



wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 1, 2016 – 10:43 AM GMT

PDB ID : 3MG6
Title : Structure of yeast 20S open-gate proteasome with Compound 6
Authors : Sintchak, M.D.
Deposited on : 2010-04-05
Resolution : 2.60 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
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) : trunk26865

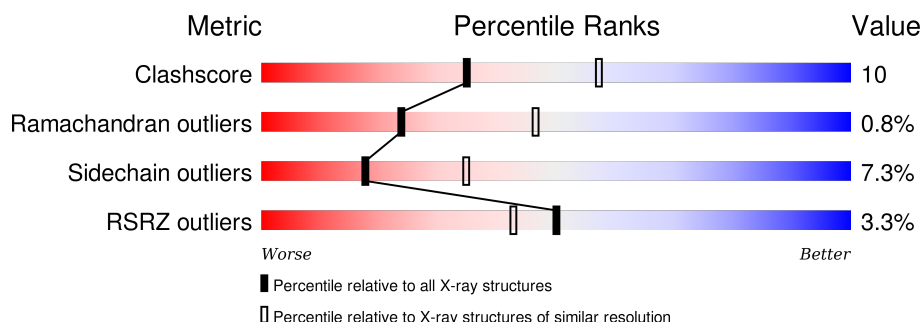
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 Å.

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






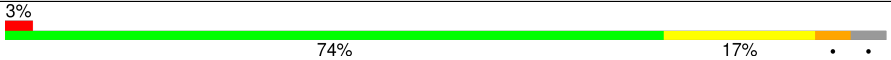
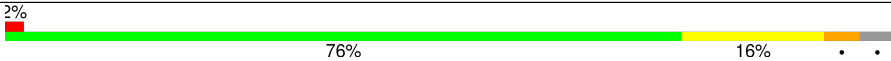
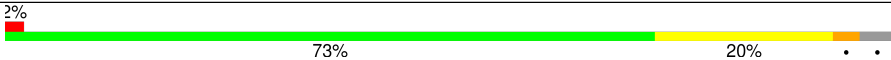
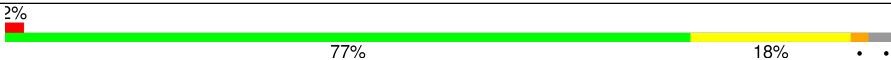
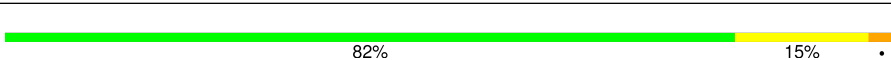
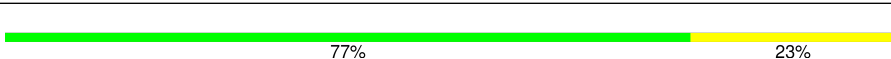
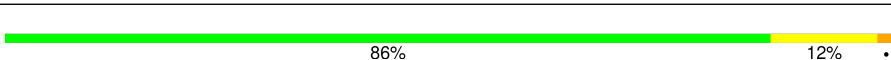



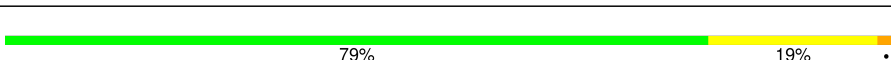
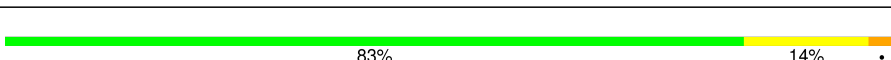
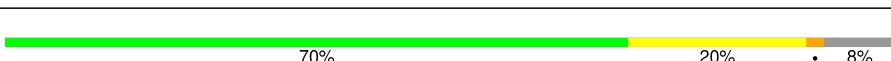
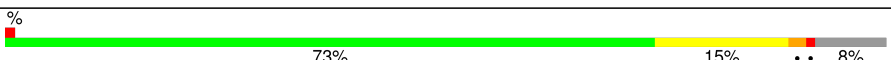
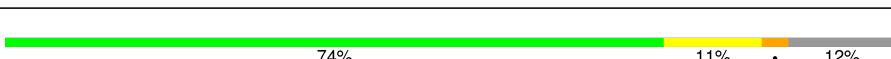
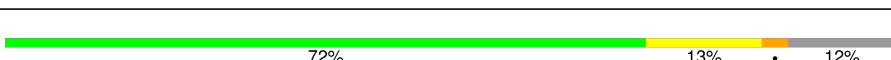
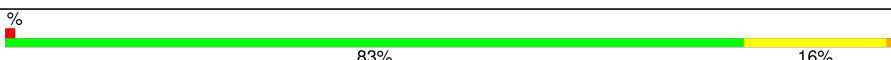
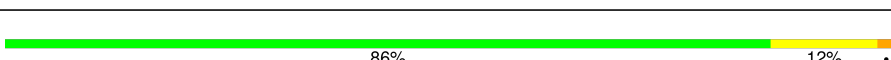
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	102246	2679 (2.60-2.60)
Ramachandran outliers	100387	2635 (2.60-2.60)
Sidechain outliers	100360	2635 (2.60-2.60)
RSRZ outliers	91569	2334 (2.60-2.60)

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>3%</div> <div>84%</div> <div>15%</div> <div>.</div> </div>
1	O	250	<div> <div>4%</div> <div>79%</div> <div>18%</div> <div>.</div> </div>
2	B	245	<div> <div>5%</div> <div>72%</div> <div>21%</div> <div>.</div> </div>
2	P	245	<div> <div>3%</div> <div>72%</div> <div>21%</div> <div>.</div> </div>
3	C	243	<div> <div>9%</div> <div>78%</div> <div>18%</div> <div>.</div> </div>
3	Q	243	<div> <div>18%</div> <div>71%</div> <div>25%</div> <div>.</div> </div>
4	D	250	<div> <div>5%</div> <div>73%</div> <div>21%</div> <div>.</div> </div>

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Mol	Chain	Length	Quality of chain
4	R	250	
5	E	234	
5	S	234	
6	F	248	
6	T	248	
7	G	252	
7	U	252	
8	H	222	
8	V	222	
9	I	205	
9	W	205	
10	J	198	
10	X	198	
11	K	212	
11	Y	212	
12	L	241	
12	Z	241	
13	1	266	
13	M	266	
14	2	196	
14	N	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	MG	F	242	-	-	-	X
15	MG	F	243	-	-	-	X

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
15	MG	I	196	-	-	-	X
15	MG	L	195	-	-	-	X
17	MES	K	214	-	-	-	X

2 Entry composition

There are 18 unique types of molecules in this entry. The entry contains 50276 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	235	Total	C	N	O	S	0	0	0
			1829	1158	303	365	3			
2	P	235	Total	C	N	O	S	0	0	0
			1829	1158	303	365	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
			1891	1181	331	375	4			
3	Q	241	Total	C	N	O	S	0	0	0
			1891	1181	331	375	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
			1862	1162	314	379	7			
4	R	242	Total	C	N	O	S	0	0	0
			1862	1162	314	379	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	237	Total	C	N	O	S	0	0	0
			1848	1175	323	346	4			
6	T	237	Total	C	N	O	S	0	0	0
			1848	1175	323	346	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
			1685	1061	293	324	7			
8	V	222	Total	C	N	O	S	0	0	0
			1685	1061	293	324	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	1	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	2	196	Total	C	N	O	S	0	0	0
			1512	955	250	300	7			

