



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 08:10 AM UTC

PDB ID : 3RJ1 / pdb_00003rj1
Title : Architecture of the Mediator Head module
Authors : Imasaki, T.; Calero, G.; Cai, G.; Tsai, K.L.; Yamada, K.; Cardelli, F.; Erdjument-Bromage, H.; Tempst, P.; Berger, I.; Kornberg, G.L.; Asturias, F.J.; Kornberg, R.D.; Takagi, Y.
Deposited on : 2011-04-14
Resolution : 4.30 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

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

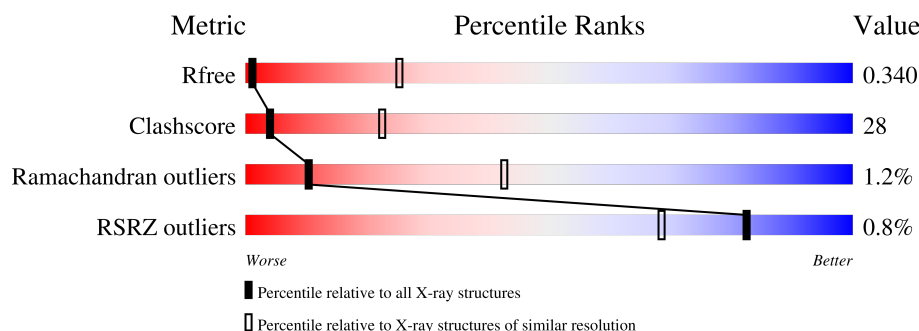
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 4.30 Å.

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



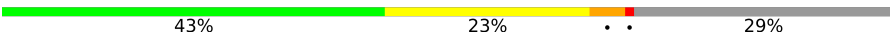
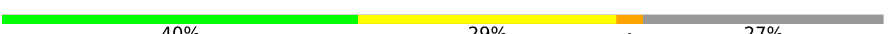
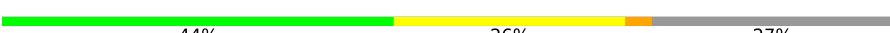
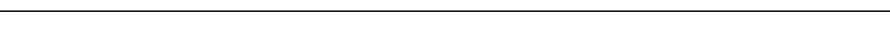
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1052 (4.70-3.90)
Clashscore	190562	1097 (4.70-3.90)
Ramachandran outliers	187476	1001 (4.70-3.90)
RSRZ outliers	180081	1049 (4.70-3.90)

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

Mol	Chain	Length	Quality of chain
1	A	131	50% 22% 28%
1	H	131	50% 22% 28%
1	O	131	51% 21% 28%
2	B	583	% 43% 11% 45%
2	I	583	43% 10% 45%
2	P	583	42% 11% 45%

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Mol	Chain	Length	Quality of chain
3	C	223	
3	J	223	
3	Q	223	
4	D	121	
4	K	121	
4	R	121	
5	E	275	
5	L	275	
5	S	275	
6	F	210	
6	M	210	
6	T	210	
7	G	295	
7	N	295	
7	U	295	

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
8	SE	B	1730	-	-	X	-
8	SE	D	1301	-	-	X	-
8	SE	D	222	-	-	X	-
8	SE	E	409	-	-	X	-
8	SE	E	410	-	-	X	-
8	SE	I	906	-	-	X	-
8	SE	J	225	-	-	X	-
8	SE	K	174	-	-	X	-
8	SE	K	175	-	-	X	-
8	SE	K	177	-	-	X	-
8	SE	L	408	-	-	X	-
8	SE	L	413	-	-	X	-

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
8	SE	L	415	-	-	X	-
8	SE	N	396	-	-	X	-
8	SE	N	398	-	-	X	-
8	SE	O	132	-	-	X	-
8	SE	P	1733	-	-	X	-
8	SE	Q	325	-	-	X	-
8	SE	S	310	-	-	X	-
8	SE	U	499	-	-	X	-

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 17945 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 Mediator of RNA polymerase II transcription subunit 11.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
1	A	94	Total	C	N	O	0	0	0
			468	280	94	94			
1	H	94	Total	C	N	O	0	0	0
			468	280	94	94			
1	O	94	Total	C	N	O	0	0	0
			468	280	94	94			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	16	GLY	THR	engineered mutation	UNP Q99278
A	17	SER	MET	engineered mutation	UNP Q99278
H	16	GLY	THR	engineered mutation	UNP Q99278
H	17	SER	MET	engineered mutation	UNP Q99278
O	16	GLY	THR	engineered mutation	UNP Q99278
O	17	SER	MET	engineered mutation	UNP Q99278

- Molecule 2 is a protein called Mediator of RNA polymerase II transcription subunit 17.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	B	318	Total	C	N	O	0	0	0
			1583	947	318	318			
2	I	318	Total	C	N	O	0	0	0
			1583	947	318	318			
2	P	318	Total	C	N	O	0	0	0
			1583	947	318	318			

There are 33 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	98	MSE	-	expression tag	UNP P32569
B	99	HIS	-	expression tag	UNP P32569

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Chain	Residue	Modelled	Actual	Comment	Reference
B	100	HIS	-	expression tag	UNP P32569
B	101	HIS	-	expression tag	UNP P32569
B	102	HIS	-	expression tag	UNP P32569
B	103	HIS	-	expression tag	UNP P32569
B	104	HIS	-	expression tag	UNP P32569
B	105	HIS	-	expression tag	UNP P32569
B	106	HIS	-	expression tag	UNP P32569
B	107	HIS	-	expression tag	UNP P32569
B	108	HIS	-	expression tag	UNP P32569
I	98	MSE	-	expression tag	UNP P32569
I	99	HIS	-	expression tag	UNP P32569
I	100	HIS	-	expression tag	UNP P32569
I	101	HIS	-	expression tag	UNP P32569
I	102	HIS	-	expression tag	UNP P32569
I	103	HIS	-	expression tag	UNP P32569
I	104	HIS	-	expression tag	UNP P32569
I	105	HIS	-	expression tag	UNP P32569
I	106	HIS	-	expression tag	UNP P32569
I	107	HIS	-	expression tag	UNP P32569
I	108	HIS	-	expression tag	UNP P32569
P	98	MSE	-	expression tag	UNP P32569
P	99	HIS	-	expression tag	UNP P32569
P	100	HIS	-	expression tag	UNP P32569
P	101	HIS	-	expression tag	UNP P32569
P	102	HIS	-	expression tag	UNP P32569
P	103	HIS	-	expression tag	UNP P32569
P	104	HIS	-	expression tag	UNP P32569
P	105	HIS	-	expression tag	UNP P32569
P	106	HIS	-	expression tag	UNP P32569
P	107	HIS	-	expression tag	UNP P32569
P	108	HIS	-	expression tag	UNP P32569

- Molecule 3 is a protein called Mediator of RNA polymerase II transcription subunit 8.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	C	159	Total	C	N	O	0	0	0
			791	473	159	159			
3	J	159	Total	C	N	O	0	0	0
			791	473	159	159			
3	Q	159	Total	C	N	O	0	0	0
			791	473	159	159			

- Molecule 4 is a protein called Mediator of RNA polymerase II transcription subunit 22.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	D	88	Total	C	N	O	0	0	0
			440	264	88	88			
4	K	88	Total	C	N	O	0	0	0
			440	264	88	88			
4	R	88	Total	C	N	O	0	0	0
			440	264	88	88			

- Molecule 5 is a protein called Mediator of RNA polymerase II transcription subunit 18.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
5	E	231	Total	C	N	O	0	0	0
			1141	679	231	231			
5	L	231	Total	C	N	O	0	0	0
			1141	679	231	231			
5	S	231	Total	C	N	O	0	0	0
			1141	679	231	231			

- Molecule 6 is a protein called Mediator of RNA polymerase II transcription subunit 20.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
6	F	205	Total	C	N	O	0	0	0
			1012	602	205	205			
6	M	205	Total	C	N	O	0	0	0
			1012	602	205	205			
6	T	205	Total	C	N	O	0	0	0
			1012	602	205	205			

- Molecule 7 is a protein called Mediator of RNA polymerase II transcription subunit 6.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
7	G	103	Total	C	N	O	0	0	0
			514	308	103	103			
7	N	103	Total	C	N	O	0	0	0
			514	308	103	103			
7	U	103	Total	C	N	O	0	0	0
			514	308	103	103			

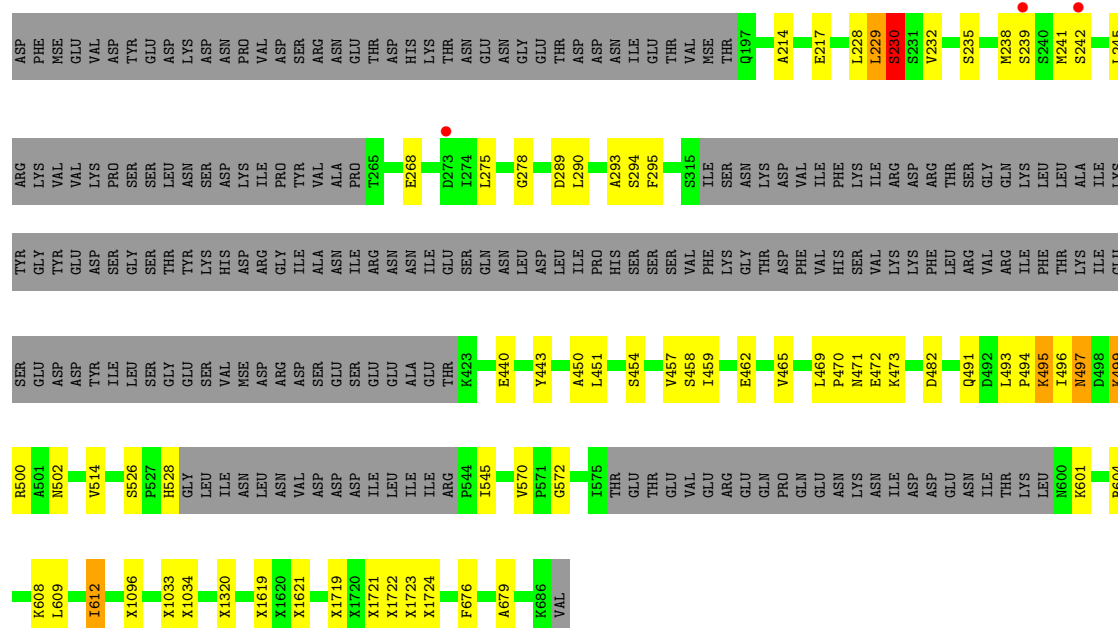
- Molecule 8 is SELENIUM ATOM (CCD ID: SE) (formula: Se).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	3	Total	Se	1	0
			3	3		

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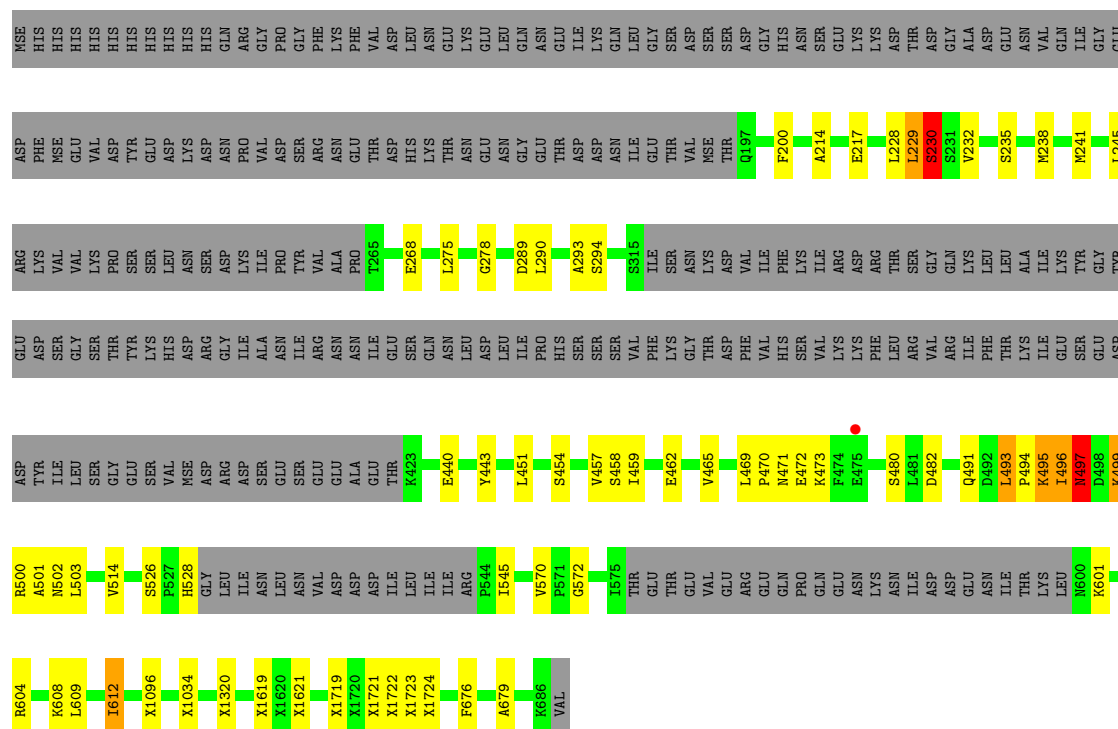
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	B	9	Total 9	Se 9	1	0
8	C	4	Total 4	Se 4	1	0
8	D	3	Total 3	Se 3	0	0
8	E	6	Total 6	Se 6	2	0
8	G	5	Total 5	Se 5	1	0
8	H	3	Total 3	Se 3	2	0
8	I	10	Total 10	Se 10	3	0
8	J	4	Total 4	Se 4	3	0
8	K	4	Total 4	Se 4	0	0
8	L	7	Total 7	Se 7	0	0
8	M	3	Total 3	Se 3	3	0
8	N	5	Total 5	Se 5	1	0
8	O	3	Total 3	Se 3	0	0
8	P	9	Total 9	Se 9	1	0
8	Q	4	Total 4	Se 4	0	0
8	R	2	Total 2	Se 2	0	0
8	S	6	Total 6	Se 6	0	0
8	T	2	Total 2	Se 2	1	0
8	U	6	Total 6	Se 6	2	0



• Molecule 2: Mediator of RNA polymerase II transcription subunit 17

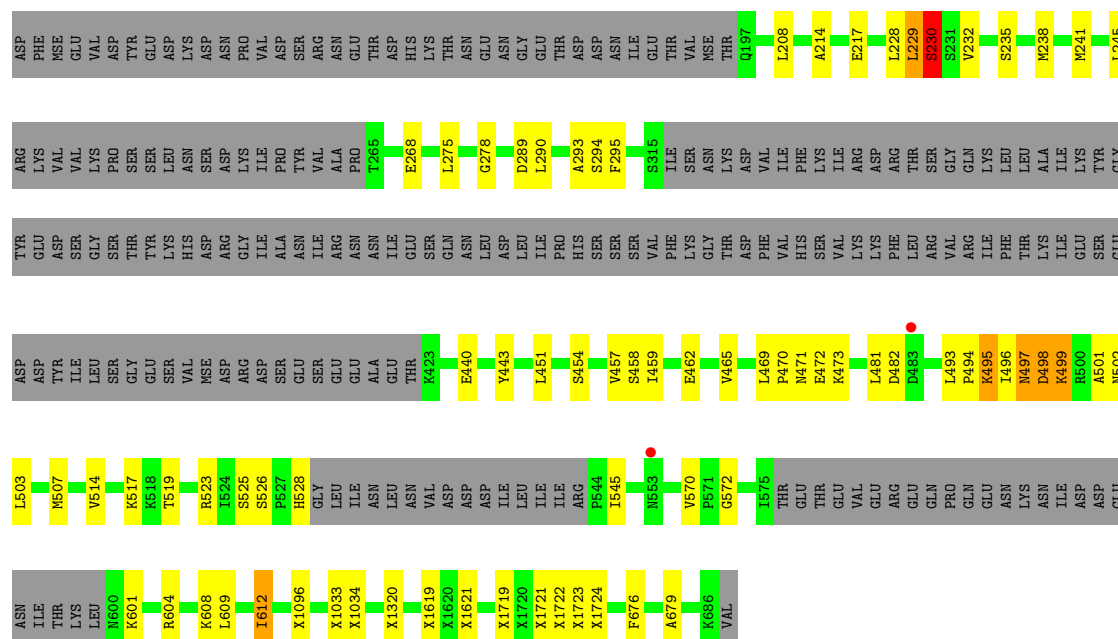
Chain I: 43% 10% 45%



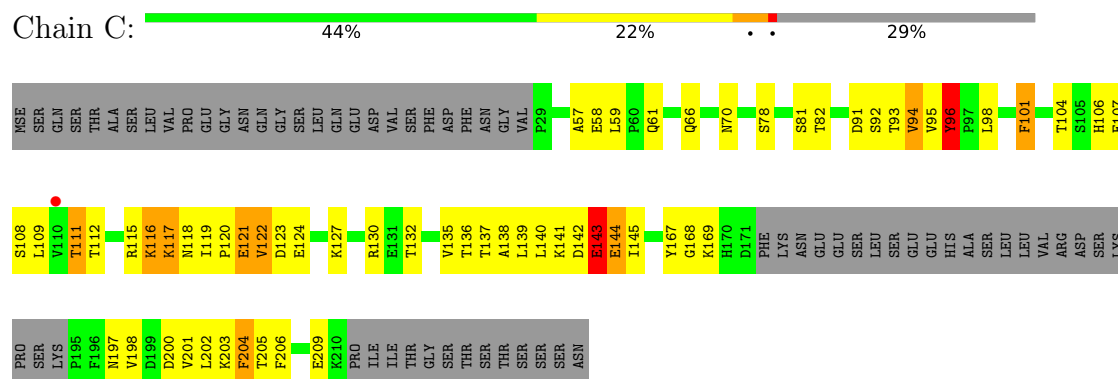
• Molecule 2: Mediator of RNA polymerase II transcription subunit 17

