



wwPDB X-ray Structure Validation Summary Report ⓘ

May 15, 2020 – 08:54 am BST

PDB ID : 4GK7
Title : yeast 20S proteasome in complex with the Syringolin-Glidobactin chimera
Authors : Groll, M.; Stein, M.L.; Bachmann, A.
Deposited on : 2012-08-10
Resolution : 2.80 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

<https://www.wwpdb.org/validation/2017/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 : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

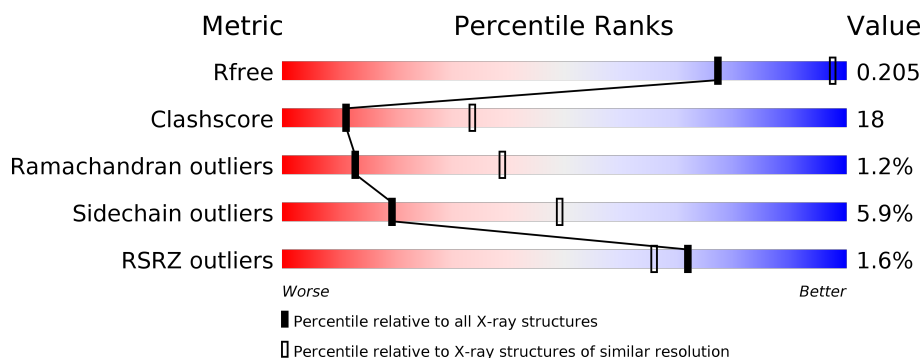
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	3140 (2.80-2.80)
Clashscore	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)
RSRZ outliers	127900	3078 (2.80-2.80)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria 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	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 0%, red 1%, green 1%, green 71%, yellow 71%, yellow 97%, orange 97%, orange 100%);"></div> <div style="display: flex; justify-content: space-between; width: 90%; margin: 0 auto;"> % 71% 27% • </div> </div>
1	O	250	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 0%, red 1%, green 1%, green 69%, yellow 69%, yellow 97%, orange 97%, orange 100%);"></div> <div style="display: flex; justify-content: space-between; width: 90%; margin: 0 auto;"> % 69% 29% • </div> </div>
2	B	244	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 0%, red 3%, green 3%, green 64%, yellow 64%, yellow 93%, orange 93%, orange 100%);"></div> <div style="display: flex; justify-content: space-between; width: 90%; margin: 0 auto;"> 3% 64% 29% 7% </div> </div>
2	P	244	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 0%, red 3%, green 3%, green 65%, yellow 65%, yellow 93%, orange 93%, orange 100%);"></div> <div style="display: flex; justify-content: space-between; width: 90%; margin: 0 auto;"> 3% 65% 28% 7% </div> </div>
3	C	241	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 0%, red 3%, green 3%, green 59%, yellow 59%, yellow 97%, orange 97%, orange 100%);"></div> <div style="display: flex; justify-content: space-between; width: 90%; margin: 0 auto;"> 3% 59% 38% • </div> </div>
3	Q	241	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 0%, red 5%, green 5%, green 59%, yellow 59%, yellow 97%, orange 97%, orange 100%);"></div> <div style="display: flex; justify-content: space-between; width: 90%; margin: 0 auto;"> 5% 59% 38% • </div> </div>

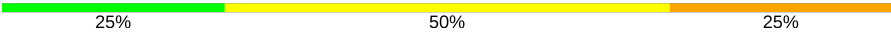


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Mol	Chain	Length	Quality of chain
4	D	242	
4	R	242	
5	E	233	
5	S	233	
6	F	244	
6	T	244	
7	G	243	
7	U	243	
8	H	222	
8	V	222	
9	I	204	
9	W	204	
10	J	198	
10	X	198	
11	K	212	
11	Y	212	
12	L	222	
12	Z	222	
13	M	233	
13	a	233	
14	N	196	
14	b	196	
15	1	4	
15	2	4	
15	3	4	

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Mol	Chain	Length	Quality of chain
15	4	4	 25% 50% 25%
15	5	4	 50% 50%
15	6	4	 75% 25%

2 Entry composition

There are 16 unique types of molecules in this entry. The entry contains 51109 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Proteasome component Y7.

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

- Molecule 2 is a protein called Proteasome component Y13.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	244	Total	C	N	O	S	0	0	0
			1905	1201	321	380	3			
2	P	244	Total	C	N	O	S	0	0	0
			1905	1201	321	380	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	244	Total	C	N	O	S	0	0	0
			1897	1205	330	358	4			
6	T	244	Total	C	N	O	S	0	0	0
			1897	1205	330	358	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	a	233	Total	C	N	O	S	0	0	0
			1824	1154	312	351	7			

- Molecule 14 is a protein called Proteasome component PRE3.

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

- Molecule 15 is a protein called SYRINGOLIN-GLIDOBACTIN CHIMERA.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
15	1	4	Total	C	N	O	0	0	0
			38	29	4	5			
15	2	4	Total	C	N	O	0	0	0
			38	29	4	5			

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
15	3	4	Total	C	N	O	0	0	0
			38	29	4	5			
15	4	4	Total	C	N	O	0	0	0
			38	29	4	5			
15	5	4	Total	C	N	O	0	0	0
			38	29	4	5			
15	6	4	Total	C	N	O	0	0	0
			38	29	4	5			

- Molecule 16 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
16	A	59	Total	O	0	0
			59	59		
16	B	35	Total	O	0	0
			35	35		
16	C	44	Total	O	0	0
			44	44		
16	D	39	Total	O	0	0
			39	39		
16	E	23	Total	O	0	0
			23	23		
16	F	49	Total	O	0	0
			49	49		
16	G	60	Total	O	0	0
			60	60		
16	H	49	Total	O	0	0
			49	49		
16	I	64	Total	O	0	0
			64	64		
16	J	53	Total	O	0	0
			53	53		
16	K	42	Total	O	0	0
			42	42		
16	L	55	Total	O	0	0
			55	55		
16	M	75	Total	O	0	0
			75	75		
16	N	56	Total	O	0	0
			56	56		
16	O	34	Total	O	0	0
			34	34		

