



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 15, 2017 – 07:29 am GMT

PDB ID : 3UN4
Title : Yeast 20S proteasome in complex with PR-957 (morpholine)
Authors : Huber, E.; Basler, M.; Schwab, R.; Heinemeyer, W.; Kirk, C.; Groettrup, M.; Groll, M.
Deposited on : 2011-11-15
Resolution : 3.40 Å(reported)

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

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

A user guide is available at

<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) : 1.9-1692
EDS : trunk28620
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : recalc28949

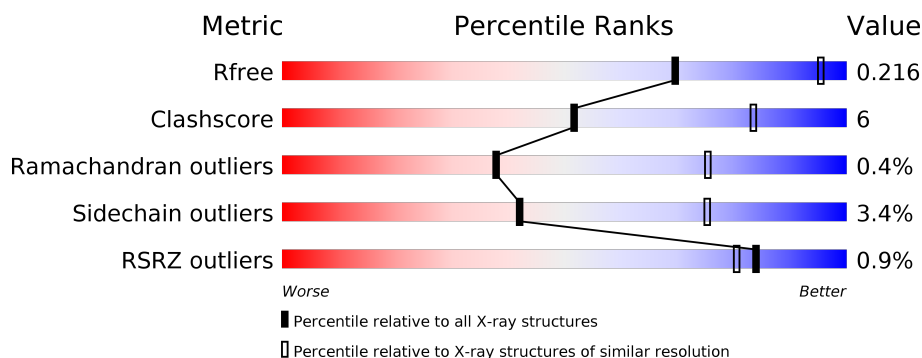
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.40 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	100719	1679 (3.50-3.30)
Clashscore	112137	1832 (3.50-3.30)
Ramachandran outliers	110173	1789 (3.50-3.30)
Sidechain outliers	110143	1789 (3.50-3.30)
RSRZ outliers	101464	1709 (3.50-3.30)

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.

Mol	Chain	Length	Quality of chain
1	A	250	<div> <div>88%</div> <div>12%</div> </div>
1	O	250	<div> <div>%</div> <div>87%</div> <div>13%</div> </div>
2	B	258	<div> <div>75%</div> <div>18%</div> <div>• 5%</div> </div>
2	P	258	<div> <div>%</div> <div>75%</div> <div>18%</div> <div>• 5%</div> </div>
3	C	254	<div> <div>%</div> <div>72%</div> <div>20%</div> <div>• 5%</div> </div>
3	Q	254	<div> <div>2%</div> <div>74%</div> <div>19%</div> <div>• 5%</div> </div>

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Mol	Chain	Length	Quality of chain
4	D	260	
4	R	260	
5	E	234	
5	S	234	
6	F	288	
6	T	288	
7	G	252	
7	U	252	
8	H	232	
8	V	232	
9	I	205	
9	W	205	
10	J	198	
10	X	198	
11	K	212	
11	Y	212	
12	L	222	
12	Z	222	
13	M	233	
13	a	233	
14	N	196	
14	b	196	

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
15	04C	N	201	-	-	-	X
15	04C	b	201	-	-	-	X

2 Entry composition

There are 16 unique types of molecules in this entry. The entry contains 51112 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 Proteasome component Y7.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	250	Total	C	N	O	S	0	0	0
			1915	1219	315	377	4			
1	O	250	Total	C	N	O	S	0	0	0
			1915	1219	315	377	4			

- Molecule 2 is a protein called Proteasome component Y13.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	244	Total	C	N	O	S	0	0	0
			1904	1201	321	379	3			
2	P	244	Total	C	N	O	S	0	0	0
			1904	1201	321	379	3			

- Molecule 3 is a protein called Proteasome component PRE6.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	241	Total	C	N	O	S	0	0	0
			1890	1181	331	374	4			
3	Q	241	Total	C	N	O	S	0	0	0
			1890	1181	331	374	4			

- Molecule 4 is a protein called Proteasome component PUP2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	242	Total	C	N	O	S	0	0	0
			1861	1162	314	378	7			
4	R	242	Total	C	N	O	S	0	0	0
			1861	1162	314	378	7			

- Molecule 5 is a protein called Proteasome component PRE5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	E	233	Total	C	N	O	S	0	0	0
			1795	1129	312	350	4			
5	S	233	Total	C	N	O	S	0	0	0
			1795	1129	312	350	4			

- Molecule 6 is a protein called Proteasome component C1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
6	F	244	Total	C	N	O	S	0	0	0
			1896	1205	330	357	4			
6	T	244	Total	C	N	O	S	0	0	0
			1896	1205	330	357	4			

- Molecule 7 is a protein called Proteasome component C7-alpha.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
7	G	243	Total	C	N	O	S	0	0	0
			1921	1221	322	370	8			
7	U	243	Total	C	N	O	S	0	0	0
			1921	1221	322	370	8			

- Molecule 8 is a protein called Proteasome component PUP1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
8	H	222	Total	C	N	O	S	0	0	0
			1684	1061	293	323	7			
8	V	222	Total	C	N	O	S	0	0	0
			1684	1061	293	323	7			

- Molecule 9 is a protein called Proteasome component PUP3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
9	I	204	Total	C	N	O	S	0	0	0
			1581	1010	258	305	8			
9	W	204	Total	C	N	O	S	0	0	0
			1581	1010	258	305	8			

- Molecule 10 is a protein called Proteasome component C11.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
10	J	198	Total	C	N	O	S	0	0	0
			1585	1005	269	305	6			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
10	X	198	Total	C	N	O	S	0	0	0
			1585	1005	269	305	6			

- Molecule 11 is a protein called Proteasome component PRE2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
11	K	212	Total	C	N	O	S	0	0	0
			1644	1045	280	312	7			
11	Y	212	Total	C	N	O	S	0	0	0
			1644	1045	280	312	7			

- Molecule 12 is a protein called Proteasome component C5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
12	L	222	Total	C	N	O	S	0	0	0
			1757	1115	303	335	4			
12	Z	222	Total	C	N	O	S	0	0	0
			1757	1115	303	335	4			

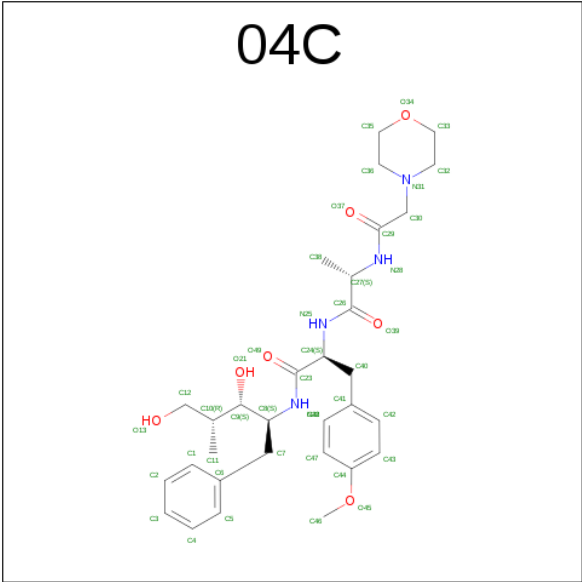
- Molecule 13 is a protein called Proteasome component PRE4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
13	M	233	Total	C	N	O	S	0	0	0
			1824	1154	312	351	7			
13	a	233	Total	C	N	O	S	0	0	0
			1824	1154	312	351	7			

- Molecule 14 is a protein called Proteasome component PRE3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
14	N	196	Total	C	N	O	S	0	0	0
			1512	955	250	300	7			
14	b	196	Total	C	N	O	S	0	0	0
			1512	955	250	300	7			

- Molecule 15 is 1,2,4-TRIDEOXY-4-METHYL-2-{[N-(MORPHOLIN-4-YLACETYL)-L-ALANYL-O-METHYL-L-TYROSYL]AMINO}-1-PHENYL-D-XYLITOL (three-letter code: 04C) (formula: C₃₁H₄₄N₄O₇).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
15	H	1	Total	C	N	O	0	0
			42	31	4	7		
15	K	1	Total	C	N	O	0	0
			42	31	4	7		
15	N	1	Total	C	N	O	0	0
			42	31	4	7		
15	V	1	Total	C	N	O	0	0
			42	31	4	7		
15	Y	1	Total	C	N	O	0	0
			42	31	4	7		
15	b	1	Total	C	N	O	0	0
			42	31	4	7		

- Molecule 16 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
16	A	56	Total	O	0	0
			56	56		
16	B	38	Total	O	0	0
			38	38		
16	C	42	Total	O	0	0
			42	42		
16	D	38	Total	O	0	0
			38	38		
16	E	23	Total	O	0	0
			23	23		
16	F	48	Total	O	0	0
			48	48		

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
16	G	61	Total O 61 61	0	0
16	H	49	Total O 49 49	0	0
16	I	62	Total O 62 62	0	0
16	J	54	Total O 54 54	0	0
16	K	46	Total O 46 46	0	0
16	L	55	Total O 55 55	0	0
16	M	72	Total O 72 72	0	0
16	N	51	Total O 51 51	0	0
16	O	32	Total O 32 32	0	0
16	P	30	Total O 30 30	0	0
16	Q	30	Total O 30 30	0	0
16	R	27	Total O 27 27	0	0
16	S	20	Total O 20 20	0	0
16	T	41	Total O 41 41	0	0
16	U	61	Total O 61 61	0	0
16	V	45	Total O 45 45	0	0
16	W	57	Total O 57 57	0	0
16	X	52	Total O 52 52	0	0
16	Y	49	Total O 49 49	0	0
16	Z	50	Total O 50 50	0	0
16	a	75	Total O 75 75	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
16	b	58	Total	O	0	0
			58	58		

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.

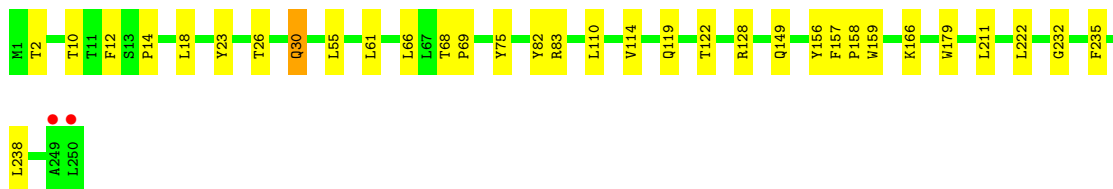
• Molecule 1: Proteasome component Y7

Chain A: 



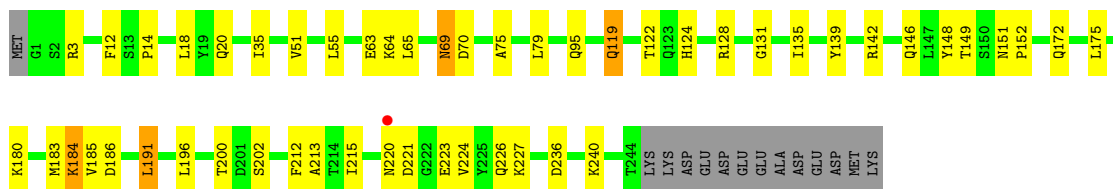
• Molecule 1: Proteasome component Y7

Chain O: 




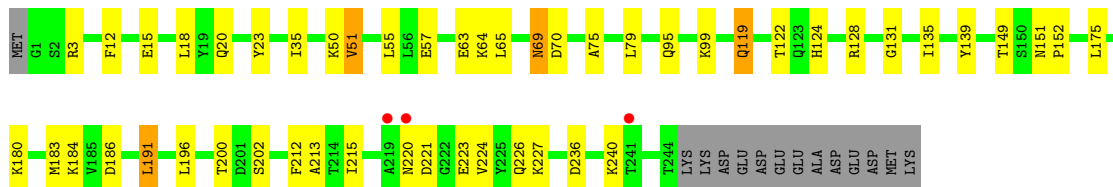
• Molecule 2: Proteasome component Y13

Chain B: 

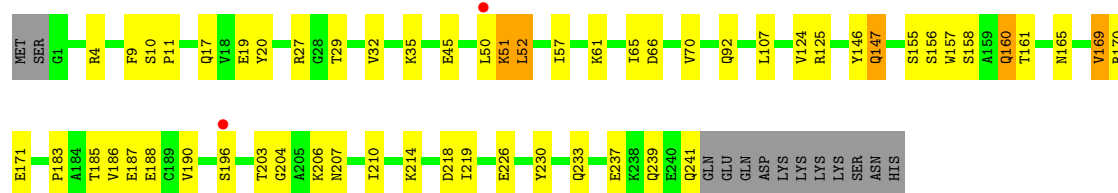
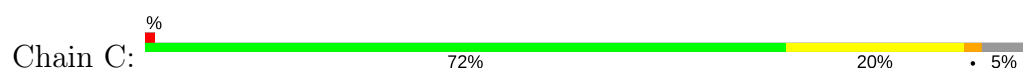


• Molecule 2: Proteasome component Y13

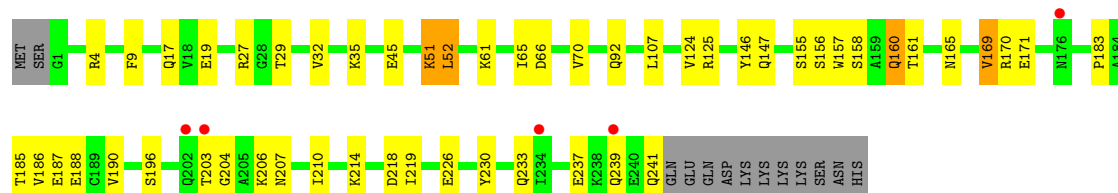
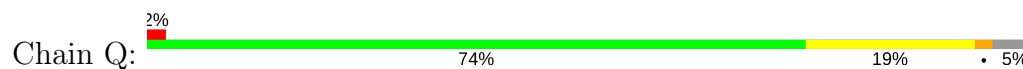
Chain P: 



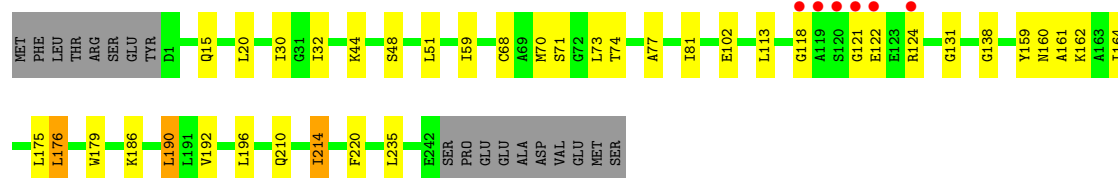
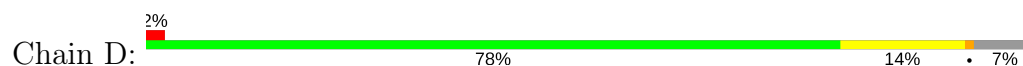
• Molecule 3: Proteasome component PRE6



