS
1	C	111	LEU
1	C	112	TRP
1	C	125	THR
1	C	128	GLU
1	C	213	ILE
1	C	214	THR
1	C	215	LEU

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Mol	Chain	Res	Type
1	C	225	THR
1	C	233	THR
1	C	234	ARG
2	D	4	THR
2	D	7	ILE
2	D	17	ASN
2	D	51	HIS
2	D	67	TYR
2	D	84	HIS
2	D	94	LYS
2	D	99	MET
3	E	2	ASP
3	E	13	LEU
3	E	25	PHE
3	E	30	ARG
3	E	32	VAL
3	E	46	LEU
3	E	72	THR
3	E	92	THR
3	E	110	GLN
3	E	132	LYS
3	E	133	ASN
3	E	148	GLU
3	E	150	THR
3	E	152	SER
3	E	153	LEU
3	E	188	VAL
3	E	189	THR
4	F	11	ILE
4	F	18	THR
4	F	34	THR
4	F	42	PRO
4	F	51	TYR
4	F	52	VAL
4	F	53	LEU
4	F	55	THR
4	F	79	SER
4	F	81	LEU
4	F	85	SER
4	F	100	GLU
4	F	106	HIS
4	F	118	ILE

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Mol	Chain	Res	Type
4	F	121	ASP
4	F	133	THR
4	F	135	PHE
4	F	144	LEU
4	F	153	LEU
4	F	155	GLU
4	F	159	PRO
4	F	160	ASP
4	F	177	SER
4	F	183	LEU
4	F	185	THR
4	F	192	PHE
4	F	212	HIS
4	F	215	ASN
4	F	221	GLN
4	F	223	ILE
4	F	224	PHE
3	G	4	VAL
3	G	11	LEU
3	G	12	SER
3	G	13	LEU
3	G	17	THR
3	G	23	CYS
3	G	25	PHE
3	G	28	THR
3	G	45	SER
3	G	64	ASP
3	G	92	THR
3	G	101	LEU
3	G	104	ASP
3	G	108	PHE
3	G	115	THR
3	G	120	SER
3	G	122	PRO
3	G	133	ASN
3	G	135	THR
3	G	140	LEU
3	G	143	ASP
3	G	163	ASP
3	G	181	GLN
3	G	191	SER
3	G	194	HIS

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Mol	Chain	Res	Type
4	H	9	ILE
4	H	14	PRO
4	H	21	ILE
4	H	34	THR
4	H	42	PRO
4	H	51	TYR
4	H	56	PRO
4	H	60	PHE
4	H	62	ASP
4	H	64	GLU
4	H	117	VAL
4	H	118	ILE
4	H	133	THR
4	H	143	ASN
4	H	144	LEU
4	H	151	LEU
4	H	161	VAL
4	H	163	ARG
4	H	165	TYR
4	H	183	LEU
4	H	184	LYS
4	H	187	ASP
4	H	189	TYR
4	H	190	MET
4	H	191	LYS
4	H	215	ASN
4	H	228	ILE
4	H	229	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (45) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	59	ASN
1	A	63	GLN
1	A	70	GLN
1	A	96	GLN
1	A	99	GLN
1	A	127	ASN
1	A	163	HIS
1	A	169	GLN
2	B	24	ASN
2	B	42	ASN

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Mol	Chain	Res	Type
1	C	32	GLN
1	C	54	GLN
1	C	83	HIS
1	C	86	ASN
1	C	96	GLN
1	C	99	GLN
1	C	108	HIS
1	C	155	HIS
1	C	218	GLN
1	C	220	ASN
2	D	24	ASN
2	D	42	ASN
2	D	83	ASN
3	E	6	GLN
3	E	33	GLN
3	E	37	GLN
3	E	74	HIS
3	E	133	ASN
4	F	6	GLN
4	F	36	HIS
4	F	44	GLN
4	F	73	ASN
4	F	186	ASN
4	F	206	HIS
4	F	214	ASN
3	G	3	GLN
3	G	6	GLN
3	G	24	ASN
3	G	33	GLN
3	G	38	ASN
3	G	136	ASN
4	H	73	ASN
4	H	106	HIS
4	H	206	HIS
4	H	221	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

13 monosaccharides are modelled in this entry.