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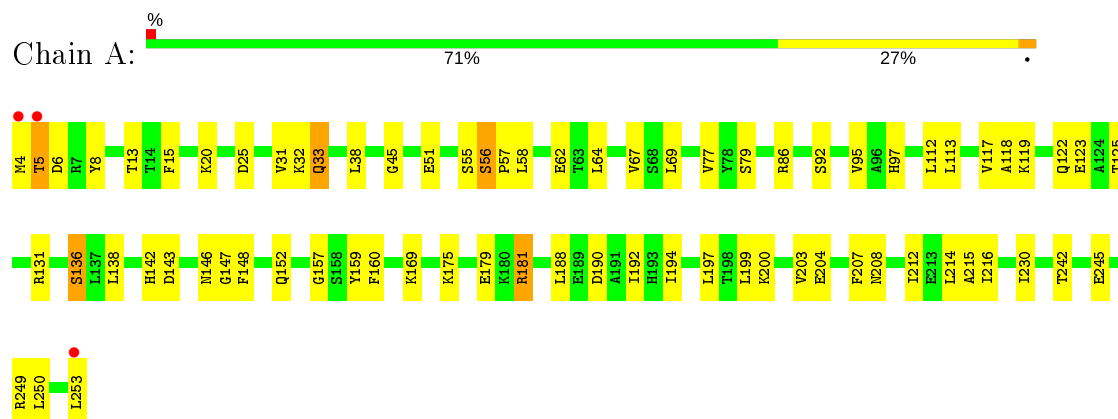
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
16	P	28	Total 28	O 28	0	0
16	Q	29	Total 29	O 29	0	0
16	R	33	Total 33	O 33	0	0
16	S	21	Total 21	O 21	0	0
16	T	39	Total 39	O 39	0	0
16	U	63	Total 63	O 63	0	0
16	V	47	Total 47	O 47	0	0
16	W	55	Total 55	O 55	0	0
16	X	46	Total 46	O 46	0	0
16	Y	50	Total 50	O 50	0	0
16	Z	51	Total 51	O 51	0	0
16	a	73	Total 73	O 73	0	0
16	b	58	Total 58	O 58	0	0
16	1	1	Total 1	O 1	0	0
16	2	1	Total 1	O 1	0	0
16	3	1	Total 1	O 1	0	0

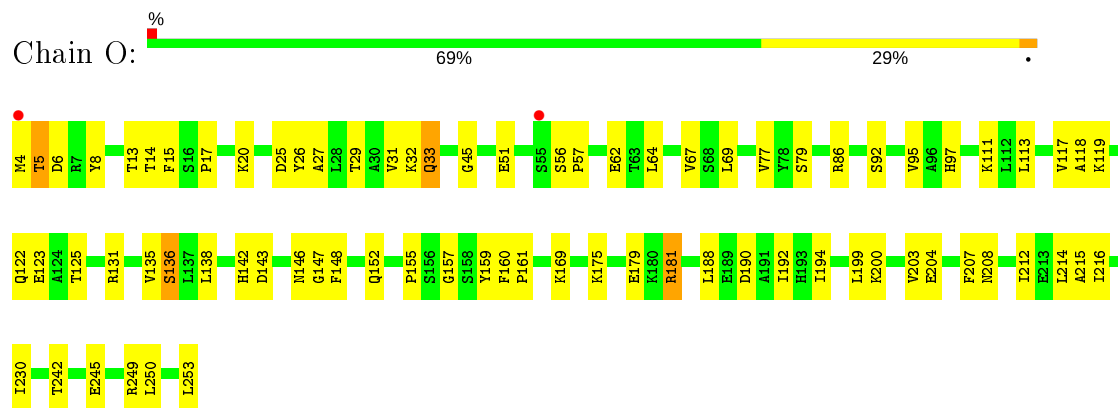
3 Residue-property plots [i](#)

