



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 04:15 PM UTC

PDB ID : 5W18 / pdb_00005w18
Title : Staphylococcus aureus ClpP in complex with (S)-N-((2R,6S,8aS,14aS,20S,23aS)-2,6-dimethyl-5,8,14,19,23-pentaoxooctadecahydro-1H,5H,14H,19H-pyrido[2,1-i]dipyrrolo[2,1-c:2',1'-l][1]oxa[4,7,10,13]tetraazacyclohexadecin-20-yl)-3-phenyl-2-(3-phenylureido)propanamide
Authors : Lee, R.E.; Griffith, E.C.
Deposited on : 2017-06-02
Resolution : 2.44 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

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

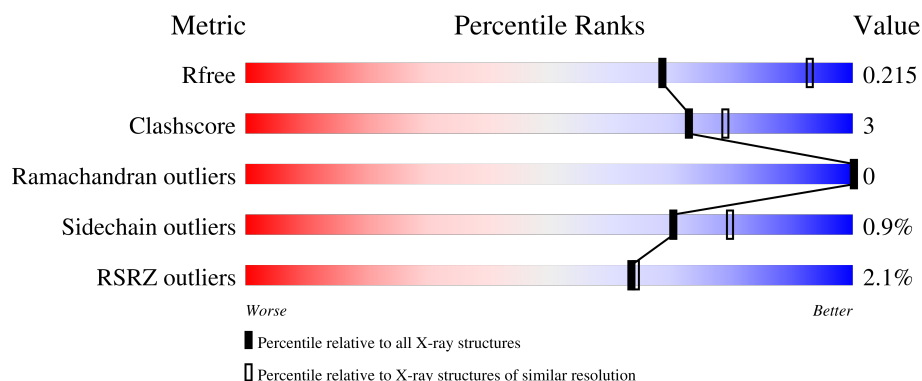
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

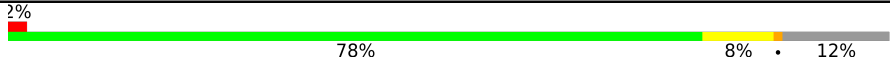
The reported resolution of this entry is 2.44 Å.

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	2340 (2.46-2.42)
Clashscore	190562	2400 (2.46-2.42)
Ramachandran outliers	187476	2379 (2.46-2.42)
Sidechain outliers	187428	2379 (2.46-2.42)
RSRZ outliers	180081	2340 (2.46-2.42)

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	203	
1	B	203	
1	C	203	
1	D	203	

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Mol	Chain	Length	Quality of chain
1	E	203	
1	F	203	
1	G	203	
1	I	203	
1	K	203	
1	L	203	
1	M	203	
1	N	203	
1	S	203	
1	T	203	
2	H	7	
2	J	7	
2	O	7	
2	P	7	
2	Q	7	
2	R	7	
2	U	7	
2	V	7	
2	X	7	
2	Y	7	
2	Z	7	
2	a	7	
2	b	7	
2	c	7	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 20484 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 ATP-dependent Clp protease proteolytic subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	179	Total	C	N	O	S	0	0	0
			1366	860	233	267	6			
1	B	180	Total	C	N	O	S	0	0	0
			1356	854	229	267	6			
1	C	179	Total	C	N	O	S	0	0	0
			1363	857	232	268	6			
1	D	180	Total	C	N	O	S	0	0	0
			1381	871	234	270	6			
1	E	178	Total	C	N	O	S	0	0	0
			1364	860	232	266	6			
1	F	179	Total	C	N	O	S	0	0	0
			1354	851	231	266	6			
1	G	180	Total	C	N	O	S	0	0	0
			1361	856	230	269	6			
1	I	179	Total	C	N	O	S	0	0	0
			1350	849	229	266	6			
1	K	180	Total	C	N	O	S	0	0	0
			1371	862	234	269	6			
1	L	180	Total	C	N	O	S	0	0	0
			1373	865	232	270	6			
1	M	179	Total	C	N	O	S	0	0	0
			1376	868	233	269	6			
1	N	179	Total	C	N	O	S	0	0	0
			1359	856	229	268	6			
1	S	178	Total	C	N	O	S	0	0	0
			1351	850	228	267	6			
1	T	178	Total	C	N	O	S	0	0	0
			1351	850	231	264	6			

There are 112 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	196	LEU	-	expression tag	UNP Q2G036

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Chain	Residue	Modelled	Actual	Comment	Reference
A	197	GLU	-	expression tag	UNP Q2G036
A	198	HIS	-	expression tag	UNP Q2G036
A	199	HIS	-	expression tag	UNP Q2G036
A	200	HIS	-	expression tag	UNP Q2G036
A	201	HIS	-	expression tag	UNP Q2G036
A	202	HIS	-	expression tag	UNP Q2G036
A	203	HIS	-	expression tag	UNP Q2G036
B	196	LEU	-	expression tag	UNP Q2G036
B	197	GLU	-	expression tag	UNP Q2G036
B	198	HIS	-	expression tag	UNP Q2G036
B	199	HIS	-	expression tag	UNP Q2G036
B	200	HIS	-	expression tag	UNP Q2G036
B	201	HIS	-	expression tag	UNP Q2G036
B	202	HIS	-	expression tag	UNP Q2G036
B	203	HIS	-	expression tag	UNP Q2G036
C	196	LEU	-	expression tag	UNP Q2G036
C	197	GLU	-	expression tag	UNP Q2G036
C	198	HIS	-	expression tag	UNP Q2G036
C	199	HIS	-	expression tag	UNP Q2G036
C	200	HIS	-	expression tag	UNP Q2G036
C	201	HIS	-	expression tag	UNP Q2G036
C	202	HIS	-	expression tag	UNP Q2G036
C	203	HIS	-	expression tag	UNP Q2G036
D	196	LEU	-	expression tag	UNP Q2G036
D	197	GLU	-	expression tag	UNP Q2G036
D	198	HIS	-	expression tag	UNP Q2G036
D	199	HIS	-	expression tag	UNP Q2G036
D	200	HIS	-	expression tag	UNP Q2G036
D	201	HIS	-	expression tag	UNP Q2G036
D	202	HIS	-	expression tag	UNP Q2G036
D	203	HIS	-	expression tag	UNP Q2G036
E	196	LEU	-	expression tag	UNP Q2G036
E	197	GLU	-	expression tag	UNP Q2G036
E	198	HIS	-	expression tag	UNP Q2G036
E	199	HIS	-	expression tag	UNP Q2G036
E	200	HIS	-	expression tag	UNP Q2G036
E	201	HIS	-	expression tag	UNP Q2G036
E	202	HIS	-	expression tag	UNP Q2G036
E	203	HIS	-	expression tag	UNP Q2G036
F	196	LEU	-	expression tag	UNP Q2G036
F	197	GLU	-	expression tag	UNP Q2G036
F	198	HIS	-	expression tag	UNP Q2G036

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Chain	Residue	Modelled	Actual	Comment	Reference
F	199	HIS	-	expression tag	UNP Q2G036
F	200	HIS	-	expression tag	UNP Q2G036
F	201	HIS	-	expression tag	UNP Q2G036
F	202	HIS	-	expression tag	UNP Q2G036
F	203	HIS	-	expression tag	UNP Q2G036
G	196	LEU	-	expression tag	UNP Q2G036
G	197	GLU	-	expression tag	UNP Q2G036
G	198	HIS	-	expression tag	UNP Q2G036
G	199	HIS	-	expression tag	UNP Q2G036
G	200	HIS	-	expression tag	UNP Q2G036
G	201	HIS	-	expression tag	UNP Q2G036
G	202	HIS	-	expression tag	UNP Q2G036
G	203	HIS	-	expression tag	UNP Q2G036
I	196	LEU	-	expression tag	UNP Q2G036
I	197	GLU	-	expression tag	UNP Q2G036
I	198	HIS	-	expression tag	UNP Q2G036
I	199	HIS	-	expression tag	UNP Q2G036
I	200	HIS	-	expression tag	UNP Q2G036
I	201	HIS	-	expression tag	UNP Q2G036
I	202	HIS	-	expression tag	UNP Q2G036
I	203	HIS	-	expression tag	UNP Q2G036
K	196	LEU	-	expression tag	UNP Q2G036
K	197	GLU	-	expression tag	UNP Q2G036
K	198	HIS	-	expression tag	UNP Q2G036
K	199	HIS	-	expression tag	UNP Q2G036
K	200	HIS	-	expression tag	UNP Q2G036
K	201	HIS	-	expression tag	UNP Q2G036
K	202	HIS	-	expression tag	UNP Q2G036
K	203	HIS	-	expression tag	UNP Q2G036
L	196	LEU	-	expression tag	UNP Q2G036
L	197	GLU	-	expression tag	UNP Q2G036
L	198	HIS	-	expression tag	UNP Q2G036
L	199	HIS	-	expression tag	UNP Q2G036
L	200	HIS	-	expression tag	UNP Q2G036
L	201	HIS	-	expression tag	UNP Q2G036
L	202	HIS	-	expression tag	UNP Q2G036
L	203	HIS	-	expression tag	UNP Q2G036
M	196	LEU	-	expression tag	UNP Q2G036
M	197	GLU	-	expression tag	UNP Q2G036
M	198	HIS	-	expression tag	UNP Q2G036
M	199	HIS	-	expression tag	UNP Q2G036
M	200	HIS	-	expression tag	UNP Q2G036

