



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 13, 2018 – 11:16 PM EDT

PDB ID : 1OCC
Title : STRUCTURE OF BOVINE HEART CYTOCHROME C OXIDASE AT THE FULLY OXIDIZED STATE
Authors : Tsukihara, T.; Aoyama, H.; Yamashita, E.; Tomizaki, T.; Yamaguchi, H.; Shinzawa-Itoh, K.; Nakashima, R.; Yaono, R.; Yoshikawa, S.
Deposited on : 1996-04-18
Resolution : 2.80 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

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : rb-20030736

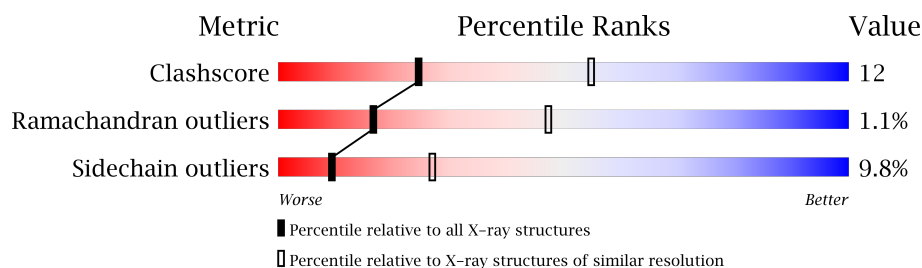
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 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	112137	3033 (2.80-2.80)
Ramachandran outliers	110173	2983 (2.80-2.80)
Sidechain outliers	110143	2985 (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. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

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	

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Mol	Chain	Length	Quality of chain
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	
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	

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
16	HEA	A	515	X	-	-	-
16	HEA	A	516	X	-	-	-
16	HEA	N	515	X	-	-	-
16	HEA	N	516	X	-	-	-

2 Entry composition

There are 17 unique types of molecules in this entry. The entry contains 28722 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 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	0	0
			1822	1184	281	339	18			
2	O	227	Total	C	N	O	S	0	0	0
			1822	1184	281	339	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			

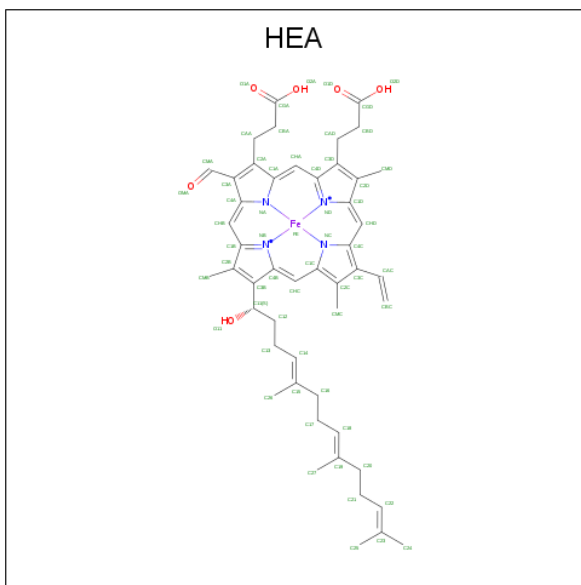
- Molecule 14 is COPPER (II) ION (three-letter code: CU) (formula: Cu).

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 HEME-A (three-letter code: HEA) (formula: $C_{49}H_{56}FeN_4O_6$).



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

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

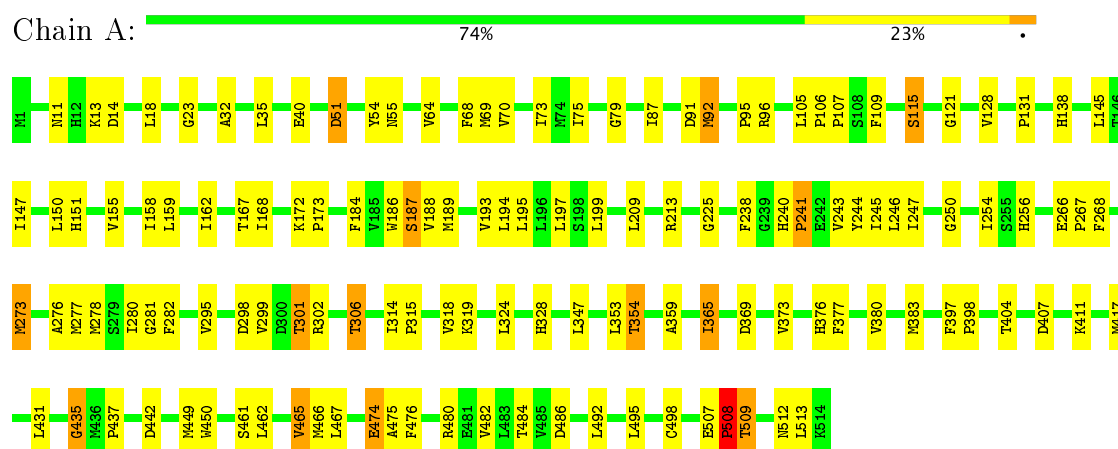
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
17	S	1	Total Zn 1 1	0	0
17	F	1	Total Zn 1 1	0	0

3 Residue-property plots

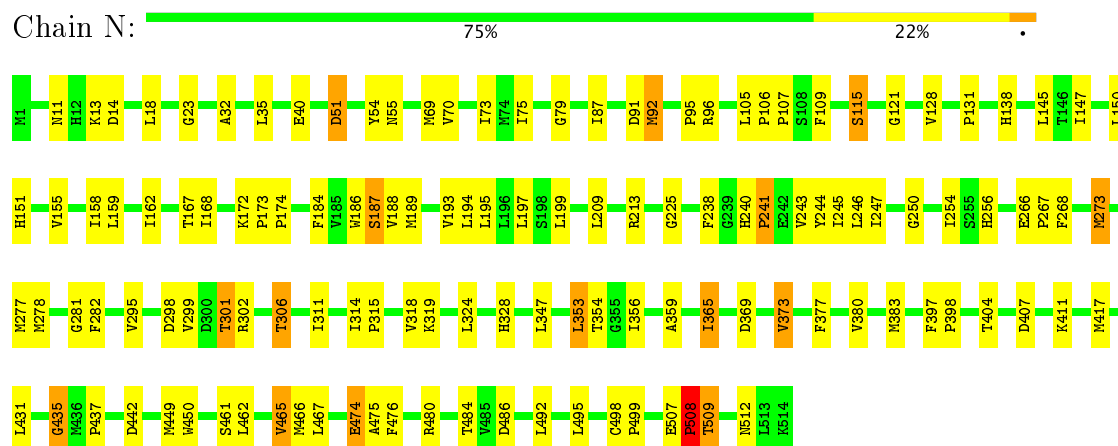
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are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

Note EDS was not executed.

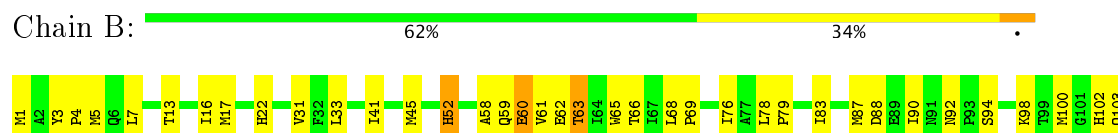
• Molecule 1: CYTOCHROME C OXIDASE

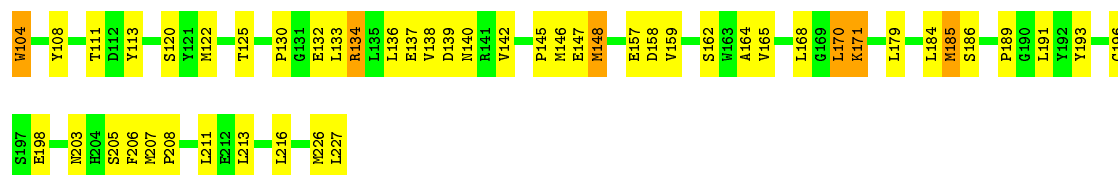


• Molecule 1: CYTOCHROME C OXIDASE

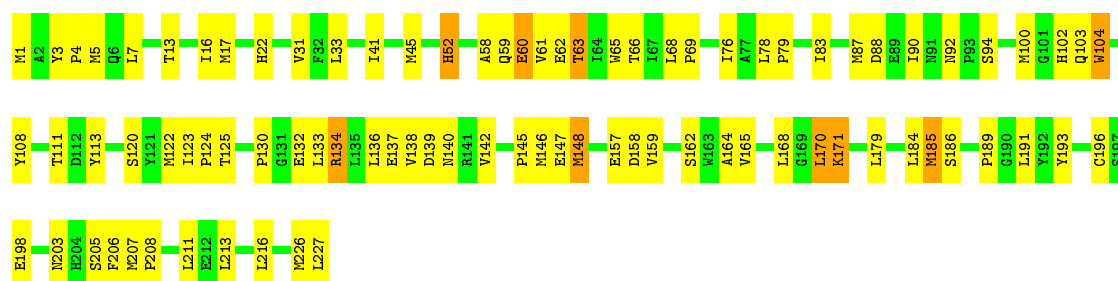


• Molecule 2: CYTOCHROME C OXIDASE

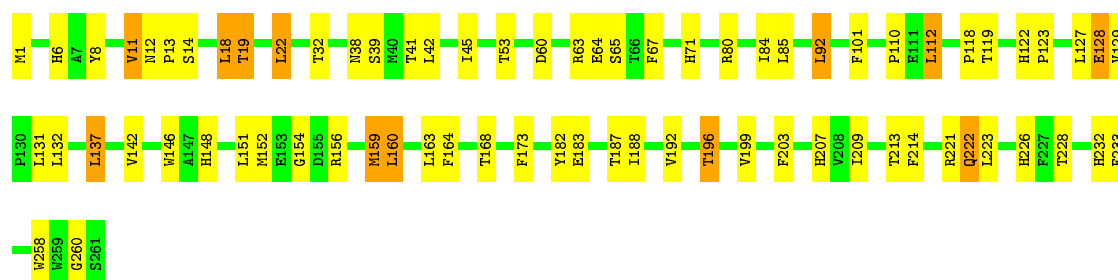




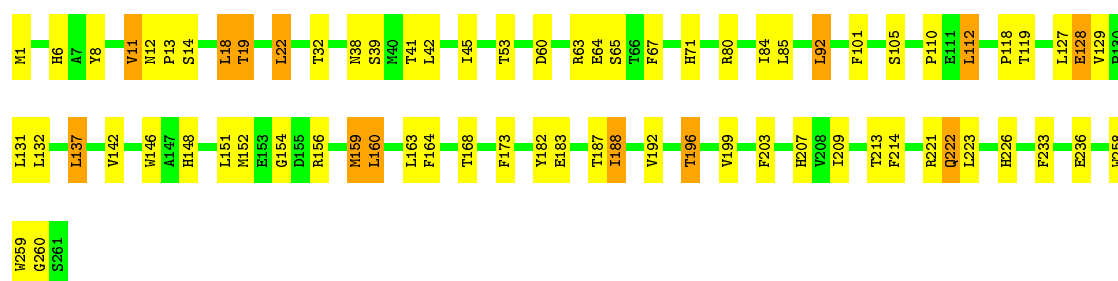
- Molecule 2: CYTOCHROME C OXIDASE



- Molecule 3: CYTOCHROME C OXIDASE



- Molecule 3: CYTOCHROME C OXIDASE



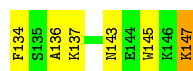
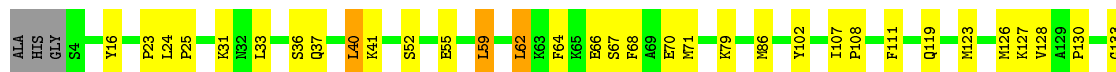
- Molecule 4: CYTOCHROME C OXIDASE





• Molecule 4: CYTOCHROME C OXIDASE

Chain Q: 71% 24% . .



• Molecule 5: CYTOCHROME C OXIDASE

Chain E: 77% 20% .



• Molecule 5: CYTOCHROME C OXIDASE

Chain R: 78% 20% .



