



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 28, 2014 – 05:42 AM GMT

PDB ID : 1OCO
Title : BOVINE HEART CYTOCHROME C OXIDASE IN CARBON MONOXIDE-
BOUND STATE
Authors : Tsukihara, T.; Yao, M.
Deposited on : 1998-07-09
Resolution : 2.80 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

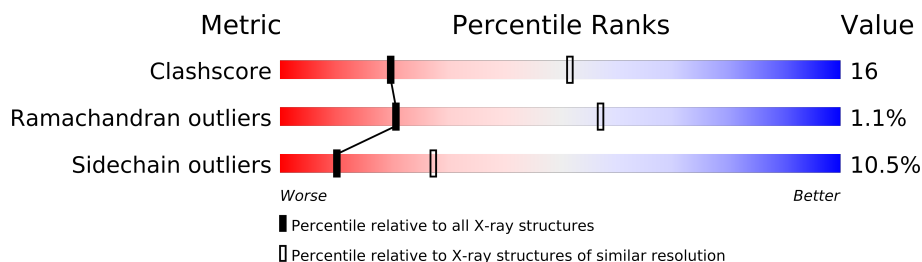
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 21963
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 2.80 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	79885	2295 (2.80-2.80)
Ramachandran outliers	78287	2252 (2.80-2.80)
Sidechain outliers	78261	2254 (2.80-2.80)







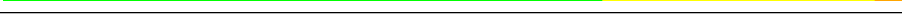

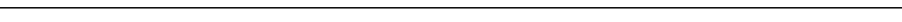
The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	514	
1	N	514	
2	B	227	
2	O	227	
3	C	261	
3	P	261	
4	D	147	
4	Q	147	
5	E	109	
5	R	109	
6	F	98	
6	S	98	
7	G	84	
7	T	84	
8	H	85	
8	U	85	
9	I	73	

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Mol	Chain	Length	Quality of chain
9	V	73	
10	J	59	
10	W	59	
11	K	56	
11	X	56	
12	L	47	
12	Y	47	
13	M	46	
13	Z	46	

2 Entry composition

There are 19 unique types of molecules in this entry. The entry contains 28810 atoms, of which 0 are hydrogen and 0 are deuterium.

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 CYTOCHROME C OXIDASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	514	Total	C	N	O	S	0	0	0
			4025	2690	623	677	35			
1	N	514	Total	C	N	O	S	0	0	0
			4025	2690	623	677	35			

- Molecule 2 is a protein called CYTOCHROME C OXIDASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	227	Total	C	N	O	S	0	5	0
			1863	1207	288	350	18			
2	O	227	Total	C	N	O	S	0	5	0
			1863	1207	288	350	18			

- Molecule 3 is a protein called CYTOCHROME C OXIDASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	261	Total	C	N	O	S	0	0	0
			2124	1420	338	353	13			
3	P	261	Total	C	N	O	S	0	0	0
			2124	1420	338	353	13			

- Molecule 4 is a protein called CYTOCHROME C OXIDASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	144	Total	C	N	O	S	0	0	0
			1195	777	196	218	4			
4	Q	144	Total	C	N	O	S	0	0	0
			1195	777	196	218	4			

- Molecule 5 is a protein called CYTOCHROME C OXIDASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	E	109	Total	C	N	O	S	0	0	0
			878	558	150	168	2			
5	R	109	Total	C	N	O	S	0	0	0
			878	558	150	168	2			

- Molecule 6 is a protein called CYTOCHROME C OXIDASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
6	F	98	Total	C	N	O	S	0	0	0
			748	464	134	145	5			
6	S	98	Total	C	N	O	S	0	0	0
			748	464	134	145	5			

- Molecule 7 is a protein called CYTOCHROME C OXIDASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
7	G	84	Total	C	N	O	S	0	0	0
			672	431	129	111	1			
7	T	84	Total	C	N	O	S	0	0	0
			672	431	129	111	1			

- Molecule 8 is a protein called CYTOCHROME C OXIDASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
8	H	75	Total	C	N	O	S	0	0	0
			628	395	114	114	5			
8	U	75	Total	C	N	O	S	0	0	0
			628	395	114	114	5			

- Molecule 9 is a protein called CYTOCHROME C OXIDASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
9	I	73	Total	C	N	O	S	0	0	0
			598	388	107	99	4			
9	V	73	Total	C	N	O	S	0	0	0
			598	388	107	99	4			

- Molecule 10 is a protein called CYTOCHROME C OXIDASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
10	J	56	Total	C	N	O	S	0	0	0
			441	285	73	80	3			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
10	W	56	Total	C	N	O	S	0	0	0
			441	285	73	80	3			

- Molecule 11 is a protein called CYTOCHROME C OXIDASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
11	K	49	Total	C	N	O	S	0	0	0
			384	250	65	67	2			
11	X	49	Total	C	N	O	S	0	0	0
			384	250	65	67	2			

- Molecule 12 is a protein called CYTOCHROME C OXIDASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
12	L	47	Total	C	N	O	S	0	0	0
			386	257	65	62	2			
12	Y	47	Total	C	N	O	S	0	0	0
			386	257	65	62	2			

- Molecule 13 is a protein called CYTOCHROME C OXIDASE.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
13	M	43	Total	C	N	O	0	0	0
			335	223	53	59			
13	Z	43	Total	C	N	O	0	0	0
			335	223	53	59			

- Molecule 14 is COPPER (II) ION (three-letter code: CU) (formula: Cu).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
14	O	2	Total	Cu	0	0
			2	2		
14	B	2	Total	Cu	0	0
			2	2		
14	A	1	Total	Cu	0	0
			1	1		
14	N	1	Total	Cu	0	0
			1	1		

- Molecule 15 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
15	A	1	Total Mg 1 1	0	0
15	N	1	Total Mg 1 1	0	0

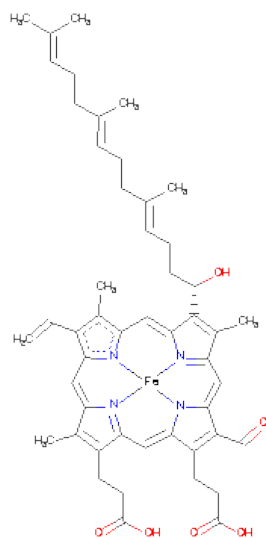
- Molecule 16 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
16	A	1	Total Na 1 1	0	0
16	N	1	Total Na 1 1	0	0

- Molecule 17 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
17	S	1	Total Zn 1 1	0	0
17	F	1	Total Zn 1 1	0	0

- Molecule 18 is HEME-A (three-letter code: HEA) (formula: C₄₉H₅₆FeN₄O₆).



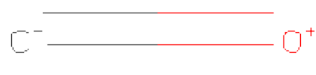
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
18	A	1	Total C Fe N O 60 49 1 4 6	0	0

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
18	A	1	Total	C	Fe	N	O	0	0
			60	49	1	4	6		
18	N	1	Total	C	Fe	N	O	0	0
			60	49	1	4	6		
18	N	1	Total	C	Fe	N	O	0	0
			60	49	1	4	6		

- Molecule 19 is CARBON MONOXIDE (three-letter code: CMO) (formula: CO).



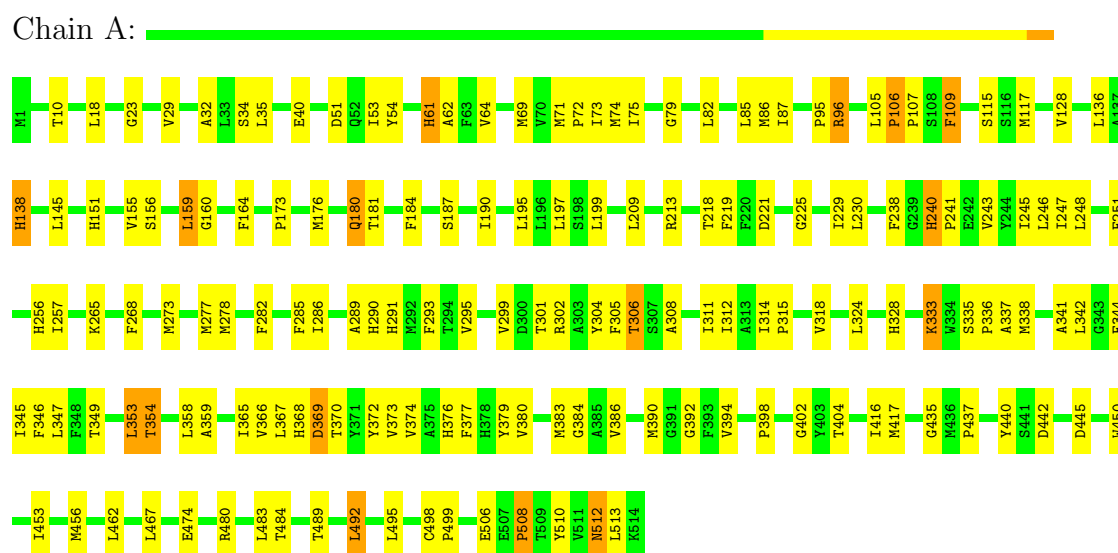
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
19	A	1	Total	C	O	0	0
			2	1	1		
19	N	1	Total	C	O	0	0
			2	1	1		

3 Residue-property plots

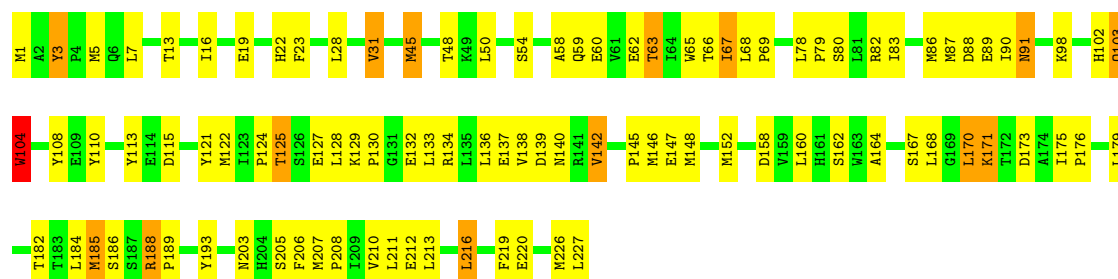
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according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

Note EDS was not executed.

• Molecule 1: CYTOCHROME C OXIDASE



Chain B: 



• Molecule 2: CYTOCHROME C OXIDASE

Chain O: 



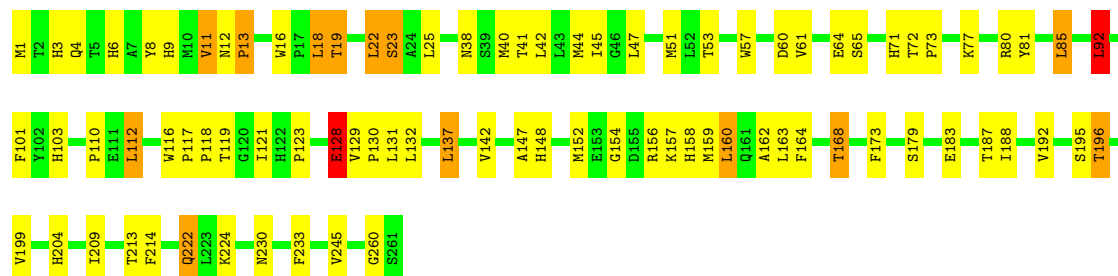
• Molecule 3: CYTOCHROME C OXIDASE

Chain C: 



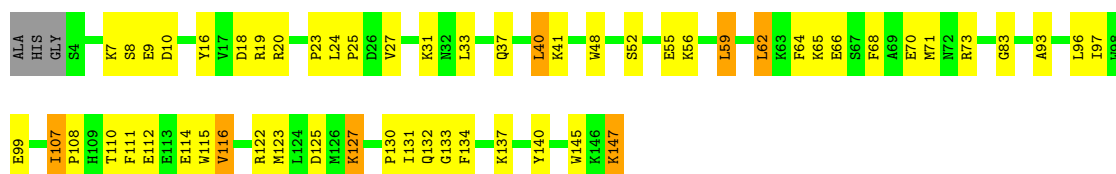
• Molecule 3: CYTOCHROME C OXIDASE

Chain P: 



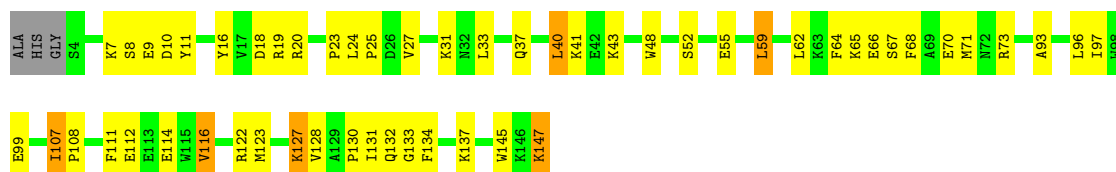
• Molecule 4: CYTOCHROME C OXIDASE

Chain D: 



• Molecule 4: CYTOCHROME C OXIDASE

Chain Q:



• Molecule 5: CYTOCHROME C OXIDASE

Chain E:



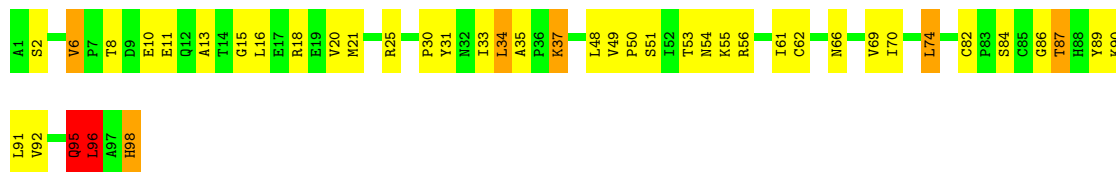
• Molecule 5: CYTOCHROME C OXIDASE

Chain R:



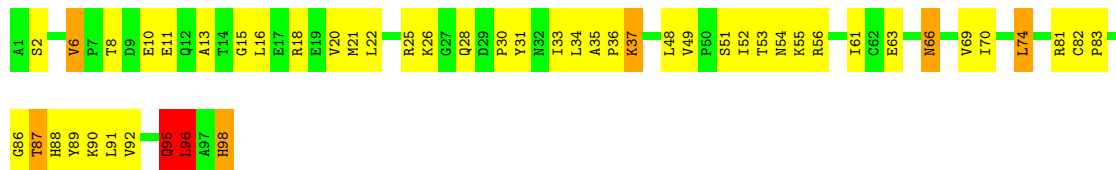
• Molecule 6: CYTOCHROME C OXIDASE

Chain F:



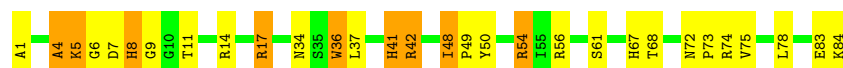
• Molecule 6: CYTOCHROME C OXIDASE

Chain S:



- Molecule 7: CYTOCHROME C OXIDASE

Chain G: 



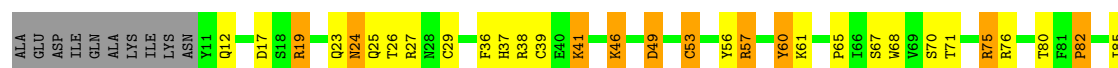
- Molecule 7: CYTOCHROME C OXIDASE

Chain T: 



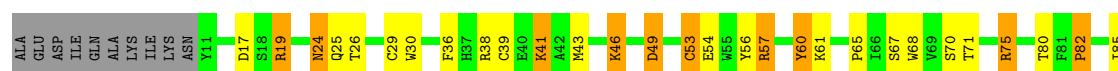
- Molecule 8: CYTOCHROME C OXIDASE

Chain H: 



- Molecule 8: CYTOCHROME C OXIDASE

Chain U: 



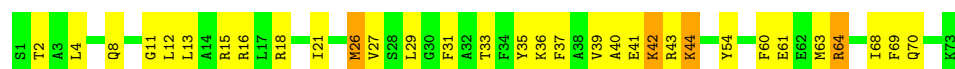
- Molecule 9: CYTOCHROME C OXIDASE

Chain I: 



- Molecule 9: CYTOCHROME C OXIDASE

Chain V: 



