



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 03:32 AM UTC

PDB ID : 6HTP / pdb_00006htp
Title : Yeast 20S proteasome with human beta2c (S171G) in complex with 7
Authors : Huber, E.M.; Groll, M.
Deposited on : 2018-10-04
Resolution : 3.00 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
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
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

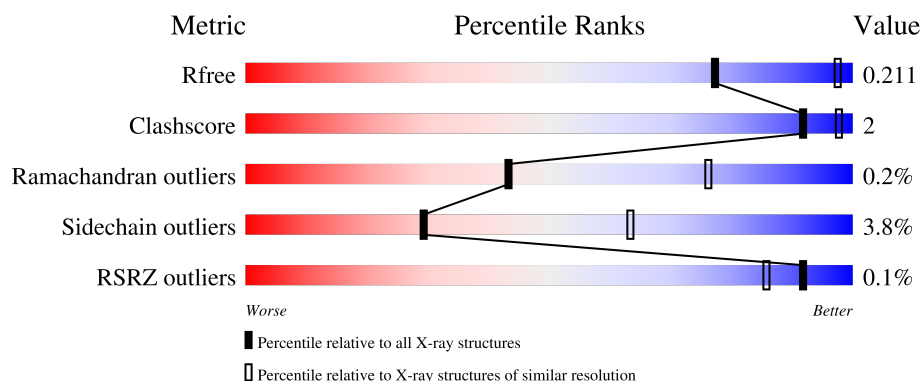
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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	2672 (3.00-3.00)
Clashscore	190562	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)
RSRZ outliers	180081	2671 (3.00-3.00)

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	250	
1	O	250	
2	B	258	
2	P	258	
3	C	254	

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Mol	Chain	Length	Quality of chain
3	Q	254	 2% 87% 7% • 6%
4	D	260	 84% 6% 10%
4	R	260	 84% 6% 10%
5	E	234	 93% 5% •
5	S	234	 93% 6% •
6	F	288	 80% • 16%
6	T	288	 80% • 16%
7	G	252	 90% 5% •
7	U	252	 89% 6% •
8	H	234	 88% • • 6%
8	V	234	 88% • • 6%
9	I	205	 92% 7%
9	W	205	 94% 5%
10	J	198	 % 91% 7% •
10	X	198	 % 91% 7% •
11	K	212	 82% 17% •
11	Y	212	 83% 16% •
12	L	222	 94% 6%
12	Z	222	 93% 6% •
13	M	246	 88% • 9%
13	a	246	 88% • 9%
14	N	196	 93% 6% •
14	b	196	 94% 5% •

2 Entry composition

There are 19 unique types of molecules in this entry. The entry contains 49342 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 subunit alpha type-2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	240	Total	C	N	O	S	0	0	0
			1842	1171	305	362	4			
1	O	240	Total	C	N	O	S	0	0	0
			1842	1171	305	362	4			

- Molecule 2 is a protein called Proteasome subunit alpha type-3.

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 subunit alpha type-4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	240	Total	C	N	O	S	0	0	0
			1881	1176	329	372	4			
3	Q	240	Total	C	N	O	S	0	0	0
			1881	1176	329	372	4			

- Molecule 4 is a protein called Proteasome subunit alpha type-5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	235	Total	C	N	O	S	0	0	0
			1813	1136	304	366	7			
4	R	235	Total	C	N	O	S	0	0	0
			1813	1136	304	366	7			

- Molecule 5 is a protein called Proteasome subunit alpha type-6.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	E	231	Total	C	N	O	S	0	0	0
			1773	1114	307	348	4			
5	S	231	Total	C	N	O	S	0	0	0
			1773	1114	307	348	4			

- Molecule 6 is a protein called Probable proteasome subunit alpha type-7.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
6	F	243	Total	C	N	O	S	0	0	0
			1892	1203	329	356	4			
6	T	243	Total	C	N	O	S	0	0	0
			1892	1203	329	356	4			

- Molecule 7 is a protein called Proteasome subunit alpha type-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
7	G	241	Total	C	N	O	S	0	0	0
			1907	1214	320	365	8			
7	U	241	Total	C	N	O	S	0	0	0
			1907	1214	320	365	8			

- Molecule 8 is a protein called Proteasome subunit beta type-7.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
8	H	219	Total	C	N	O	S	0	0	0
			1648	1038	282	316	12			
8	V	219	Total	C	N	O	S	0	0	0
			1648	1038	282	316	12			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
H	171	GLY	SER	engineered mutation	UNP Q99436
V	171	GLY	SER	engineered mutation	UNP Q99436

- Molecule 9 is a protein called Proteasome subunit beta type-3.

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 subunit beta type-4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
10	J	195	Total	C	N	O	S	0	0	0
			1561	992	264	299	6			
10	X	195	Total	C	N	O	S	0	0	0
			1561	992	264	299	6			

- Molecule 11 is a protein called Proteasome subunit beta type-5.

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 subunit beta type-6.

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 subunit beta type-7.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
13	M	224	Total	C	N	O	S	0	0	0
			1753	1108	300	338	7			
13	a	224	Total	C	N	O	S	0	0	0
			1753	1108	300	338	7			

- Molecule 14 is a protein called Proteasome subunit beta type-1.

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			

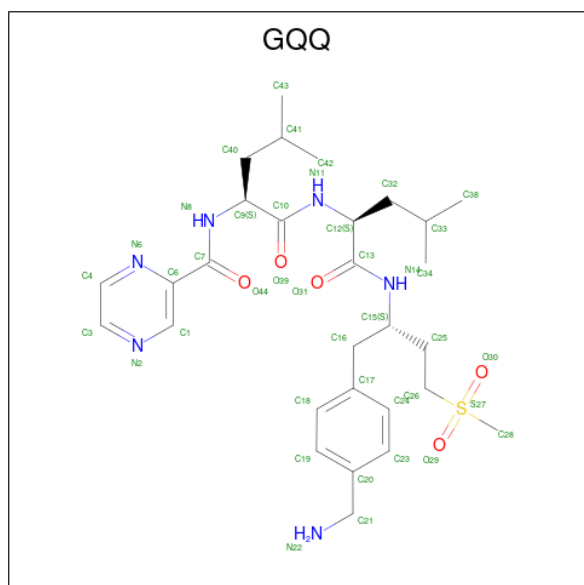
- Molecule 15 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
15	G	1	Total Mg 1 1	0	0
15	I	2	Total Mg 2 2	0	0
15	K	1	Total Mg 1 1	0	0
15	L	1	Total Mg 1 1	0	0
15	N	1	Total Mg 1 1	0	0
15	W	1	Total Mg 1 1	0	0
15	Z	1	Total Mg 1 1	0	0

- Molecule 16 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

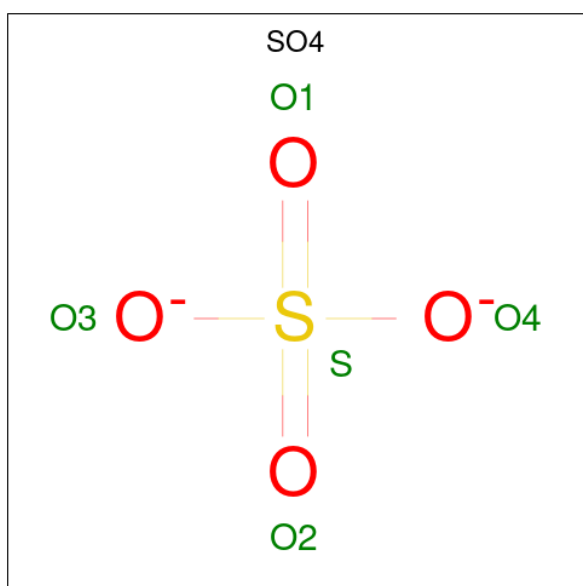
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
16	G	1	Total Cl 1 1	0	0
16	U	1	Total Cl 1 1	0	0

- Molecule 17 is {N}-(2 {S})-1-[(2 {S})-1-[(2 {S})-1-[4-(aminomethyl)phenyl]-4-methylsulfonyl-butan-2-yl]amino]-4-methyl-1-oxidanylidene-pentan-2-yl]amino]-4-methyl-1-oxidanylidene-pentan-2-yl]pyrazine-2-carboxamide (CCD ID: GQQ) (formula: C₂₉H₄₄N₆O₅S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
17	H	1	Total	C	N	O	S	0	0
			41	29	6	5	1		
17	K	1	Total	C	N	O	S	0	0
			41	29	6	5	1		
17	V	1	Total	C	N	O	S	0	0
			41	29	6	5	1		
17	Y	1	Total	C	N	O	S	0	0
			41	29	6	5	1		

- Molecule 18 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
18	N	1	Total	O	S	0	0
			5	4	1		
18	b	1	Total	O	S	0	0
			5	4	1		

- Molecule 19 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
19	A	5	Total	O	0	0
			5	5		
19	B	9	Total	O	0	0
			9	9		
19	C	7	Total	O	0	0
			7	7		
19	D	4	Total	O	0	0
			4	4		

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
19	E	5	Total O 5 5	0	0
19	F	13	Total O 13 13	0	0
19	G	8	Total O 8 8	0	0
19	H	5	Total O 5 5	0	0
19	I	4	Total O 4 4	0	0
19	J	11	Total O 11 11	0	0
19	K	7	Total O 7 7	0	0
19	L	8	Total O 8 8	0	0
19	M	16	Total O 16 16	0	0
19	N	10	Total O 10 10	0	0
19	O	2	Total O 2 2	0	0
19	P	8	Total O 8 8	0	0
19	Q	8	Total O 8 8	0	0
19	R	3	Total O 3 3	0	0
19	S	2	Total O 2 2	0	0
19	T	9	Total O 9 9	0	0
19	U	12	Total O 12 12	0	0
19	V	11	Total O 11 11	0	0
19	W	4	Total O 4 4	0	0
19	X	10	Total O 10 10	0	0
19	Y	10	Total O 10 10	0	0

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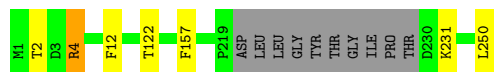
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
19	Z	9	Total 9	O 9	0	0
19	a	14	Total 14	O 14	0	0
19	b	8	Total 8	O 8	0	0

3 Residue-property plots [i](#)

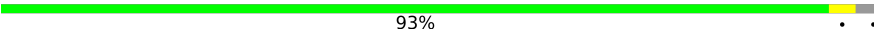
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density.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

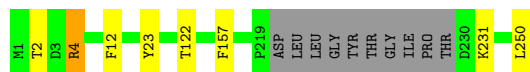
- Molecule 1: Proteasome subunit alpha type-2

Chain A:  93%



- Molecule 1: Proteasome subunit alpha type-2

Chain O:  93%



- Molecule 2: Proteasome subunit alpha type-3

Chain B:  87% 7% 5%



- Molecule 2: Proteasome subunit alpha type-3

Chain P:  88% 6% 5%



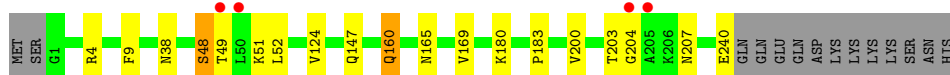
- Molecule 3: Proteasome subunit alpha type-4

Chain C:  86% 7% 6%

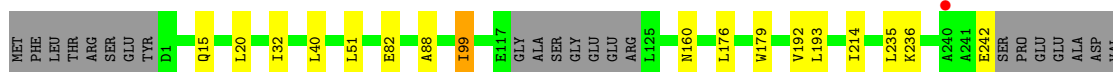
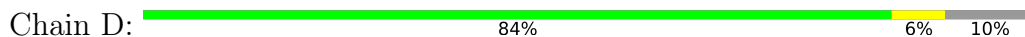


