Full wwPDB X-ray Structure Validation Report (i)

Jun 23, 2024 – 02:00 AM EDT

PDB ID	:	6JPL
Title	:	The X-ray structure of yeast tRNA methyltransferase Trm7-Trm734 in complex with S-adenosyl-L-methionine
Authors	:	Hirata, A.; Okada, K.; Yoshii, K.; Shiraisi, H.; Saijo, S.; Yonezawa, K.; Shimizu, N.; Hori, H.
Deposited on	:	2019-03-27
Resolution	:	2.32 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 (i) 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 (i)) were used in the production of this report:

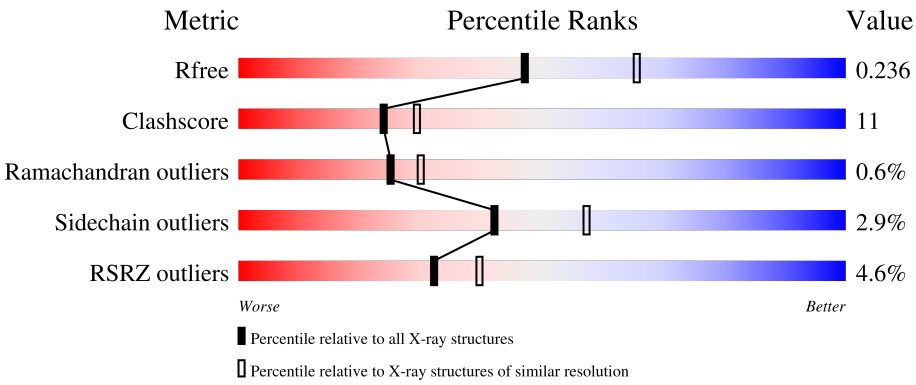
MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	2.37.1
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.37.1

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.32 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	5974 (2.34-2.30)
Clashscore	141614	6604 (2.34-2.30)
Ramachandran outliers	138981	6523 (2.34-2.30)
Sidechain outliers	138945	6523 (2.34-2.30)
RSRZ outliers	127900	5855 (2.34-2.30)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments 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 <=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	1028	82% 15% • •
1	C	1028	% 82% 14% • •
2	B	325	15% 41% 26% • 29%
2	D	325	16% 42% 24% • • 29%

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
5	SAM	D	401	-	-	-	X

2 Entry composition (i)

There are 6 unique types of molecules in this entry. The entry contains 20155 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 tRNA (guanosine(34)-2'-O)-methyltransferase non-catalytic subunit TRM734.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	998	7951	5074	1328	1517	32	0	0	0
1	C	998	7951	5074	1328	1517	32	0	0	0

There are 30 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1014	LEU	-	expression tag	UNP Q08924
A	1015	GLU	-	expression tag	UNP Q08924
A	1016	VAL	-	expression tag	UNP Q08924
A	1017	LEU	-	expression tag	UNP Q08924
A	1018	PHE	-	expression tag	UNP Q08924
A	1019	GLN	-	expression tag	UNP Q08924
A	1020	GLY	-	expression tag	UNP Q08924
A	1021	PRO	-	expression tag	UNP Q08924
A	1022	SER	-	expression tag	UNP Q08924
A	1023	HIS	-	expression tag	UNP Q08924
A	1024	HIS	-	expression tag	UNP Q08924
A	1025	HIS	-	expression tag	UNP Q08924
A	1026	HIS	-	expression tag	UNP Q08924
A	1027	HIS	-	expression tag	UNP Q08924
A	1028	HIS	-	expression tag	UNP Q08924
C	1014	LEU	-	expression tag	UNP Q08924
C	1015	GLU	-	expression tag	UNP Q08924
C	1016	VAL	-	expression tag	UNP Q08924
C	1017	LEU	-	expression tag	UNP Q08924
C	1018	PHE	-	expression tag	UNP Q08924
C	1019	GLN	-	expression tag	UNP Q08924
C	1020	GLY	-	expression tag	UNP Q08924
C	1021	PRO	-	expression tag	UNP Q08924
C	1022	SER	-	expression tag	UNP Q08924

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Chain	Residue	Modelled	Actual	Comment	Reference
C	1023	HIS	-	expression tag	UNP Q08924
C	1024	HIS	-	expression tag	UNP Q08924
C	1025	HIS	-	expression tag	UNP Q08924
C	1026	HIS	-	expression tag	UNP Q08924
C	1027	HIS	-	expression tag	UNP Q08924
C	1028	HIS	-	expression tag	UNP Q08924

- Molecule 2 is a protein called tRNA (cytidine(34)/guanosine(34)-2'-O)-methyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	231	1829	1158	318	342	11	0	0	0
2	D	231	1829	1158	318	342	11	0	0	0

There are 30 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	311	LEU	-	expression tag	UNP P38238
B	312	GLU	-	expression tag	UNP P38238
B	313	VAL	-	expression tag	UNP P38238
B	314	LEU	-	expression tag	UNP P38238
B	315	PHE	-	expression tag	UNP P38238
B	316	GLN	-	expression tag	UNP P38238
B	317	GLY	-	expression tag	UNP P38238
B	318	PRO	-	expression tag	UNP P38238
B	319	SER	-	expression tag	UNP P38238
B	320	HIS	-	expression tag	UNP P38238
B	321	HIS	-	expression tag	UNP P38238
B	322	HIS	-	expression tag	UNP P38238
B	323	HIS	-	expression tag	UNP P38238
B	324	HIS	-	expression tag	UNP P38238
B	325	HIS	-	expression tag	UNP P38238
D	311	LEU	-	expression tag	UNP P38238
D	312	GLU	-	expression tag	UNP P38238
D	313	VAL	-	expression tag	UNP P38238
D	314	LEU	-	expression tag	UNP P38238
D	315	PHE	-	expression tag	UNP P38238
D	316	GLN	-	expression tag	UNP P38238
D	317	GLY	-	expression tag	UNP P38238
D	318	PRO	-	expression tag	UNP P38238
D	319	SER	-	expression tag	UNP P38238

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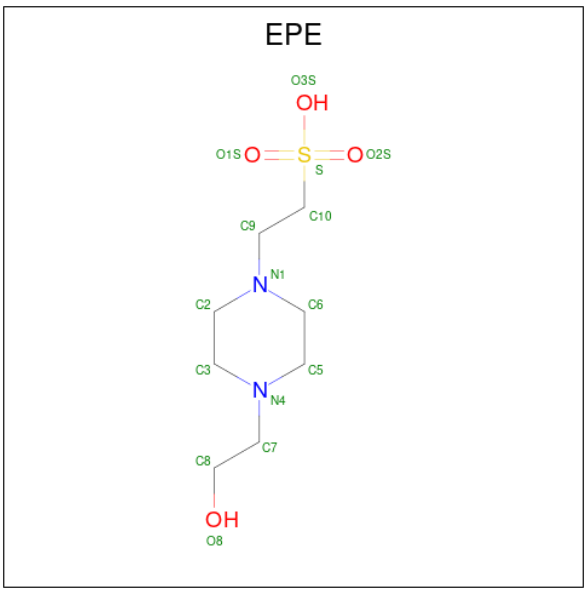
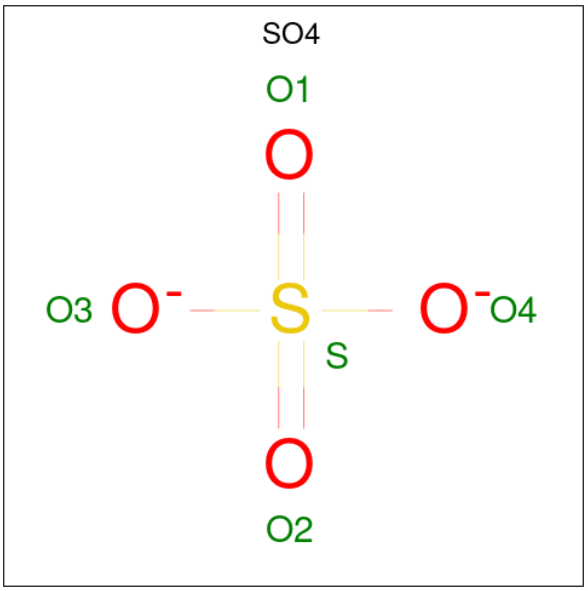
Chain	Residue	Modelled	Actual	Comment	Reference
D	320	HIS	-	expression tag	UNP P38238
D	321	HIS	-	expression tag	UNP P38238
D	322	HIS	-	expression tag	UNP P38238
D	323	HIS	-	expression tag	UNP P38238
D	324	HIS	-	expression tag	UNP P38238
D	325	HIS	-	expression tag	UNP P38238

- Molecule 3 is 4-(2-HYDROXYETHYL)-1-PIPERAZINE ETHANESULFONIC ACID (three-letter code: EPE) (formula: $C_8H_{18}N_2O_4S$).



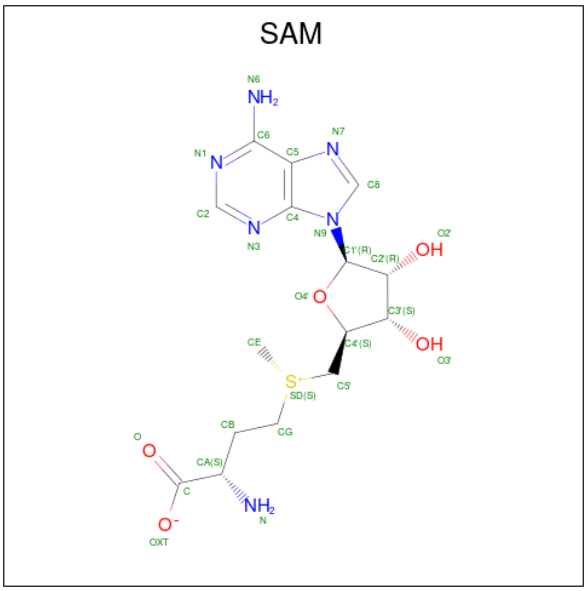
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	S		
3	A	1	15	8	2	4	1	0	0
			Total	C	N	O	S		
3	C	1	15	8	2	4	1	0	0

- Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O_4S).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	O	S			
4	A	1	5	4	1		0	0
			Total	O	S			
4	A	1	5	4	1		0	0
			Total	O	S			
4	A	1	5	4	1		0	0
			Total	O	S			
4	A	1	5	4	1		0	0
			Total	O	S			
4	A	1	5	4	1		0	0
			Total	O	S			
4	C	1	5	4	1		0	0
			Total	O	S			
4	C	1	5	4	1		0	0
			Total	O	S			
4	C	1	5	4	1		0	0
			Total	O	S			
4	C	1	5	4	1		0	0
			Total	O	S			
4	C	1	5	4	1		0	0

- Molecule 5 is S-ADENOSYLMETHIONINE (three-letter code: SAM) (formula: $C_{15}H_{22}N_6O_5S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	S		
5	B	1	27	15	6	5	1	0	0
			Total	C	N	O	S		
5	D	1	27	15	6	5	1	0	0

- Molecule 6 is water.