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

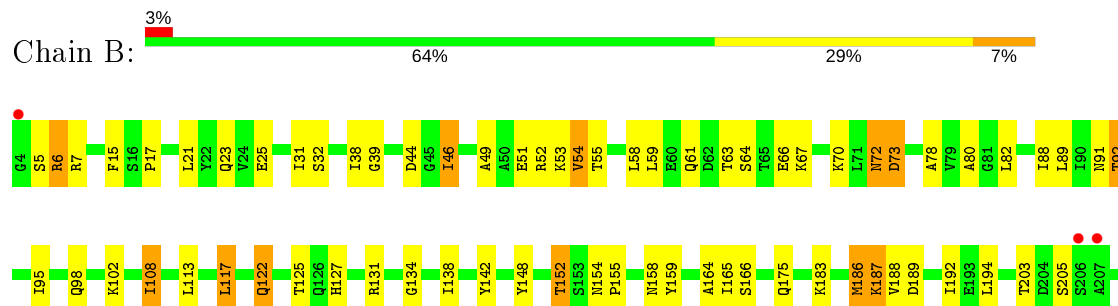
• Molecule 1: Proteasome component Y7



• Molecule 1: Proteasome component Y7

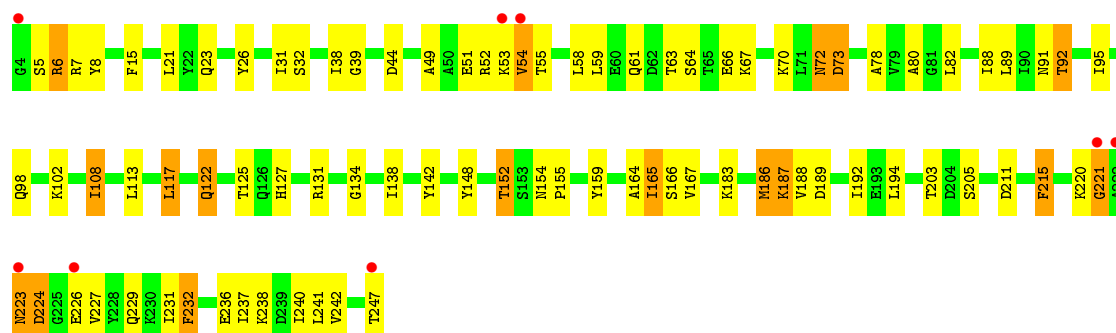


• Molecule 2: Proteasome component Y13

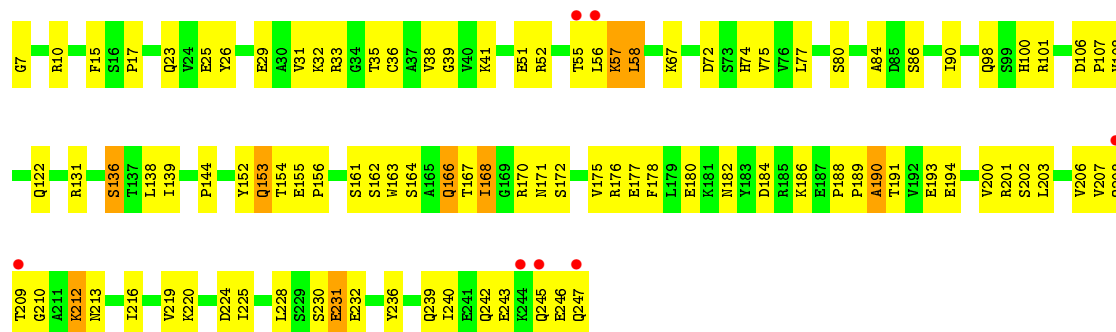




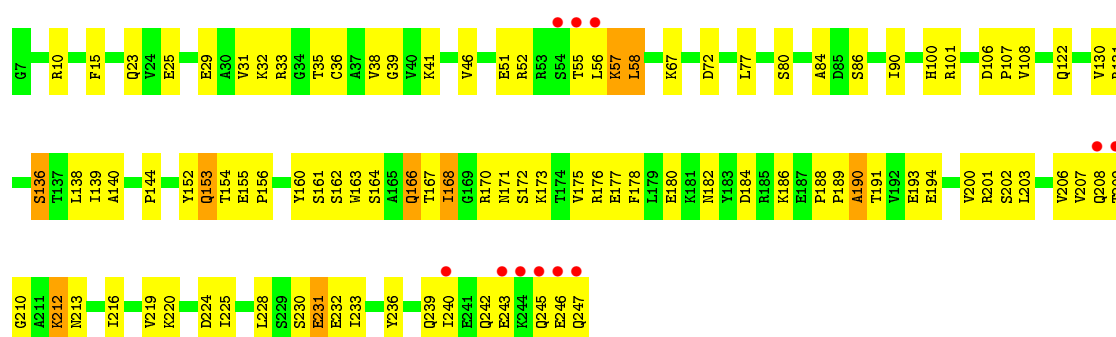
• Molecule 2: Proteasome component Y13



• Molecule 3: Proteasome component PRE6

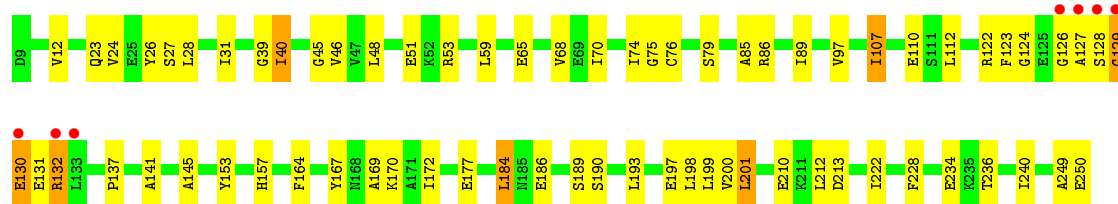


• Molecule 3: Proteasome component PRE6

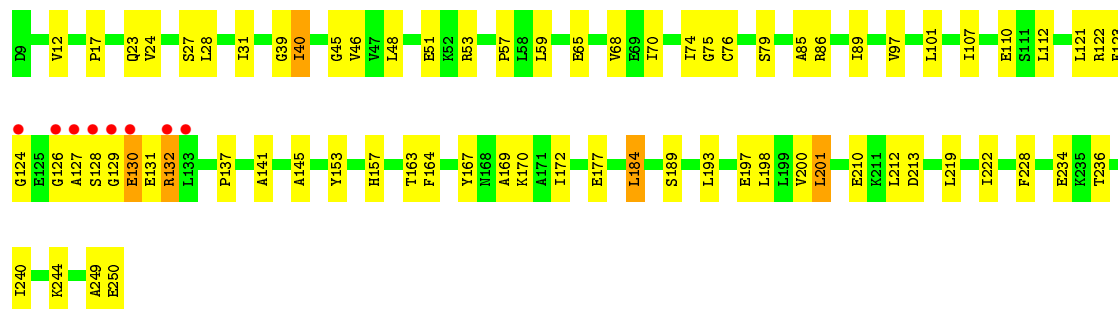


• Molecule 4: Proteasome component PUP2

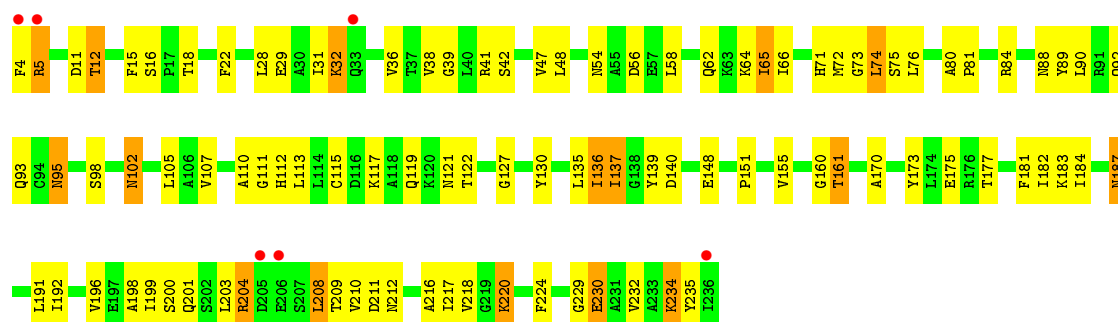