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Chain	Residue	Modelled	Actual	Comment	Reference
M	201	HIS	-	expression tag	UNP Q2G036
M	202	HIS	-	expression tag	UNP Q2G036
M	203	HIS	-	expression tag	UNP Q2G036
N	196	LEU	-	expression tag	UNP Q2G036
N	197	GLU	-	expression tag	UNP Q2G036
N	198	HIS	-	expression tag	UNP Q2G036
N	199	HIS	-	expression tag	UNP Q2G036
N	200	HIS	-	expression tag	UNP Q2G036
N	201	HIS	-	expression tag	UNP Q2G036
N	202	HIS	-	expression tag	UNP Q2G036
N	203	HIS	-	expression tag	UNP Q2G036
S	196	LEU	-	expression tag	UNP Q2G036
S	197	GLU	-	expression tag	UNP Q2G036
S	198	HIS	-	expression tag	UNP Q2G036
S	199	HIS	-	expression tag	UNP Q2G036
S	200	HIS	-	expression tag	UNP Q2G036
S	201	HIS	-	expression tag	UNP Q2G036
S	202	HIS	-	expression tag	UNP Q2G036
S	203	HIS	-	expression tag	UNP Q2G036
T	196	LEU	-	expression tag	UNP Q2G036
T	197	GLU	-	expression tag	UNP Q2G036
T	198	HIS	-	expression tag	UNP Q2G036
T	199	HIS	-	expression tag	UNP Q2G036
T	200	HIS	-	expression tag	UNP Q2G036
T	201	HIS	-	expression tag	UNP Q2G036
T	202	HIS	-	expression tag	UNP Q2G036
T	203	HIS	-	expression tag	UNP Q2G036

- Molecule 2 is a protein called 9V7-PHE-SER-PRO-YCP-ALA-MP8.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	H	7	Total	C	N	O	0	0	0
			54	39	7	8			
2	J	7	Total	C	N	O	0	0	0
			54	39	7	8			
2	O	7	Total	C	N	O	0	0	0
			54	39	7	8			
2	P	7	Total	C	N	O	0	0	0
			54	39	7	8			
2	Q	7	Total	C	N	O	0	0	0
			54	39	7	8			
2	R	7	Total	C	N	O	0	0	0
			54	39	7	8			

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	U	7	Total	C	N	O	0	0	0
			54	39	7	8			
2	V	7	Total	C	N	O	0	0	0
			54	39	7	8			
2	X	7	Total	C	N	O	0	0	0
			54	39	7	8			
2	Y	7	Total	C	N	O	0	0	0
			54	39	7	8			
2	Z	7	Total	C	N	O	0	0	0
			54	39	7	8			
2	a	7	Total	C	N	O	0	0	0
			54	39	7	8			
2	b	7	Total	C	N	O	0	0	0
			54	39	7	8			
2	c	7	Total	C	N	O	0	0	0
			54	39	7	8			

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	33	Total	O	0	0
			33	33		
3	B	40	Total	O	0	0
			40	40		
3	C	56	Total	O	0	0
			56	56		
3	D	45	Total	O	0	0
			45	45		
3	E	65	Total	O	0	0
			65	65		
3	F	54	Total	O	0	0
			54	54		
3	G	45	Total	O	0	0
			45	45		
3	I	26	Total	O	0	0
			26	26		
3	K	41	Total	O	0	0
			41	41		
3	L	62	Total	O	0	0
			62	62		
3	M	57	Total	O	0	0
			57	57		

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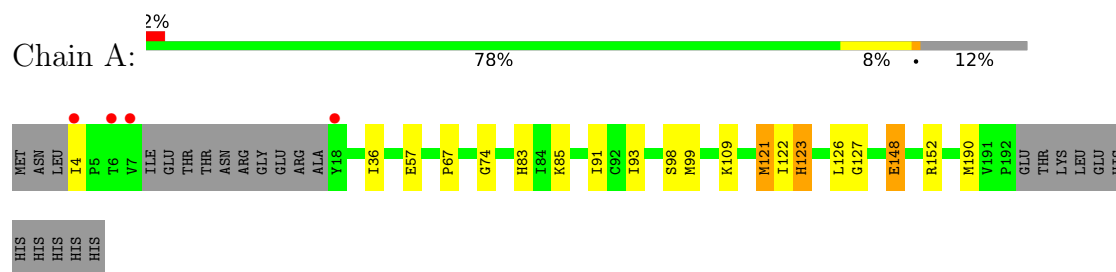
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	N	62	Total 62	O 62	0	0
3	S	33	Total 33	O 33	0	0
3	T	30	Total 30	O 30	0	0
3	Y	1	Total 1	O 1	0	0
3	Z	1	Total 1	O 1	0	0
3	a	1	Total 1	O 1	0	0

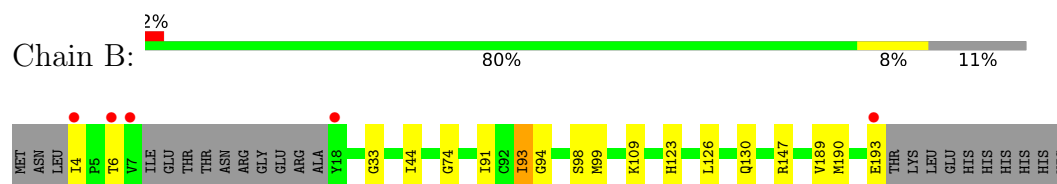
3 Residue-property plots

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.

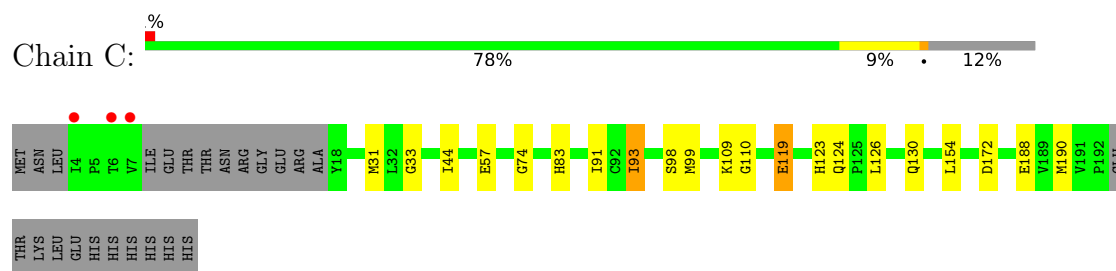
- Molecule 1: ATP-dependent Clp protease proteolytic subunit



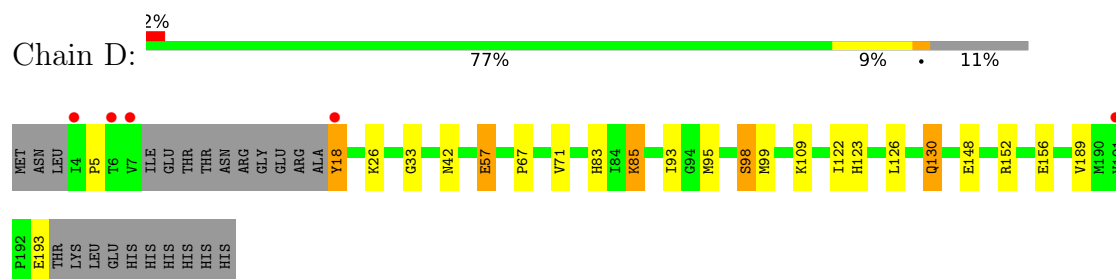
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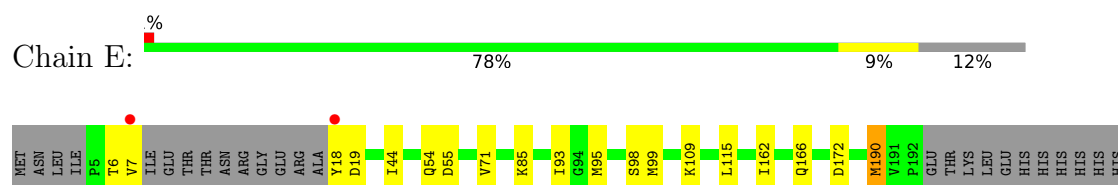
- Molecule 1: ATP-dependent Clp protease proteolytic subunit



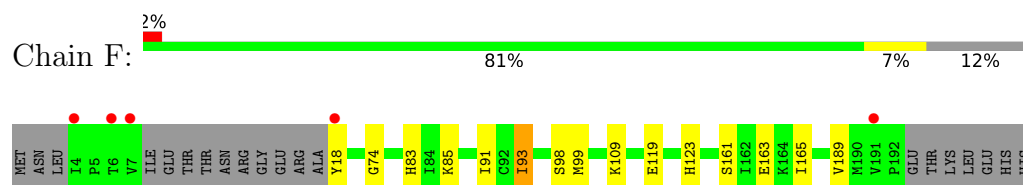
- Molecule 1: ATP-dependent Clp protease proteolytic subunit



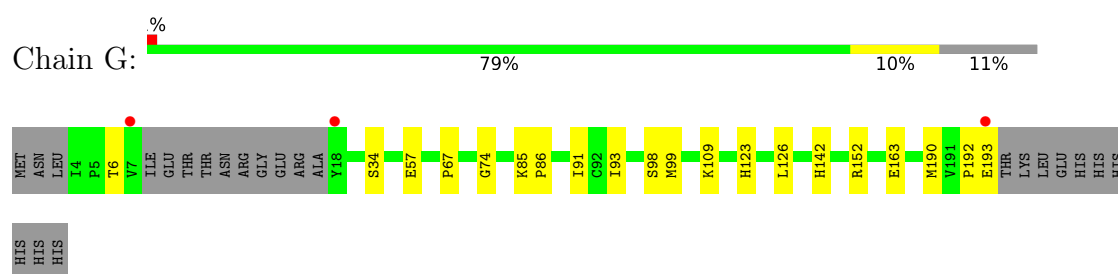
- Molecule 1: ATP-dependent Clp protease proteolytic subunit