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

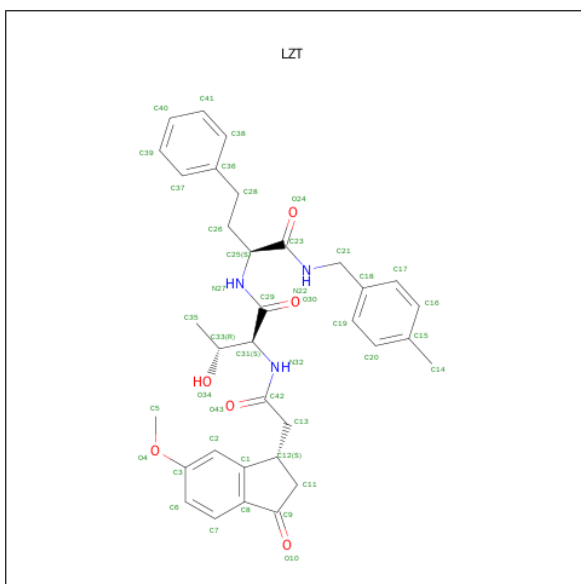
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
15	G	1	Total	Mg	0	0
			1	1		
15	K	1	Total	Mg	0	0
			1	1		

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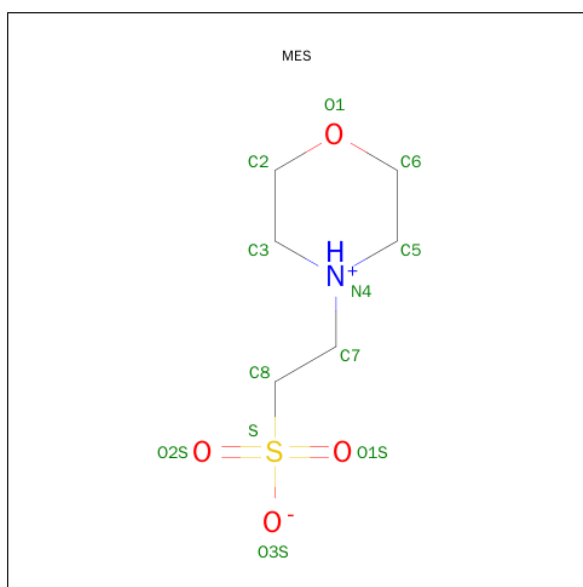
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
15	H	1	Total	Mg	0	0
			1	1		
15	I	2	Total	Mg	0	0
			2	2		
15	N	1	Total	Mg	0	0
			1	1		
15	L	2	Total	Mg	0	0
			2	2		
15	F	2	Total	Mg	0	0
			2	2		

- Molecule 16 is N 2 -{[(1S)-6-METHOXY-3-OXO-2,3-DIHYDRO-1H-INDEN-1-YL]ACETYL}-N-{[(1S)-1-[(4-METHYLBENZYL)CARBAMOYL]-3-PHENYLPROPYL]-L-THREONINAMIDE (three-letter code: LZT) (formula: C₃₄H₃₉N₃O₆).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
16	K	1	Total	C	N	O	0	0
			43	34	3	6		
16	Y	1	Total	C	N	O	0	0
			43	34	3	6		

- Molecule 17 is 2-(N-MORPHOLINO)-ETHANESULFONIC ACID (three-letter code: MES) (formula: C₆H₁₃NO₄S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
17	K	1	Total	C	N	O	S	0	0
			12	6	1	4	1		
17	Y	1	Total	C	N	O	S	0	0
			12	6	1	4	1		

- Molecule 18 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
18	A	27	Total	O	0	0
			27	27		
18	B	28	Total	O	0	0
			28	28		
18	C	27	Total	O	0	0
			27	27		
18	D	24	Total	O	0	0
			24	24		
18	E	21	Total	O	0	0
			21	21		
18	F	38	Total	O	0	0
			38	38		
18	G	44	Total	O	0	0
			44	44		
18	H	30	Total	O	0	0
			30	30		
18	I	37	Total	O	0	0
			37	37		
18	J	33	Total	O	0	0
			33	33		

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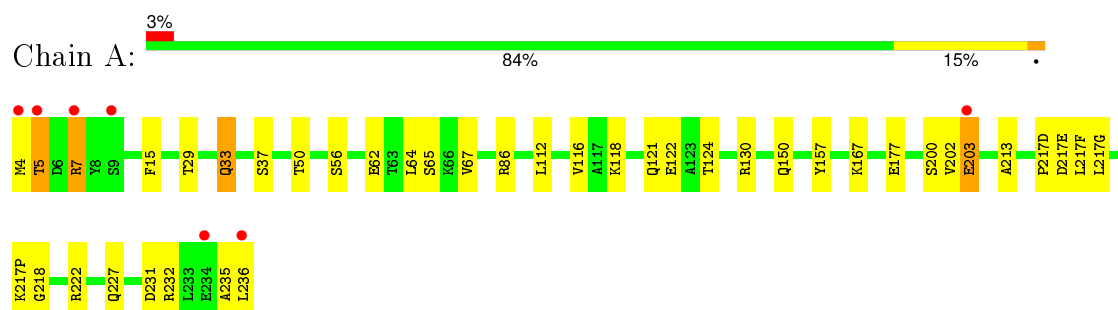
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
18	K	39	Total 39	O 39	0	0
18	L	34	Total 34	O 34	0	0
18	M	35	Total 35	O 35	0	0
18	N	29	Total 29	O 29	0	0
18	O	23	Total 23	O 23	0	0
18	P	17	Total 17	O 17	0	0
18	Q	32	Total 32	O 32	0	0
18	R	21	Total 21	O 21	0	0
18	S	23	Total 23	O 23	0	0
18	T	24	Total 24	O 24	0	0
18	U	34	Total 34	O 34	0	0
18	V	27	Total 27	O 27	0	0
18	W	32	Total 32	O 32	0	0
18	X	38	Total 38	O 38	0	0
18	Y	32	Total 32	O 32	0	0
18	Z	31	Total 31	O 31	0	0
18	1	46	Total 46	O 46	0	0
18	2	32	Total 32	O 32	0	0

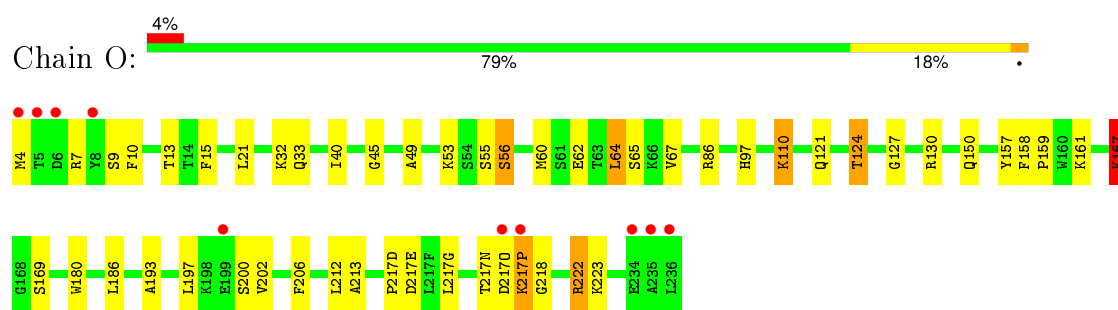
3 Residue-property plots [i](#)