• Molecule 6: CYTOCHROME C OXIDASE

Chain F: 61% 33% 6%



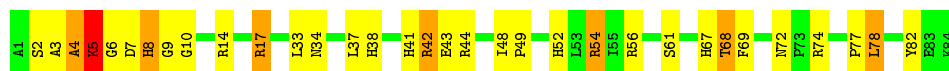
• Molecule 6: CYTOCHROME C OXIDASE

Chain S: 64% 30% 6%

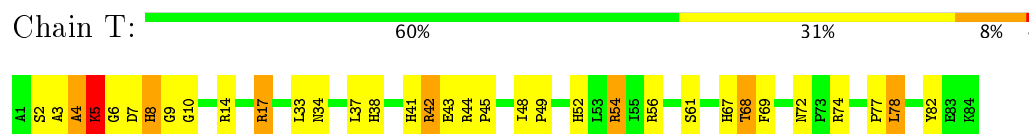


• Molecule 7: CYTOCHROME C OXIDASE

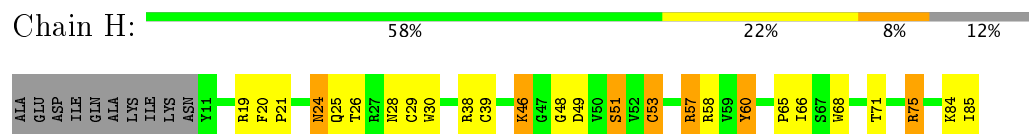
Chain G: 61% 30% 8%



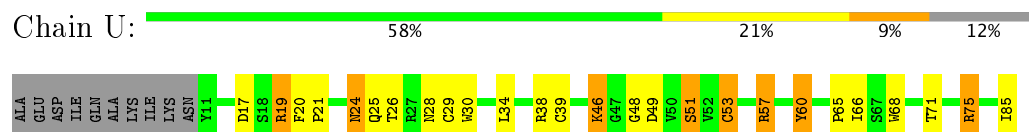
• Molecule 7: CYTOCHROME C OXIDASE



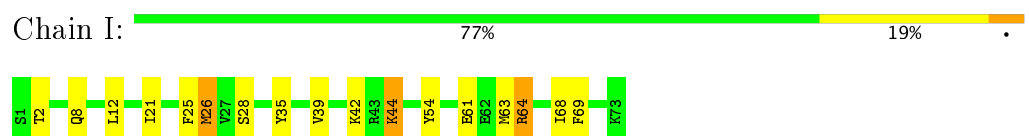
- Molecule 8: CYTOCHROME C OXIDASE



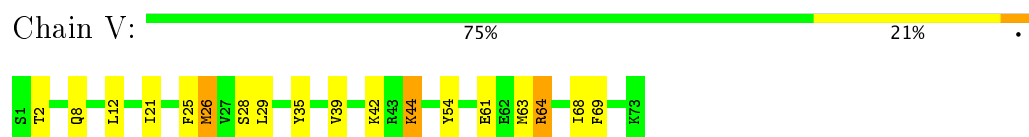
- Molecule 8: CYTOCHROME C OXIDASE



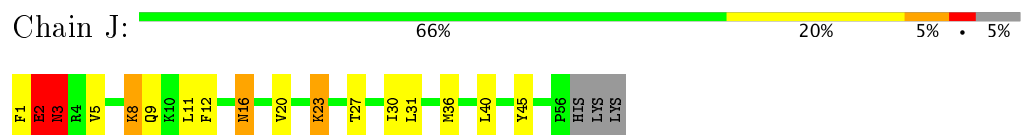
- Molecule 9: CYTOCHROME C OXIDASE



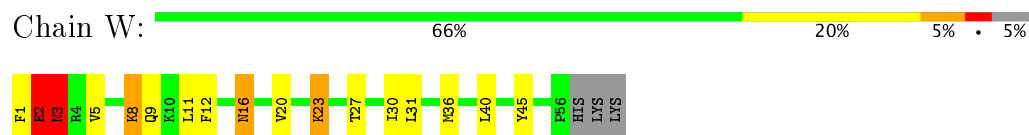
- Molecule 9: CYTOCHROME C OXIDASE



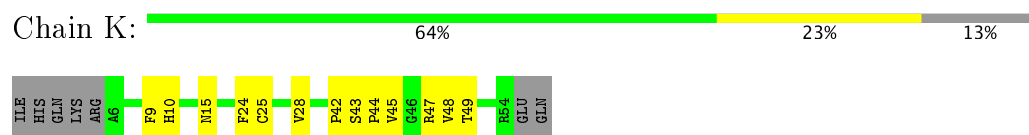
- Molecule 10: CYTOCHROME C OXIDASE



- Molecule 10: CYTOCHROME C OXIDASE



- Molecule 11: CYTOCHROME C OXIDASE




- Molecule 11: CYTOCHROME C OXIDASE

Chain X:  64% 23% 13%




- Molecule 12: CYTOCHROME C OXIDASE

Chain L:  81% 17% .



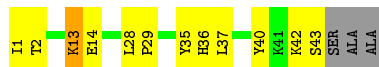
- Molecule 12: CYTOCHROME C OXIDASE

Chain Y:  81% 17% .



- Molecule 13: CYTOCHROME C OXIDASE

Chain M:  67% 24% . 7%



- Molecule 13: CYTOCHROME C OXIDASE

Chain Z:  70% 22% . 7%



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	189.10Å 210.50Å 178.60Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	10.00 – 2.80	Depositor
% Data completeness (in resolution range)	88.7 (10.00-2.80)	Depositor
R_{merge}	0.78	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR 3.1	Depositor
R, R_{free}	0.201 , 0.252	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	28722	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: MG, ZN, CU, HEA

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	0.60	0/4164	0.76	1/5688 (0.0%)
1	N	0.60	0/4164	0.76	1/5688 (0.0%)
2	B	0.57	0/1868	0.79	0/2544
2	O	0.57	0/1868	0.79	0/2544
3	C	0.56	0/2211	0.68	0/3023
3	P	0.56	0/2211	0.68	0/3023
4	D	0.56	0/1229	0.64	1/1658 (0.1%)
4	Q	0.56	0/1229	0.64	1/1658 (0.1%)
5	E	0.50	0/898	0.66	0/1218
5	R	0.50	0/898	0.66	0/1218
6	F	0.56	0/765	0.81	0/1038
6	S	0.56	0/765	0.81	0/1038
7	G	0.54	0/699	0.73	1/950 (0.1%)
7	T	0.55	0/699	0.73	1/950 (0.1%)
8	H	0.55	0/648	0.73	0/877
8	U	0.55	0/648	0.73	0/877
9	I	0.60	0/611	0.64	0/810
9	V	0.60	0/611	0.64	0/810
10	J	0.61	0/451	0.72	0/610
10	W	0.61	0/451	0.72	0/610
11	K	0.57	0/398	0.66	0/546
11	X	0.57	0/398	0.66	0/546
12	L	0.63	0/399	0.62	0/534
12	Y	0.63	0/399	0.62	0/534
13	M	0.51	0/345	0.65	0/470
13	Z	0.51	0/345	0.65	0/470
All	All	0.57	0/29372	0.72	6/39932 (0.0%)

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	Q	133	GLY	N-CA-C	6.44	129.21	113.10
4	D	133	GLY	N-CA-C	6.44	129.21	113.10
7	G	5	LYS	N-CA-C	5.22	125.10	111.00
7	T	5	LYS	N-CA-C	5.22	125.10	111.00
1	A	435	GLY	N-CA-C	5.18	126.04	113.10
1	N	435	GLY	N-CA-C	5.17	126.02	113.10

There are no chirality outliers.

There are no planarity outliers.

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	4025	0	4003	91	0
1	N	4025	0	4003	90	0
2	B	1822	0	1834	72	0
2	O	1822	0	1834	73	0
3	C	2124	0	2044	50	0
3	P	2124	0	2044	51	0
4	D	1195	0	1183	32	0
4	Q	1195	0	1183	33	0
5	E	878	0	868	22	0
5	R	878	0	868	20	0
6	F	748	0	728	27	0
6	S	748	0	728	24	0
7	G	672	0	645	40	0
7	T	672	0	645	42	0
8	H	628	0	582	29	0
8	U	628	0	582	30	0
9	I	598	0	612	15	0
9	V	598	0	612	14	0
10	J	441	0	439	14	0
10	W	441	0	439	14	0
11	K	384	0	366	9	0
11	X	384	0	366	10	0
12	L	386	0	388	8	0
12	Y	386	0	388	8	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
13	M	335	0	352	12	0
13	Z	335	0	352	10	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	120	0	108	8	0
16	N	120	0	108	6	0
17	F	1	0	0	0	0
17	S	1	0	0	0	0
All	All	28722	0	28304	697	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 12.