- Molecule 10: CYTOCHROME C OXIDASE

Chain J: 



- Molecule 10: CYTOCHROME C OXIDASE

Chain W: 



- Molecule 11: CYTOCHROME C OXIDASE

Chain K:



- Molecule 11: CYTOCHROME C OXIDASE

Chain X:



- Molecule 12: CYTOCHROME C OXIDASE

Chain L:



- Molecule 12: CYTOCHROME C OXIDASE

Chain Y:



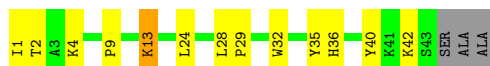
- Molecule 13: CYTOCHROME C OXIDASE

Chain M:



- Molecule 13: CYTOCHROME C OXIDASE

Chain Z:



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	189.10Å 210.50Å 178.60Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	7.00 – 2.80	Depositor
% Data completeness (in resolution range)	97.3 (7.00-2.80)	Depositor
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR 3.84	Depositor
R, R_{free}	0.213 , 0.256	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	28810	wwPDB-VP
Average B, all atoms (Å ²)	31.0	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: CMO, MG, NA, ZN, HEA, CU

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	1/4164 (0.0%)	0.86	2/5688 (0.0%)
1	N	0.69	2/4164 (0.0%)	0.84	1/5688 (0.0%)
2	B	0.70	0/1909	0.91	1/2601 (0.0%)
2	O	0.65	1/1909 (0.1%)	0.88	1/2601 (0.0%)
3	C	0.67	0/2211	0.77	1/3023 (0.0%)
3	P	0.66	0/2211	0.77	1/3023 (0.0%)
4	D	0.65	0/1229	0.72	1/1658 (0.1%)
4	Q	0.61	0/1229	0.70	1/1658 (0.1%)
5	E	0.59	0/898	0.72	0/1218
5	R	0.56	0/898	0.72	0/1218
6	F	0.68	0/765	0.86	0/1038
6	S	0.64	0/765	0.87	0/1038
7	G	0.66	0/699	0.85	1/950 (0.1%)
7	T	0.63	0/699	0.86	1/950 (0.1%)
8	H	0.67	0/648	0.78	0/877
8	U	0.61	0/648	0.77	0/877
9	I	0.67	0/611	0.73	0/810
9	V	0.71	0/611	0.73	0/810
10	J	0.64	0/451	0.76	0/610
10	W	0.66	0/451	0.75	0/610
11	K	0.66	0/398	0.70	0/546
11	X	0.58	0/398	0.67	0/546
12	L	0.70	0/399	0.65	0/534
12	Y	0.66	0/399	0.64	0/534
13	M	0.60	0/345	0.74	0/470
13	Z	0.58	0/345	0.70	0/470
All	All	0.67	4/29454 (0.0%)	0.80	11/40046 (0.0%)

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
1	A	0	2
1	N	0	2
2	B	0	2
2	O	0	1
3	C	0	1
3	P	0	1
5	E	0	1
5	R	0	1
All	All	0	11

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	O	198	GLU	CG-CD	6.00	1.60	1.51
1	N	61	HIS	CG-CD2	5.63	1.45	1.35
1	N	378	HIS	CG-CD2	5.43	1.45	1.35
1	A	61	HIS	CG-CD2	5.18	1.44	1.35

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	133	GLY	N-CA-C	6.20	128.60	113.10
3	C	92	LEU	CA-CB-CG	-6.11	101.25	115.30
1	N	61	HIS	CG-ND1-CE1	-5.94	97.98	105.70
3	P	92	LEU	CA-CB-CG	-5.85	101.85	115.30
4	Q	133	GLY	N-CA-C	5.72	127.41	113.10
2	B	170	LEU	CA-CB-CG	5.72	128.45	115.30
7	G	9	GLY	N-CA-C	-5.53	99.28	113.10
7	T	9	GLY	N-CA-C	-5.46	99.46	113.10
1	A	435	GLY	N-CA-C	5.27	126.28	113.10
1	A	96	ARG	NE-CZ-NH2	-5.21	117.70	120.30
2	O	134	ARG	N-CA-C	5.18	124.98	111.00

There are no chirality outliers.

All (11) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	240	HIS	Sidechain
1	A	304	TYR	Sidechain
2	B	110	TYR	Sidechain

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Mol	Chain	Res	Type	Group
2	B	121	TYR	Sidechain
3	C	8	TYR	Sidechain
5	E	82	TYR	Sidechain
1	N	240	HIS	Sidechain
1	N	304	TYR	Sidechain
2	O	121	TYR	Sidechain
3	P	8	TYR	Sidechain
5	R	82	TYR	Sidechain

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4025	0	4002	128	0
1	N	4025	0	4002	125	0
2	B	1863	0	1867	82	0
2	O	1863	0	1867	81	0
3	C	2124	0	2044	64	0
3	P	2124	0	2044	74	0
4	D	1195	0	1183	44	0
4	Q	1195	0	1183	44	0
5	E	878	0	868	33	0
5	R	878	0	868	30	0
6	F	748	0	728	39	0
6	S	748	0	728	48	0
7	G	672	0	645	34	0
7	T	672	0	645	33	0
8	H	628	0	582	46	0
8	U	628	0	582	45	0
9	I	598	0	612	28	0
9	V	598	0	612	28	0
10	J	441	0	439	13	0
10	W	441	0	439	11	0
11	K	384	0	366	12	0
11	X	384	0	366	15	0
12	L	386	0	388	14	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
12	Y	386	0	388	13	0
13	M	335	0	352	15	0
13	Z	335	0	352	15	0
14	A	1	0	0	0	0
14	B	2	0	0	0	0
14	N	1	0	0	0	0
14	O	2	0	0	0	0
15	A	1	0	0	0	0
15	N	1	0	0	0	0
16	A	1	0	0	0	0
16	N	1	0	0	0	0
17	F	1	0	0	0	0
17	S	1	0	0	0	0
18	A	120	0	107	12	0
18	N	120	0	107	8	0
19	A	2	0	0	0	0
19	N	2	0	0	0	0
All	All	28810	0	28366	920	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 16.

All (920) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
8:H:57:ARG:HH11	8:H:57:ARG:HB3	1.23	1.03
8:U:57:ARG:HB3	8:U:57:ARG:HH11	1.20	1.01
2:B:86:MET:O	2:B:89[B]:GLU:HG2	1.65	0.96
1:A:35:LEU:HD11	1:A:462:LEU:HD13	1.53	0.90
2:O:86:MET:O	2:O:89[B]:GLU:HG2	1.71	0.89
2:O:78:LEU:HB2	2:O:79:PRO:HD3	1.55	0.88
2:B:78:LEU:HB2	2:B:79:PRO:HD3	1.55	0.87
8:H:39:CYS:SG	8:H:53:CYS:CB	2.64	0.86
8:U:39:CYS:SG	8:U:53:CYS:CB	2.65	0.84
8:H:39:CYS:SG	8:H:53:CYS:HB3	2.17	0.84
1:A:282:PHE:HB2	7:T:5:LYS:HB3	1.59	0.83
1:A:365:ILE:HG13	2:B:87:MET:HE1	1.60	0.83
7:G:5:LYS:HB3	1:N:282:PHE:HB2	1.60	0.82
3:P:209:ILE:O	3:P:213:THR:HG23	1.80	0.82
8:U:39:CYS:SG	8:U:53:CYS:HB3	2.20	0.82
1:N:184:PHE:H	1:N:256:HIS:HE1	1.27	0.81
2:O:103:GLN:HG2	2:O:104:TRP:CE2	2.15	0.81
2:B:103:GLN:HG2	2:B:104:TRP:CE2	2.16	0.80

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
7:G:1:ALA:HB1	1:N:286:ILE:HG22	1.63	0.80
1:A:286:ILE:HG22	7:T:1:ALA:HB1	1.64	0.79
7:G:6:GLY:HA3	1:N:190:ILE:HG12	1.63	0.78
1:N:506:GLU:HG3	3:P:1:MET:SD	2.24	0.78
1:N:35:LEU:HD11	1:N:462:LEU:HD13	1.65	0.77
3:C:209:ILE:O	3:C:213:THR:HG23	1.84	0.77
3:P:18:LEU:HD22	3:P:22:LEU:HD22	1.67	0.77
18:A:515:HEA:HBC1	18:A:515:HEA:HMC1	1.65	0.77
2:O:13:THR:HB	2:O:168:LEU:HD23	1.67	0.77
1:A:190:ILE:HG12	7:T:6:GLY:HA3	1.66	0.76
6:F:33:ILE:HG22	6:F:34:LEU:HD12	1.67	0.76
1:A:506:GLU:HG3	3:C:1:MET:SD	2.24	0.76
4:Q:147:LYS:HA	4:Q:147:LYS:HE3	1.68	0.75
2:B:13:THR:HB	2:B:168:LEU:HD23	1.68	0.75
2:B:59:GLN:H	2:B:62:GLU:HG3	1.52	0.74
4:D:147:LYS:HE3	4:D:147:LYS:HA	1.68	0.74
2:B:1:MET:SD	2:B:133:LEU:CD1	2.76	0.74
2:O:152:MET:HB2	2:O:182:THR:HG23	1.70	0.73
6:F:30:PRO:HB3	6:F:96:LEU:HB3	1.70	0.73
2:B:188:ARG:HG3	9:I:54:TYR:OH	1.89	0.73
8:H:39:CYS:HG	8:H:53:CYS:HG	0.78	0.73
10:W:29:ASN:O	10:W:33:ARG:HG3	1.89	0.72
13:M:13:LYS:HD3	13:M:13:LYS:H	1.54	0.72
1:A:184:PHE:H	1:A:256:HIS:HE1	1.38	0.71
1:A:306:THR:HB	1:A:359:ALA:O	1.90	0.71
8:H:53:CYS:HA	8:U:46:LYS:NZ	2.06	0.71
1:N:306:THR:HB	1:N:359:ALA:O	1.90	0.71
3:P:60:ASP:O	3:P:64:GLU:HG3	1.91	0.71
6:S:30:PRO:HB3	6:S:96:LEU:HB3	1.71	0.71
2:B:59:GLN:HA	2:B:62:GLU:HB2	1.72	0.70
2:B:152:MET:HB2	2:B:182:THR:HG23	1.73	0.70
1:A:506:GLU:HB2	3:C:1:MET:SD	2.32	0.70
18:N:515:HEA:HMC1	18:N:515:HEA:HBC1	1.73	0.70
8:H:46:LYS:NZ	8:U:53:CYS:HA	2.07	0.70
1:N:506:GLU:HB2	3:P:1:MET:HA	1.74	0.70
9:I:27:VAL:HG13	9:I:31:PHE:HE2	1.57	0.69
3:C:60:ASP:O	3:C:64:GLU:HG3	1.91	0.69
2:O:188:ARG:HG3	9:V:54:TYR:OH	1.92	0.69
2:O:59:GLN:HA	2:O:62:GLU:HB2	1.74	0.69
8:U:71:THR:O	8:U:75:ARG:HD3	1.92	0.69
4:D:23:PRO:O	5:E:66:ARG:HD3	1.93	0.69
1:A:506:GLU:HB2	3:C:1:MET:HA	1.74	0.69

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:O:186:SER:HB3	2:O:213:LEU:HD22	1.75	0.69
10:J:29:ASN:O	10:J:33:ARG:HG3	1.93	0.68
8:U:39:CYS:HG	8:U:53:CYS:HG	0.69	0.68
8:H:71:THR:O	8:H:75:ARG:HD3	1.92	0.68
2:O:1:MET:SD	2:O:133:LEU:CD1	2.82	0.68
7:G:4:ALA:HB2	1:N:197:LEU:HD12	1.76	0.68
2:O:59:GLN:H	2:O:62:GLU:HG3	1.58	0.67
8:H:41:LYS:HB3	8:H:41:LYS:NZ	2.09	0.67
1:A:197:LEU:HD12	7:T:4:ALA:HB2	1.77	0.66
2:O:132:GLU:HB3	2:O:137:GLU:HG3	1.78	0.66
9:V:27:VAL:HG13	9:V:31:PHE:HE2	1.60	0.66
5:E:82:TYR:HB3	5:E:83:PRO:HD3	1.76	0.66
1:N:184:PHE:H	1:N:256:HIS:CE1	2.11	0.66
6:S:33:ILE:HG22	6:S:34:LEU:HD12	1.76	0.66
2:B:103:GLN:HG2	2:B:104:TRP:CZ2	2.30	0.66
2:O:5:MET:HE3	11:X:43:SER:H	1.59	0.66
8:H:39:CYS:HG	8:H:53:CYS:CB	2.08	0.66
7:G:54:ARG:N	7:G:54:ARG:HD3	2.10	0.66
3:C:19:THR:HG23	3:C:53:THR:OG1	1.96	0.66
2:O:1:MET:SD	2:O:133:LEU:HD13	2.36	0.65
8:U:41:LYS:NZ	8:U:41:LYS:HB3	2.12	0.65
3:C:164:PHE:O	3:C:168:THR:HG23	1.97	0.65
7:G:4:ALA:CB	1:N:197:LEU:HD12	2.26	0.65
4:D:24:LEU:HB3	5:E:30:ARG:HG2	1.78	0.65
4:Q:24:LEU:HB3	5:R:30:ARG:HG2	1.79	0.65
4:D:40:LEU:HD22	4:D:59:LEU:HD13	1.77	0.65
13:Z:13:LYS:H	13:Z:13:LYS:HD3	1.61	0.64
4:Q:68:PHE:HA	4:Q:71:MET:HG2	1.78	0.64
8:H:17:ASP:OD1	8:H:19:ARG:HG2	1.98	0.64
3:P:164:PHE:O	3:P:168:THR:HG23	1.97	0.64
1:A:184:PHE:H	1:A:256:HIS:CE1	2.15	0.64
8:U:65:PRO:HG2	8:U:68:TRP:CG	2.33	0.64
7:G:4:ALA:HB2	1:N:197:LEU:CD1	2.28	0.64
2:O:103:GLN:HG2	2:O:104:TRP:CZ2	2.33	0.64
13:M:32:TRP:O	13:M:36:HIS:HD2	1.80	0.64
13:Z:32:TRP:O	13:Z:36:HIS:HD2	1.79	0.64
3:C:18:LEU:HD22	3:C:22:LEU:HD22	1.79	0.64
2:O:90[A]:ILE:H	2:O:90[A]:ILE:HD12	1.63	0.64
2:B:1:MET:SD	2:B:133:LEU:HD11	2.38	0.64
1:A:510:TYR:HB2	6:F:56:ARG:NH1	2.12	0.64
1:N:390:MET:CE	18:N:515:HEA:H242	2.27	0.64
1:A:197:LEU:HD12	7:T:4:ALA:CB	2.28	0.63