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

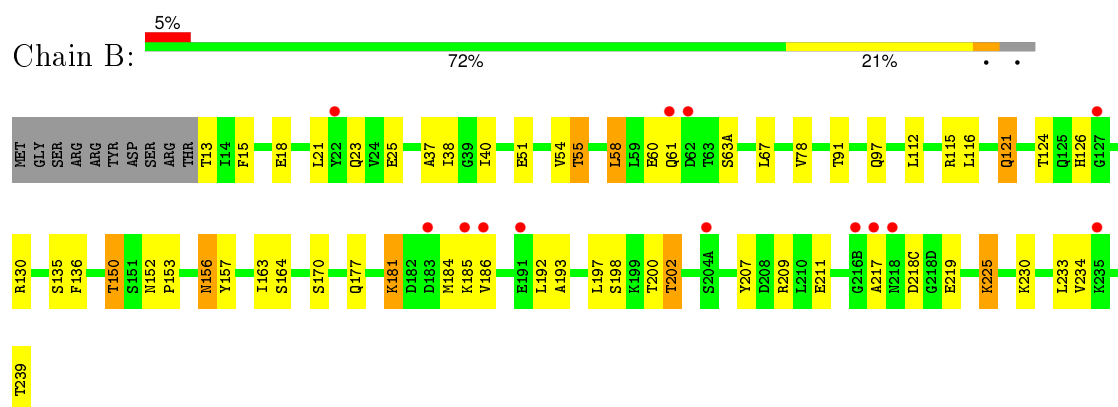
• Molecule 1: Proteasome component Y7



• Molecule 1: Proteasome component Y7

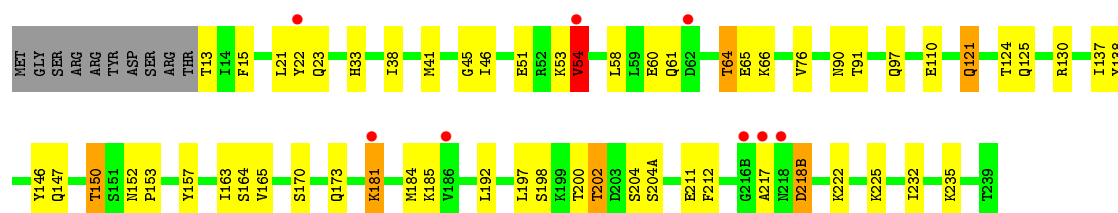


• Molecule 2: Proteasome component Y13

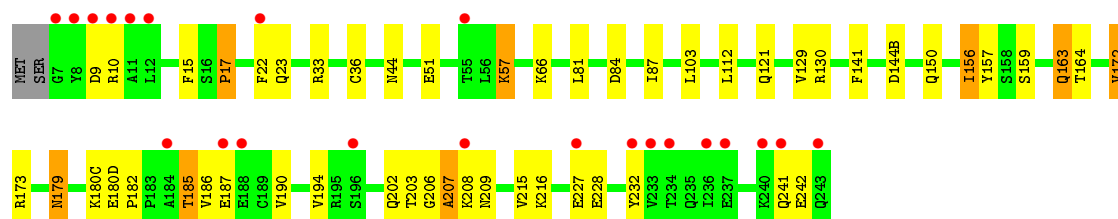
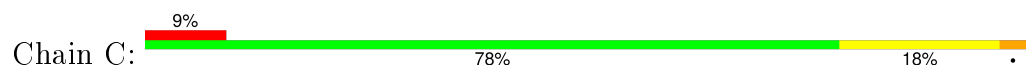


• Molecule 2: Proteasome component Y13

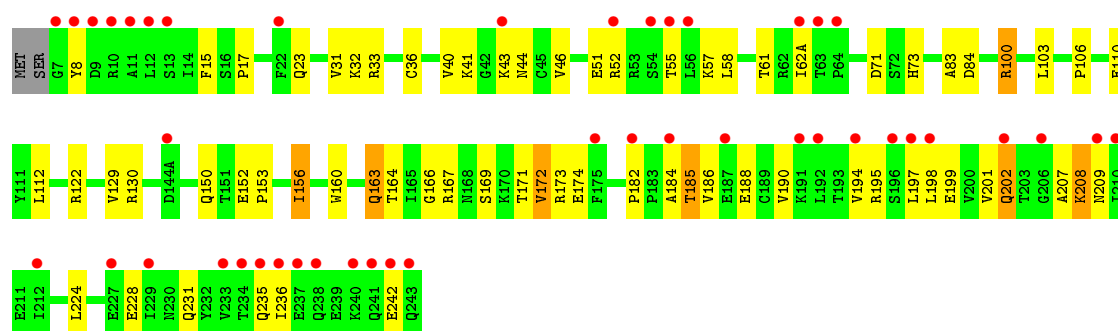




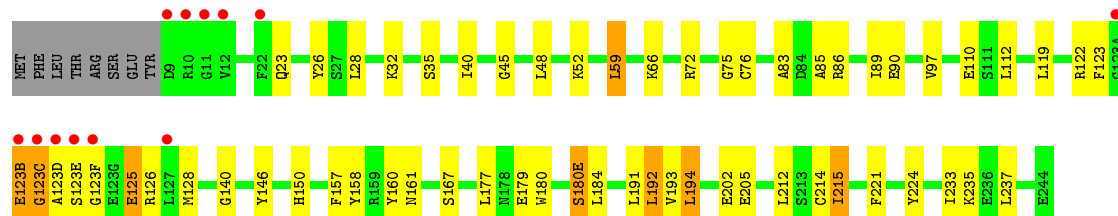
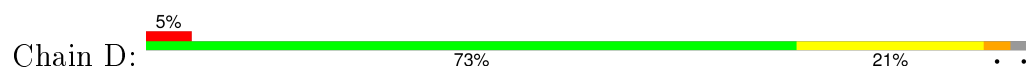
• Molecule 3: Proteasome component PRE6