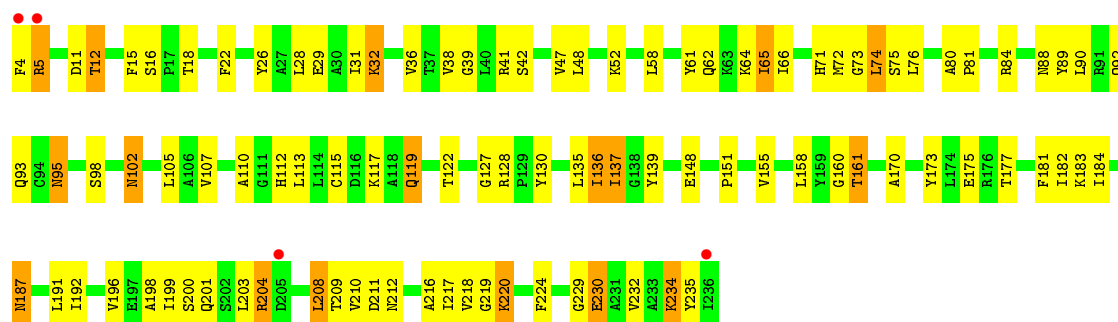
• Molecule 4: Proteasome component PUP2



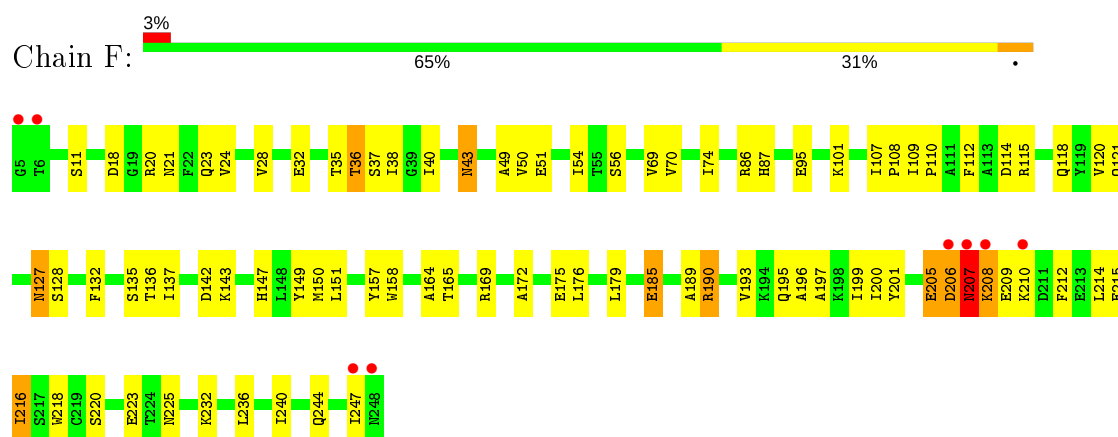
• Molecule 5: Proteasome component PRE5



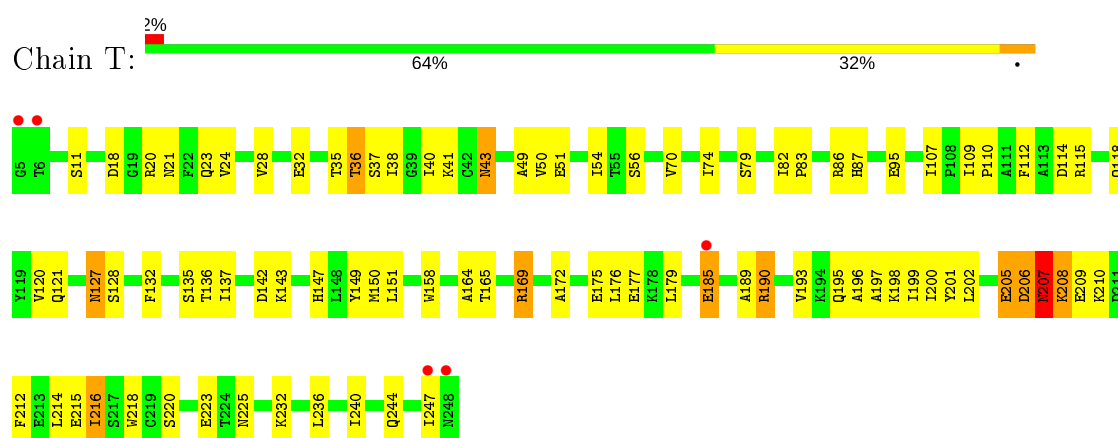
• Molecule 5: Proteasome component PRE5



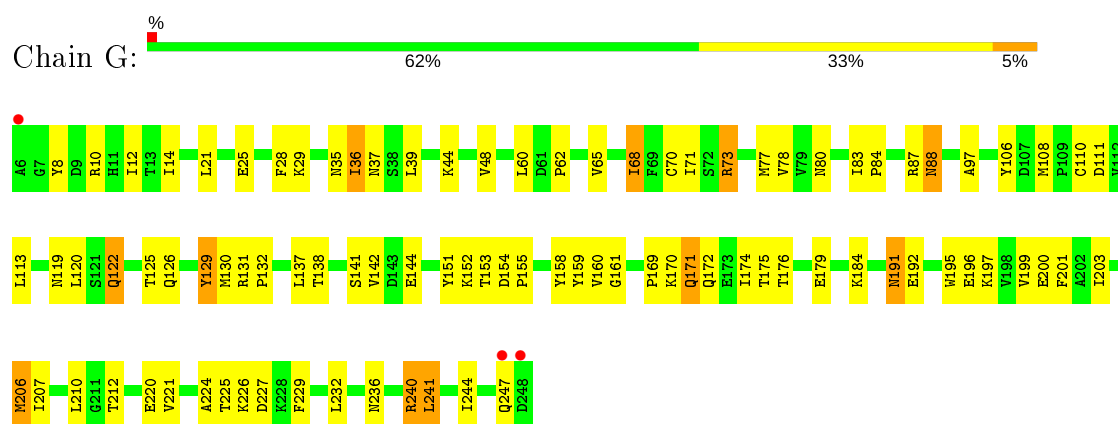
• Molecule 6: Proteasome component C1



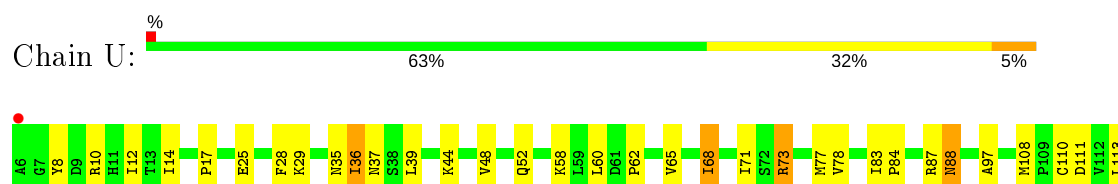
• Molecule 6: Proteasome component C1

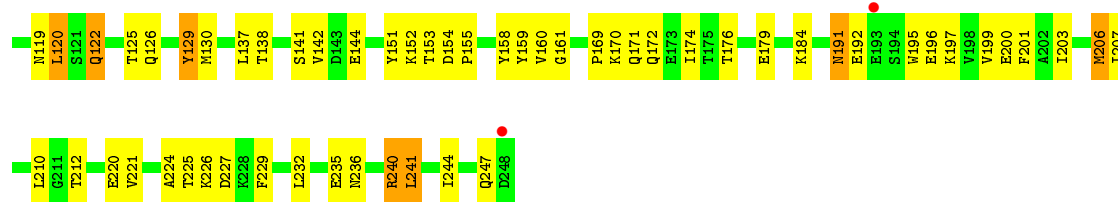


• Molecule 7: Proteasome component C7-alpha

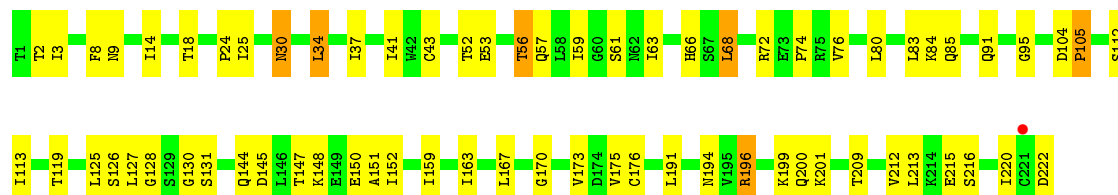


• Molecule 7: Proteasome component C7-alpha