All (697) 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)
7:T:5:LYS:HZ2	7:T:6:GLY:H	1.13	0.97
1:A:365:ILE:HD12	2:B:87:MET:HE1	1.52	0.90
1:N:365:ILE:HD12	2:O:87:MET:HE1	1.51	0.90
7:G:5:LYS:HZ2	7:G:6:GLY:H	1.14	0.89
3:P:112:LEU:HG	3:P:118:PRO:HB3	1.55	0.88
3:C:112:LEU:HG	3:C:118:PRO:HB3	1.55	0.87
2:B:78:LEU:HB2	2:B:79:PRO:HD3	1.58	0.86
2:O:78:LEU:HB2	2:O:79:PRO:HD3	1.58	0.83
7:G:5:LYS:HB3	7:G:5:LYS:HZ2	1.45	0.82
1:A:193:VAL:HG11	7:T:5:LYS:O	1.78	0.81
1:A:281:GLY:HA3	7:T:5:LYS:HE3	1.62	0.81
7:G:5:LYS:NZ	7:G:6:GLY:H	1.78	0.81
7:T:5:LYS:NZ	7:T:6:GLY:H	1.78	0.80
7:T:5:LYS:HZ2	7:T:5:LYS:HB3	1.45	0.80
8:H:39:CYS:SG	8:H:53:CYS:HB3	2.22	0.79
8:U:39:CYS:SG	8:U:53:CYS:HB3	2.22	0.79
8:H:75:ARG:HG2	8:H:75:ARG:HH11	1.48	0.79
8:H:39:CYS:SG	8:H:53:CYS:CB	2.70	0.79
7:G:5:LYS:O	1:N:193:VAL:HG11	1.82	0.79
8:U:39:CYS:SG	8:U:53:CYS:CB	2.70	0.79
8:U:39:CYS:HG	8:U:53:CYS:CB	1.96	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:U:75:ARG:HH11	8:U:75:ARG:HG2	1.48	0.78
7:G:5:LYS:HE3	1:N:281:GLY:HA3	1.65	0.77
8:H:49:ASP:HB2	8:U:49:ASP:HB2	1.67	0.76
2:B:59:GLN:H	2:B:62:GLU:HG3	1.51	0.75
2:O:59:GLN:H	2:O:62:GLU:HG3	1.51	0.75
2:O:184:LEU:HD11	2:O:211:LEU:HD21	1.69	0.75
2:B:184:LEU:HD11	2:B:211:LEU:HD21	1.69	0.75
3:P:187:THR:HG21	7:T:68:THR:HG21	1.68	0.74
7:G:54:ARG:N	7:G:54:ARG:HD3	2.04	0.73
7:T:54:ARG:N	7:T:54:ARG:HD3	2.04	0.73
4:D:23:PRO:O	5:E:66:ARG:HD3	1.89	0.73
3:C:221:ARG:HE	3:C:226:HIS:HD1	1.37	0.73
3:C:187:THR:HG21	7:G:68:THR:HG21	1.68	0.73
4:Q:23:PRO:O	5:R:66:ARG:HD3	1.89	0.73
4:D:147:LYS:HE3	4:D:147:LYS:HA	1.72	0.72
1:A:194:LEU:HD11	7:T:5:LYS:HD3	1.72	0.71
3:C:12:ASN:HD22	10:J:20:VAL:H	1.39	0.71
8:H:48:GLY:O	8:U:51:SER:HB3	1.91	0.71
2:B:13:THR:HB	2:B:168:LEU:HD23	1.74	0.70
3:P:221:ARG:HE	3:P:226:HIS:HD1	1.37	0.70
2:O:13:THR:HB	2:O:168:LEU:HD23	1.74	0.70
8:H:51:SER:HB3	8:U:48:GLY:O	1.92	0.70
3:P:12:ASN:HD22	10:W:20:VAL:H	1.39	0.70
1:A:11:ASN:HD22	1:A:14:ASP:H	1.39	0.70
3:P:156:ARG:HG3	3:P:156:ARG:HH11	1.56	0.70
3:C:156:ARG:HH11	3:C:156:ARG:HG3	1.56	0.69
4:Q:147:LYS:HE3	4:Q:147:LYS:HA	1.71	0.69
1:N:11:ASN:HD22	1:N:14:ASP:H	1.39	0.69
2:B:186:SER:HB3	2:B:213:LEU:HD22	1.75	0.69
2:O:186:SER:HB3	2:O:213:LEU:HD22	1.75	0.68
7:G:5:LYS:HD3	1:N:194:LEU:HD11	1.76	0.68
3:C:154:GLY:HA2	6:F:6:VAL:HG22	1.75	0.68
5:E:43:PRO:HB2	5:E:48:ILE:HD11	1.76	0.68
3:P:154:GLY:HA2	6:S:6:VAL:HG22	1.75	0.68
8:H:38:ARG:HG2	8:H:85:ILE:HG23	1.77	0.67
5:R:43:PRO:HB2	5:R:48:ILE:HD11	1.76	0.67
8:U:38:ARG:HG2	8:U:85:ILE:HG23	1.77	0.67
12:L:19:TRP:HZ2	13:M:14:GLU:HG2	1.59	0.67
1:N:151:HIS:O	1:N:155:VAL:HG23	1.95	0.67
3:P:209:ILE:O	3:P:213:THR:HG23	1.95	0.67
12:Y:19:TRP:HZ2	13:Z:14:GLU:HG2	1.58	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:151:HIS:O	1:A:155:VAL:HG23	1.95	0.67
1:N:298:ASP:HB2	1:N:301:THR:HG23	1.77	0.67
3:C:209:ILE:O	3:C:213:THR:HG23	1.95	0.66
1:A:298:ASP:HB2	1:A:301:THR:HG23	1.77	0.66
3:C:160:LEU:HD13	3:C:222:GLN:HG2	1.78	0.66
3:C:101:PHE:HZ	3:C:260:GLY:HA3	1.61	0.66
4:D:24:LEU:H	5:E:34:ASN:HD21	1.44	0.65
2:O:120:SER:OG	2:O:138:VAL:HG21	1.97	0.65
12:L:45:LEU:HD21	13:M:40:TYR:HA	1.78	0.65
1:A:172:LYS:NZ	1:A:172:LYS:HB2	2.12	0.65
3:P:101:PHE:HZ	3:P:260:GLY:HA3	1.61	0.65
2:B:120:SER:OG	2:B:138:VAL:HG21	1.97	0.65
4:Q:24:LEU:H	5:R:34:ASN:HD21	1.44	0.65
1:N:195:LEU:HD23	1:N:245:ILE:HD13	1.79	0.64
12:Y:45:LEU:HD21	13:Z:40:TYR:HA	1.78	0.64
1:N:172:LYS:HB2	1:N:172:LYS:NZ	2.12	0.64
3:P:160:LEU:HD13	3:P:222:GLN:HG2	1.78	0.64
7:G:5:LYS:HG3	1:N:282:PHE:N	2.13	0.64
1:A:195:LEU:HD23	1:A:245:ILE:HD13	1.79	0.63
2:B:122:MET:HB2	2:B:208:PRO:HD2	1.81	0.62
4:Q:16:TYR:CE1	4:Q:25:PRO:HG3	2.34	0.62
8:U:57:ARG:HH11	8:U:57:ARG:HB3	1.64	0.62
13:Z:13:LYS:HD3	13:Z:13:LYS:H	1.65	0.62
4:D:16:TYR:CE1	4:D:25:PRO:HG3	2.34	0.62
1:A:11:ASN:HD21	1:A:13:LYS:HB2	1.64	0.62
2:B:16:ILE:HD12	2:B:17:MET:N	2.15	0.62
5:E:82:TYR:HB3	5:E:83:PRO:HD3	1.81	0.62
1:N:11:ASN:HD21	1:N:13:LYS:HB2	1.64	0.61
1:A:282:PHE:N	7:T:5:LYS:HG3	2.15	0.61
6:F:13:ALA:O	6:F:18:ARG:HD2	2.00	0.61
2:O:16:ILE:HD12	2:O:17:MET:N	2.15	0.61
8:H:53:CYS:HA	8:U:46:LYS:NZ	2.15	0.61
6:S:13:ALA:O	6:S:18:ARG:HD2	2.00	0.61
3:C:192:VAL:O	3:C:196:THR:HB	2.01	0.61
13:M:13:LYS:H	13:M:13:LYS:HD3	1.65	0.61
3:C:6:HIS:HD2	3:C:8:TYR:H	1.49	0.61
3:P:154:GLY:HA2	6:S:6:VAL:CG2	2.30	0.61
1:A:128:VAL:O	1:A:128:VAL:HG22	1.99	0.61
8:H:39:CYS:HG	8:H:53:CYS:HG	0.67	0.61
2:B:59:GLN:HA	2:B:62:GLU:HB2	1.83	0.60
1:A:92:MET:CE	1:A:167:THR:HG21	2.31	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:H:57:ARG:HH11	8:H:57:ARG:HB3	1.64	0.60
3:P:192:VAL:O	3:P:196:THR:HB	2.01	0.60
3:C:154:GLY:HA2	6:F:6:VAL:CG2	2.30	0.60
2:O:122:MET:HB2	2:O:208:PRO:HD2	1.81	0.60
2:O:59:GLN:HA	2:O:62:GLU:HB2	1.83	0.60
8:H:46:LYS:NZ	8:U:53:CYS:HA	2.16	0.60
6:F:62:CYS:SG	6:F:84:SER:OG	2.60	0.60
1:N:128:VAL:HG22	1:N:128:VAL:O	1.99	0.60
16:A:516:HEA:HMC1	16:A:516:HEA:HBC1	1.83	0.60
3:C:119:THR:O	7:G:52:HIS:HE1	1.84	0.60
13:M:28:LEU:HB2	13:M:29:PRO:HD3	1.84	0.60
3:P:148:HIS:O	3:P:152:MET:HG3	2.01	0.60
2:B:226:MET:O	2:B:227:LEU:HB2	2.02	0.60
1:N:92:MET:CE	1:N:167:THR:HG21	2.31	0.60
13:Z:28:LEU:HB2	13:Z:29:PRO:HD3	1.84	0.60
3:C:148:HIS:O	3:C:152:MET:HG3	2.01	0.60
6:S:62:CYS:SG	6:S:84:SER:OG	2.60	0.60
5:R:82:TYR:HB3	5:R:83:PRO:HD3	1.81	0.60
1:N:92:MET:HE2	1:N:167:THR:HG21	1.84	0.59
3:P:6:HIS:HD2	3:P:8:TYR:H	1.49	0.59
16:N:516:HEA:HBC1	16:N:516:HEA:HMC1	1.83	0.59
2:O:193:TYR:HE1	4:Q:126:MET:HE1	1.68	0.59
1:A:92:MET:HE2	1:A:167:THR:HG21	1.84	0.59
1:A:194:LEU:CD1	7:T:5:LYS:HD3	2.31	0.59
2:O:226:MET:O	2:O:227:LEU:HB2	2.02	0.59
2:O:179:LEU:HD21	8:U:65:PRO:HD3	1.85	0.58
3:P:119:THR:O	7:T:52:HIS:HE1	1.85	0.58
12:Y:46:LYS:O	12:Y:47:LYS:HB2	2.02	0.58
2:B:179:LEU:HD21	8:H:65:PRO:HD3	1.85	0.58
12:L:46:LYS:O	12:L:47:LYS:HB2	2.02	0.58
8:U:57:ARG:HG3	8:U:60:TYR:CE2	2.38	0.58
11:X:43:SER:OG	11:X:45:VAL:HG12	2.03	0.58
11:K:43:SER:OG	11:K:45:VAL:HG12	2.03	0.58
1:A:184:PHE:H	1:A:256:HIS:HE1	1.52	0.58
1:A:404:THR:O	1:A:480:ARG:NH1	2.37	0.58
2:B:193:TYR:HE1	4:D:126:MET:HE1	1.69	0.58
3:C:203:PHE:O	3:C:207:HIS:HD2	1.88	0.57
8:H:57:ARG:HG3	8:H:60:TYR:CE2	2.38	0.57
10:J:3:ASN:C	10:J:3:ASN:HD22	2.08	0.57
7:G:5:LYS:HG3	1:N:281:GLY:C	2.25	0.57
1:N:404:THR:O	1:N:480:ARG:NH1	2.37	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:281:GLY:C	7:T:5:LYS:HG3	2.25	0.57
12:L:41:ARG:HG3	13:M:40:TYR:CE2	2.40	0.57
2:O:132:GLU:HB3	2:O:137:GLU:HG3	1.87	0.57
2:B:189:PRO:HD2	9:I:54:TYR:OH	2.05	0.57
1:A:95:PRO:HB2	3:C:11:VAL:HG13	1.87	0.57
8:H:71:THR:O	8:H:75:ARG:HD3	2.05	0.57
8:U:39:CYS:HG	8:U:53:CYS:HG	0.79	0.57
12:Y:41:ARG:HG3	13:Z:40:TYR:CE2	2.40	0.57
8:U:71:THR:O	8:U:75:ARG:HD3	2.05	0.56
7:T:5:LYS:HD2	7:T:6:GLY:N	2.21	0.56
1:N:184:PHE:H	1:N:256:HIS:HE1	1.52	0.56
3:P:203:PHE:O	3:P:207:HIS:HD2	1.88	0.56
10:W:3:ASN:HD22	10:W:3:ASN:C	2.08	0.56
3:C:42:LEU:HD13	10:J:45:TYR:CD2	2.41	0.56
1:N:95:PRO:HB2	3:P:11:VAL:HG13	1.87	0.56
5:R:78:HIS:ND1	9:V:12:LEU:HD22	2.21	0.56
1:A:273:MET:HG3	1:A:319:LYS:HZ1	1.71	0.56
5:E:78:HIS:ND1	9:I:12:LEU:HD22	2.21	0.56
7:G:5:LYS:HD3	1:N:194:LEU:CD1	2.35	0.56
2:O:189:PRO:HD2	9:V:54:TYR:OH	2.05	0.56
2:B:132:GLU:HB3	2:B:137:GLU:HG3	1.87	0.55
3:P:42:LEU:HD13	10:W:45:TYR:CD2	2.41	0.55
2:B:111:THR:HG21	8:H:66:ILE:HD11	1.88	0.55
8:U:75:ARG:HG2	8:U:75:ARG:NH1	2.19	0.55
7:G:5:LYS:HD2	7:G:6:GLY:N	2.21	0.55
7:G:6:GLY:HA3	1:N:278:MET:HG2	1.87	0.55
3:P:222:GLN:HE21	3:P:222:GLN:HA	1.72	0.55
4:D:108:PRO:HG2	4:D:111:PHE:CE2	2.42	0.55
2:O:203:ASN:HD22	2:O:206:PHE:HD2	1.55	0.55
3:C:222:GLN:HE21	3:C:222:GLN:HA	1.72	0.55
6:S:82:CYS:N	6:S:86:GLY:O	2.40	0.55
2:O:111:THR:HG21	8:U:66:ILE:HD11	1.88	0.55
3:C:151:LEU:HB2	3:C:159:MET:HG3	1.89	0.54
1:A:281:GLY:CA	7:T:5:LYS:HE3	2.36	0.54
1:A:407:ASP:O	1:A:411:LYS:HG3	2.07	0.54
6:F:82:CYS:N	6:F:86:GLY:O	2.41	0.54
1:N:407:ASP:O	1:N:411:LYS:HG3	2.07	0.54
4:Q:108:PRO:HG2	4:Q:111:PHE:CE2	2.42	0.54
2:B:203:ASN:HD22	2:B:206:PHE:HD2	1.55	0.54
2:O:145:PRO:CG	2:O:148:MET:HE2	2.38	0.54
3:C:156:ARG:NH1	3:C:156:ARG:HG3	2.23	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:O:165:VAL:HB	2:O:170:LEU:HD12	1.90	0.54
2:O:136:LEU:HB3	2:O:193:TYR:HD2	1.73	0.54
4:Q:40:LEU:HD13	4:Q:59:LEU:HD13	1.90	0.54
6:S:55:LYS:HA	6:S:74:LEU:O	2.08	0.54
4:Q:52:SER:OG	4:Q:55:GLU:HG3	2.08	0.54
5:R:41:LEU:HD12	5:R:41:LEU:O	2.08	0.54
1:A:398:PRO:O	1:A:498:CYS:HB3	2.08	0.53
2:B:165:VAL:HB	2:B:170:LEU:HD12	1.90	0.53
5:E:80:GLU:H	5:E:80:GLU:CD	2.10	0.53
6:F:55:LYS:HA	6:F:74:LEU:O	2.08	0.53
3:P:156:ARG:HG3	3:P:156:ARG:NH1	2.22	0.53
7:T:2:SER:OG	7:T:3:ALA:N	2.41	0.53
6:F:48:LEU:O	6:F:50:PRO:HD3	2.08	0.53
9:V:63:MET:HB3	9:V:68:ILE:HD11	1.89	0.53
4:D:40:LEU:HD13	4:D:59:LEU:HD13	1.90	0.53
9:I:63:MET:HB3	9:I:68:ILE:HD11	1.89	0.53
4:Q:67:SER:OG	4:Q:70:GLU:HG3	2.08	0.53
4:D:23:PRO:HD2	5:E:34:ASN:HD22	1.74	0.53
2:B:145:PRO:CG	2:B:148:MET:HE2	2.38	0.53
5:E:41:LEU:O	5:E:41:LEU:HD12	2.08	0.53
1:N:11:ASN:ND2	1:N:14:ASP:H	2.07	0.53
6:S:48:LEU:O	6:S:50:PRO:HD3	2.08	0.53
6:S:51:SER:HB2	6:S:91:LEU:HD11	1.90	0.53
5:R:80:GLU:CD	5:R:80:GLU:H	2.10	0.53
2:B:102:HIS:O	2:B:104:TRP:N	2.42	0.53
2:B:136:LEU:HB3	2:B:193:TYR:HD2	1.73	0.53
4:D:52:SER:OG	4:D:55:GLU:HG3	2.08	0.53
6:F:51:SER:HB2	6:F:91:LEU:HD11	1.90	0.53
7:G:8:HIS:O	7:G:10:GLY:N	2.42	0.53
7:G:2:SER:OG	7:G:3:ALA:N	2.41	0.53
3:P:151:LEU:HB2	3:P:159:MET:HG3	1.89	0.53
4:Q:128:VAL:O	4:Q:134:PHE:HB3	2.09	0.53
1:N:302:ARG:O	1:N:306:THR:HG23	2.09	0.53
7:T:8:HIS:O	7:T:10:GLY:N	2.42	0.53
1:A:278:MET:HG2	7:T:6:GLY:HA3	1.91	0.53
1:A:302:ARG:O	1:A:306:THR:HG23	2.09	0.52
4:D:128:VAL:O	4:D:134:PHE:HB3	2.09	0.52
6:S:8:THR:OG1	6:S:11:GLU:HG2	2.10	0.52
1:A:268:PHE:CZ	2:B:58:ALA:HA	2.44	0.52
1:N:268:PHE:CZ	2:O:58:ALA:HA	2.44	0.52
7:T:42:ARG:HH11	7:T:42:ARG:HG3	1.75	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:225:GLY:HA3	3:C:112:LEU:HD13	1.92	0.52
2:B:33:LEU:HD23	9:I:28:SER:HB2	1.91	0.52
2:O:104:TRP:CG	2:O:203:ASN:HB2	2.44	0.52
3:C:18:LEU:HD22	3:C:22:LEU:HD22	1.92	0.52
4:D:67:SER:OG	4:D:70:GLU:HG3	2.08	0.52
1:N:398:PRO:O	1:N:498:CYS:HB3	2.08	0.52
4:Q:23:PRO:HD2	5:R:34:ASN:HD22	1.74	0.52