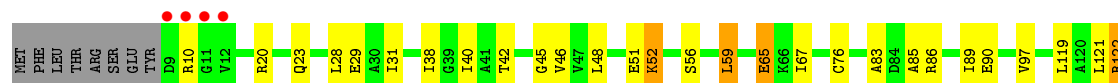
• Molecule 3: Proteasome component PRE6

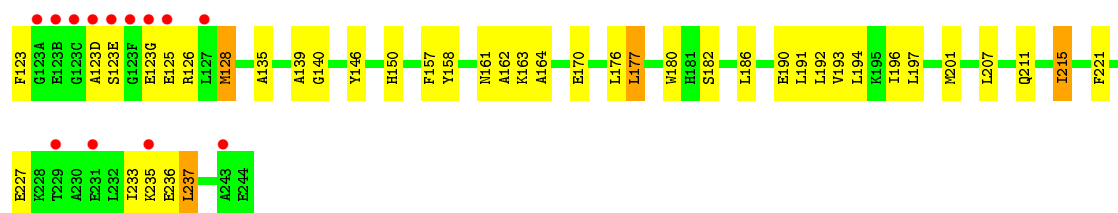


• Molecule 4: Proteasome component PUP2

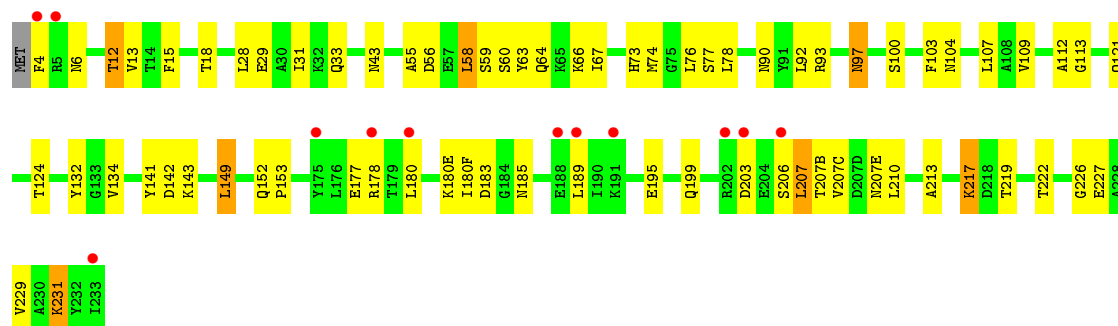


• Molecule 4: Proteasome component PUP2

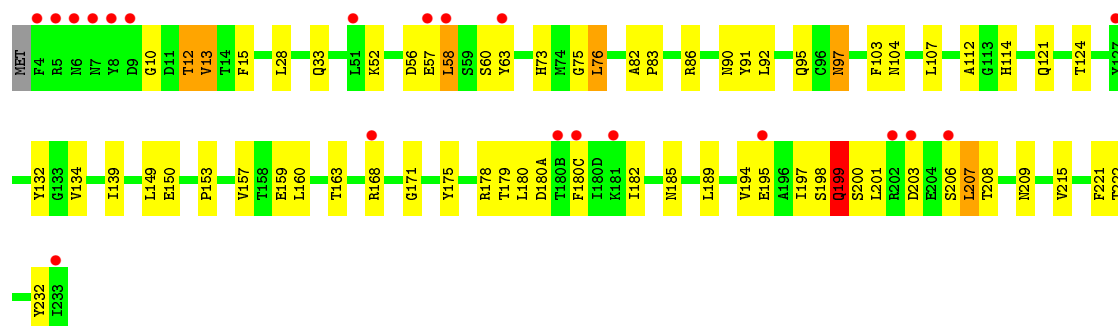




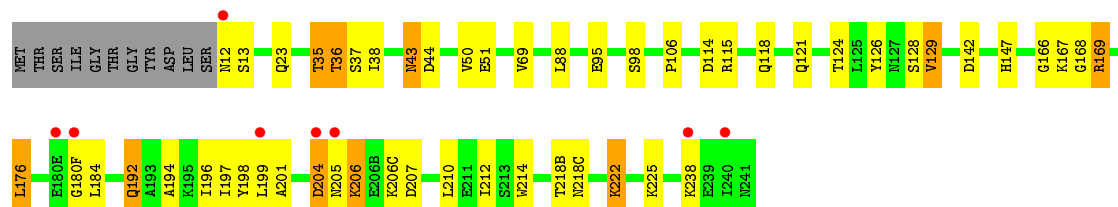
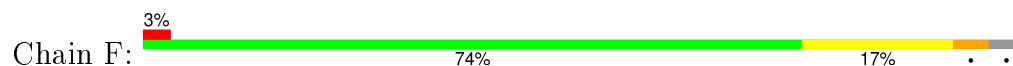
• Molecule 5: Proteasome component PRE5



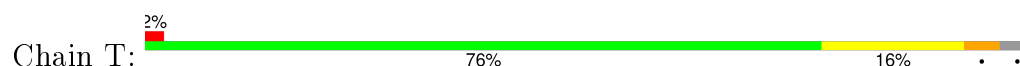
• Molecule 5: Proteasome component PRE5

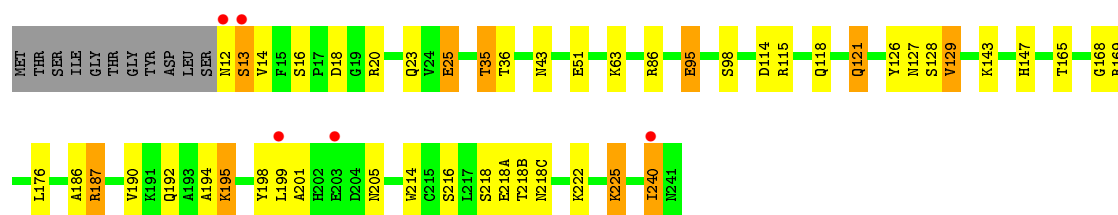


• Molecule 6: Proteasome component C1

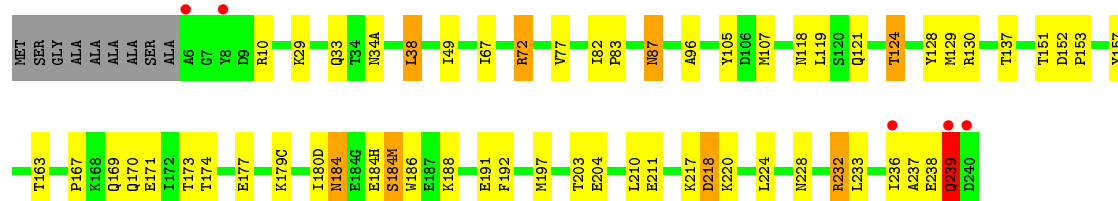


• Molecule 6: Proteasome component C1

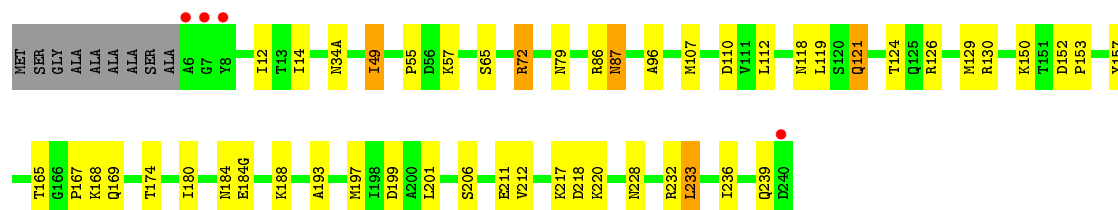
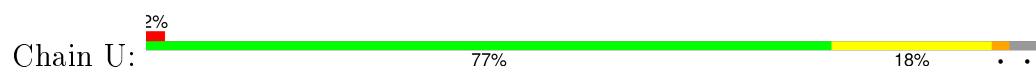