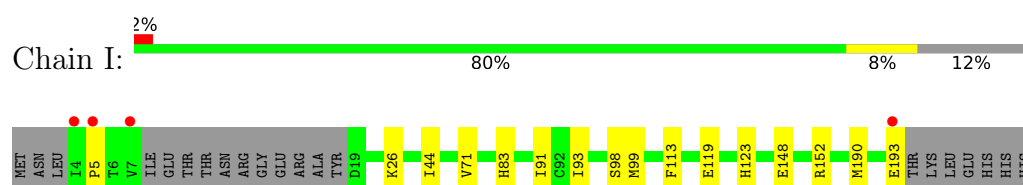
- Molecule 1: ATP-dependent Clp protease proteolytic subunit



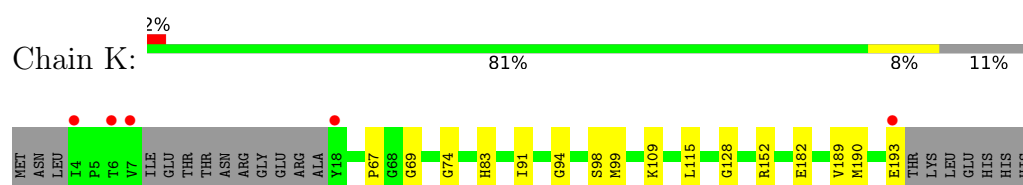
- Molecule 1: ATP-dependent Clp protease proteolytic subunit



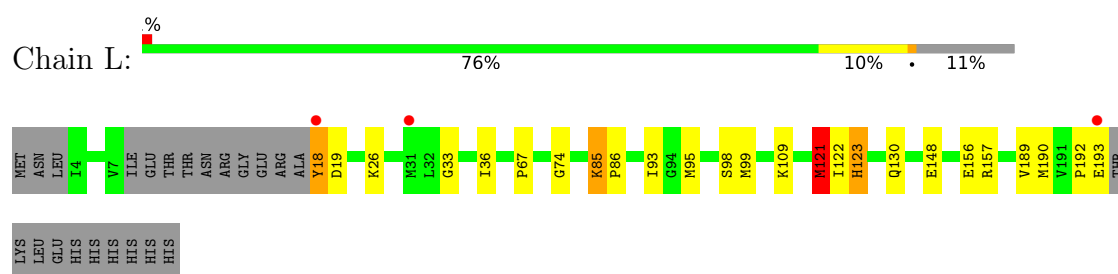
- Molecule 1: ATP-dependent Clp protease proteolytic subunit



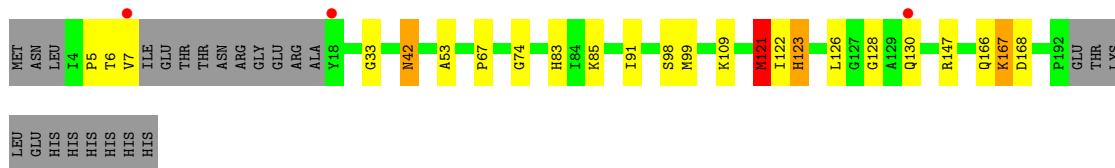
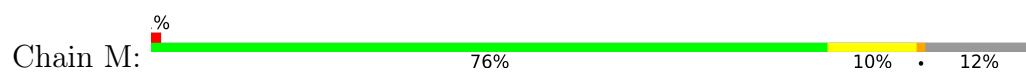
- Molecule 1: ATP-dependent Clp protease proteolytic subunit



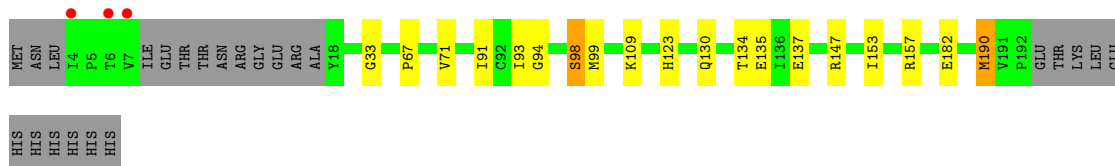
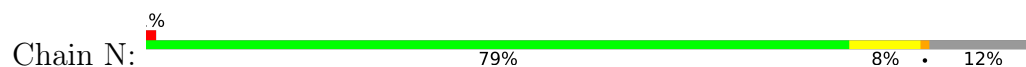
- Molecule 1: ATP-dependent Clp protease proteolytic subunit



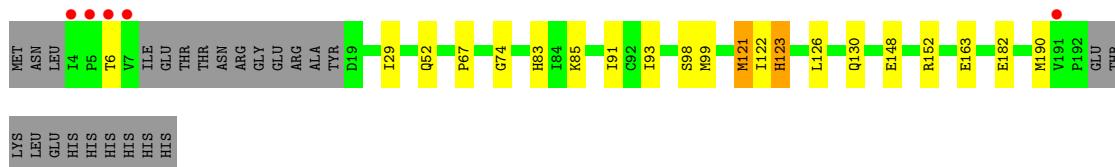
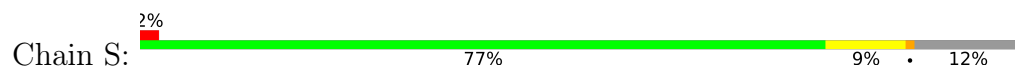
- Molecule 1: ATP-dependent Clp protease proteolytic subunit



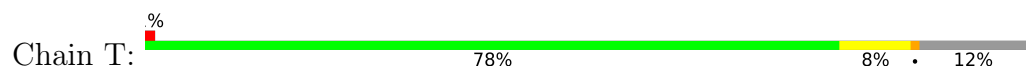
- Molecule 1: ATP-dependent Clp protease proteolytic subunit



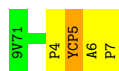
- Molecule 1: ATP-dependent Clp protease proteolytic subunit



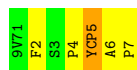
- Molecule 1: ATP-dependent Clp protease proteolytic subunit



- Molecule 2: 9V7-PHE-SER-PRO-YCP-ALA-MP8



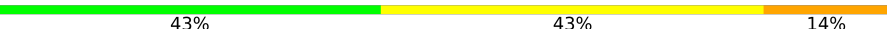
- Molecule 2: 9V7-PHE-SER-PRO-YCP-ALA-MP8



● Molecule 2: 9V7-PHE-SER-PRO-YCP-ALA-MP8

Chain O:  43% 29% 29%

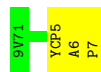
● Molecule 2: 9V7-PHE-SER-PRO-YCP-ALA-MP8

Chain P:  43% 43% 14%

● Molecule 2: 9V7-PHE-SER-PRO-YCP-ALA-MP8

Chain Q:  43% 29% 29%

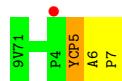
● Molecule 2: 9V7-PHE-SER-PRO-YCP-ALA-MP8

Chain R:  57% 43%

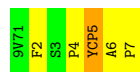
● Molecule 2: 9V7-PHE-SER-PRO-YCP-ALA-MP8

Chain U:  29% 57% 14%

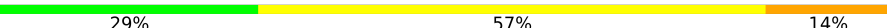
● Molecule 2: 9V7-PHE-SER-PRO-YCP-ALA-MP8

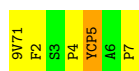
Chain V:  14% 57% 29% 14%

● Molecule 2: 9V7-PHE-SER-PRO-YCP-ALA-MP8

Chain X:  29% 57% 14%

● Molecule 2: 9V7-PHE-SER-PRO-YCP-ALA-MP8

Chain Y:  29% 57% 14%



- Molecule 2: 9V7-PHE-SER-PRO-YCP-ALA-MP8



- Molecule 2: 9V7-PHE-SER-PRO-YCP-ALA-MP8



- Molecule 2: 9V7-PHE-SER-PRO-YCP-ALA-MP8



- Molecule 2: 9V7-PHE-SER-PRO-YCP-ALA-MP8



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	94.92Å 126.38Å 146.06Å 90.00° 93.42° 90.00°	Depositor
Resolution (Å)	50.00 – 2.44 50.00 – 2.44	Depositor EDS
% Data completeness (in resolution range)	99.5 (50.00-2.44) 99.8 (50.00-2.44)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.26 (at 2.42Å)	Xtriage
Refinement program	REFMAC 5.8.0158	Depositor
R, R_{free}	0.189 , 0.210 0.193 , 0.215	Depositor DCC
R_{free} test set	6556 reflections (5.10%)	wwPDB-VP
Wilson B-factor (Å ²)	33.6	Xtriage
Anisotropy	0.167	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 22.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	20484	wwPDB-VP
Average B, all atoms (Å ²)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.76% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MP8, 9V7, YCP