5:R:63:SER:O	5:R:67:ILE:HG13	2.10	0.52
5:E:63:SER:O	5:E:67:ILE:HG13	2.10	0.52
2:O:68:LEU:HB3	2:O:69:PRO:HD3	1.91	0.52
2:B:100:MET:CE	2:B:157:GLU:HG3	2.40	0.52
2:B:68:LEU:HB3	2:B:69:PRO:HD3	1.91	0.52
3:C:119:THR:O	7:G:52:HIS:CE1	2.63	0.52
1:N:240:HIS:HB3	1:N:241:PRO:HD3	1.92	0.52
2:B:104:TRP:CG	2:B:203:ASN:HB2	2.44	0.52
6:F:8:THR:OG1	6:F:11:GLU:HG2	2.10	0.52
2:O:33:LEU:HD23	9:V:28:SER:HB2	1.91	0.52
2:B:186:SER:CB	2:B:213:LEU:HD22	2.40	0.51
8:H:75:ARG:HG2	8:H:75:ARG:NH1	2.19	0.51
2:O:1:MET:SD	2:O:133:LEU:CD1	2.98	0.51
1:A:250:GLY:O	1:A:254:ILE:HG12	2.11	0.51
2:O:102:HIS:O	2:O:104:TRP:N	2.42	0.51
7:G:42:ARG:HG3	7:G:42:ARG:HH11	1.75	0.51
7:G:5:LYS:NZ	7:G:6:GLY:N	2.54	0.51
2:O:100:MET:CE	2:O:157:GLU:HG3	2.40	0.51
1:A:11:ASN:ND2	1:A:14:ASP:H	2.07	0.51
1:N:225:GLY:HA3	3:P:112:LEU:HD13	1.92	0.51
4:Q:40:LEU:HD22	4:Q:59:LEU:HD13	1.91	0.51
1:A:240:HIS:HB3	1:A:241:PRO:HD3	1.92	0.51
1:A:281:GLY:HA3	7:T:5:LYS:CE	2.39	0.51
2:B:139:ASP:OD1	2:B:140:ASN:N	2.44	0.51
1:N:75:ILE:O	1:N:79:GLY:HA3	2.11	0.51
3:P:119:THR:O	7:T:52:HIS:CE1	2.63	0.51
2:B:59:GLN:N	2:B:62:GLU:HG3	2.24	0.51
7:G:5:LYS:HE3	1:N:281:GLY:CA	2.38	0.51
1:A:75:ILE:O	1:A:79:GLY:HA3	2.11	0.51
2:B:136:LEU:HB3	2:B:193:TYR:CD2	2.46	0.51
4:D:40:LEU:HD22	4:D:59:LEU:HD13	1.92	0.51
2:O:136:LEU:HB3	2:O:193:TYR:CD2	2.46	0.51
1:N:377:PHE:HA	1:N:380:VAL:HG22	1.93	0.51
1:A:377:PHE:HA	1:A:380:VAL:HG22	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:1:MET:SD	2:B:133:LEU:CD1	2.98	0.51
1:A:40:GLU:HG2	1:A:54:TYR:CD1	2.47	0.50
4:Q:64:PHE:CE1	5:R:66:ARG:HD2	2.47	0.50
1:A:184:PHE:H	1:A:256:HIS:CE1	2.29	0.50
1:N:273:MET:HG3	1:N:319:LYS:HZ1	1.75	0.50
2:O:139:ASP:OD1	2:O:140:ASN:N	2.44	0.50
7:T:5:LYS:C	7:T:5:LYS:HD2	2.32	0.50
1:A:509:THR:HG1	6:F:71:TRP:HZ3	1.58	0.50
2:O:59:GLN:N	2:O:62:GLU:HG3	2.24	0.50
3:P:18:LEU:HD22	3:P:22:LEU:HD22	1.92	0.50
6:S:37:LYS:N	6:S:37:LYS:HD3	2.27	0.50
1:A:168:ILE:HG21	1:A:189:MET:HG3	1.94	0.50
7:G:42:ARG:NH1	7:G:74:ARG:HH21	2.10	0.50
1:N:184:PHE:H	1:N:256:HIS:CE1	2.29	0.50
1:A:145:LEU:HD21	3:C:32:THR:HG21	1.93	0.50
2:B:13:THR:HG22	2:B:13:THR:O	2.12	0.50
4:D:64:PHE:CE1	5:E:66:ARG:HD2	2.47	0.50
1:N:172:LYS:HZ2	1:N:172:LYS:HB2	1.75	0.50
2:B:16:ILE:HD12	2:B:17:MET:H	1.78	0.49
6:F:37:LYS:N	6:F:37:LYS:HD3	2.27	0.49
2:O:186:SER:HB3	2:O:213:LEU:CD2	2.42	0.49
3:P:42:LEU:HD13	10:W:45:TYR:HD2	1.77	0.49
7:T:5:LYS:NZ	7:T:6:GLY:N	2.54	0.49
1:N:145:LEU:HD21	3:P:32:THR:HG21	1.93	0.49
1:N:365:ILE:HD12	2:O:87:MET:CE	2.34	0.49
1:A:365:ILE:HD12	2:B:87:MET:CE	2.34	0.49
3:C:42:LEU:HD13	10:J:45:TYR:HD2	1.77	0.49
7:G:5:LYS:C	7:G:5:LYS:HD2	2.32	0.49
2:O:13:THR:O	2:O:13:THR:HG22	2.11	0.49
1:N:250:GLY:O	1:N:254:ILE:HG12	2.11	0.49
1:N:40:GLU:HG2	1:N:54:TYR:CD1	2.47	0.49
7:G:42:ARG:O	7:G:42:ARG:HD2	2.12	0.49
1:N:509:THR:HG1	6:S:71:TRP:HZ3	1.61	0.49
1:A:172:LYS:HZ2	1:A:172:LYS:HB2	1.78	0.49
6:F:49:VAL:HG21	6:F:74:LEU:HD12	1.95	0.49
1:N:266:GLU:HB2	1:N:267:PRO:HD2	1.94	0.49
2:O:90:ILE:HD12	2:O:90:ILE:H	1.77	0.49
7:T:42:ARG:O	7:T:42:ARG:HD2	2.13	0.49
7:T:44:ARG:HD2	7:T:74:ARG:O	2.12	0.49
8:H:46:LYS:HZ3	8:U:53:CYS:HA	1.76	0.49
1:A:187:SER:HB2	1:A:277:MET:HE1	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:90:ILE:H	2:B:90:ILE:HD12	1.77	0.49
1:N:508:PRO:HG3	3:P:6:HIS:HB3	1.94	0.49
8:U:24:ASN:ND2	8:U:26:THR:H	2.11	0.49
1:A:87:ILE:O	1:A:173:PRO:HD3	2.13	0.49
7:G:44:ARG:HD2	7:G:74:ARG:O	2.12	0.49
7:G:5:LYS:CE	1:N:281:GLY:HA3	2.41	0.49
2:O:186:SER:CB	2:O:213:LEU:HD22	2.40	0.49
4:Q:68:PHE:HA	4:Q:71:MET:HG2	1.95	0.48
1:A:508:PRO:HG3	3:C:6:HIS:HB3	1.94	0.48
3:C:65:SER:HB3	3:C:71:HIS:CE1	2.48	0.48
7:T:42:ARG:NH1	7:T:74:ARG:HH21	2.10	0.48
1:N:476:PHE:CD1	12:Y:15:VAL:HG21	2.48	0.48
1:A:115:SER:O	1:A:121:GLY:HA2	2.14	0.48
1:A:397:PHE:HB3	1:A:398:PRO:HD3	1.95	0.48
1:N:197:LEU:O	3:P:92:LEU:HD22	2.13	0.48
2:B:186:SER:HB3	2:B:213:LEU:CD2	2.42	0.48
1:A:449:MET:SD	2:B:5:MET:HE2	2.54	0.48
1:N:11:ASN:ND2	1:N:13:LYS:HB2	2.28	0.48
1:N:87:ILE:O	1:N:173:PRO:HD3	2.13	0.48
2:O:59:GLN:O	2:O:62:GLU:HB2	2.13	0.48
3:P:65:SER:HB3	3:P:71:HIS:CE1	2.48	0.48
2:B:22:HIS:CE1	9:I:44:LYS:HD3	2.49	0.48
8:H:24:ASN:ND2	8:H:26:THR:H	2.11	0.48
9:V:39:VAL:O	9:V:42:LYS:HE2	2.14	0.48
1:N:168:ILE:HG21	1:N:189:MET:HG3	1.94	0.48
10:J:16:ASN:H	10:J:16:ASN:ND2	2.12	0.48
1:A:476:PHE:CD1	12:L:15:VAL:HG21	2.48	0.48
1:A:449:MET:SD	2:B:5:MET:CE	3.02	0.48
1:N:397:PHE:HB3	1:N:398:PRO:HD3	1.95	0.48
4:D:68:PHE:HA	4:D:71:MET:HG2	1.95	0.48
1:N:449:MET:SD	2:O:5:MET:CE	3.02	0.48
9:I:39:VAL:O	9:I:42:LYS:HE2	2.14	0.48
1:A:197:LEU:O	3:C:92:LEU:HD22	2.13	0.47
10:J:8:LYS:HB3	10:J:8:LYS:NZ	2.29	0.47
6:S:49:VAL:HG21	6:S:74:LEU:HD12	1.95	0.47
2:B:63:THR:O	2:B:66:THR:HG22	2.15	0.47
10:W:8:LYS:HB3	10:W:8:LYS:NZ	2.29	0.47
2:B:59:GLN:O	2:B:62:GLU:HB2	2.13	0.47
6:F:31:TYR:HE1	6:F:98:HIS:HE1	1.62	0.47
3:P:80:ARG:O	3:P:84:ILE:HG12	2.14	0.47
2:O:4:PRO:HB2	11:X:43:SER:HA	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:266:GLU:HB2	1:A:267:PRO:HD2	1.94	0.47
8:H:58:ARG:HD2	8:H:58:ARG:HA	1.74	0.47
2:O:22:HIS:CE1	9:V:44:LYS:HD3	2.49	0.47
1:A:69:MET:HE1	1:A:70:VAL:HG23	1.96	0.47
2:B:5:MET:CE	11:K:42:PRO:HA	2.45	0.47
2:B:83:ILE:O	2:B:87:MET:HG3	2.14	0.47
7:G:67:HIS:CD2	7:G:78:LEU:HD11	2.50	0.47
2:B:4:PRO:HB2	11:K:43:SER:HA	1.96	0.47
2:O:63:THR:O	2:O:66:THR:HG22	2.15	0.47
2:O:83:ILE:O	2:O:87:MET:HG3	2.14	0.47
3:P:164:PHE:O	3:P:168:THR:HG23	2.14	0.47
4:Q:41:LYS:HD3	4:Q:62:LEU:HD23	1.96	0.47
7:T:67:HIS:CD2	7:T:78:LEU:HD11	2.50	0.47
9:V:21:ILE:HA	9:V:21:ILE:HD12	1.82	0.47
3:C:164:PHE:O	3:C:168:THR:HG23	2.14	0.47
3:C:80:ARG:O	3:C:84:ILE:HG12	2.14	0.47
1:N:115:SER:O	1:N:121:GLY:HA2	2.14	0.47
10:W:16:ASN:H	10:W:16:ASN:ND2	2.12	0.47
4:D:23:PRO:HB3	5:E:70:VAL:CG2	2.45	0.47
2:O:16:ILE:HD12	2:O:17:MET:H	1.78	0.47
4:Q:23:PRO:HB3	5:R:70:VAL:CG2	2.45	0.47
4:Q:33:LEU:HD22	4:Q:37:GLN:HB3	1.97	0.47
8:U:60:TYR:CD1	8:U:60:TYR:C	2.88	0.47
2:O:5:MET:CE	11:X:42:PRO:HA	2.45	0.47
1:A:197:LEU:HD11	7:T:4:ALA:HB3	1.96	0.47
1:A:195:LEU:CD2	1:A:245:ILE:HD13	2.45	0.47
2:B:108:TYR:N	2:B:108:TYR:CD1	2.83	0.47
4:Q:108:PRO:HG2	4:Q:111:PHE:CD2	2.50	0.47
10:W:11:LEU:O	10:W:11:LEU:HD23	2.15	0.47
1:A:306:THR:HB	1:A:359:ALA:O	2.15	0.46
10:J:11:LEU:O	10:J:11:LEU:HD23	2.15	0.46
2:O:108:TYR:CD1	2:O:108:TYR:N	2.83	0.46
1:A:484:THR:HB	13:M:2:THR:OG1	2.16	0.46
4:D:41:LYS:HD3	4:D:62:LEU:HD23	1.97	0.46
4:D:102:TYR:CD2	13:M:35:TYR:HE1	2.33	0.46
1:N:507:GLU:O	1:N:508:PRO:O	2.32	0.46
4:Q:33:LEU:HD13	4:Q:41:LYS:HG3	1.97	0.46
2:B:191:LEU:HB2	4:D:126:MET:HE1	1.97	0.46
4:D:33:LEU:HD13	4:D:41:LYS:HG3	1.97	0.46
1:N:273:MET:HG3	1:N:319:LYS:NZ	2.31	0.46
2:O:58:ALA:O	2:O:60:GLU:HG3	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:1:MET:SD	2:B:133:LEU:HD11	2.56	0.46
4:D:108:PRO:HG2	4:D:111:PHE:CD2	2.50	0.46
8:H:60:TYR:C	8:H:60:TYR:CD1	2.88	0.46
6:S:31:TYR:HE1	6:S:98:HIS:HE1	1.62	0.46
1:A:186:TRP:HH2	7:T:8:HIS:HB3	1.81	0.46
1:A:299:VAL:HA	1:A:302:ARG:NH1	2.31	0.46
4:D:33:LEU:HD22	4:D:37:GLN:HB3	1.97	0.46
7:G:4:ALA:HB3	1:N:197:LEU:HD11	1.98	0.46
11:K:9:PHE:HD1	11:K:10:HIS:HD2	1.64	0.46
1:N:240:HIS:NE2	1:N:244:TYR:CE2	2.81	0.46
1:N:484:THR:HB	13:Z:2:THR:OG1	2.15	0.46
4:D:66:GLU:O	5:E:66:ARG:NH2	2.49	0.46
13:Z:1:ILE:HG23	13:Z:1:ILE:O	2.15	0.46
1:A:273:MET:HG3	1:A:319:LYS:NZ	2.31	0.46
2:B:76:ILE:O	2:B:79:PRO:HD2	2.16	0.46
6:F:40:SER:OG	6:F:45:ASP:HB3	2.16	0.46
1:N:306:THR:HB	1:N:359:ALA:O	2.15	0.46
2:O:78:LEU:HB2	2:O:79:PRO:CD	2.38	0.46
8:H:53:CYS:HA	8:U:46:LYS:HZ3	1.80	0.46
1:A:147:ILE:HD11	1:A:209:LEU:HD23	1.98	0.46
2:B:59:GLN:CA	2:B:62:GLU:HB2	2.46	0.46
1:N:195:LEU:CD2	1:N:245:ILE:HD13	2.45	0.46
1:N:435:GLY:O	1:N:437:PRO:HD3	2.15	0.46
6:S:40:SER:OG	6:S:45:ASP:HB3	2.16	0.46
1:A:240:HIS:NE2	1:A:244:TYR:CE2	2.81	0.46
11:X:9:PHE:HD1	11:X:10:HIS:HD2	1.64	0.46
1:A:240:HIS:O	1:A:243:VAL:HG22	2.16	0.45
7:G:8:HIS:HB3	1:N:186:TRP:HH2	1.80	0.45
1:N:32:ALA:HB3	12:Y:36:PRO:HG2	1.98	0.45
2:O:1:MET:SD	2:O:133:LEU:HD11	2.56	0.45
1:A:507:GLU:O	1:A:508:PRO:O	2.32	0.45
1:N:240:HIS:O	1:N:243:VAL:HG22	2.16	0.45
2:O:41:ILE:O	2:O:45:MET:HG2	2.16	0.45
6:S:98:HIS:N	6:S:98:HIS:ND1	2.64	0.45
4:Q:102:TYR:CD2	13:Z:35:TYR:HE1	2.33	0.45
1:A:32:ALA:HB3	12:L:36:PRO:HG2	1.98	0.45
2:B:41:ILE:O	2:B:45:MET:HG2	2.16	0.45
13:M:37:LEU:HD23	13:M:37:LEU:HA	1.83	0.45
7:G:5:LYS:HZ3	1:N:278:MET:HG2	1.81	0.45
6:S:33:ILE:HG22	6:S:34:LEU:HD12	1.98	0.45
1:A:435:GLY:O	1:A:437:PRO:HD3	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:158:ILE:O	1:A:162:ILE:HG13	2.17	0.45
6:F:98:HIS:N	6:F:98:HIS:ND1	2.64	0.45
2:O:76:ILE:O	2:O:79:PRO:HD2	2.16	0.45
16:A:515:HEA:CBC	16:A:515:HEA:HMC1	2.46	0.45
1:N:299:VAL:HA	1:N:302:ARG:NH1	2.31	0.45
2:O:122:MET:HB2	2:O:208:PRO:CD	2.47	0.45
6:S:31:TYR:HE1	6:S:98:HIS:CE1	2.34	0.45
4:Q:79:LYS:NZ	11:X:15:ASN:HD21	2.15	0.45
12:Y:35:ALA:HB3	12:Y:36:PRO:HD3	1.99	0.45
1:A:11:ASN:ND2	1:A:13:LYS:HB2	2.28	0.45
1:A:495:LEU:HD23	1:A:495:LEU:HA	1.68	0.45
10:J:2:GLU:CG	10:J:3:ASN:H	2.30	0.45
13:M:1:ILE:HG23	13:M:1:ILE:O	2.15	0.45
1:N:158:ILE:O	1:N:162:ILE:HG13	2.17	0.45
4:Q:40:LEU:HD22	4:Q:59:LEU:CD1	2.47	0.45
4:Q:66:GLU:O	5:R:66:ARG:NH2	2.49	0.45
1:A:431:LEU:HD21	1:A:450:TRP:HB2	1.99	0.45
4:D:40:LEU:HD22	4:D:59:LEU:CD1	2.47	0.45
16:N:515:HEA:HMC1	16:N:515:HEA:CBC	2.46	0.45
3:C:128:GLU:HB3	3:C:129:VAL:H	1.56	0.45
3:C:42:LEU:HA	3:C:42:LEU:HD12	1.66	0.45
1:A:513:LEU:HD12	1:A:513:LEU:HA	1.65	0.45
16:A:516:HEA:HHC	16:A:516:HEA:O11	2.17	0.45
2:B:162:SER:HB2	2:B:198:GLU:HB2	1.99	0.45
5:E:78:HIS:CE1	9:I:12:LEU:HD22	2.52	0.45
6:F:31:TYR:HE1	6:F:98:HIS:CE1	2.34	0.45
7:G:42:ARG:HG3	7:G:42:ARG:NH1	2.32	0.45
8:H:57:ARG:NH1	8:H:57:ARG:HB3	2.31	0.45
1:N:187:SER:HB2	1:N:277:MET:HE1	1.99	0.45
1:N:431:LEU:HD21	1:N:450:TRP:HB2	1.99	0.45
5:R:70:VAL:HG12	5:R:71:VAL:N	2.31	0.45
3:C:137:LEU:HD12	3:C:137:LEU:HA	1.69	0.44
5:E:70:VAL:HG12	5:E:71:VAL:N	2.31	0.44
3:P:42:LEU:HA	3:P:42:LEU:HD12	1.66	0.44
10:W:30:ILE:HG13	10:W:31:LEU:N	2.32	0.44
1:N:70:VAL:HG11	1:N:246:LEU:HD22	2.00	0.44
7:G:5:LYS:HG3	1:N:282:PHE:CA	2.47	0.44
5:R:78:HIS:CE1	9:V:12:LEU:HD22	2.52	0.44