- Molecule 3: Proteasome subunit alpha type-4

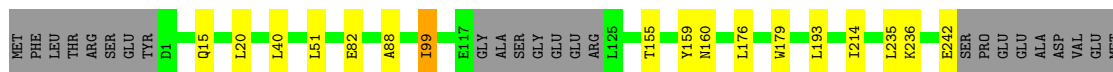
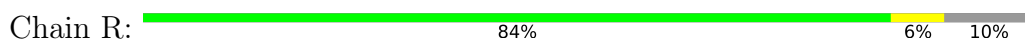
Chain Q:  87% 7% 6%



- Molecule 4: Proteasome subunit alpha type-5



- Molecule 4: Proteasome subunit alpha type-5



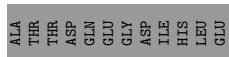
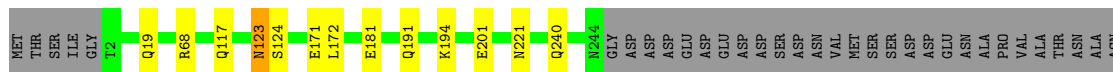
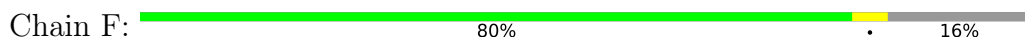
- Molecule 5: Proteasome subunit alpha type-6



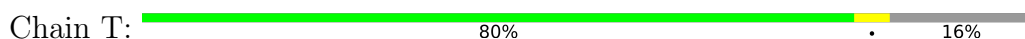
- Molecule 5: Proteasome subunit alpha type-6

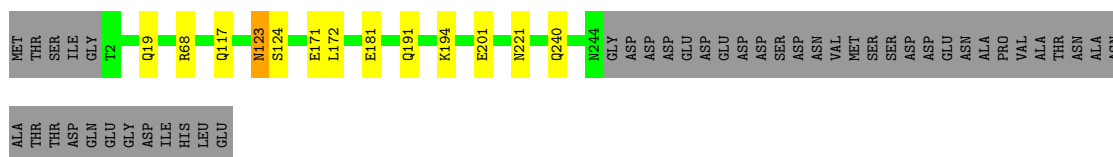


- Molecule 6: Probable proteasome subunit alpha type-7



- Molecule 6: Probable proteasome subunit alpha type-7





- Molecule 7: Proteasome subunit alpha type-1

Chain G: 90% 5% . .



- Molecule 7: Proteasome subunit alpha type-1

Chain U: 89% 6% .



- Molecule 8: Proteasome subunit beta type-7

Chain H: 88% . . 6%



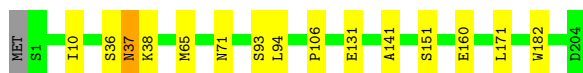
- Molecule 8: Proteasome subunit beta type-7

Chain V: 88% . . 6%



- Molecule 9: Proteasome subunit beta type-3

Chain I: 92% 7%

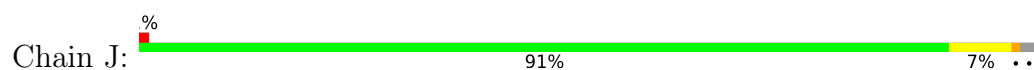


- Molecule 9: Proteasome subunit beta type-3

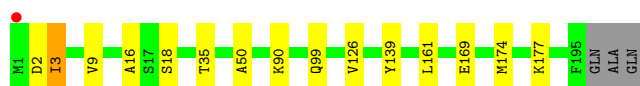
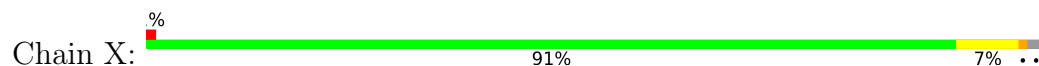
Chain W: 94% 5%



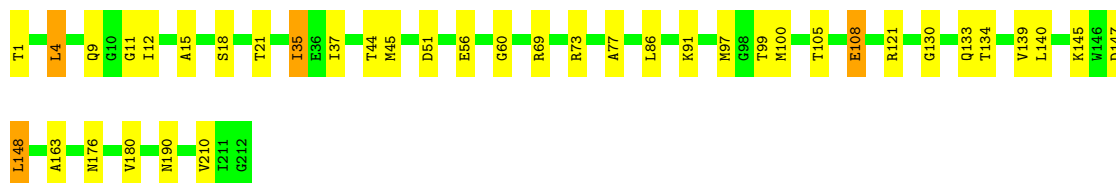
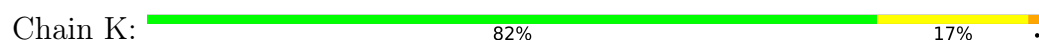
- Molecule 10: Proteasome subunit beta type-4



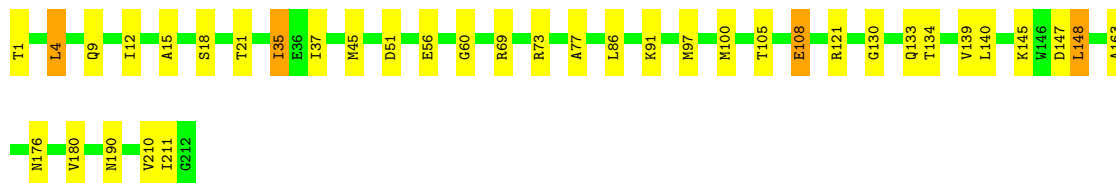
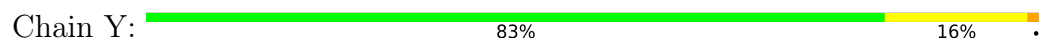
- Molecule 10: Proteasome subunit beta type-4



- Molecule 11: Proteasome subunit beta type-5



- Molecule 11: Proteasome subunit beta type-5



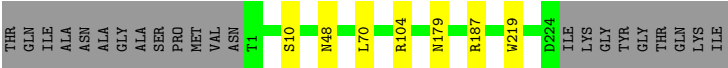
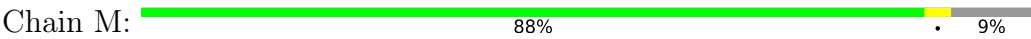
- Molecule 12: Proteasome subunit beta type-6



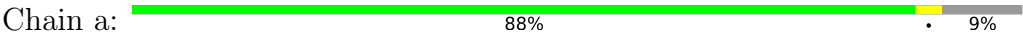
- Molecule 12: Proteasome subunit beta type-6



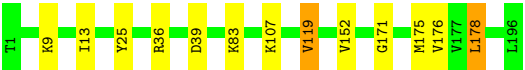
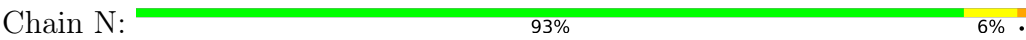
- Molecule 13: Proteasome subunit beta type-7



• Molecule 13: Proteasome subunit beta type-7



• Molecule 14: Proteasome subunit beta type-1



• Molecule 14: Proteasome subunit beta type-1



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	133.91Å 300.42Å 144.04Å 90.00° 112.84° 90.00°	Depositor
Resolution (Å)	15.00 – 3.00 15.00 – 3.00	Depositor EDS
% Data completeness (in resolution range)	95.6 (15.00-3.00) 94.9 (15.00-3.00)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.03 (at 3.01Å)	Xtriage
Refinement program	REFMAC 5.8.0158	Depositor
R, R_{free}	0.177 , 0.206 0.185 , 0.211	Depositor DCC
R_{free} test set	9979 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	69.4	Xtriage
Anisotropy	0.257	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 63.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	49342	wwPDB-VP
Average B, all atoms (Å ²)	83.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.57% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: SO4, MG, CL, GQQ

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	0.53	0/1876	0.88	3/2535 (0.1%)
1	O	0.54	0/1876	0.90	4/2535 (0.2%)
2	B	0.55	0/1934	0.84	0/2618
2	P	0.56	0/1934	0.84	0/2618
3	C	0.55	0/1910	0.83	0/2586
3	Q	0.55	0/1910	0.83	0/2586
4	D	0.52	0/1837	0.81	0/2475
4	R	0.52	0/1837	0.81	0/2475
5	E	0.55	0/1800	0.80	0/2433
5	S	0.53	0/1800	0.80	0/2433
6	F	0.53	0/1932	0.81	0/2609
6	T	0.53	0/1932	0.81	0/2609
7	G	0.52	0/1945	0.85	3/2634 (0.1%)
7	U	0.52	0/1945	0.86	2/2634 (0.1%)
8	H	0.53	0/1675	0.90	3/2267 (0.1%)
8	V	0.53	0/1675	0.88	4/2267 (0.2%)
9	I	0.50	0/1611	0.78	0/2174
9	W	0.51	0/1611	0.79	0/2174
10	J	0.52	0/1589	0.78	0/2142
10	X	0.52	0/1589	0.80	0/2142
11	K	0.49	0/1681	0.82	1/2274 (0.0%)
11	Y	0.49	0/1681	0.82	0/2274
12	L	0.51	0/1795	0.78	0/2420
12	Z	0.51	0/1795	0.77	0/2420
13	M	0.53	0/1783	0.82	0/2420
13	a	0.54	0/1783	0.83	0/2420
14	N	0.51	0/1541	0.78	0/2087
14	b	0.51	0/1541	0.78	0/2087
All	All	0.53	0/49818	0.82	20/67348 (0.0%)

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

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	O	0	1
7	U	0	1
8	V	0	1
All	All	0	4

There are no bond length outliers.

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	U	68	ARG	NE-CZ-NH2	-11.36	108.98	119.20
1	O	4	ARG	NE-CZ-NH1	-11.32	110.18	121.50
7	U	68	ARG	CD-NE-CZ	11.27	140.18	124.40
8	H	143	ARG	NE-CZ-NH1	-10.75	110.75	121.50
8	H	143	ARG	NE-CZ-NH2	10.56	128.71	119.20
1	A	4	ARG	NE-CZ-NH2	-10.44	109.81	119.20
7	G	68	ARG	NE-CZ-NH1	-9.94	111.56	121.50
1	O	4	ARG	NE-CZ-NH2	9.92	128.13	119.20
8	V	143	ARG	NE-CZ-NH2	-9.83	110.36	119.20
7	G	68	ARG	NE-CZ-NH2	9.69	127.92	119.20
1	O	4	ARG	CD-NE-CZ	9.55	137.77	124.40
1	A	4	ARG	CD-NE-CZ	9.41	137.57	124.40
8	V	143	ARG	CD-NE-CZ	9.37	137.52	124.40
8	H	143	ARG	CD-NE-CZ	7.56	134.99	124.40
7	G	68	ARG	CD-NE-CZ	6.96	134.15	124.40
1	A	4	ARG	NE-CZ-NH1	6.12	127.62	121.50
8	V	143	ARG	NE-CZ-NH1	5.63	127.13	121.50
11	K	11	GLY	N-CA-C	5.45	116.88	111.21
1	O	4	ARG	CG-CD-NE	5.22	123.49	112.00
8	V	54	MET	CG-SD-CE	5.22	112.39	100.90

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	4	ARG	Sidechain
1	O	4	ARG	Sidechain
7	U	68	ARG	Sidechain

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Mol	Chain	Res	Type	Group
8	V	143	ARG	Sidechain