Chain P: 42% 11% 45%

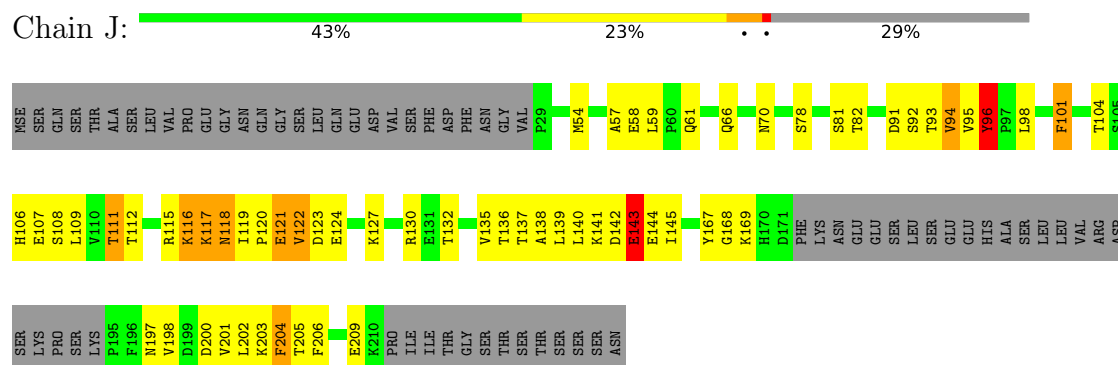




- Molecule 3: Mediator of RNA polymerase II transcription subunit 8

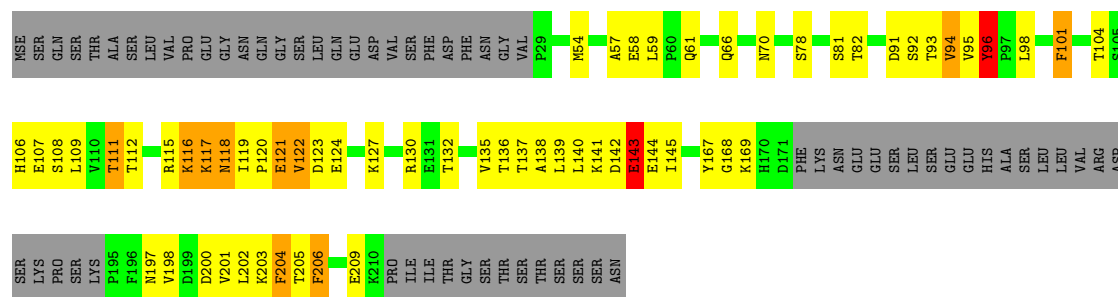


- Molecule 3: Mediator of RNA polymerase II transcription subunit 8

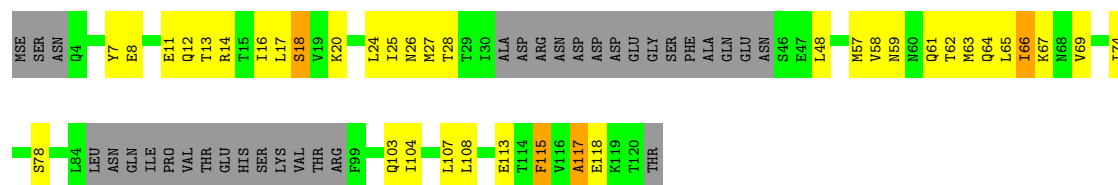


- Molecule 3: Mediator of RNA polymerase II transcription subunit 8

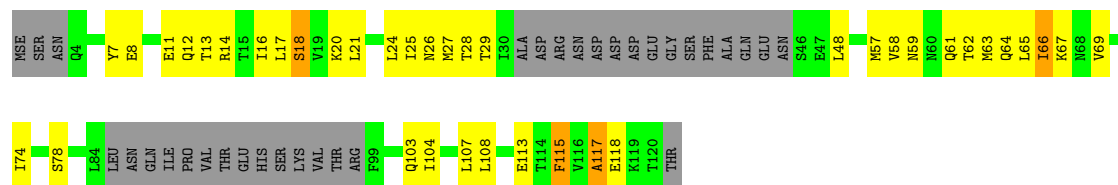




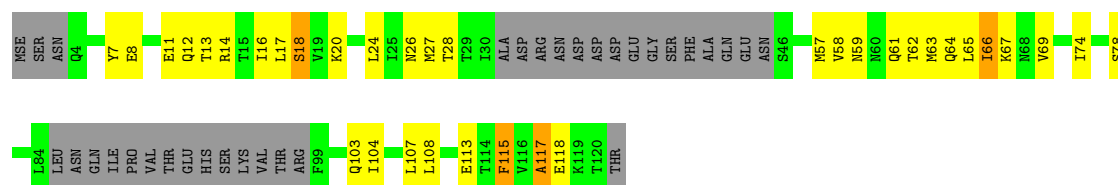
- Molecule 4: Mediator of RNA polymerase II transcription subunit 22



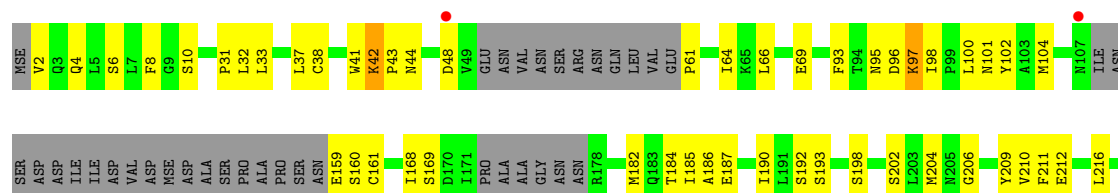
- Molecule 4: Mediator of RNA polymerase II transcription subunit 22



- Molecule 4: Mediator of RNA polymerase II transcription subunit 22



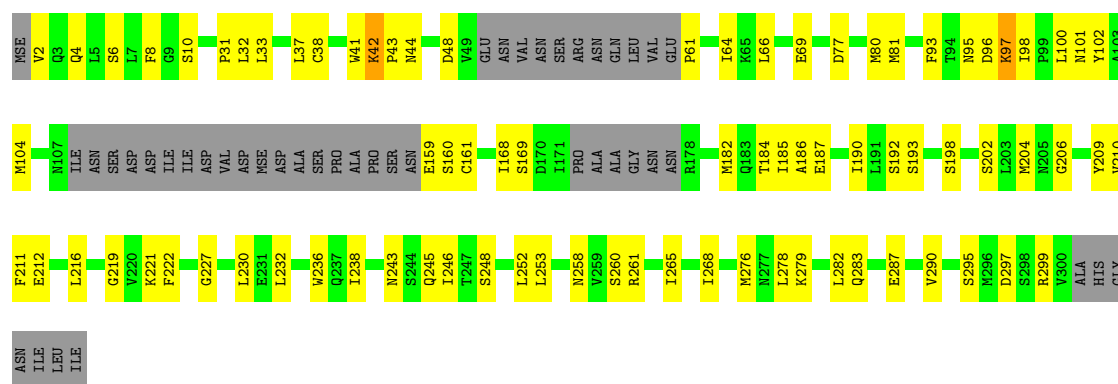
- Molecule 5: Mediator of RNA polymerase II transcription subunit 18





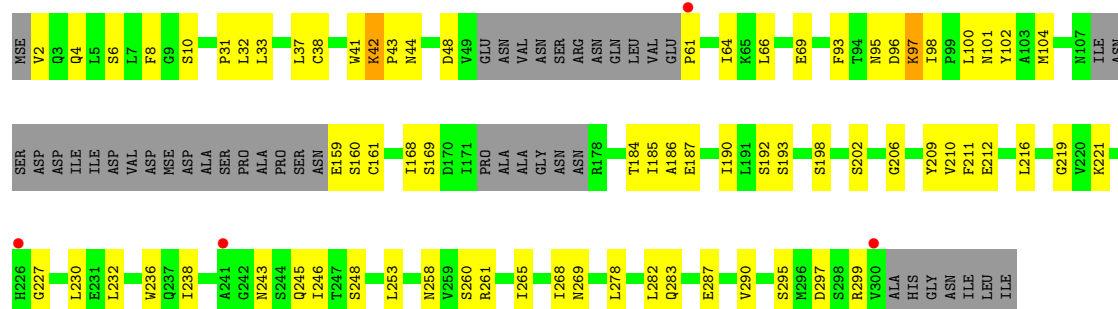
• Molecule 5: Mediator of RNA polymerase II transcription subunit 18

Chain L: 54% 29% 16%



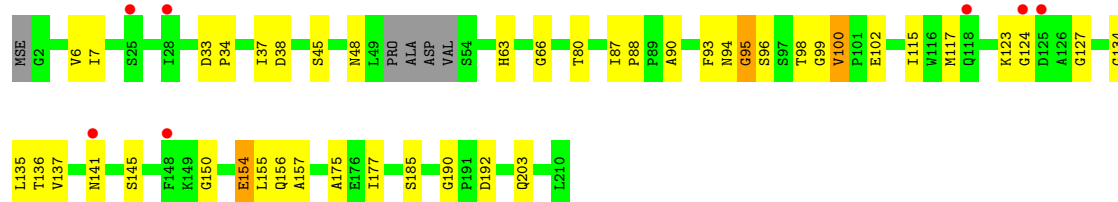
• Molecule 5: Mediator of RNA polymerase II transcription subunit 18

Chain S: 57% 26% 16%



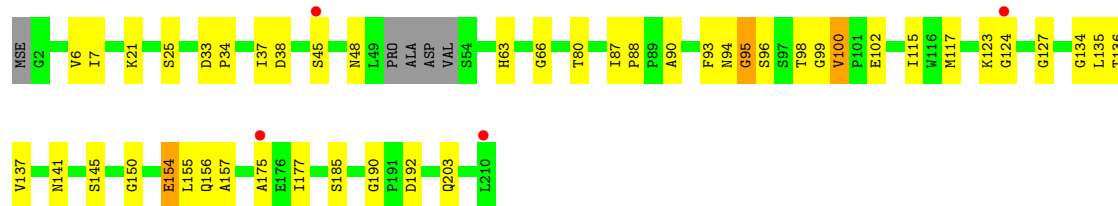
• Molecule 6: Mediator of RNA polymerase II transcription subunit 20

Chain F: 3% 77% 20% 2%

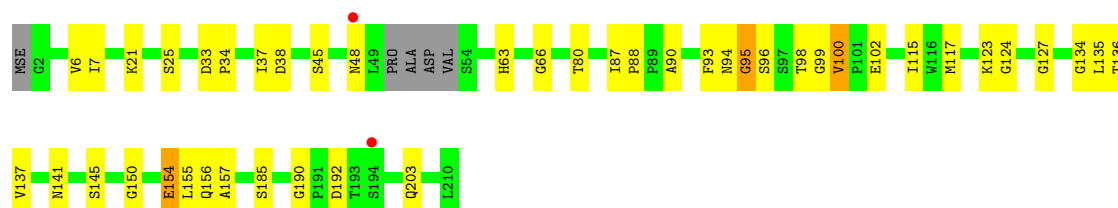
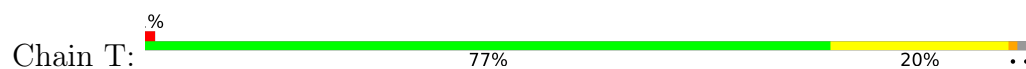


• Molecule 6: Mediator of RNA polymerase II transcription subunit 20

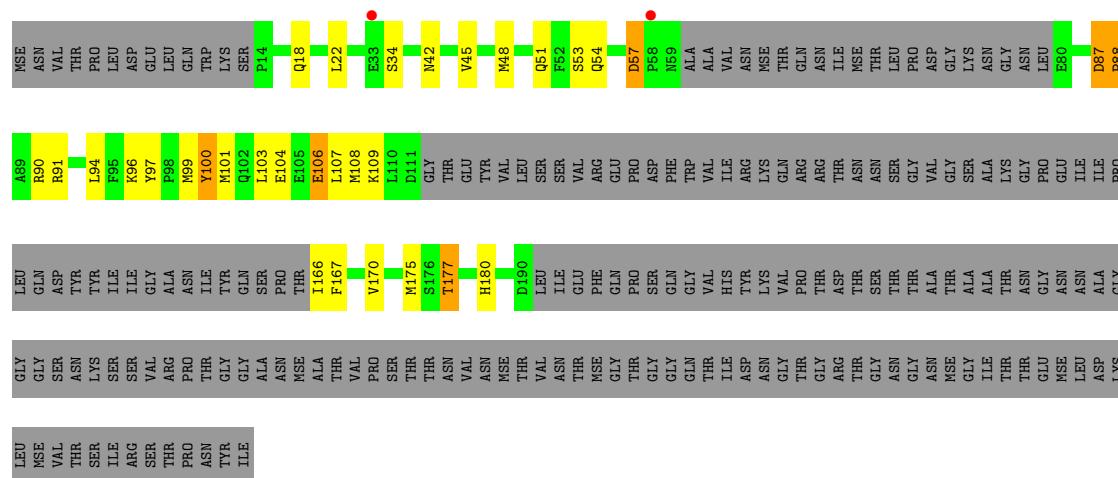
Chain M: 2% 76% 20% 2%



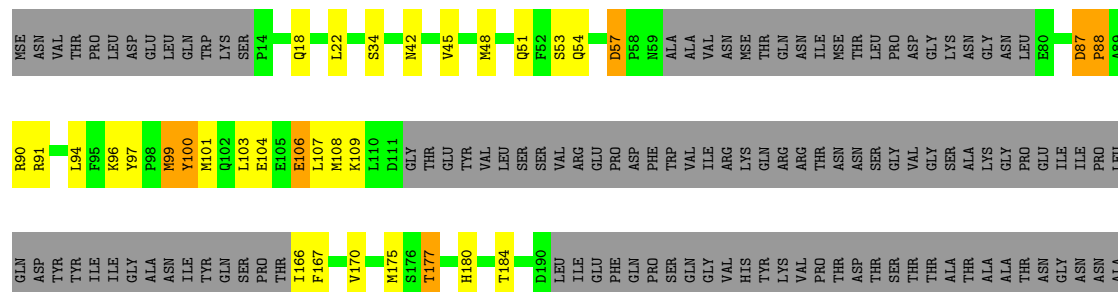
• Molecule 6: Mediator of RNA polymerase II transcription subunit 20