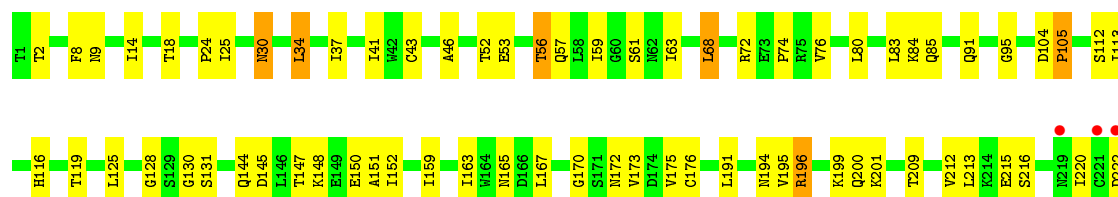




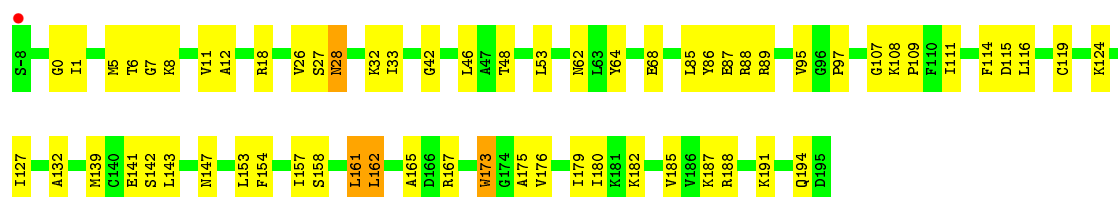
• Molecule 8: Proteasome component PUP1



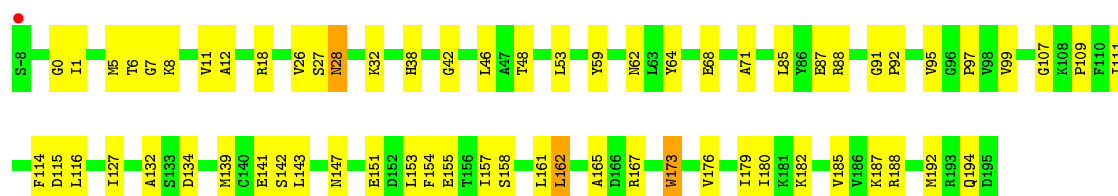
• Molecule 8: Proteasome component PUP1



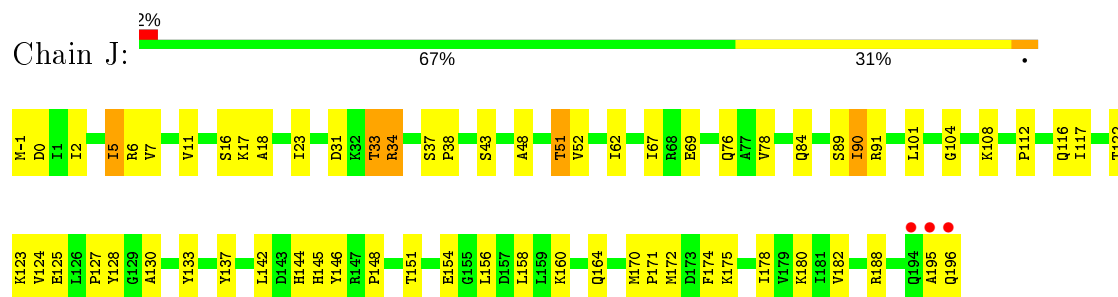
• Molecule 9: Proteasome component PUP3



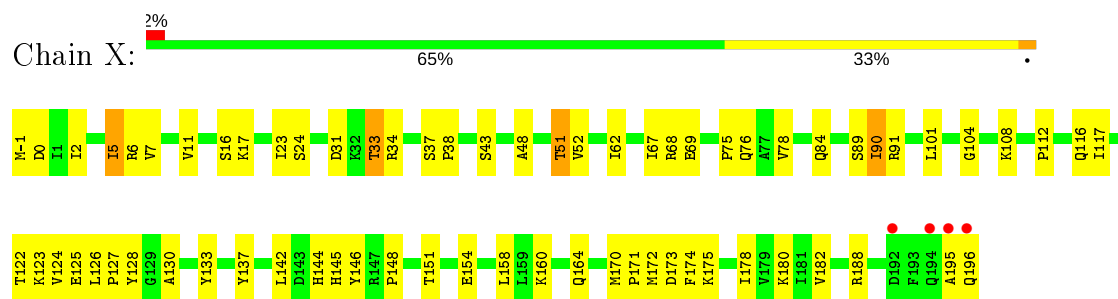
• Molecule 9: Proteasome component PUP3



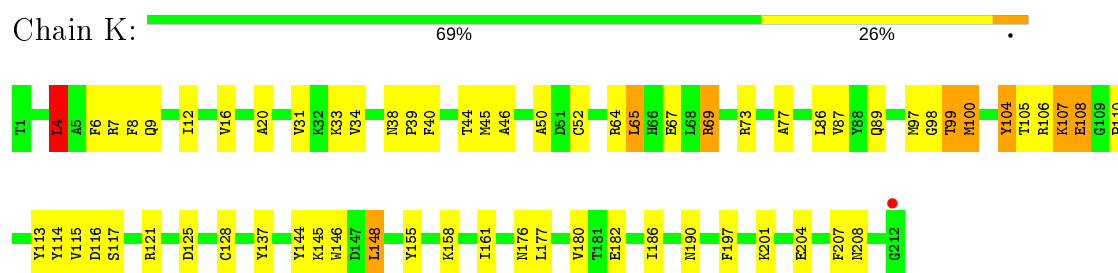
• Molecule 10: Proteasome component C11



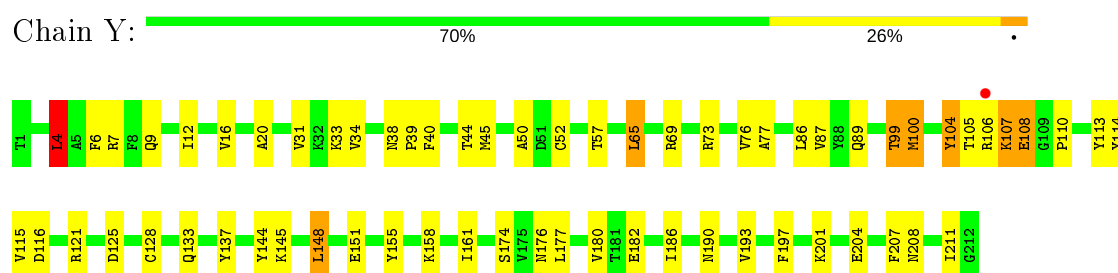
- Molecule 10: Proteasome component C11



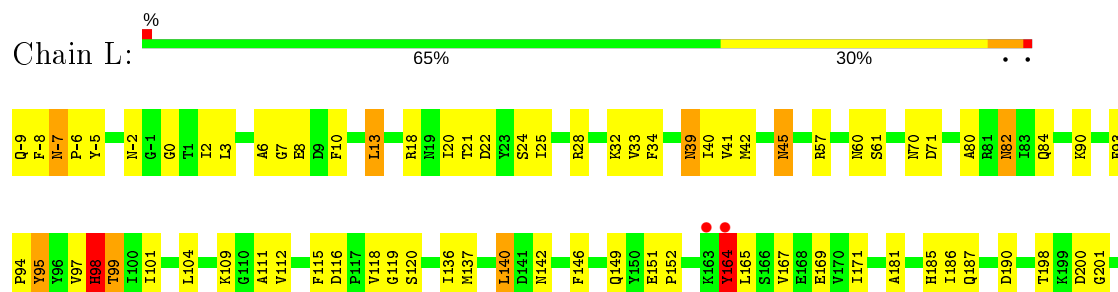
- Molecule 11: Proteasome component PRE2