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1842	0	1855	1	0
1	O	1842	0	1855	2	0
2	B	1904	0	1904	7	0
2	P	1904	0	1904	6	0
3	C	1881	0	1895	9	0
3	Q	1881	0	1895	8	0
4	D	1813	0	1797	5	0
4	R	1813	0	1797	6	0
5	E	1773	0	1775	8	0
5	S	1773	0	1775	9	0
6	F	1892	0	1883	4	0
6	T	1892	0	1883	4	0
7	G	1907	0	1901	5	0
7	U	1907	0	1901	7	0
8	H	1648	0	1672	5	0
8	V	1648	0	1672	4	0
9	I	1581	0	1574	8	0
9	W	1581	0	1574	5	0
10	J	1561	0	1569	8	0
10	X	1561	0	1569	8	0
11	K	1644	0	1594	24	0
11	Y	1644	0	1594	25	0
12	L	1757	0	1711	7	0
12	Z	1757	0	1711	8	0
13	M	1753	0	1754	1	0
13	a	1753	0	1754	1	0
14	N	1512	0	1481	8	0
14	b	1512	0	1481	7	0
15	G	1	0	0	0	0
15	I	2	0	0	0	0
15	K	1	0	0	0	0
15	L	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
15	N	1	0	0	0	0
15	W	1	0	0	0	0
15	Z	1	0	0	0	0
16	G	1	0	0	0	0
16	U	1	0	0	0	0
17	H	41	0	0	1	0
17	K	41	0	0	1	0
17	V	41	0	0	1	0
17	Y	41	0	0	1	0
18	N	5	0	0	0	0
18	b	5	0	0	0	0
19	A	5	0	0	0	1
19	B	9	0	0	0	0
19	C	7	0	0	0	0
19	D	4	0	0	0	0
19	E	5	0	0	1	0
19	F	13	0	0	0	0
19	G	8	0	0	0	0
19	H	5	0	0	0	0
19	I	4	0	0	0	0
19	J	11	0	0	1	0
19	K	7	0	0	0	0
19	L	8	0	0	0	0
19	M	16	0	0	0	0
19	N	10	0	0	1	0
19	O	2	0	0	0	0
19	P	8	0	0	0	0
19	Q	8	0	0	1	0
19	R	3	0	0	0	0
19	S	2	0	0	0	0
19	T	9	0	0	0	0
19	U	12	0	0	1	1
19	V	11	0	0	0	0
19	W	4	0	0	0	0
19	X	10	0	0	1	0
19	Y	10	0	0	0	0
19	Z	9	0	0	0	0
19	a	14	0	0	0	0
19	b	8	0	0	0	0
All	All	49342	0	48730	159	1

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 2.

All (159) 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:E:92:ASN:HD21	12:L:70:ASN:HD21	1.14	0.90
5:S:92:ASN:HD21	12:Z:70:ASN:HD21	1.14	0.89
3:C:160:GLN:HE21	3:C:160:GLN:HA	1.52	0.75
3:Q:160:GLN:HE21	3:Q:160:GLN:HA	1.53	0.73
11:Y:77:ALA:O	11:Y:121:ARG:NH1	2.23	0.71
11:K:77:ALA:O	11:K:121:ARG:NH1	2.22	0.71
5:E:92:ASN:HD21	12:L:70:ASN:ND2	1.91	0.66
5:E:92:ASN:ND2	12:L:70:ASN:HD21	1.91	0.64
5:S:92:ASN:ND2	12:Z:70:ASN:HD21	1.92	0.64
7:G:68:ARG:NH1	14:N:39:ASP:OD2	2.30	0.64
11:K:51:ASP:HB3	11:K:97:MET:HE2	1.79	0.63
11:Y:51:ASP:HB3	11:Y:97:MET:HE2	1.80	0.63
5:S:92:ASN:HD21	12:Z:70:ASN:ND2	1.91	0.61
14:b:152:VAL:HA	14:b:175:MET:HE1	1.81	0.61
1:O:12:PHE:H	2:P:20:GLN:HE22	1.49	0.60
12:L:31:THR:HG23	12:L:36:ASN:HD21	1.68	0.59
14:N:152:VAL:HA	14:N:175:MET:HE1	1.85	0.59
7:G:23:PHE:O	7:G:26:THR:HB	2.02	0.59
5:S:12:PHE:H	6:T:19:GLN:HE22	1.50	0.59
7:G:99:TYR:O	8:H:78:THR:HB	2.04	0.58
5:E:12:PHE:H	6:F:19:GLN:HE22	1.51	0.58
12:L:31:THR:CG2	12:L:36:ASN:HD21	2.17	0.58
12:Z:31:THR:CG2	12:Z:36:ASN:HD21	2.17	0.58
12:Z:31:THR:HG23	12:Z:36:ASN:HD21	1.69	0.56
11:Y:4:LEU:HD11	11:Y:15:ALA:HB3	1.87	0.56
11:K:35:ILE:HB	11:K:45:MET:HE3	1.88	0.56
11:K:4:LEU:HD23	11:K:140:LEU:HD22	1.88	0.56
8:V:53:ASP:OD1	17:V:301:GQQ:N22	2.40	0.55
7:U:68:ARG:NH1	14:b:39:ASP:OD2	2.39	0.55
11:Y:35:ILE:HB	11:Y:45:MET:HE3	1.88	0.55
11:K:4:LEU:HD11	11:K:15:ALA:HB3	1.88	0.55
6:T:123:ASN:HD22	6:T:124:SER:N	2.06	0.54
8:H:53:ASP:OD1	17:H:301:GQQ:N22	2.40	0.54
11:K:1:THR:HB	17:K:301:GQQ:O30	2.07	0.54
11:Y:4:LEU:HD23	11:Y:140:LEU:HD22	1.90	0.54
11:K:130:GLY:O	11:K:133:GLN:HG2	2.08	0.54
7:U:23:PHE:O	7:U:26:THR:HB	2.07	0.53
6:F:123:ASN:HD22	6:F:124:SER:N	2.07	0.53
1:A:12:PHE:H	2:B:20:GLN:HE22	1.56	0.53
11:Y:130:GLY:O	11:Y:133:GLN:HG2	2.09	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
11:K:210:VAL:HA	9:W:37:ASN:OD1	2.08	0.53
2:B:12:PHE:H	3:C:17:GLN:HE22	1.55	0.52
5:E:87:LEU:HD21	5:E:107:ALA:HB1	1.92	0.52
11:K:4:LEU:CD1	11:K:15:ALA:HB3	2.39	0.52
5:S:87:LEU:HD21	5:S:107:ALA:HB1	1.91	0.52
11:Y:1:THR:HB	17:Y:301:GQQ:O30	2.09	0.52
11:Y:4:LEU:CD1	11:Y:15:ALA:HB3	2.40	0.52
6:F:123:ASN:HD22	6:F:123:ASN:C	2.17	0.52
9:I:37:ASN:OD1	11:Y:210:VAL:HA	2.10	0.52
11:Y:105:THR:OG1	11:Y:108:GLU:HG3	2.11	0.51
11:K:51:ASP:HB3	11:K:97:MET:CE	2.40	0.51
10:X:16:ALA:HB2	10:X:161:LEU:HD21	1.92	0.51
11:Y:145:LYS:HG3	11:Y:148:LEU:HD12	1.93	0.51
6:T:123:ASN:HD22	6:T:123:ASN:C	2.18	0.51
4:R:82:GLU:OE2	11:Y:69:ARG:NH1	2.44	0.51
3:C:48:SER:HB2	3:C:207:ASN:HD21	1.76	0.50
7:U:200:HIS:CE1	19:U:402:HOH:O	2.64	0.50
11:K:145:LYS:HG3	11:K:148:LEU:HD12	1.93	0.50
10:J:16:ALA:HB2	10:J:161:LEU:HD21	1.92	0.50
11:K:176:ASN:ND2	11:K:190:ASN:HD22	2.10	0.49
14:N:13:ILE:HG21	14:N:175:MET:HE2	1.94	0.49
11:K:105:THR:OG1	11:K:108:GLU:HG3	2.11	0.49
14:N:83:LYS:HG3	14:N:119:VAL:CG2	2.43	0.49
10:X:3:ILE:HG23	10:X:18:SER:HB3	1.95	0.49
14:b:83:LYS:HG3	14:b:119:VAL:CG2	2.43	0.48
10:J:3:ILE:HG23	10:J:18:SER:HB3	1.95	0.48
11:Y:51:ASP:HB3	11:Y:97:MET:CE	2.42	0.48
9:I:38:LYS:HG2	11:Y:211:ILE:HD11	1.96	0.48
7:G:167:GLN:HE21	7:G:171:THR:HG23	1.78	0.48
14:N:171:GLY:HA2	13:a:219:TRP:CH2	2.49	0.47
11:Y:176:ASN:ND2	11:Y:190:ASN:HD22	2.13	0.47
3:Q:48:SER:HB2	3:Q:207:ASN:HD21	1.78	0.47
3:C:99:GLU:CD	11:K:121:ARG:HH22	2.23	0.47
12:Z:42:LYS:HD2	12:Z:55:ASN:HD22	1.80	0.47
12:L:42:LYS:HD2	12:L:55:ASN:HD22	1.81	0.46
14:b:13:ILE:HG21	14:b:175:MET:HE2	1.97	0.46
7:U:99:TYR:O	8:V:78:THR:HB	2.16	0.46
7:U:167:GLN:HE21	7:U:171:THR:HG23	1.80	0.45
8:V:132:LEU:HD22	14:b:25:TYR:CE1	2.51	0.45
10:X:50:ALA:O	11:Y:91:LYS:NZ	2.49	0.45
11:Y:97:MET:HE3	11:Y:97:MET:HB2	1.91	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:R:88:ALA:HA	4:R:99:ILE:HG21	1.99	0.45
10:J:139:TYR:CE1	11:Y:134:THR:HG22	2.51	0.45
13:M:219:TRP:CH2	14:b:171:GLY:HA2	2.52	0.45
11:Y:35:ILE:HG21	11:Y:56:GLU:HB3	1.99	0.45
8:H:132:LEU:HD22	14:N:25:TYR:CE1	2.52	0.44
3:Q:165:ASN:HB2	3:Q:200:VAL:HG11	1.99	0.44
11:K:35:ILE:HG21	11:K:56:GLU:HB3	1.99	0.44
3:Q:9:PHE:H	4:R:15:GLN:HE22	1.66	0.44
3:C:165:ASN:HB2	3:C:200:VAL:HG11	1.99	0.44
10:X:174:MET:HB2	19:X:205:HOH:O	2.16	0.44
2:P:30:HIS:O	2:P:50:LYS:NZ	2.46	0.44
2:B:217:LYS:O	2:B:219:ALA:N	2.51	0.44
4:D:88:ALA:HA	4:D:99:ILE:HG21	1.99	0.44
11:K:12:ILE:HB	11:K:180:VAL:HB	2.00	0.44
3:C:48:SER:HB2	3:C:207:ASN:ND2	2.33	0.44
9:W:37:ASN:HB3	9:W:182:TRP:CE3	2.53	0.44
2:B:95:GLN:HE22	9:I:71:ASN:HD22	1.66	0.43
11:K:134:THR:HG22	10:X:139:TYR:CE1	2.52	0.43
3:C:203:THR:HG22	3:C:204:GLY:H	1.83	0.43
8:H:184:ASP:HB3	8:H:186:LEU:CD1	2.48	0.43
2:P:217:LYS:O	2:P:219:ALA:N	2.52	0.43
3:Q:203:THR:HG22	3:Q:204:GLY:H	1.83	0.43
8:V:184:ASP:HB3	8:V:186:LEU:CD1	2.49	0.43
10:J:177:LYS:NZ	10:X:169:GLU:O	2.51	0.43
3:Q:48:SER:HB2	3:Q:207:ASN:ND2	2.34	0.43
7:G:106:ASP:HB3	7:G:146:TYR:CZ	2.54	0.43
7:U:106:ASP:HB3	7:U:146:TYR:CZ	2.54	0.43
5:E:118:ASN:HD22	5:E:118:ASN:N	2.17	0.42
10:J:50:ALA:O	11:K:91:LYS:NZ	2.52	0.42
11:Y:12:ILE:HB	11:Y:180:VAL:HB	2.01	0.42
6:F:191:GLN:HE22	6:F:194:LYS:HE2	1.84	0.42
4:R:155:THR:HG23	5:S:59:GLN:HE22	1.84	0.42
2:B:139:TYR:CD1	2:B:224:VAL:HG21	2.54	0.42
9:I:36:SER:HB2	10:J:126:VAL:HG11	2.01	0.42
2:P:139:TYR:CD1	2:P:224:VAL:HG21	2.54	0.42
5:E:68:HIS:HE1	5:E:102:LEU:O	2.03	0.42
11:Y:37:ILE:HG23	11:Y:60:GLY:HA2	2.01	0.42
11:Y:86:LEU:C	11:Y:86:LEU:HD13	2.45	0.42
10:J:174:MET:HB2	19:J:205:HOH:O	2.19	0.42
9:I:37:ASN:HB3	9:I:182:TRP:CE3	2.54	0.42
5:S:68:HIS:HE1	5:S:102:LEU:O	2.03	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:H:120:ASP:HB3	8:H:122:LEU:HD22	2.02	0.42
2:P:52:THR:HG21	2:P:209:ARG:HG3	2.02	0.42
3:Q:52:LEU:N	19:Q:301:HOH:O	2.39	0.42
11:Y:139:VAL:HG21	11:Y:163:ALA:HB2	2.03	0.41
9:I:94:LEU:HD11	9:I:106:PRO:HG2	2.02	0.41
12:L:8:ASN:HA	12:L:30:ILE:O	2.20	0.41
5:E:147:GLN:HG2	19:E:305:HOH:O	2.20	0.41
11:K:86:LEU:HD13	11:K:86:LEU:C	2.45	0.41
4:R:160:ASN:HB3	4:R:179:TRP:CE2	2.56	0.41
11:Y:139:VAL:HG21	11:Y:163:ALA:CB	2.51	0.41
11:K:139:VAL:HG21	11:K:163:ALA:HB2	2.02	0.41
5:S:118:ASN:N	5:S:118:ASN:HD22	2.19	0.41
9:W:10:ILE:HG21	9:W:141:ALA:HB3	2.02	0.41
11:Y:35:ILE:HB	11:Y:45:MET:CE	2.50	0.41
4:D:82:GLU:OE2	11:K:69:ARG:NH1	2.53	0.41
9:I:10:ILE:HG21	9:I:141:ALA:HB3	2.03	0.41
11:K:44:THR:O	11:K:99:THR:OG1	2.35	0.41
11:K:139:VAL:HG21	11:K:163:ALA:CB	2.51	0.41
14:N:36:ARG:HD2	19:N:302:HOH:O	2.20	0.41
6:T:191:GLN:HE22	6:T:194:LYS:HE2	1.85	0.41
2:B:124:HIS:HB3	3:C:124:VAL:HG12	2.03	0.41
14:b:176:VAL:HG12	14:b:178:LEU:HD13	2.03	0.41
2:B:52:THR:HG21	2:B:209:ARG:HG3	2.02	0.41
3:C:9:PHE:H	4:D:15:GLN:HE22	1.69	0.41
9:W:36:SER:HB2	10:X:126:VAL:HG11	2.03	0.40
10:J:169:GLU:O	10:X:177:LYS:NZ	2.54	0.40
14:N:176:VAL:HG12	14:N:178:LEU:HD13	2.03	0.40
12:Z:3:ASN:HD22	12:Z:4:PRO:HD2	1.86	0.40
9:I:65:MET:HE1	9:I:93:SER:HB3	2.04	0.40
4:R:159:TYR:CE2	5:S:56:SER:HB3	2.56	0.40
4:D:32:ILE:HD12	4:D:192:VAL:HG23	2.02	0.40
4:D:160:ASN:HB3	4:D:179:TRP:CE2	2.57	0.40
11:K:37:ILE:HG23	11:K:60:GLY:HA2	2.02	0.40
12:Z:146:ILE:HG22	12:Z:150:LEU:HD22	2.04	0.40
1:O:23:TYR:CD1	7:U:12:PRO:HA	2.56	0.40
2:P:124:HIS:HB3	3:Q:124:VAL:HG12	2.04	0.40
9:W:94:LEU:HD11	9:W:106:PRO:HG2	2.02	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
19:A:305:HOH:O	19:U:402:HOH:O[2_656]	2.11	0.09