• Molecule 7: Mediator of RNA polymerase II transcription subunit 6



• Molecule 7: Mediator of RNA polymerase II transcription subunit 6



4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	265.91Å 265.91Å 319.95Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	49.91 – 4.30 49.91 – 4.30	Depositor EDS
% Data completeness (in resolution range)	92.6 (49.91-4.30) 92.6 (49.91-4.30)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.82 (at 4.29Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7_650)	Depositor
R, R_{free}	0.345 , 0.373 0.341 , 0.340	Depositor DCC
R_{free} test set	4144 reflections (4.65%)	wwPDB-VP
Wilson B-factor (Å ²)	145.9	Xtriage
Anisotropy	0.048	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.21 , 795.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.29$, $\langle L^2 \rangle = 0.13$	Xtriage
Estimated twinning fraction	0.257 for -h,-k,l	Xtriage
F_o, F_c correlation	0.86	EDS
Total number of atoms	17945	wwPDB-VP
Average B, all atoms (Å ²)	130.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.99% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: SE

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	0.61	0/464	1.21	4/642 (0.6%)
1	H	0.61	0/464	1.21	4/642 (0.6%)
1	O	0.61	0/464	1.21	4/642 (0.6%)
2	B	0.65	0/1346	1.57	22/1859 (1.2%)
2	I	0.65	0/1346	1.58	21/1859 (1.1%)
2	P	0.63	0/1346	1.59	24/1859 (1.3%)
3	C	0.71	0/786	1.90	24/1090 (2.2%)
3	J	0.71	0/786	1.90	24/1090 (2.2%)
3	Q	0.71	0/786	1.90	24/1090 (2.2%)
4	D	0.69	0/433	1.30	8/595 (1.3%)
4	K	0.69	0/433	1.30	8/595 (1.3%)
4	R	0.69	0/433	1.30	8/595 (1.3%)
5	E	0.64	0/1129	1.44	15/1553 (1.0%)
5	L	0.64	0/1129	1.44	15/1553 (1.0%)
5	S	0.64	0/1129	1.44	15/1553 (1.0%)
6	F	0.65	0/1007	1.41	14/1394 (1.0%)
6	M	0.65	0/1007	1.41	14/1394 (1.0%)
6	T	0.65	0/1007	1.41	14/1394 (1.0%)
7	G	0.60	0/506	1.84	13/695 (1.9%)
7	N	0.59	0/506	1.84	13/695 (1.9%)
7	U	0.60	0/506	1.84	13/695 (1.9%)
All	All	0.65	0/17013	1.55	301/23484 (1.3%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
2	B	0	2
2	I	0	2

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Mol	Chain	#Chirality outliers	#Planarity outliers
2	P	0	2
3	C	0	4
3	J	0	4
3	Q	0	4
5	E	0	1
5	L	0	1
5	S	0	1
7	G	0	2
7	N	0	2
7	U	0	2
All	All	0	27

There are no bond length outliers.

All (301) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	98	LEU	CA-C-N	23.71	149.48	119.84
3	C	98	LEU	C-N-CA	23.71	149.48	119.84
3	J	98	LEU	CA-C-N	23.70	149.46	119.84
3	J	98	LEU	C-N-CA	23.70	149.46	119.84
3	Q	98	LEU	CA-C-N	23.66	149.41	119.84
3	Q	98	LEU	C-N-CA	23.66	149.41	119.84
7	U	87	ASP	CA-C-N	19.86	144.67	119.84
7	U	87	ASP	C-N-CA	19.86	144.67	119.84
7	N	87	ASP	CA-C-N	19.86	144.67	119.84
7	N	87	ASP	C-N-CA	19.86	144.67	119.84
7	G	87	ASP	CA-C-N	19.84	144.64	119.84
7	G	87	ASP	C-N-CA	19.84	144.64	119.84
2	P	526	SER	CA-C-N	18.43	138.17	119.56
2	P	526	SER	C-N-CA	18.43	138.17	119.56
2	I	526	SER	CA-C-N	18.41	138.16	119.56
2	I	526	SER	C-N-CA	18.41	138.16	119.56
2	B	526	SER	CA-C-N	18.40	138.15	119.56
2	B	526	SER	C-N-CA	18.40	138.15	119.56
3	C	121	GLU	N-CA-C	-14.45	95.53	111.28
3	J	121	GLU	N-CA-C	-14.43	95.55	111.28
3	Q	121	GLU	N-CA-C	-14.43	95.55	111.28
5	E	290	VAL	CA-C-N	13.98	133.87	120.03
5	E	290	VAL	C-N-CA	13.98	133.87	120.03
5	S	290	VAL	CA-C-N	13.95	133.84	120.03
5	S	290	VAL	C-N-CA	13.95	133.84	120.03
5	L	290	VAL	CA-C-N	13.94	133.83	120.03

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	L	290	VAL	C-N-CA	13.94	133.83	120.03
2	I	495	LYS	N-CA-C	-13.42	95.10	111.33
2	B	570	VAL	CA-C-N	13.38	133.78	119.87
2	B	570	VAL	C-N-CA	13.38	133.78	119.87
2	I	570	VAL	CA-C-N	13.38	133.78	119.87
2	I	570	VAL	C-N-CA	13.38	133.78	119.87
2	P	570	VAL	CA-C-N	13.34	133.75	119.87
2	P	570	VAL	C-N-CA	13.34	133.75	119.87
2	P	495	LYS	N-CA-C	-12.80	97.33	111.28
5	L	98	ILE	CA-C-N	12.60	132.59	118.85
5	L	98	ILE	C-N-CA	12.60	132.59	118.85
5	S	98	ILE	CA-C-N	12.60	132.58	118.85
5	S	98	ILE	C-N-CA	12.60	132.58	118.85
5	E	98	ILE	CA-C-N	12.59	132.57	118.85
5	E	98	ILE	C-N-CA	12.59	132.57	118.85
2	B	612	ILE	CA-C-N	12.26	131.72	118.97
2	B	612	ILE	C-N-CA	12.26	131.72	118.97
2	P	612	ILE	CA-C-N	12.26	131.72	118.97
2	P	612	ILE	C-N-CA	12.26	131.72	118.97
2	I	612	ILE	CA-C-N	12.25	131.71	118.97
2	I	612	ILE	C-N-CA	12.25	131.71	118.97
2	B	495	LYS	N-CA-C	-11.77	98.45	111.28
6	F	33	ASP	CA-C-N	11.12	131.58	119.90
6	F	33	ASP	C-N-CA	11.12	131.58	119.90
6	T	33	ASP	CA-C-N	11.11	131.57	119.90
6	T	33	ASP	C-N-CA	11.11	131.57	119.90
3	J	59	LEU	CA-C-N	11.09	131.54	119.90
3	J	59	LEU	C-N-CA	11.09	131.54	119.90
6	M	33	ASP	CA-C-N	11.08	131.54	119.90
6	M	33	ASP	C-N-CA	11.08	131.54	119.90
3	C	59	LEU	CA-C-N	11.07	131.53	119.90
3	C	59	LEU	C-N-CA	11.07	131.53	119.90
3	Q	59	LEU	CA-C-N	11.05	131.51	119.90
3	Q	59	LEU	C-N-CA	11.05	131.51	119.90
7	U	100	TYR	N-CA-C	-10.33	100.16	113.17
7	N	100	TYR	N-CA-C	-10.31	100.17	113.17
7	G	100	TYR	N-CA-C	-10.30	100.20	113.17
2	P	498	ASP	N-CA-C	-9.84	101.21	113.02
7	U	57	ASP	CA-C-N	9.17	128.91	119.56
7	U	57	ASP	C-N-CA	9.17	128.91	119.56
7	G	57	ASP	CA-C-N	9.15	128.90	119.56
7	G	57	ASP	C-N-CA	9.15	128.90	119.56

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	N	57	ASP	CA-C-N	9.14	128.88	119.56
7	N	57	ASP	C-N-CA	9.14	128.88	119.56
6	T	150	GLY	N-CA-C	9.03	121.84	110.29
6	M	150	GLY	N-CA-C	9.02	121.84	110.29
6	F	150	GLY	N-CA-C	9.01	121.82	110.29
5	S	42	LYS	CA-C-N	-8.88	110.42	119.92
5	S	42	LYS	C-N-CA	-8.88	110.42	119.92
5	E	42	LYS	CA-C-N	-8.87	110.43	119.92
5	E	42	LYS	C-N-CA	-8.87	110.43	119.92
3	Q	93	THR	N-CA-C	-8.86	100.42	111.75
5	L	42	LYS	CA-C-N	-8.83	110.47	119.92
5	L	42	LYS	C-N-CA	-8.83	110.47	119.92
3	C	93	THR	N-CA-C	-8.82	100.46	111.75
3	J	93	THR	N-CA-C	-8.82	100.46	111.75
6	T	100	VAL	CA-C-N	8.63	128.67	119.78
6	T	100	VAL	C-N-CA	8.63	128.67	119.78
6	M	100	VAL	CA-C-N	8.62	128.66	119.78
6	M	100	VAL	C-N-CA	8.62	128.66	119.78
6	F	100	VAL	CA-C-N	8.57	128.61	119.78
6	F	100	VAL	C-N-CA	8.57	128.61	119.78
7	U	167	PHE	N-CA-C	-7.97	103.56	113.20
7	N	167	PHE	N-CA-C	-7.94	103.59	113.20
7	G	167	PHE	N-CA-C	-7.92	103.62	113.20
7	U	97	TYR	CA-C-N	7.71	126.98	118.97
7	U	97	TYR	C-N-CA	7.71	126.98	118.97
7	G	97	TYR	CA-C-N	7.69	126.97	118.97
7	G	97	TYR	C-N-CA	7.69	126.97	118.97
7	N	97	TYR	CA-C-N	7.66	126.94	118.97
7	N	97	TYR	C-N-CA	7.66	126.94	118.97
2	P	482	ASP	N-CA-C	-7.63	102.61	112.23
2	I	482	ASP	N-CA-C	-7.62	102.62	112.23
2	B	482	ASP	N-CA-C	-7.61	102.64	112.23
3	Q	143	GLU	CB-CA-C	-7.24	108.23	116.63
5	E	6	SER	N-CA-C	7.24	119.86	108.79
5	L	6	SER	N-CA-C	7.23	119.86	108.79
5	S	6	SER	N-CA-C	7.23	119.85	108.79
3	C	143	GLU	CB-CA-C	-7.22	108.25	116.63
3	J	143	GLU	CB-CA-C	-7.21	108.27	116.63
3	J	101	PHE	CA-C-N	7.17	128.81	119.84
3	J	101	PHE	C-N-CA	7.17	128.81	119.84
3	C	101	PHE	CA-C-N	7.14	128.76	119.84
3	C	101	PHE	C-N-CA	7.14	128.76	119.84

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	Q	101	PHE	CA-C-N	7.12	128.74	119.84
3	Q	101	PHE	C-N-CA	7.12	128.74	119.84
7	G	96	LYS	N-CA-C	7.01	121.43	113.02
2	P	545	ILE	N-CA-C	6.99	118.86	112.29
7	N	96	LYS	N-CA-C	6.97	121.39	113.02
2	I	545	ILE	N-CA-C	6.97	118.84	112.29
1	O	126	ASP	CA-C-N	6.97	126.22	118.97
1	O	126	ASP	C-N-CA	6.97	126.22	118.97
7	U	96	LYS	N-CA-C	6.97	121.38	113.02
1	A	126	ASP	CA-C-N	6.96	126.21	118.97
1	A	126	ASP	C-N-CA	6.96	126.21	118.97
1	H	126	ASP	CA-C-N	6.96	126.21	118.97
1	H	126	ASP	C-N-CA	6.96	126.21	118.97
2	B	545	ILE	N-CA-C	6.92	118.79	112.29
3	J	206	PHE	N-CA-C	-6.90	104.86	113.28
3	Q	206	PHE	N-CA-C	-6.89	104.87	113.28
3	C	206	PHE	N-CA-C	-6.88	104.89	113.28
4	R	13	THR	N-CA-C	-6.82	104.99	113.38
7	G	34	SER	CA-C-N	6.81	126.41	119.19
7	G	34	SER	C-N-CA	6.81	126.41	119.19
4	K	13	THR	N-CA-C	-6.81	105.00	113.38
3	J	111	THR	N-CA-C	-6.81	104.97	113.28
3	Q	111	THR	N-CA-C	-6.79	105.00	113.28
4	D	13	THR	N-CA-C	-6.78	105.04	113.38
3	C	111	THR	N-CA-C	-6.77	105.02	113.28
7	N	34	SER	CA-C-N	6.75	126.34	119.19
7	N	34	SER	C-N-CA	6.75	126.34	119.19
7	U	34	SER	CA-C-N	6.74	126.34	119.19
7	U	34	SER	C-N-CA	6.74	126.34	119.19
2	B	232	VAL	CA-C-N	6.66	134.01	122.56
2	B	232	VAL	C-N-CA	6.66	134.01	122.56
2	B	572	GLY	N-CA-C	6.66	122.22	114.16
2	P	572	GLY	N-CA-C	6.65	122.20	114.16
2	P	232	VAL	CA-C-N	6.64	133.99	122.56
2	P	232	VAL	C-N-CA	6.64	133.99	122.56
2	I	232	VAL	CA-C-N	6.64	133.98	122.56
2	I	232	VAL	C-N-CA	6.64	133.98	122.56
3	C	117	LYS	CB-CA-C	-6.63	108.22	117.23
2	I	572	GLY	N-CA-C	6.63	122.18	114.16
2	B	493	LEU	CA-C-N	6.62	126.21	119.19
2	B	493	LEU	C-N-CA	6.62	126.21	119.19
5	S	8	PHE	N-CA-C	6.61	118.24	108.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	J	117	LYS	CB-CA-C	-6.60	108.25	117.23
5	L	8	PHE	N-CA-C	6.59	118.23	108.60
5	E	8	PHE	N-CA-C	6.59	118.22	108.60
3	Q	117	LYS	CB-CA-C	-6.58	108.28	117.23
3	J	118	ASN	CA-C-N	-6.50	117.67	123.33
3	J	118	ASN	C-N-CA	-6.50	117.67	123.33
3	C	118	ASN	CA-C-N	-6.46	117.71	123.33
3	C	118	ASN	C-N-CA	-6.46	117.71	123.33
3	Q	118	ASN	CA-C-N	-6.45	117.72	123.33
3	Q	118	ASN	C-N-CA	-6.45	117.72	123.33
2	I	499	LYS	N-CA-C	-6.37	105.16	113.12
4	R	7	TYR	N-CA-C	-6.34	105.03	112.89
4	D	7	TYR	N-CA-C	-6.32	105.05	112.89
4	K	7	TYR	N-CA-C	-6.30	105.08	112.89
2	B	499	LYS	N-CA-C	-6.27	103.54	112.13
6	M	48	ASN	N-CA-C	6.26	119.13	108.67
6	T	48	ASN	N-CA-C	6.26	119.13	108.67
6	F	48	ASN	N-CA-C	6.26	119.12	108.67
3	J	121	GLU	CA-C-N	-6.14	110.91	121.97
3	J	121	GLU	C-N-CA	-6.14	110.91	121.97
3	Q	121	GLU	CA-C-N	-6.14	110.92	121.97
3	Q	121	GLU	C-N-CA	-6.14	110.92	121.97
3	C	121	GLU	CA-C-N	-6.12	110.95	121.97
3	C	121	GLU	C-N-CA	-6.12	110.95	121.97
2	P	493	LEU	CA-C-N	6.07	126.18	119.87
2	P	493	LEU	C-N-CA	6.07	126.18	119.87
2	I	497	ASN	N-CA-C	6.06	123.71	110.80
3	Q	116	LYS	CA-C-N	-6.04	114.22	123.04
3	Q	116	LYS	C-N-CA	-6.04	114.22	123.04
6	F	145	SER	N-CA-C	-6.04	105.87	113.18
5	S	96	ASP	N-CA-C	-6.04	106.17	112.93
6	M	145	SER	N-CA-C	-6.03	105.89	113.18
3	J	116	LYS	CA-C-N	-6.03	114.24	123.04
3	J	116	LYS	C-N-CA	-6.03	114.24	123.04
6	T	145	SER	N-CA-C	-6.02	105.89	113.18
5	L	96	ASP	N-CA-C	-6.02	106.19	112.93
5	E	96	ASP	N-CA-C	-6.01	106.20	112.93
3	C	116	LYS	CA-C-N	-6.00	114.28	123.04
3	C	116	LYS	C-N-CA	-6.00	114.28	123.04
3	C	119	ILE	N-CA-C	6.00	114.75	107.61
2	P	497	ASN	N-CA-C	5.97	119.88	112.24
2	P	499	LYS	N-CA-C	-5.97	104.78	111.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	Q	119	ILE	N-CA-C	5.96	114.71	107.61
3	J	119	ILE	N-CA-C	5.96	114.70	107.61
4	K	117	ALA	N-CA-C	-5.93	105.70	113.17
4	D	117	ALA	N-CA-C	-5.93	105.70	113.17
4	R	117	ALA	N-CA-C	-5.93	105.70	113.17
6	F	95	GLY	N-CA-C	-5.91	107.79	115.59
6	M	95	GLY	N-CA-C	-5.90	107.81	115.59
6	T	95	GLY	N-CA-C	-5.89	107.81	115.59
3	Q	119	ILE	CA-C-O	5.86	123.14	119.51
4	R	18	SER	N-CA-C	-5.85	105.03	111.82
6	T	117	MSE	N-CA-C	-5.85	101.07	110.14
4	K	18	SER	N-CA-C	-5.83	105.05	111.82
6	F	117	MSE	N-CA-C	-5.83	101.10	110.14
6	M	117	MSE	N-CA-C	-5.83	101.11	110.14
3	J	119	ILE	CA-C-O	5.82	123.12	119.51
4	D	18	SER	N-CA-C	-5.82	105.07	111.82
4	R	115	PHE	N-CA-C	5.81	119.56	112.23
3	C	119	ILE	CA-C-O	5.80	123.11	119.51
4	K	115	PHE	N-CA-C	5.79	119.52	112.23
4	D	115	PHE	N-CA-C	5.79	119.52	112.23
2	P	499	LYS	CA-C-N	-5.77	109.76	121.18
2	P	499	LYS	C-N-CA	-5.77	109.76	121.18
6	T	99	GLY	N-CA-C	-5.62	107.98	115.40
2	I	496	ILE	N-CA-C	-5.60	100.25	108.54
6	M	99	GLY	N-CA-C	-5.60	108.01	115.40
2	B	469	LEU	CA-C-N	5.59	126.07	120.14
2	B	469	LEU	C-N-CA	5.59	126.07	120.14
2	P	469	LEU	CA-C-N	5.58	126.06	120.14
2	P	469	LEU	C-N-CA	5.58	126.06	120.14
6	F	99	GLY	N-CA-C	-5.58	108.04	115.40
5	E	104	MSE	CB-CA-C	-5.57	110.17	116.63
2	I	469	LEU	CA-C-N	5.57	126.04	120.14
2	I	469	LEU	C-N-CA	5.57	126.04	120.14
5	L	104	MSE	CB-CA-C	-5.57	110.17	116.63
5	S	104	MSE	CB-CA-C	-5.53	110.22	116.63
3	Q	61	GLN	N-CA-C	5.49	118.36	108.48
6	T	7	ILE	N-CA-C	5.49	115.54	107.75
3	C	61	GLN	N-CA-C	5.48	118.35	108.48
3	J	61	GLN	N-CA-C	5.48	118.34	108.48
6	M	6	VAL	N-CA-C	5.48	115.53	107.75
6	T	6	VAL	N-CA-C	5.47	115.52	107.75
6	F	6	VAL	N-CA-C	5.47	115.51	107.75