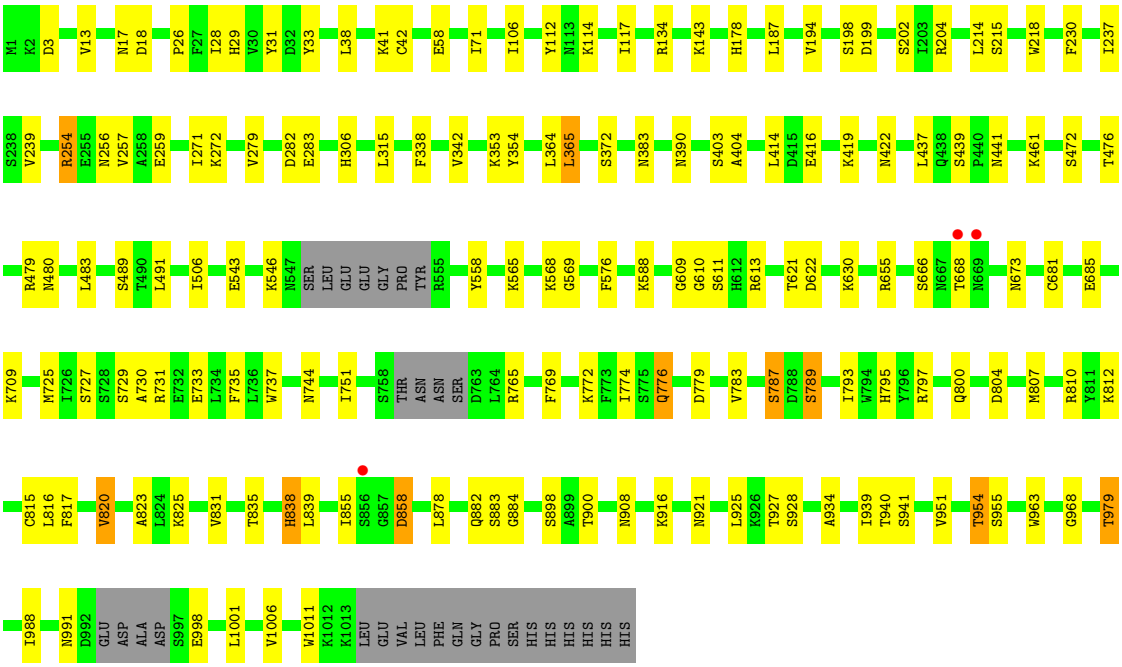
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
6	A	213	213	213	0	0
			Total	O		
6	B	19	19	19	0	0
			Total	O		
6	C	208	208	208	0	0
			Total	O		
6	D	21	21	21	0	0

3 Residue-property plots (i)

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

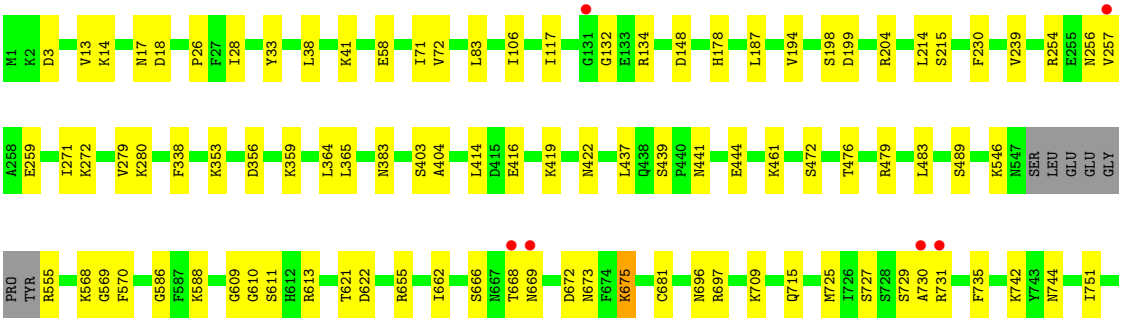
• Molecule 1: tRNA (guanosine(34)-2'-O)-methyltransferase non-catalytic subunit TRM734

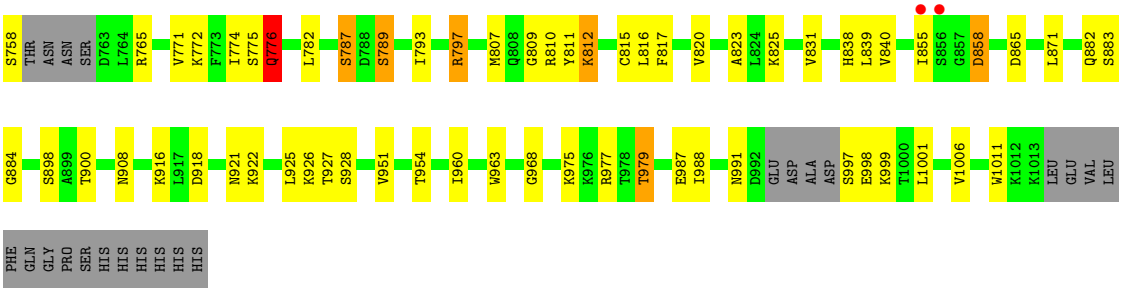
Chain A:



• Molecule 1: tRNA (guanosine(34)-2'-O)-methyltransferase non-catalytic subunit TRM734

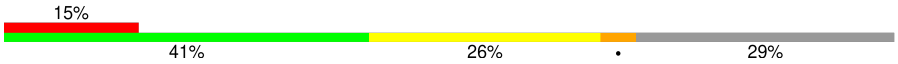
Chain C:

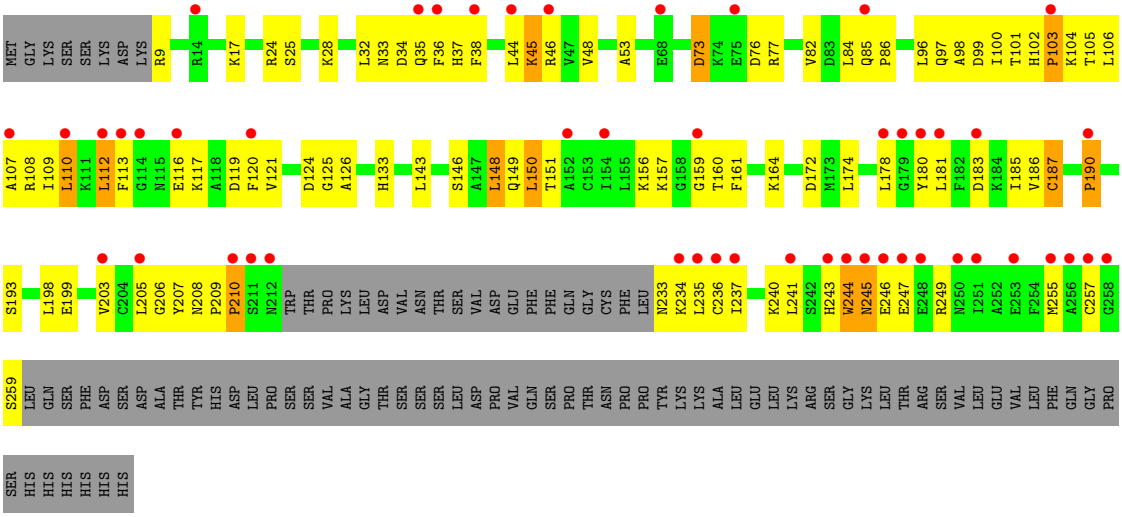




• Molecule 2: tRNA (cytidine(34)/guanosine(34)-2'-O)-methyltransferase

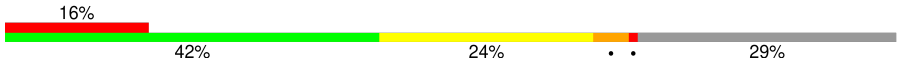
Chain B:

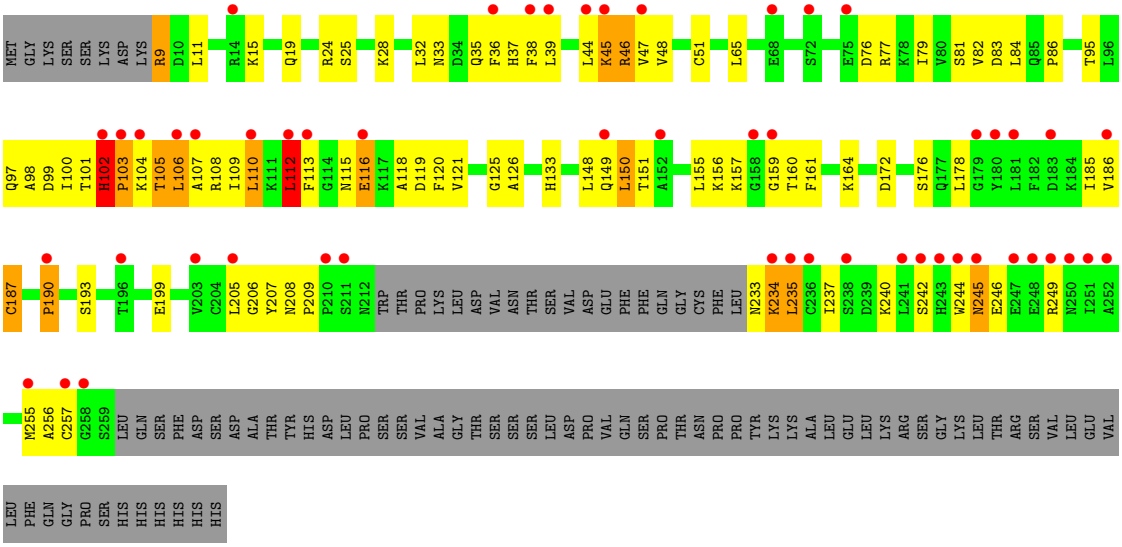




• Molecule 2: tRNA (cytidine(34)/guanosine(34)-2'-O)-methyltransferase

Chain D:





4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	253.56Å 110.67Å 133.56Å 90.00° 90.01° 90.00°	Depositor
Resolution (Å)	47.41 – 2.32 47.41 – 2.32	Depositor EDS
% Data completeness (in resolution range)	99.8 (47.41-2.32) 99.8 (47.41-2.32)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.69 (at 2.32Å)	Xtriage
Refinement program	PHENIX 1.10.1_2155	Depositor
R, R_{free}	0.203 , 0.235 0.206 , 0.236	Depositor DCC
R_{free} test set	7996 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å^2)	42.0	Xtriage
Anisotropy	0.436	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.34 , 30.7	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	0.487 for -h,-k,l	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	20155	wwPDB-VP
Average B, all atoms (Å^2)	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.96% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ 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: SAM, SO4, EPE

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	0.48	1/8110 (0.0%)	0.67	0/10972
1	C	0.48	1/8110 (0.0%)	0.69	3/10972 (0.0%)
2	B	0.52	0/1865	0.88	4/2516 (0.2%)
2	D	0.63	3/1865 (0.2%)	1.04	13/2516 (0.5%)
All	All	0.50	5/19950 (0.0%)	0.74	20/26976 (0.1%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2
1	C	0	2
2	B	0	1
2	D	0	2
All	All	0	7

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	116	GLU	CB-CG	-9.80	1.33	1.52
2	D	46	ARG	CB-CG	-6.13	1.35	1.52
1	A	681	CYS	CB-SG	-5.54	1.72	1.81
2	D	46	ARG	NE-CZ	-5.16	1.26	1.33
1	C	681	CYS	CB-SG	-5.07	1.73	1.81

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	D	106	LEU	CB-CG-CD1	-18.14	80.17	111.00
2	D	46	ARG	NE-CZ-NH1	-13.90	113.35	120.30
1	C	675	LYS	CD-CE-NZ	-10.02	88.66	111.70
2	D	110	LEU	CB-CG-CD2	-9.87	94.22	111.00
2	D	235	LEU	CA-CB-CG	9.47	137.07	115.30
2	B	112	LEU	CA-CB-CG	8.84	135.63	115.30
2	B	110	LEU	CA-CB-CG	8.42	134.66	115.30
2	D	105	THR	C-N-CA	-8.33	100.87	121.70
2	D	106	LEU	CB-CG-CD2	7.62	123.96	111.00
2	B	150	LEU	CB-CG-CD1	-7.24	98.70	111.00
2	B	235	LEU	CA-CB-CG	7.10	131.63	115.30
1	C	675	LYS	CB-CG-CD	6.87	129.46	111.60
2	D	110	LEU	CA-CB-CG	6.62	130.51	115.30
2	D	46	ARG	NE-CZ-NH2	6.33	123.46	120.30
2	D	149	GLN	CA-CB-CG	5.86	126.28	113.40
2	D	112	LEU	CB-CG-CD1	5.63	120.58	111.00
2	D	150	LEU	CA-CB-CG	5.43	127.78	115.30
1	C	675	LYS	CA-CB-CG	-5.40	101.53	113.40
2	D	45	LYS	CB-CA-C	5.23	120.85	110.40
2	D	112	LEU	CA-CB-CG	5.09	127.00	115.30

There are no chirality outliers.

All (7) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	776	GLN	Peptide
1	A	855	ILE	Peptide
2	B	245	ASN	Peptide
1	C	776	GLN	Peptide
1	C	855	ILE	Peptide
2	D	102	HIS	Peptide
2	D	245	ASN	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	7951	0	7901	123	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	7951	0	7901	110	1
2	B	1829	0	1829	104	1
2	D	1829	0	1829	116	0
3	A	15	0	17	3	0
3	C	15	0	17	0	0
4	A	25	0	0	1	0
4	C	25	0	0	1	0
5	B	27	0	22	5	0
5	D	27	0	22	6	0
6	A	213	0	0	12	0
6	B	19	0	0	1	0
6	C	208	0	0	5	0
6	D	21	0	0	2	0
All	All	20155	0	19538	436	1

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 11.

All (436) 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:B:102:HIS:ND1	2:B:103:PRO:CD	1.97	1.26
2:B:246:GLU:HG3	2:B:247:GLU:OE2	1.33	1.23
2:B:102:HIS:ND1	2:B:103:PRO:HD3	1.56	1.18
2:B:102:HIS:ND1	2:B:103:PRO:HD2	1.66	1.10
1:A:479:ARG:HG3	1:A:480:ASN:H	1.12	1.08
2:D:104:LYS:O	2:D:108:ARG:N	1.92	1.02
2:B:246:GLU:CG	2:B:247:GLU:OE2	2.08	1.02
2:D:45:LYS:NZ	2:D:46:ARG:HH12	1.61	0.98
2:D:100:ILE:O	2:D:150:LEU:HD21	1.65	0.97
2:D:45:LYS:HZ1	2:D:46:ARG:HH12	1.12	0.94
2:D:106:LEU:HD12	2:D:106:LEU:C	1.87	0.94
1:A:730:ALA:CB	1:A:765:ARG:HG2	1.99	0.93
2:B:35:GLN:NE2	2:B:257:CYS:SG	2.41	0.93
2:B:104:LYS:O	2:B:108:ARG:N	2.03	0.90
1:C:882:GLN:HE21	2:D:256:ALA:H	1.13	0.89
1:A:28:ILE:HB	6:A:1273:HOH:O	1.74	0.87
1:A:479:ARG:HG3	1:A:480:ASN:N	1.90	0.86
1:A:479:ARG:CG	1:A:480:ASN:H	1.89	0.85
2:B:46:ARG:HH12	2:B:76:ASP:HB3	1.41	0.83
2:D:110:LEU:HD12	2:D:115:ASN:OD1	1.77	0.83