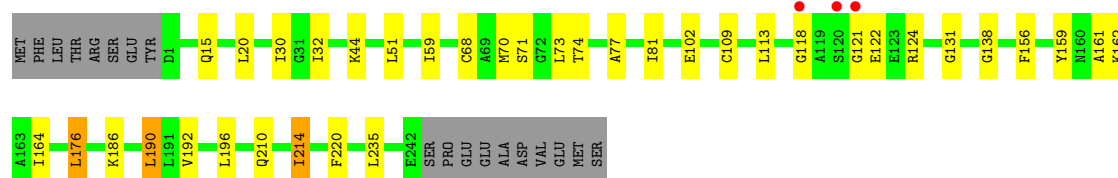
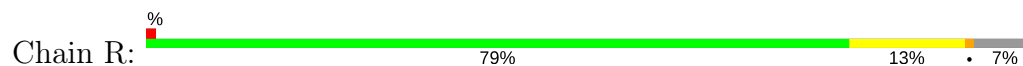
• Molecule 3: Proteasome component PRE6



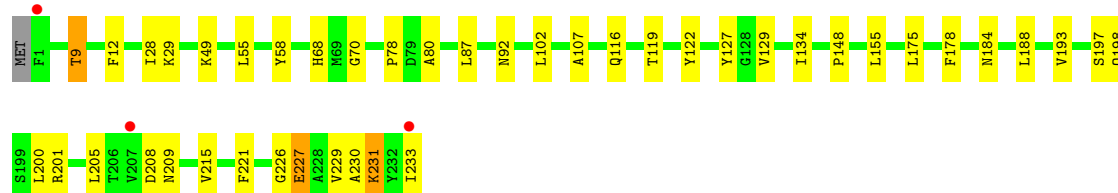
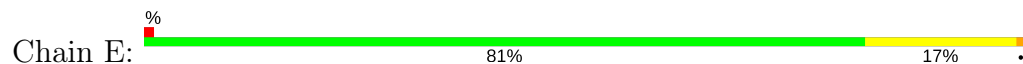
• Molecule 4: Proteasome component PUP2



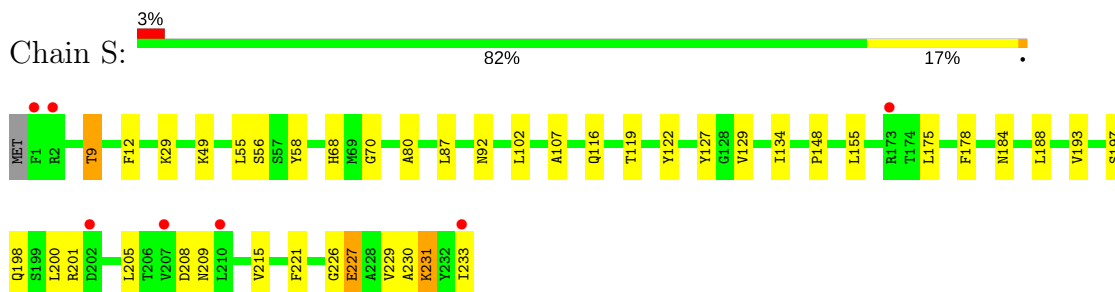
• Molecule 4: Proteasome component PUP2



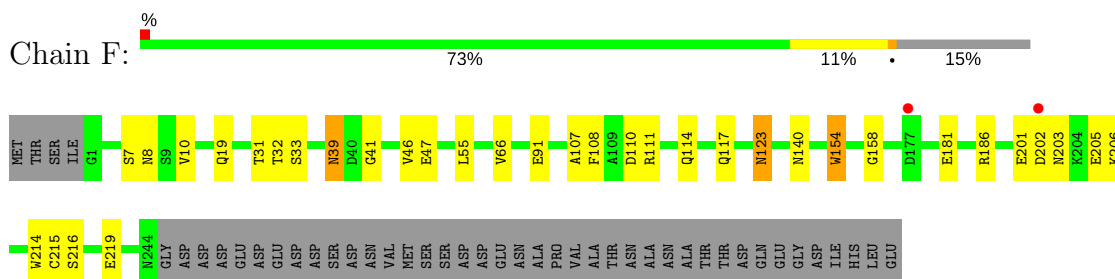
• Molecule 5: Proteasome component PRE5



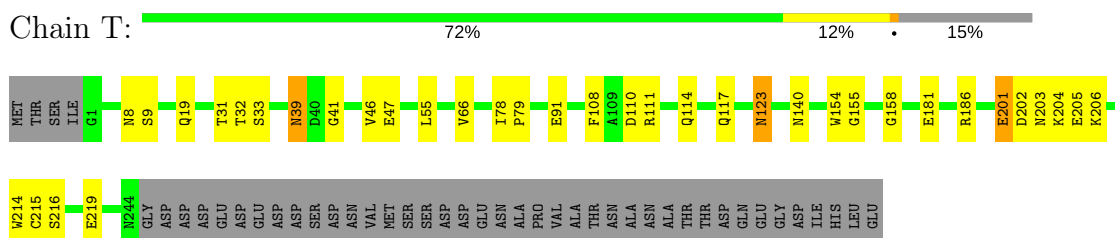
- Molecule 5: Proteasome component PRE5



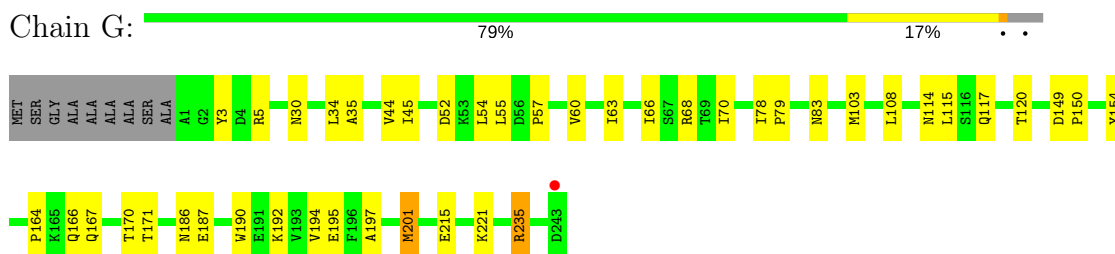
- Molecule 6: Proteasome component C1



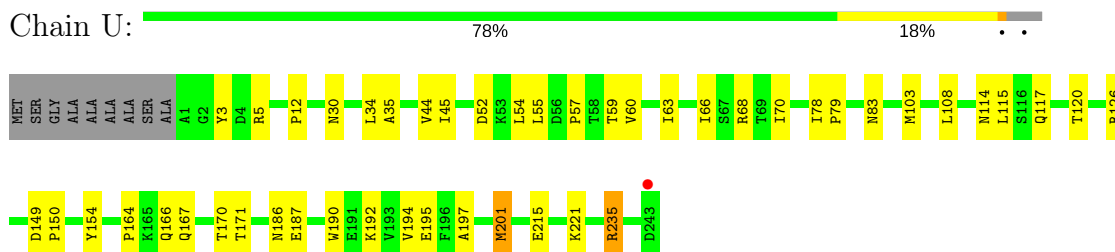
- Molecule 6: Proteasome component C1



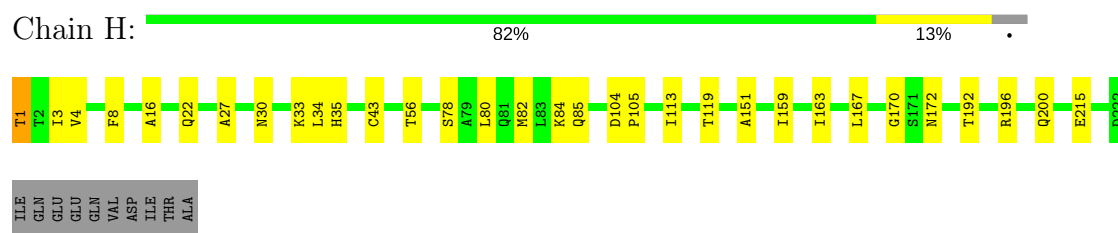
- Molecule 7: Proteasome component C7-alpha



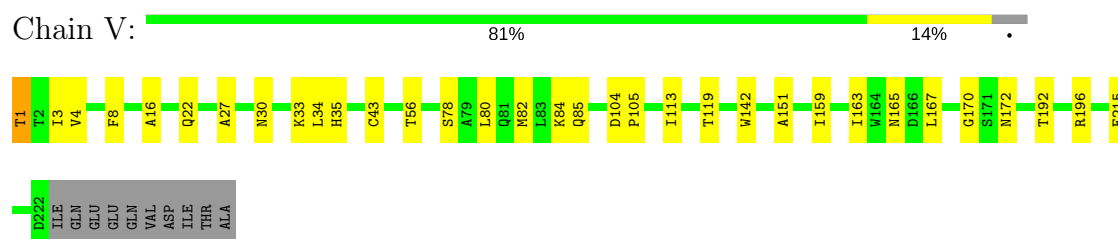
- Molecule 7: Proteasome component C7-alpha



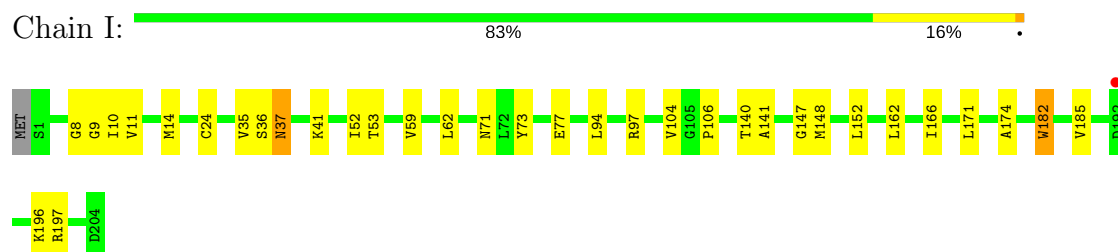
- Molecule 8: Proteasome component PUP1



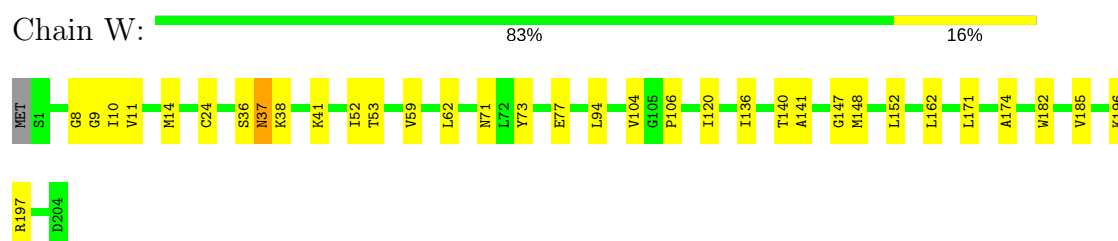
- Molecule 8: Proteasome component PUP1



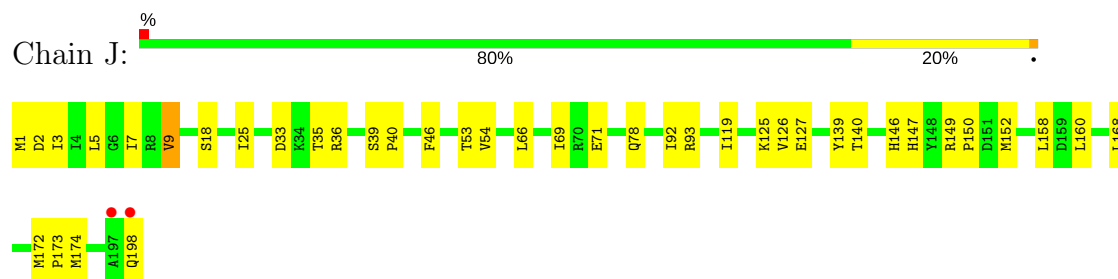
- Molecule 9: Proteasome component PUP3



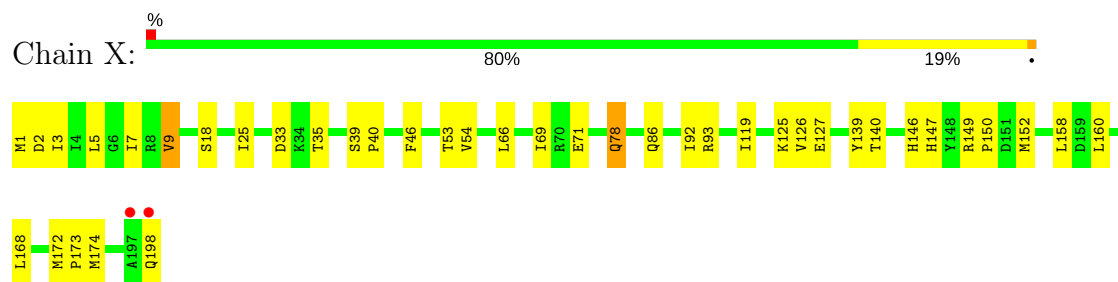
- Molecule 9: Proteasome component PUP3



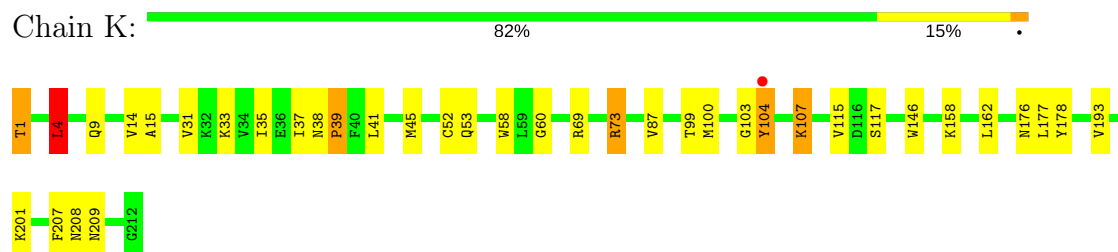
- Molecule 10: Proteasome component C11



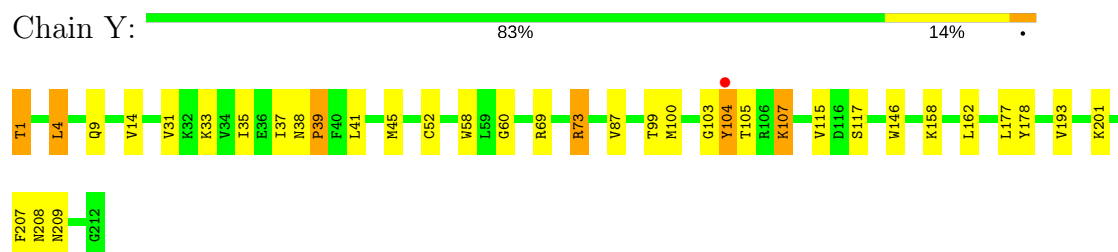
- Molecule 10: Proteasome component C11



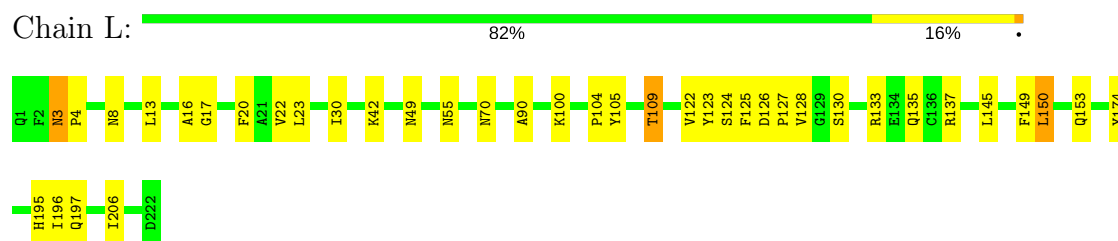
- Molecule 11: Proteasome component PRE2