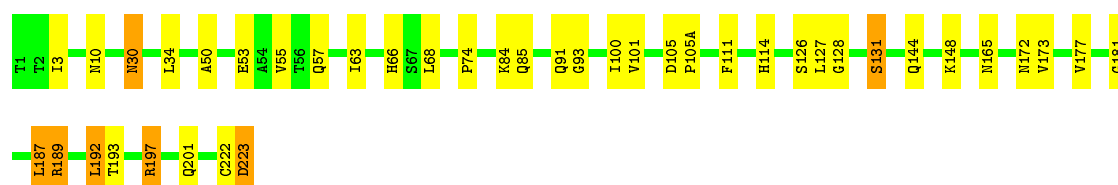
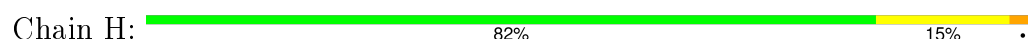
• Molecule 7: Proteasome component C7-alpha



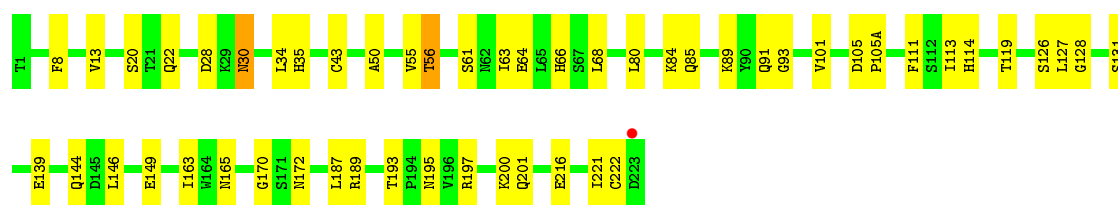
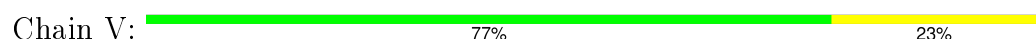
• Molecule 7: Proteasome component C7-alpha




• Molecule 8: Proteasome component PUP1



• Molecule 8: Proteasome component PUP1




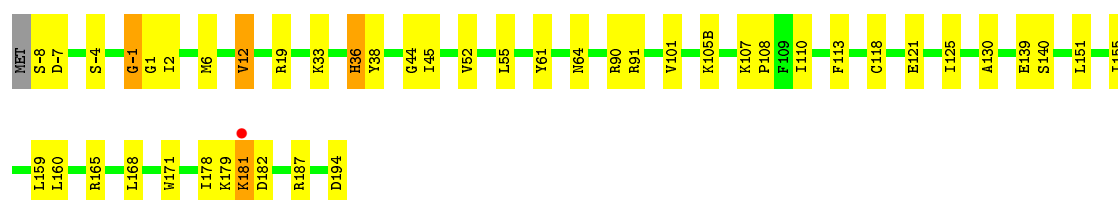
• Molecule 9: Proteasome component PUP3

Chain I:  86% 12%




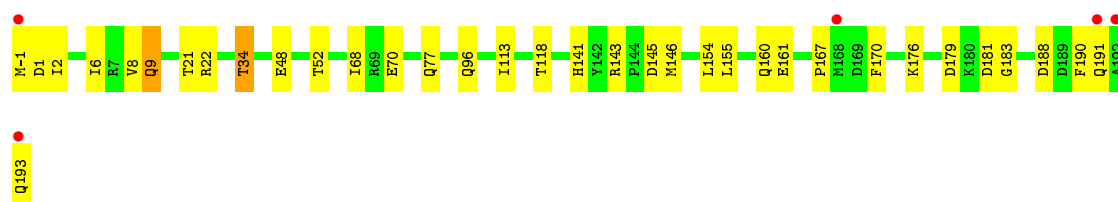
• Molecule 9: Proteasome component PUP3

Chain W:  78% 20%




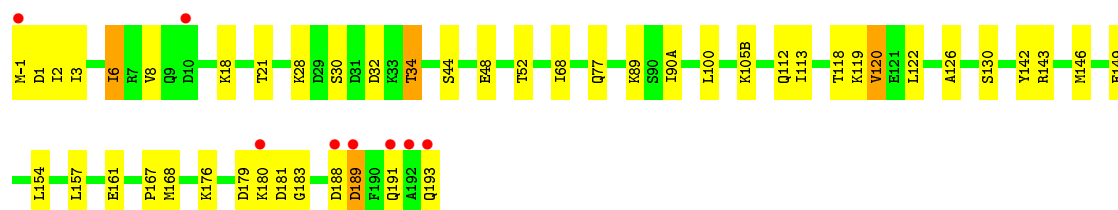
• Molecule 10: Proteasome component C11

Chain J:  3% 82% 17%




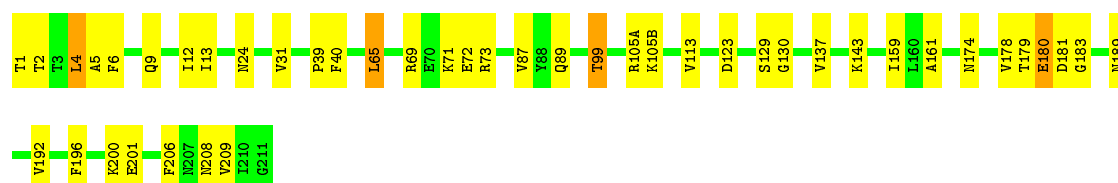
• Molecule 10: Proteasome component C11

Chain X:  4% 76% 22%




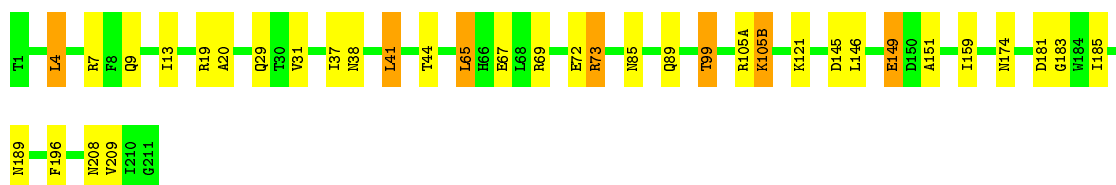
• Molecule 11: Proteasome component PRE2

Chain K:  79% 19%



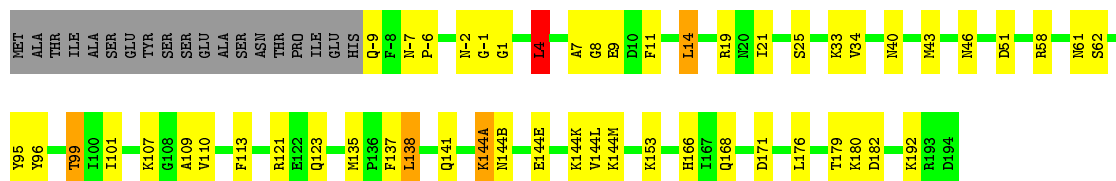
• Molecule 11: Proteasome component PRE2