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	168	GLY	N-CA-C	-5.47	108.53	115.31
6	M	7	ILE	N-CA-C	5.45	115.49	107.75
5	E	97	LYS	N-CA-C	5.44	119.41	112.34
6	F	7	ILE	N-CA-C	5.43	115.47	107.75
3	Q	168	GLY	N-CA-C	-5.43	108.58	115.31
3	J	168	GLY	N-CA-C	-5.42	108.58	115.31
2	B	499	LYS	CA-C-N	-5.42	111.19	121.54
2	B	499	LYS	C-N-CA	-5.42	111.19	121.54
5	L	97	LYS	N-CA-C	5.42	119.38	112.34
5	S	97	LYS	N-CA-C	5.41	119.37	112.34
2	P	465	VAL	N-CA-C	5.40	116.66	109.37
3	C	204	PHE	N-CA-C	-5.38	106.22	112.89
3	Q	204	PHE	N-CA-C	-5.37	106.23	112.89
2	I	465	VAL	N-CA-C	5.37	116.62	109.37
2	B	465	VAL	N-CA-C	5.37	116.62	109.37
3	J	204	PHE	N-CA-C	-5.37	106.23	112.89
6	F	48	ASN	CB-CA-C	-5.36	109.94	117.23
6	M	48	ASN	CB-CA-C	-5.34	109.97	117.23
3	Q	132	THR	N-CA-C	-5.31	105.53	112.23
6	T	48	ASN	CB-CA-C	-5.31	110.00	117.23
3	J	132	THR	N-CA-C	-5.31	105.54	112.23
5	L	93	PHE	N-CA-C	5.31	117.06	108.41
3	C	132	THR	N-CA-C	-5.30	105.56	112.23
6	T	154	GLU	N-CA-C	5.30	117.73	109.52
6	M	154	GLU	N-CA-C	5.29	117.72	109.52
5	E	93	PHE	N-CA-C	5.29	117.03	108.41
5	S	93	PHE	N-CA-C	5.28	117.02	108.41
6	F	154	GLU	N-CA-C	5.28	117.70	109.52
2	P	268	GLU	N-CA-C	5.27	115.08	108.45
2	B	268	GLU	N-CA-C	5.26	115.08	108.45
2	I	457	VAL	N-CA-C	5.26	115.67	107.99
1	A	53	LEU	N-CA-C	-5.25	106.94	113.19
2	B	457	VAL	N-CA-C	5.25	115.66	107.99
2	P	457	VAL	N-CA-C	5.25	115.66	107.99
2	I	268	GLU	N-CA-C	5.24	115.05	108.45
1	H	53	LEU	N-CA-C	-5.23	106.97	113.19
2	I	493	LEU	O-C-N	5.22	124.50	120.38
4	K	58	VAL	N-CA-C	-5.22	105.31	111.00
4	K	103	GLN	CB-CA-C	-5.22	110.58	116.63
5	S	10	SER	N-CA-C	5.21	116.99	109.07
4	D	103	GLN	CB-CA-C	-5.21	110.59	116.63
1	O	53	LEU	N-CA-C	-5.20	107.00	113.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	R	58	VAL	N-CA-C	-5.20	105.33	111.00
5	E	10	SER	N-CA-C	5.20	116.97	109.07
5	L	10	SER	N-CA-C	5.20	116.97	109.07
4	D	58	VAL	N-CA-C	-5.19	105.34	111.00
4	R	103	GLN	CB-CA-C	-5.18	110.62	116.63
5	L	48	ASP	N-CA-C	5.13	116.92	110.91
5	S	48	ASP	N-CA-C	5.13	116.92	110.91
5	S	187	GLU	N-CA-C	5.13	117.08	109.69
5	E	48	ASP	N-CA-C	5.13	116.91	110.91
5	L	187	GLU	N-CA-C	5.12	117.07	109.69
5	E	187	GLU	N-CA-C	5.12	117.06	109.69
1	A	117	GLN	N-CA-C	-5.10	106.83	113.16
7	N	106	GLU	N-CA-C	-5.10	107.06	113.28
7	U	106	GLU	N-CA-C	-5.10	107.06	113.28
1	H	117	GLN	N-CA-C	-5.10	106.84	113.16
7	G	106	GLU	N-CA-C	-5.09	107.07	113.28
1	O	117	GLN	N-CA-C	-5.08	106.86	113.16
4	D	66	ILE	N-CA-C	-5.07	106.97	112.80
4	R	66	ILE	N-CA-C	-5.06	106.98	112.80
4	K	66	ILE	N-CA-C	-5.05	106.99	112.80
7	U	177	THR	N-CA-C	-5.04	106.64	112.89
7	N	177	THR	N-CA-C	-5.03	106.65	112.89
7	G	177	THR	N-CA-C	-5.03	106.65	112.89

There are no chirality outliers.

All (27) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	1722	UNK	Peptide
2	B	230	SER	Mainchain
3	C	101	PHE	Peptide
3	C	104	THR	Peptide
3	C	143	GLU	Peptide
3	C	96	TYR	Peptide
5	E	69	GLU	Peptide
7	G	87	ASP	Peptide
7	G	99	MSE	Peptide
2	I	1722	UNK	Peptide
2	I	230	SER	Mainchain
3	J	101	PHE	Peptide
3	J	104	THR	Peptide
3	J	143	GLU	Peptide

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Mol	Chain	Res	Type	Group
3	J	96	TYR	Peptide
5	L	69	GLU	Peptide
7	N	87	ASP	Peptide
7	N	99	MSE	Peptide
2	P	1722	UNK	Peptide
2	P	230	SER	Mainchain
3	Q	101	PHE	Peptide
3	Q	104	THR	Peptide
3	Q	143	GLU	Peptide
3	Q	96	TYR	Peptide
5	S	69	GLU	Peptide
7	U	87	ASP	Peptide
7	U	99	MSE	Peptide

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	468	0	192	20	0
1	H	468	0	192	19	0
1	O	468	0	192	18	0
2	B	1583	0	628	39	0
2	I	1583	0	629	46	0
2	P	1583	0	628	49	0
3	C	791	0	333	56	0
3	J	791	0	333	56	0
3	Q	791	0	333	59	0
4	D	440	0	188	26	0
4	K	440	0	188	32	0
4	R	440	0	188	21	0
5	E	1141	0	481	53	0
5	L	1141	0	481	63	0
5	S	1141	0	481	53	0
6	F	1012	0	469	21	0
6	M	1012	0	469	22	0
6	T	1012	0	469	21	0
7	G	514	0	209	21	0
7	N	514	0	209	26	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
7	U	514	0	209	23	0
8	A	3	0	0	1	0
8	B	9	0	0	4	0
8	C	4	0	0	0	0
8	D	3	0	0	5	0
8	E	6	0	0	6	0
8	G	5	0	0	0	0
8	H	3	0	0	1	0
8	I	10	0	0	4	0
8	J	4	0	0	2	0
8	K	4	0	0	10	0
8	L	7	0	0	14	0
8	M	3	0	0	0	0
8	N	5	0	0	5	0
8	O	3	0	0	3	0
8	P	9	0	0	6	0
8	Q	4	0	0	3	0
8	R	2	0	0	0	0
8	S	6	0	0	2	0
8	T	2	0	0	0	0
8	U	6	0	0	3	0
All	All	17945	0	7501	706	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 28.

All (706) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:Q:54:MSE:HA	8:Q:325:SE:SE	1.32	1.75
3:Q:54:MSE:CA	8:Q:325:SE:SE	2.29	1.27
4:K:28:THR:HA	8:K:177:SE:SE	1.90	1.22
3:C:143:GLU:O	3:C:145:ILE:N	1.74	1.21
3:J:143:GLU:O	3:J:145:ILE:N	1.74	1.18
3:Q:143:GLU:O	3:Q:145:ILE:N	1.74	1.18
5:E:182:MSE:O	8:E:409:SE:SE	2.16	1.13
2:P:498:ASP:CB	2:P:501:ALA:HB3	1.78	1.13
3:C:136:THR:CB	3:C:137:THR:CB	2.29	1.10
3:Q:136:THR:CB	3:Q:137:THR:CB	2.29	1.10
3:J:136:THR:CB	3:J:137:THR:CB	2.29	1.09
4:D:28:THR:HA	8:D:222:SE:SE	2.03	1.09

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:Q:138:ALA:HA	3:Q:139:LEU:C	1.75	1.09
3:J:138:ALA:HA	3:J:140:LEU:N	1.69	1.08
3:C:138:ALA:HA	3:C:140:LEU:N	1.69	1.08
3:J:138:ALA:HA	3:J:139:LEU:C	1.75	1.08
3:C:138:ALA:HA	3:C:139:LEU:C	1.75	1.07
5:S:2:VAL:HA	5:S:261:ARG:HA	1.32	1.07
5:E:2:VAL:HA	5:E:261:ARG:HA	1.32	1.05
3:Q:138:ALA:HA	3:Q:140:LEU:N	1.69	1.04
5:L:2:VAL:HA	5:L:261:ARG:HA	1.32	1.04
3:J:54:MSE:HA	8:J:225:SE:SE	2.07	1.04
3:Q:135:VAL:HA	3:Q:136:THR:CB	1.89	1.03
3:J:121:GLU:O	3:J:124:GLU:N	1.94	1.01
3:Q:121:GLU:O	3:Q:124:GLU:N	1.94	1.01
3:C:121:GLU:O	3:C:124:GLU:N	1.94	1.01
3:J:135:VAL:HA	3:J:136:THR:CB	1.89	1.00
3:Q:121:GLU:O	3:Q:123:ASP:N	1.95	0.99
3:J:121:GLU:O	3:J:123:ASP:N	1.95	0.99
3:C:121:GLU:O	3:C:123:ASP:N	1.95	0.99
2:P:494:PRO:O	2:P:496:ILE:O	1.80	0.99
3:C:135:VAL:HA	3:C:136:THR:CB	1.89	0.98
2:P:498:ASP:CA	2:P:501:ALA:HB3	1.93	0.97
2:B:499:LYS:O	2:B:502:ASN:N	1.96	0.97
7:N:184:THR:O	8:N:396:SE:SE	2.32	0.96
3:Q:78:SER:CB	8:Q:327:SE:SE	2.63	0.96
2:I:480:SER:HA	8:I:906:SE:SE	2.16	0.96
4:D:48:LEU:O	8:D:1301:SE:SE	2.35	0.95
4:D:57:MSE:CB	8:D:222:SE:SE	2.65	0.94
2:P:498:ASP:C	2:P:501:ALA:HB3	1.94	0.93
3:C:121:GLU:O	3:C:122:VAL:C	2.12	0.92
1:O:39:MSE:O	8:O:132:SE:SE	2.37	0.92
3:J:57:ALA:HB1	3:J:58:GLU:HA	1.52	0.92
3:Q:121:GLU:O	3:Q:122:VAL:C	2.12	0.91
5:L:80:MSE:HA	8:L:408:SE:SE	2.20	0.91
6:M:135:LEU:HA	6:M:156:GLN:O	1.71	0.91
6:T:135:LEU:HA	6:T:156:GLN:O	1.71	0.91
6:F:135:LEU:HA	6:F:156:GLN:O	1.71	0.91
3:C:57:ALA:HB1	3:C:58:GLU:HA	1.52	0.91
3:Q:57:ALA:HB1	3:Q:58:GLU:HA	1.52	0.91
3:C:141:LYS:CB	3:C:145:ILE:CB	2.50	0.90
5:E:41:TRP:CB	8:E:410:SE:SE	2.69	0.90
3:J:141:LYS:CB	3:J:145:ILE:CB	2.50	0.89