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	1.52	9/1383 (0.7%)	1.14	3/1867 (0.2%)
1	B	1.47	12/1373 (0.9%)	1.10	1/1859 (0.1%)
1	C	1.57	11/1380 (0.8%)	1.11	2/1865 (0.1%)
1	D	1.70	17/1399 (1.2%)	1.16	0/1890
1	E	1.62	8/1382 (0.6%)	1.15	4/1866 (0.2%)
1	F	1.58	9/1371 (0.7%)	1.14	2/1855 (0.1%)
1	G	1.52	9/1378 (0.7%)	1.09	1/1864 (0.1%)
1	I	1.46	8/1367 (0.6%)	1.08	0/1849
1	K	1.58	11/1388 (0.8%)	1.15	0/1875
1	L	1.69	16/1391 (1.2%)	1.12	1/1882 (0.1%)
1	M	1.64	13/1394 (0.9%)	1.15	0/1883
1	N	1.56	10/1376 (0.7%)	1.13	3/1861 (0.2%)
1	S	1.46	10/1368 (0.7%)	1.10	0/1850
1	T	1.46	10/1368 (0.7%)	1.12	2/1848 (0.1%)
2	H	1.26	0/29	0.89	0/37
2	J	1.24	0/29	1.07	0/37
2	O	1.19	0/29	1.13	0/37
2	P	1.33	0/29	1.32	1/37 (2.7%)
2	Q	1.73	1/29 (3.4%)	1.34	0/37
2	R	1.40	0/29	0.97	0/37
2	U	1.04	0/29	0.97	0/37
2	V	1.29	0/29	1.19	0/37
2	X	1.23	0/29	0.92	0/37
2	Y	1.25	0/29	1.09	0/37
2	Z	1.39	1/29 (3.4%)	1.29	0/37
2	a	1.54	1/29 (3.4%)	0.79	0/37
2	b	1.22	0/29	0.91	0/37
2	c	1.33	0/29	1.15	0/37
All	All	1.56	156/19724 (0.8%)	1.12	20/26632 (0.1%)

All (156) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	I	5	PRO	C-O	10.38	1.37	1.23
1	G	193	GLU	N-CA	9.48	1.64	1.46
1	C	109	LYS	C-O	-9.30	1.12	1.23
1	D	109	LYS	C-O	-9.28	1.12	1.23
1	M	123	HIS	C-O	-9.21	1.13	1.23
1	T	123	HIS	C-O	-9.14	1.13	1.23
1	G	109	LYS	C-O	-9.04	1.13	1.23
1	B	123	HIS	C-O	-9.04	1.14	1.24
1	C	123	HIS	C-O	-9.03	1.14	1.23
1	D	5	PRO	C-O	8.64	1.33	1.23
1	L	123	HIS	C-O	-8.47	1.14	1.23
1	K	109	LYS	C-O	-8.47	1.13	1.23
1	L	109	LYS	C-O	-8.46	1.13	1.23
1	M	98	SER	C-O	-8.31	1.16	1.24
1	K	182	GLU	CD-OE1	8.19	1.41	1.25
1	D	123	HIS	C-O	-8.03	1.15	1.23
1	I	193	GLU	N-CA	8.03	1.61	1.46
1	F	109	LYS	C-O	-7.88	1.14	1.23
1	A	127	GLY	C-O	-7.87	1.15	1.23
1	N	109	LYS	C-O	-7.82	1.14	1.23
1	L	98	SER	C-O	-7.79	1.17	1.24
1	C	119	GLU	CD-OE1	7.46	1.39	1.25
1	D	57	GLU	CD-OE2	7.41	1.39	1.25
1	M	121	MET	SD-CE	-7.33	1.61	1.79
1	D	193	GLU	N-CA	7.32	1.60	1.46
1	G	57	GLU	CD-OE2	7.30	1.39	1.25
1	I	123	HIS	C-O	-7.29	1.16	1.24
1	L	189	VAL	C-O	-7.27	1.16	1.24
1	T	98	SER	C-O	-7.25	1.18	1.24
1	B	193	GLU	N-CA	7.07	1.59	1.46
1	A	123	HIS	C-O	-7.04	1.16	1.23
2	Q	4	PRO	N-CD	6.97	1.57	1.47
1	M	109	LYS	C-O	-6.94	1.15	1.23
1	C	98	SER	C-O	-6.88	1.18	1.24
1	E	98	SER	C-O	-6.88	1.18	1.24
1	T	163	GLU	CD-OE1	6.88	1.38	1.25
1	S	98	SER	C-O	-6.85	1.18	1.24
1	L	85	LYS	C-O	-6.75	1.18	1.24
1	B	193	GLU	C-O	6.74	1.37	1.23
1	L	121	MET	SD-CE	-6.65	1.62	1.79
1	G	152	ARG	CZ-NH2	6.65	1.42	1.33
1	D	98	SER	C-O	-6.62	1.18	1.24
1	E	172	ASP	C-O	-6.61	1.15	1.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	L	19	ASP	CG-OD2	6.59	1.37	1.25
1	D	152	ARG	CD-NE	6.55	1.55	1.46
1	E	54	GLN	CD-OE1	6.50	1.35	1.23
1	B	130	GLN	CD-OE1	6.47	1.35	1.23
1	S	123	HIS	C-O	-6.46	1.16	1.23
1	N	130	GLN	CD-OE1	6.45	1.35	1.23
1	S	152	ARG	CZ-NH2	6.39	1.41	1.33
1	M	85	LYS	C-O	-6.39	1.18	1.24
1	K	193	GLU	N-CA	6.38	1.58	1.46
1	D	130	GLN	CD-NE2	6.34	1.46	1.33
1	M	42	ASN	CG-OD1	6.33	1.35	1.23
1	K	98	SER	C-O	-6.32	1.18	1.24
1	B	126	LEU	C-O	-6.30	1.16	1.23
1	N	98	SER	C-O	-6.26	1.18	1.24
1	F	98	SER	C-O	-6.26	1.18	1.24
1	E	95	MET	SD-CE	6.05	1.94	1.79
1	F	85	LYS	C-O	-6.04	1.18	1.24
1	G	6	THR	C-O	6.04	1.31	1.23
1	F	161	SER	CA-C	6.02	1.60	1.52
1	A	36	ILE	C-O	-6.02	1.16	1.23
1	D	126	LEU	C-O	-6.01	1.16	1.23
1	I	148	GLU	CD-OE1	6.01	1.36	1.25
1	E	6	THR	CA-C	5.99	1.59	1.52
1	D	85	LYS	C-O	-5.98	1.18	1.24
1	D	42	ASN	CG-OD1	5.97	1.34	1.23
1	K	193	GLU	C-O	5.97	1.35	1.23
1	L	193	GLU	N-CA	5.94	1.57	1.46
1	N	94	GLY	N-CA	5.94	1.53	1.45
1	C	130	GLN	CD-NE2	5.92	1.45	1.33
1	D	67	PRO	C-O	-5.89	1.16	1.24
1	D	122	ILE	C-O	-5.89	1.18	1.24
1	F	123	HIS	C-O	-5.88	1.17	1.23
1	A	98	SER	C-O	-5.84	1.19	1.24
1	M	5	PRO	C-O	5.83	1.31	1.23
1	A	126	LEU	C-O	-5.82	1.16	1.23
1	E	6	THR	C-O	5.79	1.31	1.24
1	S	126	LEU	C-O	-5.79	1.16	1.23
2	a	4	PRO	N-CD	5.77	1.55	1.47
1	S	130	GLN	CD-OE1	5.74	1.34	1.23
1	A	109	LYS	C-O	-5.72	1.17	1.23
1	G	123	HIS	C-O	-5.71	1.17	1.23
1	D	189	VAL	C-O	-5.67	1.17	1.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	I	98	SER	C-O	-5.64	1.19	1.24
1	L	156	GLU	CD-OE1	5.63	1.36	1.25
1	M	33	GLY	C-O	-5.62	1.16	1.23
1	K	69	GLY	C-O	-5.61	1.19	1.23
1	I	152	ARG	CZ-NH2	5.61	1.40	1.33
1	L	67	PRO	C-O	-5.59	1.17	1.24
1	M	126	LEU	C-O	-5.57	1.17	1.23
1	T	95	MET	SD-CE	5.56	1.93	1.79
1	M	130	GLN	CD-OE1	5.56	1.34	1.23
1	B	98	SER	C-O	-5.53	1.19	1.24
1	C	172	ASP	C-O	-5.52	1.17	1.23
1	B	94	GLY	N-CA	5.51	1.53	1.45
1	A	148	GLU	CD-OE2	5.50	1.35	1.25
1	I	193	GLU	C-O	5.50	1.34	1.23
1	S	148	GLU	CD-OE2	5.50	1.35	1.25
1	G	193	GLU	C-O	5.45	1.34	1.23
1	D	156	GLU	CD-OE2	5.44	1.35	1.25
1	F	189	VAL	N-CA	-5.42	1.39	1.46
1	C	188	GLU	CD-OE1	5.42	1.35	1.25
1	B	33	GLY	C-O	-5.42	1.16	1.23
1	N	67	PRO	C-O	-5.41	1.17	1.24
1	K	152	ARG	CD-NE	5.40	1.53	1.46
1	T	36	ILE	C-O	-5.39	1.17	1.23
1	K	94	GLY	N-CA	5.38	1.53	1.45
1	E	109	LYS	C-O	-5.37	1.17	1.23
1	F	165	ILE	CA-C	5.36	1.59	1.52
1	N	182	GLU	C-O	-5.36	1.17	1.24
1	M	128	GLY	C-O	-5.34	1.18	1.24
1	N	123	HIS	C-O	-5.34	1.18	1.23
1	E	19	ASP	CG-OD2	5.33	1.35	1.25
1	B	109	LYS	C-O	-5.31	1.17	1.23
1	F	163	GLU	CD-OE1	5.29	1.35	1.25
1	F	119	GLU	CD-OE2	5.28	1.35	1.25
1	T	109	LYS	C-O	-5.27	1.17	1.23
1	A	67	PRO	C-O	-5.26	1.17	1.24
1	M	6	THR	C-O	5.25	1.30	1.24
1	A	4	ILE	N-CA	5.23	1.56	1.46
1	C	126	LEU	C-O	-5.22	1.17	1.23
1	L	193	GLU	C-O	5.22	1.33	1.23
1	C	57	GLU	CD-OE2	5.22	1.35	1.25
1	K	189	VAL	C-O	-5.21	1.18	1.24
1	L	36	ILE	C-O	-5.21	1.17	1.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	33	GLY	C-O	-5.21	1.16	1.23
1	I	119	GLU	CD-OE2	5.21	1.35	1.25
1	G	126	LEU	C-O	-5.20	1.17	1.23
1	M	67	PRO	C-O	-5.17	1.17	1.24
1	S	182	GLU	C-O	-5.17	1.18	1.24
1	T	157	ARG	CD-NE	-5.17	1.39	1.46
1	T	172	ASP	C-O	-5.17	1.17	1.23
2	Z	4	PRO	N-CD	5.16	1.54	1.47
1	S	6	THR	C-O	5.16	1.30	1.24
1	G	67	PRO	C-O	-5.15	1.17	1.24
1	S	163	GLU	CD-OE1	5.14	1.35	1.25
1	K	67	PRO	C-O	-5.13	1.17	1.24
1	N	135	GLU	C-O	-5.13	1.18	1.24
1	L	148	GLU	CD-OE2	5.12	1.35	1.25
1	L	95	MET	CG-SD	5.11	1.93	1.80
1	N	137	GLU	CD-OE1	5.11	1.35	1.25
1	T	161	SER	CA-C	5.11	1.60	1.53
1	K	128	GLY	C-O	-5.10	1.18	1.24
1	S	67	PRO	C-O	-5.10	1.17	1.24
1	L	157	ARG	NE-CZ	5.10	1.38	1.33
1	T	119	GLU	CD-OE2	5.09	1.35	1.25
1	C	33	GLY	C-O	-5.08	1.17	1.23
1	D	148	GLU	CD-OE1	5.08	1.35	1.25
1	B	6	THR	CA-C	5.08	1.58	1.52
1	B	189	VAL	C-O	-5.07	1.18	1.24
1	C	110	GLY	C-O	-5.06	1.17	1.24
1	L	130	GLN	CD-OE1	5.06	1.33	1.23
1	N	33	GLY	C-O	-5.05	1.17	1.23
1	B	147	ARG	CZ-NH1	5.03	1.39	1.32