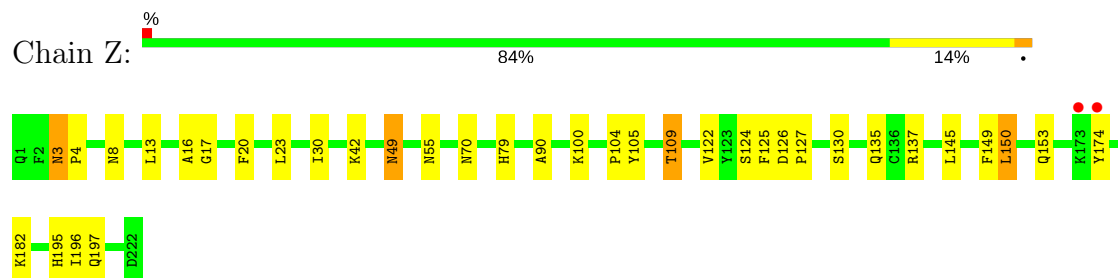
- Molecule 11: Proteasome component PRE2



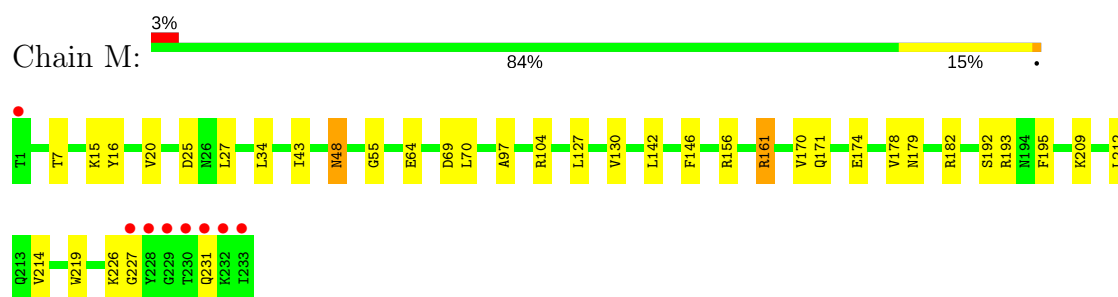
- Molecule 12: Proteasome component C5



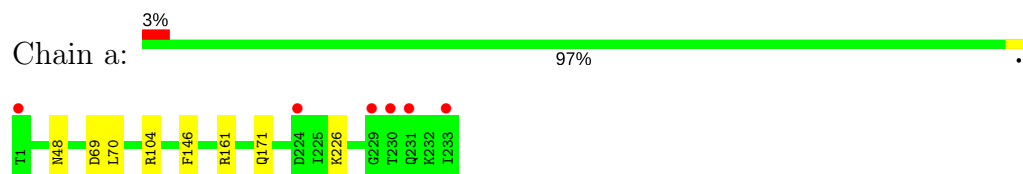
- Molecule 12: Proteasome component C5



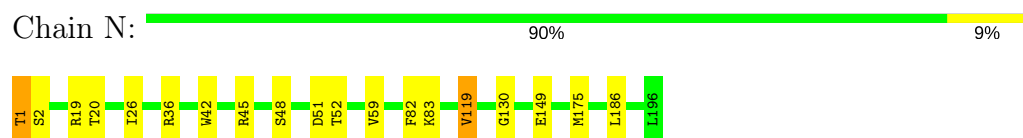
- Molecule 13: Proteasome component PRE4



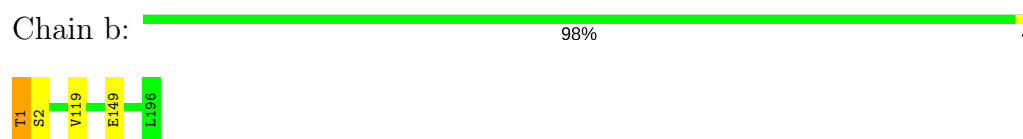
- Molecule 13: Proteasome component PRE4



- Molecule 14: Proteasome component PRE3



- Molecule 14: Proteasome component PRE3



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	134.41Å 300.78Å 143.82Å 90.00° 112.75° 90.00°	Depositor
Resolution (Å)	15.00 – 3.40 49.31 – 3.40	Depositor EDS
% Data completeness (in resolution range)	98.3 (15.00-3.40) 98.3 (49.31-3.40)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.67 (at 3.40Å)	Xtriage
Refinement program	REFMAC 5.6.0119	Depositor
R, R_{free}	0.179 , 0.219 0.176 , 0.216	Depositor DCC
R_{free} test set	7082 reflections (5.26%)	DCC
Wilson B-factor (Å ²)	66.8	Xtriage
Anisotropy	0.786	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 64.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	51112	wwPDB-VP
Average B, all atoms (Å ²)	80.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.25% 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: 04C

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	0.36	2/1952 (0.1%)	0.48	0/2642
1	O	0.36	2/1952 (0.1%)	0.48	0/2642
2	B	0.33	0/1934	0.48	0/2618
2	P	0.33	0/1934	0.48	0/2618
3	C	0.34	0/1919	0.49	0/2598
3	Q	0.33	0/1919	0.49	0/2598
4	D	0.36	0/1886	0.50	0/2541
4	R	0.35	0/1886	0.50	0/2541
5	E	0.31	0/1823	0.48	0/2463
5	S	0.30	0/1823	0.48	0/2463
6	F	0.41	1/1936 (0.1%)	0.47	0/2614
6	T	0.40	0/1936	0.47	0/2614
7	G	0.34	0/1959	0.48	0/2652
7	U	0.34	0/1959	0.48	0/2652
8	H	0.46	1/1715 (0.1%)	0.48	0/2326
8	V	0.46	2/1715 (0.1%)	0.49	0/2326
9	I	0.35	1/1611 (0.1%)	0.49	0/2174
9	W	0.34	0/1611	0.49	0/2174
10	J	0.31	0/1613	0.48	0/2173
10	X	0.31	0/1613	0.48	0/2173
11	K	0.52	3/1681 (0.2%)	0.52	1/2274 (0.0%)
11	Y	0.52	3/1681 (0.2%)	0.52	1/2274 (0.0%)
12	L	0.36	0/1795	0.50	0/2420
12	Z	0.36	0/1795	0.50	0/2420
13	M	0.36	1/1855 (0.1%)	0.50	0/2514
13	a	0.36	0/1855	0.50	0/2514
14	N	0.43	1/1541 (0.1%)	0.49	1/2087 (0.0%)
14	b	0.43	1/1541 (0.1%)	0.51	1/2087 (0.0%)
All	All	0.38	18/50440 (0.0%)	0.49	4/68192 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if

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

Mol	Chain	#Chirality outliers	#Planarity outliers
14	b	0	1

All (18) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
14	b	1	THR	C-N	6.64	1.49	1.34
14	N	1	THR	C-N	6.25	1.48	1.34
8	V	1	THR	C-N	5.61	1.47	1.34
11	K	1	THR	C-N	5.37	1.46	1.34
8	H	1	THR	C-N	5.34	1.46	1.34
11	Y	1	THR	C-N	5.12	1.45	1.34
6	F	154	TRP	CD2-CE2	5.07	1.47	1.41
13	M	219	TRP	CD2-CE2	5.05	1.47	1.41
1	O	159	TRP	CD2-CE2	5.04	1.47	1.41
8	V	142	TRP	CD2-CE2	5.04	1.47	1.41
1	A	159	TRP	CD2-CE2	5.04	1.47	1.41
11	Y	146	TRP	CD2-CE2	5.02	1.47	1.41
11	Y	58	TRP	CD2-CE2	5.02	1.47	1.41
1	A	179	TRP	CD2-CE2	5.02	1.47	1.41
1	O	179	TRP	CD2-CE2	5.01	1.47	1.41
9	I	182	TRP	CD2-CE2	5.01	1.47	1.41
11	K	58	TRP	CD2-CE2	5.01	1.47	1.41
11	K	146	TRP	CD2-CE2	5.00	1.47	1.41

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
14	b	1	THR	N-CA-C	6.01	127.24	111.00
11	Y	4	LEU	CA-CB-CG	5.22	127.30	115.30
11	K	4	LEU	CA-CB-CG	5.21	127.27	115.30
14	N	1	THR	N-CA-C	5.11	124.79	111.00

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
14	b	1	THR	Mainchain

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	1915	0	1929	17	0
1	O	1915	0	1929	20	0
2	B	1904	0	1904	33	0
2	P	1904	0	1904	31	0
3	C	1890	0	1903	40	0
3	Q	1890	0	1903	34	0
4	D	1861	0	1839	19	0
4	R	1861	0	1839	20	0
5	E	1795	0	1800	31	0
5	S	1795	0	1800	30	0
6	F	1896	0	1889	21	0
6	T	1896	0	1889	22	0
7	G	1921	0	1913	25	0
7	U	1921	0	1913	30	0
8	H	1684	0	1686	18	0
8	V	1684	0	1686	18	0
9	I	1581	0	1574	25	0
9	W	1581	0	1574	23	0
10	J	1585	0	1590	41	0
10	X	1585	0	1590	38	0
11	K	1644	0	1593	24	0
11	Y	1644	0	1593	20	0
12	L	1757	0	1711	28	0
12	Z	1757	0	1711	29	0
13	M	1824	0	1832	24	0
13	a	1824	0	1832	0	0
14	N	1512	0	1479	11	0
14	b	1512	0	1479	0	0
15	H	42	0	42	0	0
15	K	42	0	42	4	0
15	N	42	0	42	4	0
15	V	42	0	42	0	0
15	Y	42	0	42	2	0
15	b	42	0	42	0	0
16	A	56	0	0	0	0
16	B	38	0	0	1	0
16	C	42	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
16	D	38	0	0	0	0
16	E	23	0	0	0	0
16	F	48	0	0	0	0
16	G	61	0	0	0	0
16	H	49	0	0	0	0
16	I	62	0	0	1	0
16	J	54	0	0	3	0
16	K	46	0	0	0	0
16	L	55	0	0	0	0
16	M	72	0	0	0	0
16	N	51	0	0	0	0
16	O	32	0	0	0	0
16	P	30	0	0	0	0
16	Q	30	0	0	0	0
16	R	27	0	0	2	0
16	S	20	0	0	0	0
16	T	41	0	0	0	0
16	U	61	0	0	0	0
16	V	45	0	0	0	0
16	W	57	0	0	0	0
16	X	52	0	0	3	0
16	Y	49	0	0	0	0
16	Z	50	0	0	0	0
16	a	75	0	0	0	0
16	b	58	0	0	0	0
All	All	51112	0	49536	582	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 6.