Chain Y:  83% 14%



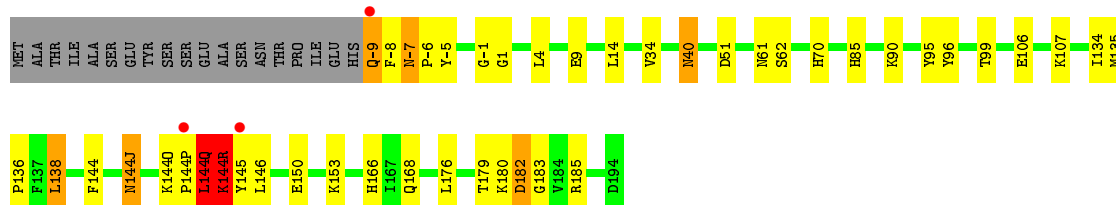
• Molecule 12: Proteasome component C5

Chain L: 70% 20% 8%



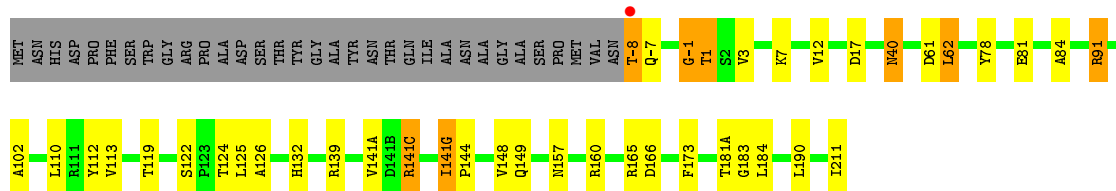
• Molecule 12: Proteasome component C5

Chain Z: 73% 15% 8%



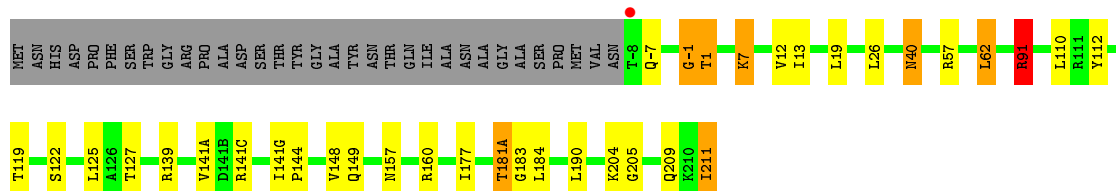
• Molecule 13: Proteasome component PRE4

Chain M: 72% 13% 12%

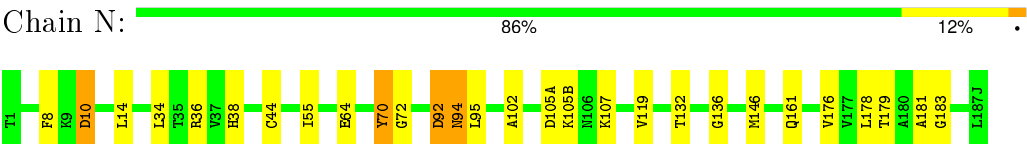


• Molecule 13: Proteasome component PRE4

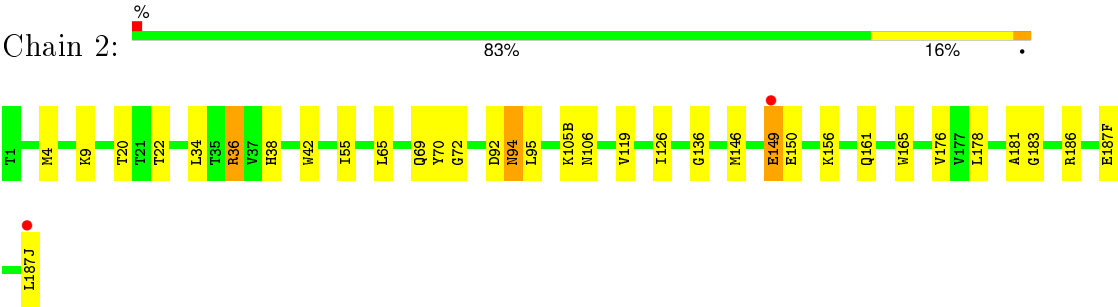
Chain 1: 74% 11% 12%



• 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	137.25Å 299.65Å 145.78Å 90.00° 113.67° 90.00°	Depositor
Resolution (Å)	50.00 – 2.60 49.94 – 2.60	Depositor EDS
% Data completeness (in resolution range)	95.4 (50.00-2.60) 95.4 (49.94-2.60)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.40 (at 2.61Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.215 , 0.249 (Not available) , (Not available)	Depositor DCC
R_{free} test set	NotAvailable	DCC
Wilson B-factor (Å ²)	53.7	Xtriage
Anisotropy	0.182	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 32.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.52$, $\langle L^2 \rangle = 0.36$	Xtriage
Outliers	0 of 314176 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	50276	wwPDB-VP
Average B, all atoms (Å ²)	55.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.17% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: MG, LZT, MES

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	0.51	0/1951	0.63	0/2639
1	O	0.54	0/1951	0.62	0/2639
2	B	0.48	0/1857	0.64	0/2513
2	P	0.50	0/1857	0.62	0/2513
3	C	0.47	0/1918	0.60	1/2591 (0.0%)
3	Q	0.51	0/1918	0.63	1/2591 (0.0%)
4	D	0.52	0/1883	0.67	1/2529 (0.0%)
4	R	0.53	0/1884	0.66	0/2532
5	E	0.48	0/1819	0.64	1/2451 (0.0%)
5	S	0.48	0/1821	0.63	1/2457 (0.0%)
6	F	0.51	0/1885	0.61	0/2540
6	T	0.54	0/1886	0.63	0/2543
7	G	0.81	1/1957 (0.1%)	0.67	2/2646 (0.1%)
7	U	0.55	0/1958	0.63	0/2649
8	H	0.84	1/1715 (0.1%)	0.78	1/2323 (0.0%)
8	V	0.55	1/1714 (0.1%)	0.65	0/2320
9	I	1.13	2/1610 (0.1%)	0.83	3/2171 (0.1%)
9	W	1.14	2/1610 (0.1%)	0.94	5/2171 (0.2%)
10	J	0.55	0/1611	0.67	0/2167
10	X	0.58	0/1611	0.68	0/2167
11	K	0.57	0/1680	0.69	0/2271
11	Y	0.53	0/1680	0.66	0/2271
12	L	0.56	0/1793	0.65	1/2414 (0.0%)
12	Z	0.55	0/1793	0.68	0/2414
13	1	0.76	1/1853 (0.1%)	0.85	3/2507 (0.1%)
13	M	1.06	2/1854 (0.1%)	0.85	5/2511 (0.2%)
14	2	0.58	0/1538	0.66	0/2078
14	N	1.17	2/1540 (0.1%)	0.82	5/2084 (0.2%)
All	All	0.67	12/50147 (0.0%)	0.69	30/67702 (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	O	0	1
8	H	0	1
All	All	0	2

The worst 5 of 12 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
14	N	70	TYR	C-N	34.46	1.95	1.33
13	M	141(G)	ILE	C-N	31.31	1.93	1.34
9	I	-1	GLY	C-N	29.40	1.85	1.33
9	W	-1	GLY	C-N	28.22	1.83	1.33
9	W	36	HIS	C-N	27.17	1.96	1.34

The worst 5 of 30 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
9	W	-1	GLY	O-C-N	-20.50	88.36	123.20
13	M	-1	GLY	O-C-N	-20.43	90.00	122.70
13	1	-1	GLY	O-C-N	-20.28	90.26	122.70
8	H	187	LEU	O-C-N	-19.97	90.75	122.70
9	I	36	HIS	O-C-N	-18.17	93.63	122.70