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	NAG	I	1	5,3	15,15,15	0.56	0	21,21,21	0.77	0
5	NAG	I	2	5	15,15,15	0.47	0	21,21,21	0.70	0
5	MAN	I	3	5	12,12,12	0.44	0	17,17,17	0.45	0
6	NAG	J	1	3,6	15,15,15	0.56	0	21,21,21	0.79	0
6	NAG	J	2	6	15,15,15	0.48	0	21,21,21	0.70	0
6	NAG	K	1	4,6	15,15,15	0.56	0	21,21,21	1.13	2 (9%)
6	NAG	K	2	6	15,15,15	0.48	0	21,21,21	0.70	0
5	NAG	L	1	5,3	15,15,15	0.57	0	21,21,21	0.97	1 (4%)
5	NAG	L	2	5	15,15,15	0.48	0	21,21,21	0.70	0
5	MAN	L	3	5	12,12,12	0.44	0	17,17,17	0.45	0
7	NAG	M	1	3,7	15,15,15	1.20	1 (6%)	21,21,21	1.79	5 (23%)
7	NAG	M	2	7	14,14,15	0.78	0	17,19,21	0.84	0
7	FUC	M	3	7	10,10,11	0.97	1 (10%)	14,14,16	2.01	2 (14%)

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	NAG	I	1	5,3	1/1/6/7	0/6/26/26	0/1/1/1
5	NAG	I	2	5	-	4/6/26/26	0/1/1/1
5	MAN	I	3	5	-	2/2/22/22	0/1/1/1
6	NAG	J	1	3,6	1/1/6/7	0/6/26/26	0/1/1/1
6	NAG	J	2	6	-	4/6/26/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	NAG	K	1	4,6	-	0/6/26/26	0/1/1/1
6	NAG	K	2	6	-	4/6/26/26	0/1/1/1
5	NAG	L	1	5,3	1/1/6/7	0/6/26/26	0/1/1/1
5	NAG	L	2	5	-	4/6/26/26	0/1/1/1
5	MAN	L	3	5	-	2/2/22/22	0/1/1/1
7	NAG	M	1	3,7	1/1/6/7	5/6/26/26	0/1/1/1
7	NAG	M	2	7	-	5/6/23/26	0/1/1/1
7	FUC	M	3	7	-	-	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	M	1	NAG	C1-C2	-3.52	1.48	1.52
7	M	3	FUC	C1-C2	2.20	1.57	1.52

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	M	3	FUC	C1-C2-C3	6.49	119.10	109.64
7	M	1	NAG	C3-C4-C5	4.12	117.70	110.23
7	M	1	NAG	C1-C2-C3	-3.60	105.63	110.54
6	K	1	NAG	O1-C1-C2	-3.27	102.42	109.22
7	M	1	NAG	C4-C3-C2	2.54	114.09	110.40
5	L	1	NAG	O1-C1-C2	-2.40	104.23	109.22
7	M	1	NAG	O4-C4-C3	-2.38	104.75	110.38
7	M	1	NAG	O5-C1-C2	-2.30	107.20	109.52
7	M	3	FUC	C2-C3-C4	2.06	114.49	110.86
6	K	1	NAG	O1-C1-O5	-2.02	104.42	110.41

All (4) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
5	I	1	NAG	C1
5	L	1	NAG	C1
6	J	1	NAG	C1
7	M	1	NAG	C1

All (30) torsion outliers are listed below:

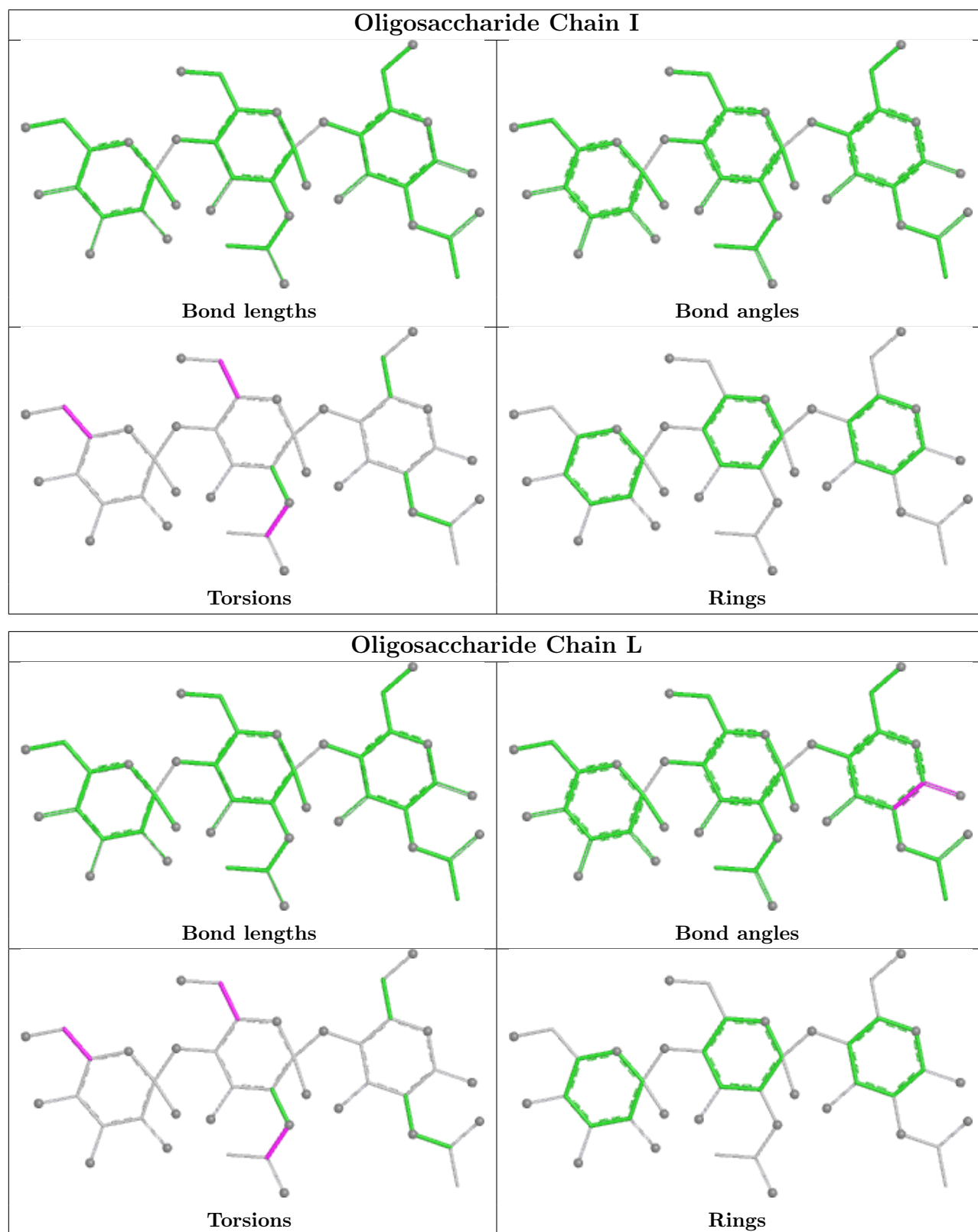
Mol	Chain	Res	Type	Atoms
5	I	2	NAG	C8-C7-N2-C2
5	I	2	NAG	O7-C7-N2-C2
5	L	2	NAG	C8-C7-N2-C2
5	L	2	NAG	O7-C7-N2-C2
6	J	2	NAG	C8-C7-N2-C2
6	J	2	NAG	O7-C7-N2-C2
6	K	2	NAG	C8-C7-N2-C2
6	K	2	NAG	O7-C7-N2-C2
7	M	1	NAG	C3-C2-N2-C7
7	M	1	NAG	C8-C7-N2-C2
7	M	1	NAG	O7-C7-N2-C2
7	M	2	NAG	C8-C7-N2-C2
7	M	2	NAG	O7-C7-N2-C2
7	M	1	NAG	C4-C5-C6-O6
7	M	1	NAG	O5-C5-C6-O6
5	I	3	MAN	C4-C5-C6-O6
5	L	3	MAN	C4-C5-C6-O6
5	I	2	NAG	C4-C5-C6-O6
5	L	2	NAG	C4-C5-C6-O6
6	J	2	NAG	C4-C5-C6-O6
6	K	2	NAG	C4-C5-C6-O6
5	I	2	NAG	O5-C5-C6-O6
5	L	2	NAG	O5-C5-C6-O6
6	J	2	NAG	O5-C5-C6-O6
6	K	2	NAG	O5-C5-C6-O6
7	M	2	NAG	O5-C5-C6-O6
7	M	2	NAG	C3-C2-N2-C7
5	L	3	MAN	O5-C5-C6-O6
5	I	3	MAN	O5-C5-C6-O6
7	M	2	NAG	C1-C2-N2-C7