All (582) 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)
5:S:92:ASN:HD21	12:Z:70:ASN:HD21	1.13	0.91
5:E:92:ASN:HD21	12:L:70:ASN:HD21	1.19	0.91
11:K:209:ASN:HB3	16:X:234:HOH:O	1.77	0.85
8:V:35:HIS:HB3	8:V:56:THR:HG21	1.63	0.80
1:O:12:PHE:H	2:P:20:GLN:HE22	1.27	0.80
8:H:35:HIS:HB3	8:H:56:THR:HG21	1.62	0.80
10:J:139:TYR:HD1	16:J:228:HOH:O	1.65	0.79
2:P:200:THR:HG22	2:P:202:SER:H	1.49	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:T:91:GLU:HG2	6:T:111:ARG:HB3	1.64	0.78
2:B:200:THR:HG22	2:B:202:SER:H	1.49	0.78
13:M:161:ARG:HH11	13:M:161:ARG:HG3	1.50	0.77
2:B:12:PHE:H	3:C:17:GLN:HE22	1.32	0.76
3:C:160:GLN:HE21	3:C:160:GLN:HA	1.50	0.76
6:F:91:GLU:HG2	6:F:111:ARG:HB3	1.66	0.76
12:Z:16:ALA:HB2	12:Z:122:VAL:HG23	1.68	0.76
3:Q:160:GLN:HE21	3:Q:160:GLN:HA	1.50	0.76
5:S:12:PHE:H	6:T:19:GLN:HE22	1.31	0.75
15:N:201:04C:H7	15:N:201:04C:H22	1.68	0.75
12:L:16:ALA:HB2	12:L:122:VAL:HG23	1.68	0.75
13:M:43:ILE:HG12	13:M:64:GLU:HG3	1.68	0.75
5:S:92:ASN:HD21	12:Z:70:ASN:ND2	1.85	0.75
12:Z:13:LEU:HD11	12:Z:150:LEU:HD21	1.69	0.74
3:Q:9:PHE:H	4:R:15:GLN:HE22	1.35	0.74
10:J:139:TYR:CE2	10:J:172:MET:HG3	2.23	0.74
1:O:128:ARG:HH21	7:U:120:THR:HG22	1.51	0.73
13:M:48:ASN:H	13:M:48:ASN:HD22	1.35	0.72
10:X:139:TYR:CE2	10:X:172:MET:HG3	2.24	0.72
12:L:13:LEU:HD11	12:L:150:LEU:HD21	1.71	0.72
4:R:161:ALA:HB3	5:S:55:LEU:HD23	1.71	0.71
5:S:87:LEU:HD11	5:S:107:ALA:HB1	1.72	0.71
5:E:12:PHE:H	6:F:19:GLN:HE22	1.37	0.71
2:P:95:GLN:HE22	9:W:71:ASN:HD22	1.39	0.71
3:C:9:PHE:H	4:D:15:GLN:HE22	1.39	0.71
5:S:92:ASN:ND2	12:Z:70:ASN:HD21	1.87	0.71
5:S:205:LEU:HA	5:S:209:ASN:HD22	1.56	0.70
5:E:205:LEU:HA	5:E:209:ASN:HD22	1.56	0.70
2:B:122:THR:HG22	3:C:125:ARG:HH21	1.55	0.70
5:E:87:LEU:HD11	5:E:107:ALA:HB1	1.71	0.70
1:A:12:PHE:H	2:B:20:GLN:HE22	1.37	0.70
10:X:39:SER:HB2	10:X:40:PRO:HD2	1.73	0.70
1:A:128:ARG:HH21	7:G:120:THR:HG22	1.56	0.70
10:J:39:SER:HB2	10:J:40:PRO:HD2	1.73	0.70
12:L:195:HIS:HD2	12:L:197:GLN:H	1.40	0.70
2:P:124:HIS:HB3	3:Q:124:VAL:HG12	1.74	0.70
3:C:160:GLN:HE22	3:C:170:ARG:HE	1.41	0.69
7:G:103:MET:HE3	7:G:108:LEU:HD13	1.75	0.69
7:U:103:MET:HE3	7:U:108:LEU:HD13	1.75	0.69
11:K:209:ASN:ND2	9:W:37:ASN:HD21	1.90	0.69
5:E:92:ASN:HD21	12:L:70:ASN:ND2	1.89	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
12:Z:195:HIS:HD2	12:Z:197:GLN:H	1.40	0.68
2:B:95:GLN:HE22	9:I:71:ASN:HD22	1.40	0.68
7:G:187:GLU:HG2	7:G:192:LYS:HB2	1.74	0.68
10:J:150:PRO:HG3	11:Y:208:ASN:HD21	1.59	0.68
7:U:187:GLU:HG2	7:U:192:LYS:HB2	1.74	0.68
11:K:73:ARG:NH2	11:K:104:TYR:O	2.27	0.67
13:M:179:ASN:HD22	13:M:182:ARG:HH11	1.40	0.67
12:Z:13:LEU:CD1	12:Z:150:LEU:HD21	2.25	0.67
8:H:104:ASP:HB2	8:H:105:PRO:HD2	1.77	0.67
11:K:208:ASN:HD21	10:X:150:PRO:HG3	1.60	0.67
11:Y:73:ARG:NH2	11:Y:104:TYR:O	2.28	0.67
2:P:122:THR:HG22	3:Q:125:ARG:HH21	1.60	0.66
9:I:148:MET:HE3	9:I:152:LEU:HD11	1.77	0.66
12:L:13:LEU:CD1	12:L:150:LEU:HD21	2.24	0.66
3:Q:160:GLN:HE22	3:Q:170:ARG:HE	1.42	0.65
8:V:104:ASP:HB2	8:V:105:PRO:HD2	1.78	0.65
8:V:35:HIS:CB	8:V:56:THR:HG21	2.27	0.65
2:B:124:HIS:HB3	3:C:124:VAL:HG12	1.77	0.65
9:I:37:ASN:HD21	11:Y:209:ASN:ND2	1.94	0.64
6:F:32:THR:HG22	6:F:47:GLU:OE2	1.97	0.64
1:A:122:THR:HG22	2:B:128:ARG:HH21	1.62	0.64
5:E:92:ASN:ND2	12:L:70:ASN:HD21	1.91	0.64
8:H:35:HIS:CB	8:H:56:THR:HG21	2.27	0.64
6:T:32:THR:HG22	6:T:47:GLU:OE2	1.98	0.64
7:G:195:GLU:HG3	7:G:235:ARG:HG3	1.78	0.63
1:O:122:THR:HG22	2:P:128:ARG:HH21	1.62	0.63
2:P:12:PHE:H	3:Q:17:GLN:HE22	1.47	0.63
7:U:195:GLU:HG3	7:U:235:ARG:HG3	1.79	0.62
13:M:48:ASN:HD22	13:M:48:ASN:N	1.97	0.62
5:E:205:LEU:HA	5:E:209:ASN:ND2	2.15	0.61
12:L:3:ASN:HD22	12:L:4:PRO:HD2	1.65	0.61
6:F:31:THR:HG21	6:F:47:GLU:O	2.01	0.61
12:Z:3:ASN:HD22	12:Z:4:PRO:HD2	1.65	0.61
14:N:1:THR:HG22	14:N:2:SER:N	2.15	0.61
6:T:31:THR:HG21	6:T:47:GLU:O	2.00	0.61
2:B:180:LYS:HG3	2:B:183:MET:HG3	1.83	0.61
5:S:205:LEU:HA	5:S:209:ASN:ND2	2.15	0.61
9:I:37:ASN:ND2	9:I:37:ASN:H	1.99	0.60
4:R:73:LEU:HD12	4:R:131:GLY:HA3	1.83	0.60
3:C:161:THR:HG21	3:C:169:VAL:HG13	1.83	0.60
14:N:175:MET:HB2	14:N:186:LEU:HB2	1.83	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:S:127:TYR:O	5:S:148:PRO:HB3	2.01	0.60
3:Q:161:THR:HG21	3:Q:169:VAL:HG13	1.82	0.60
2:B:63:GLU:HG3	2:B:64:LYS:HG3	1.82	0.60
10:J:149:ARG:HB2	10:J:152:MET:HG3	1.84	0.60
11:K:87:VAL:HG11	11:K:117:SER:HA	1.83	0.60
11:K:208:ASN:ND2	10:X:150:PRO:HG3	2.16	0.60
2:P:63:GLU:HG3	2:P:64:LYS:HG3	1.82	0.60
3:Q:169:VAL:HG23	3:Q:196:SER:HB2	1.83	0.60
5:E:127:TYR:O	5:E:148:PRO:HB3	2.01	0.60
9:W:148:MET:HE3	9:W:152:LEU:HD11	1.84	0.60
10:J:1:MET:HG2	10:J:2:ASP:H	1.67	0.59
4:D:73:LEU:HD12	4:D:131:GLY:HA3	1.84	0.59
1:A:30:GLN:HE21	1:A:30:GLN:HA	1.67	0.59
10:J:150:PRO:HG3	11:Y:208:ASN:ND2	2.16	0.59
10:X:149:ARG:HB2	10:X:152:MET:HG3	1.84	0.59
7:G:117:GLN:O	7:G:120:THR:HB	2.02	0.59
9:I:36:SER:HB2	10:J:126:VAL:HG21	1.84	0.59
6:T:91:GLU:HG3	6:T:111:ARG:HH11	1.67	0.59
12:Z:126:ASP:HB2	12:Z:130:SER:HB3	1.85	0.59
3:C:169:VAL:HG23	3:C:196:SER:HB2	1.83	0.59
10:X:3:ILE:HD13	10:X:168:LEU:HD13	1.83	0.59
2:B:18:LEU:HD13	2:B:122:THR:HG23	1.83	0.59
6:F:91:GLU:HG3	6:F:111:ARG:HH11	1.66	0.59
13:M:179:ASN:HD22	13:M:182:ARG:NH1	1.99	0.59
2:P:18:LEU:HD13	2:P:122:THR:HG23	1.85	0.58
5:E:197:SER:HA	5:E:200:LEU:HG	1.84	0.58
11:K:38:ASN:HB2	11:K:39:PRO:HD2	1.85	0.58
1:O:30:GLN:HE21	1:O:30:GLN:HA	1.68	0.58
3:C:204:GLY:HA3	3:C:207:ASN:HB2	1.86	0.58
10:X:1:MET:HG2	10:X:2:ASP:H	1.69	0.58
11:Y:87:VAL:HG11	11:Y:117:SER:HA	1.83	0.58
11:Y:38:ASN:HB2	11:Y:39:PRO:HD2	1.86	0.58
10:J:3:ILE:HD13	10:J:168:LEU:HD13	1.85	0.58
12:L:126:ASP:HB2	12:L:130:SER:HB3	1.86	0.58
7:U:117:GLN:O	7:U:120:THR:HB	2.04	0.58
2:P:180:LYS:HG3	2:P:183:MET:HG3	1.84	0.57
5:S:197:SER:HA	5:S:200:LEU:HG	1.85	0.57
9:W:37:ASN:ND2	9:W:37:ASN:H	2.01	0.57
2:B:151:ASN:HB2	2:B:152:PRO:HD2	1.84	0.57
13:M:227:GLY:HA3	13:M:231:GLN:HB3	1.85	0.57
1:O:119:GLN:O	1:O:122:THR:HB	2.05	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:P:151:ASN:HB2	2:P:152:PRO:HD2	1.85	0.57
9:W:36:SER:HB2	10:X:126:VAL:HG21	1.86	0.57
11:K:87:VAL:CG1	11:K:117:SER:HA	2.35	0.57
2:B:215:ILE:HG12	2:B:226:GLN:HG2	1.87	0.57
2:B:122:THR:CG2	3:C:125:ARG:HH21	2.18	0.57
9:I:97:ARG:HD2	16:I:360:HOH:O	2.05	0.57
9:I:94:LEU:HD11	9:I:106:PRO:HG2	1.87	0.56
10:X:147:HIS:HB2	10:X:160:LEU:HD11	1.87	0.56
1:A:119:GLN:O	1:A:122:THR:HB	2.05	0.56
11:Y:87:VAL:CG1	11:Y:117:SER:HA	2.34	0.56
3:Q:204:GLY:HA3	3:Q:207:ASN:HB2	1.87	0.56
10:J:25:ILE:O	10:X:139:TYR:OH	2.23	0.56
2:P:215:ILE:HG12	2:P:226:GLN:HG2	1.87	0.56
4:D:176:LEU:HD22	5:E:55:LEU:HD13	1.88	0.56
2:P:75:ALA:HB3	2:P:135:ILE:HB	1.88	0.55
10:J:139:TYR:OH	10:X:25:ILE:O	2.24	0.55
9:W:14:MET:HB3	9:W:162:LEU:HD11	1.87	0.55
2:B:3:ARG:HB2	5:E:122:TYR:OH	2.06	0.55
14:N:1:THR:CG2	14:N:2:SER:N	2.70	0.55
9:W:9:GLY:HA3	9:W:41:LYS:HE2	1.88	0.55
7:G:167:GLN:HE21	7:G:171:THR:HG23	1.71	0.55
9:W:52:ILE:HB	9:W:59:VAL:HG13	1.89	0.55
11:Y:107:LYS:H	11:Y:107:LYS:HD2	1.71	0.55
10:X:46:PHE:HD1	10:X:53:THR:HG1	1.55	0.55
11:Y:45:MET:HG2	11:Y:52:CYS:HB3	1.87	0.55
15:N:201:04C:H8	15:N:201:04C:H24	1.88	0.55
9:I:9:GLY:HA3	9:I:41:LYS:HE2	1.88	0.55
10:J:147:HIS:HB2	10:J:160:LEU:HD11	1.89	0.55
9:W:94:LEU:HD11	9:W:106:PRO:HG2	1.88	0.55
2:B:151:ASN:HB2	2:B:152:PRO:CD	2.38	0.54
9:I:14:MET:HB3	9:I:162:LEU:HD11	1.88	0.54
7:U:167:GLN:HE21	7:U:171:THR:HG23	1.72	0.54
3:Q:185:THR:HB	3:Q:188:GLU:HG2	1.90	0.54
10:X:7:ILE:HD11	10:X:160:LEU:HD23	1.89	0.54
11:K:107:LYS:H	11:K:107:LYS:HD2	1.73	0.54
3:C:185:THR:HB	3:C:188:GLU:HG2	1.89	0.54
13:M:209:LYS:HB3	13:M:212:LEU:HD11	1.90	0.54
1:A:222:LEU:HD13	1:A:232:GLY:HA2	1.88	0.54
10:J:46:PHE:HD1	10:J:53:THR:HG1	1.55	0.54
14:N:36:ARG:HG3	14:N:42:TRP:CE2	2.42	0.54
1:A:55:LEU:HD12	7:G:170:THR:HG23	1.89	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
13:M:27:LEU:HD21	13:M:34:LEU:HD22	1.89	0.53
2:B:75:ALA:HB3	2:B:135:ILE:HB	1.90	0.53
5:E:12:PHE:HB2	6:F:19:GLN:HE22	1.73	0.53
11:K:45:MET:HG2	11:K:52:CYS:HB3	1.90	0.53
2:P:151:ASN:HB2	2:P:152:PRO:CD	2.38	0.53
10:J:7:ILE:HD11	10:J:160:LEU:HD23	1.90	0.53
1:O:222:LEU:HD13	1:O:232:GLY:HA2	1.90	0.53
3:C:157:TRP:CE2	4:D:51:LEU:HD23	2.44	0.53
5:S:205:LEU:HD23	5:S:205:LEU:H	1.74	0.53
10:J:139:TYR:HE1	16:X:236:HOH:O	1.91	0.53
6:T:154:TRP:CZ3	7:U:60:VAL:HA	2.44	0.53
9:I:52:ILE:HB	9:I:59:VAL:HG13	1.92	0.52
4:R:159:TYR:CE2	5:S:56:SER:HB3	2.44	0.52
8:V:113:ILE:HG12	8:V:119:THR:HG22	1.92	0.52
1:O:14:PRO:HA	2:P:23:TYR:CD1	2.44	0.52
8:V:163:ILE:HG23	8:V:170:GLY:HA2	1.91	0.52
2:B:69:ASN:ND2	2:B:70:ASP:H	2.07	0.52
9:I:35:VAL:HG13	16:J:242:HOH:O	2.08	0.52
12:L:42:LYS:HD2	12:L:55:ASN:HD22	1.73	0.52
5:E:205:LEU:HD23	5:E:205:LEU:H	1.75	0.52
1:O:68:THR:HB	1:O:69:PRO:HD2	1.92	0.52
8:H:113:ILE:HG12	8:H:119:THR:HG22	1.92	0.52
14:N:19:ARG:HG3	14:N:26:ILE:HG23	1.92	0.51
1:A:68:THR:HB	1:A:69:PRO:HD2	1.91	0.51
3:C:35:LYS:HG2	3:C:158:SER:O	2.10	0.51
3:C:160:GLN:HE21	3:C:160:GLN:CA	2.22	0.51
13:M:15:LYS:HB3	13:M:20:VAL:HG12	1.91	0.51
12:Z:90:ALA:HA	12:Z:125:PHE:HZ	1.75	0.51
6:F:33:SER:HB3	6:F:46:VAL:HG23	1.92	0.51
8:H:1:THR:HG23	8:H:33:LYS:HD3	1.93	0.51
6:T:31:THR:HG23	6:T:47:GLU:HB3	1.91	0.51
12:Z:8:ASN:HA	12:Z:30:ILE:O	2.10	0.51
3:C:66:ASP:HA	10:J:69:ILE:HD13	1.93	0.51
3:Q:157:TRP:CE2	4:R:51:LEU:HD23	2.46	0.51
5:S:226:GLY:O	5:S:229:VAL:HG22	2.11	0.51
2:B:139:TYR:CD1	2:B:224:VAL:HG21	2.46	0.51
12:L:135:GLN:HG3	12:L:174:TYR:OH	2.11	0.51
12:L:8:ASN:HA	12:L:30:ILE:O	2.10	0.51
13:M:161:ARG:NH1	13:M:161:ARG:HG3	2.24	0.51
14:N:45:ARG:HD2	14:N:52:THR:HB	1.93	0.51
3:Q:27:ARG:HB2	3:Q:27:ARG:HH11	1.76	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:35:ILE:HD12	2:B:196:LEU:HG	1.92	0.51
10:X:119:ILE:HG12	10:X:125:LYS:HG3	1.92	0.51
6:F:216:SER:HB3	6:F:219:GLU:HB2	1.93	0.50
6:F:31:THR:HG23	6:F:47:GLU:HB3	1.93	0.50
6:T:216:SER:HB3	6:T:219:GLU:HB2	1.93	0.50
3:C:29:THR:HB	3:C:45:GLU:HG3	1.93	0.50
2:P:69:ASN:ND2	2:P:70:ASP:H	2.08	0.50
12:Z:126:ASP:CB	12:Z:130:SER:HB3	2.41	0.50
13:M:97:ALA:HA	13:M:130:VAL:HG21	1.93	0.50
2:P:139:TYR:CD1	2:P:224:VAL:HG21	2.46	0.50
11:Y:99:THR:HG22	11:Y:115:VAL:O	2.11	0.50
10:J:174:MET:HE2	10:X:174:MET:HE2	1.92	0.50
6:T:158:GLY:O	7:U:54:LEU:HB3	2.11	0.50
2:B:119:GLN:O	2:B:122:THR:HB	2.11	0.50
12:Z:135:GLN:HG3	12:Z:174:TYR:OH	2.11	0.50
12:L:126:ASP:CB	12:L:130:SER:HB3	2.42	0.50
1:A:110:LEU:O	1:A:114:VAL:HG23	2.11	0.50
5:E:200:LEU:HD11	5:E:205:LEU:HD22	1.94	0.50
7:G:78:ILE:N	7:G:79:PRO:HD2	2.26	0.50
13:M:27:LEU:HB2	13:M:192:SER:HB2	1.93	0.50
5:S:230:ALA:HA	5:S:233:ILE:HD12	1.94	0.50
7:U:78:ILE:N	7:U:79:PRO:HD2	2.26	0.50
3:C:214:LYS:HB2	3:C:218:ASP:HB3	1.94	0.50
5:S:70:GLY:HA3	5:S:221:PHE:CE2	2.47	0.50
12:Z:42:LYS:HD2	12:Z:55:ASN:HD22	1.77	0.50
3:C:27:ARG:HB2	3:C:27:ARG:HH11	1.77	0.49
13:M:16:TYR:CE2	13:M:170:VAL:HG22	2.47	0.49
2:P:119:GLN:O	2:P:122:THR:HB	2.12	0.49
2:P:175:LEU:HD23	2:P:191:LEU:HD22	1.93	0.49
3:Q:27:ARG:HB2	3:Q:27:ARG:NH1	2.27	0.49
10:X:7:ILE:CD1	10:X:160:LEU:HD23	2.42	0.49
5:E:226:GLY:O	5:E:229:VAL:HG22	2.11	0.49
7:G:187:GLU:HG2	7:G:192:LYS:CB	2.42	0.49
12:L:90:ALA:HA	12:L:125:PHE:HZ	1.76	0.49
7:U:63:ILE:HD12	7:U:215:GLU:HG2	1.94	0.49
3:C:65:ILE:HG21	3:C:107:LEU:HD21	1.95	0.49
6:F:8:ASN:HB3	6:F:123:ASN:HA	1.94	0.49
5:S:68:HIS:HE1	5:S:102:LEU:O	1.95	0.49
5:E:193:VAL:HG13	5:E:205:LEU:HD11	1.94	0.49
7:G:63:ILE:HD12	7:G:215:GLU:HG2	1.95	0.49
13:M:193:ARG:HG3	13:M:214:VAL:HB	1.93	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:Q:214:LYS:HB2	3:Q:218:ASP:HB3	1.94	0.49
3:Q:65:ILE:HG21	3:Q:107:LEU:HD21	1.95	0.49
5:S:9:THR:HG21	5:S:119:THR:HA	1.95	0.49
6:T:33:SER:HB3	6:T:46:VAL:HG23	1.93	0.49
2:B:223:GLU:HG2	2:B:224:VAL:H	1.77	0.49
8:H:163:ILE:HG23	8:H:170:GLY:HA2	1.93	0.49
10:J:119:ILE:HG12	10:J:125:LYS:HG3	1.94	0.49
1:A:211:LEU:HD22	1:A:238:LEU:HD12	1.95	0.49
3:C:27:ARG:NH1	3:C:27:ARG:HB2	2.28	0.49
2:P:99:LYS:HG2	10:X:86:GLN:HE22	1.78	0.49
9:I:10:ILE:HG21	9:I:141:ALA:HB3	1.94	0.49
8:H:84:LYS:HG3	8:H:85:GLN:N	2.28	0.49
1:O:110:LEU:O	1:O:114:VAL:HG23	2.13	0.49
3:Q:185:THR:HG22	3:Q:187:GLU:H	1.78	0.49
12:L:109:THR:HG23	12:L:125:PHE:HB2	1.94	0.49
10:J:139:TYR:OH	10:X:25:ILE:HG12	2.13	0.48
3:C:155:SER:HB2	4:D:51:LEU:HD21	1.95	0.48
10:J:7:ILE:CD1	10:J:160:LEU:HD23	2.43	0.48
5:E:68:HIS:HE1	5:E:102:LEU:O	1.95	0.48
7:G:30:ASN:HD22	7:G:164:PRO:HG2	1.78	0.48
5:S:193:VAL:HG13	5:S:205:LEU:HD11	1.95	0.48
1:O:211:LEU:HD22	1:O:238:LEU:HD12	1.95	0.48
2:P:122:THR:CG2	3:Q:125:ARG:HH21	2.26	0.48
5:S:134:ILE:HD12	5:S:215:VAL:HG12	1.94	0.48
2:B:175:LEU:HD23	2:B:191:LEU:HD22	1.94	0.48
5:E:230:ALA:HA	5:E:233:ILE:HD12	1.95	0.48
5:E:80:ALA:HB2	5:E:129:VAL:HG21	1.96	0.48
2:P:35:ILE:HD12	2:P:196:LEU:HG	1.95	0.48
3:Q:29:THR:HB	3:Q:45:GLU:HG3	1.94	0.48
9:W:10:ILE:HG21	9:W:141:ALA:HB3	1.96	0.48
10:X:139:TYR:HE2	10:X:172:MET:HG3	1.75	0.48
12:L:145:LEU:O	9:W:147:GLY:HA3	2.14	0.48
5:S:200:LEU:HD11	5:S:205:LEU:HD22	1.95	0.48
5:S:80:ALA:HB2	5:S:129:VAL:HG21	1.95	0.48
3:C:185:THR:HG22	3:C:187:GLU:H	1.78	0.48
5:S:49:LYS:HB3	5:S:58:TYR:HB3	1.96	0.48
2:B:148:TYR:OH	3:C:57:ILE:HB	2.14	0.47
12:L:124:SER:CB	12:L:137:ARG:HG2	2.43	0.47
1:O:23:TYR:CD1	7:U:12:PRO:HA	2.49	0.47
8:V:3:ILE:HG22	8:V:16:ALA:HB2	1.96	0.47
9:W:62:LEU:CD1	9:W:104:VAL:HG21	2.43	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
10:J:174:MET:HE3	10:X:174:MET:HE3	1.95	0.47
2:B:142:ARG:HB2	16:B:332:HOH:O	2.14	0.47
2:B:172:GLN:HG2	3:C:50:LEU:HD12	1.96	0.47
4:D:138:GLY:HA2	4:D:214:ILE:HG12	1.95	0.47
9:I:62:LEU:CD1	9:I:104:VAL:HG21	2.45	0.47
13:M:156:ARG:HH11	8:V:165:ASN:HD22	1.60	0.47