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
8	H	181	GLY	Peptide
1	O	217(N)	THR	Peptide

5.2 Too-close contacts [i](#)

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	1925	31	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	O	1915	0	1925	45	0
2	B	1829	0	1828	39	0
2	P	1829	0	1828	36	0
3	C	1891	0	1899	38	0
3	Q	1891	0	1899	62	0
4	D	1862	0	1832	42	0
4	R	1862	0	1833	45	0
5	E	1795	0	1793	43	0
5	S	1795	0	1795	43	0
6	F	1848	0	1842	45	0
6	T	1848	0	1843	36	0
7	G	1921	0	1907	53	0
7	U	1921	0	1909	41	0
8	H	1685	0	1686	29	0
8	V	1685	0	1686	36	0
9	I	1581	0	1574	26	0
9	W	1581	0	1574	37	0
10	J	1585	0	1591	24	0
10	X	1585	0	1591	35	0
11	K	1644	0	1594	35	0
11	Y	1644	0	1594	30	0
12	L	1757	0	1712	44	0
12	Z	1757	0	1712	58	0
13	1	1824	0	1833	34	0
13	M	1824	0	1833	39	0
14	2	1512	0	1478	33	0
14	N	1512	0	1478	24	0
15	F	2	0	0	0	0
15	G	1	0	0	0	0
15	H	1	0	0	0	0
15	I	2	0	0	0	0
15	K	1	0	0	0	0
15	L	2	0	0	0	0
15	N	1	0	0	0	0
16	K	43	0	39	5	0
16	Y	43	0	39	4	0
17	K	12	0	13	2	0
17	Y	12	0	13	1	0
18	1	46	0	0	4	0
18	2	32	0	0	6	0
18	A	27	0	0	4	0
18	B	28	0	0	9	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
18	C	27	0	0	7	0
18	D	24	0	0	6	0
18	E	21	0	0	5	0
18	F	38	0	0	10	0
18	G	44	0	0	6	0
18	H	30	0	0	4	0
18	I	37	0	0	1	0
18	J	33	0	0	1	0
18	K	39	0	0	9	0
18	L	34	0	0	5	0
18	M	35	0	0	5	0
18	N	29	0	0	5	0
18	O	23	0	0	9	0
18	P	17	0	0	4	0
18	Q	32	0	0	19	0
18	R	21	0	0	10	0
18	S	23	0	0	4	0
18	T	24	0	0	7	0
18	U	34	0	0	7	0
18	V	27	0	0	4	0
18	W	32	0	0	3	0
18	X	38	0	0	5	0
18	Y	32	0	0	6	0
18	Z	31	0	0	6	0
All	All	50276	0	49098	979	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 10.

The worst 5 of 979 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)
13:1:-1:GLY:C	13:1:1:THR:H1	1.21	1.40
14:N:92:ASP:C	14:N:94:ASN:N	1.76	1.39
9:W:-1:GLY:C	9:W:1:GLY:H1	1.24	1.38
8:H:91:GLN:O	8:H:93:GLY:N	1.59	1.32
9:I:-1:GLY:C	9:I:1:GLY:N	1.86	1.28

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	246/250 (98%)	237 (96%)	5 (2%)	4 (2%)	12	24
1	O	246/250 (98%)	233 (95%)	10 (4%)	3 (1%)	16	33
2	B	231/245 (94%)	215 (93%)	14 (6%)	2 (1%)	21	42
2	P	231/245 (94%)	213 (92%)	13 (6%)	5 (2%)	8	15
3	C	235/243 (97%)	226 (96%)	7 (3%)	2 (1%)	21	42
3	Q	235/243 (97%)	223 (95%)	10 (4%)	2 (1%)	21	42
4	D	232/250 (93%)	215 (93%)	13 (6%)	4 (2%)	11	22
4	R	234/250 (94%)	220 (94%)	9 (4%)	5 (2%)	9	16
5	E	223/234 (95%)	211 (95%)	5 (2%)	7 (3%)	5	8
5	S	227/234 (97%)	213 (94%)	12 (5%)	2 (1%)	21	42
6	F	231/248 (93%)	219 (95%)	12 (5%)	0	100	100
6	T	233/248 (94%)	219 (94%)	12 (5%)	2 (1%)	21	42
7	G	237/252 (94%)	228 (96%)	8 (3%)	1 (0%)	39	65
7	U	239/252 (95%)	235 (98%)	2 (1%)	2 (1%)	24	46
8	H	218/222 (98%)	209 (96%)	8 (4%)	1 (0%)	34	60
8	V	216/222 (97%)	207 (96%)	9 (4%)	0	100	100
9	I	200/205 (98%)	194 (97%)	6 (3%)	0	100	100
9	W	200/205 (98%)	191 (96%)	8 (4%)	1 (0%)	34	60
10	J	193/198 (98%)	185 (96%)	7 (4%)	1 (0%)	34	60
10	X	193/198 (98%)	185 (96%)	6 (3%)	2 (1%)	19	39
11	K	208/212 (98%)	202 (97%)	5 (2%)	1 (0%)	34	60
11	Y	208/212 (98%)	201 (97%)	6 (3%)	1 (0%)	34	60
12	L	216/241 (90%)	206 (95%)	10 (5%)	0	100	100
12	Z	216/241 (90%)	206 (95%)	8 (4%)	2 (1%)	21	42
13	1	227/266 (85%)	219 (96%)	7 (3%)	1 (0%)	39	65

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
13	M	229/266 (86%)	218 (95%)	10 (4%)	1 (0%)	39	65
14	2	188/196 (96%)	182 (97%)	6 (3%)	0	100	100
14	N	192/196 (98%)	187 (97%)	5 (3%)	0	100	100
All	All	6184/6524 (95%)	5899 (95%)	233 (4%)	52 (1%)	24	46

5 of 52 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	54	VAL
2	B	217	ALA
3	C	207	ALA
4	D	123(E)	SER
5	E	6	ASN

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%)	199 (95%)	10 (5%)	31	58
1	O	209/209 (100%)	198 (95%)	11 (5%)	28	53
2	B	195/204 (96%)	177 (91%)	18 (9%)	11	21
2	P	195/204 (96%)	174 (89%)	21 (11%)	8	15
3	C	213/215 (99%)	193 (91%)	20 (9%)	11	20
3	Q	213/215 (99%)	199 (93%)	14 (7%)	21	40
4	D	198/206 (96%)	181 (91%)	17 (9%)	13	25
4	R	198/206 (96%)	181 (91%)	17 (9%)	13	25
5	E	192/193 (100%)	168 (88%)	24 (12%)	6	10
5	S	192/193 (100%)	174 (91%)	18 (9%)	11	20
6	F	196/205 (96%)	179 (91%)	17 (9%)	13	24
6	T	196/205 (96%)	177 (90%)	19 (10%)	10	19
7	G	207/210 (99%)	187 (90%)	20 (10%)	10	19