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	236/250 (94%)	226 (96%)	9 (4%)	1 (0%)	30	65
1	O	236/250 (94%)	227 (96%)	8 (3%)	1 (0%)	30	65
2	B	242/258 (94%)	234 (97%)	6 (2%)	2 (1%)	16	50
2	P	242/258 (94%)	234 (97%)	6 (2%)	2 (1%)	16	50
3	C	238/254 (94%)	236 (99%)	1 (0%)	1 (0%)	30	65
3	Q	238/254 (94%)	236 (99%)	1 (0%)	1 (0%)	30	65
4	D	231/260 (89%)	226 (98%)	5 (2%)	0	100	100
4	R	231/260 (89%)	225 (97%)	6 (3%)	0	100	100
5	E	229/234 (98%)	219 (96%)	10 (4%)	0	100	100
5	S	229/234 (98%)	219 (96%)	10 (4%)	0	100	100
6	F	241/288 (84%)	238 (99%)	3 (1%)	0	100	100
6	T	241/288 (84%)	238 (99%)	3 (1%)	0	100	100
7	G	239/252 (95%)	237 (99%)	2 (1%)	0	100	100
7	U	239/252 (95%)	236 (99%)	3 (1%)	0	100	100
8	H	217/234 (93%)	215 (99%)	2 (1%)	0	100	100
8	V	217/234 (93%)	215 (99%)	2 (1%)	0	100	100
9	I	202/205 (98%)	193 (96%)	9 (4%)	0	100	100
9	W	202/205 (98%)	193 (96%)	9 (4%)	0	100	100
10	J	193/198 (98%)	189 (98%)	4 (2%)	0	100	100
10	X	193/198 (98%)	189 (98%)	3 (2%)	1 (0%)	24	60
11	K	210/212 (99%)	199 (95%)	9 (4%)	2 (1%)	12	45

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
11	Y	210/212 (99%)	199 (95%)	9 (4%)	2 (1%)	12	45
12	L	220/222 (99%)	212 (96%)	8 (4%)	0	100	100
12	Z	220/222 (99%)	213 (97%)	7 (3%)	0	100	100
13	M	222/246 (90%)	213 (96%)	9 (4%)	0	100	100
13	a	222/246 (90%)	213 (96%)	9 (4%)	0	100	100
14	N	194/196 (99%)	190 (98%)	4 (2%)	0	100	100
14	b	194/196 (99%)	190 (98%)	4 (2%)	0	100	100
All	All	6228/6618 (94%)	6054 (97%)	161 (3%)	13 (0%)	43	76

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	231	LYS
2	B	52	THR
2	B	218	GLY
1	O	231	LYS
2	P	52	THR
2	P	218	GLY
11	K	147	ASP
11	Y	147	ASP
11	K	148	LEU
11	Y	148	LEU
3	Q	183	PRO
3	C	183	PRO
10	X	9	VAL

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	201/209 (96%)	197 (98%)	4 (2%)	48	76
1	O	201/209 (96%)	197 (98%)	4 (2%)	48	76

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	B	203/216 (94%)	194 (96%)	9 (4%)	25	60
2	P	203/216 (94%)	195 (96%)	8 (4%)	28	62
3	C	212/226 (94%)	201 (95%)	11 (5%)	21	55
3	Q	212/226 (94%)	202 (95%)	10 (5%)	23	58
4	D	194/215 (90%)	184 (95%)	10 (5%)	21	55
4	R	194/215 (90%)	184 (95%)	10 (5%)	21	55
5	E	190/193 (98%)	184 (97%)	6 (3%)	34	67
5	S	190/193 (98%)	184 (97%)	6 (3%)	34	67
6	F	201/239 (84%)	192 (96%)	9 (4%)	24	59
6	T	201/239 (84%)	192 (96%)	9 (4%)	24	59
7	G	206/210 (98%)	198 (96%)	8 (4%)	28	62
7	U	206/210 (98%)	198 (96%)	8 (4%)	28	62
8	H	179/194 (92%)	170 (95%)	9 (5%)	22	56
8	V	179/194 (92%)	170 (95%)	9 (5%)	22	56
9	I	172/173 (99%)	167 (97%)	5 (3%)	37	70
9	W	172/173 (99%)	167 (97%)	5 (3%)	37	70
10	J	173/175 (99%)	168 (97%)	5 (3%)	37	70
10	X	173/175 (99%)	168 (97%)	5 (3%)	37	70
11	K	169/169 (100%)	161 (95%)	8 (5%)	23	58
11	Y	169/169 (100%)	161 (95%)	8 (5%)	23	58
12	L	185/185 (100%)	177 (96%)	8 (4%)	26	60
12	Z	185/185 (100%)	177 (96%)	8 (4%)	26	60
13	M	192/208 (92%)	186 (97%)	6 (3%)	35	68
13	a	192/208 (92%)	186 (97%)	6 (3%)	35	68
14	N	162/162 (100%)	158 (98%)	4 (2%)	42	72
14	b	162/162 (100%)	158 (98%)	4 (2%)	42	72
All	All	5278/5548 (95%)	5076 (96%)	202 (4%)	29	63

All (202) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	THR
1	A	122	THR

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Mol	Chain	Res	Type
1	A	157	PHE
1	A	250	LEU
2	B	50	LYS
2	B	51	VAL
2	B	52	THR
2	B	55	LEU
2	B	79	LEU
2	B	102	ASN
2	B	113	ARG
2	B	191	LEU
2	B	238	LEU
3	C	4	ARG
3	C	38	ASN
3	C	48	SER
3	C	49	THR
3	C	51	LYS
3	C	52	LEU
3	C	147	GLN
3	C	160	GLN
3	C	169	VAL
3	C	180	LYS
3	C	240	GLU
4	D	20	LEU
4	D	40	LEU
4	D	51	LEU
4	D	99	ILE
4	D	176	LEU
4	D	193	LEU
4	D	214	ILE
4	D	235	LEU
4	D	236	LYS
4	D	242	GLU
5	E	9	THR
5	E	29	LYS
5	E	55	LEU
5	E	71	LEU
5	E	118	ASN
5	E	188	LEU
6	F	68	ARG
6	F	117	GLN
6	F	123	ASN
6	F	171	GLU

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Mol	Chain	Res	Type
6	F	172	LEU
6	F	181	GLU
6	F	201	GLU
6	F	221	ASN
6	F	240	GLN
7	G	13	GLU
7	G	26	THR
7	G	83	ASN
7	G	115	LEU
7	G	122	ARG
7	G	125	MET
7	G	235	ARG
7	G	236	LEU
8	H	14	LEU
8	H	24	MET
8	H	52	THR
8	H	78	THR
8	H	98	LEU
8	H	122	LEU
8	H	132	LEU
8	H	143	ARG
8	H	201	ARG
9	I	37	ASN
9	I	131	GLU
9	I	151	SER
9	I	160	GLU
9	I	171	LEU
10	J	2	ASP
10	J	3	ILE
10	J	35	THR
10	J	90	LYS
10	J	99	GLN
11	K	4	LEU
11	K	9	GLN
11	K	18	SER
11	K	21	THR
11	K	35	ILE
11	K	73	ARG
11	K	100	MET
11	K	108	GLU
12	L	23	LEU
12	L	31	THR