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
6:S:8:THR:OG1	6:S:11:GLU:HG2	1.98	0.63
2:B:90[A]:ILE:HD12	2:B:90[A]:ILE:H	1.63	0.63
1:A:197:LEU:CD1	7:T:4:ALA:HB2	2.29	0.63
7:T:67:HIS:CD2	7:T:78:LEU:HD11	2.34	0.63
4:Q:40:LEU:HD22	4:Q:59:LEU:HD13	1.80	0.62
2:B:22:HIS:CE1	9:I:44:LYS:HD3	2.34	0.62
7:G:67:HIS:CD2	7:G:78:LEU:HD11	2.34	0.62
2:B:186:SER:HB3	2:B:213:LEU:HD22	1.81	0.62
2:B:50:LEU:HD21	5:E:77:PRO:HD2	1.82	0.62
10:J:50:LEU:HD22	10:J:50:LEU:O	1.99	0.62
4:Q:7:LYS:O	4:Q:10:ASP:HB2	2.00	0.62
6:S:10:GLU:HG2	6:S:25:ARG:HH22	1.63	0.62
6:F:13:ALA:O	6:F:18:ARG:HD2	1.99	0.62
6:F:10:GLU:HG2	6:F:25:ARG:HH22	1.64	0.62
2:O:136:LEU:HB3	2:O:193:TYR:CD2	2.35	0.62
8:U:57:ARG:HH11	8:U:57:ARG:CB	2.05	0.62
6:S:53:THR:HG22	6:S:54:ASN:H	1.64	0.62
3:C:154:GLY:HA2	6:F:6:VAL:HG22	1.82	0.62
1:N:390:MET:HE2	18:N:515:HEA:H242	1.80	0.61
2:B:132:GLU:HB3	2:B:137:GLU:HG3	1.82	0.61
6:F:8:THR:OG1	6:F:11:GLU:HG2	2.00	0.61
10:J:12:PHE:O	10:J:23:LYS:HE2	2.01	0.61
2:B:5:MET:HE3	11:K:43:SER:H	1.65	0.61
7:T:54:ARG:HD3	7:T:54:ARG:N	2.15	0.61
3:P:19:THR:HG23	3:P:53:THR:OG1	2.00	0.61
6:F:53:THR:HG22	6:F:54:ASN:H	1.64	0.61
4:Q:23:PRO:O	5:R:66:ARG:HD3	2.00	0.61
1:N:219:PHE:HZ	3:P:199:VAL:HG21	1.63	0.61
9:I:39:VAL:O	9:I:42:LYS:HE2	2.00	0.61
1:A:333:LYS:HE2	1:A:335:SER:HB3	1.81	0.61
1:N:510:TYR:HB2	6:S:56:ARG:NH1	2.16	0.61
7:G:42:ARG:CZ	7:G:74:ARG:HH21	2.13	0.61
10:W:12:PHE:O	10:W:23:LYS:HE2	2.01	0.61
1:N:128:VAL:O	1:N:128:VAL:HG22	2.00	0.61
3:C:222:GLN:HE21	3:C:222:GLN:HA	1.65	0.60
1:N:333:LYS:HE2	1:N:335:SER:HB3	1.84	0.60
2:O:90[B]:ILE:HG23	2:O:91[B]:ASN:H	1.65	0.60
5:R:82:TYR:HB3	5:R:83:PRO:HD3	1.82	0.60
7:T:42:ARG:CZ	7:T:74:ARG:HH21	2.15	0.60
10:W:27:THR:O	10:W:30:ILE:HG12	2.01	0.60
8:U:38:ARG:HG2	8:U:85:ILE:HG23	1.83	0.60
6:S:13:ALA:O	6:S:18:ARG:HD2	2.02	0.60

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:O:179:LEU:HD21	8:U:65:PRO:HD3	1.83	0.59
1:N:219:PHE:CZ	3:P:199:VAL:HG21	2.37	0.59
3:P:222:GLN:HA	3:P:222:GLN:HE21	1.66	0.59
8:U:17:ASP:OD1	8:U:19:ARG:HG2	2.02	0.59
2:B:90[B]:ILE:HG23	2:B:91[B]:ASN:H	1.68	0.59
4:D:68:PHE:HA	4:D:71:MET:HG2	1.84	0.59
1:N:225:GLY:HA3	3:P:112:LEU:CD1	2.33	0.59
2:O:226:MET:O	2:O:227:LEU:HB2	2.02	0.59
8:H:65:PRO:HG2	8:H:68:TRP:CG	2.38	0.59
1:N:506:GLU:HB2	3:P:1:MET:SD	2.42	0.59
8:H:57:ARG:NH1	8:H:57:ARG:HB3	2.08	0.58
1:A:225:GLY:HA3	3:C:112:LEU:CD1	2.33	0.58
3:C:148:HIS:O	3:C:152:MET:HG3	2.02	0.58
4:Q:130:PRO:HG2	4:Q:131:ILE:HG13	1.85	0.58
2:B:179:LEU:HD21	8:H:65:PRO:HD3	1.84	0.58
2:B:48:THR:HB	9:I:16:ARG:CZ	2.33	0.58
1:A:128:VAL:O	1:A:128:VAL:HG22	2.03	0.58
10:W:50:LEU:HD22	10:W:50:LEU:O	2.02	0.58
2:O:76:ILE:O	2:O:79:PRO:HD2	2.04	0.58
2:B:226:MET:O	2:B:227:LEU:HB2	2.03	0.58
2:B:83:ILE:O	2:B:87:MET:HG3	2.02	0.58
1:A:82:LEU:O	1:A:86:MET:HG3	2.03	0.58
4:D:64:PHE:CE1	5:E:66:ARG:HD2	2.39	0.58
4:D:130:PRO:HG2	4:D:131:ILE:HG13	1.85	0.58
2:O:22:HIS:CE1	9:V:44:LYS:HD3	2.38	0.58
2:B:152:MET:HB2	2:B:182:THR:CG2	2.33	0.58
4:D:70:GLU:HA	5:E:109:VAL:HG12	1.84	0.58
1:N:87:ILE:O	1:N:173:PRO:HD3	2.04	0.58
9:V:39:VAL:O	9:V:42:LYS:HE2	2.03	0.58
9:V:29:LEU:O	9:V:33:THR:HG23	2.03	0.58
8:U:57:ARG:HA	8:U:60:TYR:CD2	2.39	0.58
1:A:87:ILE:O	1:A:173:PRO:HD3	2.03	0.58
1:N:219:PHE:HZ	3:P:199:VAL:CG2	2.17	0.58
2:O:146:MET:SD	2:O:189:PRO:HB3	2.43	0.58
9:I:15:ARG:HD2	9:I:18:ARG:NH2	2.19	0.58
8:U:57:ARG:O	8:U:61:LYS:HB2	2.04	0.57
5:R:7:THR:OG1	5:R:10:GLU:HG3	2.04	0.57
5:R:6:GLU:HB2	5:R:10:GLU:HB2	1.86	0.57
3:P:154:GLY:HA2	6:S:6:VAL:HG22	1.86	0.57
2:B:1:MET:SD	2:B:133:LEU:HD13	2.44	0.57
6:F:53:THR:HG22	6:F:54:ASN:OD1	2.05	0.57
4:Q:52:SER:OG	4:Q:55:GLU:HG3	2.04	0.57

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:O:125:THR:HA	2:O:128:LEU:HG	1.86	0.57
2:B:63:THR:O	2:B:66:THR:HG22	2.04	0.57
1:A:75:ILE:O	1:A:79:GLY:HA3	2.05	0.57
8:H:57:ARG:HA	8:H:60:TYR:CD2	2.39	0.57
1:N:29:VAL:HG13	12:Y:36:PRO:HG3	1.87	0.57
8:U:39:CYS:CB	8:U:53:CYS:SG	2.93	0.57
9:I:15:ARG:HD2	9:I:18:ARG:HH22	1.69	0.57
4:Q:70:GLU:HA	5:R:109:VAL:HG12	1.85	0.57
1:N:492:LEU:HD23	1:N:495:LEU:HD12	1.86	0.57
1:A:398:PRO:O	1:A:498:CYS:HB3	2.05	0.57
1:A:506:GLU:CG	3:C:1:MET:SD	2.93	0.57
10:W:29:ASN:ND2	10:W:33:ARG:HD2	2.19	0.57
10:W:36:MET:O	10:W:40:LEU:HG	2.05	0.57
2:O:83:ILE:O	2:O:87:MET:HG3	2.04	0.57
9:I:29:LEU:O	9:I:33:THR:HG23	2.04	0.57
2:O:48:THR:HB	9:V:16:ARG:CZ	2.35	0.57
1:N:32:ALA:HB3	12:Y:36:PRO:HG2	1.86	0.57
2:B:102:HIS:O	2:B:104:TRP:N	2.38	0.56
4:Q:137:LYS:O	4:Q:145:TRP:HE3	1.88	0.56
1:A:240:HIS:O	1:A:243:VAL:HG22	2.05	0.56
12:L:41:ARG:HD2	13:M:40:TYR:CZ	2.39	0.56
12:Y:41:ARG:HD2	13:Z:40:TYR:CZ	2.40	0.56
2:B:125:THR:HA	2:B:128:LEU:HG	1.87	0.56
2:O:3:TYR:N	2:O:3:TYR:CD1	2.73	0.56
1:A:506:GLU:CB	3:C:1:MET:SD	2.93	0.56
1:A:53:ILE:HD11	12:L:43:GLN:HB3	1.86	0.56
1:N:53:ILE:HD11	12:Y:43:GLN:HB3	1.88	0.56
2:B:86:MET:O	2:B:89[B]:GLU:CG	2.49	0.56
12:L:41:ARG:HG3	13:M:40:TYR:CE2	2.41	0.56
1:N:398:PRO:O	1:N:498:CYS:HB3	2.04	0.56
3:P:154:GLY:HA2	6:S:6:VAL:CG2	2.36	0.56
4:D:137:LYS:O	4:D:145:TRP:HE3	1.89	0.56
8:H:39:CYS:CB	8:H:53:CYS:SG	2.93	0.56
9:V:15:ARG:HD2	9:V:18:ARG:HH22	1.71	0.56
4:D:48:TRP:CH2	5:E:56:ARG:HA	2.41	0.56
4:Q:8:SER:OG	13:Z:4:LYS:HE3	2.06	0.56
1:A:35:LEU:CD1	1:A:462:LEU:HD13	2.33	0.55
1:N:506:GLU:CG	3:P:1:MET:SD	2.93	0.55
1:N:35:LEU:CD1	1:N:462:LEU:HD13	2.35	0.55
4:Q:64:PHE:CE1	5:R:66:ARG:HD2	2.40	0.55
4:D:7:LYS:O	4:D:10:ASP:HB2	2.07	0.55
3:P:148:HIS:O	3:P:152:MET:HG3	2.05	0.55

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:N:290:HIS:CD2	1:N:291:HIS:CD2	2.94	0.55
8:H:49:ASP:HB2	8:U:49:ASP:HB2	1.88	0.55
1:N:115:SER:HA	1:N:145:LEU:HD12	1.89	0.55
2:O:63:THR:O	2:O:66:THR:HG22	2.07	0.55
7:T:42:ARG:HG3	7:T:42:ARG:HH11	1.72	0.55
6:F:16:LEU:O	6:F:20:VAL:HG13	2.06	0.55
8:U:57:ARG:HB3	8:U:57:ARG:NH1	2.05	0.55
2:O:152:MET:HB2	2:O:182:THR:CG2	2.37	0.55
3:C:121:ILE:HG22	3:C:123:PRO:HD3	1.89	0.55
13:Z:28:LEU:HB2	13:Z:29:PRO:HD3	1.89	0.55
2:B:78:LEU:HB2	2:B:79:PRO:CD	2.34	0.55
12:Y:41:ARG:HG3	13:Z:40:TYR:CE2	2.42	0.55
2:O:142:VAL:HG13	2:O:210:VAL:O	2.06	0.55
3:P:154:GLY:C	6:S:6:VAL:HG13	2.27	0.55
1:N:230:LEU:HB2	3:P:103:HIS:CD2	2.42	0.55
2:B:136:LEU:HB3	2:B:193:TYR:CD2	2.42	0.55
3:C:156:ARG:HH11	3:C:156:ARG:HG3	1.72	0.55
9:V:15:ARG:HD2	9:V:18:ARG:NH2	2.21	0.54
11:X:24:PHE:O	11:X:28:VAL:HG23	2.07	0.54
1:N:328:HIS:HB2	2:O:45:MET:HE3	1.88	0.54
7:G:42:ARG:NH1	7:G:74:ARG:HH21	2.04	0.54
3:P:137:LEU:HD13	3:P:173:PHE:CD2	2.43	0.54
8:H:53:CYS:HA	8:U:46:LYS:HZ3	1.72	0.54
1:A:390:MET:CE	18:A:515:HEA:H242	2.38	0.54
1:N:33:LEU:O	1:N:37:ILE:HG13	2.07	0.54
2:O:132:GLU:HA	4:Q:122:ARG:NH1	2.22	0.54
7:G:41:HIS:CD2	7:G:42:ARG:HH12	2.26	0.54
3:P:154:GLY:HA2	6:S:6:VAL:HG13	1.89	0.54
4:D:52:SER:OG	4:D:55:GLU:HG3	2.06	0.54
2:B:28:LEU:HD23	9:I:35:TYR:CZ	2.43	0.54
2:O:184:LEU:HD11	2:O:211:LEU:HD21	1.89	0.54
2:O:164:ALA:HB2	2:O:171:LYS:HD3	1.89	0.54
6:F:62:CYS:SG	6:F:84:SER:OG	2.65	0.54
1:A:115:SER:HA	1:A:145:LEU:HD12	1.89	0.54
5:E:6:GLU:HB2	5:E:10:GLU:HB2	1.88	0.54
1:A:219:PHE:HZ	3:C:199:VAL:HG21	1.73	0.54
2:B:3:TYR:CD1	2:B:3:TYR:N	2.75	0.54
8:H:57:ARG:CB	8:H:57:ARG:HH11	2.08	0.54
4:Q:127:LYS:HD3	9:V:63:MET:HE2	1.90	0.54
2:O:66:THR:HG23	2:O:67:ILE:HD13	1.90	0.54
2:O:139:ASP:OD1	2:O:140:ASN:N	2.40	0.54
1:A:225:GLY:HA3	3:C:112:LEU:HD11	1.90	0.54

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
9:V:63:MET:HB3	9:V:68:ILE:HD11	1.89	0.54
3:P:81:TYR:O	3:P:85:LEU:HD22	2.08	0.54
7:T:49:PRO:HG2	8:U:80:THR:HG21	1.90	0.54
3:P:154:GLY:HA2	6:S:6:VAL:CG1	2.37	0.54
2:B:23:PHE:CZ	2:B:80:SER:HB2	2.43	0.54
10:J:36:MET:O	10:J:40:LEU:HG	2.07	0.54
1:N:225:GLY:HA3	3:P:112:LEU:HD11	1.89	0.54
8:U:65:PRO:HG2	8:U:68:TRP:CD1	2.43	0.53
6:S:87:THR:HG22	6:S:89:TYR:CE2	2.43	0.53
1:N:75:ILE:O	1:N:79:GLY:HA3	2.08	0.53
3:P:16:TRP:CE3	3:P:19:THR:HG21	2.43	0.53
13:Z:1:ILE:O	13:Z:1:ILE:HG23	2.08	0.53
3:P:4:GLN:NE2	3:P:6:HIS:O	2.41	0.53
7:T:42:ARG:NH1	7:T:74:ARG:HH21	2.06	0.53
1:A:219:PHE:HZ	3:C:199:VAL:CG2	2.22	0.53
1:A:32:ALA:HB3	12:L:36:PRO:HG2	1.91	0.53
10:J:27:THR:O	10:J:30:ILE:HG12	2.09	0.53
5:E:12:ASP:O	5:E:16:VAL:HG23	2.08	0.53
4:Q:33:LEU:HD13	4:Q:41:LYS:HG3	1.90	0.53
9:I:27:VAL:HG13	9:I:31:PHE:CE2	2.41	0.53
8:H:38:ARG:HG2	8:H:85:ILE:HG23	1.90	0.53
1:N:483:LEU:HB2	13:Z:2:THR:OG1	2.08	0.53
6:S:53:THR:HG22	6:S:54:ASN:OD1	2.08	0.53
2:B:146:MET:SD	2:B:189:PRO:HB3	2.49	0.53
6:S:49:VAL:O	6:S:91:LEU:HD12	2.09	0.53
4:D:40:LEU:CD2	4:D:59:LEU:HD13	2.39	0.53
5:E:7:THR:OG1	5:E:10:GLU:HG3	2.08	0.53
7:T:36:TRP:O	7:T:36:TRP:HD1	1.92	0.53
1:N:484:THR:HB	13:Z:2:THR:CG2	2.39	0.53
2:B:164:ALA:HB2	2:B:171:LYS:HD3	1.91	0.53
10:J:29:ASN:ND2	10:J:33:ARG:HD2	2.24	0.52
1:A:176:MET:CE	1:A:181:THR:HG22	2.39	0.52
8:U:24:ASN:ND2	8:U:26:THR:H	2.07	0.52
1:N:107:PRO:HB3	3:P:25:LEU:HB2	1.91	0.52
11:X:47:ARG:HG3	11:X:48:VAL:HG23	1.92	0.52
1:N:367:LEU:O	1:N:370:THR:HG23	2.09	0.52
5:R:95:GLU:HG2	5:R:96:LEU:HD23	1.92	0.52
3:C:154:GLY:HA2	6:F:6:VAL:CG2	2.39	0.52
1:N:240:HIS:O	1:N:243:VAL:HG22	2.09	0.52
1:A:513:LEU:HD22	6:F:35:ALA:HB3	1.91	0.52
6:F:49:VAL:O	6:F:91:LEU:HD12	2.09	0.52
1:N:402:GLY:HA3	1:N:499:PRO:HD3	1.91	0.52