1:A:51:ASP:O	1:A:55:ASN:HB2	2.17	0.44
4:D:79:LYS:NZ	11:K:15:ASN:HD21	2.15	0.44
1:N:495:LEU:HD23	1:N:495:LEU:HA	1.68	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:58:ALA:O	2:B:60:GLU:HG3	2.16	0.44
6:F:33:ILE:HG22	6:F:34:LEU:HD12	1.98	0.44
7:G:3:ALA:O	7:G:4:ALA:HB2	2.17	0.44
1:N:147:ILE:HD11	1:N:209:LEU:HD23	1.98	0.44
3:P:128:GLU:HB3	3:P:129:VAL:H	1.56	0.44
4:Q:37:GLN:O	4:Q:41:LYS:HG2	2.18	0.44
3:P:183:GLU:O	7:T:42:ARG:NH2	2.50	0.44
10:W:31:LEU:HA	10:W:31:LEU:HD23	1.83	0.44
10:W:3:ASN:ND2	10:W:5:VAL:HG13	2.33	0.44
2:B:189:PRO:HG2	9:I:63:MET:HE3	1.99	0.44
16:N:516:HEA:O11	16:N:516:HEA:HHC	2.17	0.44
2:O:59:GLN:CA	2:O:62:GLU:HB2	2.46	0.44
2:O:63:THR:HA	2:O:66:THR:HG22	2.00	0.44
4:Q:79:LYS:HZ3	11:X:15:ASN:HD21	1.65	0.44
3:C:63:ARG:HA	3:C:67:PHE:CD2	2.53	0.44
8:H:84:LYS:HA	8:H:84:LYS:HD2	1.83	0.44
2:O:162:SER:HB2	2:O:198:GLU:HB2	1.99	0.44
7:T:42:ARG:NH1	7:T:42:ARG:HG3	2.32	0.44
10:W:2:GLU:CG	10:W:3:ASN:H	2.30	0.44
2:O:1:MET:SD	2:O:133:LEU:HD13	2.58	0.44
1:N:449:MET:SD	2:O:5:MET:HE1	2.58	0.44
3:P:63:ARG:HA	3:P:67:PHE:CD2	2.53	0.44
2:O:191:LEU:HB2	4:Q:126:MET:HE1	1.99	0.44
2:B:1:MET:SD	2:B:133:LEU:HD13	2.58	0.44
3:C:19:THR:CG2	3:C:53:THR:OG1	2.66	0.44
4:D:37:GLN:O	4:D:41:LYS:HG2	2.18	0.44
9:I:61:GLU:OE1	9:I:64:ARG:NH1	2.51	0.44
12:L:35:ALA:HB3	12:L:36:PRO:HD3	1.99	0.44
2:B:145:PRO:HG3	2:B:148:MET:HE2	2.00	0.43
2:B:63:THR:HA	2:B:66:THR:HG22	2.00	0.43
10:J:36:MET:O	10:J:40:LEU:HG	2.18	0.43
3:P:60:ASP:O	3:P:64:GLU:HG3	2.18	0.43
9:V:68:ILE:HG13	9:V:69:PHE:N	2.33	0.43
5:E:43:PRO:HB2	5:E:48:ILE:CD1	2.48	0.43
3:P:19:THR:CG2	3:P:53:THR:OG1	2.66	0.43
3:P:154:GLY:CA	6:S:6:VAL:HG22	2.47	0.43
9:V:26:MET:SD	9:V:26:MET:N	2.91	0.43
3:C:41:THR:O	3:C:45:ILE:HG13	2.18	0.43
3:C:19:THR:HG22	3:C:53:THR:OG1	2.18	0.43
10:J:3:ASN:ND2	10:J:5:VAL:HG13	2.33	0.43
1:N:131:PRO:HB2	2:O:159:VAL:HA	1.99	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:P:41:THR:O	3:P:45:ILE:HG13	2.18	0.43
10:W:36:MET:O	10:W:40:LEU:HG	2.18	0.43
1:A:282:PHE:CA	7:T:5:LYS:HG3	2.49	0.43
1:N:311:ILE:HG21	1:N:311:ILE:HD13	1.75	0.43
7:T:3:ALA:O	7:T:4:ALA:HB2	2.17	0.43
9:V:21:ILE:O	9:V:25:PHE:HD2	2.02	0.43
2:B:52:HIS:CD2	5:E:40:ASP:HB2	2.54	0.43
2:B:78:LEU:HB2	2:B:79:PRO:CD	2.38	0.43
3:C:183:GLU:O	7:G:42:ARG:NH2	2.50	0.43
1:N:353:LEU:HA	1:N:353:LEU:HD12	1.91	0.43
1:N:356:ILE:HD13	1:N:356:ILE:HA	1.90	0.43
1:N:51:ASP:O	1:N:55:ASN:HB2	2.17	0.43
1:A:131:PRO:HB2	2:B:159:VAL:HA	1.99	0.43
3:C:60:ASP:O	3:C:64:GLU:HG3	2.19	0.43
10:J:30:ILE:HG13	10:J:31:LEU:N	2.32	0.43
9:V:61:GLU:OE1	9:V:64:ARG:NH1	2.51	0.43
3:C:173:PHE:C	3:C:173:PHE:CD1	2.92	0.43
9:I:42:LYS:HB3	9:I:42:LYS:HE2	1.83	0.43
9:I:68:ILE:HG13	9:I:69:PHE:N	2.33	0.43
12:L:41:ARG:HD2	13:M:40:TYR:CZ	2.54	0.43
1:N:247:ILE:HB	16:N:516:HEA:HBC1	2.01	0.43
2:O:3:TYR:N	2:O:3:TYR:CD1	2.87	0.43
5:R:27:TRP:CH2	6:S:86:GLY:HA2	2.54	0.43
8:H:49:ASP:CB	8:U:49:ASP:HB2	2.44	0.43
12:Y:41:ARG:HD2	13:Z:40:TYR:CZ	2.54	0.43
1:A:278:MET:HG2	7:T:5:LYS:HZ3	1.84	0.43
1:A:377:PHE:CD2	16:A:516:HEA:HAD1	2.53	0.43
3:C:80:ARG:HG2	3:C:233:PHE:CE1	2.54	0.43
4:D:126:MET:HG3	4:D:128:VAL:HG23	2.00	0.43
3:C:110:PRO:HB3	8:H:30:TRP:CD2	2.54	0.43
1:N:106:PRO:HB2	1:N:107:PRO:HD3	2.01	0.43
6:S:53:THR:HG22	6:S:54:ASN:OD1	2.19	0.43
2:B:3:TYR:N	2:B:3:TYR:CD1	2.87	0.43
11:K:44:PRO:HA	11:K:47:ARG:NH1	2.34	0.43
3:P:80:ARG:HG2	3:P:233:PHE:CE1	2.54	0.43
4:Q:126:MET:HG3	4:Q:128:VAL:HG23	2.00	0.43
5:E:27:TRP:CH2	6:F:86:GLY:HA2	2.54	0.42
9:I:35:TYR:O	9:I:39:VAL:HB	2.19	0.42
7:T:5:LYS:CD	7:T:6:GLY:N	2.81	0.42
11:X:44:PRO:HA	11:X:47:ARG:NH1	2.34	0.42
1:A:70:VAL:HG11	1:A:246:LEU:HD22	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:98:LYS:HB2	2:B:98:LYS:HE3	1.80	0.42
9:I:26:MET:SD	9:I:26:MET:N	2.91	0.42
10:J:31:LEU:HD23	10:J:31:LEU:HA	1.83	0.42
1:N:377:PHE:CD2	16:N:516:HEA:HAD1	2.53	0.42
2:O:145:PRO:HG3	2:O:148:MET:HE2	2.00	0.42
3:P:110:PRO:HB3	8:U:30:TRP:CD2	2.54	0.42
7:T:78:LEU:HA	7:T:78:LEU:HD12	1.75	0.42
3:P:105:SER:HG	3:P:259:TRP:HZ3	1.65	0.42
3:P:173:PHE:C	3:P:173:PHE:CD1	2.92	0.42
7:T:44:ARG:HA	7:T:45:PRO:HD2	1.75	0.42
9:V:35:TYR:O	9:V:39:VAL:HB	2.19	0.42
1:A:462:LEU:O	1:A:466:MET:HG3	2.20	0.42
6:F:86:GLY:O	6:F:87:THR:O	2.37	0.42
7:G:78:LEU:HD12	7:G:78:LEU:HA	1.75	0.42
1:N:462:LEU:O	1:N:466:MET:HG3	2.20	0.42
3:P:19:THR:HG22	3:P:53:THR:OG1	2.18	0.42
6:S:86:GLY:O	6:S:87:THR:O	2.37	0.42
8:U:57:ARG:NH1	8:U:57:ARG:HB3	2.31	0.42
16:A:515:HEA:H172	16:A:515:HEA:H272	1.90	0.42
7:G:5:LYS:CD	7:G:6:GLY:N	2.81	0.42
7:G:77:PRO:HA	7:G:82:TYR:HA	2.02	0.42
2:O:52:HIS:CD2	5:R:40:ASP:HB2	2.54	0.42
8:U:34:LEU:HA	8:U:34:LEU:HD23	1.93	0.42
1:A:106:PRO:HB2	1:A:107:PRO:HD3	2.01	0.42
1:N:498:CYS:HA	1:N:499:PRO:HA	1.78	0.42
2:O:94:SER:OG	2:O:148:MET:HE3	2.19	0.42
5:R:21:LYS:HA	5:R:22:PRO:HD2	1.84	0.42
6:F:53:THR:HG22	6:F:54:ASN:OD1	2.19	0.42
2:O:146:MET:SD	2:O:189:PRO:HB3	2.60	0.42
5:R:21:LYS:O	5:R:57:ARG:NH1	2.53	0.42
1:N:474:GLU:HG3	1:N:475:ALA:N	2.34	0.42
1:A:474:GLU:HG3	1:A:475:ALA:N	2.34	0.42
6:F:6:VAL:HA	6:F:7:PRO:HD3	1.91	0.42
8:U:57:ARG:HA	8:U:60:TYR:CD2	2.55	0.42
9:V:29:LEU:HD23	9:V:29:LEU:HA	1.85	0.42
3:P:188:ILE:H	3:P:188:ILE:HG13	1.62	0.42
4:Q:127:LYS:O	4:Q:130:PRO:HD3	2.20	0.42
3:P:146:TRP:NE1	7:T:17:ARG:HB2	2.35	0.42
1:A:247:ILE:HB	16:A:516:HEA:HBC1	2.01	0.41
2:B:185:MET:C	2:B:185:MET:SD	2.99	0.41
2:B:193:TYR:N	2:B:193:TYR:CD1	2.88	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:H:24:ASN:HD22	8:H:25:GLN:N	2.18	0.41
8:H:65:PRO:HG2	8:H:68:TRP:CG	2.56	0.41
9:I:21:ILE:O	9:I:25:PHE:HD2	2.02	0.41
7:T:42:ARG:NH1	7:T:74:ARG:NH2	2.68	0.41
1:A:276:ALA:O	1:A:280:ILE:HG13	2.21	0.41
1:A:314:ILE:HB	1:A:315:PRO:CD	2.50	0.41
5:E:21:LYS:O	5:E:57:ARG:NH1	2.53	0.41
8:H:57:ARG:HA	8:H:60:TYR:CD2	2.55	0.41
2:O:193:TYR:CE1	4:Q:126:MET:HE1	2.54	0.41
3:P:146:TRP:CE2	7:T:17:ARG:HB2	2.56	0.41
8:U:24:ASN:HD22	8:U:25:GLN:N	2.18	0.41
4:D:130:PRO:O	4:D:136:ALA:HB2	2.20	0.41
3:P:223:LEU:HD23	3:P:223:LEU:HA	1.79	0.41
3:P:6:HIS:CD2	3:P:8:TYR:HB2	2.55	0.41
6:S:16:LEU:O	6:S:20:VAL:HG13	2.21	0.41
7:T:77:PRO:HA	7:T:82:TYR:HA	2.02	0.41
1:A:187:SER:HB2	1:A:277:MET:CE	2.50	0.41
1:A:298:ASP:HB2	1:A:301:THR:CG2	2.49	0.41
16:A:515:HEA:H122	16:A:515:HEA:HHC	2.02	0.41
2:B:122:MET:HB2	2:B:208:PRO:CD	2.47	0.41
2:B:94:SER:OG	2:B:148:MET:HE3	2.19	0.41
10:J:12:PHE:O	10:J:23:LYS:HE2	2.21	0.41
10:J:5:VAL:O	10:J:9:GLN:HG3	2.21	0.41
2:O:185:MET:C	2:O:185:MET:SD	2.99	0.41
13:Z:37:LEU:HD23	13:Z:37:LEU:HA	1.83	0.41
2:B:146:MET:SD	2:B:189:PRO:HB3	2.60	0.41
4:D:127:LYS:O	4:D:130:PRO:HD3	2.20	0.41
3:C:146:TRP:CE2	7:G:17:ARG:HB2	2.56	0.41
1:N:173:PRO:HA	1:N:174:PRO:HD3	1.96	0.41
5:R:76:GLY:HA3	5:R:77:PRO:HD2	1.85	0.41
10:W:5:VAL:O	10:W:9:GLN:HG3	2.21	0.41
2:O:5:MET:HE2	11:X:42:PRO:HA	2.01	0.41
3:C:228:THR:O	3:C:232:HIS:HD2	2.04	0.41
4:D:137:LYS:O	4:D:145:TRP:HE3	2.03	0.41
3:P:233:PHE:HA	3:P:236:GLU:HG3	2.03	0.41
8:U:20:PHE:N	8:U:21:PRO:HD3	2.35	0.41
10:W:12:PHE:O	10:W:23:LYS:HE2	2.21	0.41
1:A:461:SER:O	1:A:465:VAL:HG13	2.21	0.41
2:B:60:GLU:CD	2:B:61:VAL:H	2.24	0.41
3:C:154:GLY:CA	6:F:6:VAL:HG22	2.47	0.41
8:H:20:PHE:N	8:H:21:PRO:HD3	2.35	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:5:MET:HE3	11:K:42:PRO:HA	2.03	0.41
13:M:35:TYR:HD2	13:M:36:HIS:CE1	2.39	0.41
1:N:461:SER:O	1:N:465:VAL:HG13	2.21	0.41
1:N:69:MET:HE1	1:N:70:VAL:HG23	2.02	0.41
4:Q:86:MET:O	11:X:25:CYS:HB2	2.20	0.41
1:A:168:ILE:CG2	1:A:189:MET:HG3	2.51	0.41
5:E:76:GLY:HA3	5:E:77:PRO:HD2	1.85	0.41
4:D:86:MET:O	11:K:25:CYS:HB2	2.20	0.41
1:N:314:ILE:HB	1:N:315:PRO:CD	2.50	0.41
2:O:100:MET:HE2	2:O:157:GLU:HG3	2.03	0.41
5:E:104:LEU:HA	5:E:104:LEU:HD23	1.87	0.41
1:N:187:SER:HB2	1:N:277:MET:CE	2.50	0.41
16:N:515:HEA:HMC1	16:N:515:HEA:HBC1	2.03	0.41
1:N:23:GLY:HA3	1:N:73:ILE:HG13	2.03	0.41
2:O:193:TYR:N	2:O:193:TYR:CD1	2.88	0.41
8:U:65:PRO:HG2	8:U:68:TRP:CG	2.55	0.41
1:A:398:PRO:HB3	1:A:482:VAL:HG21	2.03	0.41
16:A:515:HEA:HMC1	16:A:515:HEA:HBC1	2.03	0.41
3:C:223:LEU:HA	3:C:223:LEU:HD23	1.79	0.41
6:F:10:GLU:HG2	6:F:25:ARG:HH22	1.86	0.41
7:G:42:ARG:NH1	7:G:74:ARG:NH2	2.68	0.41
2:O:164:ALA:HB2	2:O:171:LYS:HD3	2.03	0.41
1:A:23:GLY:HA3	1:A:73:ILE:HG13	2.03	0.41
3:C:6:HIS:CD2	3:C:8:TYR:HB2	2.55	0.41
6:F:16:LEU:O	6:F:20:VAL:HG13	2.21	0.41
6:F:49:VAL:O	6:F:91:LEU:HD12	2.21	0.41
1:N:69:MET:CE	1:N:70:VAL:HG23	2.51	0.41
3:P:137:LEU:HD12	3:P:137:LEU:HA	1.69	0.41
3:P:182:TYR:O	7:T:72:ASN:HB2	2.21	0.41
6:F:19:GLU:HB3	6:F:98:HIS:CE1	2.56	0.40
2:O:123:ILE:HA	2:O:124:PRO:HD3	1.96	0.40
2:O:60:GLU:CD	2:O:61:VAL:H	2.24	0.40
4:Q:137:LYS:O	4:Q:145:TRP:HE3	2.03	0.40
11:X:24:PHE:CE1	11:X:28:VAL:HG21	2.56	0.40
3:C:122:HIS:HA	3:C:123:PRO:HD2	1.91	0.40
5:E:79:LYS:H	5:E:79:LYS:HD3	1.87	0.40
6:F:77:GLY:O	6:F:90:LYS:NZ	2.54	0.40
1:N:442:ASP:OD2	2:O:134:ARG:NH2	2.54	0.40
8:U:17:ASP:OD1	8:U:19:ARG:HG2	2.21	0.40
1:A:172:LYS:HZ3	1:A:172:LYS:HB2	1.85	0.40
3:C:146:TRP:NE1	7:G:17:ARG:HB2	2.35	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:182:TYR:O	7:G:72:ASN:HB2	2.21	0.40
9:I:21:ILE:HD12	9:I:21:ILE:HA	1.82	0.40
4:Q:119:GLN:O	4:Q:123:MET:HG3	2.22	0.40
6:S:10:GLU:HG2	6:S:25:ARG:HH22	1.86	0.40
1:A:354:THR:HG21	1:A:376:HIS:HA	2.03	0.40
1:A:64:VAL:HA	1:A:68:PHE:HD2	1.87	0.40
2:B:100:MET:HE2	2:B:157:GLU:HG3	2.04	0.40
1:A:442:ASP:OD2	2:B:134:ARG:NH2	2.54	0.40
2:B:164:ALA:HB2	2:B:171:LYS:HD3	2.03	0.40
4:D:24:LEU:HB3	5:E:30:ARG:HG2	2.04	0.40
11:K:24:PHE:CE1	11:K:28:VAL:HG21	2.56	0.40
4:Q:130:PRO:O	4:Q:136:ALA:HB2	2.21	0.40
5:R:43:PRO:HB2	5:R:48:ILE:CD1	2.48	0.40
6:S:19:GLU:HB3	6:S:98:HIS:CE1	2.57	0.40
7:T:54:ARG:CD	7:T:54:ARG:N	2.80	0.40
2:B:196:CYS:HB2	2:B:207:MET:HG3	2.04	0.40
4:D:102:TYR:HD2	13:M:35:TYR:HE1	1.68	0.40
1:N:373:VAL:O	1:N:377:PHE:CD2	2.75	0.40
2:O:168:LEU:HD13	2:O:184:LEU:HG	2.04	0.40
2:O:196:CYS:HB2	2:O:207:MET:HG3	2.04	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	512/514 (100%)	479 (94%)	29 (6%)	4 (1%)	22	55
1	N	512/514 (100%)	479 (94%)	29 (6%)	4 (1%)	22	55
2	B	225/227 (99%)	203 (90%)	19 (8%)	3 (1%)	14	41
2	O	225/227 (99%)	203 (90%)	19 (8%)	3 (1%)	14	41