4:R:32:ILE:HD12	4:R:192:VAL:HG23	1.97	0.47
5:S:227:GLU:CD	5:S:227:GLU:H	2.16	0.47
7:U:187:GLU:HG2	7:U:192:LYS:CB	2.43	0.47
8:V:84:LYS:HG3	8:V:85:GLN:N	2.29	0.47
3:C:92:GLN:HG3	10:J:66:LEU:HB2	1.97	0.47
15:K:301:04C:H7	15:K:301:04C:H24	1.97	0.47
5:E:9:THR:HG21	5:E:119:THR:HA	1.95	0.47
4:R:138:GLY:HA2	4:R:214:ILE:HG12	1.95	0.47
5:E:70:GLY:HA3	5:E:221:PHE:CE2	2.49	0.47
5:E:227:GLU:CD	5:E:227:GLU:H	2.17	0.47
3:Q:35:LYS:HG2	3:Q:158:SER:O	2.14	0.47
4:R:44:LYS:HE3	4:R:210:GLN:HB2	1.96	0.47
6:T:155:GLY:HA3	7:U:59:THR:HG21	1.96	0.47
12:Z:124:SER:CB	12:Z:137:ARG:HG2	2.44	0.47
9:I:53:THR:O	9:I:104:VAL:HA	2.15	0.47
11:K:99:THR:HG22	11:K:115:VAL:O	2.15	0.47
12:L:42:LYS:HD2	12:L:55:ASN:ND2	2.30	0.47
7:U:70:ILE:HG21	7:U:108:LEU:HD23	1.97	0.47
9:I:10:ILE:HD11	9:I:174:ALA:HB2	1.97	0.46
4:R:113:LEU:HB2	16:R:315:HOH:O	2.15	0.46
6:T:110:ASP:O	6:T:114:GLN:HG2	2.15	0.46
12:L:124:SER:HB3	12:L:137:ARG:HG2	1.97	0.46
13:M:161:ARG:HH11	13:M:161:ARG:CG	2.26	0.46
9:W:10:ILE:HD11	9:W:174:ALA:HB2	1.97	0.46
5:E:49:LYS:HB3	5:E:58:TYR:HB3	1.96	0.46
2:P:223:GLU:HG2	2:P:224:VAL:H	1.79	0.46
2:P:79:LEU:HD23	2:P:131:GLY:HA3	1.97	0.46
4:R:77:ALA:O	4:R:81:ILE:HG12	2.15	0.46
9:W:185:VAL:HG21	9:W:196:LYS:HE3	1.97	0.46
12:Z:109:THR:HG23	12:Z:125:PHE:HB2	1.97	0.46
12:Z:17:GLY:HA3	12:Z:20:PHE:CE2	2.51	0.46
14:N:20:THR:HG22	15:N:201:04C:C5	2.46	0.46
12:L:126:ASP:OD2	12:L:127:PRO:HD2	2.15	0.46
1:O:26:THR:O	1:O:30:GLN:HG2	2.15	0.46
6:T:8:ASN:HB3	6:T:123:ASN:HA	1.96	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
12:Z:124:SER:HB3	12:Z:137:ARG:HG2	1.98	0.46
5:E:134:ILE:HD12	5:E:215:VAL:HG12	1.96	0.46
9:I:147:GLY:HA3	12:Z:145:LEU:O	2.15	0.46
7:U:55:LEU:O	7:U:57:PRO:HD3	2.16	0.46
8:V:22:GLN:HG3	8:V:27:ALA:HB2	1.96	0.46
7:U:30:ASN:HD22	7:U:164:PRO:HG2	1.79	0.46
2:B:79:LEU:HD23	2:B:131:GLY:HA3	1.98	0.46
4:D:44:LYS:HE3	4:D:210:GLN:HB2	1.97	0.46
13:M:27:LEU:HD11	13:M:34:LEU:HB3	1.98	0.46
14:N:48:SER:HB3	14:N:51:ASP:HB2	1.98	0.46
5:E:231:LYS:H	5:E:231:LYS:HD2	1.81	0.45
10:J:53:THR:HG23	10:J:54:VAL:N	2.32	0.45
12:Z:126:ASP:OD2	12:Z:127:PRO:HD2	2.17	0.45
3:Q:160:GLN:HE21	3:Q:160:GLN:CA	2.22	0.45
12:Z:49:ASN:HD22	12:Z:49:ASN:HA	1.63	0.45
8:H:22:GLN:HG3	8:H:27:ALA:HB2	1.97	0.45
12:L:17:GLY:HA3	12:L:20:PHE:CE2	2.52	0.45
11:Y:201:LYS:HG3	11:Y:207:PHE:HB2	1.99	0.45
6:F:110:ASP:O	6:F:114:GLN:HG2	2.17	0.45
7:G:70:ILE:HG21	7:G:108:LEU:HD23	1.99	0.45
1:A:149:GLN:O	1:A:156:TYR:HA	2.17	0.45
8:H:3:ILE:HG22	8:H:16:ALA:HB2	1.99	0.45
11:K:162:LEU:HD12	11:K:193:VAL:HG13	1.98	0.45
8:H:104:ASP:HB2	8:H:105:PRO:CD	2.46	0.45
1:O:149:GLN:O	1:O:156:TYR:HA	2.16	0.45
4:R:71:SER:HB3	4:R:164:ILE:HD12	1.98	0.45
5:S:231:LYS:HD2	5:S:231:LYS:H	1.82	0.45
1:A:26:THR:O	1:A:30:GLN:HG2	2.17	0.45
4:D:186:LYS:O	4:D:190:LEU:HD22	2.16	0.45
4:D:32:ILE:HD12	4:D:192:VAL:HG23	1.99	0.45
6:F:66:VAL:HG11	6:F:108:PHE:CE1	2.52	0.45
10:J:146:HIS:HD2	10:J:147:HIS:CE1	2.35	0.45
3:Q:190:VAL:HG13	3:Q:210:ILE:HG21	1.99	0.45
3:Q:51:LYS:O	3:Q:52:LEU:HB2	2.17	0.45
4:R:186:LYS:O	4:R:190:LEU:HD22	2.17	0.45
10:X:5:LEU:HD21	10:X:140:THR:HG21	1.99	0.45
5:E:12:PHE:HB2	6:F:19:GLN:NE2	2.32	0.44
7:G:55:LEU:O	7:G:57:PRO:HD3	2.17	0.44
10:J:172:MET:HA	10:J:173:PRO:HD3	1.79	0.44
4:R:159:TYR:CG	4:R:162:LYS:HB2	2.51	0.44
5:S:155:LEU:HD23	6:T:55:LEU:HA	1.98	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:W:53:THR:O	9:W:104:VAL:HA	2.17	0.44
3:C:51:LYS:O	3:C:52:LEU:HB2	2.16	0.44
10:J:1:MET:HG2	10:J:2:ASP:N	2.32	0.44
11:K:31:VAL:HG11	15:K:301:04C:C4	2.47	0.44
11:K:45:MET:HE1	11:K:53:GLN:HG3	2.00	0.44
5:S:70:GLY:HA3	5:S:221:PHE:CZ	2.52	0.44
3:C:190:VAL:HG13	3:C:210:ILE:HG21	1.99	0.44
7:G:52:ASP:HB3	7:G:55:LEU:HG	1.99	0.44
10:J:158:LEU:HD13	10:J:198:GLN:HE22	1.82	0.44
9:W:73:TYR:CE1	9:W:77:GLU:HG3	2.52	0.44
10:J:174:MET:HE1	10:X:173:PRO:HB2	1.99	0.44
11:Y:162:LEU:HD12	11:Y:193:VAL:HG13	1.99	0.44
11:Y:31:VAL:HG11	15:Y:301:04C:C4	2.47	0.44
9:I:8:GLY:HA2	9:I:140:THR:HG21	1.99	0.44
10:J:25:ILE:HG12	10:X:139:TYR:OH	2.18	0.44
1:O:75:TYR:HB3	1:O:82:TYR:CD1	2.52	0.44
1:A:75:TYR:HB3	1:A:82:TYR:CD1	2.52	0.44
3:C:239:GLN:C	3:C:241:GLN:H	2.21	0.44
8:H:78:SER:O	8:H:82:MET:HG3	2.17	0.44
9:I:37:ASN:N	9:I:37:ASN:ND2	2.64	0.44
6:T:205:GLU:HG3	6:T:206:LYS:HG3	2.00	0.44
10:X:33:ASP:OD2	10:X:35:THR:HG22	2.18	0.44
4:D:71:SER:HB3	4:D:164:ILE:HD12	1.99	0.44
11:K:201:LYS:HG3	11:K:207:PHE:HB2	2.00	0.44
4:R:59:ILE:HG22	4:R:220:PHE:HZ	1.83	0.44
7:U:45:ILE:HD13	7:U:197:ALA:CB	2.48	0.44
8:V:8:PHE:HB3	8:V:151:ALA:HB2	2.00	0.44
10:X:158:LEU:HD13	10:X:198:GLN:HE22	1.82	0.44
11:Y:158:LYS:HB2	11:Y:177:LEU:HD11	1.99	0.44
6:F:158:GLY:O	7:G:54:LEU:HB3	2.18	0.44
7:G:66:ILE:HG21	7:G:108:LEU:HD21	1.99	0.44
2:P:3:ARG:HB2	5:S:122:TYR:OH	2.17	0.44
6:T:39:ASN:HD22	6:T:39:ASN:N	2.16	0.44
8:V:78:SER:O	8:V:82:MET:HG3	2.17	0.44
10:X:172:MET:HA	10:X:173:PRO:HD3	1.79	0.44
10:X:53:THR:HG23	10:X:54:VAL:N	2.33	0.44
3:Q:70:VAL:HG13	3:Q:219:ILE:HD13	1.99	0.44
8:H:8:PHE:HB3	8:H:151:ALA:HB2	2.00	0.44
14:N:2:SER:OG	14:N:130:GLY:HA3	2.17	0.44
12:Z:42:LYS:HD2	12:Z:55:ASN:ND2	2.33	0.44
1:A:10:THR:HG22	1:A:18:LEU:HD22	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:70:VAL:HG13	3:C:219:ILE:HD13	2.00	0.43
11:K:158:LYS:HB2	11:K:177:LEU:HD11	2.01	0.43
1:A:66:LEU:HD12	1:A:235:PHE:CD2	2.52	0.43
4:D:59:ILE:HG22	4:D:220:PHE:HZ	1.82	0.43
6:F:39:ASN:N	6:F:39:ASN:HD22	2.16	0.43
12:L:100:LYS:HD3	12:L:105:TYR:CE1	2.53	0.43
14:N:59:VAL:HG11	14:N:82:PHE:CE2	2.53	0.43
6:T:66:VAL:HG11	6:T:108:PHE:CE1	2.53	0.43
8:H:167:LEU:HD22	12:Z:196:ILE:O	2.18	0.43
4:D:159:TYR:CG	4:D:162:LYS:HB2	2.53	0.43
3:Q:239:GLN:C	3:Q:241:GLN:H	2.21	0.43
3:Q:92:GLN:HG3	10:X:66:LEU:HB2	1.99	0.43
7:U:66:ILE:HG21	7:U:108:LEU:HD21	2.00	0.43
8:V:80:LEU:HD12	8:V:113:ILE:HD11	2.01	0.43
2:B:213:ALA:HA	2:B:227:LYS:O	2.19	0.43
2:B:223:GLU:HG2	2:B:224:VAL:N	2.34	0.43
10:J:173:PRO:HB2	10:X:174:MET:HE1	2.01	0.43
10:X:3:ILE:HB	10:X:18:SER:HB3	2.00	0.43
9:I:73:TYR:CE1	9:I:77:GLU:HG3	2.53	0.43
3:Q:155:SER:HB2	4:R:51:LEU:HD21	2.00	0.43
13:M:25:ASP:HA	13:M:195:PHE:HA	2.00	0.43
3:C:32:VAL:HG23	3:C:161:THR:HB	2.00	0.43
7:G:34:LEU:HD23	7:G:201:MET:HE3	2.00	0.43
8:V:215:GLU:HG3	9:W:197:ARG:HG2	2.01	0.43
5:E:70:GLY:HA3	5:E:221:PHE:CZ	2.53	0.43
7:G:3:TYR:C	7:G:5:ARG:H	2.22	0.43
5:S:175:LEU:HA	5:S:178:PHE:CE2	2.54	0.43
8:V:104:ASP:HB2	8:V:105:PRO:CD	2.47	0.43
13:M:174:GLU:O	13:M:178:VAL:HG23	2.19	0.43
7:U:70:ILE:HD11	7:U:103:MET:O	2.18	0.43
7:G:45:ILE:HD13	7:G:197:ALA:CB	2.49	0.42
1:O:55:LEU:HD12	7:U:170:THR:HG23	2.01	0.42
7:U:34:LEU:HD23	7:U:201:MET:HE3	2.01	0.42
13:M:179:ASN:ND2	13:M:182:ARG:HH11	2.12	0.42
2:P:213:ALA:HA	2:P:227:LYS:O	2.19	0.42
3:C:226:GLU:O	3:C:230:TYR:HD1	2.03	0.42
5:E:175:LEU:HA	5:E:178:PHE:CE2	2.55	0.42
8:H:215:GLU:HG3	9:I:197:ARG:HG2	2.02	0.42
9:I:185:VAL:HG21	9:I:196:LYS:HE3	2.00	0.42
2:P:236:ASP:O	2:P:240:LYS:HG2	2.19	0.42
11:Y:45:MET:HE2	15:Y:301:04C:H44	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:77:ALA:O	4:D:81:ILE:HG12	2.19	0.42
4:R:159:TYR:CE2	4:R:162:LYS:HD3	2.54	0.42
7:U:149:ASP:HB2	7:U:150:PRO:HD2	2.01	0.42
9:W:62:LEU:HD12	9:W:104:VAL:HG21	2.00	0.42
10:X:146:HIS:HD2	10:X:147:HIS:CE1	2.37	0.42
15:K:301:04C:H11	12:L:128:VAL:HG22	2.02	0.42
10:J:9:VAL:HB	10:J:152:MET:O	2.20	0.42
1:O:66:LEU:HD12	1:O:235:PHE:CD2	2.54	0.42
6:F:7:SER:HB3	6:F:10:VAL:HG23	2.02	0.42
10:J:139:TYR:HE2	10:J:172:MET:HG3	1.76	0.42
7:U:149:ASP:HB2	7:U:150:PRO:CD	2.49	0.42
7:U:52:ASP:HB3	7:U:55:LEU:HG	2.00	0.42
12:Z:100:LYS:O	12:Z:104:PRO:HA	2.20	0.42
12:L:100:LYS:O	12:L:104:PRO:HA	2.20	0.42
10:X:78:GLN:HB2	16:X:213:HOH:O	2.20	0.42
12:Z:195:HIS:CD2	12:Z:197:GLN:H	2.29	0.42
11:K:1:THR:HG23	11:K:33:LYS:HD3	2.02	0.42
2:P:223:GLU:HG2	2:P:224:VAL:N	2.35	0.42
3:Q:165:ASN:O	3:Q:169:VAL:HG12	2.20	0.42
4:R:30:ILE:HD12	4:R:196:LEU:HG	2.02	0.42
10:J:33:ASP:OD2	10:J:35:THR:HG22	2.19	0.42
6:T:78:ILE:HB	6:T:79:PRO:HD3	2.02	0.42
11:K:209:ASN:O	9:W:38:LYS:NZ	2.53	0.42
10:X:92:ILE:HG23	10:X:93:ARG:HG3	2.00	0.42
11:Y:1:THR:HG23	11:Y:33:LYS:HD3	2.01	0.42
6:F:205:GLU:HG3	6:F:206:LYS:HG3	2.01	0.41
12:L:149:PHE:CE1	12:L:153:GLN:HG3	2.54	0.41
6:T:201:GLU:O	6:T:204:LYS:HD2	2.20	0.41
8:V:4:VAL:HG22	8:V:159:ILE:HD11	2.02	0.41
12:L:196:ILE:O	8:V:167:LEU:HD22	2.19	0.41
8:V:1:THR:HA	8:V:33:LYS:HZ3	1.85	0.41
9:W:8:GLY:HA2	9:W:140:THR:HG21	2.02	0.41
1:A:30:GLN:CA	1:A:30:GLN:HE21	2.33	0.41
6:F:154:TRP:CZ3	7:G:60:VAL:HA	2.55	0.41
12:L:22:VAL:HG12	12:L:206:ILE:HG13	2.02	0.41
15:N:201:04C:C35	15:N:201:04C:H22	2.46	0.41
4:R:109:CYS:SG	4:R:156:PHE:HB3	2.60	0.41
3:Q:66:ASP:HA	10:X:69:ILE:HD13	2.02	0.41
2:B:146:GLN:HG2	3:C:57:ILE:HG21	2.01	0.41
4:D:70:MET:HG3	4:D:74:THR:HG22	2.02	0.41
9:I:36:SER:CB	10:J:126:VAL:HG21	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
10:J:3:ILE:HB	10:J:18:SER:HB3	2.02	0.41
7:U:35:ALA:HB2	7:U:44:VAL:HG12	2.03	0.41
7:G:149:ASP:HB2	7:G:150:PRO:CD	2.50	0.41
9:I:11:VAL:HG22	9:I:24:CYS:HB3	2.01	0.41
10:J:139:TYR:CD1	16:J:228:HOH:O	2.53	0.41
12:L:123:TYR:CE1	12:L:133:ARG:HB2	2.56	0.41
3:Q:226:GLU:O	3:Q:230:TYR:HD1	2.03	0.41
3:Q:233:GLN:O	3:Q:237:GLU:HG2	2.20	0.41
3:C:165:ASN:O	3:C:169:VAL:HG12	2.20	0.41
10:J:36:ARG:HD3	10:J:36:ARG:HA	1.91	0.41
11:K:176:ASN:HD22	11:K:176:ASN:HA	1.74	0.41
1:O:10:THR:HG22	1:O:18:LEU:HD22	2.03	0.41
9:W:11:VAL:HG22	9:W:24:CYS:HB3	2.02	0.41
11:Y:37:ILE:HG23	11:Y:60:GLY:HA2	2.02	0.41
2:B:236:ASP:O	2:B:240:LYS:HG2	2.19	0.41
4:D:160:ASN:HB3	4:D:179:TRP:CE2	2.56	0.41
4:D:113:LEU:HD12	5:E:78:PRO:HB2	2.01	0.41
8:H:172:ASN:HD22	8:H:192:THR:HA	1.84	0.41
9:I:62:LEU:HD12	9:I:104:VAL:HG21	2.02	0.41
11:K:14:VAL:HB	11:K:178:TYR:HB2	2.01	0.41
11:K:37:ILE:HG23	11:K:60:GLY:HA2	2.03	0.41
1:O:83:ARG:HH21	7:U:114:ASN:HD22	1.68	0.41
5:S:12:PHE:HB2	6:T:19:GLN:HE22	1.85	0.41
7:U:190:TRP:O	7:U:194:VAL:HG23	2.20	0.41
8:V:172:ASN:HD22	8:V:192:THR:HA	1.86	0.41
10:X:1:MET:HG2	10:X:2:ASP:N	2.34	0.41
6:F:41:GLY:HA3	6:F:215:CYS:O	2.20	0.41
7:G:190:TRP:O	7:G:194:VAL:HG23	2.21	0.41
8:H:4:VAL:HG22	8:H:159:ILE:HD11	2.02	0.41
10:J:53:THR:HG23	10:J:54:VAL:HG23	2.02	0.41
1:O:158:PRO:HB2	2:P:57:GLU:HB3	2.02	0.41
6:T:41:GLY:HA3	6:T:215:CYS:O	2.20	0.41
2:B:180:LYS:CG	2:B:183:MET:HG3	2.49	0.41
3:C:233:GLN:O	3:C:237:GLU:HG2	2.21	0.41
6:F:107:ALA:O	6:F:111:ARG:HG2	2.21	0.41
9:I:14:MET:HE3	9:I:166:ILE:HG13	2.03	0.41
4:R:70:MET:HG3	4:R:74:THR:HG22	2.02	0.41
3:C:147:GLN:HE21	3:C:147:GLN:HB3	1.61	0.41
13:M:48:ASN:ND2	13:M:48:ASN:N	2.68	0.41
1:O:83:ARG:HE	7:U:114:ASN:HD21	1.69	0.41
4:R:176:LEU:HD22	5:S:55:LEU:HD13	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
12:Z:100:LYS:HD3	12:Z:105:TYR:CE1	2.56	0.41
13:M:127:LEU:HG	13:M:142:LEU:HD12	2.01	0.41
3:Q:186:VAL:O	3:Q:190:VAL:HG23	2.21	0.41
3:C:10:SER:HB2	3:C:11:PRO:HD2	2.03	0.41
4:D:161:ALA:HB1	4:D:175:LEU:HD22	2.02	0.41
2:P:15:GLU:O	3:Q:27:ARG:NH1	2.54	0.41
6:T:9:SER:O	7:U:126:ARG:HD3	2.21	0.41
11:Y:14:VAL:HB	11:Y:178:TYR:HB2	2.03	0.41
2:B:14:PRO:HA	3:C:20:TYR:CD1	2.56	0.40
2:B:184:LYS:HD3	2:B:185:VAL:N	2.36	0.40
3:C:146:TYR:CE1	3:C:156:SER:HB3	2.56	0.40
4:D:30:ILE:HD12	4:D:196:LEU:HG	2.02	0.40
7:G:35:ALA:HB2	7:G:44:VAL:HG12	2.04	0.40
13:M:7:THR:HA	13:M:55:GLY:O	2.21	0.40
7:U:3:TYR:C	7:U:5:ARG:H	2.24	0.40
10:X:53:THR:HG23	10:X:54:VAL:HG23	2.03	0.40
8:H:200:GLN:HG2	12:Z:182:LYS:HA	2.03	0.40
8:H:80:LEU:HD12	8:H:113:ILE:HD11	2.02	0.40
11:K:31:VAL:HG11	15:K:301:04C:C3	2.51	0.40
3:Q:32:VAL:HG23	3:Q:161:THR:HB	2.02	0.40
9:W:73:TYR:CZ	9:W:77:GLU:HG3	2.56	0.40
3:C:157:TRP:CZ3	4:D:48:SER:HB3	2.56	0.40
7:G:70:ILE:HD11	7:G:103:MET:O	2.20	0.40
11:K:41:LEU:HD23	11:K:103:GLY:HA3	2.02	0.40
14:N:83:LYS:HG3	14:N:119:VAL:HG22	2.03	0.40
2:P:50:LYS:HG2	2:P:51:VAL:HG23	2.03	0.40
12:Z:149:PHE:CE1	12:Z:153:GLN:HG3	2.56	0.40
10:J:5:LEU:HD21	10:J:140:THR:HG21	2.02	0.40
3:Q:146:TYR:CE1	3:Q:156:SER:HB3	2.57	0.40
10:X:9:VAL:HB	10:X:152:MET:O	2.22	0.40
11:Y:41:LEU:HD23	11:Y:103:GLY:HA3	2.03	0.40
11:Y:105:THR:HB	11:Y:107:LYS:HD2	2.03	0.40
1:A:83:ARG:HH21	7:G:114:ASN:HD22	1.70	0.40
3:C:186:VAL:O	3:C:190:VAL:HG23	2.21	0.40
5:E:28:ILE:HD11	5:E:148:PRO:CD	2.52	0.40
5:E:155:LEU:HD23	6:F:55:LEU:HA	2.03	0.40
10:J:174:MET:HE3	10:X:174:MET:CE	2.51	0.40
10:J:92:ILE:HG23	10:J:93:ARG:HG3	2.02	0.40
11:K:4:LEU:CD1	11:K:15:ALA:HB3	2.52	0.40
9:W:120:ILE:HD12	9:W:136:ILE:HG12	2.02	0.40
16:R:318:HOH:O	12:Z:79:HIS:HE1	2.04	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	248/250 (99%)	239 (96%)	7 (3%)	2 (1%)	22	62
1	O	248/250 (99%)	239 (96%)	7 (3%)	2 (1%)	22	62
2	B	242/258 (94%)	230 (95%)	10 (4%)	2 (1%)	22	62
2	P	242/258 (94%)	231 (96%)	9 (4%)	2 (1%)	22	62
3	C	239/254 (94%)	227 (95%)	9 (4%)	3 (1%)	14	53
3	Q	239/254 (94%)	227 (95%)	9 (4%)	3 (1%)	14	53
4	D	240/260 (92%)	230 (96%)	7 (3%)	3 (1%)	14	53
4	R	240/260 (92%)	230 (96%)	7 (3%)	3 (1%)	14	53
5	E	231/234 (99%)	218 (94%)	12 (5%)	1 (0%)	38	75
5	S	231/234 (99%)	218 (94%)	12 (5%)	1 (0%)	38	75
6	F	242/288 (84%)	232 (96%)	10 (4%)	0	100	100
6	T	242/288 (84%)	232 (96%)	10 (4%)	0	100	100
7	G	241/252 (96%)	232 (96%)	9 (4%)	0	100	100
7	U	241/252 (96%)	231 (96%)	10 (4%)	0	100	100
8	H	220/232 (95%)	212 (96%)	8 (4%)	0	100	100
8	V	220/232 (95%)	210 (96%)	10 (4%)	0	100	100
9	I	202/205 (98%)	196 (97%)	6 (3%)	0	100	100
9	W	202/205 (98%)	196 (97%)	6 (3%)	0	100	100
10	J	196/198 (99%)	187 (95%)	8 (4%)	1 (0%)	32	71
10	X	196/198 (99%)	187 (95%)	8 (4%)	1 (0%)	32	71
11	K	210/212 (99%)	202 (96%)	7 (3%)	1 (0%)	32	71
11	Y	210/212 (99%)	201 (96%)	8 (4%)	1 (0%)	32	71