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:N:437:PRO:HG2	1:N:440:TYR:CE2	2.45	0.52
1:A:367:LEU:O	1:A:370:THR:HG23	2.10	0.52
2:B:108:TYR:CD1	2:B:108:TYR:N	2.77	0.52
6:F:51:SER:HB2	6:F:91:LEU:HD11	1.90	0.52
11:K:13:TYR:O	11:K:17:VAL:HG23	2.09	0.52
2:B:184:LEU:HD11	2:B:211:LEU:HD21	1.92	0.52
2:O:102:HIS:O	2:O:104:TRP:N	2.43	0.52
10:W:30:ILE:O	10:W:34:VAL:HG23	2.10	0.52
4:Q:48:TRP:CH2	5:R:56:ARG:HA	2.45	0.52
1:N:187:SER:HB2	1:N:277:MET:CE	2.40	0.52
7:G:36:TRP:O	7:G:36:TRP:HD1	1.93	0.52
1:A:219:PHE:CZ	3:C:199:VAL:HG21	2.44	0.52
13:M:1:ILE:O	13:M:1:ILE:HG23	2.09	0.52
1:A:117:MET:HE3	12:L:39:ILE:HG23	1.91	0.52
7:G:49:PRO:HG2	8:H:80:THR:HG21	1.91	0.52
8:H:65:PRO:HG2	8:H:68:TRP:CD1	2.45	0.52
2:O:68:LEU:HB3	2:O:69:PRO:HD3	1.92	0.52
2:B:216:LEU:O	2:B:219:PHE:HB3	2.10	0.52
6:F:33:ILE:HG22	6:F:34:LEU:CD1	2.37	0.51
10:W:27:THR:HG22	10:W:30:ILE:HD11	1.91	0.51
8:H:57:ARG:O	8:H:61:LYS:HB2	2.09	0.51
2:O:13:THR:HG22	2:O:13:THR:O	2.10	0.51
3:P:16:TRP:HE3	3:P:19:THR:HG21	1.73	0.51
7:T:41:HIS:CD2	7:T:42:ARG:HH12	2.28	0.51
7:T:42:ARG:HG3	7:T:42:ARG:NH1	2.25	0.51
2:B:68:LEU:HB3	2:B:69:PRO:HD3	1.91	0.51
4:D:73:ARG:CZ	4:D:73:ARG:HB3	2.40	0.51
1:A:333:LYS:CE	1:A:335:SER:HB3	2.40	0.51
1:N:176:MET:CE	1:N:181:THR:HG22	2.40	0.51
1:A:390:MET:HE2	18:A:515:HEA:H242	1.93	0.51
13:M:35:TYR:HD2	13:M:36:HIS:CD2	2.28	0.51
7:G:48:ILE:HD13	7:G:50:TYR:CZ	2.46	0.51
2:O:50:LEU:HD21	5:R:77:PRO:HD2	1.93	0.51
4:Q:41:LYS:HD3	4:Q:62:LEU:HD23	1.93	0.51
5:R:12:ASP:O	5:R:16:VAL:HG23	2.11	0.51
2:O:28:LEU:HD23	9:V:35:TYR:CZ	2.46	0.51
1:A:107:PRO:HB3	3:C:25:LEU:HB2	1.91	0.51
2:B:5:MET:CE	11:K:42:PRO:HA	2.40	0.51
2:B:139:ASP:OD1	2:B:140:ASN:N	2.43	0.51
4:Q:114:GLU:HG3	11:X:51:LYS:HE2	1.93	0.51
5:R:80:GLU:CD	5:R:80:GLU:H	2.13	0.51
1:A:483:LEU:HB2	13:M:2:THR:OG1	2.11	0.51

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:40:GLU:HG2	1:A:54:TYR:CD1	2.46	0.51
2:B:132:GLU:HA	4:D:122:ARG:NH1	2.24	0.50
2:O:108:TYR:N	2:O:108:TYR:CD1	2.79	0.50
1:N:506:GLU:OE2	3:P:1:MET:HG3	2.11	0.50
8:H:41:LYS:HB3	8:H:41:LYS:HZ2	1.74	0.50
9:I:40:ALA:O	9:I:44:LYS:HE2	2.11	0.50
7:G:42:ARG:HG3	7:G:42:ARG:HH11	1.75	0.50
5:R:20:ASN:HA	5:R:57:ARG:HH22	1.76	0.50
4:Q:16:TYR:CE1	4:Q:25:PRO:HG3	2.46	0.50
6:F:87:THR:HG22	6:F:89:TYR:CE2	2.45	0.50
10:J:30:ILE:O	10:J:34:VAL:HG23	2.11	0.50
1:A:187:SER:HB2	1:A:277:MET:HE1	1.92	0.50
2:O:5:MET:HE3	11:X:43:SER:N	2.27	0.50
1:N:290:HIS:HA	1:N:293:PHE:CZ	2.46	0.50
6:S:16:LEU:O	6:S:20:VAL:HG13	2.12	0.50
2:O:78:LEU:HB2	2:O:79:PRO:CD	2.35	0.50
4:D:40:LEU:HD22	4:D:59:LEU:CD1	2.41	0.50
5:E:21:LYS:O	5:E:57:ARG:NH1	2.45	0.50
5:E:20:ASN:HA	5:E:57:ARG:HH22	1.76	0.50
4:Q:128:VAL:O	4:Q:134:PHE:HB3	2.11	0.50
1:A:230:LEU:HB2	3:C:103:HIS:CD2	2.46	0.50
4:Q:37:GLN:O	4:Q:41:LYS:HG2	2.12	0.50
8:U:17:ASP:OD2	8:U:19:ARG:NH1	2.44	0.50
3:P:156:ARG:HG3	3:P:156:ARG:HH11	1.77	0.50
2:B:5:MET:HE3	11:K:43:SER:N	2.26	0.50
1:A:248:LEU:O	1:A:251:PHE:HB2	2.12	0.50
5:R:43:PRO:HB2	5:R:48:ILE:HD11	1.93	0.50
12:Y:35:ALA:HB3	12:Y:36:PRO:HD3	1.94	0.50
1:N:176:MET:HE3	1:N:181:THR:HG22	1.93	0.50
2:O:162:SER:HA	2:O:173:ASP:HA	1.93	0.50
2:B:66:THR:HG23	2:B:67:ILE:HD13	1.93	0.49
2:B:189:PRO:HG2	9:I:63:MET:HE3	1.94	0.49
3:C:154:GLY:C	6:F:6:VAL:HG13	2.33	0.49
1:A:268:PHE:CZ	2:B:58:ALA:HA	2.47	0.49
1:N:268:PHE:CZ	2:O:58:ALA:HA	2.47	0.49
6:F:95:GLN:NE2	6:F:96:LEU:H	2.09	0.49
8:U:41:LYS:HB3	8:U:41:LYS:HZ2	1.77	0.49
3:C:160:LEU:HD13	3:C:222:GLN:HG2	1.94	0.49
1:N:85:LEU:O	1:N:492:LEU:HD13	2.12	0.49
8:U:24:ASN:HD22	8:U:25:GLN:N	2.10	0.49
2:O:188:ARG:HH11	2:O:188:ARG:HG2	1.77	0.49
1:N:40:GLU:HG2	1:N:54:TYR:CD1	2.47	0.49

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
6:S:37:LYS:N	6:S:37:LYS:HD3	2.28	0.49
1:A:484:THR:HB	13:M:2:THR:CG2	2.41	0.49
2:O:216:LEU:O	2:O:219:PHE:HB3	2.12	0.49
1:A:10:THR:HG21	3:C:11:VAL:HG22	1.95	0.49
8:H:37:HIS:CD2	8:H:76:ARG:NH1	2.80	0.49
7:G:42:ARG:HG3	7:G:42:ARG:NH1	2.28	0.49
2:B:125:THR:HB	2:B:134:ARG:HG3	1.95	0.49
4:D:107:ILE:HB	4:D:108:PRO:HD2	1.95	0.49
5:E:78:HIS:CE1	9:I:12:LEU:HD22	2.48	0.49
1:N:358:LEU:HD21	1:N:372:TYR:HD2	1.77	0.49
8:H:46:LYS:HZ3	8:U:53:CYS:HA	1.77	0.49
2:O:5:MET:CE	11:X:42:PRO:HA	2.42	0.49
2:O:125:THR:HB	2:O:134:ARG:HG3	1.95	0.49
8:H:17:ASP:OD2	8:H:19:ARG:NH1	2.45	0.49
6:S:18:ARG:NH1	6:S:21:MET:HE1	2.28	0.49
1:N:513:LEU:HD22	6:S:35:ALA:HB3	1.95	0.49
1:N:74:MET:HE1	1:N:246:LEU:O	2.12	0.49
1:N:506:GLU:CB	3:P:1:MET:SD	3.00	0.49
2:O:189:PRO:HG2	9:V:63:MET:HE3	1.94	0.49
10:J:27:THR:HG22	10:J:30:ILE:HD11	1.94	0.49
1:N:71:MET:CE	1:N:195:LEU:HD21	2.43	0.49
5:E:80:GLU:H	5:E:80:GLU:CD	2.17	0.49
11:X:13:TYR:O	11:X:17:VAL:HG23	2.13	0.49
4:D:37:GLN:O	4:D:41:LYS:HG2	2.12	0.49
4:D:33:LEU:HD13	4:D:41:LYS:HG3	1.94	0.48
6:S:48:LEU:HB3	6:S:92:VAL:HG21	1.94	0.48
2:O:23:PHE:CZ	2:O:80:SER:HB2	2.48	0.48
9:I:36:LYS:O	9:I:40:ALA:HB3	2.12	0.48
7:T:42:ARG:NH2	7:T:75:VAL:HG11	2.28	0.48
2:O:116:LEU:HD11	2:O:226:MET:HG3	1.94	0.48
3:C:228:THR:O	3:C:230:ASN:N	2.46	0.48
6:S:81:ARG:HG2	6:S:88:HIS:CD2	2.48	0.48
1:A:358:LEU:HD21	1:A:372:TYR:HD2	1.77	0.48
3:P:160:LEU:HD13	3:P:222:GLN:HG2	1.94	0.48
3:C:66:THR:HG22	3:C:67:PHE:CD1	2.49	0.48
1:A:379:TYR:O	1:A:383:MET:HB2	2.13	0.48
6:F:30:PRO:CB	6:F:96:LEU:HB3	2.42	0.48
9:V:27:VAL:HG13	9:V:31:PHE:CE2	2.44	0.48
7:G:42:ARG:NH2	7:G:75:VAL:HG11	2.29	0.48
5:R:87:GLN:OE1	9:V:4:LEU:HD23	2.13	0.48
4:D:114:GLU:HG3	11:K:51:LYS:HE2	1.96	0.48
1:N:333:LYS:CE	1:N:335:SER:HB3	2.43	0.48

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
4:Q:67:SER:OG	4:Q:70:GLU:HG3	2.13	0.48
6:F:55:LYS:HA	6:F:74:LEU:O	2.13	0.48
3:C:65:SER:HB3	3:C:71:HIS:CE1	2.48	0.48
3:P:128:GLU:HB3	3:P:129:VAL:H	1.56	0.48
8:U:38:ARG:HG2	8:U:85:ILE:HA	1.95	0.48
11:K:47:ARG:HG3	11:K:48:VAL:HG23	1.95	0.48
3:C:147:ALA:HB2	3:C:162:ALA:HB3	1.95	0.48
2:B:142:VAL:HG13	2:B:210:VAL:O	2.13	0.48
2:B:124:PRO:HB2	2:B:127:GLU:HG2	1.96	0.48
6:S:33:ILE:HG22	6:S:34:LEU:CD1	2.43	0.48
3:P:16:TRP:HA	3:P:19:THR:HG22	1.96	0.48
4:Q:132:GLN:HE22	9:V:43:ARG:HB2	1.78	0.48
1:N:117:MET:HA	1:N:117:MET:HE2	1.95	0.48
1:A:69:MET:O	1:A:72:PRO:HD2	2.13	0.48
13:Z:35:TYR:HD2	13:Z:36:HIS:CD2	2.31	0.48
3:C:57:TRP:O	3:C:61:VAL:HG23	2.14	0.48
3:C:77:LYS:HE2	3:C:81:TYR:OH	2.14	0.48
6:F:61:ILE:HG12	6:F:69:VAL:HG22	1.95	0.48
7:T:47:PHE:CD2	7:T:77:PRO:HB2	2.49	0.48
6:S:61:ILE:HG12	6:S:69:VAL:HG22	1.94	0.48
5:R:81:ILE:O	5:R:85:VAL:HG23	2.14	0.48
7:G:5:LYS:O	1:N:278:MET:HG2	2.13	0.48
5:R:80:GLU:O	5:R:83:PRO:HD2	2.14	0.48
1:A:513:LEU:HD12	1:A:513:LEU:HA	1.61	0.48
9:V:26:MET:HA	9:V:26:MET:CE	2.43	0.48
1:N:328:HIS:HB2	2:O:45:MET:CE	2.44	0.47
8:H:38:ARG:HG2	8:H:85:ILE:HA	1.96	0.47
6:S:31:TYR:HE2	6:S:98:HIS:CE1	2.32	0.47
2:B:59:GLN:N	2:B:62:GLU:HG3	2.26	0.47
2:B:188:ARG:HH11	2:B:188:ARG:HG2	1.79	0.47
3:P:195:SER:O	3:P:199:VAL:HG22	2.14	0.47
3:P:77:LYS:HE2	3:P:81:TYR:OH	2.14	0.47
1:N:337:ALA:HB2	1:N:394:VAL:HG23	1.96	0.47
1:A:160:GLY:O	1:A:164:PHE:HD2	1.97	0.47
1:N:248:LEU:O	1:N:251:PHE:HB2	2.14	0.47
5:E:87:GLN:OE1	9:I:4:LEU:HD23	2.14	0.47
1:A:290:HIS:HA	1:A:293:PHE:CZ	2.50	0.47
1:N:365:ILE:HD13	1:N:365:ILE:HA	1.66	0.47
2:O:122:MET:HB2	2:O:208:PRO:HD2	1.94	0.47
2:O:86:MET:O	2:O:89[B]:GLU:CG	2.55	0.47
1:A:365:ILE:HA	1:A:365:ILE:HD13	1.71	0.47
6:S:82:CYS:N	6:S:86:GLY:O	2.48	0.47