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Mol	Chain	Res	Type
12	L	49	ASN
12	L	108	HIS
12	L	130	SER
12	L	136	CYS
12	L	150	LEU
12	L	161	GLU
13	M	10	SER
13	M	48	ASN
13	M	70	LEU
13	M	104	ARG
13	M	179	ASN
13	M	187	ARG
14	N	9	LYS
14	N	107	LYS
14	N	119	VAL
14	N	178	LEU
1	O	2	THR
1	O	122	THR
1	O	157	PHE
1	O	250	LEU
2	P	50	LYS
2	P	52	THR
2	P	55	LEU
2	P	79	LEU
2	P	102	ASN
2	P	113	ARG
2	P	191	LEU
2	P	238	LEU
3	Q	4	ARG
3	Q	38	ASN
3	Q	48	SER
3	Q	49	THR
3	Q	51	LYS
3	Q	147	GLN
3	Q	160	GLN
3	Q	169	VAL
3	Q	180	LYS
3	Q	240	GLU
4	R	20	LEU
4	R	40	LEU
4	R	51	LEU
4	R	99	ILE

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Mol	Chain	Res	Type
4	R	176	LEU
4	R	193	LEU
4	R	214	ILE
4	R	235	LEU
4	R	236	LYS
4	R	242	GLU
5	S	9	THR
5	S	29	LYS
5	S	55	LEU
5	S	71	LEU
5	S	118	ASN
5	S	188	LEU
6	T	68	ARG
6	T	117	GLN
6	T	123	ASN
6	T	171	GLU
6	T	172	LEU
6	T	181	GLU
6	T	201	GLU
6	T	221	ASN
6	T	240	GLN
7	U	13	GLU
7	U	26	THR
7	U	83	ASN
7	U	115	LEU
7	U	122	ARG
7	U	125	MET
7	U	235	ARG
7	U	236	LEU
8	V	14	LEU
8	V	24	MET
8	V	52	THR
8	V	54	MET
8	V	78	THR
8	V	98	LEU
8	V	132	LEU
8	V	143	ARG
8	V	201	ARG
9	W	37	ASN
9	W	131	GLU
9	W	151	SER
9	W	160	GLU

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Mol	Chain	Res	Type
9	W	171	LEU
10	X	2	ASP
10	X	3	ILE
10	X	35	THR
10	X	90	LYS
10	X	99	GLN
11	Y	4	LEU
11	Y	9	GLN
11	Y	18	SER
11	Y	21	THR
11	Y	35	ILE
11	Y	73	ARG
11	Y	100	MET
11	Y	108	GLU
12	Z	23	LEU
12	Z	31	THR
12	Z	49	ASN
12	Z	108	HIS
12	Z	130	SER
12	Z	136	CYS
12	Z	150	LEU
12	Z	161	GLU
13	a	10	SER
13	a	48	ASN
13	a	70	LEU
13	a	104	ARG
13	a	179	ASN
13	a	187	ARG
14	b	9	LYS
14	b	107	LYS
14	b	119	VAL
14	b	178	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (166) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	143	ASN
2	B	20	GLN
2	B	119	GLN
2	B	123	GLN
2	B	155	ASN
2	B	167	ASN

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Mol	Chain	Res	Type
2	B	176	GLN
3	C	17	GLN
3	C	38	ASN
3	C	77	ASN
3	C	116	GLN
3	C	120	GLN
3	C	147	GLN
3	C	160	GLN
3	C	165	ASN
3	C	207	ASN
3	C	233	GLN
4	D	15	GLN
4	D	91	HIS
4	D	106	GLN
4	D	160	ASN
4	D	198	GLN
4	D	225	ASN
5	E	68	HIS
5	E	92	ASN
5	E	99	ASN
5	E	116	GLN
5	E	118	ASN
5	E	120	GLN
5	E	151	ASN
5	E	184	ASN
6	F	19	GLN
6	F	86	ASN
6	F	117	GLN
6	F	123	ASN
6	F	191	GLN
6	F	203	ASN
6	F	240	GLN
7	G	83	ASN
7	G	114	ASN
7	G	117	GLN
7	G	121	GLN
7	G	167	GLN
7	G	175	ASN
7	G	184	HIS
8	H	30	ASN
8	H	80	ASN
8	H	153	ASN

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Mol	Chain	Res	Type
9	I	71	ASN
9	I	88	GLN
10	J	10	GLN
10	J	55	GLN
10	J	63	ASN
10	J	78	GLN
10	J	146	HIS
10	J	147	HIS
11	K	9	GLN
11	K	85	ASN
11	K	143	ASN
11	K	176	ASN
11	K	190	ASN
11	K	208	ASN
11	K	209	ASN
12	L	3	ASN
12	L	29	ASN
12	L	36	ASN
12	L	49	ASN
12	L	55	ASN
12	L	80	ASN
12	L	152	ASN
12	L	155	ASN
12	L	158	ASN
13	M	2	GLN
13	M	18	ASN
13	M	48	ASN
13	M	102	GLN
13	M	179	ASN
13	M	194	ASN
14	N	141	ASN
14	N	161	GLN
1	O	143	ASN
2	P	20	GLN
2	P	95	GLN
2	P	119	GLN
2	P	123	GLN
2	P	146	GLN
2	P	167	ASN
2	P	176	GLN
3	Q	17	GLN
3	Q	38	ASN

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Mol	Chain	Res	Type
3	Q	77	ASN
3	Q	116	GLN
3	Q	120	GLN
3	Q	147	GLN
3	Q	160	GLN
3	Q	165	ASN
3	Q	207	ASN
3	Q	233	GLN
4	R	15	GLN
4	R	91	HIS
4	R	106	GLN
4	R	198	GLN
4	R	207	ASN
4	R	225	ASN
5	S	59	GLN
5	S	68	HIS
5	S	92	ASN
5	S	99	ASN
5	S	116	GLN
5	S	118	ASN
5	S	120	GLN
5	S	151	ASN
5	S	184	ASN
6	T	19	GLN
6	T	86	ASN
6	T	117	GLN
6	T	123	ASN
6	T	191	GLN
6	T	203	ASN
6	T	240	GLN
7	U	83	ASN
7	U	114	ASN
7	U	117	GLN
7	U	121	GLN
7	U	167	GLN
7	U	175	ASN
7	U	184	HIS
8	V	30	ASN
8	V	80	ASN
8	V	116	HIS
8	V	153	ASN
9	W	63	ASN

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Mol	Chain	Res	Type
9	W	71	ASN
9	W	88	GLN
9	W	203	GLN
10	X	10	GLN
10	X	55	GLN
10	X	63	ASN
10	X	78	GLN
10	X	86	GLN
10	X	112	ASN
10	X	146	HIS
10	X	147	HIS
11	Y	9	GLN
11	Y	62	GLN
11	Y	85	ASN
11	Y	143	ASN
11	Y	176	ASN
11	Y	190	ASN
11	Y	209	ASN
12	Z	3	ASN
12	Z	29	ASN
12	Z	36	ASN
12	Z	49	ASN
12	Z	55	ASN
12	Z	80	ASN
12	Z	155	ASN
12	Z	158	ASN
13	a	2	GLN
13	a	18	ASN
13	a	48	ASN
13	a	102	GLN
13	a	108	ASN
13	a	179	ASN
14	b	141	ASN
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 ⓘ

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 16 ligands modelled in this entry, 10 are monoatomic - leaving 6 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
17	GQQ	V	301	8	42,42,42	1.97	10 (23%)	55,57,57	1.50	10 (18%)
17	GQQ	H	301	8	42,42,42	1.92	10 (23%)	55,57,57	1.51	11 (20%)
17	GQQ	K	301	11	42,42,42	2.00	10 (23%)	55,57,57	1.76	11 (20%)
18	SO4	N	202	-	4,4,4	0.40	0	6,6,6	0.21	0
17	GQQ	Y	301	11	42,42,42	2.07	10 (23%)	55,57,57	1.72	12 (21%)
18	SO4	b	201	-	4,4,4	0.40	0	6,6,6	0.19	0

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
17	GQQ	V	301	8	-	3/44/44/44	0/2/2/2
17	GQQ	K	301	11	-	5/44/44/44	0/2/2/2
17	GQQ	Y	301	11	-	5/44/44/44	0/2/2/2
17	GQQ	H	301	8	-	3/44/44/44	0/2/2/2

All (40) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
17	Y	301	GQQ	C16-C17	-5.68	1.37	1.51
17	K	301	GQQ	C16-C17	-5.45	1.38	1.51
17	H	301	GQQ	C16-C17	-5.17	1.39	1.51
17	Y	301	GQQ	O29-S27	5.17	1.54	1.44

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
17	V	301	GQQ	C16-C17	-5.07	1.39	1.51
17	K	301	GQQ	O29-S27	4.75	1.53	1.44
17	Y	301	GQQ	O30-S27	4.63	1.53	1.44
17	H	301	GQQ	O29-S27	4.60	1.53	1.44
17	V	301	GQQ	O29-S27	4.50	1.53	1.44
17	V	301	GQQ	O30-S27	4.39	1.52	1.44
17	V	301	GQQ	C26-S27	4.35	1.83	1.78
17	K	301	GQQ	O30-S27	4.34	1.52	1.44
17	Y	301	GQQ	C1-N2	4.33	1.43	1.34
17	K	301	GQQ	C1-N2	4.15	1.43	1.34
17	H	301	GQQ	C26-S27	4.15	1.83	1.78
17	V	301	GQQ	C6-C7	-4.12	1.39	1.50
17	K	301	GQQ	C4-N6	4.08	1.43	1.34
17	H	301	GQQ	C6-C7	-4.01	1.39	1.50
17	Y	301	GQQ	C4-N6	4.01	1.43	1.34
17	H	301	GQQ	O30-S27	3.91	1.52	1.44
17	Y	301	GQQ	C6-C7	-3.76	1.40	1.50
17	K	301	GQQ	C6-C7	-3.68	1.40	1.50
17	K	301	GQQ	C3-N2	3.57	1.43	1.33
17	V	301	GQQ	C1-N2	3.51	1.41	1.34
17	H	301	GQQ	C1-N2	3.31	1.41	1.34
17	Y	301	GQQ	C3-N2	3.31	1.43	1.33
17	H	301	GQQ	C21-C20	-3.21	1.40	1.51
17	Y	301	GQQ	C25-C26	3.19	1.55	1.52
17	V	301	GQQ	C3-N2	3.15	1.42	1.33
17	H	301	GQQ	C3-N2	3.03	1.42	1.33
17	V	301	GQQ	C21-C20	-2.99	1.40	1.51
17	V	301	GQQ	C4-N6	2.97	1.40	1.34
17	K	301	GQQ	C25-C26	2.82	1.55	1.52
17	H	301	GQQ	C4-N6	2.77	1.40	1.34
17	Y	301	GQQ	C21-C20	-2.73	1.41	1.51
17	K	301	GQQ	C21-C20	-2.71	1.41	1.51
17	Y	301	GQQ	C26-S27	2.64	1.81	1.78
17	H	301	GQQ	C25-C26	2.57	1.55	1.52
17	V	301	GQQ	C25-C26	2.55	1.55	1.52
17	K	301	GQQ	C26-S27	2.52	1.81	1.78