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:730:ALA:CB	1:C:765:ARG:HG2	2.09	0.83
2:B:102:HIS:CG	2:B:103:PRO:HD2	2.14	0.82
2:D:103:PRO:HG2	2:D:104:LYS:H	1.45	0.81
1:A:898:SER:HB3	1:A:916:LYS:HG3	1.60	0.81
2:D:107:ALA:HA	2:D:110:LEU:HD21	1.62	0.80
2:B:45:LYS:HZ3	2:B:119:ASP:N	1.80	0.80
2:D:104:LYS:HA	2:D:107:ALA:HB3	1.64	0.80
1:A:673:ASN:ND2	1:A:744:ASN:O	2.15	0.80
1:C:898:SER:HB3	1:C:916:LYS:HG3	1.64	0.79
1:C:416:GLU:OE2	1:C:461:LYS:NZ	2.14	0.77
1:C:673:ASN:ND2	1:C:744:ASN:O	2.19	0.76
2:B:100:ILE:O	2:B:150:LEU:HD11	1.85	0.76
2:D:103:PRO:CG	2:D:104:LYS:H	1.94	0.76
2:B:73:ASP:OD1	2:B:73:ASP:N	2.18	0.75
2:D:45:LYS:HZ1	2:D:46:ARG:NH1	1.84	0.75
1:C:422:ASN:ND2	1:C:472:SER:O	2.21	0.74
1:C:730:ALA:HB2	1:C:765:ARG:HG2	1.70	0.74
1:C:729:SER:OG	1:C:730:ALA:N	2.21	0.74
2:D:103:PRO:HD2	2:D:105:THR:HG23	1.68	0.74
2:B:104:LYS:HA	2:B:107:ALA:HB3	1.68	0.74
1:A:729:SER:O	6:A:1201:HOH:O	2.06	0.73
2:B:103:PRO:HG2	2:B:104:LYS:H	1.54	0.73
1:A:988:ILE:HG12	1:A:1001:LEU:HD22	1.69	0.73
1:A:134:ARG:HG2	2:B:233:ASN:HA	1.69	0.72
1:A:630:LYS:NZ	6:A:1202:HOH:O	2.21	0.72
1:C:825:LYS:NZ	1:C:921:ASN:OD1	2.19	0.72
2:D:108:ARG:O	2:D:112:LEU:HD13	1.90	0.72
1:C:882:GLN:HE22	2:D:255:MET:HA	1.55	0.71
2:D:39:LEU:HD23	2:D:44:LEU:CD2	2.20	0.71
1:A:730:ALA:HB2	1:A:765:ARG:HG2	1.72	0.70
2:D:106:LEU:HD12	2:D:106:LEU:O	1.91	0.70
1:C:730:ALA:CB	1:C:765:ARG:CG	2.70	0.70
1:A:282:ASP:OD2	1:A:283:GLU:HG3	1.92	0.69
2:B:107:ALA:O	2:B:110:LEU:HB3	1.91	0.69
1:C:672:ASP:O	1:C:675:LYS:HG2	1.92	0.69
1:C:882:GLN:NE2	2:D:255:MET:HA	2.07	0.69
2:D:39:LEU:HD23	2:D:44:LEU:HD21	1.73	0.69
1:C:882:GLN:NE2	2:D:256:ALA:H	1.89	0.68
1:C:999:LYS:NZ	4:C:1104:SO4:O2	2.26	0.68
1:C:988:ILE:HG12	1:C:1001:LEU:HD22	1.75	0.68
2:B:102:HIS:CG	2:B:103:PRO:CD	2.74	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:134:ARG:HG3	2:D:233:ASN:HA	1.76	0.67
2:D:104:LYS:HG2	2:D:108:ARG:HB2	1.75	0.67
2:B:85:GLN:NE2	2:B:86:PRO:O	2.28	0.67
1:C:26:PRO:HD3	1:C:979:THR:HG21	1.76	0.67
5:B:401:SAM:N	6:B:501:HOH:O	2.28	0.66
2:D:36:PHE:HE2	2:D:257:CYS:SG	2.18	0.66
2:D:107:ALA:HA	2:D:110:LEU:CD2	2.26	0.66
2:D:36:PHE:HE2	2:D:257:CYS:HG	1.42	0.66
1:C:669:ASN:OD1	1:C:742:LYS:NZ	2.29	0.66
1:C:925:LEU:O	1:C:926:LYS:HD2	1.96	0.66
2:B:9:ARG:HH22	2:D:9:ARG:NH2	1.94	0.65
1:A:271:ILE:HG13	1:A:272:LYS:HB3	1.78	0.65
2:D:99:ASP:OD1	2:D:101:THR:OG1	2.15	0.65
1:A:422:ASN:ND2	1:A:472:SER:O	2.30	0.65
1:A:282:ASP:OD2	1:A:283:GLU:N	2.29	0.64
1:A:609:GLY:O	1:A:611:SER:N	2.30	0.64
2:D:33:ASN:ND2	2:D:39:LEU:H	1.96	0.64
2:D:208:ASN:HB3	2:D:209:PRO:HD2	1.78	0.64
2:B:208:ASN:HB3	2:B:209:PRO:HD2	1.79	0.64
2:D:103:PRO:HD2	2:D:105:THR:H	1.63	0.64
1:A:730:ALA:HB1	1:A:765:ARG:HG2	1.80	0.63
2:D:103:PRO:CG	2:D:104:LYS:N	2.61	0.63
2:B:104:LYS:HG2	2:B:108:ARG:HB2	1.80	0.63
1:A:729:SER:OG	1:A:730:ALA:N	2.25	0.63
1:C:898:SER:CB	1:C:916:LYS:HG3	2.30	0.62
1:A:26:PRO:HD3	1:A:979:THR:HG21	1.81	0.62
2:B:125:GLY:HA2	5:B:401:SAM:HG1	1.82	0.62
2:B:208:ASN:HB3	2:B:209:PRO:CD	2.29	0.62
2:D:208:ASN:HB3	2:D:209:PRO:CD	2.29	0.62
1:A:991:ASN:ND2	1:A:998:GLU:HG2	2.14	0.62
1:A:882:GLN:NE2	2:B:255:MET:HA	2.14	0.62
2:B:45:LYS:HE2	2:B:46:ARG:HB2	1.82	0.62
1:C:730:ALA:O	1:C:731:ARG:HB2	1.99	0.62
2:B:35:GLN:C	2:B:37:HIS:H	2.04	0.61
2:D:237:ILE:HG13	2:D:240:LYS:HD3	1.81	0.61
1:A:568:LYS:HD2	1:A:569:GLY:N	2.15	0.61
2:D:46:ARG:NH1	2:D:116:GLU:OE2	2.34	0.61
2:B:246:GLU:CD	2:B:247:GLU:OE2	2.38	0.61
1:A:735:PHE:HB3	1:A:751:ILE:HD11	1.83	0.61
2:B:174:LEU:O	2:B:178:LEU:HB2	2.01	0.61
1:C:725:MET:HE1	1:C:727:SER:HB3	1.83	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:45:LYS:HG2	2:D:46:ARG:NH1	2.17	0.60
2:D:46:ARG:HG2	2:D:113:PHE:CD1	2.37	0.60
2:B:99:ASP:OD1	2:B:101:THR:OG1	2.21	0.59
2:D:106:LEU:HB3	2:D:150:LEU:HD13	1.84	0.59
2:B:46:ARG:HH12	2:B:76:ASP:CB	2.11	0.59
1:C:38:LEU:HD21	1:C:41:LYS:HG3	1.84	0.59
2:B:190:PRO:HD2	2:B:193:SER:HB2	1.85	0.59
1:A:611:SER:HA	1:C:810:ARG:CZ	2.32	0.59
1:A:797:ARG:NH1	1:A:804:ASP:OD1	2.35	0.59
1:A:611:SER:HA	1:C:810:ARG:NH2	2.17	0.59
1:A:991:ASN:HD21	1:A:998:GLU:HG2	1.68	0.58
1:C:991:ASN:OD1	1:C:998:GLU:HG2	2.03	0.58
1:C:772:LYS:HE2	1:C:823:ALA:HB2	1.83	0.58
2:D:11:LEU:CD2	2:D:15:LYS:HD3	2.34	0.58
1:A:414:LEU:HD12	1:A:437:LEU:HD23	1.85	0.58
1:C:730:ALA:HB1	1:C:765:ARG:HG3	1.84	0.58
2:D:242:SER:O	2:D:246:GLU:HB2	2.04	0.58
1:A:898:SER:CB	1:A:916:LYS:HG3	2.33	0.57
1:C:609:GLY:O	1:C:611:SER:N	2.37	0.57
2:D:98:ALA:HB1	2:D:105:THR:HG21	1.86	0.57
1:A:795:HIS:CE1	1:A:797:ARG:HD3	2.39	0.57
2:D:45:LYS:NZ	2:D:46:ARG:NH1	2.43	0.57
1:C:730:ALA:HB1	1:C:765:ARG:CG	2.33	0.57
1:C:271:ILE:HG13	1:C:272:LYS:HB3	1.86	0.57
1:A:772:LYS:HE2	1:A:823:ALA:HB2	1.86	0.56
1:A:858:ASP:OD2	1:A:858:ASP:N	2.38	0.56
1:C:815:CYS:N	6:C:1207:HOH:O	2.38	0.56
1:C:882:GLN:HE21	2:D:256:ALA:N	1.95	0.56
2:B:102:HIS:CB	2:B:103:PRO:HD2	2.36	0.56
2:B:99:ASP:HB3	2:B:102:HIS:HB2	1.87	0.56
1:C:71:ILE:HD11	1:C:106:ILE:HD13	1.88	0.56
2:D:125:GLY:HA2	5:D:401:SAM:HG1	1.88	0.56
2:D:190:PRO:HD2	2:D:193:SER:HB2	1.88	0.56
1:A:71:ILE:HD11	1:A:106:ILE:HD13	1.88	0.55
1:A:38:LEU:HD21	1:A:41:LYS:HG3	1.89	0.55
1:A:797:ARG:HH21	1:A:800:GLN:HE22	1.53	0.55
1:A:198:SER:OG	1:A:199:ASP:N	2.38	0.55
1:A:416:GLU:OE2	1:A:461:LYS:NZ	2.26	0.55
1:C:882:GLN:HE22	2:D:255:MET:CA	2.19	0.55
1:A:441:ASN:HB2	6:A:1215:HOH:O	2.07	0.55
2:D:11:LEU:HD21	2:D:15:LYS:HD3	1.88	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:143:LYS:HZ2	1:A:187:LEU:HB2	1.72	0.55
2:B:151:THR:HG21	2:B:161:PHE:CD1	2.41	0.55
2:B:46:ARG:NH1	2:B:76:ASP:O	2.40	0.54
2:B:109:ILE:HG23	2:B:113:PHE:HE1	1.72	0.54
1:C:669:ASN:OD1	1:C:669:ASN:N	2.40	0.54
2:D:100:ILE:O	2:D:150:LEU:CD2	2.48	0.54
5:D:401:SAM:N	6:D:501:HOH:O	2.33	0.54
1:C:72:VAL:CG2	1:C:83:LEU:HD13	2.36	0.54
2:D:47:VAL:HG12	2:D:120:PHE:HD1	1.70	0.54
2:B:119:ASP:HA	2:B:156:LYS:HE2	1.89	0.54
2:B:146:SER:O	2:B:149:GLN:NE2	2.41	0.54
1:C:198:SER:OG	1:C:199:ASP:N	2.39	0.54
2:B:45:LYS:NZ	2:B:119:ASP:OD2	2.36	0.54
1:C:441:ASN:HB2	6:C:1302:HOH:O	2.08	0.54
1:C:414:LEU:HD12	1:C:437:LEU:HD23	1.89	0.54