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
12	L	220/222 (99%)	210 (96%)	10 (4%)	0	100	100
12	Z	220/222 (99%)	210 (96%)	10 (4%)	0	100	100
13	M	231/233 (99%)	221 (96%)	10 (4%)	0	100	100
13	a	231/233 (99%)	221 (96%)	10 (4%)	0	100	100
14	N	194/196 (99%)	187 (96%)	7 (4%)	0	100	100
14	b	194/196 (99%)	186 (96%)	7 (4%)	1 (0%)	32	71
All	All	6312/6588 (96%)	6042 (96%)	243 (4%)	27 (0%)	38	75

All (27) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	C	52	LEU
3	Q	52	LEU
1	A	166	LYS
3	C	203	THR
4	D	122	GLU
5	E	201	ARG
1	O	166	LYS
3	Q	203	THR
4	R	122	GLU
5	S	201	ARG
14	b	2	SER
1	A	2	THR
2	B	221	ASP
4	D	121	GLY
11	K	39	PRO
1	O	2	THR
2	P	221	ASP
4	R	121	GLY
11	Y	39	PRO
2	B	51	VAL
3	C	183	PRO
2	P	51	VAL
3	Q	183	PRO
10	J	9	VAL
10	X	9	VAL
4	D	118	GLY
4	R	118	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.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	209/209 (100%)	206 (99%)	3 (1%)	71	88
1	O	209/209 (100%)	206 (99%)	3 (1%)	71	88
2	B	203/216 (94%)	193 (95%)	10 (5%)	29	66
2	P	203/216 (94%)	193 (95%)	10 (5%)	29	66
3	C	213/226 (94%)	204 (96%)	9 (4%)	34	70
3	Q	213/226 (94%)	204 (96%)	9 (4%)	34	70
4	D	198/215 (92%)	190 (96%)	8 (4%)	36	71
4	R	198/215 (92%)	190 (96%)	8 (4%)	36	71
5	E	192/193 (100%)	183 (95%)	9 (5%)	30	68
5	S	192/193 (100%)	183 (95%)	9 (5%)	30	68
6	F	201/239 (84%)	191 (95%)	10 (5%)	28	65
6	T	201/239 (84%)	191 (95%)	10 (5%)	28	65
7	G	207/210 (99%)	198 (96%)	9 (4%)	33	70
7	U	207/210 (99%)	198 (96%)	9 (4%)	33	70
8	H	181/190 (95%)	177 (98%)	4 (2%)	57	83
8	V	181/190 (95%)	177 (98%)	4 (2%)	57	83
9	I	172/173 (99%)	169 (98%)	3 (2%)	66	86
9	W	172/173 (99%)	169 (98%)	3 (2%)	66	86
10	J	175/175 (100%)	172 (98%)	3 (2%)	66	86
10	X	175/175 (100%)	172 (98%)	3 (2%)	66	86
11	K	169/169 (100%)	161 (95%)	8 (5%)	30	68
11	Y	169/169 (100%)	161 (95%)	8 (5%)	30	68
12	L	185/185 (100%)	180 (97%)	5 (3%)	50	80
12	Z	185/185 (100%)	180 (97%)	5 (3%)	50	80
13	M	199/199 (100%)	191 (96%)	8 (4%)	36	71
13	a	199/199 (100%)	191 (96%)	8 (4%)	36	71