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	190	MET	CG-SD-CE	-6.43	86.75	100.90
1	L	192	PRO	CA-C-O	-6.21	113.95	121.03
1	G	192	PRO	CA-C-O	-5.88	114.32	121.03
1	F	93	ILE	O-C-N	-5.84	117.04	123.18
1	N	93	ILE	O-C-N	-5.58	117.32	123.18
1	A	93	ILE	O-C-N	-5.51	117.40	123.18
1	T	93	ILE	O-C-N	-5.48	117.42	123.18
1	E	55	ASP	CB-CG-OD2	-5.47	105.82	118.40
1	E	85	LYS	CG-CD-CE	-5.35	99.00	111.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	T	190	MET	CB-CA-C	-5.32	101.56	110.29
1	C	124	GLN	CB-CA-C	5.29	117.05	109.11
1	B	93	ILE	O-C-N	-5.13	117.79	123.18
1	E	7	VAL	CA-CB-CG2	5.13	119.12	110.40
1	F	18	TYR	N-CA-CB	-5.08	101.86	110.50
1	N	134	THR	OG1-CB-CG2	-5.08	99.13	109.30
1	A	190	MET	CB-CA-C	-5.08	101.96	110.29
1	N	190	MET	CB-CA-C	-5.06	102.66	110.26
2	P	3	SER	CA-CB-OG	-5.03	101.04	111.10
1	C	93	ILE	O-C-N	-5.02	117.91	123.18
1	E	93	ILE	O-C-N	-5.02	117.84	123.26

There are no chirality outliers.

There are no planarity outliers.

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	1366	0	1374	10	0
1	B	1356	0	1343	7	0
1	C	1363	0	1360	11	0
1	D	1381	0	1387	11	0
1	E	1364	0	1371	10	0
1	F	1354	0	1343	4	0
1	G	1361	0	1349	11	0
1	I	1350	0	1336	12	0
1	K	1371	0	1371	10	0
1	L	1373	0	1365	10	0
1	M	1376	0	1385	10	0
1	N	1359	0	1356	7	0
1	S	1351	0	1345	13	0
1	T	1351	0	1350	10	0
2	H	54	0	43	2	0
2	J	54	0	43	3	0
2	O	54	0	43	3	0
2	P	54	0	43	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	Q	54	0	43	6	0
2	R	54	0	43	1	0
2	U	54	0	43	3	0
2	V	54	0	43	2	0
2	X	54	0	43	3	0
2	Y	54	0	43	3	0
2	Z	54	0	43	1	0
2	a	54	0	43	2	0
2	b	54	0	43	4	0
2	c	54	0	43	4	0
3	A	33	0	0	1	0
3	B	40	0	0	0	0
3	C	56	0	0	1	0
3	D	45	0	0	2	0
3	E	65	0	0	2	0
3	F	54	0	0	0	0
3	G	45	0	0	2	0
3	I	26	0	0	1	0
3	K	41	0	0	0	0
3	L	62	0	0	0	0
3	M	57	0	0	0	0
3	N	62	0	0	3	0
3	S	33	0	0	0	0
3	T	30	0	0	0	0
3	Y	1	0	0	0	0
3	Z	1	0	0	0	0
3	a	1	0	0	0	0
All	All	20484	0	19637	128	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 3.

All (128) 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)
1:D:95:MET:HG3	3:D:301:HOH:O	1.59	1.00
1:L:190:MET:HE1	2:Y:2:PHE:HB3	1.49	0.93
1:S:190:MET:CE	1:T:83:HIS:CE1	2.55	0.89
1:A:83:HIS:CE1	1:G:190:MET:HE3	2.09	0.87
1:A:148:GLU:OE1	1:A:152:ARG:NH2	2.09	0.84
1:D:71:VAL:HG22	1:D:99:MET:HE3	1.58	0.83