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:Q:141:LYS:CB	3:Q:145:ILE:CB	2.50	0.89
7:N:107:LEU:O	8:N:400:SE:SE	2.40	0.88
7:U:104:GLU:O	8:U:499:SE:SE	2.43	0.87
4:K:28:THR:CA	8:K:177:SE:SE	2.72	0.86
3:J:121:GLU:O	3:J:122:VAL:C	2.12	0.86
2:P:498:ASP:O	2:P:502:ASN:N	2.06	0.84
3:J:112:THR:O	3:J:116:LYS:CB	2.26	0.83
3:Q:112:THR:O	3:Q:116:LYS:CB	2.26	0.83
4:K:57:MSE:CB	8:K:177:SE:SE	2.77	0.82
3:C:112:THR:O	3:C:116:LYS:CB	2.26	0.82
2:I:480:SER:CA	8:I:906:SE:SE	2.79	0.80
5:L:77:ASP:HA	8:L:408:SE:SE	2.32	0.80
2:I:496:ILE:O	2:I:497:ASN:CB	2.30	0.80
2:B:229:LEU:CB	8:B:1729:SE:SE	2.80	0.79
6:M:124:GLY:HA3	6:M:141:ASN:O	1.83	0.79
3:Q:135:VAL:CB	3:Q:136:THR:O	2.32	0.78
3:C:135:VAL:CB	3:C:136:THR:O	2.32	0.78
6:F:124:GLY:HA3	6:F:141:ASN:O	1.83	0.78
2:I:491:GLN:O	2:I:495:LYS:CB	2.32	0.78
2:P:498:ASP:CB	2:P:501:ALA:CB	2.61	0.78
3:J:135:VAL:CB	3:J:136:THR:O	2.32	0.78
3:C:137:THR:C	3:C:140:LEU:CB	2.57	0.78
3:Q:137:THR:C	3:Q:140:LEU:CB	2.57	0.78
3:J:137:THR:C	3:J:140:LEU:CB	2.57	0.78
5:S:186:ALA:HB1	6:T:102:GLU:CB	2.15	0.77
6:T:124:GLY:HA3	6:T:141:ASN:O	1.83	0.77
3:J:138:ALA:CA	3:J:139:LEU:C	2.56	0.77
2:P:208:LEU:HA	8:P:1728:SE:SE	2.35	0.77
5:L:186:ALA:HB1	6:M:102:GLU:CB	2.15	0.77
5:E:186:ALA:HB1	6:F:102:GLU:CB	2.15	0.76
6:F:134:GLY:O	6:F:157:ALA:HA	1.85	0.76
5:E:41:TRP:O	5:E:64:ILE:N	2.19	0.75
5:L:41:TRP:O	5:L:64:ILE:N	2.19	0.75
6:M:134:GLY:O	6:M:157:ALA:HA	1.85	0.75
3:J:117:LYS:CB	7:N:170:VAL:HA	2.17	0.75
3:Q:117:LYS:CB	7:U:170:VAL:HA	2.17	0.75
3:C:117:LYS:CB	7:G:170:VAL:HA	2.17	0.74
2:I:494:PRO:O	2:I:496:ILE:C	2.29	0.74
5:S:41:TRP:O	5:S:64:ILE:N	2.19	0.74
6:T:134:GLY:O	6:T:157:ALA:HA	1.85	0.74
7:U:189:TYR:CB	8:U:496:SE:SE	2.85	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:S:169:SER:HA	5:S:184:THR:O	1.88	0.74
5:E:182:MSE:CB	8:E:409:SE:SE	2.85	0.73
4:K:14:ARG:HA	4:K:17:LEU:CB	2.18	0.73
5:E:169:SER:HA	5:E:184:THR:O	1.88	0.73
5:L:169:SER:HA	5:L:184:THR:O	1.88	0.73
3:J:121:GLU:C	3:J:123:ASP:N	2.47	0.73
5:L:80:MSE:CA	8:L:408:SE:SE	2.86	0.73
4:R:14:ARG:HA	4:R:17:LEU:CB	2.18	0.72
3:Q:121:GLU:C	3:Q:123:ASP:N	2.47	0.72
4:D:14:ARG:HA	4:D:17:LEU:CB	2.18	0.72
2:I:493:LEU:O	2:I:496:ILE:O	2.07	0.72
2:B:241:MSE:O	2:B:245:LEU:CB	2.38	0.72
2:P:241:MSE:O	2:P:245:LEU:CB	2.38	0.72
3:Q:138:ALA:CA	3:Q:139:LEU:C	2.56	0.72
1:A:114:MSE:O	8:A:133:SE:SE	2.57	0.71
2:I:241:MSE:O	2:I:245:LEU:CB	2.38	0.71
5:L:221:LYS:HA	5:L:230:LEU:O	1.90	0.71
5:S:160:SER:HA	5:S:193:SER:O	1.90	0.71
5:E:160:SER:HA	5:E:193:SER:O	1.90	0.71
5:E:221:LYS:HA	5:E:230:LEU:O	1.90	0.71
5:S:41:TRP:HA	5:S:211:PHE:HA	1.71	0.71
5:E:41:TRP:HA	5:E:211:PHE:HA	1.71	0.71
5:L:160:SER:HA	5:L:193:SER:O	1.90	0.71
7:G:88:PRO:HA	7:G:91:ARG:CB	2.21	0.71
5:S:221:LYS:HA	5:S:230:LEU:O	1.90	0.71
7:N:88:PRO:HA	7:N:91:ARG:CB	2.21	0.70
3:C:121:GLU:C	3:C:123:ASP:N	2.47	0.70
5:S:33:LEU:HA	5:S:219:GLY:O	1.91	0.70
5:E:33:LEU:HA	5:E:219:GLY:O	1.91	0.70
5:L:33:LEU:HA	5:L:219:GLY:O	1.91	0.70
2:B:239:SER:HA	8:B:1730:SE:SE	2.42	0.70
7:N:99:MSE:CB	8:N:398:SE:SE	2.90	0.70
2:P:1621:UNK:O	2:P:1721:UNK:HA	1.92	0.69
5:L:41:TRP:HA	5:L:211:PHE:HA	1.71	0.69
3:J:200:ASP:O	5:L:295:SER:CB	2.41	0.69
2:P:601:LYS:HA	2:P:604:ARG:CB	2.23	0.69
3:Q:200:ASP:O	5:S:295:SER:CB	2.41	0.69
2:B:1621:UNK:O	2:B:1721:UNK:HA	1.92	0.69
5:L:37:LEU:HA	5:L:216:LEU:HA	1.74	0.69
1:O:92:ASN:CB	8:O:313:SE:SE	2.91	0.69
7:U:88:PRO:HA	7:U:91:ARG:CB	2.21	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:200:ASP:O	5:E:295:SER:CB	2.41	0.68
8:K:175:SE:SE	7:N:175:MSE:CB	2.91	0.68
2:I:494:PRO:O	2:I:495:LYS:C	2.36	0.68
2:B:601:LYS:HA	2:B:604:ARG:CB	2.23	0.68
2:I:1621:UNK:O	2:I:1721:UNK:HA	1.92	0.68
5:E:37:LEU:HA	5:E:216:LEU:HA	1.74	0.68
2:I:601:LYS:HA	2:I:604:ARG:CB	2.23	0.68
5:S:37:LEU:HA	5:S:216:LEU:HA	1.74	0.68
2:P:458:SER:H	2:P:462:GLU:CB	2.08	0.67
2:B:217:GLU:CB	7:G:177:THR:HA	2.25	0.67
5:L:80:MSE:H	8:L:408:SE:SE	2.27	0.67
2:P:217:GLU:CB	7:U:177:THR:HA	2.25	0.66
2:I:458:SER:H	2:I:462:GLU:CB	2.08	0.66
2:B:458:SER:H	2:B:462:GLU:CB	2.08	0.66
3:C:138:ALA:CA	3:C:139:LEU:C	2.56	0.66
1:A:65:GLU:O	1:A:69:LYS:N	2.29	0.66
1:O:65:GLU:O	1:O:69:LYS:N	2.29	0.66
1:H:65:GLU:O	1:H:69:LYS:N	2.29	0.66
2:B:496:ILE:O	2:B:497:ASN:CB	2.44	0.66
2:B:289:ASP:O	2:B:293:ALA:N	2.29	0.66
7:U:100:TYR:O	7:U:104:GLU:N	2.29	0.66
7:G:100:TYR:O	7:G:104:GLU:N	2.29	0.65
2:I:217:GLU:CB	7:N:177:THR:HA	2.25	0.65
5:E:236:TRP:HA	5:E:245:GLN:HA	1.79	0.65
2:I:289:ASP:O	2:I:293:ALA:N	2.29	0.64
5:L:204:MSE:HA	8:L:415:SE:SE	2.47	0.64
7:N:100:TYR:O	7:N:104:GLU:N	2.29	0.64
5:S:236:TRP:HA	5:S:245:GLN:HA	1.79	0.64
7:G:104:GLU:HA	7:G:107:LEU:CB	2.28	0.64
5:L:278:LEU:O	5:L:282:LEU:N	2.26	0.64
7:N:104:GLU:HA	7:N:107:LEU:CB	2.28	0.64
4:D:11:GLU:HA	4:D:14:ARG:CB	2.28	0.64
2:I:228:LEU:O	2:I:229:LEU:C	2.41	0.64
5:L:77:ASP:C	8:L:408:SE:SE	2.90	0.64
5:L:80:MSE:N	8:L:408:SE:SE	2.81	0.64
5:L:236:TRP:HA	5:L:245:GLN:HA	1.79	0.64
5:E:278:LEU:O	5:E:282:LEU:N	2.26	0.64
4:K:11:GLU:HA	4:K:14:ARG:CB	2.28	0.63
2:B:228:LEU:O	2:B:229:LEU:C	2.41	0.63
5:L:43:PRO:HA	5:L:209:TYR:HA	1.80	0.63
7:U:104:GLU:HA	7:U:107:LEU:CB	2.28	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:108:SER:O	3:C:112:THR:N	2.31	0.63
2:P:229:LEU:O	2:P:230:SER:O	2.17	0.63
3:C:106:HIS:O	3:C:108:SER:N	2.31	0.63
2:B:1619:UNK:O	2:B:1719:UNK:HA	1.99	0.63
2:P:499:LYS:O	2:P:503:LEU:N	2.30	0.63
2:P:1619:UNK:O	2:P:1719:UNK:HA	1.99	0.63
4:D:59:ASN:O	4:D:63:MSE:CB	2.47	0.63
5:E:43:PRO:HA	5:E:209:TYR:HA	1.80	0.63
2:I:1619:UNK:O	2:I:1719:UNK:HA	1.99	0.63
4:R:26:ASN:C	4:R:28:THR:H	2.07	0.63
3:C:111:THR:O	3:C:115:ARG:N	2.32	0.63
2:I:229:LEU:O	2:I:230:SER:O	2.17	0.63
4:K:59:ASN:O	4:K:63:MSE:CB	2.47	0.63
3:Q:108:SER:O	3:Q:112:THR:N	2.31	0.63
3:Q:204:PHE:O	3:Q:209:GLU:N	2.32	0.63
3:J:108:SER:O	3:J:112:THR:N	2.31	0.63
2:P:228:LEU:O	2:P:229:LEU:C	2.41	0.63
3:Q:106:HIS:O	3:Q:108:SER:N	2.31	0.63
2:I:499:LYS:O	2:I:502:ASN:N	2.31	0.62
3:J:106:HIS:O	3:J:108:SER:N	2.31	0.62
3:J:204:PHE:O	3:J:209:GLU:N	2.32	0.62
4:R:11:GLU:HA	4:R:14:ARG:CB	2.28	0.62
5:E:2:VAL:HA	5:E:261:ARG:CA	2.19	0.62
3:J:111:THR:O	3:J:115:ARG:N	2.32	0.62
4:K:26:ASN:C	4:K:28:THR:H	2.07	0.62
2:B:229:LEU:O	2:B:230:SER:O	2.17	0.62
2:P:289:ASP:O	2:P:293:ALA:N	2.29	0.62
1:H:112:GLU:O	1:H:114:MSE:N	2.33	0.62
4:R:59:ASN:O	4:R:63:MSE:CB	2.47	0.61
5:S:209:TYR:CB	8:S:310:SE:SE	2.98	0.61
1:O:112:GLU:O	1:O:114:MSE:N	2.33	0.61
5:S:2:VAL:HA	5:S:261:ARG:CA	2.19	0.61
5:L:2:VAL:HA	5:L:261:ARG:CA	2.19	0.61
4:R:104:ILE:O	4:R:108:LEU:N	2.33	0.61
5:S:43:PRO:HA	5:S:209:TYR:HA	1.81	0.61
4:D:26:ASN:C	4:D:28:THR:H	2.07	0.61
1:A:112:GLU:O	1:A:114:MSE:N	2.33	0.61
3:C:204:PHE:O	3:C:209:GLU:N	2.32	0.61
2:P:498:ASP:CA	2:P:501:ALA:CB	2.74	0.61
2:P:497:ASN:C	2:P:499:LYS:H	2.09	0.60
7:U:104:GLU:O	7:U:108:MSE:N	2.35	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:N:104:GLU:O	7:N:108:MSE:N	2.35	0.60
5:E:182:MSE:C	8:E:409:SE:SE	2.92	0.60
7:G:104:GLU:O	7:G:108:MSE:N	2.35	0.60
6:F:136:THR:O	6:F:155:LEU:HA	2.02	0.60
5:L:81:MSE:CB	8:L:409:SE:SE	3.00	0.60
4:K:104:ILE:O	4:K:108:LEU:N	2.33	0.60
3:J:91:ASP:C	3:J:94:VAL:HA	2.27	0.60
6:T:136:THR:O	6:T:155:LEU:HA	2.02	0.60
3:C:91:ASP:C	3:C:94:VAL:HA	2.27	0.60
6:F:137:VAL:HA	6:F:154:GLU:O	2.02	0.60
5:S:278:LEU:O	5:S:282:LEU:N	2.26	0.60
3:C:78:SER:HA	3:C:81:SER:CB	2.32	0.59
5:L:61:PRO:O	5:L:299:ARG:HA	2.02	0.59
6:M:136:THR:O	6:M:155:LEU:HA	2.02	0.59
5:S:61:PRO:O	5:S:299:ARG:HA	2.02	0.59
6:M:137:VAL:HA	6:M:154:GLU:O	2.02	0.59
3:C:136:THR:CB	3:C:137:THR:CA	2.80	0.59
5:E:61:PRO:O	5:E:299:ARG:HA	2.02	0.59
3:J:136:THR:CB	3:J:137:THR:CA	2.80	0.59
3:Q:91:ASP:C	3:Q:94:VAL:HA	2.27	0.59
6:T:137:VAL:HA	6:T:154:GLU:O	2.02	0.59
3:Q:111:THR:O	3:Q:115:ARG:N	2.32	0.58
3:Q:78:SER:O	3:Q:82:THR:N	2.36	0.58
5:S:44:ASN:N	5:S:209:TYR:HA	2.19	0.58
3:J:78:SER:O	3:J:82:THR:N	2.36	0.58
3:Q:78:SER:HA	3:Q:81:SER:CB	2.32	0.58
3:C:78:SER:O	3:C:82:THR:N	2.36	0.58
7:U:103:LEU:O	7:U:107:LEU:N	2.37	0.58
4:D:66:ILE:HA	4:D:69:VAL:CB	2.34	0.58
3:J:78:SER:HA	3:J:81:SER:CB	2.32	0.58
4:R:66:ILE:HA	4:R:69:VAL:CB	2.34	0.58
7:G:103:LEU:O	7:G:107:LEU:N	2.37	0.58
3:J:120:PRO:O	3:J:121:GLU:C	2.46	0.58
5:L:44:ASN:N	5:L:209:TYR:HA	2.19	0.58
2:P:494:PRO:O	2:P:495:LYS:C	2.47	0.58
2:B:491:GLN:O	2:B:495:LYS:CB	2.51	0.57
4:D:104:ILE:O	4:D:108:LEU:N	2.33	0.57
4:K:66:ILE:HA	4:K:69:VAL:CB	2.34	0.57
5:S:168:ILE:O	5:S:185:ILE:HA	2.05	0.57
5:L:77:ASP:CA	8:L:408:SE:SE	3.03	0.57
7:N:103:LEU:O	7:N:107:LEU:N	2.37	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:Q:136:THR:CB	3:Q:137:THR:CA	2.80	0.57
5:E:168:ILE:O	5:E:185:ILE:HA	2.05	0.57
5:E:204:MSE:CB	8:E:410:SE:SE	3.03	0.57
4:K:62:THR:C	8:K:174:SE:SE	2.97	0.57
5:L:168:ILE:O	5:L:185:ILE:HA	2.05	0.57
4:K:62:THR:CB	8:K:174:SE:SE	3.03	0.57
5:E:44:ASN:N	5:E:209:TYR:HA	2.19	0.56
4:K:62:THR:HA	4:K:65:LEU:CB	2.35	0.56
2:B:494:PRO:O	2:B:495:LYS:C	2.48	0.56
4:D:62:THR:HA	4:D:65:LEU:CB	2.36	0.56
5:L:278:LEU:O	5:L:282:LEU:CB	2.54	0.56
5:E:278:LEU:O	5:E:282:LEU:CB	2.54	0.56
1:A:55:ARG:HA	1:H:57:ASN:CB	2.36	0.56
3:J:54:MSE:CA	8:J:225:SE:SE	2.95	0.56
2:P:235:SER:HA	2:P:238:MSE:CB	2.36	0.56
3:Q:135:VAL:CA	3:Q:136:THR:CB	2.69	0.56
3:C:167:TYR:C	3:C:169:LYS:H	2.14	0.56
2:I:1096:UNK:CB	2:I:1621:UNK:HA	2.36	0.56
3:J:92:SER:C	3:J:94:VAL:N	2.60	0.56
4:R:62:THR:HA	4:R:65:LEU:CB	2.35	0.56