All (44) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
17	H	301	GQQ	O30-S27-O29	-4.48	108.81	117.22
17	Y	301	GQQ	C40-C9-C10	-4.33	100.33	110.59

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
17	K	301	GQQ	C4-N6-C6	4.32	122.50	116.93
17	K	301	GQQ	O30-S27-O29	-4.32	109.12	117.22
17	K	301	GQQ	C40-C9-C10	-4.26	100.50	110.59
17	V	301	GQQ	O30-S27-O29	-4.24	109.26	117.22
17	Y	301	GQQ	O30-S27-O29	-4.13	109.47	117.22
17	Y	301	GQQ	C4-N6-C6	3.73	121.74	116.93
17	K	301	GQQ	C25-C15-N14	-3.47	104.81	110.50
17	H	301	GQQ	O30-S27-C26	3.44	110.81	108.33
17	H	301	GQQ	C4-N6-C6	3.28	121.15	116.93
17	H	301	GQQ	C32-C12-C13	-3.26	102.86	110.59
17	V	301	GQQ	C32-C12-C13	-3.16	103.10	110.59
17	K	301	GQQ	C32-C12-C13	-3.16	103.10	110.59
17	K	301	GQQ	C16-C17-C18	-3.15	115.05	120.90
17	K	301	GQQ	C16-C17-C24	3.10	126.67	120.90
17	Y	301	GQQ	C25-C15-N14	-3.08	105.45	110.50
17	K	301	GQQ	C4-C3-N2	-3.07	118.14	121.96
17	Y	301	GQQ	C16-C17-C18	-3.04	115.25	120.90
17	V	301	GQQ	O30-S27-C26	3.04	110.52	108.33
17	Y	301	GQQ	C16-C17-C24	3.01	126.51	120.90
17	Y	301	GQQ	C4-C3-N2	-2.92	118.33	121.96
17	V	301	GQQ	C16-C17-C18	-2.88	115.54	120.90
17	V	301	GQQ	C4-N6-C6	2.86	120.61	116.93
17	Y	301	GQQ	C32-C12-C13	-2.84	103.86	110.59
17	V	301	GQQ	C12-C13-N14	-2.81	110.63	116.63
17	V	301	GQQ	C16-C17-C24	2.81	126.13	120.90
17	H	301	GQQ	C16-C17-C24	2.73	125.98	120.90
17	Y	301	GQQ	O29-S27-C26	-2.70	106.38	108.33
17	H	301	GQQ	C12-C13-N14	-2.69	110.89	116.63
17	Y	301	GQQ	C40-C9-N8	2.61	116.46	110.58
17	K	301	GQQ	C40-C9-N8	2.61	116.46	110.58
17	V	301	GQQ	C28-S27-C26	2.60	114.93	105.28
17	H	301	GQQ	C16-C17-C18	-2.58	116.09	120.90
17	H	301	GQQ	C28-S27-C26	2.35	114.03	105.28
17	Y	301	GQQ	C28-S27-C26	2.29	113.79	105.28
17	K	301	GQQ	C3-N2-C1	2.21	120.72	116.85
17	H	301	GQQ	C40-C9-N8	-2.18	105.67	110.58
17	K	301	GQQ	C28-S27-C26	2.16	113.31	105.28
17	H	301	GQQ	C6-C1-N2	-2.12	119.19	121.97
17	H	301	GQQ	C3-C4-N6	-2.07	119.27	122.19
17	V	301	GQQ	C40-C9-N8	-2.05	105.95	110.58
17	Y	301	GQQ	C12-C13-N14	-2.01	112.34	116.63
17	V	301	GQQ	C6-C1-N2	-2.00	119.34	121.97

There are no chirality outliers.

All (16) torsion outliers are listed below:

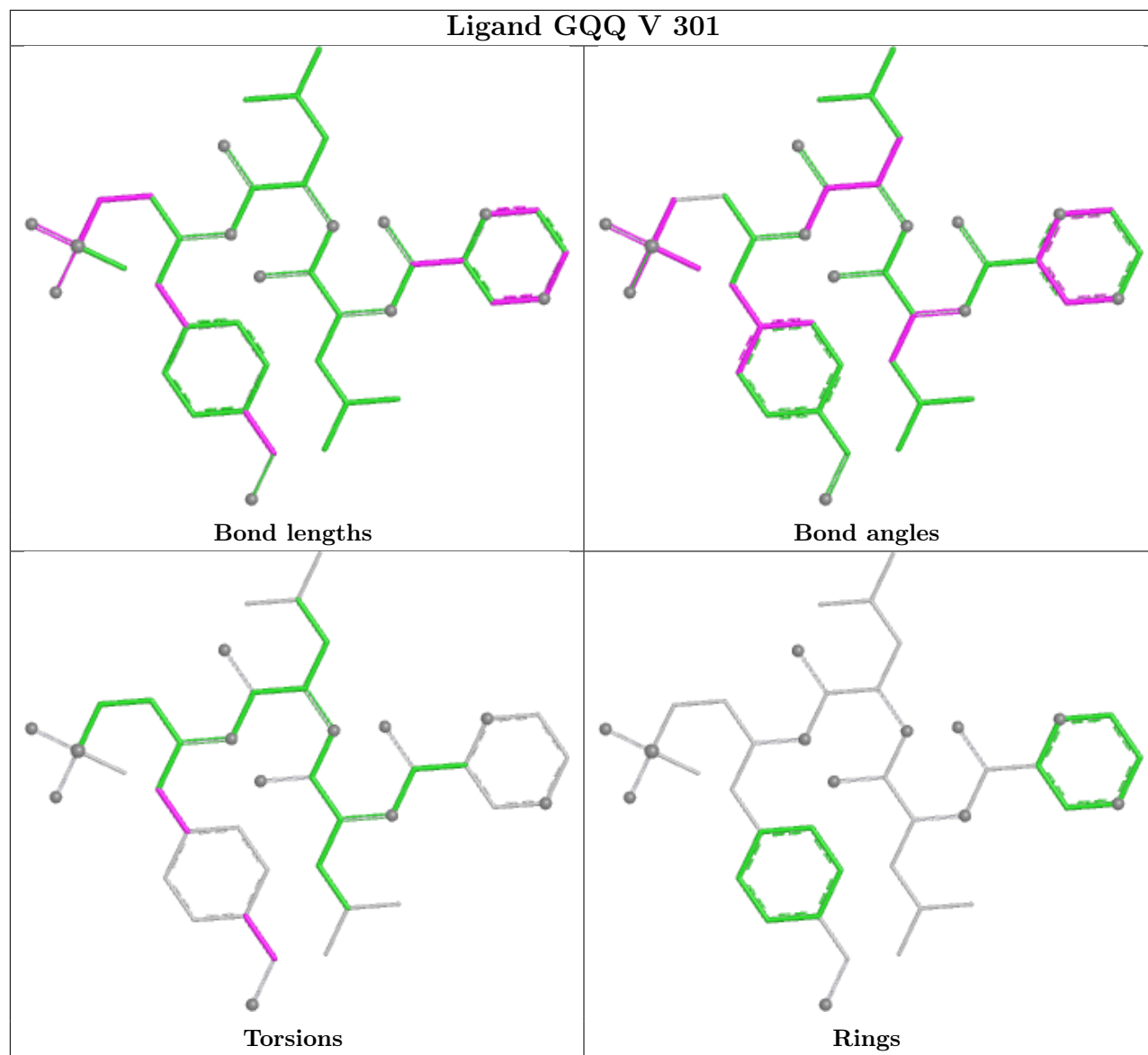
Mol	Chain	Res	Type	Atoms
17	K	301	GQQ	C25-C26-S27-C28
17	K	301	GQQ	C25-C26-S27-O29
17	K	301	GQQ	C25-C26-S27-O30
17	Y	301	GQQ	C25-C26-S27-C28
17	Y	301	GQQ	C25-C26-S27-O29
17	Y	301	GQQ	C25-C26-S27-O30
17	H	301	GQQ	C15-C16-C17-C24
17	V	301	GQQ	C15-C16-C17-C24
17	H	301	GQQ	C15-C16-C17-C18
17	Y	301	GQQ	C15-C16-C17-C24
17	K	301	GQQ	C15-C16-C17-C24
17	V	301	GQQ	C15-C16-C17-C18
17	K	301	GQQ	C15-C16-C17-C18
17	Y	301	GQQ	C15-C16-C17-C18
17	H	301	GQQ	C23-C20-C21-N22
17	V	301	GQQ	C23-C20-C21-N22

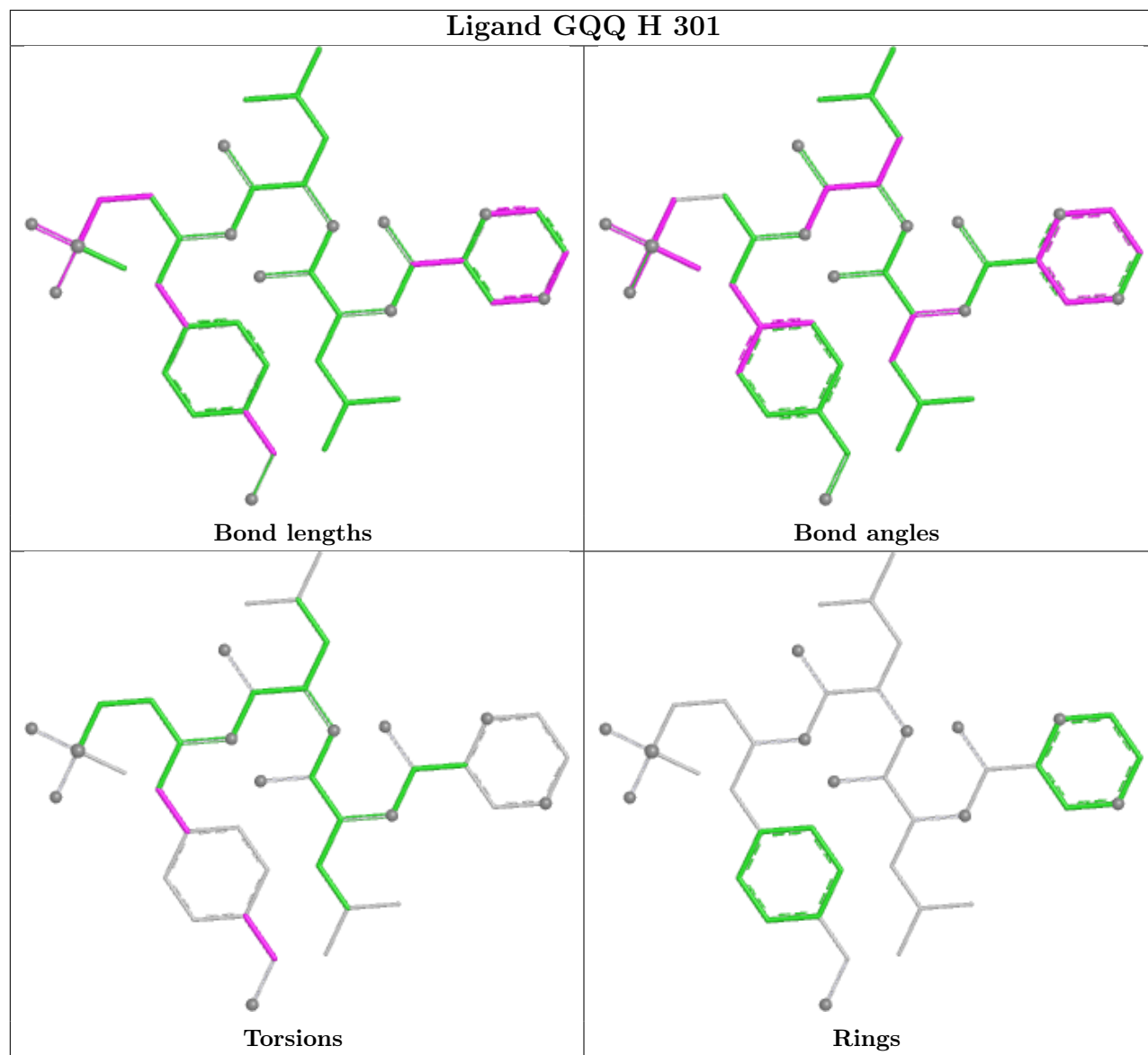
There are no ring outliers.