2:B:84:LEU:HD21	2:B:99:ASP:HB2	1.89	0.53
2:B:246:GLU:OE1	2:B:246:GLU:HA	2.07	0.53
1:C:132:GLY:O	2:D:233:ASN:HB2	2.09	0.53
1:C:18:ASP:H	1:C:280:LYS:NZ	2.06	0.53
2:D:102:HIS:CG	2:D:103:PRO:HD3	2.44	0.53
2:B:45:LYS:HD3	2:B:46:ARG:H	1.73	0.53
1:A:730:ALA:HB2	6:A:1201:HOH:O	2.07	0.53
1:A:254:ARG:H	1:A:254:ARG:CD	2.21	0.53
1:C:928:SER:OG	1:C:968:GLY:O	2.25	0.53
2:D:84:LEU:HD21	2:D:99:ASP:HB2	1.90	0.52
1:A:254:ARG:H	1:A:254:ARG:HD2	1.75	0.52
2:B:35:GLN:HG3	2:B:257:CYS:O	2.10	0.52
2:D:120:PHE:HD2	2:D:160:THR:OG1	1.91	0.52
1:A:733:GLU:HG2	1:A:735:PHE:CZ	2.45	0.52
1:A:568:LYS:HD2	1:A:569:GLY:H	1.75	0.52
2:D:33:ASN:HD21	2:D:39:LEU:H	1.57	0.52
1:A:338:PHE:CD1	1:A:383:ASN:HA	2.45	0.52
1:C:134:ARG:HG2	2:D:233:ASN:OD1	2.10	0.52
2:D:19:GLN:HG2	6:D:505:HOH:O	2.10	0.52
2:D:48:VAL:HB	2:D:121:VAL:HG22	1.92	0.52
2:B:86:PRO:HG3	2:B:97:GLN:OE1	2.11	0.51
2:D:193:SER:OG	2:D:199:GLU:OE1	2.21	0.51
1:A:820:VAL:HG12	1:A:831:VAL:HB	1.93	0.51
1:C:568:LYS:HG3	1:C:588:LYS:HG2	1.92	0.51
1:C:858:ASP:N	1:C:858:ASP:OD2	2.43	0.51
2:D:100:ILE:HG12	5:D:401:SAM:N1	2.26	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:730:ALA:HB1	1:A:765:ARG:CG	2.41	0.51
1:C:215:SER:HB2	1:C:259:GLU:HA	1.93	0.51
1:C:730:ALA:HB2	1:C:765:ARG:CG	2.36	0.51
1:C:568:LYS:HD2	1:C:569:GLY:N	2.26	0.51
2:D:35:GLN:C	2:D:37:HIS:H	2.15	0.51
1:A:134:ARG:NH1	2:B:236:CYS:SG	2.83	0.50
1:A:735:PHE:HB3	1:A:751:ILE:CD1	2.41	0.50
1:C:353:LYS:HE3	1:C:404:ALA:HB3	1.93	0.50
2:D:35:GLN:HB2	2:D:257:CYS:O	2.12	0.50
2:B:103:PRO:CG	2:B:104:LYS:H	2.13	0.50
1:C:134:ARG:CG	2:D:233:ASN:HA	2.41	0.50
1:C:838:HIS:HD2	6:C:1259:HOH:O	1.94	0.50
2:D:39:LEU:HD23	2:D:44:LEU:HD22	1.94	0.50
2:B:36:PHE:HZ	2:B:186:VAL:HG11	1.76	0.50
2:D:246:GLU:OE1	2:D:246:GLU:HA	2.11	0.50
1:A:565:LYS:NZ	6:A:1214:HOH:O	2.44	0.49
1:C:709:LYS:HD3	1:C:735:PHE:CZ	2.47	0.49
2:D:45:LYS:HA	2:D:77:ARG:HG2	1.94	0.49
1:A:838:HIS:HD2	6:A:1292:HOH:O	1.94	0.49
2:B:45:LYS:HD3	2:B:119:ASP:OD1	2.13	0.49
2:D:36:PHE:HD1	2:D:205:LEU:HD11	1.76	0.49
1:A:928:SER:OG	1:A:968:GLY:O	2.29	0.49
1:A:883:SER:OG	1:A:884:GLY:N	2.45	0.49
1:A:364:LEU:HB3	1:A:365:LEU:HD23	1.94	0.49
2:B:45:LYS:HA	2:B:77:ARG:HG2	1.95	0.49
1:C:26:PRO:HD3	1:C:979:THR:CG2	2.42	0.49
2:D:186:VAL:CG1	2:D:255:MET:HB2	2.42	0.49
1:A:42:CYS:HB3	6:A:1273:HOH:O	2.12	0.49
1:A:882:GLN:HE21	2:B:255:MET:HA	1.78	0.49
2:B:45:LYS:O	2:B:77:ARG:HA	2.12	0.49
2:B:86:PRO:HB2	2:D:86:PRO:HB2	1.95	0.49
1:A:951:VAL:HB	1:A:963:TRP:HB2	1.95	0.49
2:B:103:PRO:HG2	2:B:104:LYS:N	2.25	0.49
1:C:666:SER:HB2	1:C:668:THR:HG22	1.95	0.49
1:C:883:SER:OG	1:C:884:GLY:N	2.46	0.49
1:A:797:ARG:HG3	1:A:797:ARG:HH11	1.79	0.48
2:D:106:LEU:O	2:D:110:LEU:HD23	2.13	0.48
1:A:178:HIS:CE1	1:A:204:ARG:HD2	2.48	0.48
2:B:25:SER:HA	2:B:28:LYS:HD2	1.94	0.48
1:A:3:ASP:HA	1:A:1011:TRP:CD1	2.49	0.48
1:A:114:LYS:NZ	2:B:233:ASN:OD1	2.45	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:48:VAL:HB	2:B:121:VAL:HG22	1.96	0.48
1:C:675:LYS:HG2	1:C:675:LYS:H	1.40	0.48
1:C:789:SER:HA	1:C:816:LEU:HG	1.96	0.48
1:A:215:SER:HB2	1:A:259:GLU:HA	1.95	0.48
2:D:36:PHE:CE2	2:D:257:CYS:SG	3.02	0.48
2:D:45:LYS:HZ3	2:D:76:ASP:HB3	1.79	0.47
1:A:576:PHE:HB2	3:A:1101:EPE:H91	1.96	0.47
2:D:98:ALA:HB1	2:D:105:THR:CG2	2.45	0.47
2:B:46:ARG:HB3	2:B:113:PHE:CE2	2.49	0.47
2:D:103:PRO:HG2	2:D:104:LYS:N	2.22	0.47
1:A:112:TYR:CG	1:A:134:ARG:NH2	2.82	0.47
1:C:793:ILE:HB	1:C:807:MET:HB3	1.96	0.47
1:A:214:LEU:HD22	1:A:257:VAL:HB	1.97	0.47
2:D:15:LYS:HB3	2:D:15:LYS:HE3	1.68	0.47
2:D:35:GLN:HE22	2:D:255:MET:CE	2.28	0.47
2:D:82:VAL:HG21	2:D:109:ILE:HD11	1.97	0.47
2:B:45:LYS:HG3	2:B:46:ARG:NH2	2.29	0.47
1:C:14:LYS:HE3	6:C:1347:HOH:O	2.13	0.47
1:C:546:LYS:HB2	1:C:546:LYS:HE3	1.72	0.47
1:C:918:ASP:OD2	1:C:922:LYS:HG2	2.14	0.47
3:A:1101:EPE:H21	6:A:1328:HOH:O	2.14	0.47
1:C:178:HIS:CE1	1:C:204:ARG:HD2	2.49	0.47
1:C:820:VAL:HG12	1:C:831:VAL:HB	1.96	0.47
2:D:47:VAL:CG2	2:D:79:ILE:HG12	2.45	0.47
2:B:117:LYS:CD	2:B:157:LYS:HZ1	2.28	0.47
2:B:126:ALA:HB2	2:B:164:LYS:HE2	1.97	0.47
1:C:730:ALA:HA	1:C:765:ARG:HG2	1.96	0.47
2:B:186:VAL:CG1	2:B:255:MET:HB2	2.45	0.47
2:B:36:PHE:CD2	2:B:203:VAL:HG11	2.50	0.46
1:A:709:LYS:HD3	1:A:735:PHE:CZ	2.51	0.46
2:B:120:PHE:HD2	2:B:160:THR:OG1	1.99	0.46
1:C:18:ASP:HA	1:C:33:TYR:CE2	2.51	0.46
1:C:364:LEU:HD12	1:C:364:LEU:HA	1.74	0.46
1:C:476:THR:HG22	1:C:483:LEU:HB2	1.96	0.46
2:B:28:LYS:O	2:B:32:LEU:HG	2.14	0.46
2:B:45:LYS:HG2	2:B:46:ARG:NE	2.30	0.46
1:C:214:LEU:HD22	1:C:257:VAL:HB	1.98	0.46
2:D:45:LYS:N	2:D:119:ASP:OD1	2.45	0.46
1:C:672:ASP:O	1:C:675:LYS:CD	2.64	0.46
1:C:774:ILE:HG22	1:C:774:ILE:O	2.14	0.46
2:B:35:GLN:C	2:B:37:HIS:N	2.68	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:45:LYS:CE	2:D:46:ARG:HH12	2.27	0.46
1:A:774:ILE:O	1:A:774:ILE:HG22	2.15	0.46
2:B:183:ASP:OD2	2:B:208:ASN:ND2	2.42	0.46
1:A:479:ARG:CG	1:A:480:ASN:N	2.61	0.46
2:B:237:ILE:HA	2:B:240:LYS:HG2	1.98	0.46
1:A:825:LYS:HZ3	1:A:921:ASN:HD21	1.63	0.46
1:C:72:VAL:HG22	1:C:83:LEU:HD13	1.96	0.46
1:C:951:VAL:HB	1:C:963:TRP:HB2	1.97	0.46
2:D:45:LYS:NZ	2:D:76:ASP:HB3	2.30	0.46
2:B:103:PRO:CG	2:B:104:LYS:N	2.78	0.46
1:A:730:ALA:CA	1:A:765:ARG:HG2	2.45	0.45
1:A:1006:VAL:HG21	2:B:133:HIS:CE1	2.51	0.45
2:B:243:HIS:ND1	2:B:244:TRP:HD1	2.14	0.45
1:C:3:ASP:HA	1:C:1011:TRP:CD1	2.51	0.45
1:A:372:SER:OG	1:A:390:ASN:OD1	2.21	0.45
1:A:793:ILE:HB	1:A:807:MET:HB3	1.98	0.45
1:A:940:THR:OG1	1:A:954:THR:HG23	2.16	0.45
2:B:36:PHE:CD1	2:B:205:LEU:HD11	2.51	0.45
2:B:96:LEU:HD22	2:B:108:ARG:CZ	2.47	0.45
2:B:98:ALA:HB1	2:B:105:THR:OG1	2.17	0.45
1:A:730:ALA:CB	1:A:765:ARG:CG	2.83	0.45
2:B:33:ASN:C	2:B:35:GLN:H	2.19	0.45
2:B:180:TYR:CE2	2:B:244:TRP:HE3	2.34	0.45
1:C:271:ILE:HA	1:C:272:LYS:HA	1.40	0.45
2:D:28:LYS:HE2	2:D:199:GLU:OE1	2.16	0.45
1:A:353:LYS:HB2	1:A:364:LEU:HD22	1.99	0.45
1:A:797:ARG:NH2	1:A:800:GLN:HE22	2.15	0.45
2:B:102:HIS:CB	2:B:103:PRO:CD	2.95	0.45
1:C:811:TYR:CE1	1:C:840:VAL:HG21	2.51	0.45
1:A:17:ASN:O	1:A:18:ASP:HB2	2.17	0.45
1:A:353:LYS:HE3	1:A:404:ALA:HB3	1.99	0.45
2:D:86:PRO:HG3	2:D:97:GLN:NE2	2.32	0.45