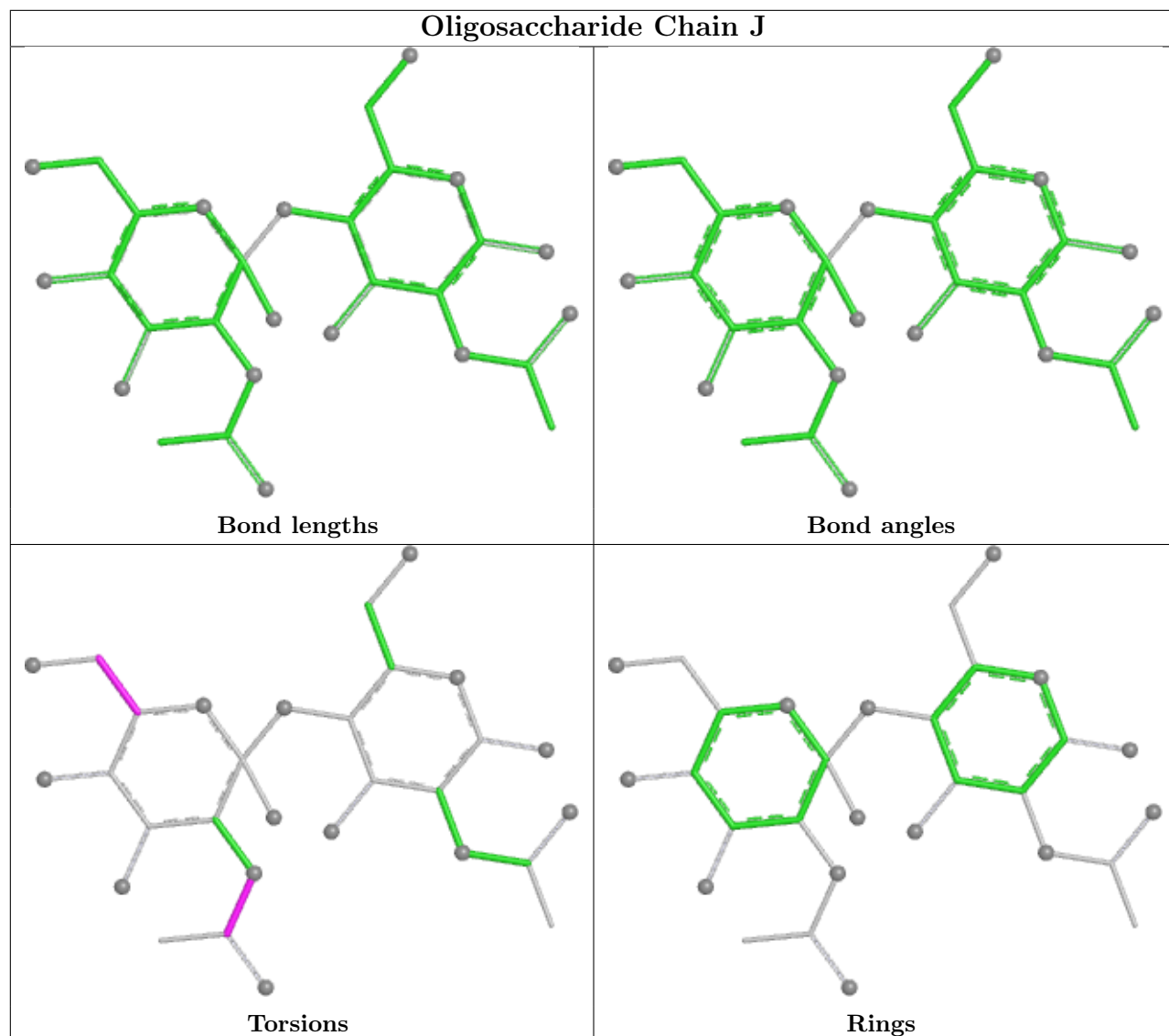
There are no ring outliers.