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:83:HIS:CE1	1:G:190:MET:CE	2.62	0.82
1:C:190:MET:HE1	2:O:2:PHE:HB3	1.66	0.77
1:S:190:MET:HE3	1:T:83:HIS:CE1	2.20	0.76
1:C:190:MET:CE	1:D:83:HIS:CE1	2.69	0.76
1:E:71:VAL:HG22	1:E:99:MET:HE3	1.68	0.74
1:G:190:MET:HE1	2:U:2:PHE:HB3	1.67	0.74
1:B:190:MET:HE2	1:C:83:HIS:CE1	2.24	0.73
1:L:190:MET:CE	1:M:83:HIS:CE1	2.72	0.72
1:T:190:MET:HE1	2:c:2:PHE:HB3	1.70	0.72
1:M:147:ARG:NH1	1:M:166:GLN:OE1	2.22	0.72
1:S:190:MET:HE1	2:b:2:PHE:HB3	1.72	0.72
1:N:71:VAL:HG22	1:N:99:MET:HE3	1.71	0.72
1:I:71:VAL:HG22	1:I:99:MET:HE3	1.72	0.70
1:B:190:MET:CE	1:C:83:HIS:CE1	2.76	0.68
1:D:95:MET:CE	3:E:301:HOH:O	2.42	0.68
1:I:190:MET:CE	1:K:83:HIS:NE2	2.58	0.66
1:D:95:MET:HE3	3:E:301:HOH:O	1.97	0.65
1:S:52:GLN:HE21	1:S:85:LYS:H	1.42	0.65
1:E:190:MET:HE1	2:Q:2:PHE:HB3	1.79	0.65
1:K:74:GLY:HA3	1:K:99:MET:HE2	1.80	0.64
1:T:153:ILE:O	1:T:157:ARG:HG2	1.99	0.63
1:M:74:GLY:HA3	1:M:99:MET:HE2	1.81	0.63
1:D:57:GLU:OE2	1:D:85:LYS:NZ	2.32	0.62
1:T:74:GLY:HA3	1:T:99:MET:HE2	1.82	0.62
1:E:190:MET:HE3	1:F:83:HIS:CE1	2.35	0.62
1:L:190:MET:HE3	1:M:83:HIS:CE1	2.35	0.61
1:D:71:VAL:HG22	1:D:99:MET:CE	2.32	0.58
1:S:121:MET:HE3	1:S:122:ILE:C	2.28	0.58
1:E:190:MET:CE	2:Q:2:PHE:HB3	2.34	0.58
1:K:190:MET:HE1	2:X:2:PHE:CG	2.39	0.58
1:N:71:VAL:HG22	1:N:99:MET:CE	2.34	0.57
1:C:190:MET:HE2	1:D:83:HIS:CE1	2.40	0.57
1:B:74:GLY:HA3	1:B:99:MET:HE2	1.85	0.57
1:M:167:LYS:NZ	1:M:168:ASP:OD1	2.37	0.57
1:E:190:MET:HE1	2:Q:2:PHE:CG	2.40	0.56
1:I:190:MET:HE1	1:K:83:HIS:NE2	2.20	0.56
1:E:190:MET:HE1	2:Q:2:PHE:CB	2.34	0.56
1:L:121:MET:HE1	1:L:123:HIS:CG	2.40	0.56
1:K:115:LEU:HD21	1:K:190:MET:HE2	1.88	0.56
1:I:71:VAL:HG22	1:I:99:MET:CE	2.36	0.55
1:K:115:LEU:CD2	1:K:190:MET:HE2	2.37	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:74:GLY:HA3	1:F:99:MET:HE2	1.89	0.55
1:A:121:MET:HE3	1:A:122:ILE:C	2.32	0.55
1:A:57:GLU:OE2	1:A:85:LYS:NZ	2.40	0.54
1:E:162:ILE:O	1:E:166:GLN:HG3	2.07	0.54
1:S:74:GLY:HA3	1:S:99:MET:HE2	1.90	0.54
1:G:74:GLY:HA3	1:G:99:MET:HE2	1.88	0.54
1:I:190:MET:HE2	1:K:83:HIS:NE2	2.23	0.54
1:C:74:GLY:HA3	1:C:99:MET:HE2	1.91	0.53
1:E:71:VAL:HG22	1:E:99:MET:CE	2.37	0.53
1:M:121:MET:HE3	1:M:122:ILE:C	2.33	0.53
1:A:74:GLY:HA3	1:A:99:MET:HE2	1.90	0.53
3:I:321:HOH:O	1:K:83:HIS:HD2	1.92	0.52
1:G:142:HIS:HD2	3:G:345:HOH:O	1.93	0.51
1:S:121:MET:HE1	1:S:123:HIS:CG	2.45	0.51
1:C:119:GLU:OE2	3:C:301:HOH:O	2.19	0.51
1:S:52:GLN:NE2	1:S:85:LYS:H	2.09	0.51
1:I:83:HIS:NE2	1:T:190:MET:HE2	2.26	0.50
1:M:53:ALA:HB2	2:Y:1:9V7:C4	2.41	0.50
1:N:153:ILE:O	1:N:157:ARG:HG2	2.11	0.50
1:A:148:GLU:CD	1:A:152:ARG:NH2	2.70	0.49
1:D:98:SER:HB3	3:D:342:HOH:O	2.11	0.49
1:M:121:MET:HE1	1:M:123:HIS:CG	2.49	0.48
1:A:121:MET:HE1	1:A:123:HIS:CG	2.49	0.47
1:L:74:GLY:HA3	1:L:99:MET:HE2	1.97	0.47
1:I:83:HIS:NE2	1:T:190:MET:CE	2.79	0.46
1:S:190:MET:HE2	1:T:83:HIS:CE1	2.48	0.46
1:E:115:LEU:CD2	1:E:190:MET:HE2	2.46	0.45
1:B:4:ILE:O	1:B:4:ILE:HG13	2.16	0.45
1:G:98:SER:HB3	3:G:332:HOH:O	2.16	0.45
1:L:121:MET:HE3	1:L:122:ILE:C	2.42	0.45
1:G:93:ILE:O	1:G:93:ILE:HG13	2.16	0.45
3:N:348:HOH:O	1:S:83:HIS:HD2	2.00	0.44
1:D:18:TYR:OH	1:D:26:LYS:HD3	2.17	0.44
2:U:4:PRO:HA	2:U:5:YCP:HA	1.77	0.44
1:C:91:ILE:HD11	2:O:6:ALA:HB1	2.00	0.44
1:I:91:ILE:HD11	2:V:6:ALA:HB1	1.99	0.44
1:K:91:ILE:HD11	2:X:6:ALA:HB1	1.99	0.44
2:J:4:PRO:HA	2:J:5:YCP:HA	1.78	0.44
1:A:91:ILE:HD11	2:H:6:ALA:HB1	1.98	0.44
1:C:93:ILE:O	1:C:93:ILE:HG13	2.16	0.44
1:D:93:ILE:O	1:D:93:ILE:HG13	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:91:ILE:HD11	2:R:6:ALA:HB1	1.99	0.44
2:O:5:YCP:HA	2:O:7:MP8:O	2.18	0.43
1:G:91:ILE:HD11	2:U:6:ALA:HB1	2.00	0.43
2:a:4:PRO:HA	2:a:5:YCP:HA	1.76	0.43
1:T:91:ILE:HD11	2:c:6:ALA:HB1	1.99	0.43
1:N:91:ILE:HD11	2:a:6:ALA:HB1	2.00	0.43
1:I:113:PHE:CZ	2:V:5:YCP:HD	2.52	0.43
1:B:93:ILE:O	1:B:93:ILE:HG13	2.18	0.43
1:S:91:ILE:HD11	2:b:6:ALA:HB1	2.01	0.43
1:L:93:ILE:O	1:L:93:ILE:HG13	2.19	0.43
1:S:29:ILE:HD11	2:b:7:MP8:HE	2.01	0.43
1:E:190:MET:HE1	2:Q:2:PHE:CD2	2.54	0.42
2:P:4:PRO:HA	2:P:5:YCP:HA	1.78	0.42
1:I:190:MET:HE2	1:K:83:HIS:CD2	2.54	0.42
1:L:33:GLY:HA3	1:M:42:ASN:ND2	2.35	0.42
1:M:91:ILE:HD11	2:Z:6:ALA:HB1	2.02	0.42
1:B:91:ILE:HD11	2:J:6:ALA:HB1	2.01	0.42
1:S:93:ILE:O	1:S:93:ILE:HG13	2.19	0.42
1:I:93:ILE:O	1:I:93:ILE:HG13	2.19	0.42
2:b:4:PRO:HA	2:b:5:YCP:HA	1.73	0.42
1:A:148:GLU:CD	1:A:152:ARG:HH21	2.28	0.41
1:B:190:MET:HE1	2:J:2:PHE:HB3	2.01	0.41
1:N:98:SER:HB3	3:N:342:HOH:O	2.20	0.41
1:G:163:GLU:H	1:G:163:GLU:CD	2.28	0.41
1:N:147:ARG:HD3	3:N:332:HOH:O	2.19	0.41
1:I:190:MET:HE2	1:I:190:MET:HB3	1.81	0.41
2:Q:4:PRO:HA	2:Q:5:YCP:HA	1.74	0.41
2:H:4:PRO:HA	2:H:5:YCP:HA	1.85	0.41
2:c:4:PRO:HA	2:c:5:YCP:HA	1.79	0.41
1:T:113:PHE:CZ	2:c:5:YCP:HD	2.56	0.41
1:C:31:MET:HE2	1:C:31:MET:HB2	1.97	0.41
1:C:190:MET:HE2	1:C:190:MET:HB3	1.96	0.41
1:G:85:LYS:N	1:G:86:PRO:CD	2.84	0.41
1:N:190:MET:HE2	1:N:190:MET:HB3	1.93	0.41
3:A:318:HOH:O	1:G:34:SER:HA	2.20	0.40
1:L:85:LYS:N	1:L:86:PRO:CD	2.84	0.40
2:X:4:PRO:HA	2:X:5:YCP:HA	1.67	0.40
2:Y:4:PRO:HA	2:Y:5:YCP:HA	1.81	0.40
1:F:93:ILE:O	1:F:93:ILE:HG13	2.20	0.40
1:L:18:TYR:OH	1:L:26:LYS:HD2	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	175/203 (86%)	172 (98%)	3 (2%)	0	100	100
1	B	176/203 (87%)	172 (98%)	4 (2%)	0	100	100
1	C	175/203 (86%)	172 (98%)	3 (2%)	0	100	100
1	D	176/203 (87%)	171 (97%)	5 (3%)	0	100	100
1	E	174/203 (86%)	170 (98%)	4 (2%)	0	100	100
1	F	175/203 (86%)	171 (98%)	4 (2%)	0	100	100
1	G	176/203 (87%)	172 (98%)	4 (2%)	0	100	100
1	I	175/203 (86%)	171 (98%)	4 (2%)	0	100	100
1	K	176/203 (87%)	171 (97%)	5 (3%)	0	100	100
1	L	176/203 (87%)	171 (97%)	5 (3%)	0	100	100
1	M	175/203 (86%)	170 (97%)	5 (3%)	0	100	100
1	N	175/203 (86%)	172 (98%)	3 (2%)	0	100	100
1	S	174/203 (86%)	170 (98%)	4 (2%)	0	100	100
1	T	174/203 (86%)	170 (98%)	4 (2%)	0	100	100
2	H	4/7 (57%)	3 (75%)	1 (25%)	0	100	100
2	J	4/7 (57%)	3 (75%)	1 (25%)	0	100	100
2	O	4/7 (57%)	3 (75%)	1 (25%)	0	100	100
2	P	4/7 (57%)	3 (75%)	1 (25%)	0	100	100
2	Q	4/7 (57%)	3 (75%)	1 (25%)	0	100	100
2	R	4/7 (57%)	3 (75%)	1 (25%)	0	100	100
2	U	4/7 (57%)	3 (75%)	1 (25%)	0	100	100
2	V	4/7 (57%)	3 (75%)	1 (25%)	0	100	100
2	X	4/7 (57%)	3 (75%)	1 (25%)	0	100	100
2	Y	4/7 (57%)	2 (50%)	2 (50%)	0	100	100
2	Z	4/7 (57%)	3 (75%)	1 (25%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	a	4/7 (57%)	3 (75%)	1 (25%)	0	100	100
2	b	4/7 (57%)	3 (75%)	1 (25%)	0	100	100
2	c	4/7 (57%)	2 (50%)	2 (50%)	0	100	100
All	All	2508/2940 (85%)	2435 (97%)	73 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	146/171 (85%)	145 (99%)	1 (1%)	76	83
1	B	143/171 (84%)	142 (99%)	1 (1%)	76	83
1	C	145/171 (85%)	143 (99%)	2 (1%)	59	70
1	D	148/171 (86%)	146 (99%)	2 (1%)	59	70
1	E	146/171 (85%)	143 (98%)	3 (2%)	47	60
1	F	143/171 (84%)	143 (100%)	0	100	100
1	G	144/171 (84%)	144 (100%)	0	100	100
1	I	142/171 (83%)	140 (99%)	2 (1%)	59	70
1	K	146/171 (85%)	146 (100%)	0	100	100
1	L	146/171 (85%)	144 (99%)	2 (1%)	59	70
1	M	148/171 (86%)	145 (98%)	3 (2%)	48	62
1	N	145/171 (85%)	145 (100%)	0	100	100
1	S	144/171 (84%)	143 (99%)	1 (1%)	76	83
1	T	143/171 (84%)	142 (99%)	1 (1%)	76	83
2	H	3/3 (100%)	3 (100%)	0	100	100
2	J	3/3 (100%)	3 (100%)	0	100	100
2	O	3/3 (100%)	3 (100%)	0	100	100
2	P	3/3 (100%)	3 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	Q	3/3 (100%)	3 (100%)	0	100	100
2	R	3/3 (100%)	3 (100%)	0	100	100
2	U	3/3 (100%)	3 (100%)	0	100	100
2	V	3/3 (100%)	3 (100%)	0	100	100
2	X	3/3 (100%)	3 (100%)	0	100	100
2	Y	3/3 (100%)	3 (100%)	0	100	100
2	Z	3/3 (100%)	3 (100%)	0	100	100
2	a	3/3 (100%)	3 (100%)	0	100	100
2	b	3/3 (100%)	3 (100%)	0	100	100
2	c	3/3 (100%)	3 (100%)	0	100	100
All	All	2071/2436 (85%)	2053 (99%)	18 (1%)	70	80