2:D:159:GLY:O	2:D:207:TYR:HB2	2.16	0.45
1:A:939:ILE:HA	1:A:955:SER:HB3	1.98	0.45
1:C:230:PHE:CD2	1:C:279:VAL:HG11	2.51	0.45
1:A:825:LYS:NZ	1:A:921:ASN:HD21	2.15	0.45
2:B:35:GLN:CG	2:B:257:CYS:O	2.65	0.45
1:C:148:ASP:OD1	6:C:1201:HOH:O	2.21	0.45
1:C:338:PHE:CD1	1:C:383:ASN:HA	2.52	0.45
1:C:765:ARG:H	1:C:787:SER:HG	1.64	0.44
2:B:33:ASN:C	2:B:35:GLN:N	2.70	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:125:GLY:HA2	5:D:401:SAM:H5'2	1.98	0.44
1:C:419:LYS:O	1:C:439:SER:HB2	2.18	0.44
1:C:839:LEU:HB3	1:C:925:LEU:HD21	2.00	0.44
1:A:479:ARG:NH1	1:A:546:LYS:HD3	2.33	0.44
2:B:157:LYS:HZ3	2:B:157:LYS:HG2	1.60	0.44
1:A:546:LYS:HE3	1:A:546:LYS:HB2	1.67	0.44
2:D:28:LYS:O	2:D:32:LEU:HG	2.16	0.44
2:D:118:ALA:O	2:D:156:LYS:HB3	2.18	0.44
1:A:730:ALA:HA	1:A:765:ARG:HG2	1.99	0.44
2:B:36:PHE:CE2	2:B:203:VAL:HG11	2.53	0.44
2:B:106:LEU:HD12	2:B:106:LEU:O	2.17	0.44
2:B:117:LYS:CE	2:B:157:LYS:HZ1	2.31	0.44
1:C:730:ALA:CA	1:C:765:ARG:HG2	2.47	0.44
2:D:176:SER:HB2	2:D:245:ASN:OD1	2.17	0.44
1:A:568:LYS:HG3	1:A:588:LYS:HG2	1.99	0.44
2:D:102:HIS:ND1	2:D:103:PRO:CD	2.81	0.44
1:A:26:PRO:HD3	1:A:979:THR:CG2	2.47	0.44
1:A:306:HIS:HE1	6:A:1322:HOH:O	2.01	0.44
1:A:543:GLU:O	1:A:558:TYR:HA	2.18	0.44
2:B:100:ILE:HG12	5:B:401:SAM:N1	2.32	0.44
1:A:187:LEU:HD22	1:A:194:VAL:HG22	2.00	0.43
1:A:611:SER:OG	1:A:613:ARG:NE	2.51	0.43
2:B:149:GLN:OE1	2:B:150:LEU:HD12	2.18	0.43
2:B:172:ASP:HB2	2:B:249:ARG:HB3	2.00	0.43
2:D:46:ARG:HG2	2:D:113:PHE:HD1	1.79	0.43
1:A:18:ASP:HA	1:A:33:TYR:CE2	2.54	0.43
1:A:613:ARG:NH2	4:A:1106:SO4:O1	2.51	0.43
1:A:825:LYS:NZ	1:A:921:ASN:ND2	2.66	0.43
2:D:45:LYS:NZ	2:D:76:ASP:CB	2.81	0.43
1:A:58:GLU:HB2	6:A:1229:HOH:O	2.18	0.43
1:A:419:LYS:O	1:A:439:SER:HB2	2.19	0.43
2:B:124:ASP:HB3	5:B:401:SAM:N	2.33	0.43
2:B:186:VAL:HG12	2:B:187:CYS:N	2.34	0.43
1:C:662:ILE:HG13	1:C:987:GLU:HG2	2.00	0.43
1:C:1006:VAL:HG21	2:D:133:HIS:CE1	2.53	0.43
1:A:117:ILE:HD12	1:A:117:ILE:N	2.33	0.43
2:D:172:ASP:HB2	2:D:249:ARG:HB3	2.01	0.43
1:A:230:PHE:CE1	1:A:237:ILE:HB	2.54	0.43
1:A:271:ILE:HA	1:A:272:LYS:HA	1.39	0.43
1:C:672:ASP:O	1:C:675:LYS:CG	2.63	0.43
1:A:779:ASP:CG	1:A:797:ARG:HA	2.39	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:441:ASN:HB3	1:C:444:GLU:HG3	2.00	0.43
2:D:45:LYS:HE3	2:D:46:ARG:HH22	1.83	0.43
1:C:797:ARG:HE	1:C:797:ARG:HB3	1.18	0.43
2:D:81:SER:OG	2:D:95:THR:HG22	2.19	0.43
2:D:44:LEU:HD23	2:D:65:LEU:HD11	1.99	0.43
2:D:86:PRO:HG3	2:D:97:GLN:HE22	1.83	0.43
1:A:991:ASN:ND2	1:A:998:GLU:O	2.51	0.43
2:B:46:ARG:CZ	2:B:116:GLU:OE1	2.66	0.43
2:D:151:THR:HG21	2:D:161:PHE:CD1	2.53	0.43
1:C:479:ARG:NH1	1:C:546:LYS:HD3	2.34	0.42
2:D:178:LEU:O	2:D:185:ILE:HD11	2.19	0.42
1:A:941:SER:N	1:A:954:THR:HG22	2.35	0.42
2:B:17:LYS:HA	2:B:17:LYS:HD3	1.88	0.42
2:B:82:VAL:HG21	2:B:109:ILE:HD11	2.01	0.42
1:A:769:PHE:HA	1:A:783:VAL:O	2.19	0.42
2:B:164:LYS:HD2	2:B:199:GLU:HG3	2.01	0.42
1:C:809:GLY:HA3	1:C:871:LEU:HB3	2.02	0.42
2:D:102:HIS:HB3	2:D:103:PRO:HD3	2.00	0.42
2:D:155:LEU:HD21	2:D:159:GLY:O	2.20	0.42
1:A:621:THR:O	1:A:622:ASP:HB2	2.19	0.42
1:C:17:ASN:O	1:C:18:ASP:HB2	2.20	0.42
1:C:666:SER:CB	1:C:668:THR:HG22	2.50	0.42
1:C:811:TYR:O	1:C:812:LYS:HG2	2.20	0.42
1:A:666:SER:HB2	1:A:668:THR:HG22	2.01	0.42
1:A:934:ALA:HB3	1:A:939:ILE:HD11	2.02	0.42
2:B:159:GLY:O	2:B:207:TYR:HB2	2.18	0.42
1:C:187:LEU:HD22	1:C:194:VAL:HG22	2.02	0.42
1:C:742:LYS:HE3	1:C:742:LYS:HB3	1.77	0.42
2:D:126:ALA:HB2	2:D:164:LYS:HE2	2.01	0.42
1:A:685:GLU:O	1:A:685:GLU:HG2	2.20	0.42
2:B:100:ILE:O	2:B:150:LEU:CD1	2.64	0.42
2:B:198:LEU:HD23	2:B:198:LEU:HA	1.91	0.42
2:D:36:PHE:CD1	2:D:205:LEU:HD11	2.54	0.42
2:D:83:ASP:O	2:D:97:GLN:HA	2.20	0.42
1:C:775:SER:OG	1:C:776:GLN:N	2.53	0.42
2:B:241:LEU:HD23	2:B:241:LEU:HA	1.87	0.42
2:D:102:HIS:CB	2:D:103:PRO:HD3	2.50	0.42
1:A:765:ARG:H	1:A:787:SER:HG	1.66	0.41
1:C:570:PHE:O	1:C:586:GLY:HA3	2.19	0.41
1:C:758:SER:O	1:C:758:SER:OG	2.37	0.41
1:C:254:ARG:HA	1:C:254:ARG:HD3	1.78	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:715:GLN:HG3	1:C:729:SER:O	2.21	0.41
3:A:1101:EPE:H101	3:A:1101:EPE:H22	1.79	0.41
1:C:977:ARG:HH11	1:C:977:ARG:HG2	1.84	0.41
1:A:315:LEU:HD12	1:A:315:LEU:HA	1.89	0.41
2:B:148:LEU:HD23	2:B:148:LEU:HA	1.85	0.41
1:A:29:HIS:HB3	1:A:31:TYR:CE1	2.55	0.41
1:C:621:THR:O	1:C:622:ASP:HB2	2.20	0.41
1:C:735:PHE:HB3	1:C:751:ILE:CD1	2.51	0.41
2:D:35:GLN:O	2:D:36:PHE:HB2	2.20	0.41
2:D:46:ARG:CZ	2:D:116:GLU:OE2	2.68	0.41
2:D:103:PRO:CD	2:D:105:THR:HG23	2.44	0.41
2:D:186:VAL:HG12	2:D:187:CYS:N	2.35	0.41
2:D:240:LYS:HE2	2:D:244:TRP:CH2	2.56	0.41
2:B:28:LYS:HE2	2:B:199:GLU:OE1	2.20	0.41
2:D:106:LEU:HB3	2:D:150:LEU:CD1	2.50	0.41
1:A:783:VAL:HB	1:A:793:ILE:HD13	2.03	0.41
1:A:491:LEU:HD22	1:A:506:ILE:HD12	2.03	0.41
2:B:53:ALA:HB3	5:B:401:SAM:O3'	2.21	0.41
1:C:117:ILE:N	1:C:117:ILE:HD12	2.36	0.41
1:C:356:ASP:O	1:C:359:LYS:NZ	2.38	0.41
1:C:555:ARG:HA	1:C:555:ARG:HD3	1.91	0.41
2:D:84:LEU:HD12	5:D:401:SAM:C6	2.51	0.41
1:A:725:MET:CE	1:A:737:TRP:CD1	3.04	0.41
1:A:815:CYS:O	1:A:835:THR:HG23	2.22	0.41
2:B:193:SER:OG	2:B:199:GLU:OE1	2.23	0.41
2:D:234:LYS:HG2	2:D:235:LEU:N	2.36	0.41
1:A:476:THR:HG22	1:A:483:LEU:HB2	2.02	0.40
1:A:839:LEU:HB3	1:A:925:LEU:HD21	2.03	0.40
1:A:878:LEU:HD21	1:A:927:THR:HB	2.02	0.40
2:B:178:LEU:O	2:B:185:ILE:HD11	2.21	0.40
2:B:209:PRO:HA	2:B:210:PRO:HD2	1.98	0.40
1:C:960:ILE:O	1:C:975:LYS:HA	2.21	0.40
1:A:342:VAL:HB	1:A:354:TYR:HB3	2.03	0.40
1:A:725:MET:HE1	1:A:727:SER:HB3	2.03	0.40
1:A:772:LYS:NZ	6:A:1226:HOH:O	2.54	0.40
2:B:44:LEU:HA	2:B:119:ASP:OD1	2.21	0.40
1:C:771:VAL:HG12	1:C:782:LEU:HD12	2.03	0.40
2:B:46:ARG:NE	2:B:116:GLU:OE1	2.54	0.40
2:D:36:PHE:HZ	2:D:186:VAL:HG11	1.87	0.40
2:D:102:HIS:ND1	2:D:103:PRO:N	2.69	0.40
2:D:234:LYS:N	2:D:234:LYS:HD3	2.36	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:202:SER:HB3	1:A:218:TRP:CZ3	2.56	0.40
1:A:789:SER:HA	1:A:816:LEU:HG	2.04	0.40
1:C:611:SER:OG	1:C:613:ARG:NH1	2.55	0.40
2:D:25:SER:HA	2:D:28:LYS:HD2	2.02	0.40
2:D:51:CYS:HB3	5:D:401:SAM:O4'	2.21	0.40
1:A:230:PHE:CD2	1:A:279:VAL:HG11	2.56	0.40