- Molecule 11: Proteasome component PRE2

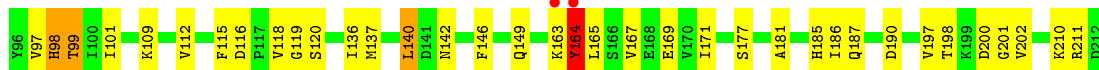


- Molecule 12: Proteasome component C5





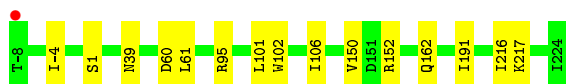
- Molecule 12: Proteasome component C5



- Molecule 13: Proteasome component PRE4



- Molecule 13: Proteasome component PRE4



- Molecule 14: Proteasome component PRE3



- Molecule 14: Proteasome component PRE3



- Molecule 15: SYRINGOLIN-GLIDOBACTIN CHIMERA

Chain 1:  50% 25% 25%



- Molecule 15: SYRINGOLIN-GLIDOBACTIN CHIMERA

Chain 2:  75% 25%



- Molecule 15: SYRINGOLIN-GLIDOBACTIN CHIMERA

Chain 3:  25% 50% 25%



- Molecule 15: SYRINGOLIN-GLIDOBACTIN CHIMERA

Chain 4:  25% 50% 25%



- Molecule 15: SYRINGOLIN-GLIDOBACTIN CHIMERA

Chain 5:  50% 50%



- Molecule 15: SYRINGOLIN-GLIDOBACTIN CHIMERA

Chain 6:  75% 25%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	135.85Å 298.45Å 145.25Å 90.00° 112.63° 90.00°	Depositor
Resolution (Å)	15.00 – 2.80 29.84 – 2.80	Depositor EDS
% Data completeness (in resolution range)	97.8 (15.00-2.80) 97.9 (29.84-2.80)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.94 (at 2.80Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.210 , 0.241 0.206 , 0.205	Depositor DCC
R_{free} test set	12775 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	50.3	Xtriage
Anisotropy	0.648	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 57.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	51109	wwPDB-VP
Average B, all atoms (Å ²)	56.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.16% 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: 0JT, LYH, MH9

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	0.37	0/1952	0.64	0/2642
1	O	0.38	0/1952	0.63	0/2642
2	B	0.38	0/1935	0.64	0/2618
2	P	0.39	0/1935	0.64	0/2618
3	C	0.36	0/1920	0.62	0/2598
3	Q	0.38	0/1920	0.62	0/2598
4	D	0.36	0/1887	0.64	0/2541
4	R	0.36	0/1887	0.64	0/2541
5	E	0.36	0/1823	0.63	0/2463
5	S	0.36	0/1823	0.63	0/2463
6	F	0.39	0/1937	0.62	0/2614
6	T	0.39	0/1937	0.62	0/2614
7	G	0.41	0/1959	0.64	0/2652
7	U	0.39	0/1959	0.63	0/2652
8	H	0.41	0/1716	0.66	0/2326
8	V	0.39	0/1716	0.66	0/2326
9	I	0.39	0/1611	0.67	0/2174
9	W	0.41	0/1611	0.67	0/2174
10	J	0.40	0/1613	0.65	0/2173
10	X	0.40	0/1613	0.65	0/2173
11	K	0.41	0/1681	0.66	1/2274 (0.0%)
11	Y	0.39	0/1681	0.66	1/2274 (0.0%)
12	L	0.40	0/1795	0.69	3/2420 (0.1%)
12	Z	0.39	0/1795	0.69	3/2420 (0.1%)
13	M	0.39	0/1855	0.67	0/2514
13	a	0.38	0/1855	0.67	1/2514 (0.0%)
14	N	0.40	0/1541	0.64	0/2087
14	b	0.39	0/1541	0.64	0/2087
15	1	1.68	0/6	1.40	0/7
15	2	1.23	0/6	1.89	0/7
15	3	1.71	0/6	1.64	0/7
15	4	1.75	0/6	1.44	0/7

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
15	5	1.31	0/6	1.84	0/7
15	6	1.69	0/6	1.68	0/7
All	All	0.39	0/50486	0.65	9/68234 (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
12	L	0	1
12	Z	0	1
All	All	0	2

There are no bond length outliers.

The worst 5 of 9 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
12	L	164	TYR	CA-CB-CG	-5.36	103.22	113.40
11	Y	4	LEU	CA-CB-CG	5.32	127.53	115.30
11	K	4	LEU	CA-CB-CG	5.29	127.48	115.30
12	Z	98	HIS	N-CA-C	-5.27	96.77	111.00
12	Z	164	TYR	CA-CB-CG	-5.27	103.39	113.40

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
12	L	164	TYR	Sidechain
12	Z	164	TYR	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	1915	0	1926	66	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	O	1915	0	1926	68	0
2	B	1905	0	1901	78	0
2	P	1905	0	1901	84	0
3	C	1891	0	1900	91	0
3	Q	1891	0	1900	88	0
4	D	1862	0	1836	52	0
4	R	1862	0	1836	55	0
5	E	1795	0	1797	95	0
5	S	1795	0	1797	97	0
6	F	1897	0	1886	77	0
6	T	1897	0	1886	79	0
7	G	1921	0	1910	84	0
7	U	1921	0	1910	87	0
8	H	1685	0	1687	57	0
8	V	1685	0	1687	58	0
9	I	1581	0	1574	59	0
9	W	1581	0	1574	59	0
10	J	1585	0	1590	73	0
10	X	1585	0	1590	73	0
11	K	1644	0	1594	55	0
11	Y	1644	0	1594	58	0
12	L	1757	0	1711	64	0
12	Z	1757	0	1711	60	0
13	M	1824	0	1832	60	0
13	a	1824	0	1832	0	0
14	N	1512	0	1480	46	0
14	b	1512	0	1480	0	0
15	1	38	0	45	4	0
15	2	38	0	45	1	0
15	3	38	0	45	3	0
15	4	38	0	45	5	0
15	5	38	0	45	2	0
15	6	38	0	45	0	0
16	1	1	0	0	0	0
16	2	1	0	0	0	0
16	3	1	0	0	0	0
16	A	59	0	0	1	0
16	B	35	0	0	0	0
16	C	44	0	0	5	0
16	D	39	0	0	4	0
16	E	23	0	0	2	0
16	F	49	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
16	G	60	0	0	6	0
16	H	49	0	0	3	0
16	I	64	0	0	2	0
16	J	53	0	0	2	0
16	K	42	0	0	3	0
16	L	55	0	0	4	0
16	M	75	0	0	3	0
16	N	56	0	0	3	0
16	O	34	0	0	1	0
16	P	28	0	0	1	0
16	Q	29	0	0	2	0
16	R	33	0	0	2	0
16	S	21	0	0	0	0
16	T	39	0	0	2	0
16	U	63	0	0	2	0
16	V	47	0	0	5	0
16	W	55	0	0	2	0
16	X	46	0	0	4	0
16	Y	50	0	0	5	0
16	Z	51	0	0	2	0
16	a	73	0	0	0	0
16	b	58	0	0	0	0
All	All	51109	0	49518	1646	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 18.