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
3:P:119:THR:HG21	8:U:82:PRO:O	2.15	0.47
3:P:183:GLU:O	7:T:42:ARG:NH2	2.47	0.47
4:Q:18:ASP:OD2	4:Q:66:GLU:HG2	2.13	0.47
2:B:162:SER:HA	2:B:173:ASP:HA	1.97	0.47
3:C:137:LEU:HD13	3:C:173:PHE:CD2	2.50	0.47
5:E:95:GLU:HG2	5:E:96:LEU:HD23	1.95	0.47
3:P:147:ALA:HB2	3:P:162:ALA:HB3	1.96	0.47
4:Q:73:ARG:HB3	4:Q:73:ARG:CZ	2.43	0.47
1:A:218:THR:HB	1:A:221:ASP:HB3	1.95	0.47
3:C:4:GLN:NE2	3:C:6:HIS:O	2.48	0.47
10:W:29:ASN:HD21	10:W:33:ARG:HD2	1.79	0.47
12:L:35:ALA:HB3	12:L:36:PRO:HD3	1.95	0.47
2:O:89[B]:GLU:HB2	2:O:91[B]:ASN:ND2	2.30	0.47
1:A:62:ALA:HB2	18:A:515:HEA:HBD1	1.96	0.47
6:S:95:GLN:NE2	6:S:96:LEU:H	2.12	0.47
4:Q:145:TRP:CE2	11:X:45:VAL:HA	2.49	0.47
4:D:24:LEU:CB	5:E:30:ARG:HG2	2.45	0.47
3:P:192:VAL:O	3:P:196:THR:HB	2.14	0.47
3:P:154:GLY:CA	6:S:6:VAL:HG13	2.45	0.47
1:A:209:LEU:O	1:A:213:ARG:HG3	2.15	0.47
4:D:93:ALA:O	4:D:97:ILE:HG13	2.15	0.47
8:H:36:PHE:HB2	8:H:56:TYR:CB	2.45	0.47
4:Q:24:LEU:CB	5:R:30:ARG:HG2	2.43	0.47
1:A:492:LEU:O	1:A:492:LEU:HD22	2.14	0.47
3:P:187:THR:HG21	7:T:68:THR:HG21	1.96	0.47
8:H:53:CYS:CA	8:U:46:LYS:HZ3	2.28	0.47
2:B:129:LYS:O	2:B:132:GLU:HG3	2.15	0.47
8:H:49:ASP:CB	8:U:49:ASP:HB2	2.45	0.47
9:I:26:MET:HA	9:I:26:MET:CE	2.45	0.47
2:B:13:THR:O	2:B:13:THR:HG22	2.15	0.47
3:P:112:LEU:HG	3:P:118:PRO:HB3	1.95	0.47
5:E:78:HIS:ND1	9:I:12:LEU:HD22	2.30	0.47
1:A:290:HIS:CD2	1:A:291:HIS:CD2	3.03	0.47
3:C:187:THR:HG21	7:G:68:THR:HG21	1.97	0.47
1:A:506:GLU:OE2	3:C:1:MET:HG3	2.15	0.46
7:T:54:ARG:HB2	7:T:54:ARG:HH11	1.80	0.46
12:L:41:ARG:HD2	13:M:40:TYR:OH	2.15	0.46
1:N:160:GLY:O	1:N:164:PHE:HD2	1.99	0.46
4:D:16:TYR:CE1	4:D:25:PRO:HG3	2.49	0.46
2:B:145:PRO:CG	2:B:148:MET:HE2	2.46	0.46
12:Y:4:GLU:HB3	12:Y:9:LYS:HD3	1.97	0.46
7:G:83:GLU:O	7:G:84:LYS:HB2	2.15	0.46

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
4:Q:93:ALA:O	4:Q:97:ILE:HG13	2.16	0.46
1:N:138:HIS:O	1:N:213:ARG:NH2	2.48	0.46
3:C:173:PHE:C	3:C:173:PHE:CD1	2.89	0.46
4:D:110:THR:HG22	4:D:115:TRP:CE2	2.50	0.46
12:L:4:GLU:HB3	12:L:9:LYS:HD3	1.97	0.46
2:B:185:MET:SD	2:B:185:MET:C	2.93	0.46
7:T:83:GLU:O	7:T:84:LYS:HB2	2.15	0.46
3:C:183:GLU:O	7:G:42:ARG:NH2	2.48	0.46
4:D:68:PHE:CZ	5:E:70:VAL:HG23	2.50	0.46
2:O:63:THR:HA	2:O:66:THR:HG22	1.98	0.46
1:N:374:VAL:HA	1:N:377:PHE:CE2	2.50	0.46
6:F:82:CYS:N	6:F:86:GLY:O	2.49	0.46
7:G:5:LYS:O	7:G:6:GLY:C	2.51	0.46
13:M:13:LYS:HD3	13:M:13:LYS:N	2.25	0.46
8:H:39:CYS:HB3	8:H:53:CYS:SG	2.56	0.46
1:N:468:MET:SD	18:N:515:HEA:H243	2.55	0.46
3:C:16:TRP:CE3	3:C:19:THR:HG21	2.51	0.46
4:Q:40:LEU:CD2	4:Q:59:LEU:HD13	2.45	0.46
5:E:49:ASP:O	5:E:53:ARG:HG3	2.16	0.46
3:C:146:TRP:CE2	7:G:17:ARG:HB2	2.51	0.46
9:I:61:GLU:OE1	9:I:64:ARG:NH1	2.48	0.46
8:H:46:LYS:HZ1	8:U:53:CYS:HA	1.81	0.46
3:C:16:TRP:HE3	3:C:19:THR:HG21	1.81	0.46
1:A:29:VAL:HG13	12:L:36:PRO:HG3	1.97	0.46
5:E:43:PRO:HB2	5:E:48:ILE:HD11	1.97	0.46
3:P:72:THR:HB	3:P:73:PRO:CD	2.46	0.46
2:O:160:LEU:HD21	2:O:175:ILE:HG23	1.98	0.46
2:B:50:LEU:CD2	5:E:77:PRO:HD2	2.45	0.46
12:Y:41:ARG:HD2	13:Z:40:TYR:OH	2.16	0.46
5:R:49:ASP:O	5:R:53:ARG:HG3	2.15	0.46
1:N:187:SER:HB2	1:N:277:MET:HE3	1.96	0.46
8:U:36:PHE:HB2	8:U:56:TYR:CB	2.44	0.46
1:N:377:PHE:CD2	18:N:516:HEA:HAD1	2.51	0.46
6:F:31:TYR:HE2	6:F:98:HIS:CE1	2.34	0.46
2:B:160:LEU:HD23	2:B:175:ILE:HA	1.97	0.46
2:B:160:LEU:CD2	2:B:175:ILE:HG12	2.46	0.46
1:A:404:THR:O	1:A:480:ARG:NH1	2.49	0.46
1:A:484:THR:HB	13:M:2:THR:HG23	1.98	0.46
1:N:71:MET:HE1	1:N:195:LEU:HD21	1.97	0.46
6:S:31:TYR:HE2	6:S:98:HIS:HE1	1.64	0.46
9:V:61:GLU:OE1	9:V:64:ARG:NH1	2.49	0.46
6:F:37:LYS:HD3	6:F:37:LYS:N	2.31	0.46

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
5:E:90:ARG:HD3	5:E:90:ARG:HA	1.62	0.46
2:B:89[B]:GLU:HB2	2:B:91[B]:ASN:ND2	2.31	0.45
13:Z:13:LYS:CD	13:Z:13:LYS:H	2.28	0.45
1:N:117:MET:HE3	12:Y:39:ILE:HG23	1.98	0.45
3:P:204:HIS:HE1	3:P:245:VAL:O	1.99	0.45
1:N:456:MET:HG2	4:Q:96:LEU:HD13	1.98	0.45
3:P:40:MET:O	3:P:44:MET:HG2	2.16	0.45
13:M:13:LYS:CD	13:M:13:LYS:H	2.23	0.45
7:G:42:ARG:CZ	7:G:74:ARG:NH2	2.80	0.45
1:A:402:GLY:HA3	1:A:499:PRO:HD3	1.97	0.45
4:D:145:TRP:CE2	11:K:45:VAL:HA	2.52	0.45
12:Y:15:VAL:HG23	12:Y:21:LEU:HD22	1.98	0.45
1:A:337:ALA:HB2	1:A:394:VAL:HG23	1.98	0.45
1:N:353:LEU:HG	2:O:31:VAL:HB	1.98	0.45
2:B:98:LYS:HB2	2:B:98:LYS:HE3	1.70	0.45
3:C:22:LEU:HD12	3:C:22:LEU:HA	1.74	0.45
2:O:116:LEU:HD11	2:O:226:MET:CG	2.46	0.45
4:D:41:LYS:HD3	4:D:62:LEU:HD23	1.97	0.45
1:A:456:MET:HG2	4:D:96:LEU:HD13	1.97	0.45
4:D:123:MET:HE2	4:D:134:PHE:CE2	2.50	0.45
1:A:71:MET:CE	1:A:195:LEU:HD21	2.47	0.45
1:N:5:ARG:HG2	1:N:6:TRP:NE1	2.32	0.45
13:Z:13:LYS:N	13:Z:13:LYS:HD3	2.29	0.45
3:P:118:PRO:HG2	3:P:121:ILE:HG13	1.99	0.45
3:C:112:LEU:HG	3:C:118:PRO:HB3	1.98	0.45
1:N:115:SER:O	1:N:121:GLY:HA2	2.16	0.45
6:S:55:LYS:HA	6:S:74:LEU:O	2.15	0.45
3:P:101:PHE:HZ	3:P:260:GLY:HA3	1.81	0.45
5:R:12:ASP:HA	5:R:47:ILE:HD11	1.99	0.45
4:D:108:PRO:HG2	4:D:111:PHE:CE2	2.52	0.45
10:W:11:LEU:C	10:W:11:LEU:HD23	2.36	0.45
1:A:64:VAL:HG22	1:A:109:PHE:HE1	1.82	0.45
7:T:5:LYS:O	7:T:6:GLY:C	2.51	0.45
3:P:1:MET:O	3:P:3:HIS:N	2.49	0.45
2:B:128:LEU:HD11	2:B:134:ARG:HA	1.98	0.45
7:T:48:ILE:HG13	8:U:80:THR:HG22	1.99	0.45
10:J:3:ASN:OD1	10:J:5:VAL:HG12	2.16	0.45
1:A:353:LEU:HG	2:B:31:VAL:HB	1.98	0.45
1:N:10:THR:HG21	3:P:11:VAL:HG22	1.98	0.45
3:P:16:TRP:HA	3:P:16:TRP:CE3	2.52	0.45
1:A:229:ILE:HD11	2:B:175:ILE:HD13	1.98	0.45
3:C:72:THR:HB	3:C:73:PRO:CD	2.47	0.45