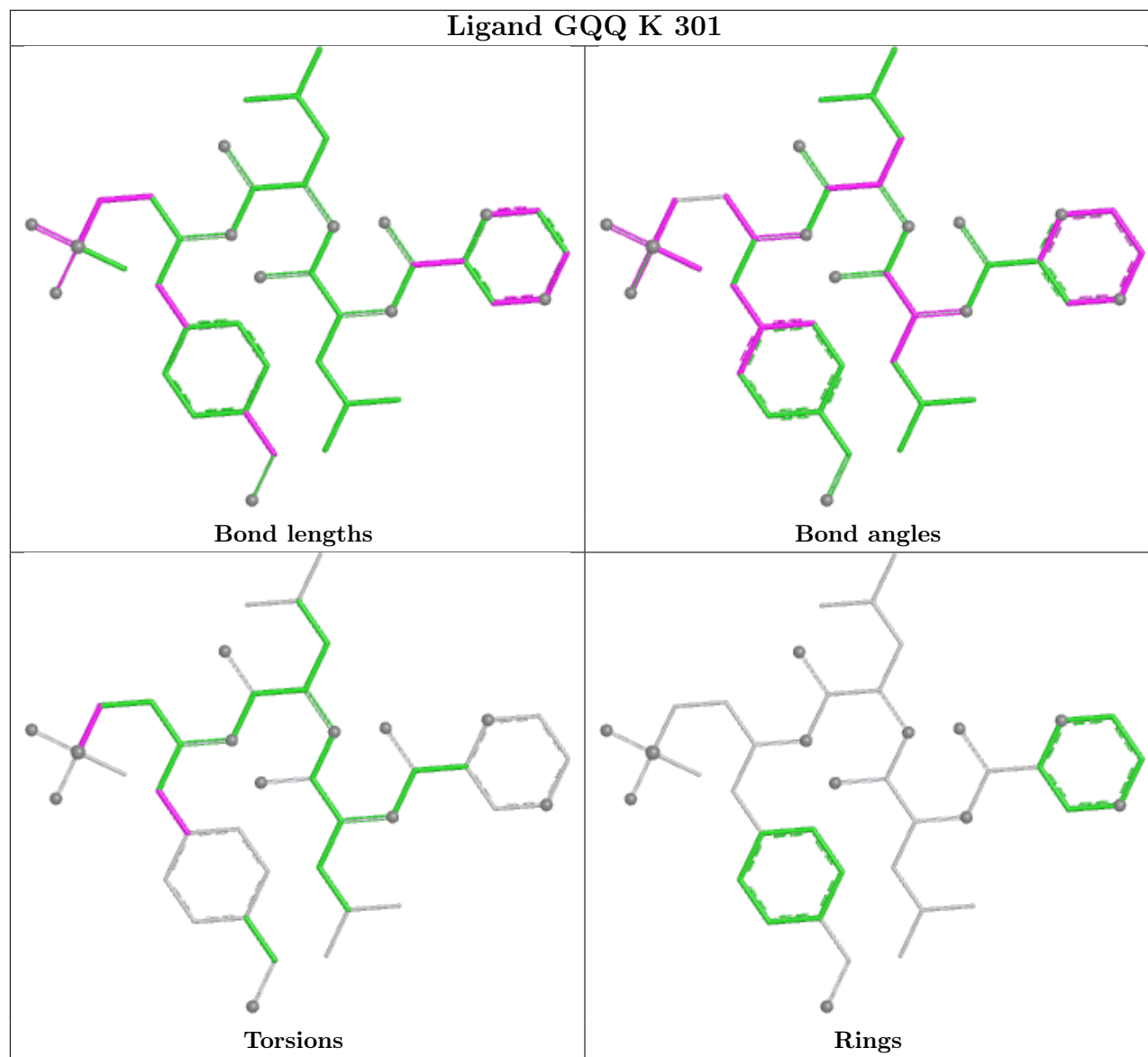
4 monomers are involved in 4 short contacts:

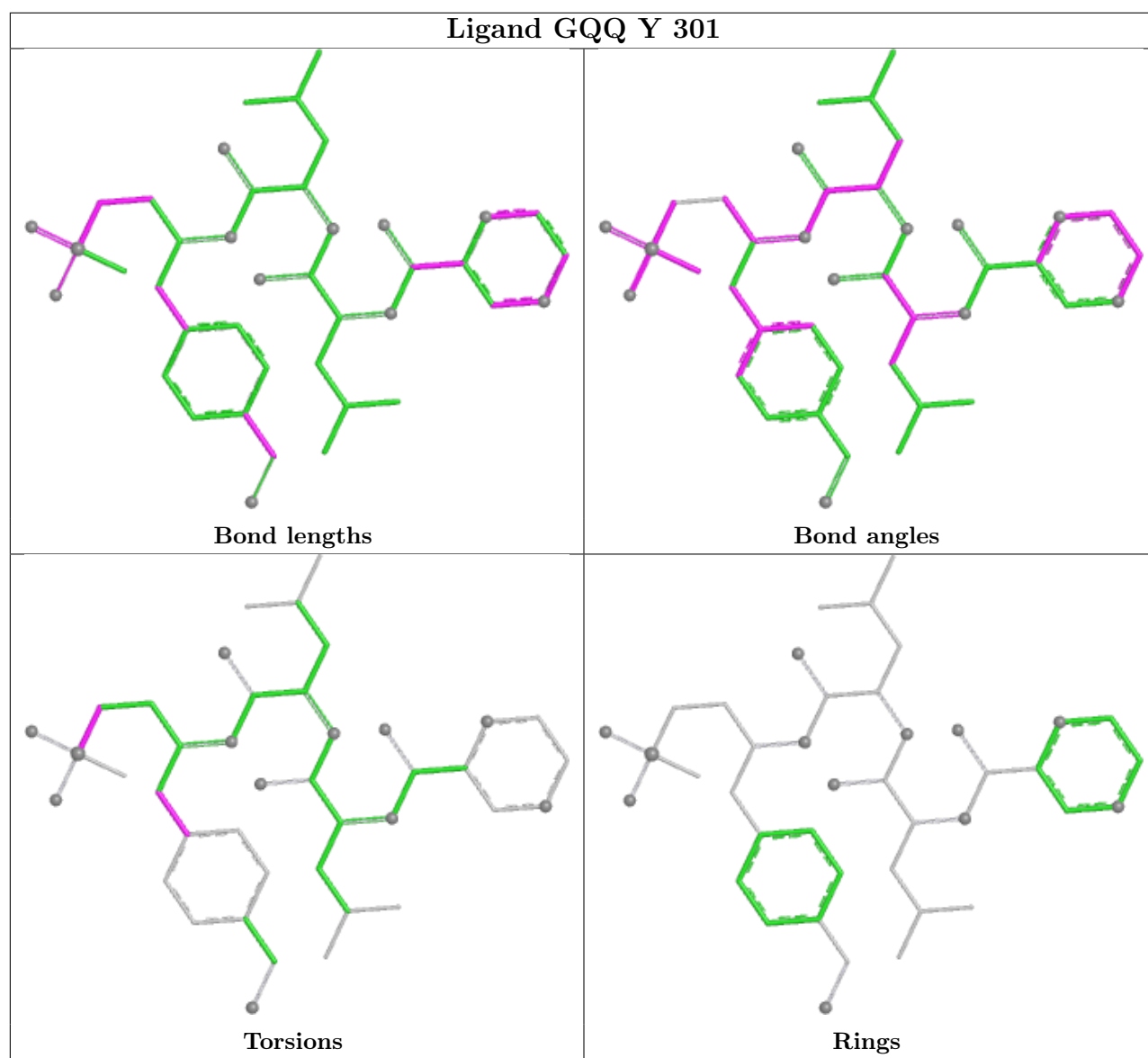
Mol	Chain	Res	Type	Clashes	Symm-Clashes
17	V	301	GQQ	1	0
17	H	301	GQQ	1	0
17	K	301	GQQ	1	0
17	Y	301	GQQ	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.









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	240/250 (96%)	-0.65	0 100 100	57, 77, 119, 154	0
1	O	240/250 (96%)	-0.55	0 100 100	62, 85, 132, 167	0
2	B	244/258 (94%)	-0.56	0 100 100	58, 81, 134, 189	0
2	P	244/258 (94%)	-0.56	0 100 100	60, 82, 130, 183	0
3	C	240/254 (94%)	-0.61	0 100 100	57, 81, 143, 169	0
3	Q	240/254 (94%)	-0.45	4 (1%) 69 45	72, 99, 175, 195	0
4	D	235/260 (90%)	-0.68	1 (0%) 88 76	60, 82, 115, 156	0
4	R	235/260 (90%)	-0.54	0 100 100	68, 91, 132, 165	0
5	E	231/234 (98%)	-0.60	0 100 100	62, 85, 125, 168	0
5	S	231/234 (98%)	-0.45	0 100 100	60, 93, 144, 186	0
6	F	243/288 (84%)	-0.62	0 100 100	53, 78, 131, 163	0
6	T	243/288 (84%)	-0.60	0 100 100	57, 87, 145, 175	0
7	G	241/252 (95%)	-0.71	0 100 100	55, 73, 110, 159	0
7	U	241/252 (95%)	-0.63	0 100 100	60, 77, 114, 155	0
8	H	219/234 (93%)	-0.56	1 (0%) 87 72	52, 71, 127, 147	0
8	V	219/234 (93%)	-0.54	0 100 100	55, 74, 128, 159	0
9	I	204/205 (99%)	-0.79	0 100 100	50, 68, 96, 127	0
9	W	204/205 (99%)	-0.79	0 100 100	51, 68, 98, 120	0
10	J	195/198 (98%)	-0.70	1 (0%) 87 72	51, 71, 99, 140	0
10	X	195/198 (98%)	-0.74	1 (0%) 87 72	51, 72, 98, 157	0
11	K	212/212 (100%)	-0.73	0 100 100	55, 72, 101, 126	0
11	Y	212/212 (100%)	-0.61	0 100 100	53, 77, 110, 136	0
12	L	222/222 (100%)	-0.64	0 100 100	49, 71, 114, 145	0
12	Z	222/222 (100%)	-0.69	0 100 100	51, 71, 113, 150	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
13	M	224/246 (91%)	-0.65	0 100 100	50, 73, 103, 147	0
13	a	224/246 (91%)	-0.70	1 (0%) 88 76	49, 72, 100, 150	0
14	N	196/196 (100%)	-0.77	0 100 100	50, 65, 99, 125	0
14	b	196/196 (100%)	-0.76	0 100 100	50, 67, 100, 126	0
All	All	6292/6618 (95%)	-0.63	9 (0%) 92 86	49, 77, 128, 195	0