The worst 5 of 1646 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)
2:P:203:THR:HG22	2:P:205:SER:H	1.08	1.13
11:K:107:LYS:H	11:K:107:LYS:HD2	1.03	1.11
11:Y:107:LYS:H	11:Y:107:LYS:HD2	1.04	1.11
2:B:203:THR:HG22	2:B:205:SER:H	1.11	1.06
2:P:72:ASN:ND2	2:P:73:ASP:H	1.56	1.04

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	248/250 (99%)	235 (95%)	10 (4%)	3 (1%)	13	39
1	O	248/250 (99%)	236 (95%)	9 (4%)	3 (1%)	13	39
2	B	242/244 (99%)	214 (88%)	22 (9%)	6 (2%)	5	19
2	P	242/244 (99%)	215 (89%)	21 (9%)	6 (2%)	5	19
3	C	239/241 (99%)	220 (92%)	14 (6%)	5 (2%)	7	23
3	Q	239/241 (99%)	221 (92%)	13 (5%)	5 (2%)	7	23
4	D	240/242 (99%)	219 (91%)	16 (7%)	5 (2%)	7	23
4	R	240/242 (99%)	221 (92%)	14 (6%)	5 (2%)	7	23
5	E	231/233 (99%)	207 (90%)	20 (9%)	4 (2%)	9	29
5	S	231/233 (99%)	207 (90%)	20 (9%)	4 (2%)	9	29
6	F	242/244 (99%)	226 (93%)	13 (5%)	3 (1%)	13	39
6	T	242/244 (99%)	226 (93%)	13 (5%)	3 (1%)	13	39
7	G	241/243 (99%)	228 (95%)	12 (5%)	1 (0%)	34	66
7	U	241/243 (99%)	228 (95%)	12 (5%)	1 (0%)	34	66
8	H	220/222 (99%)	206 (94%)	10 (4%)	4 (2%)	8	28
8	V	220/222 (99%)	207 (94%)	9 (4%)	4 (2%)	8	28
9	I	202/204 (99%)	195 (96%)	7 (4%)	0	100	100
9	W	202/204 (99%)	195 (96%)	7 (4%)	0	100	100
10	J	196/198 (99%)	184 (94%)	9 (5%)	3 (2%)	10	33
10	X	196/198 (99%)	184 (94%)	9 (5%)	3 (2%)	10	33
11	K	210/212 (99%)	199 (95%)	10 (5%)	1 (0%)	29	61
11	Y	210/212 (99%)	199 (95%)	10 (5%)	1 (0%)	29	61
12	L	220/222 (99%)	206 (94%)	13 (6%)	1 (0%)	29	61
12	Z	220/222 (99%)	208 (94%)	10 (4%)	2 (1%)	17	46
13	M	231/233 (99%)	215 (93%)	14 (6%)	2 (1%)	17	46

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
13	a	231/233 (99%)	215 (93%)	14 (6%)	2 (1%)	17	46
14	N	194/196 (99%)	187 (96%)	7 (4%)	0	100	100
14	b	194/196 (99%)	187 (96%)	7 (4%)	0	100	100
15	1	1/4 (25%)	0	1 (100%)	0	100	100
15	2	1/4 (25%)	1 (100%)	0	0	100	100
15	3	1/4 (25%)	1 (100%)	0	0	100	100
15	4	1/4 (25%)	0	1 (100%)	0	100	100
15	5	1/4 (25%)	1 (100%)	0	0	100	100
15	6	1/4 (25%)	1 (100%)	0	0	100	100
All	All	6318/6392 (99%)	5894 (93%)	347 (6%)	77 (1%)	13	39

5 of 77 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	5	THR
1	A	169	LYS
2	B	186	MET
2	B	224	ASP
3	C	58	LEU

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%)	204 (98%)	5 (2%)	49	81
1	O	209/209 (100%)	204 (98%)	5 (2%)	49	81
2	B	203/203 (100%)	186 (92%)	17 (8%)	11	31
2	P	203/203 (100%)	188 (93%)	15 (7%)	13	37
3	C	213/213 (100%)	201 (94%)	12 (6%)	21	51
3	Q	213/213 (100%)	201 (94%)	12 (6%)	21	51
4	D	198/198 (100%)	188 (95%)	10 (5%)	24	55

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
4	R	198/198 (100%)	188 (95%)	10 (5%)	24	55
5	E	192/192 (100%)	173 (90%)	19 (10%)	8	23
5	S	192/192 (100%)	173 (90%)	19 (10%)	8	23
6	F	201/201 (100%)	185 (92%)	16 (8%)	12	34
6	T	201/201 (100%)	185 (92%)	16 (8%)	12	34
7	G	207/207 (100%)	195 (94%)	12 (6%)	20	50
7	U	207/207 (100%)	195 (94%)	12 (6%)	20	50
8	H	181/181 (100%)	175 (97%)	6 (3%)	38	72
8	V	181/181 (100%)	175 (97%)	6 (3%)	38	72
9	I	172/172 (100%)	167 (97%)	5 (3%)	42	76
9	W	172/172 (100%)	168 (98%)	4 (2%)	50	82
10	J	175/175 (100%)	165 (94%)	10 (6%)	20	50
10	X	175/175 (100%)	165 (94%)	10 (6%)	20	50
11	K	169/169 (100%)	157 (93%)	12 (7%)	14	39
11	Y	169/169 (100%)	158 (94%)	11 (6%)	17	44
12	L	185/185 (100%)	170 (92%)	15 (8%)	11	33
12	Z	185/185 (100%)	171 (92%)	14 (8%)	13	36
13	M	199/199 (100%)	188 (94%)	11 (6%)	21	52
13	a	199/199 (100%)	187 (94%)	12 (6%)	19	48
14	N	162/162 (100%)	154 (95%)	8 (5%)	25	57
14	b	162/162 (100%)	153 (94%)	9 (6%)	21	51
15	1	1/1 (100%)	0	1 (100%)	0	0
15	2	1/1 (100%)	1 (100%)	0	100	100
15	3	1/1 (100%)	0	1 (100%)	0	0
15	4	1/1 (100%)	0	1 (100%)	0	0
15	5	1/1 (100%)	1 (100%)	0	100	100
15	6	1/1 (100%)	0	1 (100%)	0	0
All	All	5338/5338 (100%)	5021 (94%)	317 (6%)	19	49

5 of 317 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
13	M	106	ILE

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Mol	Chain	Res	Type
2	P	232	PHE
13	a	-4	ILE
13	M	191	ILE
1	O	136	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 202 such sidechains are listed below:

Mol	Chain	Res	Type
13	M	39	ASN
3	Q	83	ASN
12	Z	143	GLN
13	M	162	GLN
1	O	97	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