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
3	C	259/261 (99%)	249 (96%)	10 (4%)	0	100	100
3	P	259/261 (99%)	249 (96%)	10 (4%)	0	100	100
4	D	142/147 (97%)	135 (95%)	7 (5%)	0	100	100
4	Q	142/147 (97%)	135 (95%)	7 (5%)	0	100	100
5	E	107/109 (98%)	104 (97%)	3 (3%)	0	100	100
5	R	107/109 (98%)	104 (97%)	3 (3%)	0	100	100
6	F	96/98 (98%)	86 (90%)	6 (6%)	4 (4%)	3	10
6	S	96/98 (98%)	86 (90%)	6 (6%)	4 (4%)	3	10
7	G	82/84 (98%)	67 (82%)	10 (12%)	5 (6%)	2	4
7	T	82/84 (98%)	67 (82%)	10 (12%)	5 (6%)	2	4
8	H	73/85 (86%)	64 (88%)	8 (11%)	1 (1%)	13	39
8	U	73/85 (86%)	64 (88%)	8 (11%)	1 (1%)	13	39
9	I	71/73 (97%)	65 (92%)	6 (8%)	0	100	100
9	V	71/73 (97%)	65 (92%)	6 (8%)	0	100	100
10	J	54/59 (92%)	48 (89%)	4 (7%)	2 (4%)	4	13
10	W	54/59 (92%)	48 (89%)	4 (7%)	2 (4%)	4	13
11	K	47/56 (84%)	41 (87%)	6 (13%)	0	100	100
11	X	47/56 (84%)	41 (87%)	6 (13%)	0	100	100
12	L	45/47 (96%)	42 (93%)	3 (7%)	0	100	100
12	Y	45/47 (96%)	42 (93%)	3 (7%)	0	100	100
13	M	41/46 (89%)	39 (95%)	2 (5%)	0	100	100
13	Z	41/46 (89%)	39 (95%)	2 (5%)	0	100	100
All	All	3508/3612 (97%)	3244 (92%)	226 (6%)	38 (1%)	17	47