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
8:H:24:ASN:ND2	8:H:26:THR:H	2.14	0.45
1:N:52:GLN:O	1:N:56:VAL:HG23	2.17	0.45
1:N:218:THR:HB	1:N:221:ASP:HB3	1.98	0.45
2:B:122:MET:HB2	2:B:208:PRO:HD2	1.99	0.45
8:U:39:CYS:HB3	8:U:53:CYS:SG	2.57	0.44
3:P:57:TRP:O	3:P:61:VAL:HG23	2.17	0.44
1:N:404:THR:O	1:N:480:ARG:NH1	2.51	0.44
4:D:112:GLU:O	4:D:116:VAL:HG13	2.17	0.44
4:Q:108:PRO:HG2	4:Q:111:PHE:CE2	2.52	0.44
1:A:354:THR:HG21	1:A:376:HIS:HA	1.99	0.44
4:D:83:GLY:HA3	11:K:18:LEU:HA	1.99	0.44
7:G:7:ASP:CG	7:G:8:HIS:H	2.21	0.44
6:S:30:PRO:CB	6:S:96:LEU:HB3	2.42	0.44
4:Q:68:PHE:CZ	5:R:70:VAL:HG23	2.52	0.44
1:N:513:LEU:HD12	1:N:513:LEU:HA	1.56	0.44
10:W:3:ASN:OD1	10:W:5:VAL:HG12	2.16	0.44
5:E:52:LEU:HA	5:E:52:LEU:HD23	1.86	0.44
2:O:1:MET:SD	2:O:133:LEU:HD11	2.55	0.44
3:P:173:PHE:C	3:P:173:PHE:CD1	2.90	0.44
6:S:31:TYR:CE2	6:S:98:HIS:HE1	2.35	0.44
1:A:71:MET:HE1	1:A:195:LEU:HD21	1.99	0.44
4:Q:107:ILE:HB	4:Q:108:PRO:HD2	2.00	0.44
1:A:34:SER:HB2	18:A:515:HEA:C2B	2.47	0.44
5:R:21:LYS:O	5:R:57:ARG:NH1	2.50	0.44
1:A:495:LEU:HA	1:A:495:LEU:HD23	1.71	0.44
2:O:160:LEU:CD2	2:O:175:ILE:HG12	2.47	0.44
1:N:315:PRO:O	1:N:318:VAL:HG22	2.18	0.44
4:D:18:ASP:OD2	4:D:66:GLU:HG2	2.17	0.44
1:A:437:PRO:HG2	1:A:440:TYR:CE2	2.52	0.44
1:A:417:MET:HG3	1:A:417:MET:O	2.18	0.44
4:Q:145:TRP:CZ2	11:X:45:VAL:HA	2.52	0.44
8:H:49:ASP:HB2	8:U:49:ASP:CB	2.46	0.44
6:F:31:TYR:HE2	6:F:98:HIS:HE1	1.65	0.44
1:A:374:VAL:HA	1:A:377:PHE:CE2	2.53	0.44
3:C:224:LYS:HB3	3:C:224:LYS:NZ	2.32	0.44
1:A:315:PRO:O	1:A:318:VAL:HG22	2.18	0.44
13:Z:32:TRP:O	13:Z:36:HIS:CD2	2.65	0.44
1:A:225:GLY:HA3	3:C:112:LEU:HD13	1.99	0.44
7:T:48:ILE:HD13	7:T:50:TYR:CZ	2.52	0.44
6:S:98:HIS:N	6:S:98:HIS:ND1	2.66	0.44
1:N:5:ARG:HG2	1:N:6:TRP:CE2	2.52	0.44
9:I:27:VAL:CG1	9:I:31:PHE:HE2	2.26	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
18:A:516:HEA:HMB1	18:A:516:HEA:H11	1.75	0.44
1:A:314:ILE:HB	1:A:315:PRO:CD	2.48	0.44
2:O:121:TYR:O	2:O:138:VAL:HA	2.18	0.44
12:L:46:LYS:O	12:L:47:LYS:HB2	2.18	0.44
7:G:6:GLY:CA	1:N:190:ILE:HG12	2.43	0.44
3:C:154:GLY:HA2	6:F:6:VAL:HG13	1.99	0.44
1:N:324:LEU:O	2:O:45:MET:HE3	2.17	0.44
7:G:48:ILE:HG13	8:H:80:THR:HG22	1.99	0.44
1:A:456:MET:CE	11:K:40:TRP:CH2	3.01	0.44
8:H:24:ASN:HD22	8:H:25:GLN:N	2.15	0.44
2:O:145:PRO:CG	2:O:148:MET:HE2	2.48	0.44
4:Q:9:GLU:CD	4:Q:9:GLU:H	2.22	0.44
1:A:386:VAL:HG11	18:A:515:HEA:H261	1.99	0.43
6:S:48:LEU:HD23	6:S:90:LYS:HB3	2.00	0.43
3:C:192:VAL:O	3:C:196:THR:HB	2.18	0.43
3:P:224:LYS:HB3	3:P:224:LYS:NZ	2.33	0.43
11:K:24:PHE:O	11:K:28:VAL:HG23	2.17	0.43
8:U:39:CYS:HG	8:U:53:CYS:CB	2.12	0.43
6:S:8:THR:H	6:S:11:GLU:HG2	1.82	0.43
4:Q:40:LEU:HD22	4:Q:59:LEU:CD1	2.47	0.43
1:N:492:LEU:HD22	1:N:492:LEU:O	2.17	0.43
1:N:484:THR:HB	13:Z:2:THR:HG23	2.00	0.43
5:R:87:GLN:HG2	5:R:88:GLU:OE2	2.18	0.43
4:D:56:LYS:HB3	5:E:61:PHE:CE2	2.53	0.43
3:P:65:SER:HB3	3:P:71:HIS:CE1	2.52	0.43
3:P:110:PRO:HB3	8:U:30:TRP:CD2	2.53	0.43
11:X:43:SER:OG	11:X:45:VAL:HG12	2.18	0.43
4:Q:123:MET:HE2	4:Q:134:PHE:CE2	2.53	0.43
1:N:159:LEU:HD12	1:N:159:LEU:HA	1.85	0.43
5:E:76:GLY:HA3	5:E:77:PRO:HD2	1.86	0.43
3:P:16:TRP:HE3	3:P:16:TRP:HA	1.82	0.43
12:Y:15:VAL:O	12:Y:15:VAL:HG13	2.18	0.43
1:A:328:HIS:HB2	2:B:45:MET:HE3	2.01	0.43
2:O:212:GLU:OE1	9:V:70:GLN:HG2	2.18	0.43
13:M:28:LEU:HB2	13:M:29:PRO:HD3	1.99	0.43
4:Q:11:TYR:CD1	4:Q:11:TYR:C	2.91	0.43
1:A:308:ALA:O	1:A:311:ILE:HG12	2.18	0.43
3:C:42:LEU:HD21	10:J:46:SER:HB3	2.01	0.43
2:O:114:GLU:HG3	2:O:115:ASP:H	1.84	0.43
3:C:1:MET:O	3:C:3:HIS:N	2.49	0.43
6:F:18:ARG:NH1	6:F:21:MET:HE1	2.34	0.43
1:N:106:PRO:HB2	1:N:107:PRO:HD3	2.01	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:106:PRO:HB2	1:A:107:PRO:HD3	2.00	0.43
1:A:492:LEU:HD23	1:A:495:LEU:HD12	2.01	0.43
1:A:299:VAL:HA	1:A:302:ARG:NH1	2.33	0.43
6:F:50:PRO:HA	6:F:92:VAL:O	2.18	0.43
6:F:8:THR:H	6:F:11:GLU:HG2	1.83	0.43
1:N:225:GLY:HA3	3:P:112:LEU:HD13	2.00	0.43
6:S:51:SER:HB2	6:S:91:LEU:HD11	1.99	0.43
2:O:136:LEU:HD22	2:O:193:TYR:HB3	2.01	0.43
2:B:63:THR:HA	2:B:66:THR:HG22	2.00	0.43
1:A:138:HIS:O	1:A:213:ARG:NH2	2.51	0.43
8:U:36:PHE:HB2	8:U:56:TYR:HB3	2.00	0.43
2:O:100:MET:SD	2:O:155:SER:HB3	2.58	0.43
4:D:132:GLN:HE22	9:I:43:ARG:HB2	1.84	0.43
5:E:41:LEU:O	5:E:41:LEU:HD12	2.18	0.43
1:N:197:LEU:O	3:P:92:LEU:HD22	2.18	0.43
8:U:24:ASN:HD22	8:U:24:ASN:C	2.22	0.43
5:R:53:ARG:O	5:R:56:ARG:HB3	2.19	0.43
1:A:368:HIS:CD2	1:A:369:ASP:HB2	2.53	0.43
11:K:9:PHE:CD1	11:K:9:PHE:C	2.92	0.43
2:B:212:GLU:OE1	9:I:70:GLN:HG2	2.19	0.43
1:A:278:MET:HG2	7:T:5:LYS:O	2.19	0.43
9:V:27:VAL:CG1	9:V:31:PHE:HE2	2.28	0.43
8:U:65:PRO:HG2	8:U:68:TRP:CD2	2.54	0.43
6:F:53:THR:HG22	6:F:54:ASN:N	2.32	0.43
1:N:508:PRO:HG3	3:P:6:HIS:HB3	2.00	0.43
2:O:216:LEU:O	2:O:219:PHE:N	2.52	0.43
7:G:84:LYS:HE3	7:G:84:LYS:HB3	1.86	0.43
1:N:456:MET:CE	11:X:40:TRP:CH2	3.02	0.43
5:R:90:ARG:HD3	5:R:90:ARG:HA	1.63	0.43
7:T:42:ARG:CZ	7:T:74:ARG:NH2	2.82	0.43
1:A:95:PRO:HB2	3:C:11:VAL:HG13	2.00	0.43
1:N:353:LEU:HD12	1:N:353:LEU:HA	1.92	0.43
1:A:450:TRP:CE3	1:A:453:ILE:HD12	2.54	0.43
1:N:82:LEU:O	1:N:86:MET:HG3	2.19	0.43
8:U:57:ARG:HA	8:U:60:TYR:CE2	2.54	0.42
1:N:397:PHE:HB3	1:N:398:PRO:HD3	2.00	0.42
4:D:108:PRO:HG2	4:D:111:PHE:CD2	2.54	0.42
1:N:311:ILE:HD13	1:N:311:ILE:HG21	1.65	0.42
1:A:159:LEU:HD12	1:A:159:LEU:HA	1.72	0.42
1:N:390:MET:HE1	18:N:515:HEA:H211	2.01	0.42
3:C:26:LEU:HD23	3:C:45:ILE:HG22	2.01	0.42
2:O:193:TYR:CD1	2:O:193:TYR:N	2.87	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:N:176:MET:HG3	1:N:180:GLN:HB3	2.01	0.42
7:T:47:PHE:CD2	7:T:77:PRO:CB	3.02	0.42
5:E:87:GLN:HG2	5:E:88:GLU:OE2	2.20	0.42
5:E:27:TRP:CH2	6:F:86:GLY:HA2	2.54	0.42
6:F:31:TYR:CE2	6:F:98:HIS:HE1	2.36	0.42
18:A:516:HEA:HHC	18:A:516:HEA:O11	2.19	0.42
5:R:42:VAL:HG13	5:R:71:VAL:HG13	1.99	0.42
1:N:198:SER:HB2	1:N:238:PHE:HA	2.01	0.42
1:A:285:PHE:CE2	7:T:4:ALA:HB3	2.54	0.42
1:A:85:LEU:O	1:A:492:LEU:HD13	2.19	0.42
18:N:516:HEA:HMB1	18:N:516:HEA:H11	1.74	0.42
1:N:308:ALA:O	1:N:311:ILE:HG12	2.19	0.42
1:A:74:MET:HE1	1:A:246:LEU:O	2.19	0.42
11:X:31:TYR:CD2	11:X:35:GLN:HB2	2.53	0.42
3:C:133:ASN:ND2	3:C:176:LEU:HD13	2.34	0.42
9:I:21:ILE:HD12	9:I:21:ILE:HA	1.74	0.42
3:C:101:PHE:HZ	3:C:260:GLY:HA3	1.84	0.42
7:G:1:ALA:HB1	1:N:286:ILE:CG2	2.42	0.42
1:A:34:SER:HB3	1:A:61:HIS:CE1	2.54	0.42
1:A:498:CYS:HA	1:A:499:PRO:HA	1.90	0.42
6:S:35:ALA:HA	6:S:36:PRO:HD3	1.89	0.42
7:T:38:HIS:C	7:T:40:GLY:H	2.22	0.42
7:G:42:ARG:NH1	7:G:74:ARG:NH2	2.66	0.42
9:I:15:ARG:NH1	9:I:18:ARG:HH22	2.18	0.42
9:V:15:ARG:NH1	9:V:18:ARG:HH22	2.18	0.42
2:B:158:ASP:O	2:B:176:PRO:HG3	2.19	0.42
3:C:118:PRO:HG2	3:C:121:ILE:HG13	2.01	0.42
5:E:53:ARG:O	5:E:56:ARG:HB3	2.20	0.42
4:D:18:ASP:OD2	4:D:65:LYS:HB3	2.20	0.42
3:P:116:TRP:HA	3:P:117:PRO:C	2.39	0.42
2:B:19:GLU:CD	2:B:82:ARG:HH22	2.23	0.42
5:E:81:ILE:O	5:E:85:VAL:HG23	2.20	0.42
2:B:104:TRP:HA	2:B:207:MET:SD	2.59	0.42
1:A:61:HIS:HE1	18:A:515:HEA:C4B	2.33	0.42
3:C:16:TRP:HA	3:C:16:TRP:CE3	2.55	0.42
5:R:53:ARG:O	5:R:56:ARG:N	2.52	0.42
1:A:187:SER:HB2	1:A:277:MET:CE	2.48	0.42
2:O:145:PRO:HG2	2:O:148:MET:HE2	2.02	0.42
2:O:141:ARG:HD2	2:O:212:GLU:OE1	2.20	0.42
3:P:42:LEU:HA	3:P:42:LEU:HD12	1.86	0.42
11:X:9:PHE:C	11:X:9:PHE:CD1	2.92	0.42
9:I:36:LYS:HD2	9:I:37:PHE:CE2	2.54	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
9:V:11:GLY:O	9:V:15:ARG:HD3	2.19	0.42
1:A:377:PHE:CD2	18:A:516:HEA:HAD1	2.55	0.42
6:F:48:LEU:HD23	6:F:90:LYS:HB3	2.01	0.42
1:N:489:THR:HB	6:S:70:ILE:HG23	2.01	0.42
1:A:342:LEU:HA	1:A:342:LEU:HD23	1.89	0.42
1:N:512:ASN:HA	1:N:512:ASN:HD22	1.69	0.42
9:V:21:ILE:HA	9:V:21:ILE:HD12	1.86	0.42
8:H:46:LYS:NZ	8:U:53:CYS:CA	2.80	0.42
7:G:54:ARG:H	7:G:54:ARG:HD3	1.82	0.42
5:R:52:LEU:HD11	5:R:68:LEU:HD21	2.01	0.42
1:A:510:TYR:HB2	6:F:56:ARG:HH11	1.83	0.42
2:B:193:TYR:CD1	2:B:193:TYR:N	2.87	0.42
5:E:81:ILE:HA	9:I:7:PRO:HG3	2.01	0.42
9:V:36:LYS:O	9:V:40:ALA:HB3	2.19	0.42
1:N:431:LEU:HD21	1:N:450:TRP:HB2	2.02	0.42
4:D:23:PRO:HD2	5:E:34:ASN:OD1	2.20	0.41
13:M:32:TRP:O	13:M:36:HIS:CD2	2.65	0.41
2:O:136:LEU:HB3	2:O:193:TYR:HD2	1.79	0.41
1:A:117:MET:HE3	12:L:39:ILE:HG12	2.01	0.41
1:A:241:PRO:O	1:A:245:ILE:HG13	2.20	0.41
2:O:191:LEU:CD2	2:O:212:GLU:HG3	2.49	0.41
1:N:23:GLY:HA3	1:N:73:ILE:HG13	2.01	0.41
7:G:72:ASN:HA	7:G:73:PRO:HD3	1.84	0.41
10:J:11:LEU:HD23	10:J:11:LEU:C	2.40	0.41
8:H:36:PHE:HB2	8:H:56:TYR:HB3	2.01	0.41
18:A:515:HEA:H272	18:A:515:HEA:H172	1.87	0.41
3:C:16:TRP:HA	3:C:16:TRP:HE3	1.86	0.41
3:P:192:VAL:HA	3:P:195:SER:HB2	2.02	0.41
2:B:179:LEU:HD21	8:H:65:PRO:CD	2.50	0.41
1:A:176:MET:HE2	1:A:181:THR:HG22	2.02	0.41
3:P:72:THR:HB	3:P:73:PRO:HD2	2.02	0.41
1:N:354:THR:HG21	1:N:376:HIS:HA	2.01	0.41
3:C:169:LEU:HA	3:C:169:LEU:HD23	1.93	0.41
12:L:15:VAL:HG23	12:L:21:LEU:HD22	2.02	0.41
3:C:154:GLY:HA2	6:F:6:VAL:CG1	2.51	0.41
7:T:54:ARG:HH11	7:T:54:ARG:CG	2.34	0.41
1:N:195:LEU:CD2	1:N:245:ILE:HD13	2.50	0.41
4:D:96:LEU:HD23	4:D:96:LEU:HA	1.92	0.41
1:N:416:ILE:HD12	1:N:467:LEU:HD12	2.02	0.41
11:X:53:TRP:O	11:X:54:ARG:CB	2.68	0.41
12:Y:46:LYS:O	12:Y:47:LYS:HB2	2.20	0.41
4:Q:43:LYS:HE2	4:Q:43:LYS:HB3	1.80	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:344:PHE:CD1	1:A:344:PHE:C	2.94	0.41
8:H:46:LYS:HZ3	8:U:53:CYS:CA	2.33	0.41
8:U:39:CYS:O	8:U:43:MET:HG2	2.21	0.41
2:O:104:TRP:HA	2:O:207:MET:SD	2.61	0.41
7:G:54:ARG:HH11	7:G:54:ARG:CG	2.33	0.41
7:T:54:ARG:HD3	7:T:54:ARG:H	1.83	0.41
4:Q:23:PRO:HD2	5:R:34:ASN:OD1	2.20	0.41
1:A:442:ASP:OD2	2:B:134:ARG:NH2	2.53	0.41
2:B:189:PRO:HG2	9:I:63:MET:CE	2.50	0.41
1:A:314:ILE:HB	1:A:315:PRO:HD2	2.02	0.41
3:C:119:THR:HG21	8:H:82:PRO:O	2.21	0.41
4:D:9:GLU:H	4:D:9:GLU:CD	2.23	0.41
2:B:203:ASN:HD22	2:B:206:PHE:HD2	1.68	0.41
8:H:57:ARG:HA	8:H:60:TYR:CE2	2.56	0.41
7:G:54:ARG:HH11	7:G:54:ARG:HB2	1.85	0.41
6:S:37:LYS:N	6:S:37:LYS:CD	2.84	0.41
1:A:346:PHE:O	1:A:349:THR:HB	2.20	0.41
1:N:266:GLU:HB2	1:N:267:PRO:HD2	2.02	0.41
3:P:22:LEU:HA	3:P:22:LEU:HD12	1.85	0.41
2:O:59:GLN:N	2:O:62:GLU:HG3	2.32	0.41
9:I:15:ARG:HH11	9:I:18:ARG:HH22	1.68	0.41
1:N:131:PRO:HB2	2:O:159:VAL:HA	2.02	0.41
4:D:8:SER:OG	13:M:4:LYS:HE3	2.20	0.41
6:S:26:LYS:HB3	6:S:28:GLN:NE2	2.35	0.41
1:A:512:ASN:HD22	1:A:512:ASN:HA	1.69	0.41
1:N:241:PRO:O	1:N:245:ILE:HG13	2.21	0.41
1:N:209:LEU:O	1:N:213:ARG:HG3	2.21	0.41
1:A:456:MET:HE1	11:K:38:ILE:HD13	2.02	0.41
1:A:341:ALA:O	1:A:345:ILE:HG13	2.21	0.41
1:N:64:VAL:HG22	1:N:109:PHE:HE1	1.85	0.41
1:N:151:HIS:O	1:N:155:VAL:HG23	2.21	0.41
1:A:257:ILE:HD11	1:A:392:GLY:HA2	2.03	0.41
1:A:416:ILE:HD12	1:A:467:LEU:HD12	2.02	0.41
8:H:36:PHE:CE1	8:H:57:ARG:HB2	2.56	0.41
6:S:22:LEU:HA	6:S:25:ARG:HG2	2.02	0.41
6:S:81:ARG:HB3	6:S:86:GLY:O	2.21	0.41
4:Q:108:PRO:HG2	4:Q:111:PHE:CD2	2.56	0.41
1:A:380:VAL:HG21	18:A:516:HEA:C3C	2.51	0.41
1:N:450:TRP:CE3	1:N:453:ILE:HD12	2.55	0.41
3:P:47:LEU:O	3:P:51:MET:HG2	2.21	0.41
1:A:489:THR:HB	6:F:70:ILE:HG23	2.01	0.41
12:Y:22:LEU:HD23	12:Y:22:LEU:HA	1.71	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:N:366:VAL:O	1:N:366:VAL:HG23	2.20	0.41
10:J:29:ASN:HD21	10:J:33:ARG:HD2	1.85	0.41
3:P:19:THR:O	3:P:23:SER:HB2	2.21	0.41
4:D:127:LYS:HG3	4:D:140:TYR:OH	2.21	0.41
6:F:49:VAL:HG21	6:F:74:LEU:HD12	2.03	0.41
6:S:31:TYR:CE2	6:S:98:HIS:CE1	3.08	0.41
1:N:294:THR:HG22	1:N:365:ILE:HD13	2.03	0.41
1:A:508:PRO:HG3	3:C:6:HIS:HB3	2.02	0.41
6:S:63:GLU:HB2	6:S:66:ASN:ND2	2.36	0.41
9:V:60:PHE:CE1	9:V:69:PHE:CE2	3.09	0.41
1:N:299:VAL:HA	1:N:302:ARG:NH1	2.36	0.41
8:U:54:GLU:CD	8:U:57:ARG:NH1	2.75	0.41
2:B:103:GLN:HE21	2:B:104:TRP:HZ2	1.69	0.41
4:Q:132:GLN:NE2	9:V:42:LYS:HE3	2.36	0.41
4:Q:18:ASP:OD2	4:Q:65:LYS:HB3	2.21	0.41
9:V:36:LYS:HD2	9:V:37:PHE:CE2	2.56	0.41
1:A:289:ALA:HB3	1:A:305:PHE:CD2	2.56	0.41
7:T:7:ASP:CG	7:T:8:HIS:H	2.23	0.41
8:H:23:GLN:HB2	8:H:23:GLN:HE21	1.70	0.41
1:N:83:VAL:HB	1:N:84:PRO:HD3	2.03	0.41
1:N:34:SER:HB2	18:N:515:HEA:C2B	2.51	0.40
3:C:22:LEU:O	3:C:26:LEU:HD13	2.22	0.40
4:D:122:ARG:O	4:D:125:ASP:HB2	2.20	0.40
1:N:128:VAL:O	1:N:128:VAL:CG2	2.67	0.40
3:P:137:LEU:HD22	3:P:173:PHE:CE2	2.56	0.40
1:N:511:VAL:O	6:S:36:PRO:HD2	2.22	0.40
1:A:328:HIS:HB2	2:B:45:MET:CE	2.51	0.40
6:S:52:ILE:O	6:S:52:ILE:HG23	2.21	0.40
2:B:59:GLN:CA	2:B:62:GLU:HB2	2.47	0.40
2:O:50:LEU:HB2	9:V:13:LEU:HD13	2.03	0.40
8:H:76:ARG:HD2	8:H:76:ARG:HH11	1.75	0.40
1:A:23:GLY:HA3	1:A:73:ILE:HG13	2.03	0.40
4:Q:112:GLU:O	4:Q:116:VAL:HG13	2.21	0.40
2:B:113:TYR:OH	8:H:12:GLN:HA	2.20	0.40
2:O:19:GLU:CD	2:O:82:ARG:HH22	2.25	0.40
3:P:157:LYS:HE3	3:P:158:HIS:CD2	2.56	0.40
5:R:78:HIS:CE1	9:V:12:LEU:HD22	2.56	0.40
7:T:42:ARG:NH1	7:T:74:ARG:NH2	2.69	0.40
10:J:30:ILE:HG13	10:J:31:LEU:N	2.36	0.40
1:A:176:MET:HG3	1:A:180:GLN:HB3	2.03	0.40
6:S:82:CYS:HA	6:S:83:PRO:HD3	1.95	0.40
3:C:81:TYR:O	3:C:85:LEU:HD22	2.22	0.40