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
7	U	207/210 (99%)	195 (94%)	12 (6%)	25	49
8	H	181/181 (100%)	171 (94%)	10 (6%)	27	51
8	V	181/181 (100%)	168 (93%)	13 (7%)	18	35
9	I	172/173 (99%)	166 (96%)	6 (4%)	43	71
9	W	172/173 (99%)	162 (94%)	10 (6%)	25	49
10	J	175/175 (100%)	165 (94%)	10 (6%)	25	49
10	X	175/175 (100%)	163 (93%)	12 (7%)	19	38
11	K	169/169 (100%)	159 (94%)	10 (6%)	24	47
11	Y	169/169 (100%)	156 (92%)	13 (8%)	16	31
12	L	185/201 (92%)	173 (94%)	12 (6%)	21	42
12	Z	185/201 (92%)	173 (94%)	12 (6%)	21	42
13	1	199/224 (89%)	187 (94%)	12 (6%)	24	47
13	M	199/224 (89%)	190 (96%)	9 (4%)	34	62
14	2	162/162 (100%)	152 (94%)	10 (6%)	23	45
14	N	162/162 (100%)	154 (95%)	8 (5%)	31	57
All	All	5306/5454 (97%)	4921 (93%)	385 (7%)	17	35

5 of 385 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
13	M	-8	THR
2	P	192	LEU
12	Z	62	SER
13	M	141(C)	ARG
1	O	110	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 165 such sidechains are listed below:

Mol	Chain	Res	Type
14	N	161	GLN
4	R	147	GLN
13	1	10	ASN
1	O	97	HIS
2	P	218	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 14 ligands modelled in this entry, 10 are monoatomic - leaving 4 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
16	LZT	K	213	-	46,46,46	1.01	1 (2%)	59,63,63	1.67	6 (10%)
17	MES	K	214	-	11,12,12	0.68	0	14,16,16	1.45	2 (14%)
16	LZT	Y	212	-	46,46,46	1.05	1 (2%)	59,63,63	1.41	5 (8%)
17	MES	Y	213	-	11,12,12	0.66	0	14,16,16	1.75	1 (7%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
16	LZT	K	213	-	-	0/36/48/48	0/4/4/4
17	MES	K	214	-	-	0/6/14/14	0/1/1/1
16	LZT	Y	212	-	-	0/36/48/48	0/4/4/4
17	MES	Y	213	-	-	0/6/14/14	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
16	Y	212	LZT	C8-C9	-5.89	1.40	1.47
16	K	213	LZT	C8-C9	-5.44	1.40	1.47

The worst 5 of 14 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
16	Y	212	LZT	C1-C8-C9	-5.65	105.49	109.58
16	K	213	LZT	C12-C11-C9	-5.54	92.48	104.87
16	K	213	LZT	C8-C1-C12	-4.83	105.88	111.55
16	K	213	LZT	C1-C8-C9	-4.58	106.26	109.58
16	Y	212	LZT	C12-C11-C9	-4.55	94.69	104.87

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
16	K	213	LZT	5	0
17	K	214	MES	2	0
16	Y	212	LZT	4	0
17	Y	213	MES	1	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 [i](#)

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

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	250/250 (100%)	-0.00	7 (2%) 56 49	38, 56, 80, 104	0
1	O	250/250 (100%)	0.00	10 (4%) 42 34	38, 57, 80, 104	0
2	B	235/245 (95%)	0.13	13 (5%) 29 21	36, 60, 90, 96	0
2	P	235/245 (95%)	0.10	8 (3%) 49 41	36, 60, 90, 96	0
3	C	241/243 (99%)	0.44	22 (9%) 11 7	43, 71, 111, 126	0
3	Q	241/243 (99%)	0.80	44 (18%) 2 1	44, 71, 111, 126	0
4	D	242/250 (96%)	0.22	12 (4%) 32 26	35, 62, 96, 103	0
4	R	242/250 (96%)	0.32	17 (7%) 19 13	36, 62, 96, 102	0
5	E	233/234 (99%)	0.18	12 (5%) 31 24	42, 62, 93, 105	0
5	S	233/234 (99%)	0.37	20 (8%) 13 8	42, 62, 92, 105	0
6	F	237/248 (95%)	-0.01	8 (3%) 49 41	35, 54, 85, 96	0
6	T	237/248 (95%)	-0.07	5 (2%) 67 61	34, 54, 82, 96	0
7	G	243/252 (96%)	-0.20	5 (2%) 67 61	33, 49, 73, 94	0
7	U	243/252 (96%)	-0.20	4 (1%) 74 69	32, 49, 72, 94	0
8	H	222/222 (100%)	-0.22	0 100 100	36, 46, 62, 89	0
8	V	222/222 (100%)	-0.29	1 (0%) 91 90	36, 47, 63, 90	0
9	I	204/205 (99%)	-0.18	1 (0%) 91 90	34, 45, 61, 77	0
9	W	204/205 (99%)	-0.12	1 (0%) 91 90	34, 45, 62, 77	0
10	J	198/198 (100%)	-0.11	5 (2%) 61 54	37, 48, 63, 120	0
10	X	198/198 (100%)	-0.11	8 (4%) 42 34	36, 49, 63, 120	0
11	K	212/212 (100%)	-0.30	0 100 100	34, 45, 63, 72	0
11	Y	212/212 (100%)	-0.27	0 100 100	34, 45, 63, 72	0
12	L	222/241 (92%)	-0.27	0 100 100	33, 46, 68, 73	0
12	Z	222/241 (92%)	-0.21	3 (1%) 78 74	32, 46, 68, 73	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
13	1	233/266 (87%)	-0.33	1 (0%) 93 91	33, 44, 57, 63	0
13	M	233/266 (87%)	-0.29	1 (0%) 93 91	34, 44, 58, 63	0
14	2	196/196 (100%)	-0.28	2 (1%) 84 81	33, 41, 58, 68	0
14	N	196/196 (100%)	-0.33	0 100 100	33, 41, 58, 69	0
All	All	6336/6524 (97%)	-0.03	210 (3%) 50 43	32, 51, 89, 126	0

The worst 5 of 210 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
4	R	123(F)	GLY	16.4
4	D	123(E)	SER	11.7
4	D	123(B)	GLU	11.2
4	R	123(E)	SER	10.9
3	C	7	GLY	10.8

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	MG	F	243	1/1	0.81	0.79	8.46	116,116,116,116	0
15	MG	I	196	1/1	0.71	0.28	6.65	65,65,65,65	0
15	MG	F	242	1/1	0.85	0.20	2.56	67,67,67,67	0
17	MES	K	214	12/12	0.93	0.20	2.49	78,83,85,85	0
15	MG	L	195	1/1	0.77	0.19	2.14	58,58,58,58	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
17	MES	Y	213	12/12	0.94	0.17	1.39	79,83,85,85	0
15	MG	L	196	1/1	0.86	0.19	0.54	57,57,57,57	0
16	LZT	K	213	43/43	0.95	0.17	0.52	32,38,45,49	0
15	MG	K	212	1/1	0.84	0.15	0.43	53,53,53,53	0
16	LZT	Y	212	43/43	0.97	0.15	0.15	34,39,45,48	0
15	MG	N	188	1/1	0.94	0.15	-0.09	40,40,40,40	0
15	MG	G	241	1/1	0.88	0.08	-1.99	66,66,66,66	0
15	MG	H	224	1/1	0.97	0.06	-4.57	50,50,50,50	0
15	MG	I	195	1/1	0.70	0.36	-	78,78,78,78	0

6.5 Other polymers [i](#)

There are no such residues in this entry.