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	121	MET
1	B	44	ILE
1	C	44	ILE
1	C	154	LEU
1	D	18	TYR
1	D	130	GLN
1	E	18	TYR
1	E	44	ILE
1	E	190	MET
1	I	26	LYS
1	I	44	ILE
1	L	18	TYR
1	L	121	MET
1	M	7	VAL
1	M	121	MET
1	M	167	LYS
1	S	121	MET
1	T	162	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (37) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	42	ASN
1	B	65	ASN
1	B	117	ASN
1	C	42	ASN
1	C	65	ASN
1	C	117	ASN
1	C	151	ASN
1	D	42	ASN
1	D	65	ASN
1	D	151	ASN
1	E	42	ASN
1	E	65	ASN
1	E	166	GLN
1	F	42	ASN
1	F	65	ASN
1	G	42	ASN
1	G	65	ASN
1	G	151	ASN
1	I	42	ASN
1	I	65	ASN
1	I	151	ASN
1	K	42	ASN
1	K	47	GLN
1	K	151	ASN
1	L	65	ASN
1	M	42	ASN
1	N	47	GLN
1	N	65	ASN
1	N	117	ASN
1	S	42	ASN
1	S	52	GLN
1	S	54	GLN
1	S	65	ASN
1	T	39	ASN
1	T	42	ASN
1	T	65	ASN
1	T	151	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

28 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$
2	MP8	Q	7	2	6,8,9	1.50	1 (16%)	3,10,12	2.05	1 (33%)
2	MP8	c	7	2	6,8,9	0.95	1 (16%)	3,10,12	1.60	1 (33%)
2	YCP	Z	5	2	6,8,9	1.20	1 (16%)	7,9,11	1.34	1 (14%)
2	YCP	O	5	2	6,8,9	1.54	2 (33%)	7,9,11	1.49	1 (14%)
2	MP8	U	7	2	6,8,9	0.69	0	3,10,12	1.83	1 (33%)
2	MP8	V	7	2	6,8,9	1.21	1 (16%)	3,10,12	1.59	1 (33%)
2	YCP	V	5	2	6,8,9	0.76	0	7,9,11	1.89	2 (28%)
2	MP8	P	7	2	6,8,9	0.95	0	3,10,12	1.89	1 (33%)
2	YCP	R	5	2	6,8,9	1.30	1 (16%)	7,9,11	1.44	1 (14%)
2	YCP	c	5	2	6,8,9	1.42	1 (16%)	7,9,11	1.39	1 (14%)
2	MP8	J	7	2	6,8,9	1.15	1 (16%)	3,10,12	1.64	1 (33%)
2	YCP	b	5	2	6,8,9	1.67	1 (16%)	7,9,11	1.44	1 (14%)
2	YCP	Y	5	2	6,8,9	1.18	1 (16%)	7,9,11	1.64	1 (14%)
2	YCP	U	5	2	6,8,9	1.67	2 (33%)	7,9,11	1.29	1 (14%)
2	YCP	P	5	2	6,8,9	1.47	2 (33%)	7,9,11	1.32	1 (14%)
2	YCP	J	5	2	6,8,9	1.64	1 (16%)	7,9,11	1.59	1 (14%)
2	MP8	O	7	2	6,8,9	0.52	0	3,10,12	1.51	1 (33%)
2	YCP	X	5	2	6,8,9	1.30	1 (16%)	7,9,11	1.52	1 (14%)
2	MP8	H	7	2	6,8,9	1.06	1 (16%)	3,10,12	1.82	1 (33%)
2	MP8	R	7	2	6,8,9	0.89	0	3,10,12	1.46	1 (33%)
2	MP8	b	7	2	6,8,9	1.35	1 (16%)	3,10,12	1.87	1 (33%)
2	YCP	a	5	2	6,8,9	1.47	1 (16%)	7,9,11	1.27	1 (14%)
2	YCP	Q	5	2	6,8,9	1.23	1 (16%)	7,9,11	1.52	1 (14%)
2	YCP	H	5	2	6,8,9	1.35	1 (16%)	7,9,11	1.09	1 (14%)
2	MP8	Y	7	2	6,8,9	0.80	0	3,10,12	1.90	1 (33%)
2	MP8	X	7	2	6,8,9	0.94	0	3,10,12	2.23	1 (33%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	MP8	Z	7	2	6,8,9	1.18	1 (16%)	3,10,12	2.03	1 (33%)
2	MP8	a	7	2	6,8,9	1.34	1 (16%)	3,10,12	1.82	1 (33%)

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
2	MP8	Q	7	2	-	0/0/11/13	0/1/1/1
2	MP8	c	7	2	-	0/0/11/13	0/1/1/1
2	YCP	Z	5	2	-	0/1/10/12	0/1/1/1
2	YCP	O	5	2	-	0/1/10/12	0/1/1/1
2	MP8	U	7	2	-	0/0/11/13	0/1/1/1
2	MP8	V	7	2	-	0/0/11/13	0/1/1/1
2	YCP	V	5	2	-	0/1/10/12	0/1/1/1
2	MP8	P	7	2	-	0/0/11/13	0/1/1/1
2	YCP	R	5	2	-	0/1/10/12	0/1/1/1
2	YCP	c	5	2	-	0/1/10/12	0/1/1/1
2	MP8	J	7	2	-	0/0/11/13	0/1/1/1
2	YCP	b	5	2	-	0/1/10/12	0/1/1/1
2	YCP	Y	5	2	-	0/1/10/12	0/1/1/1
2	YCP	U	5	2	-	0/1/10/12	0/1/1/1
2	YCP	P	5	2	-	0/1/10/12	0/1/1/1
2	YCP	J	5	2	-	0/1/10/12	0/1/1/1
2	MP8	O	7	2	-	0/0/11/13	0/1/1/1
2	YCP	X	5	2	-	0/1/10/12	0/1/1/1
2	MP8	H	7	2	-	0/0/11/13	0/1/1/1
2	MP8	R	7	2	-	0/0/11/13	0/1/1/1
2	MP8	b	7	2	-	0/0/11/13	0/1/1/1
2	YCP	a	5	2	-	0/1/10/12	0/1/1/1
2	YCP	Q	5	2	-	0/1/10/12	0/1/1/1
2	YCP	H	5	2	-	0/1/10/12	0/1/1/1
2	MP8	Y	7	2	-	0/0/11/13	0/1/1/1
2	MP8	X	7	2	-	0/0/11/13	0/1/1/1
2	MP8	Z	7	2	-	0/0/11/13	0/1/1/1
2	MP8	a	7	2	-	0/0/11/13	0/1/1/1

All (24) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	J	5	YCP	CG-CD	3.54	1.64	1.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	Q	7	MP8	CB-CA	-3.21	1.47	1.54
2	b	5	YCP	CG-CD	3.10	1.63	1.51
2	H	5	YCP	CD-CE	3.03	1.63	1.49
2	O	5	YCP	CG-CD	2.95	1.62	1.51
2	U	5	YCP	CG-CD	2.94	1.62	1.51
2	a	5	YCP	CG-CD	2.91	1.62	1.51
2	Z	7	MP8	CB-CA	-2.65	1.48	1.54
2	a	7	MP8	CB-CA	-2.60	1.48	1.54
2	b	7	MP8	CA-N	-2.58	1.44	1.48
2	V	7	MP8	CB-CA	-2.53	1.48	1.54
2	U	5	YCP	CD-CE	2.47	1.61	1.49
2	R	5	YCP	CG-CD	2.38	1.60	1.51
2	c	5	YCP	CG-CD	2.37	1.60	1.51
2	Q	5	YCP	CG-CD	2.34	1.60	1.51
2	P	5	YCP	CD-CE	2.34	1.60	1.49
2	J	7	MP8	CA-N	-2.22	1.44	1.48
2	Y	5	YCP	CD-CE	2.22	1.60	1.49
2	H	7	MP8	CA-N	-2.22	1.44	1.48
2	X	5	YCP	CD-CE	2.17	1.59	1.49
2	Z	5	YCP	CG-CD	2.15	1.59	1.51
2	P	5	YCP	CG-CD	2.05	1.59	1.51
2	c	7	MP8	CB-CA	-2.04	1.49	1.54
2	O	5	YCP	CD-CE	2.02	1.59	1.49