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
14	N	162/162 (100%)	160 (99%)	2 (1%)	75	89
14	b	162/162 (100%)	160 (99%)	2 (1%)	75	89
All	All	5332/5522 (97%)	5150 (97%)	182 (3%)	42	75

All (182) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	30	GLN
1	A	61	LEU
1	A	157	PHE
2	B	55	LEU
2	B	65	LEU
2	B	69	ASN
2	B	119	GLN
2	B	149	THR
2	B	184	LYS
2	B	186	ASP
2	B	191	LEU
2	B	212	PHE
2	B	220	ASN
3	C	4	ARG
3	C	19	GLU
3	C	51	LYS
3	C	61	LYS
3	C	147	GLN
3	C	160	GLN
3	C	169	VAL
3	C	171	GLU
3	C	206	LYS
4	D	20	LEU
4	D	68	CYS
4	D	102	GLU
4	D	124	ARG
4	D	176	LEU
4	D	190	LEU
4	D	214	ILE
4	D	235	LEU
5	E	9	THR
5	E	29	LYS
5	E	116	GLN
5	E	184	ASN

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Mol	Chain	Res	Type
5	E	188	LEU
5	E	198	GLN
5	E	208	ASP
5	E	227	GLU
5	E	231	LYS
6	F	39	ASN
6	F	117	GLN
6	F	123	ASN
6	F	140	ASN
6	F	181	GLU
6	F	186	ARG
6	F	201	GLU
6	F	202	ASP
6	F	203	ASN
6	F	214	TRP
7	G	68	ARG
7	G	83	ASN
7	G	115	LEU
7	G	154	TYR
7	G	166	GLN
7	G	186	ASN
7	G	201	MET
7	G	221	LYS
7	G	235	ARG
8	H	30	ASN
8	H	34	LEU
8	H	43	CYS
8	H	196	ARG
9	I	37	ASN
9	I	171	LEU
9	I	182	TRP
10	J	71	GLU
10	J	78	GLN
10	J	127	GLU
11	K	4	LEU
11	K	9	GLN
11	K	35	ILE
11	K	69	ARG
11	K	73	ARG
11	K	100	MET
11	K	104	TYR
11	K	107	LYS

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Mol	Chain	Res	Type
12	L	3	ASN
12	L	23	LEU
12	L	49	ASN
12	L	109	THR
12	L	150	LEU
13	M	48	ASN
13	M	69	ASP
13	M	70	LEU
13	M	104	ARG
13	M	146	PHE
13	M	161	ARG
13	M	171	GLN
13	M	226	LYS
14	N	119	VAL
14	N	149	GLU
1	O	30	GLN
1	O	61	LEU
1	O	157	PHE
2	P	55	LEU
2	P	65	LEU
2	P	69	ASN
2	P	119	GLN
2	P	149	THR
2	P	184	LYS
2	P	186	ASP
2	P	191	LEU
2	P	212	PHE
2	P	220	ASN
3	Q	4	ARG
3	Q	19	GLU
3	Q	51	LYS
3	Q	61	LYS
3	Q	147	GLN
3	Q	160	GLN
3	Q	169	VAL
3	Q	171	GLU
3	Q	206	LYS
4	R	20	LEU
4	R	68	CYS
4	R	102	GLU
4	R	124	ARG
4	R	176	LEU

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Mol	Chain	Res	Type
4	R	190	LEU
4	R	214	ILE
4	R	235	LEU
5	S	9	THR
5	S	29	LYS
5	S	116	GLN
5	S	184	ASN
5	S	188	LEU
5	S	198	GLN
5	S	208	ASP
5	S	227	GLU
5	S	231	LYS
6	T	39	ASN
6	T	117	GLN
6	T	123	ASN
6	T	140	ASN
6	T	181	GLU
6	T	186	ARG
6	T	201	GLU
6	T	202	ASP
6	T	203	ASN
6	T	214	TRP
7	U	68	ARG
7	U	83	ASN
7	U	115	LEU
7	U	154	TYR
7	U	166	GLN
7	U	186	ASN
7	U	201	MET
7	U	221	LYS
7	U	235	ARG
8	V	30	ASN
8	V	34	LEU
8	V	43	CYS
8	V	196	ARG
9	W	37	ASN
9	W	171	LEU
9	W	182	TRP
10	X	71	GLU
10	X	78	GLN
10	X	127	GLU
11	Y	4	LEU

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Mol	Chain	Res	Type
11	Y	9	GLN
11	Y	35	ILE
11	Y	69	ARG
11	Y	73	ARG
11	Y	100	MET
11	Y	104	TYR
11	Y	107	LYS
12	Z	3	ASN
12	Z	23	LEU
12	Z	49	ASN
12	Z	109	THR
12	Z	150	LEU
13	a	48	ASN
13	a	69	ASP
13	a	70	LEU
13	a	104	ARG
13	a	146	PHE
13	a	161	ARG
13	a	171	GLN
13	a	226	LYS
14	b	119	VAL
14	b	149	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (179) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	30	GLN
1	A	94	HIS
2	B	20	GLN
2	B	69	ASN
2	B	95	GLN
2	B	119	GLN
2	B	123	GLN
2	B	155	ASN
2	B	176	GLN
2	B	220	ASN
3	C	17	GLN
3	C	77	ASN
3	C	116	GLN
3	C	120	GLN
3	C	147	GLN
3	C	160	GLN