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

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:259:SER:OG	1:C:58:GLU:OE2[4_445]	2.19	0.01

5.3 Torsion angles (i)

5.3.1 Protein backbone (i)

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	990/1028 (96%)	944 (95%)	44 (4%)	2 (0%)	47 58
1	C	990/1028 (96%)	942 (95%)	46 (5%)	2 (0%)	47 58
2	B	227/325 (70%)	209 (92%)	12 (5%)	6 (3%)	5 3
2	D	227/325 (70%)	210 (92%)	13 (6%)	4 (2%)	8 7
All	All	2434/2706 (90%)	2305 (95%)	115 (5%)	14 (1%)	25 30

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	610	GLY
2	B	103	PRO
1	C	610	GLY
2	D	103	PRO

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Mol	Chain	Res	Type
2	B	34	ASP
2	B	210	PRO
2	B	245	ASN
1	A	812	LYS
2	B	190	PRO
1	C	812	LYS
2	D	190	PRO
2	B	206	GLY
2	D	102	HIS
2	D	206	GLY

5.3.2 Protein sidechains (i)

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	900/927 (97%)	879 (98%)	21 (2%)	50	66
1	C	900/927 (97%)	877 (97%)	23 (3%)	46	62
2	B	203/289 (70%)	192 (95%)	11 (5%)	22	30
2	D	203/289 (70%)	195 (96%)	8 (4%)	32	45
All	All	2206/2432 (91%)	2143 (97%)	63 (3%)	42	57

All (63) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	13	VAL
1	A	239	VAL
1	A	254	ARG
1	A	256	ASN
1	A	365	LEU
1	A	403	SER
1	A	489	SER
1	A	655	ARG
1	A	731	ARG
1	A	776	GLN
1	A	787	SER

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Mol	Chain	Res	Type
1	A	789	SER
1	A	810	ARG
1	A	817	PHE
1	A	820	VAL
1	A	838	HIS
1	A	858	ASP
1	A	900	THR
1	A	908	ASN
1	A	954	THR
1	A	979	THR
2	B	24	ARG
2	B	38	PHE
2	B	45	LYS
2	B	73	ASP
2	B	112	LEU
2	B	143	LEU
2	B	148	LEU
2	B	181	LEU
2	B	187	CYS
2	B	234	LYS
2	B	244	TRP
1	C	13	VAL
1	C	28	ILE
1	C	239	VAL
1	C	256	ASN
1	C	365	LEU
1	C	403	SER
1	C	489	SER
1	C	655	ARG
1	C	696	ASN
1	C	697	ARG
1	C	776	GLN
1	C	787	SER
1	C	789	SER
1	C	797	ARG
1	C	817	PHE
1	C	858	ASP
1	C	865	ASP
1	C	900	THR
1	C	908	ASN
1	C	927	THR
1	C	954	THR

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Mol	Chain	Res	Type
1	C	979	THR
1	C	997	SER
2	D	9	ARG
2	D	24	ARG
2	D	38	PHE
2	D	112	LEU
2	D	148	LEU
2	D	157	LYS
2	D	187	CYS
2	D	234	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	454	GLN
1	A	678	HIS
1	A	696	ASN
1	A	800	GLN
1	A	882	GLN
1	A	921	ASN
1	A	991	ASN
2	B	35	GLN
2	B	250	ASN
1	C	702	GLN
1	C	776	GLN
1	C	882	GLN
2	D	33	ASN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

14 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
4	SO4	C	1104	-	4,4,4	0.26	0	6,6,6	0.24	0
4	SO4	A	1104	-	4,4,4	0.24	0	6,6,6	0.33	0
5	SAM	B	401	-	23,29,29	1.37	2 (8%)	20,42,42	2.06	4 (20%)
4	SO4	A	1102	-	4,4,4	0.28	0	6,6,6	0.15	0
4	SO4	A	1103	-	4,4,4	0.36	0	6,6,6	0.30	0
3	EPE	C	1101	-	15,15,15	0.75	1 (6%)	19,20,20	1.90	4 (21%)
4	SO4	A	1106	-	4,4,4	0.32	0	6,6,6	0.09	0
5	SAM	D	401	-	23,29,29	1.37	3 (13%)	20,42,42	2.05	5 (25%)
4	SO4	C	1103	-	4,4,4	0.17	0	6,6,6	0.44	0
3	EPE	A	1101	-	15,15,15	0.66	0	19,20,20	1.65	2 (10%)
4	SO4	C	1105	-	4,4,4	0.27	0	6,6,6	0.18	0
4	SO4	C	1106	-	4,4,4	0.26	0	6,6,6	0.21	0
4	SO4	C	1102	-	4,4,4	0.25	0	6,6,6	0.31	0
4	SO4	A	1105	-	4,4,4	0.34	0	6,6,6	0.12	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
5	SAM	D	401	-	-	2/13/33/33	0/3/3/3
5	SAM	B	401	-	-	4/13/33/33	0/3/3/3
3	EPE	A	1101	-	-	6/9/19/19	0/1/1/1
3	EPE	C	1101	-	-	6/9/19/19	0/1/1/1

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	D	401	SAM	C2-N3	4.66	1.39	1.32
5	B	401	SAM	C2-N3	4.65	1.39	1.32
5	D	401	SAM	C2-N1	2.84	1.39	1.33
5	B	401	SAM	C2-N1	2.73	1.38	1.33
3	C	1101	EPE	C10-S	2.49	1.81	1.77
5	D	401	SAM	OXT-C	-2.01	1.24	1.30

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
5	B	401	SAM	N3-C2-N1	-6.11	120.38	128.67
5	D	401	SAM	N3-C2-N1	-5.99	120.54	128.67
3	C	1101	EPE	C5-N4-C3	5.16	119.95	108.84
5	B	401	SAM	O4'-C1'-N9	4.63	114.89	108.75
3	A	1101	EPE	C5-N4-C3	4.53	118.60	108.84
5	D	401	SAM	O4'-C1'-N9	4.47	114.68	108.75
3	C	1101	EPE	C7-N4-C3	3.17	119.69	111.24
3	C	1101	EPE	C7-N4-C5	3.17	119.67	111.24
5	D	401	SAM	C4'-O4'-C1'	-2.74	107.41	109.92
3	C	1101	EPE	O1S-S-C10	2.48	110.47	106.73
5	B	401	SAM	OXT-C-O	-2.47	118.47	124.08
5	D	401	SAM	OXT-C-O	-2.33	118.78	124.08
3	A	1101	EPE	C9-N1-C6	-2.33	105.02	111.24
5	B	401	SAM	C4-C5-N7	-2.33	106.88	109.34
5	D	401	SAM	C4-C5-N7	-2.32	106.89	109.34

There are no chirality outliers.

All (18) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	1101	EPE	C9-C10-S-O1S
3	A	1101	EPE	N4-C7-C8-O8
3	C	1101	EPE	C9-C10-S-O3S
3	C	1101	EPE	C8-C7-N4-C3
3	A	1101	EPE	C10-C9-N1-C2
3	C	1101	EPE	C10-C9-N1-C2
3	A	1101	EPE	C9-C10-S-O2S
3	C	1101	EPE	C9-C10-S-O1S
3	C	1101	EPE	C9-C10-S-O2S
3	A	1101	EPE	S-C10-C9-N1
5	B	401	SAM	CA-CB-CG-SD
5	B	401	SAM	O4'-C4'-C5'-SD

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Mol	Chain	Res	Type	Atoms
3	C	1101	EPE	C10-C9-N1-C6
3	A	1101	EPE	C9-C10-S-O3S
5	D	401	SAM	O-C-CA-N
5	D	401	SAM	OXT-C-CA-N
5	B	401	SAM	OXT-C-CA-N
5	B	401	SAM	O-C-CA-N

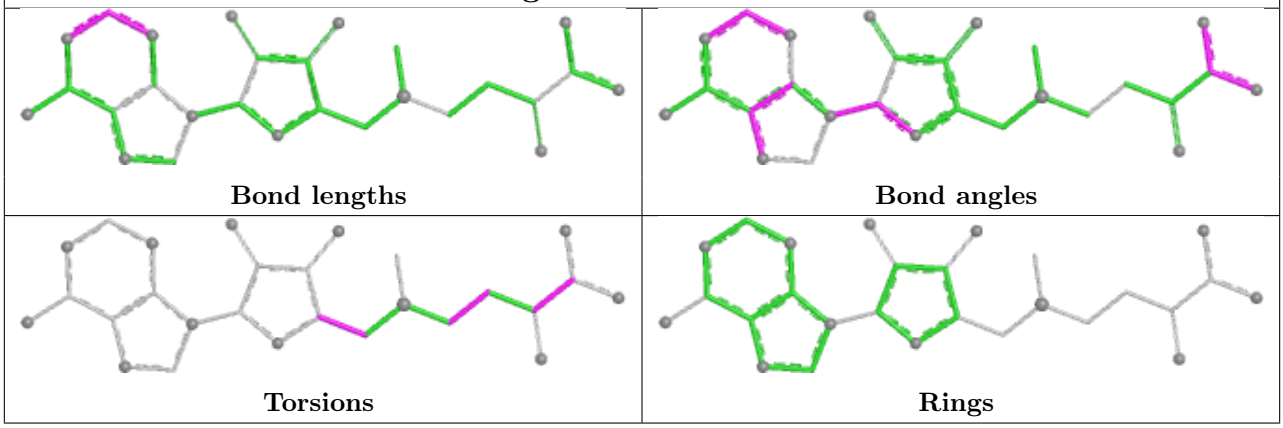
There are no ring outliers.

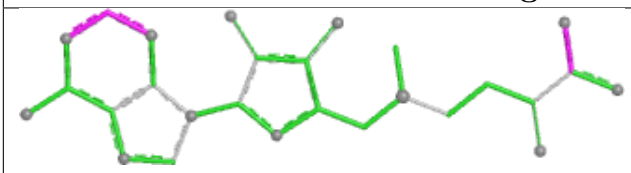
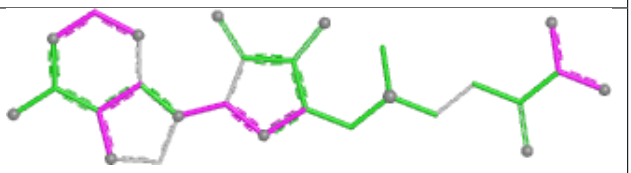
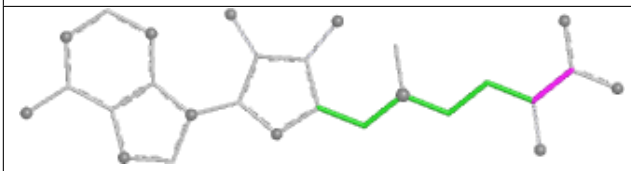
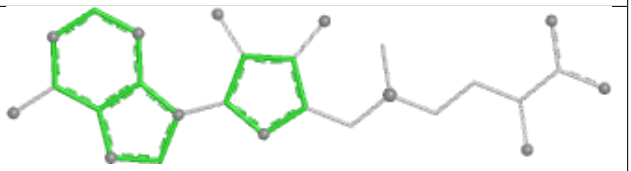
5 monomers are involved in 16 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	C	1104	SO4	1	0
5	B	401	SAM	5	0
4	A	1106	SO4	1	0
5	D	401	SAM	6	0
3	A	1101	EPE	3	0