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:151:HIS:O	1:A:155:VAL:HG23	2.21	0.40
12:L:22:LEU:HA	12:L:22:LEU:HD23	1.74	0.40
3:C:252:LEU:HD23	3:C:252:LEU:HA	1.86	0.40
1:N:141:ALA:O	1:N:145:LEU:HG	2.21	0.40
3:P:137:LEU:HD12	3:P:137:LEU:HA	1.91	0.40
6:F:98:HIS:ND1	6:F:98:HIS:N	2.68	0.40
1:A:195:LEU:HG	1:A:245:ILE:HD13	2.03	0.40
3:P:80:ARG:HG2	3:P:233:PHE:CE1	2.56	0.40
3:P:12:ASN:O	3:P:13:PRO:C	2.60	0.40
1:A:247:ILE:HA	1:A:247:ILE:HD12	1.94	0.40
3:P:41:THR:O	3:P:45:ILE:HG13	2.21	0.40
6:S:8:THR:H	6:S:11:GLU:CG	2.35	0.40
6:S:53:THR:HG22	6:S:54:ASN:N	2.33	0.40
3:P:121:ILE:HG22	3:P:123:PRO:HD3	2.04	0.40
1:A:445:ASP:OD1	2:B:134:ARG:NH1	2.54	0.40
4:D:127:LYS:HD3	9:I:63:MET:HE2	2.02	0.40
2:O:160:LEU:HD23	2:O:175:ILE:HA	2.04	0.40
11:X:31:TYR:CE2	11:X:35:GLN:CB	3.04	0.40
12:L:25:MET:O	12:L:29:PHE:HD1	2.04	0.40
1:N:220:PHE:CD1	1:N:231:TYR:HB2	2.57	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all 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	512/514 (100%)	477 (93%)	33 (6%)	2 (0%)	43	80
1	N	512/514 (100%)	476 (93%)	36 (7%)	0	100	100
2	B	230/227 (101%)	202 (88%)	24 (10%)	4 (2%)	14	42
2	O	230/227 (101%)	202 (88%)	23 (10%)	5 (2%)	10	32
3	C	259/261 (99%)	249 (96%)	8 (3%)	2 (1%)	27	65
3	P	259/261 (99%)	250 (96%)	8 (3%)	1 (0%)	43	80

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
4	D	142/147 (97%)	134 (94%)	7 (5%)	1 (1%)	30	69
4	Q	142/147 (97%)	133 (94%)	8 (6%)	1 (1%)	30	69
5	E	107/109 (98%)	103 (96%)	4 (4%)	0	100	100
5	R	107/109 (98%)	103 (96%)	4 (4%)	0	100	100
6	F	96/98 (98%)	85 (88%)	6 (6%)	5 (5%)	3	9
6	S	96/98 (98%)	86 (90%)	5 (5%)	5 (5%)	3	9
7	G	82/84 (98%)	65 (79%)	14 (17%)	3 (4%)	5	16
7	T	82/84 (98%)	64 (78%)	16 (20%)	2 (2%)	9	29
8	H	73/85 (86%)	64 (88%)	8 (11%)	1 (1%)	16	49
8	U	73/85 (86%)	62 (85%)	10 (14%)	1 (1%)	16	49
9	I	71/73 (97%)	66 (93%)	5 (7%)	0	100	100
9	V	71/73 (97%)	66 (93%)	5 (7%)	0	100	100
10	J	54/59 (92%)	49 (91%)	2 (4%)	3 (6%)	3	7
10	W	54/59 (92%)	50 (93%)	1 (2%)	3 (6%)	3	7
11	K	47/56 (84%)	40 (85%)	7 (15%)	0	100	100
11	X	47/56 (84%)	41 (87%)	6 (13%)	0	100	100
12	L	45/47 (96%)	41 (91%)	4 (9%)	0	100	100
12	Y	45/47 (96%)	42 (93%)	3 (7%)	0	100	100
13	M	41/46 (89%)	40 (98%)	1 (2%)	0	100	100
13	Z	41/46 (89%)	40 (98%)	0	1 (2%)	9	29
All	All	3518/3612 (97%)	3230 (92%)	248 (7%)	40 (1%)	21	57

All (40) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	91[A]	ASN
2	B	91[B]	ASN
4	D	20	ARG
6	F	87	THR
6	F	96	LEU
7	G	11	THR
10	J	2	GLU
10	J	3	ASN
2	O	91[A]	ASN
2	O	91[B]	ASN

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Mol	Chain	Res	Type
4	Q	20	ARG
6	S	87	THR
6	S	96	LEU
7	T	11	THR
10	W	2	GLU
10	W	3	ASN
2	B	104	TRP
6	F	2	SER
6	F	15	GLY
6	F	95	GLN
7	G	4	ALA
8	H	46	LYS
2	O	104	TRP
6	S	2	SER
6	S	15	GLY
6	S	95	GLN
7	T	4	ALA
8	U	46	LYS
1	A	508	PRO
3	C	128	GLU
3	P	128	GLU
10	W	23	LYS
2	B	103	GLN
2	O	103	GLN
3	C	229	SER
10	J	23	LYS
7	G	61	SER
1	A	384	GLY
2	O	159	VAL
13	Z	9	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	427/427 (100%)	395 (92%)	32 (8%)	19	47
1	N	427/427 (100%)	395 (92%)	32 (8%)	19	47

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	B	216/211 (102%)	189 (88%)	27 (12%)	7	19
2	O	216/211 (102%)	191 (88%)	25 (12%)	8	23
3	C	226/226 (100%)	199 (88%)	27 (12%)	8	22
3	P	226/226 (100%)	199 (88%)	27 (12%)	8	22
4	D	128/129 (99%)	117 (91%)	11 (9%)	15	40
4	Q	128/129 (99%)	118 (92%)	10 (8%)	18	45
5	E	95/95 (100%)	86 (90%)	9 (10%)	12	33
5	R	95/95 (100%)	87 (92%)	8 (8%)	16	41
6	F	81/81 (100%)	73 (90%)	8 (10%)	11	31
6	S	81/81 (100%)	74 (91%)	7 (9%)	15	40
7	G	68/68 (100%)	56 (82%)	12 (18%)	3	8
7	T	68/68 (100%)	55 (81%)	13 (19%)	2	6
8	H	67/75 (89%)	54 (81%)	13 (19%)	2	6
8	U	67/75 (89%)	55 (82%)	12 (18%)	2	7
9	I	58/58 (100%)	51 (88%)	7 (12%)	7	21
9	V	58/58 (100%)	51 (88%)	7 (12%)	7	21
10	J	47/50 (94%)	40 (85%)	7 (15%)	4	12
10	W	47/50 (94%)	40 (85%)	7 (15%)	4	12
11	K	39/46 (85%)	37 (95%)	2 (5%)	33	69
11	X	39/46 (85%)	37 (95%)	2 (5%)	33	69
12	L	40/40 (100%)	35 (88%)	5 (12%)	7	19
12	Y	40/40 (100%)	34 (85%)	6 (15%)	4	12
13	M	37/38 (97%)	34 (92%)	3 (8%)	17	43
13	Z	37/38 (97%)	34 (92%)	3 (8%)	17	43
All	All	3058/3088 (99%)	2736 (90%)	322 (10%)	10	27

All (322) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	18	LEU
1	A	51	ASP
1	A	96	ARG
1	A	105	LEU
1	A	106	PRO

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Mol	Chain	Res	Type
1	A	109	PHE
1	A	136	LEU
1	A	138	HIS
1	A	156	SER
1	A	159	LEU
1	A	180	GLN
1	A	199	LEU
1	A	238	PHE
1	A	265	LYS
1	A	273	MET
1	A	295	VAL
1	A	301	THR
1	A	306	THR
1	A	312	ILE
1	A	324	LEU
1	A	333	LYS
1	A	336	PRO
1	A	338	MET
1	A	347	LEU
1	A	353	LEU
1	A	354	THR
1	A	366	VAL
1	A	369	ASP
1	A	373	VAL
1	A	474	GLU
1	A	492	LEU
1	A	512	ASN
2	B	3	TYR
2	B	7	LEU
2	B	16	ILE
2	B	31	VAL
2	B	45	MET
2	B	54	SER
2	B	60	GLU
2	B	63	THR
2	B	65	TRP
2	B	67	ILE
2	B	88[A]	ASP
2	B	88[B]	ASP
2	B	104	TRP
2	B	115	ASP
2	B	125	THR

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Mol	Chain	Res	Type
2	B	130	PRO
2	B	138	VAL
2	B	142	VAL
2	B	147	GLU
2	B	167	SER
2	B	170	LEU
2	B	171	LYS
2	B	185	MET
2	B	188	ARG
2	B	205	SER
2	B	216	LEU
2	B	220	GLU
3	C	9	HIS
3	C	11	VAL
3	C	13	PRO
3	C	18	LEU
3	C	19	THR
3	C	22	LEU
3	C	23	SER
3	C	38	ASN
3	C	85	LEU
3	C	92	LEU
3	C	112	LEU
3	C	128	GLU
3	C	131	LEU
3	C	132	LEU
3	C	137	LEU
3	C	142	VAL
3	C	159	MET
3	C	160	LEU
3	C	163	LEU
3	C	168	THR
3	C	179	SER
3	C	188	ILE
3	C	192	VAL
3	C	196	THR
3	C	214	PHE
3	C	222	GLN
3	C	230	ASN
4	D	19	ARG
4	D	27	VAL
4	D	31	LYS

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Mol	Chain	Res	Type
4	D	40	LEU
4	D	59	LEU
4	D	62	LEU
4	D	99	GLU
4	D	107	ILE
4	D	116	VAL
4	D	127	LYS
4	D	147	LYS
5	E	29	LEU
5	E	37	VAL
5	E	41	LEU
5	E	70	VAL
5	E	79	LYS
5	E	80	GLU
5	E	99	SER
5	E	101	PRO
5	E	109	VAL
6	F	6	VAL
6	F	34	LEU
6	F	37	LYS
6	F	66	ASN
6	F	74	LEU
6	F	95	GLN
6	F	96	LEU
6	F	98	HIS
7	G	5	LYS
7	G	8	HIS
7	G	14	ARG
7	G	17	ARG
7	G	34	ASN
7	G	36	TRP
7	G	37	LEU
7	G	41	HIS
7	G	42	ARG
7	G	48	ILE
7	G	54	ARG
7	G	56	ARG
8	H	19	ARG
8	H	24	ASN
8	H	27	ARG
8	H	29	CYS
8	H	41	LYS

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Mol	Chain	Res	Type
8	H	49	ASP
8	H	53	CYS
8	H	57	ARG
8	H	60	TYR
8	H	67	SER
8	H	70	SER
8	H	75	ARG
8	H	82	PRO
9	I	2	THR
9	I	8	GLN
9	I	26	MET
9	I	41	GLU
9	I	42	LYS
9	I	44	LYS
9	I	64	ARG
10	J	1	PHE
10	J	2	GLU
10	J	8	LYS
10	J	10	LYS
10	J	27	THR
10	J	47	LEU
10	J	50	LEU
11	K	45	VAL
11	K	49	THR
12	L	15	VAL
12	L	20	ARG
12	L	22	LEU
12	L	26	THR
12	L	27	LEU
13	M	13	LYS
13	M	24	LEU
13	M	42	LYS
1	N	18	LEU
1	N	51	ASP
1	N	96	ARG
1	N	105	LEU
1	N	109	PHE
1	N	136	LEU
1	N	138	HIS
1	N	156	SER
1	N	159	LEU
1	N	180	GLN

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Mol	Chain	Res	Type
1	N	188	VAL
1	N	199	LEU
1	N	238	PHE
1	N	265	LYS
1	N	295	VAL
1	N	301	THR
1	N	306	THR
1	N	312	ILE
1	N	324	LEU
1	N	333	LYS
1	N	336	PRO
1	N	338	MET
1	N	347	LEU
1	N	353	LEU
1	N	354	THR
1	N	366	VAL
1	N	369	ASP
1	N	373	VAL
1	N	444	PRO
1	N	474	GLU
1	N	492	LEU
1	N	512	ASN
2	O	3	TYR
2	O	7	LEU
2	O	16	ILE
2	O	31	VAL
2	O	54	SER
2	O	60	GLU
2	O	63	THR
2	O	65	TRP
2	O	67	ILE
2	O	88[A]	ASP
2	O	88[B]	ASP
2	O	104	TRP
2	O	115	ASP
2	O	125	THR
2	O	130	PRO
2	O	142	VAL
2	O	147	GLU
2	O	167	SER
2	O	170	LEU
2	O	171	LYS

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Mol	Chain	Res	Type
2	O	185	MET
2	O	188	ARG
2	O	205	SER
2	O	216	LEU
2	O	220	GLU
3	P	9	HIS
3	P	11	VAL
3	P	13	PRO
3	P	18	LEU
3	P	19	THR
3	P	22	LEU
3	P	23	SER
3	P	38	ASN
3	P	85	LEU
3	P	92	LEU
3	P	112	LEU
3	P	128	GLU
3	P	130	PRO
3	P	131	LEU
3	P	132	LEU
3	P	137	LEU
3	P	142	VAL
3	P	159	MET
3	P	160	LEU
3	P	163	LEU
3	P	168	THR
3	P	179	SER
3	P	188	ILE
3	P	196	THR
3	P	214	PHE
3	P	222	GLN
3	P	230	ASN
4	Q	19	ARG
4	Q	27	VAL
4	Q	31	LYS
4	Q	40	LEU
4	Q	59	LEU
4	Q	99	GLU
4	Q	107	ILE
4	Q	116	VAL
4	Q	127	LYS
4	Q	147	LYS

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Mol	Chain	Res	Type
5	R	29	LEU
5	R	37	VAL
5	R	70	VAL
5	R	79	LYS
5	R	80	GLU
5	R	99	SER
5	R	101	PRO
5	R	109	VAL
6	S	6	VAL
6	S	37	LYS
6	S	66	ASN
6	S	74	LEU
6	S	95	GLN
6	S	96	LEU
6	S	98	HIS
7	T	5	LYS
7	T	8	HIS
7	T	14	ARG
7	T	17	ARG
7	T	26	PRO
7	T	34	ASN
7	T	36	TRP
7	T	37	LEU
7	T	41	HIS
7	T	42	ARG
7	T	48	ILE
7	T	54	ARG
7	T	56	ARG
8	U	19	ARG
8	U	24	ASN
8	U	29	CYS
8	U	41	LYS
8	U	49	ASP
8	U	53	CYS
8	U	57	ARG
8	U	60	TYR
8	U	67	SER
8	U	70	SER
8	U	75	ARG
8	U	82	PRO
9	V	2	THR
9	V	8	GLN