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Mol	Chain	Res	Type
3	C	241	GLN
4	D	15	GLN
4	D	91	HIS
4	D	100	ASN
4	D	160	ASN
4	D	210	GLN
4	D	225	ASN
5	E	4	ASN
5	E	68	HIS
5	E	99	ASN
5	E	116	GLN
5	E	120	GLN
5	E	151	ASN
5	E	184	ASN
5	E	198	GLN
5	E	209	ASN
6	F	19	GLN
6	F	39	ASN
6	F	86	ASN
6	F	117	GLN
6	F	123	ASN
6	F	191	GLN
7	G	30	ASN
7	G	83	ASN
7	G	114	ASN
7	G	117	GLN
7	G	121	GLN
7	G	166	GLN
7	G	167	GLN
7	G	175	ASN
7	G	186	ASN
7	G	231	ASN
8	H	30	ASN
8	H	66	HIS
8	H	144	GLN
8	H	165	ASN
8	H	172	ASN
8	H	189	ASN
9	I	37	ASN
9	I	88	GLN
9	I	156	ASN
10	J	37	GLN

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Mol	Chain	Res	Type
10	J	55	GLN
10	J	78	GLN
10	J	118	GLN
10	J	146	HIS
10	J	191	GLN
10	J	198	GLN
11	K	9	GLN
11	K	85	ASN
11	K	176	ASN
11	K	208	ASN
12	L	1	GLN
12	L	3	ASN
12	L	49	ASN
12	L	70	ASN
12	L	76	HIS
12	L	80	ASN
12	L	135	GLN
12	L	158	ASN
12	L	165	ASN
12	L	195	HIS
13	M	18	ASN
13	M	26	ASN
13	M	48	ASN
13	M	102	GLN
13	M	108	ASN
13	M	171	GLN
13	M	179	ASN
13	M	213	GLN
14	N	69	GLN
14	N	145	ASN
14	N	157	HIS
14	N	161	GLN
1	O	30	GLN
1	O	94	HIS
2	P	20	GLN
2	P	69	ASN
2	P	95	GLN
2	P	119	GLN
2	P	123	GLN
2	P	155	ASN
2	P	176	GLN
2	P	220	ASN

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Mol	Chain	Res	Type
3	Q	17	GLN
3	Q	77	ASN
3	Q	116	GLN
3	Q	120	GLN
3	Q	147	GLN
3	Q	160	GLN
3	Q	241	GLN
4	R	15	GLN
4	R	91	HIS
4	R	100	ASN
4	R	160	ASN
4	R	210	GLN
4	R	225	ASN
5	S	4	ASN
5	S	68	HIS
5	S	99	ASN
5	S	116	GLN
5	S	120	GLN
5	S	184	ASN
5	S	198	GLN
5	S	209	ASN
6	T	19	GLN
6	T	39	ASN
6	T	86	ASN
6	T	117	GLN
6	T	123	ASN
6	T	191	GLN
7	U	30	ASN
7	U	83	ASN
7	U	114	ASN
7	U	117	GLN
7	U	121	GLN
7	U	166	GLN
7	U	167	GLN
7	U	175	ASN
7	U	186	ASN
7	U	231	ASN
8	V	30	ASN
8	V	66	HIS
8	V	144	GLN
8	V	165	ASN
8	V	172	ASN

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Mol	Chain	Res	Type
8	V	189	ASN
9	W	37	ASN
9	W	88	GLN
9	W	156	ASN
10	X	37	GLN
10	X	55	GLN
10	X	78	GLN
10	X	86	GLN
10	X	118	GLN
10	X	146	HIS
10	X	191	GLN
10	X	198	GLN
11	Y	9	GLN
11	Y	85	ASN
11	Y	176	ASN
11	Y	208	ASN
12	Z	1	GLN
12	Z	3	ASN
12	Z	49	ASN
12	Z	70	ASN
12	Z	76	HIS
12	Z	80	ASN
12	Z	135	GLN
12	Z	158	ASN
12	Z	165	ASN
12	Z	195	HIS
13	a	18	ASN
13	a	48	ASN
13	a	102	GLN
13	a	108	ASN
13	a	171	GLN
13	a	179	ASN
13	a	213	GLN
14	b	69	GLN
14	b	145	ASN
14	b	157	HIS
14	b	161	GLN

5.3.3 RNA ⓘ

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

6 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
15	04C	H	301	8	44,44,44	1.35	3 (6%)	55,58,58	0.95	5 (9%)
15	04C	K	301	11	44,44,44	1.26	2 (4%)	55,58,58	1.17	4 (7%)
15	04C	N	201	14	44,44,44	1.43	3 (6%)	55,58,58	1.17	2 (3%)
15	04C	V	301	8	44,44,44	1.40	3 (6%)	55,58,58	0.85	3 (5%)
15	04C	Y	301	11	44,44,44	1.36	2 (4%)	55,58,58	1.11	4 (7%)
15	04C	b	201	14	44,44,44	1.49	5 (11%)	55,58,58	1.42	7 (12%)

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
15	04C	H	301	8	-	0/44/52/52	0/3/3/3
15	04C	K	301	11	-	0/44/52/52	0/3/3/3
15	04C	N	201	14	-	0/44/52/52	0/3/3/3
15	04C	V	301	8	-	0/44/52/52	0/3/3/3
15	04C	Y	301	11	-	0/44/52/52	0/3/3/3
15	04C	b	201	14	-	0/44/52/52	0/3/3/3

All (18) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
15	N	201	04C	C2-C1	2.04	1.42	1.38
15	H	301	04C	C9-C8	2.16	1.57	1.53
15	V	301	04C	C9-C8	2.18	1.57	1.53
15	b	201	04C	O45-C44	2.22	1.42	1.37
15	b	201	04C	C9-C8	2.49	1.58	1.53
15	b	201	04C	C7-C8	2.55	1.59	1.53
15	H	301	04C	C10-C9	2.99	1.59	1.53
15	V	301	04C	C10-C9	3.09	1.59	1.53
15	K	301	04C	C10-C9	3.22	1.59	1.53
15	Y	301	04C	C10-C9	3.27	1.59	1.53
15	N	201	04C	C12-C10	3.51	1.56	1.52
15	H	301	04C	C12-C10	3.58	1.56	1.52
15	V	301	04C	C12-C10	3.63	1.57	1.52
15	N	201	04C	C10-C9	3.64	1.60	1.53
15	b	201	04C	C10-C9	3.82	1.60	1.53
15	K	301	04C	C12-C10	3.87	1.57	1.52
15	b	201	04C	C12-C10	4.14	1.57	1.52
15	Y	301	04C	C12-C10	4.41	1.57	1.52

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
15	b	201	04C	C11-C10-C12	-5.66	102.74	109.87
15	N	201	04C	C11-C10-C12	-5.12	103.41	109.87
15	K	301	04C	C11-C10-C12	-4.17	104.62	109.87
15	b	201	04C	C35-C36-N31	-3.97	104.55	110.11
15	Y	301	04C	C7-C8-C9	-3.22	104.26	111.07
15	K	301	04C	C7-C8-C9	-2.83	105.08	111.07
15	V	301	04C	C11-C10-C12	-2.81	106.32	109.87
15	H	301	04C	C11-C10-C12	-2.62	106.57	109.87
15	Y	301	04C	C7-C8-N22	-2.50	107.24	110.17
15	b	201	04C	C33-C32-N31	-2.37	106.78	110.11
15	b	201	04C	C29-C30-N31	-2.35	107.75	113.26
15	H	301	04C	C7-C8-N22	-2.28	107.50	110.17
15	V	301	04C	C32-N31-C36	2.09	113.59	108.87
15	H	301	04C	C35-C36-N31	2.19	113.17	110.11
15	Y	301	04C	C30-N31-C32	2.20	114.63	111.09
15	H	301	04C	C6-C7-C8	2.41	117.73	113.36
15	b	201	04C	C33-O34-C35	2.42	118.09	109.89
15	V	301	04C	C6-C7-C8	2.56	118.01	113.36
15	K	301	04C	C30-N31-C36	2.77	115.54	111.09
15	H	301	04C	C32-N31-C36	2.78	115.17	108.87
15	b	201	04C	C30-N31-C36	3.34	116.45	111.09

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
15	b	201	04C	C30-N31-C32	3.46	116.64	111.09
15	N	201	04C	C6-C7-C8	3.51	119.72	113.36
15	K	301	04C	C6-C7-C8	3.75	120.17	113.36
15	Y	301	04C	C6-C7-C8	4.14	120.88	113.36

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
15	K	301	04C	4	0
15	N	201	04C	4	0
15	Y	301	04C	2	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 ⓘ

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	250/250 (100%)	-0.47	1 (0%) 92 90	46, 72, 111, 148	0
1	O	250/250 (100%)	-0.29	2 (0%) 86 82	59, 80, 119, 157	0
2	B	244/258 (94%)	-0.33	1 (0%) 92 90	54, 80, 132, 166	0
2	P	244/258 (94%)	-0.25	3 (1%) 79 75	55, 80, 123, 174	0
3	C	241/254 (94%)	-0.26	2 (0%) 86 82	50, 79, 133, 170	0
3	Q	241/254 (94%)	-0.03	5 (2%) 64 60	62, 94, 160, 208	0
4	D	242/260 (93%)	-0.32	6 (2%) 58 53	58, 82, 128, 160	0
4	R	242/260 (93%)	-0.27	3 (1%) 79 75	66, 92, 141, 182	0
5	E	233/234 (99%)	-0.13	3 (1%) 77 73	60, 88, 131, 155	0
5	S	233/234 (99%)	-0.02	7 (3%) 51 47	67, 98, 146, 184	0
6	F	244/288 (84%)	-0.36	2 (0%) 86 82	61, 83, 132, 154	0
6	T	244/288 (84%)	-0.26	0 100 100	63, 91, 142, 177	0
7	G	243/252 (96%)	-0.42	1 (0%) 92 90	53, 76, 118, 184	0
7	U	243/252 (96%)	-0.41	1 (0%) 92 90	59, 78, 113, 163	0
8	H	222/232 (95%)	-0.48	0 100 100	18, 70, 104, 147	0
8	V	222/232 (95%)	-0.50	0 100 100	10, 72, 105, 148	0
9	I	204/205 (99%)	-0.43	1 (0%) 90 88	43, 65, 96, 117	0
9	W	204/205 (99%)	-0.51	0 100 100	46, 64, 98, 117	0
10	J	198/198 (100%)	-0.45	2 (1%) 82 78	44, 64, 94, 164	0
10	X	198/198 (100%)	-0.43	2 (1%) 82 78	50, 67, 95, 168	0
11	K	212/212 (100%)	-0.40	1 (0%) 90 88	18, 65, 100, 116	0
11	Y	212/212 (100%)	-0.45	1 (0%) 90 88	14, 69, 102, 126	0
12	L	222/222 (100%)	-0.38	0 100 100	48, 69, 104, 134	0
12	Z	222/222 (100%)	-0.30	2 (0%) 84 80	50, 69, 106, 132	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
13	M	233/233 (100%)	-0.35	8 (3%) 46 41	49, 72, 107, 149	0
13	a	233/233 (100%)	-0.38	6 (2%) 56 52	47, 68, 100, 142	0
14	N	196/196 (100%)	-0.49	0 100 100	11, 66, 97, 117	0
14	b	196/196 (100%)	-0.50	0 100 100	6, 66, 96, 115	0
All	All	6368/6588 (96%)	-0.35	60 (0%) 84 80	6, 75, 126, 208	0

All (60) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
10	X	198	GLN	5.9
4	R	121	GLY	5.4
13	M	231	GLN	5.3
4	D	121	GLY	5.2
4	D	118	GLY	5.1
13	M	1	THR	4.8
4	D	119	ALA	4.7
4	D	120	SER	4.6
13	M	230	THR	4.6
10	J	198	GLN	4.4
2	B	220	ASN	4.2
13	M	233	ILE	4.0
2	P	220	ASN	3.9
13	M	227	GLY	3.9
10	J	197	ALA	3.9
5	S	2	ARG	3.6
3	Q	234	ILE	3.5
5	S	1	PHE	3.3
13	M	232	LYS	3.3
2	P	219	ALA	3.2
1	O	250	LEU	3.2
5	E	233	ILE	3.2
13	a	230	THR	3.1
13	a	229	GLY	2.9
13	a	1	THR	2.9
5	S	210	LEU	2.8
3	Q	203	THR	2.6
5	S	207	VAL	2.6
13	M	228	TYR	2.6
5	S	173	ARG	2.6
1	O	249	ALA	2.5
5	S	233	ILE	2.5

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Mol	Chain	Res	Type	RSRZ
4	R	120	SER	2.5
10	X	197	ALA	2.4
5	E	1	PHE	2.4
5	S	202	ASP	2.3
4	D	122	GLU	2.3
3	Q	202	GLN	2.3
4	D	124	ARG	2.3
12	Z	174	TYR	2.3
6	F	202	ASP	2.3
3	Q	239	GLN	2.3
3	Q	176	ASN	2.2
12	Z	173	LYS	2.2
2	P	241	THR	2.2
5	E	207	VAL	2.2
3	C	50	LEU	2.2
13	a	231	GLN	2.2
1	A	1	MET	2.2
4	R	118	GLY	2.1
11	K	104	TYR	2.1
7	U	243	ASP	2.1
13	M	229	GLY	2.1
3	C	196	SER	2.1
6	F	177	ASP	2.1
11	Y	104	TYR	2.0
7	G	243	ASP	2.0
13	a	224	ASP	2.0
13	a	233	ILE	2.0
9	I	192	ASP	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 carbohydrates 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. LLDF column lists the quality of electron

density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
15	04C	N	201	42/42	0.95	0.24	3.70	6,10,19,25	0
15	04C	b	201	42/42	0.95	0.23	2.35	4,11,27,40	0
15	04C	V	301	42/42	0.95	0.26	1.75	9,13,54,141	0
15	04C	K	301	42/42	0.94	0.22	1.41	17,31,47,55	0
15	04C	Y	301	42/42	0.95	0.20	0.77	11,31,53,59	0
15	04C	H	301	42/42	0.95	0.20	0.40	13,26,45,51	0

6.5 Other polymers [i](#)

There are no such residues in this entry.