2:P:1096:UNK:CB	2:P:1621:UNK:HA	2.36	0.56
2:B:1096:UNK:CB	2:B:1621:UNK:HA	2.36	0.56
3:C:120:PRO:O	3:C:121:GLU:C	2.46	0.56
3:Q:120:PRO:O	3:Q:121:GLU:C	2.46	0.56
3:Q:167:TYR:C	3:Q:169:LYS:H	2.14	0.56
2:B:235:SER:HA	2:B:238:MSE:CB	2.36	0.55
3:J:167:TYR:C	3:J:169:LYS:H	2.14	0.55
4:D:16:ILE:O	4:D:20:LYS:N	2.39	0.55
2:I:499:LYS:C	2:I:501:ALA:N	2.64	0.55
3:C:203:LYS:O	5:E:295:SER:O	2.25	0.55
3:J:203:LYS:O	5:L:295:SER:O	2.25	0.55
4:K:16:ILE:O	4:K:20:LYS:N	2.39	0.55
5:S:278:LEU:O	5:S:282:LEU:CB	2.54	0.55
3:J:203:LYS:C	5:L:295:SER:O	2.50	0.55
4:K:113:GLU:O	4:K:117:ALA:HB3	2.07	0.55
3:Q:203:LYS:C	5:S:295:SER:O	2.50	0.55
4:D:8:GLU:O	4:D:12:GLN:CB	2.55	0.55
6:M:94:ASN:C	6:M:96:SER:H	2.15	0.55
2:I:235:SER:HA	2:I:238:MSE:CB	2.36	0.55
4:R:16:ILE:O	4:R:20:LYS:N	2.39	0.55
4:K:8:GLU:O	4:K:12:GLN:CB	2.55	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:Q:92:SER:C	3:Q:94:VAL:N	2.60	0.54
3:Q:203:LYS:O	5:S:295:SER:O	2.25	0.54
4:R:8:GLU:O	4:R:12:GLN:CB	2.55	0.54
2:P:1619:UNK:C	2:P:1719:UNK:HA	2.38	0.54
2:B:1619:UNK:C	2:B:1719:UNK:HA	2.38	0.54
4:D:113:GLU:O	4:D:117:ALA:HB3	2.07	0.54
2:I:1619:UNK:C	2:I:1719:UNK:HA	2.38	0.54
6:T:94:ASN:C	6:T:96:SER:H	2.15	0.54
2:B:229:LEU:HA	2:B:238:MSE:CB	2.38	0.54
5:E:209:TYR:O	8:E:410:SE:SE	2.76	0.54
4:K:26:ASN:C	4:K:28:THR:N	2.66	0.54
4:R:64:GLN:HA	4:R:67:LYS:CB	2.38	0.54
3:C:203:LYS:C	5:E:295:SER:O	2.50	0.54
4:K:64:GLN:HA	4:K:67:LYS:CB	2.38	0.54
5:E:31:PRO:HA	5:E:221:LYS:O	2.08	0.54
2:P:229:LEU:HA	2:P:238:MSE:CB	2.38	0.54
6:T:87:ILE:O	6:T:88:PRO:C	2.51	0.54
5:L:31:PRO:HA	5:L:221:LYS:O	2.08	0.53
4:R:113:GLU:O	4:R:117:ALA:HB3	2.07	0.53
2:I:229:LEU:HA	2:I:238:MSE:CB	2.38	0.53
4:R:26:ASN:C	4:R:28:THR:N	2.66	0.53
2:I:480:SER:CB	8:I:906:SE:SE	3.07	0.53
5:S:31:PRO:HA	5:S:221:LYS:O	2.08	0.53
5:S:190:ILE:HA	6:T:80:THR:HA	1.91	0.53
6:F:87:ILE:O	6:F:88:PRO:C	2.51	0.53
3:C:94:VAL:C	3:C:96:TYR:N	2.67	0.53
4:D:64:GLN:HA	4:D:67:LYS:CB	2.38	0.53
6:F:94:ASN:C	6:F:96:SER:H	2.15	0.53
3:J:94:VAL:C	3:J:96:TYR:N	2.67	0.53
5:E:190:ILE:HA	6:F:80:THR:HA	1.91	0.53
6:F:34:PRO:HA	6:F:127:GLY:O	2.10	0.52
4:K:62:THR:O	8:K:174:SE:SE	2.77	0.52
6:T:34:PRO:HA	6:T:127:GLY:O	2.10	0.52
7:U:104:GLU:C	8:U:499:SE:SE	3.02	0.52
6:M:87:ILE:O	6:M:88:PRO:C	2.51	0.52
5:L:190:ILE:HA	6:M:80:THR:HA	1.91	0.52
2:I:676:PHE:HA	2:I:679:ALA:HB3	1.92	0.52
2:P:676:PHE:HA	2:P:679:ALA:HB3	1.92	0.52
2:P:494:PRO:O	2:P:496:ILE:C	2.50	0.52
6:T:185:SER:O	6:T:203:GLN:CB	2.58	0.52
6:F:185:SER:O	6:F:203:GLN:CB	2.58	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:Q:143:GLU:C	3:Q:145:ILE:N	2.60	0.52
2:I:494:PRO:C	2:I:496:ILE:O	2.53	0.52
6:M:34:PRO:HA	6:M:127:GLY:O	2.10	0.51
7:G:45:VAL:C	7:G:48:MSE:H	2.19	0.51
2:I:499:LYS:O	2:I:503:LEU:N	2.42	0.51
3:Q:94:VAL:C	3:Q:96:TYR:N	2.67	0.51
7:U:45:VAL:C	7:U:48:MSE:H	2.19	0.51
2:B:1619:UNK:HA	2:B:1719:UNK:N	2.26	0.51
2:B:676:PHE:HA	2:B:679:ALA:HB3	1.92	0.51
4:D:74:ILE:O	4:D:78:SER:CB	2.58	0.51
7:G:53:SER:O	7:G:57:ASP:CB	2.59	0.51
3:J:127:LYS:C	3:J:130:ARG:H	2.19	0.51
4:K:74:ILE:O	4:K:78:SER:CB	2.58	0.51
2:P:507:MSE:HA	8:P:1733:SE:SE	2.61	0.51
7:N:53:SER:O	7:N:57:ASP:CB	2.59	0.51
3:C:127:LYS:C	3:C:130:ARG:H	2.19	0.51
4:R:59:ASN:O	4:R:63:MSE:N	2.42	0.51
4:R:74:ILE:O	4:R:78:SER:CB	2.58	0.51
7:U:53:SER:O	7:U:57:ASP:CB	2.59	0.51
6:M:98:THR:C	6:M:100:VAL:H	2.19	0.51
2:B:1033:UNK:CB	8:B:902:SE:SE	3.09	0.50
2:I:1619:UNK:HA	2:I:1719:UNK:N	2.26	0.50
3:Q:127:LYS:C	3:Q:130:ARG:H	2.19	0.50
5:L:276:MSE:HA	8:L:413:SE:SE	2.62	0.50
6:M:185:SER:O	6:M:203:GLN:CB	2.58	0.50
7:N:45:VAL:C	7:N:48:MSE:H	2.19	0.50
7:N:99:MSE:O	8:N:398:SE:SE	2.79	0.50
1:O:40:LEU:O	1:O:43:ALA:HB3	2.12	0.50
5:S:265:ILE:C	5:S:268:ILE:H	2.20	0.50
7:U:90:ARG:O	7:U:94:LEU:CB	2.60	0.50
3:C:92:SER:C	3:C:94:VAL:N	2.60	0.50
6:F:98:THR:C	6:F:100:VAL:H	2.19	0.50
4:K:28:THR:CB	8:K:177:SE:SE	3.10	0.50
3:Q:142:ASP:O	3:Q:143:GLU:C	2.53	0.50
2:P:1033:UNK:CB	8:P:900:SE:SE	3.10	0.50
3:Q:137:THR:O	3:Q:139:LEU:O	2.30	0.50
2:P:451:LEU:HA	2:P:454:SER:O	2.11	0.50
2:P:1619:UNK:HA	2:P:1719:UNK:N	2.26	0.50
5:S:32:LEU:H	5:S:221:LYS:H	1.60	0.50
3:C:116:LYS:O	3:C:117:LYS:C	2.54	0.50
4:D:104:ILE:HA	4:D:107:LEU:CB	2.42	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:K:48:LEU:O	8:K:175:SE:SE	2.80	0.50
1:A:40:LEU:O	1:A:43:ALA:HB3	2.12	0.50
7:G:90:ARG:O	7:G:94:LEU:CB	2.60	0.50
4:K:59:ASN:O	4:K:63:MSE:N	2.42	0.50
4:K:104:ILE:HA	4:K:107:LEU:CB	2.42	0.50
2:I:451:LEU:HA	2:I:454:SER:O	2.11	0.50
4:K:27:MSE:O	4:K:28:THR:C	2.55	0.50
3:J:143:GLU:C	3:J:145:ILE:N	2.59	0.49
3:C:66:GLN:O	3:C:70:ASN:N	2.42	0.49
3:J:116:LYS:O	3:J:117:LYS:C	2.54	0.49
7:N:90:ARG:O	7:N:94:LEU:CB	2.60	0.49
3:C:137:THR:O	3:C:139:LEU:O	2.30	0.49
5:L:265:ILE:C	5:L:268:ILE:H	2.20	0.49
2:P:214:ALA:HB2	7:U:180:HIS:C	2.38	0.49
2:I:494:PRO:O	2:I:496:ILE:O	2.30	0.49
5:S:97:LYS:CB	5:S:245:GLN:H	2.26	0.49
2:B:470:PRO:HA	2:B:471:ASN:C	2.38	0.49
4:K:14:ARG:C	4:K:17:LEU:H	2.21	0.49
2:P:1034:UNK:CB	2:P:1320:UNK:HA	2.42	0.49
5:E:265:ILE:C	5:E:268:ILE:H	2.20	0.49
1:H:40:LEU:O	1:H:43:ALA:HB3	2.12	0.49
4:R:115:PHE:HA	4:R:118:GLU:CB	2.43	0.49
5:L:32:LEU:H	5:L:221:LYS:H	1.60	0.49
4:R:104:ILE:HA	4:R:107:LEU:CB	2.42	0.49
2:B:1034:UNK:CB	2:B:1320:UNK:HA	2.42	0.49
5:E:97:LYS:CB	5:E:245:GLN:H	2.26	0.49
5:L:204:MSE:CA	8:L:415:SE:SE	3.11	0.49
4:D:26:ASN:C	4:D:28:THR:N	2.66	0.49
5:L:97:LYS:CB	5:L:245:GLN:H	2.26	0.49
2:B:451:LEU:HA	2:B:454:SER:O	2.11	0.49
4:D:27:MSE:O	4:D:28:THR:C	2.55	0.49
3:Q:116:LYS:O	3:Q:117:LYS:C	2.54	0.49
2:I:1034:UNK:CB	2:I:1320:UNK:HA	2.42	0.48
4:R:27:MSE:O	4:R:28:THR:C	2.55	0.48
6:T:98:THR:C	6:T:100:VAL:H	2.19	0.48
3:J:137:THR:O	3:J:139:LEU:O	2.30	0.48
2:I:214:ALA:HB2	7:N:180:HIS:C	2.38	0.48
2:I:470:PRO:HA	2:I:471:ASN:C	2.38	0.48
4:D:115:PHE:HA	4:D:118:GLU:CB	2.43	0.48
4:K:115:PHE:HA	4:K:118:GLU:CB	2.42	0.48
3:Q:66:GLN:O	3:Q:70:ASN:N	2.42	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:R:57:MSE:O	4:R:61:GLN:CB	2.61	0.48
5:S:100:LEU:C	5:S:102:TYR:H	2.22	0.48
4:R:14:ARG:C	4:R:17:LEU:H	2.21	0.48
1:A:62:PRO:O	1:A:66:ASN:CB	2.62	0.48
2:B:214:ALA:HB2	7:G:180:HIS:C	2.38	0.48
4:D:59:ASN:O	4:D:63:MSE:N	2.42	0.48
5:E:32:LEU:H	5:E:221:LYS:H	1.60	0.48
7:G:42:ASN:C	7:G:45:VAL:H	2.22	0.48
1:H:62:PRO:O	1:H:66:ASN:CB	2.62	0.48
1:O:62:PRO:O	1:O:66:ASN:CB	2.62	0.48
2:P:470:PRO:HA	2:P:471:ASN:C	2.38	0.48
4:D:14:ARG:C	4:D:17:LEU:H	2.21	0.48
4:D:57:MSE:O	4:D:61:GLN:CB	2.61	0.48
7:N:42:ASN:C	7:N:45:VAL:H	2.22	0.48
2:I:200:PHE:CB	8:I:904:SE:SE	3.12	0.48
4:K:57:MSE:O	4:K:61:GLN:CB	2.61	0.48
5:S:232:LEU:HA	5:S:253:LEU:HA	1.96	0.48
2:I:1723:UNK:CB	2:I:1724:UNK:CA	2.91	0.48
5:S:42:LYS:N	5:S:210:VAL:O	2.47	0.48
3:C:121:GLU:C	3:C:123:ASP:H	2.21	0.47
5:L:100:LEU:C	5:L:102:TYR:H	2.21	0.47
8:D:1301:SE:SE	7:G:175:MSE:CB	3.12	0.47
1:O:39:MSE:C	8:O:132:SE:SE	3.05	0.47
2:P:481:LEU:HA	8:P:902:SE:SE	2.64	0.47
7:U:42:ASN:C	7:U:45:VAL:H	2.22	0.47
5:E:42:LYS:N	5:E:210:VAL:O	2.47	0.47
5:L:42:LYS:N	5:L:210:VAL:O	2.47	0.47
5:L:232:LEU:HA	5:L:253:LEU:HA	1.96	0.47
3:J:142:ASP:O	3:J:143:GLU:C	2.53	0.47
3:C:200:ASP:C	5:E:295:SER:CB	2.87	0.47
5:E:100:LEU:C	5:E:102:TYR:H	2.22	0.47
2:I:494:PRO:C	2:I:496:ILE:N	2.65	0.47
3:J:200:ASP:C	5:L:295:SER:CB	2.87	0.47
3:Q:200:ASP:C	5:S:295:SER:CB	2.87	0.47
2:B:440:GLU:O	2:B:443:TYR:N	2.48	0.47
1:A:57:ASN:CB	1:H:55:ARG:HA	2.45	0.47
3:C:142:ASP:O	3:C:143:GLU:C	2.53	0.47
3:Q:120:PRO:C	3:Q:122:VAL:N	2.68	0.47
3:J:66:GLN:O	3:J:70:ASN:N	2.42	0.47
3:J:91:ASP:CA	3:J:94:VAL:HA	2.45	0.47
7:N:101:MSE:HA	7:N:104:GLU:CB	2.45	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:P:1723:UNK:CB	2:P:1724:UNK:CA	2.91	0.47
2:B:1723:UNK:CB	2:B:1724:UNK:CA	2.91	0.46
3:C:201:VAL:O	3:C:205:THR:N	2.49	0.46
5:L:32:LEU:H	5:L:221:LYS:N	2.13	0.46
2:P:440:GLU:O	2:P:443:TYR:N	2.48	0.46
2:I:440:GLU:O	2:I:443:TYR:N	2.48	0.46
3:J:167:TYR:C	3:J:169:LYS:N	2.73	0.46
3:J:201:VAL:O	3:J:205:THR:N	2.49	0.46
3:Q:167:TYR:C	3:Q:169:LYS:N	2.73	0.46
5:S:32:LEU:H	5:S:221:LYS:N	2.13	0.46
5:E:232:LEU:HA	5:E:253:LEU:HA	1.96	0.46
7:G:101:MSE:HA	7:G:104:GLU:CB	2.45	0.46
3:C:91:ASP:CA	3:C:94:VAL:HA	2.45	0.46
3:Q:91:ASP:CA	3:Q:94:VAL:HA	2.45	0.46
3:Q:121:GLU:C	3:Q:123:ASP:H	2.22	0.46
2:B:472:GLU:HA	2:B:473:LYS:CB	2.46	0.46
3:Q:201:VAL:O	3:Q:205:THR:N	2.49	0.46
3:C:116:LYS:O	3:C:117:LYS:CB	2.61	0.46
3:C:143:GLU:C	3:C:145:ILE:N	2.59	0.46
5:E:297:ASP:C	5:E:299:ARG:H	2.23	0.46
1:H:112:GLU:C	1:H:114:MSE:N	2.73	0.46
5:L:297:ASP:C	5:L:299:ARG:H	2.23	0.46
5:S:297:ASP:C	5:S:299:ARG:H	2.23	0.46
5:E:161:CYS:O	5:E:192:SER:HA	2.16	0.46
8:P:1730:SE:SE	8:P:1731:SE:SE	3.33	0.46
1:A:56:GLY:O	1:H:57:ASN:HA	2.15	0.46
2:B:1034:UNK:O	2:B:1320:UNK:CB	2.64	0.46
5:L:43:PRO:O	5:L:44:ASN:C	2.59	0.46
3:Q:116:LYS:O	3:Q:117:LYS:CB	2.61	0.46
5:S:38:CYS:HA	5:S:66:LEU:O	2.16	0.46
1:H:44:SER:C	1:H:46:VAL:N	2.74	0.46
2:I:1034:UNK:O	2:I:1320:UNK:CB	2.64	0.46
5:L:279:LYS:CB	8:L:413:SE:SE	3.14	0.46
7:U:101:MSE:HA	7:U:104:GLU:CB	2.45	0.46
6:F:93:PHE:C	6:F:95:GLY:H	2.24	0.46
6:M:93:PHE:C	6:M:95:GLY:H	2.24	0.46
5:E:38:CYS:HA	5:E:66:LEU:O	2.16	0.45
5:S:161:CYS:O	5:S:192:SER:HA	2.16	0.45
5:E:265:ILE:HA	5:E:268:ILE:CB	2.46	0.45
7:N:42:ASN:O	7:N:45:VAL:N	2.49	0.45
2:P:1034:UNK:O	2:P:1320:UNK:CB	2.64	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:S:95:ASN:C	5:S:97:LYS:H	2.25	0.45