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	J	5	YCP	O-C-CA	-4.06	114.33	124.77
2	X	5	YCP	O-C-CA	-3.97	114.56	124.77
2	Q	5	YCP	O-C-CA	-3.75	115.11	124.77
2	Y	5	YCP	O-C-CA	-3.67	115.34	124.77
2	V	5	YCP	O-C-CA	-3.61	115.48	124.77
2	O	5	YCP	O-C-CA	-3.50	115.77	124.77
2	X	7	MP8	CG-CD-N	-3.45	102.50	106.65
2	Q	7	MP8	CG-CD-N	-3.35	102.61	106.65
2	Z	7	MP8	CG-CD-N	-3.31	102.67	106.65
2	Z	5	YCP	O-C-CA	-3.29	116.32	124.77
2	b	5	YCP	O-C-CA	-3.23	116.47	124.77
2	U	7	MP8	CG-CD-N	-3.12	102.89	106.65
2	P	5	YCP	O-C-CA	-3.03	116.98	124.77
2	U	5	YCP	O-C-CA	-3.01	117.04	124.77
2	R	5	YCP	O-C-CA	-3.00	117.06	124.77

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	a	5	YCP	O-C-CA	-2.96	117.16	124.77
2	c	5	YCP	O-C-CA	-2.94	117.22	124.77
2	V	5	YCP	CB-CA-N	-2.91	104.03	109.23
2	H	7	MP8	CG-CD-N	-2.80	103.27	106.65
2	Y	7	MP8	CG-CD-N	-2.80	103.28	106.65
2	b	7	MP8	CG-CD-N	-2.74	103.35	106.65
2	a	7	MP8	CG-CD-N	-2.74	103.35	106.65
2	V	7	MP8	CG-CD-N	-2.63	103.49	106.65
2	H	5	YCP	O-C-CA	-2.55	118.21	124.77
2	P	7	MP8	CG-CD-N	-2.52	103.61	106.65
2	c	7	MP8	CG-CD-N	-2.52	103.61	106.65
2	J	7	MP8	CG-CD-N	-2.37	103.80	106.65
2	O	7	MP8	CG-CD-N	-2.22	103.97	106.65
2	R	7	MP8	CG-CD-N	-2.16	104.05	106.65

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

14 monomers are involved in 14 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	O	5	YCP	1	0
2	V	5	YCP	1	0
2	c	5	YCP	2	0
2	b	5	YCP	1	0
2	Y	5	YCP	1	0
2	U	5	YCP	1	0
2	P	5	YCP	1	0
2	J	5	YCP	1	0
2	O	7	MP8	1	0
2	X	5	YCP	1	0
2	b	7	MP8	1	0
2	a	5	YCP	1	0
2	Q	5	YCP	1	0
2	H	5	YCP	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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	179/203 (88%)	-0.12	4 (2%) 62 63	30, 35, 44, 53	0
1	B	180/203 (88%)	-0.11	5 (2%) 55 54	30, 35, 43, 52	0
1	C	179/203 (88%)	-0.22	3 (1%) 69 70	26, 32, 41, 48	0
1	D	180/203 (88%)	-0.24	5 (2%) 55 54	24, 30, 39, 46	0
1	E	178/203 (87%)	-0.25	2 (1%) 78 79	24, 30, 38, 49	0
1	F	179/203 (88%)	-0.21	5 (2%) 55 54	24, 31, 42, 50	0
1	G	180/203 (88%)	-0.12	3 (1%) 69 70	30, 35, 43, 51	0
1	I	179/203 (88%)	0.01	4 (2%) 62 63	32, 37, 44, 52	0
1	K	180/203 (88%)	-0.15	5 (2%) 55 54	28, 33, 42, 50	0
1	L	180/203 (88%)	-0.20	3 (1%) 69 70	25, 31, 38, 47	0
1	M	179/203 (88%)	-0.15	3 (1%) 69 70	26, 31, 41, 48	0
1	N	179/203 (88%)	-0.22	3 (1%) 69 70	26, 32, 40, 51	0
1	S	178/203 (87%)	0.07	5 (2%) 55 54	30, 37, 45, 51	0
1	T	178/203 (87%)	0.06	3 (1%) 69 70	33, 38, 46, 50	0
2	H	4/7 (57%)	-0.16	0 100 100	37, 38, 41, 44	0
2	J	4/7 (57%)	0.11	0 100 100	36, 39, 40, 43	0
2	O	4/7 (57%)	-0.26	0 100 100	33, 36, 38, 41	0
2	P	4/7 (57%)	-0.46	0 100 100	30, 32, 34, 37	0
2	Q	4/7 (57%)	-0.10	0 100 100	32, 33, 35, 39	0
2	R	4/7 (57%)	-0.07	0 100 100	35, 37, 38, 40	0
2	U	4/7 (57%)	0.02	0 100 100	38, 40, 41, 44	0
2	V	4/7 (57%)	0.78	1 (25%) 2 1	40, 42, 44, 48	0
2	X	4/7 (57%)	-0.03	0 100 100	35, 35, 38, 41	0
2	Y	4/7 (57%)	-0.05	0 100 100	34, 36, 36, 41	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å ²)	Q<0.9
2	Z	4/7 (57%)	-0.59	0	100	100	32, 35, 38, 41	0
2	a	4/7 (57%)	0.29	0	100	100	36, 37, 40, 41	0
2	b	4/7 (57%)	0.27	0	100	100	40, 43, 43, 44	0
2	c	4/7 (57%)	0.38	0	100	100	41, 43, 44, 46	0
All	All	2564/2940 (87%)	-0.13	54 (2%)	63	64	24, 34, 43, 53	0

All (54) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	S	6	THR	5.2
1	I	193	GLU	4.2
1	B	193	GLU	4.2
1	S	7	VAL	4.2
1	D	18	TYR	4.1
1	N	7	VAL	4.0
1	K	193	GLU	4.0
1	B	18	TYR	3.7
1	B	7	VAL	3.7
1	G	193	GLU	3.5
1	G	7	VAL	3.4
1	L	193	GLU	3.3
1	A	18	TYR	3.3
1	D	6	THR	3.2
1	F	18	TYR	3.1
1	K	7	VAL	3.1
1	E	18	TYR	3.0
1	M	7	VAL	3.0
1	K	6	THR	3.0
1	I	7	VAL	2.9
1	T	4	ILE	2.9
1	A	7	VAL	2.9
1	N	6	THR	2.9
1	F	7	VAL	2.8
1	T	18	TYR	2.8
1	M	130	GLN	2.8
1	F	4	ILE	2.7
1	B	4	ILE	2.7
1	M	18	TYR	2.6
1	A	4	ILE	2.6
1	K	4	ILE	2.6
1	S	5	PRO	2.5

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Mol	Chain	Res	Type	RSRZ
1	S	4	ILE	2.5
1	C	6	THR	2.5
1	C	4	ILE	2.5
1	K	18	TYR	2.4
1	L	18	TYR	2.4
1	B	6	THR	2.3
1	D	4	ILE	2.3
1	D	7	VAL	2.3
1	E	7	VAL	2.3
1	S	191	VAL	2.3
1	G	18	TYR	2.3
1	A	6	THR	2.2
1	F	191	VAL	2.2
1	F	6	THR	2.2
1	T	6	THR	2.2
1	N	4	ILE	2.2
1	C	7	VAL	2.2
1	D	191	VAL	2.2
2	V	4	PRO	2.2
1	I	5	PRO	2.1
1	L	31	MET	2.1
1	I	4	ILE	2.0

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
2	MP8	c	7	8/9	0.85	0.13	42,43,43,44	0
2	MP8	Y	7	8/9	0.88	0.12	35,36,36,38	0
2	MP8	b	7	8/9	0.88	0.13	40,42,44,45	0
2	YCP	b	5	8/9	0.88	0.12	43,43,44,44	0
2	MP8	O	7	8/9	0.89	0.13	35,36,37,39	0
2	MP8	a	7	8/9	0.89	0.12	36,37,38,40	0
2	MP8	U	7	8/9	0.90	0.11	39,40,41,42	0
2	MP8	R	7	8/9	0.90	0.11	37,38,39,41	0
2	MP8	Z	7	8/9	0.90	0.11	34,35,37,38	0
2	MP8	X	7	8/9	0.91	0.11	35,36,37,39	0
2	YCP	J	5	8/9	0.91	0.10	40,41,41,42	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	YCP	X	5	8/9	0.91	0.10	37,38,39,40	0
2	MP8	Q	7	8/9	0.92	0.10	33,34,34,37	0
2	YCP	V	5	8/9	0.92	0.10	43,44,45,45	0
2	MP8	J	7	8/9	0.92	0.10	38,39,40,40	0
2	MP8	V	7	8/9	0.92	0.11	39,40,42,46	0
2	YCP	H	5	8/9	0.92	0.09	39,40,41,41	0
2	YCP	P	5	8/9	0.93	0.10	34,35,36,36	0
2	YCP	Y	5	8/9	0.93	0.09	37,38,39,40	0
2	YCP	O	5	8/9	0.93	0.10	38,38,39,40	0
2	YCP	c	5	8/9	0.93	0.09	43,44,44,44	0
2	YCP	Q	5	8/9	0.94	0.09	33,35,36,36	0
2	YCP	R	5	8/9	0.94	0.09	37,39,40,40	0
2	MP8	H	7	8/9	0.94	0.08	37,38,39,41	0
2	YCP	U	5	8/9	0.94	0.09	41,42,43,44	0
2	YCP	Z	5	8/9	0.94	0.08	37,38,38,38	0
2	YCP	a	5	8/9	0.95	0.08	37,37,38,38	0
2	MP8	P	7	8/9	0.95	0.09	31,32,34,34	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides 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.