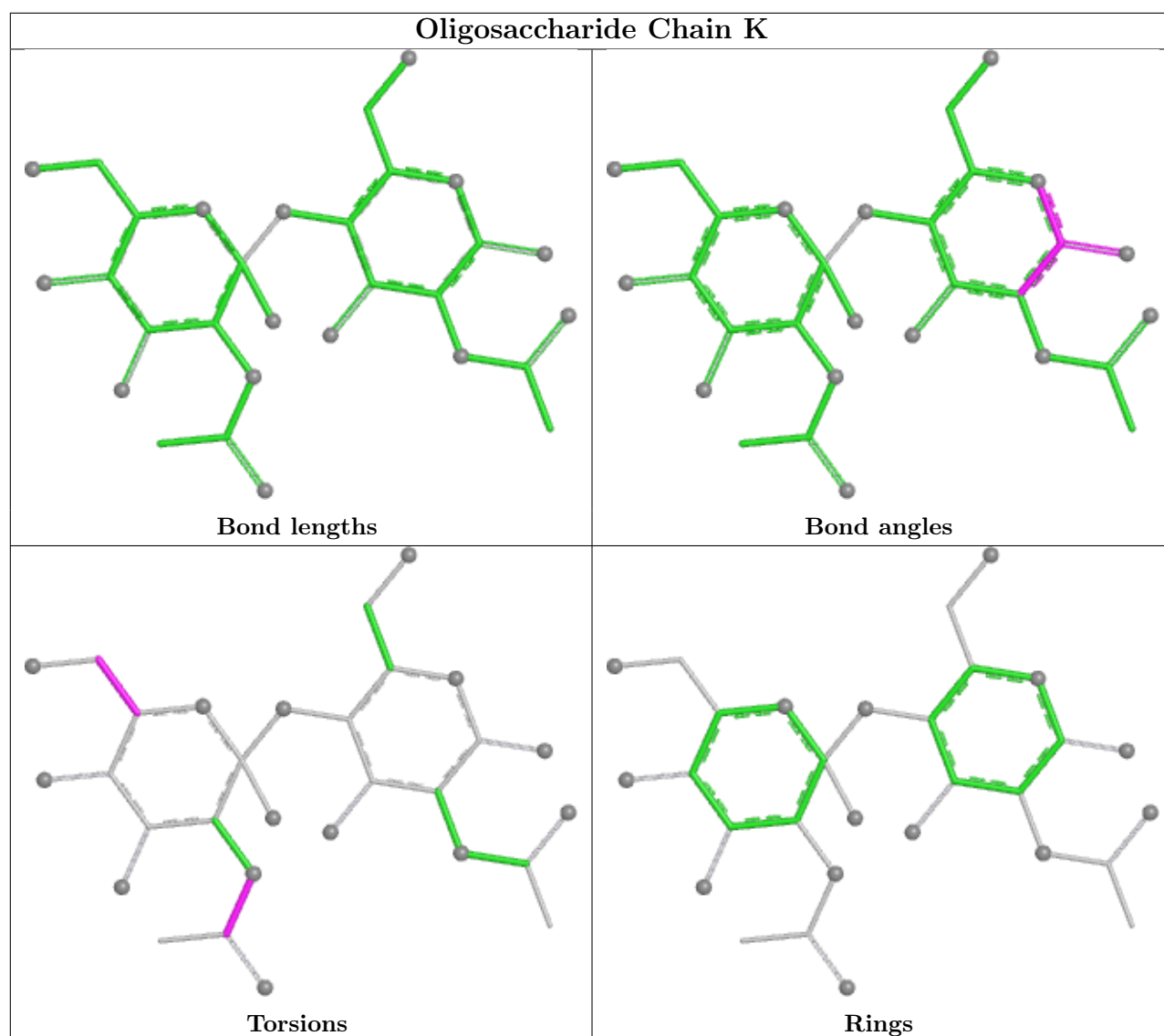
8 monomers are involved in 60 short contacts:

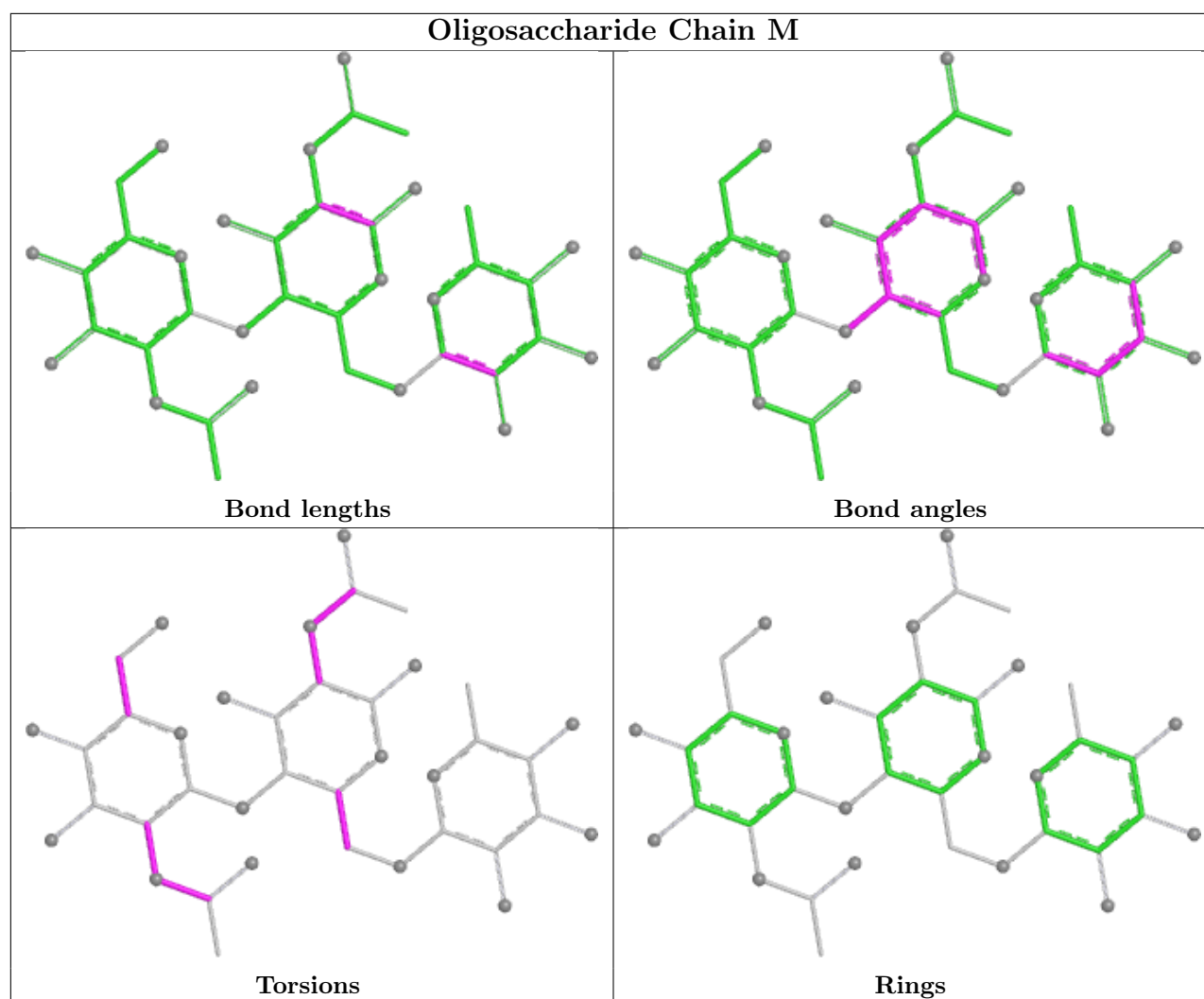
Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	J	2	NAG	1	0
5	I	3	MAN	3	0
7	M	2	NAG	2	0
6	K	1	NAG	22	0
5	L	1	NAG	7	0
6	J	1	NAG	8	0
7	M	1	NAG	10	0
5	I	1	NAG	7	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.









5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.