All (38) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	328	HIS
1	A	508	PRO
6	F	2	SER
6	F	87	THR
6	F	95	GLN
7	G	4	ALA
7	G	9	GLY

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Mol	Chain	Res	Type
8	H	46	LYS
10	J	2	GLU
1	N	328	HIS
1	N	508	PRO
6	S	2	SER
6	S	87	THR
6	S	95	GLN
7	T	4	ALA
7	T	9	GLY
8	U	46	LYS
10	W	2	GLU
7	G	5	LYS
7	T	5	LYS
2	B	104	TRP
7	G	61	SER
2	O	104	TRP
7	T	61	SER
1	A	51	ASP
1	N	51	ASP
2	B	103	GLN
10	J	3	ASN
2	O	103	GLN
10	W	3	ASN
1	A	91	ASP
2	B	158	ASP
7	G	49	PRO
1	N	91	ASP
2	O	158	ASP
7	T	49	PRO
6	F	15	GLY
6	S	15	GLY

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.

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	427/427 (100%)	389 (91%)	38 (9%)	11	32
1	N	427/427 (100%)	389 (91%)	38 (9%)	11	32
2	B	211/211 (100%)	191 (90%)	20 (10%)	10	28
2	O	211/211 (100%)	191 (90%)	20 (10%)	10	28
3	C	226/226 (100%)	199 (88%)	27 (12%)	6	18
3	P	226/226 (100%)	199 (88%)	27 (12%)	6	18
4	D	128/129 (99%)	120 (94%)	8 (6%)	21	51
4	Q	128/129 (99%)	120 (94%)	8 (6%)	21	51
5	E	95/95 (100%)	89 (94%)	6 (6%)	21	51
5	R	95/95 (100%)	89 (94%)	6 (6%)	21	51
6	F	81/81 (100%)	76 (94%)	5 (6%)	21	52
6	S	81/81 (100%)	76 (94%)	5 (6%)	21	52
7	G	68/68 (100%)	50 (74%)	18 (26%)	0	1
7	T	68/68 (100%)	50 (74%)	18 (26%)	0	1
8	H	67/75 (89%)	58 (87%)	9 (13%)	4	13
8	U	67/75 (89%)	58 (87%)	9 (13%)	4	13
9	I	58/58 (100%)	53 (91%)	5 (9%)	12	34
9	V	58/58 (100%)	53 (91%)	5 (9%)	12	34
10	J	47/50 (94%)	40 (85%)	7 (15%)	3	10
10	W	47/50 (94%)	40 (85%)	7 (15%)	3	10
11	K	39/46 (85%)	37 (95%)	2 (5%)	28	61
11	X	39/46 (85%)	37 (95%)	2 (5%)	28	61
12	L	40/40 (100%)	38 (95%)	2 (5%)	28	62
12	Y	40/40 (100%)	38 (95%)	2 (5%)	28	62
13	M	37/38 (97%)	34 (92%)	3 (8%)	14	37
13	Z	37/38 (97%)	34 (92%)	3 (8%)	14	37
All	All	3048/3088 (99%)	2748 (90%)	300 (10%)	9	27

All (300) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	18	LEU
1	A	35	LEU
1	A	92	MET
1	A	96	ARG
1	A	105	LEU
1	A	109	PHE
1	A	115	SER
1	A	138	HIS
1	A	150	LEU
1	A	159	LEU
1	A	187	SER
1	A	188	VAL
1	A	199	LEU
1	A	213	ARG
1	A	238	PHE
1	A	241	PRO
1	A	273	MET
1	A	295	VAL
1	A	301	THR
1	A	306	THR
1	A	318	VAL
1	A	324	LEU
1	A	347	LEU
1	A	353	LEU
1	A	354	THR
1	A	365	ILE
1	A	369	ASP
1	A	373	VAL
1	A	383	MET
1	A	417	MET
1	A	465	VAL
1	A	467	LEU
1	A	474	GLU
1	A	486	ASP
1	A	492	LEU
1	A	508	PRO
1	A	509	THR
1	A	512	ASN
2	B	7	LEU
2	B	31	VAL
2	B	52	HIS
2	B	60	GLU
2	B	63	THR

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Mol	Chain	Res	Type
2	B	65	TRP
2	B	88	ASP
2	B	92	ASN
2	B	113	TYR
2	B	125	THR
2	B	130	PRO
2	B	134	ARG
2	B	142	VAL
2	B	147	GLU
2	B	148	MET
2	B	170	LEU
2	B	171	LYS
2	B	185	MET
2	B	205	SER
2	B	216	LEU
3	C	1	MET
3	C	11	VAL
3	C	13	PRO
3	C	14	SER
3	C	18	LEU
3	C	19	THR
3	C	22	LEU
3	C	38	ASN
3	C	39	SER
3	C	85	LEU
3	C	92	LEU
3	C	112	LEU
3	C	127	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	188	ILE
3	C	196	THR
3	C	199	VAL
3	C	214	PHE
3	C	222	GLN
3	C	258	TRP