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

Ligand SAM B 401



Ligand SAM D 401	
 Bond lengths	 Bond angles
 Torsions	 Rings

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	998/1028 (97%)	0.02	3 (0%)	94	96	29, 44, 67, 95	0
1	C	998/1028 (97%)	0.03	8 (0%)	86	89	29, 43, 67, 94	0
2	B	231/325 (71%)	1.02	49 (21%)	0	1	34, 72, 97, 106	0
2	D	231/325 (71%)	1.09	52 (22%)	0	1	34, 71, 97, 107	0
All	All	2458/2706 (90%)	0.22	112 (4%)	32	40	29, 46, 84, 107	0

All (112) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	856	SER	7.4
2	D	235	LEU	7.3
2	B	257	CYS	7.2
2	B	36	PHE	6.8
2	D	211	SER	5.7
1	A	856	SER	5.6
2	B	211	SER	5.2
2	D	258	GLY	5.0
2	B	210	PRO	4.9
2	D	245	ASN	4.9
2	B	243	HIS	4.9
2	D	257	CYS	4.6
2	B	244	TRP	4.5
2	D	243	HIS	4.5
2	D	180	TYR	4.5
2	D	210	PRO	4.4
2	B	152	ALA	4.3
2	D	244	TRP	4.2
2	D	251	ILE	4.2
2	D	36	PHE	4.1
2	D	236	CYS	4.0

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Mol	Chain	Res	Type	RSRZ
2	B	159	GLY	4.0
1	A	669	ASN	3.9
2	D	205	LEU	3.9
2	D	159	GLY	3.8
2	D	14	ARG	3.8
2	B	113	PHE	3.8
1	A	668	THR	3.8
2	B	205	LEU	3.8
2	B	110	LEU	3.8
2	D	181	LEU	3.7
1	C	669	ASN	3.7
2	D	107	ALA	3.7
2	D	250	ASN	3.6
2	B	251	ILE	3.6
2	B	255	MET	3.5
2	B	180	TYR	3.5
2	B	256	ALA	3.5
2	B	235	LEU	3.4
2	D	106	LEU	3.3
2	B	247	GLU	3.3
2	D	112	LEU	3.3
2	D	241	LEU	3.3
2	B	203	VAL	3.2
2	D	113	PHE	3.2
2	D	110	LEU	3.1
2	B	116	GLU	3.0
2	D	238	SER	3.0
2	B	107	ALA	2.9
2	B	46	ARG	2.9
2	D	39	LEU	2.9
2	B	181	LEU	2.9
2	B	179	GLY	2.8
2	B	236	CYS	2.8
2	B	250	ASN	2.8
2	D	103	PRO	2.8
2	B	44	LEU	2.8
1	C	668	THR	2.8
2	D	203	VAL	2.7
2	D	242	SER	2.7
2	B	237	ILE	2.7
2	B	35	GLN	2.7
2	B	38	PHE	2.7

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Mol	Chain	Res	Type	RSRZ
2	D	47	VAL	2.7
2	D	255	MET	2.7
2	B	68	GLU	2.7
2	D	104	LYS	2.6
2	B	112	LEU	2.6
1	C	730	ALA	2.6
2	D	72	SER	2.6
2	B	245	ASN	2.6
2	D	44	LEU	2.6
2	B	248	GLU	2.6
2	D	149	GLN	2.5
2	D	68	GLU	2.5
2	D	252	ALA	2.5
1	C	731	ARG	2.4
1	C	257	VAL	2.4
2	D	38	PHE	2.4
1	C	131	GLY	2.4
2	D	45	LYS	2.4
2	B	14	ARG	2.4
1	C	855	ILE	2.3
2	D	152	ALA	2.3
2	D	190	PRO	2.3
2	D	248	GLU	2.3
2	D	196	THR	2.3
2	D	116	GLU	2.3
2	D	183	ASP	2.3
2	D	247	GLU	2.3
2	B	178	LEU	2.3
2	B	258	GLY	2.3
2	B	183	ASP	2.3
2	B	154	ILE	2.3
2	B	253	GLU	2.2
2	B	246	GLU	2.2
2	B	75	GLU	2.2
2	D	102	HIS	2.2
2	D	75	GLU	2.2
2	B	234	LYS	2.1
2	D	234	LYS	2.1
2	B	212	ASN	2.1
2	D	186	VAL	2.1
2	B	241	LEU	2.1
2	B	114	GLY	2.1

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Mol	Chain	Res	Type	RSRZ
2	B	85	GLN	2.1
2	B	103	PRO	2.1
2	D	179	GLY	2.1
2	D	158	GLY	2.0
2	B	190	PRO	2.0
2	B	120	PHE	2.0
2	D	249	ARG	2.0

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

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

6.3 Carbohydrates (i)

There are no monosaccharides 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. 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.

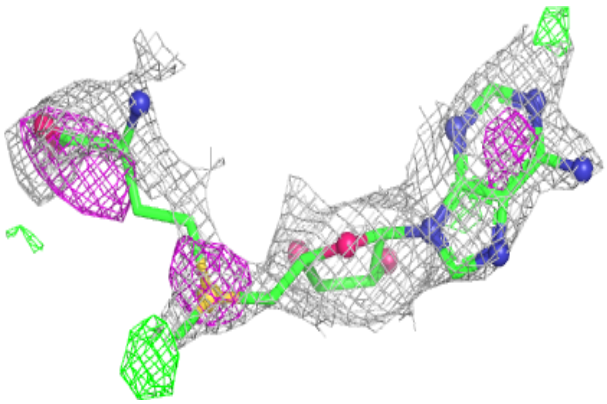
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
5	SAM	B	401	27/27	0.74	0.32	55,74,82,87	0
5	SAM	D	401	27/27	0.76	0.40	55,73,83,85	0
4	SO4	C	1105	5/5	0.90	0.12	76,77,82,85	0
4	SO4	A	1106	5/5	0.93	0.13	71,73,80,84	0
3	EPE	A	1101	15/15	0.93	0.20	51,62,84,91	0
4	SO4	C	1106	5/5	0.93	0.14	71,73,83,86	0
4	SO4	A	1104	5/5	0.93	0.13	77,78,81,83	0
4	SO4	A	1105	5/5	0.93	0.16	71,76,81,84	0
3	EPE	C	1101	15/15	0.95	0.19	54,65,87,90	0
4	SO4	C	1103	5/5	0.95	0.08	59,62,66,74	0
4	SO4	A	1103	5/5	0.97	0.07	58,64,72,76	0
4	SO4	C	1104	5/5	0.97	0.14	73,75,83,86	0
4	SO4	A	1102	5/5	0.98	0.14	38,43,47,47	0
4	SO4	C	1102	5/5	0.99	0.15	37,43,47,50	0

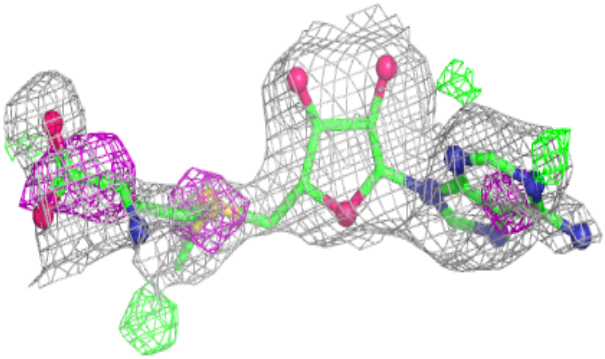
The following is a graphical depiction of the model fit to experimental electron density of all

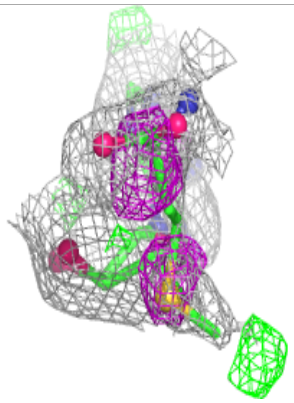
instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around SAM B 401:

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

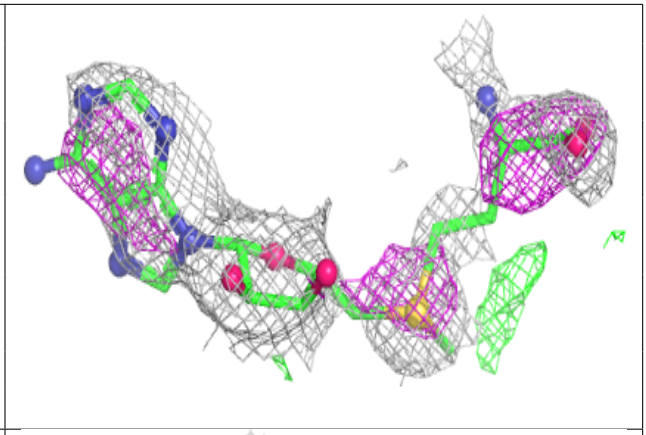


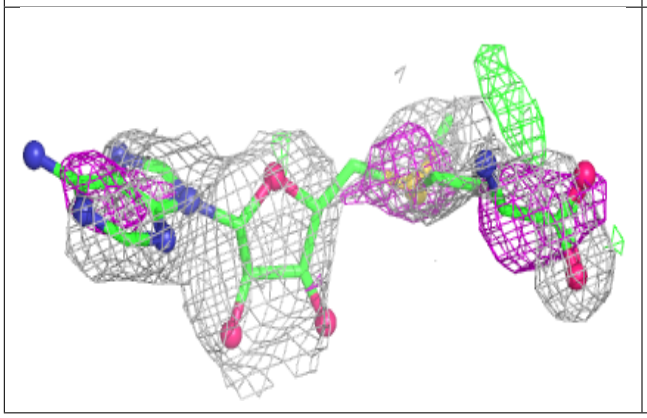


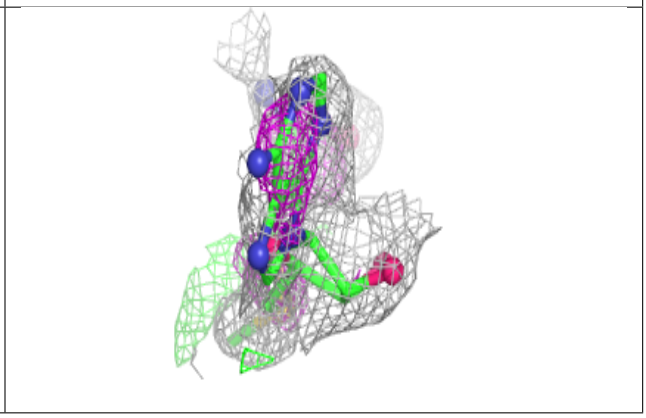


Electron density around SAM D 401:

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







6.5 Other polymers (i)

There are no such residues in this entry.