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Mol	Chain	Res	Type
9	V	26	MET
9	V	41	GLU
9	V	42	LYS
9	V	44	LYS
9	V	64	ARG
10	W	1	PHE
10	W	2	GLU
10	W	8	LYS
10	W	10	LYS
10	W	27	THR
10	W	47	LEU
10	W	50	LEU
11	X	45	VAL
11	X	49	THR
12	Y	11	ILE
12	Y	15	VAL
12	Y	20	ARG
12	Y	22	LEU
12	Y	26	THR
12	Y	27	LEU
13	Z	13	LYS
13	Z	24	LEU
13	Z	42	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (36) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	256	HIS
1	A	413	HIS
1	A	512	ASN
3	C	71	HIS
3	C	133	ASN
3	C	204	HIS
3	C	222	GLN
4	D	132	GLN
6	F	66	ASN
6	F	94	HIS
6	F	95	GLN
7	G	41	HIS
8	H	23	GLN
8	H	24	ASN
9	I	20	HIS

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Mol	Chain	Res	Type
10	J	21	HIS
11	K	10	HIS
13	M	36	HIS
13	M	39	ASN
1	N	256	HIS
1	N	512	ASN
3	P	71	HIS
3	P	133	ASN
3	P	204	HIS
3	P	222	GLN
4	Q	132	GLN
4	Q	143	ASN
6	S	66	ASN
6	S	94	HIS
6	S	95	GLN
7	T	41	HIS
8	U	23	GLN
8	U	24	ASN
11	X	10	HIS
13	Z	36	HIS
13	Z	39	ASN

5.3.3 RNA ⓘ

There are no RNA chains in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 18 ligands modelled in this entry, 12 are monoatomic - leaving 6 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link

column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
18	HEA	A	515	1	67,67,67	1.79	13 (19%)	80,103,103	2.70	31 (38%)
18	HEA	A	516	19,1	67,67,67	1.99	15 (22%)	80,103,103	2.11	29 (36%)
19	CMO	A	520	18,14	0,1,1	0.00	-	0,0,0	0.00	-
18	HEA	N	515	1	67,67,67	1.69	9 (13%)	80,103,103	2.53	30 (37%)
18	HEA	N	516	19,1	67,67,67	1.99	12 (17%)	80,103,103	2.15	27 (33%)
19	CMO	N	520	18,14	0,1,1	0.00	-	0,0,0	0.00	-

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
18	HEA	A	515	1	-	0/30/76/76	0/0/8/8
18	HEA	A	516	19,1	-	0/30/76/76	0/0/8/8
19	CMO	A	520	18,14	-	0/0/0/0	0/0/0/0
18	HEA	N	515	1	-	0/30/76/76	0/0/8/8
18	HEA	N	516	19,1	-	0/30/76/76	0/0/8/8
19	CMO	N	520	18,14	-	0/0/0/0	0/0/0/0

All (49) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
18	A	516	HEA	C1B-C2B	7.36	1.49	1.40
18	N	516	HEA	C3C-C2C	-6.76	1.29	1.41
18	N	516	HEA	C1B-C2B	6.66	1.48	1.40
18	A	515	HEA	C4B-NB	6.10	1.45	1.37
18	N	516	HEA	C1C-C2C	5.74	1.47	1.40
18	A	515	HEA	C1A-NA	5.36	1.44	1.37
18	A	516	HEA	C3C-C2C	-5.29	1.32	1.41
18	N	515	HEA	C1C-C2C	5.20	1.46	1.40
18	N	515	HEA	C3B-C11	-5.10	1.47	1.52
18	N	516	HEA	C3C-C4C	5.07	1.48	1.40
18	N	515	HEA	C3C-C4C	5.03	1.48	1.40
18	A	516	HEA	C3A-C2A	-4.95	1.32	1.41
18	A	515	HEA	C3B-C11	-4.27	1.48	1.52
18	N	515	HEA	C1B-C2B	3.94	1.45	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
18	A	515	HEA	C3A-CMA	-3.74	1.37	1.46
18	N	515	HEA	C4B-NB	3.67	1.42	1.37
18	A	516	HEA	C3C-C4C	3.63	1.46	1.40
18	A	516	HEA	C3A-CMA	-3.62	1.38	1.46
18	N	515	HEA	C3A-C2A	-3.58	1.35	1.41
18	A	515	HEA	C3C-C4C	3.53	1.45	1.40
18	A	516	HEA	C3B-C2B	-3.48	1.27	1.40
18	A	515	HEA	C1B-C2B	3.48	1.44	1.40
18	N	516	HEA	C1A-NA	3.45	1.42	1.37
18	A	515	HEA	C1C-C2C	3.37	1.44	1.40
18	A	516	HEA	C1C-NC	3.32	1.43	1.36
18	N	515	HEA	C3A-CMA	-3.30	1.38	1.46
18	A	515	HEA	C1C-NC	3.29	1.43	1.36
18	N	516	HEA	C3A-CMA	-3.15	1.39	1.46
18	A	515	HEA	C4C-NC	3.12	1.42	1.36
18	N	516	HEA	CAA-C2A	3.03	1.58	1.52
18	A	516	HEA	CAA-C2A	2.97	1.58	1.52
18	A	515	HEA	C4D-ND	2.94	1.42	1.36
18	A	516	HEA	C3B-C11	-2.93	1.49	1.52
18	A	516	HEA	FE-NA	2.93	2.05	1.92
18	N	515	HEA	C1A-NA	2.70	1.40	1.37
18	A	516	HEA	C1C-C2C	2.49	1.43	1.40
18	A	516	HEA	C4C-CHD	-2.49	1.33	1.39
18	A	515	HEA	C1B-CHB	-2.42	1.33	1.39
18	A	516	HEA	CMC-C2C	-2.36	1.46	1.51
18	N	515	HEA	C4C-NC	2.33	1.41	1.36
18	N	516	HEA	C26-C15	-2.29	1.45	1.50
18	N	516	HEA	C3B-C11	-2.27	1.50	1.52
18	A	515	HEA	C3C-CAC	-2.21	1.47	1.49
18	N	516	HEA	C3B-C2B	-2.14	1.32	1.40
18	A	516	HEA	CAD-C3D	-2.12	1.48	1.52
18	A	515	HEA	C27-C19	-2.10	1.45	1.50
18	N	516	HEA	C4A-C3A	2.10	1.44	1.41
18	A	516	HEA	FE-ND	-2.07	1.83	1.92
18	N	516	HEA	C1D-ND	2.06	1.40	1.37

All (117) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
18	A	515	HEA	C4B-C3B-C2B	-10.21	99.73	106.87
18	N	515	HEA	C4B-C3B-C2B	-8.24	101.11	106.87
18	A	515	HEA	C1B-C2B-C3B	7.01	111.87	107.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
18	N	515	HEA	C2C-C1C-NC	5.82	113.81	109.41
18	N	515	HEA	C13-C14-C15	-5.79	115.31	127.80
18	N	516	HEA	C1D-C2D-C3D	-5.67	103.29	106.89
18	A	515	HEA	C13-C14-C15	-5.56	115.81	127.80
18	A	515	HEA	C3B-C4B-NB	5.38	115.12	109.90
18	A	515	HEA	C4B-C3B-C11	5.03	133.74	124.67
18	A	515	HEA	C17-C18-C19	-4.71	117.63	127.80
18	A	515	HEA	CMB-C2B-C1B	-4.70	121.40	128.62
18	N	516	HEA	C2C-C1C-NC	4.64	112.91	109.41
18	N	515	HEA	C1B-C2B-C3B	4.62	110.21	107.00
18	A	515	HEA	C4B-NB-C1B	-4.61	100.69	106.76
18	A	515	HEA	C2C-C1C-NC	4.58	112.87	109.41
18	A	516	HEA	C4B-NB-C1B	-4.51	100.82	106.76
18	N	515	HEA	C3B-C4B-NB	4.45	114.21	109.90
18	N	516	HEA	C4C-NC-C1C	-4.42	100.95	106.76
18	A	516	HEA	C2B-C1B-NB	4.41	112.74	109.41
18	A	516	HEA	C2D-C1D-ND	4.38	112.72	109.41
18	N	515	HEA	C1A-C2A-C3A	-4.38	103.17	106.80
18	A	516	HEA	CBA-CAA-C2A	4.34	120.46	112.35
18	N	516	HEA	C3C-C4C-NC	4.32	112.84	108.64
18	N	516	HEA	C1D-CHD-C4C	-4.23	121.91	127.47
18	N	516	HEA	CBA-CAA-C2A	4.12	120.05	112.35
18	N	515	HEA	C12-C11-C3B	4.08	123.61	112.64
18	N	515	HEA	CMB-C2B-C1B	-4.04	122.40	128.62
18	N	516	HEA	C2A-C1A-NA	4.04	112.67	109.64
18	N	516	HEA	C4B-CHC-C1C	-4.01	122.19	127.47
18	A	516	HEA	C4C-NC-C1C	-4.00	101.49	106.76
18	N	516	HEA	C2D-C1D-ND	3.97	112.41	109.41
18	A	515	HEA	CMC-C2C-C1C	-3.93	122.58	128.62
18	A	516	HEA	C4B-CHC-C1C	-3.90	122.34	127.47
18	A	516	HEA	C1D-C2D-C3D	-3.88	104.43	106.89
18	A	515	HEA	C2B-C1B-NB	3.87	112.33	109.41
18	A	515	HEA	C2D-C1D-ND	3.85	112.32	109.41
18	N	515	HEA	C3C-C4C-NC	3.85	112.38	108.64
18	N	515	HEA	C4B-C3B-C11	3.82	131.57	124.67
18	A	515	HEA	C4C-NC-C1C	-3.79	101.78	106.76
18	N	515	HEA	C4B-CHC-C1C	-3.74	122.55	127.47
18	A	516	HEA	C3B-C4B-NB	3.72	113.51	109.90
18	N	515	HEA	C4A-C3A-C2A	3.70	109.35	106.78
18	A	515	HEA	C26-C15-C14	-3.67	116.26	123.52
18	N	515	HEA	CMC-C2C-C1C	-3.65	123.00	128.62
18	N	515	HEA	CMC-C2C-C3C	3.63	130.69	124.97

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
18	A	515	HEA	C12-C11-C3B	3.58	122.26	112.64
18	A	516	HEA	C2A-C1A-NA	3.57	112.31	109.64
18	N	516	HEA	C20-C19-C18	-3.56	114.23	121.08
18	N	516	HEA	C1A-CHA-C4D	-3.50	122.86	127.47
18	N	515	HEA	C2D-C1D-ND	3.49	112.05	109.41
18	N	516	HEA	C25-C23-C24	3.49	123.53	114.62
18	A	515	HEA	CMC-C2C-C3C	3.43	130.37	124.97
18	A	516	HEA	C3C-C4C-NC	3.40	111.95	108.64
18	A	516	HEA	C2C-C1C-NC	3.37	111.96	109.41
18	N	515	HEA	C2C-C1C-CHC	-3.32	119.70	126.00
18	A	516	HEA	C2C-C3C-CAC	-3.29	120.60	127.33
18	A	515	HEA	C1D-ND-C4D	-3.28	102.44	106.76
18	A	515	HEA	C1A-CHA-C4D	-3.27	123.17	127.47
18	N	515	HEA	C4B-NB-C1B	-3.27	102.46	106.76
18	N	515	HEA	C2A-C1A-NA	3.24	112.07	109.64
18	N	516	HEA	C2B-C1B-NB	3.19	111.82	109.41
18	A	515	HEA	C3C-C4C-NC	3.17	111.73	108.64
18	A	515	HEA	C26-C15-C16	3.14	120.16	115.39
18	N	515	HEA	C26-C15-C14	-3.10	117.38	123.52
18	N	515	HEA	C2B-C1B-NB	3.09	111.74	109.41
18	N	515	HEA	C4C-NC-C1C	-3.08	102.70	106.76
18	A	516	HEA	C20-C19-C18	-3.07	115.18	121.08
18	A	515	HEA	O11-C11-C12	-3.05	97.32	110.45
18	N	516	HEA	C4B-NB-C1B	-3.03	102.77	106.76
18	A	515	HEA	C4B-CHC-C1C	-3.01	123.51	127.47
18	A	516	HEA	CMB-C2B-C3B	-2.99	119.30	124.94
18	N	516	HEA	C3B-C4B-NB	2.98	112.78	109.90
18	N	515	HEA	C12-C13-C14	-2.91	104.29	112.46
18	N	516	HEA	C4D-C3D-C2D	2.91	109.93	106.92
18	N	515	HEA	C27-C19-C18	-2.87	117.83	123.52
18	A	516	HEA	CHD-C4C-NC	-2.84	119.83	124.58
18	A	515	HEA	C1D-CHD-C4C	-2.84	123.73	127.47
18	N	516	HEA	C1D-ND-C4D	-2.81	103.07	106.76
18	A	515	HEA	C12-C13-C14	-2.79	104.63	112.46
18	A	516	HEA	C17-C16-C15	2.78	121.93	112.74
18	N	516	HEA	C4B-C3B-C11	2.75	129.62	124.67
18	A	515	HEA	CBA-CAA-C2A	2.74	117.48	112.35
18	N	515	HEA	C17-C18-C19	-2.74	121.89	127.80
18	A	516	HEA	C1A-C2A-C3A	-2.74	104.53	106.80
18	A	516	HEA	C25-C23-C24	2.72	121.55	114.62
18	N	515	HEA	C1D-CHD-C4C	-2.69	123.93	127.47
18	A	516	HEA	C4B-C3B-C2B	-2.69	104.99	106.87

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
18	A	515	HEA	C20-C19-C18	2.67	126.22	121.08
18	N	516	HEA	CAA-C2A-C1A	2.65	129.44	124.67
18	A	516	HEA	C1A-CHA-C4D	-2.62	124.02	127.47
18	A	516	HEA	C1D-ND-C4D	-2.59	103.35	106.76
18	N	516	HEA	C17-C16-C15	2.59	121.31	112.74
18	N	515	HEA	C20-C19-C18	2.58	126.05	121.08
18	A	516	HEA	C4A-C3A-C2A	2.55	108.55	106.78
18	N	515	HEA	C1D-ND-C4D	-2.53	103.43	106.76
18	A	516	HEA	C3D-C4D-CHA	-2.46	121.34	126.00
18	N	515	HEA	C1A-CHA-C4D	-2.42	124.29	127.47
18	A	515	HEA	C27-C19-C18	-2.40	118.76	123.52
18	N	516	HEA	CAD-C3D-C2D	-2.38	122.20	129.00
18	A	516	HEA	C4A-NA-C1A	-2.38	102.02	105.58
18	N	516	HEA	CMB-C2B-C3B	-2.37	120.48	124.94
18	A	516	HEA	CMD-C2D-C1D	2.34	130.79	126.16
18	A	516	HEA	CAA-C2A-C1A	2.32	128.85	124.67
18	N	516	HEA	C4B-C3B-C2B	-2.32	105.25	106.87
18	A	516	HEA	CAD-C3D-C2D	-2.31	122.40	129.00
18	A	516	HEA	C27-C19-C20	2.29	118.87	115.39
18	N	515	HEA	CBA-CAA-C2A	2.29	116.62	112.35
18	N	515	HEA	O11-C11-C12	-2.28	100.61	110.45
18	N	516	HEA	C1A-C2A-C3A	-2.23	104.95	106.80
18	N	516	HEA	C3D-C4D-CHA	-2.18	121.87	126.00
18	A	515	HEA	C4A-CHB-C1B	-2.17	124.61	127.47
18	A	515	HEA	C16-C17-C18	-2.14	105.49	111.62
18	N	516	HEA	C27-C19-C18	2.08	127.65	123.52
18	A	516	HEA	CHA-C4D-ND	2.08	128.05	124.58
18	N	516	HEA	C2C-C3C-CAC	-2.08	123.08	127.33
18	A	515	HEA	C2C-C1C-CHC	-2.04	122.14	126.00
18	A	515	HEA	C4C-C3C-CAC	-2.02	123.08	127.18

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

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 will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates ⓘ

EDS was not executed - this section will therefore be empty.

6.4 Ligands ⓘ

EDS was not executed - this section will therefore be empty.

6.5 Other polymers ⓘ

EDS was not executed - this section will therefore be empty.