5:S:265:ILE:HA	5:S:268:ILE:CB	2.47	0.45
5:L:161:CYS:O	5:L:192:SER:HA	2.16	0.45
2:P:528:HIS:C	5:S:227:GLY:HA2	2.42	0.45
3:C:167:TYR:C	3:C:169:LYS:N	2.73	0.45
3:J:127:LYS:HA	3:J:130:ARG:CB	2.46	0.45
5:L:265:ILE:HA	5:L:268:ILE:CB	2.46	0.45
7:U:42:ASN:O	7:U:45:VAL:N	2.49	0.45
5:E:32:LEU:H	5:E:221:LYS:N	2.13	0.45
7:G:18:GLN:O	7:G:22:LEU:CB	2.65	0.45
2:I:472:GLU:HA	2:I:473:LYS:CB	2.46	0.45
2:I:528:HIS:C	5:L:227:GLY:HA2	2.42	0.45
1:O:58:GLU:C	1:O:60:VAL:H	2.25	0.45
6:T:90:ALA:HA	6:T:93:PHE:CB	2.47	0.45
1:A:58:GLU:C	1:A:60:VAL:H	2.25	0.45
1:H:109:GLN:HA	1:H:110:ASP:HA	1.71	0.45
5:L:38:CYS:HA	5:L:66:LEU:O	2.16	0.45
1:O:84:LYS:O	1:O:88:LEU:N	2.50	0.45
1:A:112:GLU:C	1:A:114:MSE:N	2.73	0.45
6:T:63:HIS:HA	6:T:66:GLY:O	2.17	0.45
6:M:90:ALA:HA	6:M:93:PHE:CB	2.47	0.45
1:O:52:GLU:C	1:O:54:LYS:H	2.24	0.45
7:U:18:GLN:O	7:U:22:LEU:CB	2.65	0.45
7:G:42:ASN:O	7:G:45:VAL:N	2.49	0.44
5:L:95:ASN:C	5:L:97:LYS:H	2.24	0.44
2:P:472:GLU:HA	2:P:473:LYS:CB	2.46	0.44
5:S:43:PRO:O	5:S:44:ASN:C	2.59	0.44
6:T:93:PHE:C	6:T:95:GLY:H	2.24	0.44
2:B:528:HIS:C	5:E:227:GLY:HA2	2.42	0.44
6:F:90:ALA:HA	6:F:93:PHE:CB	2.47	0.44
3:C:127:LYS:HA	3:C:130:ARG:CB	2.46	0.44
5:E:95:ASN:C	5:E:97:LYS:H	2.24	0.44
3:J:116:LYS:O	3:J:117:LYS:CB	2.61	0.44
1:A:57:ASN:HA	1:H:56:GLY:O	2.16	0.44
6:F:63:HIS:HA	6:F:66:GLY:O	2.17	0.44
1:H:52:GLU:C	1:H:54:LYS:H	2.24	0.44
6:M:63:HIS:HA	6:M:66:GLY:O	2.17	0.44
1:O:109:GLN:HA	1:O:110:ASP:HA	1.71	0.44
3:J:197:ASN:O	3:J:198:VAL:C	2.60	0.44
1:O:44:SER:C	1:O:46:VAL:N	2.75	0.44
3:Q:120:PRO:O	3:Q:123:ASP:N	2.39	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:Q:127:LYS:HA	3:Q:130:ARG:CB	2.46	0.44
1:A:44:SER:C	1:A:46:VAL:N	2.74	0.44
5:E:159:GLU:O	5:E:198:SER:CB	2.66	0.44
7:N:18:GLN:O	7:N:22:LEU:CB	2.65	0.44
3:J:120:PRO:O	3:J:123:ASP:N	2.39	0.44
3:J:135:VAL:CA	3:J:136:THR:CB	2.69	0.44
1:O:58:GLU:C	1:O:60:VAL:N	2.76	0.44
6:T:94:ASN:C	6:T:96:SER:N	2.76	0.44
2:B:450:ALA:O	2:B:454:SER:N	2.46	0.44
5:E:43:PRO:O	5:E:44:ASN:C	2.59	0.44
2:I:290:LEU:O	2:I:294:SER:CB	2.66	0.44
6:T:38:ASP:HA	6:T:123:LYS:CB	2.48	0.44
7:U:51:GLN:HA	7:U:54:GLN:CB	2.48	0.44
1:A:84:LYS:O	1:A:88:LEU:N	2.50	0.44
3:J:121:GLU:C	3:J:123:ASP:H	2.22	0.44
5:L:159:GLU:O	5:L:198:SER:CB	2.66	0.44
7:N:51:GLN:HA	7:N:54:GLN:CB	2.48	0.44
1:O:112:GLU:C	1:O:114:MSE:N	2.73	0.44
7:U:106:GLU:O	7:U:109:LYS:O	2.36	0.44
7:G:106:GLU:O	7:G:109:LYS:O	2.36	0.43
1:H:58:GLU:C	1:H:60:VAL:H	2.25	0.43
7:N:106:GLU:O	7:N:109:LYS:O	2.36	0.43
5:S:260:SER:O	5:S:261:ARG:C	2.61	0.43
3:C:143:GLU:O	3:C:144:GLU:C	2.53	0.43
6:F:94:ASN:C	6:F:96:SER:N	2.76	0.43
5:S:43:PRO:HA	5:S:209:TYR:CA	2.48	0.43
5:S:159:GLU:O	5:S:198:SER:CB	2.66	0.43
1:A:109:GLN:HA	1:A:110:ASP:HA	1.71	0.43
2:B:242:SER:CB	8:B:1730:SE:SE	3.16	0.43
2:B:275:LEU:C	2:B:278:GLY:H	2.27	0.43
3:C:197:ASN:O	3:C:198:VAL:C	2.60	0.43
6:F:38:ASP:HA	6:F:123:LYS:CB	2.48	0.43
6:F:93:PHE:C	6:F:95:GLY:N	2.77	0.43
5:S:202:SER:O	5:S:206:GLY:N	2.50	0.43
1:A:52:GLU:C	1:A:54:LYS:H	2.24	0.43
7:G:104:GLU:C	7:G:107:LEU:H	2.27	0.43
2:P:290:LEU:O	2:P:294:SER:CB	2.66	0.43
6:F:190:GLY:C	6:F:192:ASP:N	2.76	0.43
7:G:51:GLN:HA	7:G:54:GLN:CB	2.48	0.43
1:A:58:GLU:C	1:A:60:VAL:N	2.76	0.43
3:C:200:ASP:O	3:C:204:PHE:N	2.47	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:M:94:ASN:C	6:M:96:SER:N	2.76	0.43
2:B:290:LEU:O	2:B:294:SER:CB	2.66	0.43
3:Q:109:LEU:HA	3:Q:112:THR:CB	2.49	0.43
1:A:36:LEU:HA	1:A:39:MSE:CB	2.49	0.43
3:C:120:PRO:C	3:C:122:VAL:N	2.68	0.43
3:C:138:ALA:N	3:C:140:LEU:CB	2.82	0.43
5:E:202:SER:O	5:E:206:GLY:N	2.50	0.43
6:M:190:GLY:C	6:M:192:ASP:N	2.76	0.43
2:P:275:LEU:C	2:P:278:GLY:H	2.27	0.43
5:S:238:ILE:HA	5:S:243:ASN:CB	2.49	0.43
5:S:246:ILE:C	5:S:248:SER:H	2.27	0.43
7:N:104:GLU:C	7:N:107:LEU:H	2.27	0.43
2:P:609:LEU:HA	2:P:612:ILE:CB	2.49	0.43
2:B:275:LEU:O	2:B:278:GLY:N	2.49	0.43
2:I:499:LYS:O	2:I:501:ALA:N	2.52	0.43
3:J:120:PRO:C	3:J:122:VAL:N	2.68	0.43
3:J:202:LEU:HA	3:J:205:THR:CB	2.49	0.43
1:O:36:LEU:HA	1:O:39:MSE:CB	2.49	0.43
5:S:209:TYR:O	8:S:310:SE:SE	2.87	0.43
3:C:202:LEU:HA	3:C:205:THR:CB	2.49	0.42
1:H:58:GLU:C	1:H:60:VAL:N	2.76	0.42
1:H:84:LYS:O	1:H:88:LEU:N	2.50	0.42
6:M:93:PHE:C	6:M:95:GLY:N	2.76	0.42
2:P:275:LEU:O	2:P:278:GLY:N	2.49	0.42
3:Q:138:ALA:N	3:Q:140:LEU:CB	2.82	0.42
3:C:120:PRO:O	3:C:123:ASP:N	2.39	0.42
3:J:109:LEU:HA	3:J:112:THR:CB	2.49	0.42
5:L:222:PHE:O	8:L:412:SE:SE	2.87	0.42
7:U:104:GLU:C	7:U:107:LEU:H	2.27	0.42
1:A:44:SER:C	1:A:46:VAL:H	2.27	0.42
2:B:609:LEU:HA	2:B:612:ILE:CB	2.49	0.42
4:D:48:LEU:C	8:D:1301:SE:SE	3.11	0.42
3:J:138:ALA:N	3:J:140:LEU:CB	2.82	0.42
5:L:101:ASN:O	5:L:102:TYR:CB	2.67	0.42
5:S:101:ASN:O	5:S:102:TYR:CB	2.67	0.42
5:E:101:ASN:O	5:E:102:TYR:CB	2.67	0.42
6:M:38:ASP:HA	6:M:123:LYS:CB	2.48	0.42
3:Q:202:LEU:HA	3:Q:205:THR:CB	2.49	0.42
5:L:238:ILE:HA	5:L:243:ASN:CB	2.49	0.42
5:E:41:TRP:HA	5:E:212:GLU:H	1.85	0.42
5:E:246:ILE:C	5:E:248:SER:H	2.27	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:E:297:ASP:C	5:E:299:ARG:N	2.78	0.42
5:L:182:MSE:CB	8:L:410:SE:SE	3.17	0.42
4:R:14:ARG:O	4:R:18:SER:N	2.51	0.42
5:E:238:ILE:HA	5:E:243:ASN:CB	2.49	0.42
1:H:36:LEU:HA	1:H:39:MSE:CB	2.49	0.42
3:J:117:LYS:O	3:J:118:ASN:CB	2.68	0.42
3:J:121:GLU:O	3:J:123:ASP:C	2.60	0.42
5:L:41:TRP:HA	5:L:212:GLU:H	1.85	0.42
3:Q:57:ALA:CB	3:Q:58:GLU:HA	2.28	0.42
6:T:190:GLY:C	6:T:192:ASP:N	2.76	0.42
3:C:121:GLU:O	3:C:123:ASP:C	2.60	0.42
2:I:609:LEU:HA	2:I:612:ILE:CB	2.49	0.42
7:N:90:ARG:O	7:N:94:LEU:N	2.53	0.42
1:H:44:SER:C	1:H:46:VAL:H	2.27	0.42
2:P:604:ARG:O	2:P:608:LYS:N	2.52	0.42
3:C:109:LEU:HA	3:C:112:THR:CB	2.49	0.42
3:C:135:VAL:CA	3:C:136:THR:CB	2.69	0.42
4:D:14:ARG:O	4:D:18:SER:N	2.51	0.42
5:L:260:SER:O	5:L:261:ARG:C	2.61	0.42
5:L:297:ASP:C	5:L:299:ARG:N	2.78	0.42
3:Q:197:ASN:O	3:Q:198:VAL:C	2.60	0.42
3:C:57:ALA:CB	3:C:58:GLU:HA	2.28	0.41
4:K:14:ARG:O	4:K:18:SER:N	2.51	0.41
1:O:113:LYS:O	1:O:116:GLU:N	2.52	0.41
3:Q:121:GLU:O	3:Q:123:ASP:C	2.60	0.41
7:N:184:THR:C	8:N:396:SE:SE	3.11	0.41
2:P:498:ASP:C	2:P:502:ASN:H	2.21	0.41
6:T:93:PHE:C	6:T:95:GLY:N	2.76	0.41
2:B:604:ARG:O	2:B:608:LYS:N	2.52	0.41
1:H:113:LYS:O	1:H:116:GLU:N	2.52	0.41
5:L:246:ILE:C	5:L:248:SER:H	2.27	0.41
5:S:44:ASN:H	5:S:209:TYR:HA	1.84	0.41
5:E:260:SER:O	5:E:261:ARG:C	2.62	0.41
2:I:604:ARG:O	2:I:608:LYS:N	2.52	0.41
5:L:202:SER:O	5:L:206:GLY:N	2.50	0.41
5:S:4:GLN:CB	5:S:258:ASN:HA	2.51	0.41
5:S:41:TRP:HA	5:S:212:GLU:H	1.85	0.41
5:L:4:GLN:CB	5:L:258:ASN:HA	2.51	0.41
2:P:507:MSE:CA	8:P:1733:SE:SE	3.18	0.41
2:P:517:LYS:C	2:P:519:THR:N	2.79	0.41
4:D:25:ILE:C	4:D:27:MSE:N	2.77	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:I:275:LEU:C	2:I:278:GLY:H	2.27	0.41
3:Q:117:LYS:O	3:Q:118:ASN:CB	2.68	0.41
7:U:90:ARG:O	7:U:94:LEU:N	2.53	0.41
5:E:4:GLN:CB	5:E:258:ASN:HA	2.51	0.41
4:K:17:LEU:O	4:K:21:LEU:N	2.54	0.41
5:E:43:PRO:HA	5:E:209:TYR:CB	2.51	0.41
1:O:44:SER:C	1:O:46:VAL:H	2.27	0.41
1:O:68:VAL:CB	2:P:295:PHE:CB	2.99	0.41
5:S:43:PRO:HA	5:S:209:TYR:CB	2.51	0.41
5:S:283:GLN:HA	5:S:287:GLU:HA	2.03	0.41
1:A:68:VAL:CB	2:B:295:PHE:CB	2.99	0.41
6:F:45:SER:CB	6:F:115:ILE:HA	2.51	0.41
2:I:275:LEU:O	2:I:278:GLY:N	2.49	0.41
4:K:29:THR:H	8:K:177:SE:SE	2.53	0.41
6:T:21:LYS:O	6:T:25:SER:N	2.54	0.41
3:C:94:VAL:CB	7:G:166:ILE:N	2.84	0.41
7:G:90:ARG:O	7:G:94:LEU:N	2.53	0.41
1:H:89:LEU:CB	8:H:313:SE:SE	3.19	0.41
5:L:43:PRO:HA	5:L:209:TYR:CB	2.51	0.41
6:M:21:LYS:O	6:M:25:SER:N	2.54	0.41
2:P:523:ARG:C	2:P:525:SER:N	2.79	0.41
4:R:24:LEU:HA	4:R:27:MSE:CB	2.51	0.41
6:F:175:ALA:C	6:F:177:ILE:N	2.79	0.40
2:I:499:LYS:O	2:I:500:ARG:C	2.60	0.40
3:J:94:VAL:CB	7:N:166:ILE:N	2.84	0.40
4:K:25:ILE:C	4:K:27:MSE:N	2.77	0.40
5:L:283:GLN:HA	5:L:287:GLU:HA	2.03	0.40
6:M:45:SER:CB	6:M:115:ILE:HA	2.51	0.40
6:M:175:ALA:C	6:M:177:ILE:N	2.79	0.40
6:T:45:SER:CB	6:T:115:ILE:HA	2.51	0.40
1:A:113:LYS:O	1:A:116:GLU:N	2.52	0.40
4:K:24:LEU:HA	4:K:27:MSE:CB	2.51	0.40
3:Q:202:LEU:O	3:Q:206:PHE:N	2.55	0.40
4:D:24:LEU:HA	4:D:27:MSE:CB	2.51	0.40
5:L:232:LEU:HA	5:L:252:LEU:O	2.22	0.40
3:Q:94:VAL:CB	7:U:166:ILE:N	2.84	0.40
5:S:265:ILE:O	5:S:269:ASN:N	2.55	0.40
5:S:297:ASP:C	5:S:299:ARG:N	2.78	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	90/131 (69%)	83 (92%)	7 (8%)	0	100	100
1	H	90/131 (69%)	83 (92%)	7 (8%)	0	100	100
1	O	90/131 (69%)	83 (92%)	7 (8%)	0	100	100
2	B	261/583 (45%)	239 (92%)	16 (6%)	6 (2%)	5	28
2	I	261/583 (45%)	239 (92%)	17 (6%)	5 (2%)	6	32
2	P	261/583 (45%)	241 (92%)	16 (6%)	4 (2%)	8	38
3	C	155/223 (70%)	135 (87%)	14 (9%)	6 (4%)	2	19
3	J	155/223 (70%)	135 (87%)	14 (9%)	6 (4%)	2	19
3	Q	155/223 (70%)	135 (87%)	14 (9%)	6 (4%)	2	19
4	D	82/121 (68%)	69 (84%)	13 (16%)	0	100	100
4	K	82/121 (68%)	69 (84%)	13 (16%)	0	100	100
4	R	82/121 (68%)	69 (84%)	13 (16%)	0	100	100
5	E	223/275 (81%)	193 (86%)	30 (14%)	0	100	100
5	L	223/275 (81%)	193 (86%)	30 (14%)	0	100	100
5	S	223/275 (81%)	193 (86%)	30 (14%)	0	100	100
6	F	201/210 (96%)	185 (92%)	15 (8%)	1 (0%)	24	62
6	M	201/210 (96%)	184 (92%)	16 (8%)	1 (0%)	24	62
6	T	201/210 (96%)	185 (92%)	15 (8%)	1 (0%)	24	62
7	G	97/295 (33%)	82 (84%)	14 (14%)	1 (1%)	12	47
7	N	97/295 (33%)	82 (84%)	14 (14%)	1 (1%)	12	47
7	U	97/295 (33%)	82 (84%)	14 (14%)	1 (1%)	12	47
All	All	3327/5514 (60%)	2959 (89%)	329 (10%)	39 (1%)	10	42