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	Q	50	LEU	4.3
13	a	1	THR	3.2
3	Q	205	ALA	2.7
3	Q	204	GLY	2.6
10	J	1	MET	2.4
3	Q	49	THR	2.3
8	H	217	THR	2.3
10	X	1	MET	2.3
4	D	240	ALA	2.2

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

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

6.3 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

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(Å ²)	Q<0.9
15	MG	W	301	1/1	0.89	0.16	83,83,83,83	0
17	GQQ	V	301	41/41	0.90	0.11	47,71,76,85	0
15	MG	I	301	1/1	0.93	0.18	89,89,89,89	0

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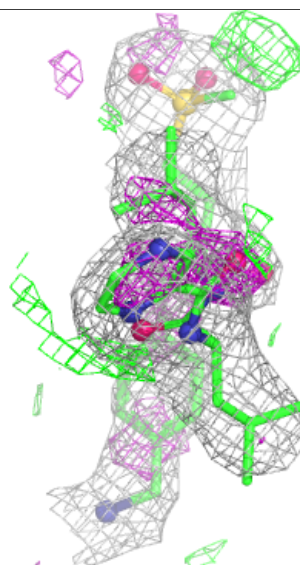
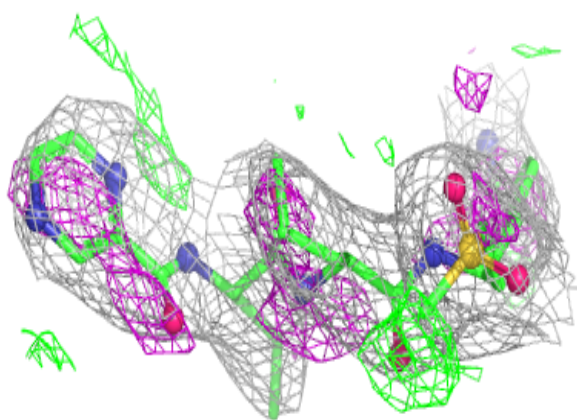
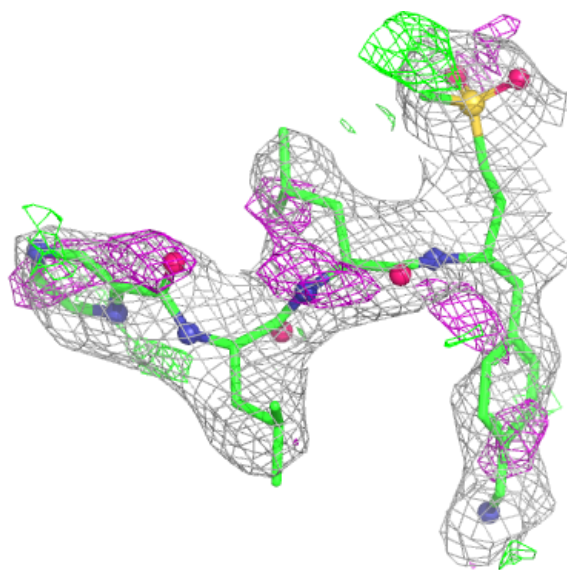
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
17	GQQ	H	301	41/41	0.93	0.10	52,76,86,91	0
15	MG	I	302	1/1	0.93	0.07	100,100,100,100	0
17	GQQ	Y	301	41/41	0.93	0.10	44,59,67,74	0
16	CL	U	301	1/1	0.94	0.15	92,92,92,92	0
15	MG	L	301	1/1	0.94	0.10	91,91,91,91	0
18	SO4	N	202	5/5	0.94	0.17	54,54,64,69	5
17	GQQ	K	301	41/41	0.95	0.07	42,56,61,71	0
16	CL	G	302	1/1	0.96	0.11	79,79,79,79	0
15	MG	Z	301	1/1	0.97	0.07	74,74,74,74	0
18	SO4	b	201	5/5	0.97	0.13	50,58,68,69	5
15	MG	K	302	1/1	0.98	0.04	80,80,80,80	0
15	MG	N	201	1/1	0.98	0.04	55,55,55,55	0
15	MG	G	301	1/1	0.99	0.03	65,65,65,65	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

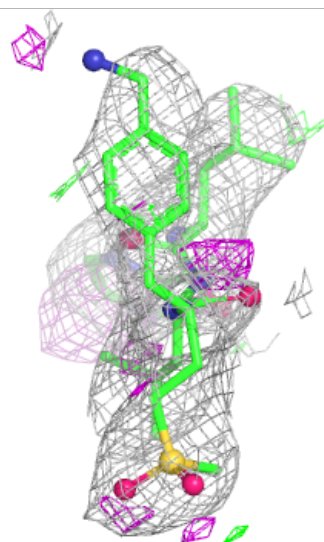
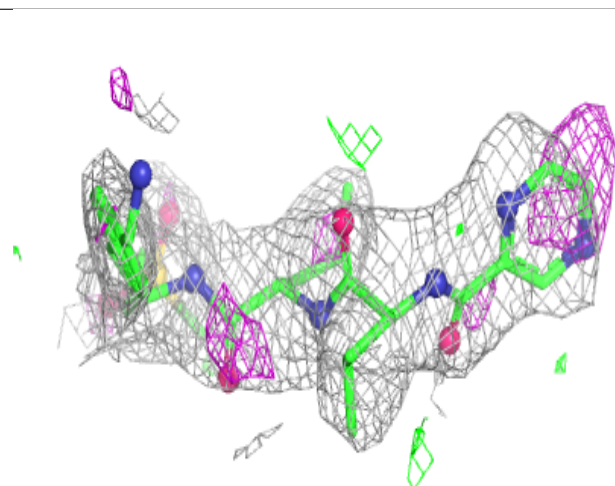
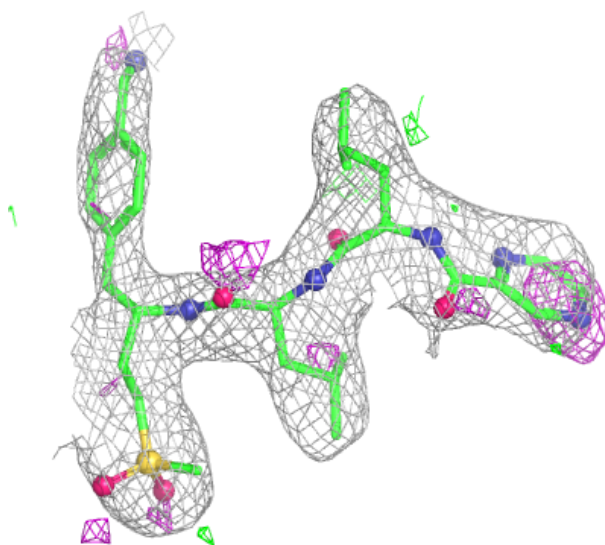
Electron density around GQQ V 301:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



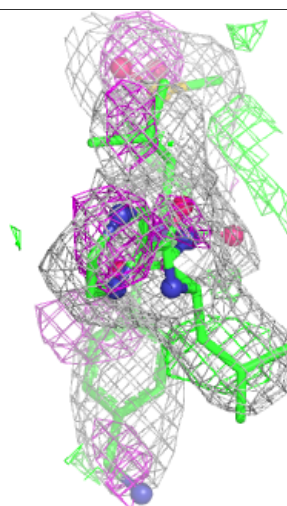
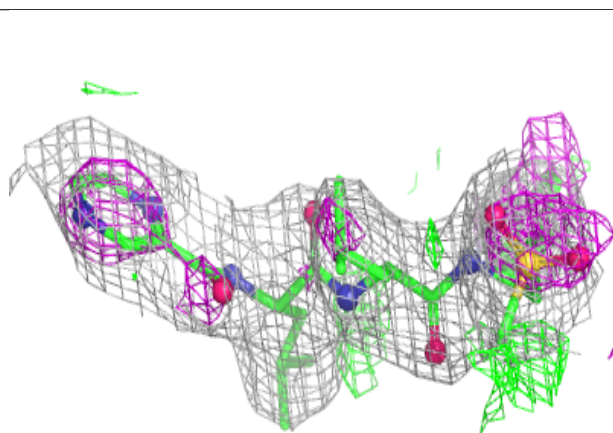
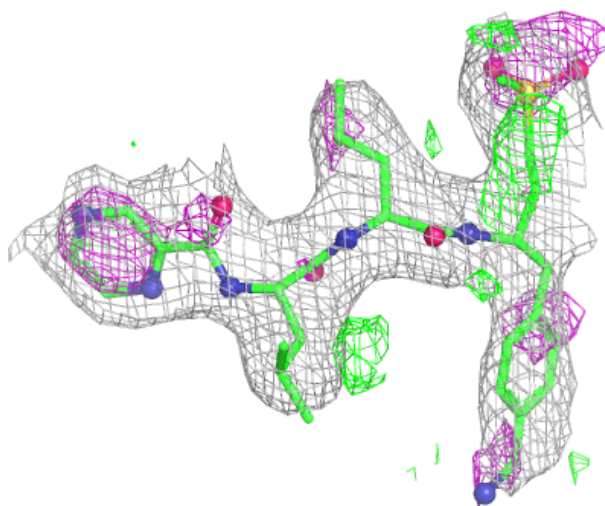
Electron density around GQQ H 301:

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and green (positive)



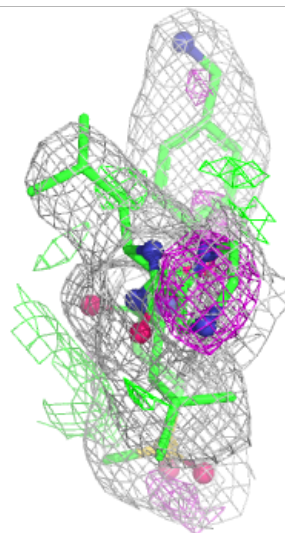
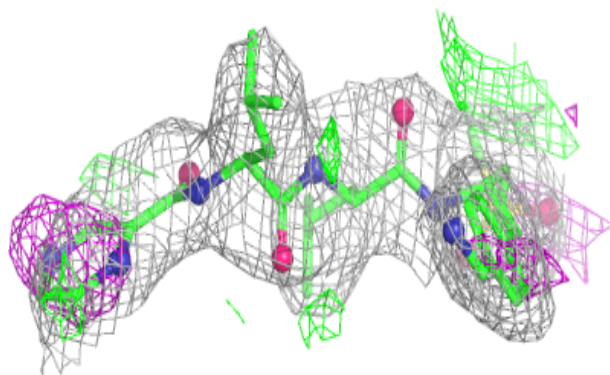
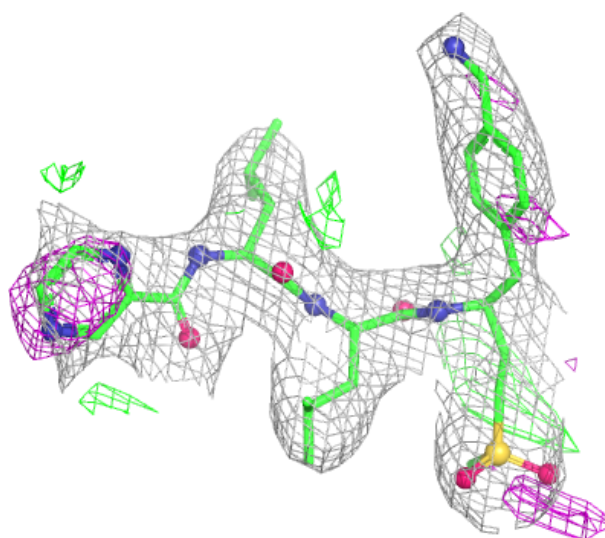
Electron density around GQQ Y 301:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around GQQ K 301:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.