12 non-standard protein/DNA/RNA residues are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
15	LYH	1	2	15	7,8,9	1.14	0	4,8,10	0.98	0
15	LYH	6	2	15	7,8,9	1.08	0	4,8,10	1.46	1 (25%)
15	LYH	4	2	15	7,8,9	1.16	1 (14%)	4,8,10	1.03	0
15	LYH	2	2	15	7,8,9	1.13	0	4,8,10	1.36	1 (25%)
15	0JT	6	3	15,14	8,8,9	1.70	2 (25%)	7,9,11	1.00	0
15	LYH	5	2	15	7,8,9	1.13	1 (14%)	4,8,10	1.41	1 (25%)
15	0JT	3	3	15,14	8,8,9	1.60	2 (25%)	7,9,11	1.06	0
15	0JT	2	3	11,15	8,8,9	1.51	1 (12%)	7,9,11	1.14	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
15	0JT	1	3	8,15	8,8,9	1.35	2 (25%)	7,9,11	1.04	0
15	LYH	3	2	15	7,8,9	1.17	0	4,8,10	1.45	1 (25%)
15	0JT	5	3	11,15	8,8,9	1.65	2 (25%)	7,9,11	1.16	1 (14%)
15	0JT	4	3	8,15	8,8,9	1.34	1 (12%)	7,9,11	1.02	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
15	LYH	1	2	15	-	1/3/7/9	-
15	LYH	6	2	15	-	1/3/7/9	-
15	LYH	4	2	15	-	1/3/7/9	-
15	LYH	2	2	15	-	1/3/7/9	-
15	0JT	6	3	15,14	-	2/7/8/9	-
15	LYH	5	2	15	-	1/3/7/9	-
15	0JT	3	3	15,14	-	2/7/8/9	-
15	0JT	2	3	11,15	-	2/7/8/9	-
15	0JT	1	3	8,15	-	5/7/8/9	-
15	LYH	3	2	15	-	1/3/7/9	-
15	0JT	5	3	11,15	-	2/7/8/9	-
15	0JT	4	3	8,15	-	5/7/8/9	-

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

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
15	6	3	0JT	C19-C18	3.69	1.58	1.53
15	5	3	0JT	C19-C18	3.60	1.58	1.53
15	3	3	0JT	C19-C18	3.42	1.58	1.53
15	2	3	0JT	C19-C18	3.34	1.58	1.53
15	4	3	0JT	C19-C18	2.81	1.57	1.53

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
15	5	2	LYH	CE-CD-CG	-2.34	107.75	113.06
15	6	2	LYH	CE-CD-CG	-2.33	107.77	113.06
15	3	2	LYH	CE-CD-CG	-2.29	107.85	113.06
15	2	2	LYH	CE-CD-CG	-2.20	108.06	113.06

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
15	5	3	0JT	C17-C16-C14	2.01	118.02	112.29

There are no chirality outliers.

5 of 24 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
15	6	2	LYH	CG-CD-CE-NZ
15	2	2	LYH	CG-CD-CE-NZ
15	5	2	LYH	CG-CD-CE-NZ
15	1	3	0JT	C14-C16-C17-C18
15	3	2	LYH	CG-CD-CE-NZ

There are no ring outliers.

5 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
15	3	3	0JT	1	0
15	2	3	0JT	1	0
15	3	2	LYH	1	0
15	5	3	0JT	1	0
15	4	3	0JT	1	0

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	250/250 (100%)	-0.45	3 (1%) 79 73	35, 55, 84, 105	0
1	O	250/250 (100%)	-0.43	2 (0%) 86 81	36, 55, 84, 105	0
2	B	244/244 (100%)	-0.25	7 (2%) 51 41	39, 56, 94, 118	0
2	P	244/244 (100%)	-0.27	8 (3%) 46 36	39, 57, 95, 118	0
3	C	241/241 (100%)	-0.21	7 (2%) 51 41	40, 59, 110, 129	0
3	Q	241/241 (100%)	-0.21	11 (4%) 32 22	42, 61, 110, 129	0
4	D	242/242 (100%)	-0.24	7 (2%) 51 41	39, 60, 92, 125	0
4	R	242/242 (100%)	-0.19	8 (3%) 46 36	41, 61, 93, 126	0
5	E	233/233 (100%)	-0.30	6 (2%) 56 46	42, 62, 87, 111	0
5	S	233/233 (100%)	-0.27	4 (1%) 70 63	42, 62, 87, 111	0
6	F	244/244 (100%)	-0.36	8 (3%) 46 36	39, 56, 92, 106	0
6	T	244/244 (100%)	-0.26	5 (2%) 65 56	39, 56, 93, 105	0
7	G	243/243 (100%)	-0.47	3 (1%) 79 73	37, 53, 80, 112	0
7	U	243/243 (100%)	-0.43	3 (1%) 79 73	36, 53, 79, 112	0
8	H	222/222 (100%)	-0.55	1 (0%) 91 88	31, 48, 68, 98	0
8	V	222/222 (100%)	-0.50	3 (1%) 75 70	33, 48, 67, 98	0
9	I	204/204 (100%)	-0.56	1 (0%) 91 88	35, 50, 67, 84	0
9	W	204/204 (100%)	-0.53	1 (0%) 91 88	37, 50, 68, 84	0
10	J	198/198 (100%)	-0.45	3 (1%) 73 68	37, 51, 68, 124	0
10	X	198/198 (100%)	-0.45	4 (2%) 65 56	38, 51, 70, 124	0
11	K	212/212 (100%)	-0.46	1 (0%) 91 88	34, 51, 75, 82	0
11	Y	212/212 (100%)	-0.50	1 (0%) 91 88	35, 51, 74, 82	0
12	L	222/222 (100%)	-0.59	2 (0%) 84 80	33, 50, 71, 96	0
12	Z	222/222 (100%)	-0.55	2 (0%) 84 80	33, 50, 71, 95	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
13	M	233/233 (100%)	-0.61	1 (0%) 92 91	35, 49, 64, 69	0
13	a	233/233 (100%)	-0.55	1 (0%) 92 91	34, 48, 64, 69	0
14	N	196/196 (100%)	-0.54	0 100 100	33, 45, 67, 77	0
14	b	196/196 (100%)	-0.53	1 (0%) 91 88	33, 46, 68, 77	0
15	1	1/4 (25%)	-0.31	0 100 100	50, 50, 50, 50	0
15	2	1/4 (25%)	-0.20	0 100 100	52, 52, 52, 52	0
15	3	1/4 (25%)	-0.23	0 100 100	57, 57, 57, 57	0
15	4	1/4 (25%)	-0.39	0 100 100	51, 51, 51, 51	0
15	5	1/4 (25%)	0.04	0 100 100	51, 51, 51, 51	0
15	6	1/4 (25%)	0.06	0 100 100	57, 57, 57, 57	0
All	All	6374/6392 (99%)	-0.41	104 (1%) 72 66	31, 53, 84, 129	0

The worst 5 of 104 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
4	D	126	GLY	9.6
3	C	55	THR	8.8
4	D	128	SER	8.5
4	R	128	SER	8.4
10	X	196	GLN	7.8

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

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
15	LYH	1	2	9/10	0.88	0.16	47,48,50,50	0
15	LYH	6	2	9/10	0.89	0.19	46,48,50,50	0
15	LYH	3	2	9/10	0.92	0.22	46,50,52,52	0
15	LYH	4	2	9/10	0.93	0.20	47,48,50,50	0
15	LYH	5	2	9/10	0.94	0.18	46,47,49,49	0
15	OJT	1	3	9/10	0.94	0.14	40,45,48,49	0
15	LYH	2	2	9/10	0.94	0.16	45,46,47,47	0
15	OJT	6	3	9/10	0.95	0.14	38,44,47,50	0
15	OJT	3	3	9/10	0.96	0.14	39,45,50,52	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
15	OJT	4	3	9/10	0.96	0.17	40,43,48,49	0
15	OJT	5	3	9/10	0.97	0.13	42,44,47,47	0
15	OJT	2	3	9/10	0.98	0.13	39,43,46,46	0

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.