All (39) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	230	SER
2	B	497	ASN
2	B	500	ARG
3	C	94	VAL
3	C	144	GLU
7	G	88	PRO
2	I	230	SER
3	J	94	VAL
3	J	144	GLU
7	N	88	PRO
2	P	230	SER
3	Q	94	VAL
3	Q	144	GLU
7	U	88	PRO
3	C	95	VAL
3	C	107	GLU
3	C	122	VAL
2	I	497	ASN
3	J	95	VAL
3	J	107	GLU
3	J	122	VAL
3	Q	95	VAL
3	Q	107	GLU
3	Q	122	VAL
2	B	459	ILE
2	I	459	ILE
2	P	459	ILE
2	B	514	VAL
2	I	514	VAL
2	P	514	VAL
2	B	229	LEU
3	C	96	TYR
6	F	37	ILE
2	I	229	LEU
3	J	96	TYR
6	M	37	ILE
2	P	229	LEU
3	Q	96	TYR
6	T	37	ILE

5.3.2 Protein sidechains ⓘ

There are no protein residues with a non-rotameric sidechain to report in this entry.

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 98 ligands modelled in this entry, 98 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
2	B	6
2	I	6
2	P	6

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	B	1727:UNK	C	669:GLU	N	41.03
1	I	1727:UNK	C	669:GLU	N	41.03
1	P	1727:UNK	C	669:GLU	N	41.03
1	B	1101:UNK	C	1028:UNK	N	37.71
1	I	1101:UNK	C	1028:UNK	N	37.71
1	P	1101:UNK	C	1028:UNK	N	37.71
1	B	1627:UNK	C	1719:UNK	N	26.03
1	I	1627:UNK	C	1719:UNK	N	26.03
1	P	1627:UNK	C	1719:UNK	N	26.03
1	B	1323:UNK	C	1619:UNK	N	14.40
1	I	1323:UNK	C	1619:UNK	N	14.40
1	P	1323:UNK	C	1619:UNK	N	14.40
1	B	616:GLU	C	1094:UNK	N	12.88
1	I	616:GLU	C	1094:UNK	N	12.88
1	P	616:GLU	C	1094:UNK	N	12.88
1	B	1038:UNK	C	1316:UNK	N	4.44
1	I	1038:UNK	C	1316:UNK	N	4.44
1	P	1038:UNK	C	1316:UNK	N	4.44

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	92/131 (70%)	-0.75	0 100 100	42, 74, 118, 119	0
1	H	92/131 (70%)	-0.96	0 100 100	28, 42, 119, 121	0
1	O	92/131 (70%)	-0.91	0 100 100	80, 92, 168, 169	0
2	B	265/583 (45%)	-0.64	3 (1%) 78 62	32, 87, 179, 181	0
2	I	265/583 (45%)	-0.67	1 (0%) 88 77	52, 71, 177, 191	0
2	P	265/583 (45%)	-0.61	2 (0%) 82 68	15, 124, 209, 210	0
3	C	156/223 (69%)	-0.70	1 (0%) 85 73	33, 106, 192, 192	0
3	J	156/223 (69%)	-0.61	0 100 100	42, 112, 192, 240	0
3	Q	156/223 (69%)	-0.57	0 100 100	61, 135, 205, 238	0
4	D	84/121 (69%)	-0.73	0 100 100	54, 66, 92, 99	0
4	K	84/121 (69%)	-0.84	0 100 100	54, 59, 113, 114	0
4	R	84/121 (69%)	-0.77	0 100 100	52, 110, 165, 167	0
5	E	223/275 (81%)	-0.35	2 (0%) 81 66	99, 184, 216, 223	0
5	L	223/275 (81%)	-0.59	0 100 100	64, 149, 196, 198	0
5	S	223/275 (81%)	-0.36	4 (1%) 67 53	103, 149, 255, 259	0
6	F	202/210 (96%)	0.00	7 (3%) 47 37	187, 252, 349, 353	0
6	M	202/210 (96%)	-0.48	4 (1%) 65 50	96, 134, 194, 199	0
6	T	202/210 (96%)	-0.25	2 (0%) 79 64	40, 182, 243, 247	0
7	G	98/295 (33%)	-0.34	2 (2%) 65 50	33, 147, 171, 172	0
7	N	98/295 (33%)	-0.85	0 100 100	35, 47, 147, 148	0
7	U	98/295 (33%)	-0.63	0 100 100	48, 119, 138, 140	0
All	All	3360/5514 (60%)	-0.55	28 (0%) 82 68	15, 129, 244, 353	0

All (28) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	239	SER	6.8
7	G	58	PRO	4.3
6	F	28	ILE	4.1
5	S	241	ALA	3.8
2	P	553	ASN	3.7
6	T	48	ASN	3.5
5	E	48	ASP	3.4
2	B	242	SER	3.4
6	F	118	GLN	3.0
6	M	124	GLY	2.9
5	S	226	HIS	2.8
5	S	300	VAL	2.7
6	M	45	SER	2.7
2	P	483	ASP	2.7
6	F	124	GLY	2.7
6	F	125	ASP	2.5
5	E	107	ASN	2.5
6	T	194	SER	2.5
6	M	210	LEU	2.5
6	M	175	ALA	2.4
2	B	273	ASP	2.4
6	F	148	PHE	2.4
6	F	25	SER	2.3
3	C	110	VAL	2.3
6	F	141	ASN	2.2
5	S	61	PRO	2.2
7	G	33	GLU	2.1
2	I	475	GLU	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum,

median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
8	SE	A	132	1/1	-	-	98,98,98,98	1
8	SE	B	1729	1/1	0.87	0.27	98,98,98,98	1
8	SE	C	225	1/1	0.90	0.37	98,98,98,98	1
8	SE	I	904	1/1	0.92	0.52	98,98,98,98	1
8	SE	Q	324	1/1	0.92	0.18	98,98,98,98	1
8	SE	Q	327	1/1	0.92	0.12	98,98,98,98	1
8	SE	B	1731	1/1	-	-	98,98,98,98	1
8	SE	P	1731	1/1	0.95	0.59	98,98,98,98	1
8	SE	P	900	1/1	0.95	0.33	98,98,98,98	1
8	SE	N	400	1/1	0.96	0.24	98,98,98,98	0
8	SE	P	1730	1/1	0.96	0.31	98,98,98,98	0
8	SE	A	313	1/1	0.96	0.12	98,98,98,98	1
8	SE	P	1733	1/1	0.96	0.38	98,98,98,98	1
8	SE	G	1297	1/1	0.96	0.14	98,98,98,98	1
8	SE	P	902	1/1	0.96	0.22	98,98,98,98	1
8	SE	C	227	1/1	-	-	98,98,98,98	1
8	SE	B	902	1/1	0.96	0.13	98,98,98,98	1
8	SE	L	414	1/1	0.96	0.20	98,98,98,98	1
8	SE	U	496	1/1	0.96	0.18	98,98,98,98	1
8	SE	E	408	1/1	-	-	98,98,98,98	1
8	SE	U	500	1/1	0.96	0.46	98,98,98,98	1
8	SE	I	1728	1/1	0.97	0.30	98,98,98,98	1
8	SE	I	901	1/1	0.97	0.20	98,98,98,98	1
8	SE	B	1732	1/1	0.97	0.16	98,98,98,98	1
8	SE	E	413	1/1	-	-	98,98,98,98	1
8	SE	K	173	1/1	0.97	0.28	98,98,98,98	1
8	SE	P	901	1/1	0.97	0.09	98,98,98,98	1
8	SE	G	1298	1/1	-	-	98,98,98,98	1
8	SE	K	175	1/1	0.97	0.27	98,98,98,98	1
8	SE	L	412	1/1	0.97	0.18	98,98,98,98	1
8	SE	H	132	1/1	-	-	98,98,98,98	1
8	SE	H	133	1/1	-	-	98,98,98,98	1
8	SE	G	1296	1/1	0.97	0.18	98,98,98,98	1
8	SE	S	308	1/1	0.97	0.13	98,98,98,98	1
8	SE	N	397	1/1	0.97	0.35	98,98,98,98	1
8	SE	I	1730	1/1	-	-	98,98,98,98	1
8	SE	I	1731	1/1	-	-	98,98,98,98	1
8	SE	A	133	1/1	0.97	0.17	98,98,98,98	1
8	SE	I	902	1/1	-	-	98,98,98,98	1
8	SE	O	133	1/1	0.98	0.23	98,98,98,98	1
8	SE	P	1729	1/1	0.98	0.42	98,98,98,98	1

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
8	SE	B	901	1/1	0.98	0.09	98,98,98,98	0
8	SE	C	226	1/1	0.98	0.14	98,98,98,98	1
8	SE	J	224	1/1	-	-	98,98,98,98	1
8	SE	D	1301	1/1	0.98	0.07	98,98,98,98	0
8	SE	J	226	1/1	-	-	98,98,98,98	1
8	SE	J	227	1/1	-	-	98,98,98,98	1
8	SE	K	177	1/1	0.98	0.15	98,98,98,98	0
8	SE	B	1730	1/1	0.98	0.29	98,98,98,98	1
8	SE	B	903	1/1	0.98	0.20	98,98,98,98	1
8	SE	L	413	1/1	0.98	0.08	98,98,98,98	0
8	SE	Q	325	1/1	0.98	0.08	98,98,98,98	1
8	SE	Q	326	1/1	0.98	0.19	98,98,98,98	1
8	SE	G	1300	1/1	0.98	0.34	98,98,98,98	0
8	SE	R	223	1/1	0.98	0.15	98,98,98,98	1
8	SE	N	396	1/1	0.98	0.11	98,98,98,98	1
8	SE	S	313	1/1	0.98	0.04	98,98,98,98	0
8	SE	H	313	1/1	0.98	0.16	98,98,98,98	1
8	SE	M	311	1/1	-	-	98,98,98,98	1
8	SE	M	312	1/1	-	-	98,98,98,98	1
8	SE	M	313	1/1	-	-	98,98,98,98	1
8	SE	B	904	1/1	0.98	0.11	98,98,98,98	1
8	SE	I	1729	1/1	0.99	0.18	98,98,98,98	1
8	SE	N	398	1/1	0.99	0.06	98,98,98,98	0
8	SE	N	399	1/1	-	-	98,98,98,98	1
8	SE	E	410	1/1	0.99	0.21	98,98,98,98	1
8	SE	O	132	1/1	0.99	0.10	98,98,98,98	1
8	SE	I	903	1/1	0.99	0.14	98,98,98,98	1
8	SE	O	313	1/1	0.99	0.05	98,98,98,98	0
8	SE	P	1728	1/1	0.99	0.10	98,98,98,98	1
8	SE	E	411	1/1	0.99	0.04	98,98,98,98	0
8	SE	I	906	1/1	0.99	0.05	98,98,98,98	1
8	SE	I	905	1/1	0.99	0.13	98,98,98,98	0
8	SE	P	1732	1/1	-	-	98,98,98,98	1
8	SE	J	225	1/1	0.99	0.23	98,98,98,98	1
8	SE	E	412	1/1	0.99	0.07	98,98,98,98	1
8	SE	K	174	1/1	0.99	0.07	98,98,98,98	1
8	SE	C	224	1/1	0.99	0.08	98,98,98,98	1
8	SE	D	222	1/1	0.99	0.12	98,98,98,98	0
8	SE	L	408	1/1	0.99	0.04	98,98,98,98	1
8	SE	L	409	1/1	0.99	0.06	98,98,98,98	1
8	SE	L	410	1/1	0.99	0.07	98,98,98,98	0
8	SE	R	222	1/1	0.99	0.26	98,98,98,98	1

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
8	SE	G	1299	1/1	0.99	0.02	98,98,98,98	1
8	SE	D	223	1/1	0.99	0.18	98,98,98,98	1
8	SE	S	309	1/1	0.99	0.19	98,98,98,98	1
8	SE	S	310	1/1	0.99	0.05	98,98,98,98	0
8	SE	S	311	1/1	0.99	0.40	98,98,98,98	1
8	SE	S	312	1/1	0.99	0.19	98,98,98,98	1
8	SE	B	1728	1/1	0.99	0.03	98,98,98,98	0
8	SE	T	311	1/1	0.99	0.14	98,98,98,98	1
8	SE	T	312	1/1	-	-	98,98,98,98	1
8	SE	U	495	1/1	0.99	0.22	98,98,98,98	0
8	SE	L	415	1/1	0.99	0.11	98,98,98,98	1
8	SE	U	497	1/1	-	-	98,98,98,98	1
8	SE	U	498	1/1	-	-	98,98,98,98	1
8	SE	U	499	1/1	0.99	0.13	98,98,98,98	1
8	SE	E	409	1/1	0.99	0.04	98,98,98,98	1

6.5 Other polymers [i](#)

There are no such residues in this entry.