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Mol	Chain	Res	Type
4	D	31	LYS
4	D	36	SER
4	D	40	LEU
4	D	59	LEU
4	D	62	LEU
4	D	107	ILE
4	D	143	ASN
4	D	147	LYS
5	E	7	THR
5	E	29	LEU
5	E	70	VAL
5	E	79	LYS
5	E	80	GLU
5	E	90	ARG
6	F	37	LYS
6	F	53	THR
6	F	74	LEU
6	F	95	GLN
6	F	98	HIS
7	G	5	LYS
7	G	7	ASP
7	G	8	HIS
7	G	14	ARG
7	G	17	ARG
7	G	33	LEU
7	G	34	ASN
7	G	37	LEU
7	G	38	HIS
7	G	41	HIS
7	G	42	ARG
7	G	43	GLU
7	G	48	ILE
7	G	54	ARG
7	G	56	ARG
7	G	68	THR
7	G	69	PHE
7	G	78	LEU
8	H	19	ARG
8	H	24	ASN
8	H	28	ASN
8	H	29	CYS
8	H	51	SER

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Mol	Chain	Res	Type
8	H	53	CYS
8	H	57	ARG
8	H	60	TYR
8	H	75	ARG
9	I	2	THR
9	I	8	GLN
9	I	26	MET
9	I	44	LYS
9	I	64	ARG
10	J	1	PHE
10	J	2	GLU
10	J	3	ASN
10	J	8	LYS
10	J	16	ASN
10	J	23	LYS
10	J	27	THR
11	K	48	VAL
11	K	49	THR
12	L	15	VAL
12	L	22	LEU
13	M	13	LYS
13	M	42	LYS
13	M	43	SER
1	N	18	LEU
1	N	35	LEU
1	N	92	MET
1	N	96	ARG
1	N	105	LEU
1	N	109	PHE
1	N	115	SER
1	N	138	HIS
1	N	150	LEU
1	N	159	LEU
1	N	187	SER
1	N	188	VAL
1	N	199	LEU
1	N	213	ARG
1	N	238	PHE
1	N	241	PRO
1	N	273	MET
1	N	295	VAL
1	N	301	THR

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Mol	Chain	Res	Type
1	N	306	THR
1	N	318	VAL
1	N	324	LEU
1	N	347	LEU
1	N	353	LEU
1	N	354	THR
1	N	365	ILE
1	N	369	ASP
1	N	373	VAL
1	N	383	MET
1	N	417	MET
1	N	465	VAL
1	N	467	LEU
1	N	474	GLU
1	N	486	ASP
1	N	492	LEU
1	N	508	PRO
1	N	509	THR
1	N	512	ASN
2	O	7	LEU
2	O	31	VAL
2	O	52	HIS
2	O	60	GLU
2	O	63	THR
2	O	65	TRP
2	O	88	ASP
2	O	92	ASN
2	O	113	TYR
2	O	125	THR
2	O	130	PRO
2	O	134	ARG
2	O	142	VAL
2	O	147	GLU
2	O	148	MET
2	O	170	LEU
2	O	171	LYS
2	O	185	MET
2	O	205	SER
2	O	216	LEU
3	P	1	MET
3	P	11	VAL
3	P	13	PRO

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Mol	Chain	Res	Type
3	P	14	SER
3	P	18	LEU
3	P	19	THR
3	P	22	LEU
3	P	38	ASN
3	P	39	SER
3	P	85	LEU
3	P	92	LEU
3	P	112	LEU
3	P	127	LEU
3	P	128	GLU
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	188	ILE
3	P	196	THR
3	P	199	VAL
3	P	214	PHE
3	P	222	GLN
3	P	258	TRP
4	Q	31	LYS
4	Q	36	SER
4	Q	40	LEU
4	Q	59	LEU
4	Q	62	LEU
4	Q	107	ILE
4	Q	143	ASN
4	Q	147	LYS
5	R	7	THR
5	R	29	LEU
5	R	70	VAL
5	R	79	LYS
5	R	80	GLU
5	R	90	ARG
6	S	37	LYS
6	S	53	THR
6	S	74	LEU
6	S	95	GLN

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Mol	Chain	Res	Type
6	S	98	HIS
7	T	5	LYS
7	T	7	ASP
7	T	8	HIS
7	T	14	ARG
7	T	17	ARG
7	T	33	LEU
7	T	34	ASN
7	T	37	LEU
7	T	38	HIS
7	T	41	HIS
7	T	42	ARG
7	T	43	GLU
7	T	48	ILE
7	T	54	ARG
7	T	56	ARG
7	T	68	THR
7	T	69	PHE
7	T	78	LEU
8	U	19	ARG
8	U	24	ASN
8	U	28	ASN
8	U	29	CYS
8	U	51	SER
8	U	53	CYS
8	U	57	ARG
8	U	60	TYR
8	U	75	ARG
9	V	2	THR
9	V	8	GLN
9	V	26	MET
9	V	44	LYS
9	V	64	ARG
10	W	1	PHE
10	W	2	GLU
10	W	3	ASN
10	W	8	LYS
10	W	16	ASN
10	W	23	LYS
10	W	27	THR
11	X	48	VAL
11	X	49	THR

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Mol	Chain	Res	Type
12	Y	15	VAL
12	Y	22	LEU
13	Z	13	LYS
13	Z	42	LYS
13	Z	43	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (66) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	11	ASN
1	A	12	HIS
1	A	43	GLN
1	A	99	ASN
1	A	170	ASN
1	A	256	HIS
1	A	360	ASN
1	A	413	HIS
1	A	512	ASN
2	B	103	GLN
2	B	203	ASN
3	C	6	HIS
3	C	12	ASN
3	C	133	ASN
3	C	148	HIS
3	C	207	HIS
3	C	222	GLN
3	C	232	HIS
4	D	109	HIS
5	E	34	ASN
6	F	66	ASN
7	G	52	HIS
8	H	24	ASN
8	H	25	GLN
8	H	28	ASN
8	H	37	HIS
9	I	8	GLN
10	J	3	ASN
10	J	16	ASN
11	K	10	HIS
11	K	15	ASN
11	K	41	ASN
13	M	39	ASN

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Mol	Chain	Res	Type
1	N	11	ASN
1	N	12	HIS
1	N	43	GLN
1	N	99	ASN
1	N	170	ASN
1	N	256	HIS
1	N	360	ASN
1	N	413	HIS
1	N	512	ASN
2	O	103	GLN
2	O	203	ASN
3	P	6	HIS
3	P	12	ASN
3	P	133	ASN
3	P	207	HIS
3	P	222	GLN
3	P	232	HIS
4	Q	109	HIS
5	R	34	ASN
6	S	66	ASN
7	T	52	HIS
8	U	23	GLN
8	U	24	ASN
8	U	25	GLN
8	U	28	ASN
8	U	37	HIS
9	V	8	GLN
10	W	3	ASN
10	W	16	ASN
11	X	10	HIS
11	X	15	ASN
11	X	41	ASN
13	Z	39	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 14 ligands modelled in this entry, 10 are monoatomic - leaving 4 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$
16	HEA	A	515	1	44,67,67	1.63	8 (18%)	37,103,103	1.84	9 (24%)
16	HEA	A	516	1	44,67,67	1.53	4 (9%)	37,103,103	1.36	7 (18%)
16	HEA	N	515	1	44,67,67	1.64	9 (20%)	37,103,103	1.84	10 (27%)
16	HEA	N	516	1	44,67,67	1.53	4 (9%)	37,103,103	1.36	7 (18%)

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
16	HEA	A	515	1	3/3/7/16	0/24/76/76	0/0/8/8
16	HEA	A	516	1	3/3/7/16	0/24/76/76	0/0/8/8
16	HEA	N	515	1	3/3/7/16	0/24/76/76	0/0/8/8
16	HEA	N	516	1	3/3/7/16	0/24/76/76	0/0/8/8

All (25) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
16	N	516	HEA	C3A-C2A	-6.81	1.31	1.40
16	A	516	HEA	C3A-C2A	-6.79	1.31	1.40
16	A	516	HEA	C3A-CMA	-3.90	1.37	1.46
16	N	516	HEA	C3A-CMA	-3.89	1.37	1.46
16	A	515	HEA	C3C-C2C	-3.36	1.35	1.40
16	N	515	HEA	C3C-C2C	-3.36	1.35	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
16	A	515	HEA	C3A-CMA	-2.97	1.39	1.46
16	N	515	HEA	C3A-CMA	-2.93	1.39	1.46
16	N	516	HEA	C3C-C2C	-2.74	1.36	1.40
16	A	516	HEA	C3C-C2C	-2.71	1.36	1.40
16	N	515	HEA	C3A-C2A	-2.30	1.37	1.40
16	A	515	HEA	C3A-C2A	-2.28	1.37	1.40
16	N	515	HEA	C3B-C2B	-2.00	1.34	1.41
16	A	515	HEA	C1C-NC	2.06	1.40	1.36
16	N	515	HEA	C1C-NC	2.07	1.40	1.36
16	N	516	HEA	C1A-NA	2.11	1.39	1.36
16	A	516	HEA	C1A-NA	2.12	1.39	1.36
16	N	515	HEA	C4C-NC	2.60	1.41	1.36
16	A	515	HEA	C4C-NC	2.61	1.41	1.36
16	N	515	HEA	C4D-ND	2.99	1.42	1.36
16	A	515	HEA	C4D-ND	3.00	1.42	1.36
16	A	515	HEA	C1A-NA	4.73	1.42	1.36
16	N	515	HEA	C1A-NA	4.81	1.42	1.36
16	A	515	HEA	C4B-NB	5.07	1.42	1.36
16	N	515	HEA	C4B-NB	5.08	1.42	1.36

All (33) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
16	N	515	HEA	C4B-C3B-C2B	-4.43	103.78	106.87
16	A	515	HEA	C4B-C3B-C2B	-4.40	103.79	106.87
16	N	515	HEA	CMB-C2B-C1B	-3.86	122.53	128.46
16	A	515	HEA	CMB-C2B-C1B	-3.86	122.54	128.46
16	A	515	HEA	C17-C18-C19	-3.82	118.09	127.68
16	N	515	HEA	C17-C18-C19	-3.81	118.12	127.68
16	A	515	HEA	C13-C14-C15	-3.06	120.00	127.68
16	N	515	HEA	C13-C14-C15	-3.05	120.01	127.68
16	N	515	HEA	C17-C16-C15	-2.34	105.03	112.93
16	A	515	HEA	C17-C16-C15	-2.34	105.03	112.93
16	A	515	HEA	C16-C17-C18	-2.27	104.19	111.97
16	N	515	HEA	C16-C17-C18	-2.26	104.21	111.97
16	N	516	HEA	C13-C14-C15	-2.22	122.09	127.68
16	A	515	HEA	C12-C13-C14	-2.21	106.51	112.33
16	N	515	HEA	C12-C13-C14	-2.21	106.52	112.33
16	A	516	HEA	C13-C14-C15	-2.21	122.14	127.68
16	A	516	HEA	C4B-C3B-C2B	-2.09	105.41	106.87
16	N	516	HEA	C4B-C3B-C2B	-2.06	105.43	106.87
16	N	515	HEA	C3C-C4C-NC	2.01	111.80	109.21

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
16	A	516	HEA	C26-C15-C16	2.11	118.95	115.29
16	N	516	HEA	C26-C15-C16	2.15	119.01	115.29
16	N	516	HEA	C25-C23-C24	2.18	119.68	114.60
16	A	516	HEA	C25-C23-C24	2.20	119.72	114.60
16	N	515	HEA	C20-C19-C18	2.33	125.86	121.10
16	A	515	HEA	C20-C19-C18	2.33	125.88	121.10
16	A	516	HEA	CMC-C2C-C3C	2.35	129.26	124.89
16	N	516	HEA	CMC-C2C-C3C	2.37	129.29	124.89
16	N	516	HEA	CBA-CAA-C2A	2.59	117.41	112.47
16	A	516	HEA	CBA-CAA-C2A	2.60	117.43	112.47
16	A	516	HEA	CBD-CAD-C3D	3.05	118.32	112.48
16	N	516	HEA	CBD-CAD-C3D	3.06	118.33	112.48
16	A	515	HEA	C1B-C2B-C3B	4.54	110.15	107.00
16	N	515	HEA	C1B-C2B-C3B	4.57	110.18	107.00

All (12) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
16	A	515	HEA	ND
16	A	515	HEA	NA
16	A	515	HEA	NB
16	N	516	HEA	ND
16	N	516	HEA	NA
16	N	516	HEA	NB
16	A	516	HEA	ND
16	A	516	HEA	NA
16	A	516	HEA	NB
16	N	515	HEA	ND
16	N	515	HEA	NA
16	N	515	HEA	NB

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 14 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
16	A	515	HEA	4	0
16	A	516	HEA	4	0
16	N	515	HEA	2	0
16	N	516	HEA	4	0

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.