



Full wwPDB X-ray Structure Validation Report

Jun 18, 2014 – 09:30 PM EDT

PDB ID : 4L6I
Title : Methylthioadenosine phosphorylase from *Schistosoma mansoni* in complex with adenine
Authors : Torini, J.R.; DeMarco, R.; Brandao-Neto, J.; Pereira, H.M.
Deposited on : 2013-06-12
Resolution : 2.10 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

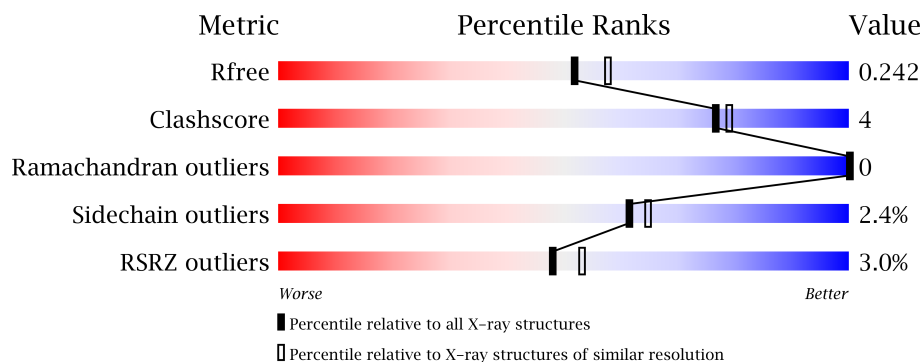
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : **FAILED**
Xtriage (Phenix) : dev-1439
EDS : stable23161
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable23161

1 Overall quality at a glance

The reported resolution of this entry is 2.10 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}	66092	3012 (2.10-2.10)
Clashscore	79885	3649 (2.10-2.10)
Ramachandran outliers	78287	3610 (2.10-2.10)
Sidechain outliers	78261	3611 (2.10-2.10)
RSRZ outliers	66119	3013 (2.10-2.10)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	320	
1	B	320	
1	C	320	
1	D	320	
1	E	320	
1	F	320	

2 Entry composition i

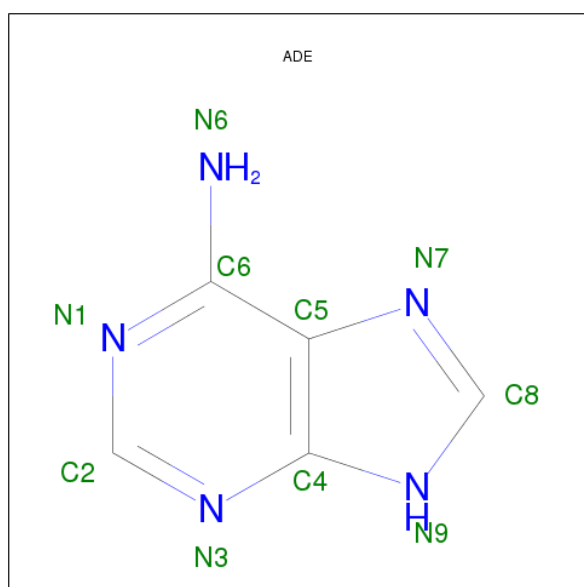
There are 4 unique types of molecules in this entry. The entry contains 13756 atoms, of which 0 are hydrogen and 0 are deuterium.

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 S-methyl-5'-thioadenosinephosphorylase.

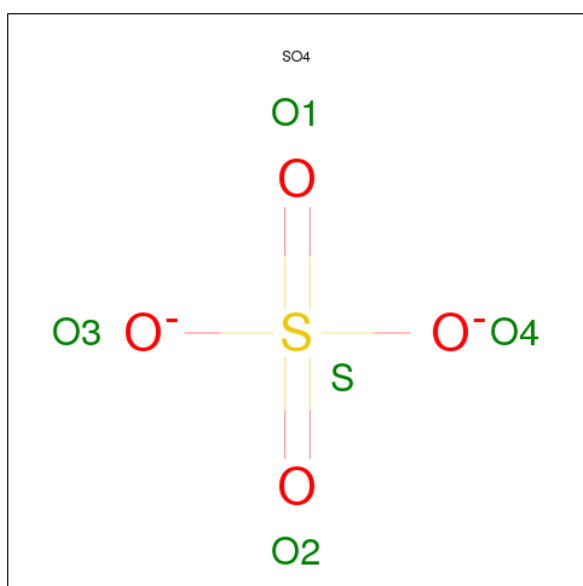
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	289	Total	C	N	O	S	0	0	0
			2197	1395	384	401	17			
1	B	290	Total	C	N	O	S	0	0	0
			2197	1390	380	409	18			
1	C	277	Total	C	N	O	S	0	0	0
			2105	1337	365	388	15			
1	D	286	Total	C	N	O	S	0	0	0
			2168	1375	379	397	17			
1	E	288	Total	C	N	O	S	0	0	0
			2202	1394	381	410	17			
1	F	288	Total	C	N	O	S	0	0	0
			2193	1388	382	408	15			

- Molecule 2 is ADENINE (three-letter code: ADE) (formula: C₅H₅N₅).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	N	0	0
			10	5	5		
2	B	1	Total	C	N	0	0
			10	5	5		
2	D	1	Total	C	N	0	0
			10	5	5		
2	E	1	Total	C	N	0	0
			10	5	5		

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		
3	E	1	Total	O	S	0	0
			5	4	1		
3	F	1	Total	O	S	0	0
			5	4	1		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	127	Total	O	0	0
			127	127		
4	B	114	Total	O	0	0
			114	114		

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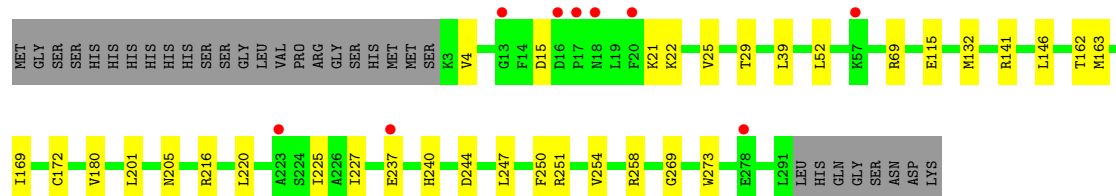
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	C	94	Total 94	O 94	0	0
4	D	80	Total 80	O 80	0	0
4	E	91	Total 91	O 91	0	0
4	F	128	Total 128	O 128	0	0

3 Residue-property plots

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

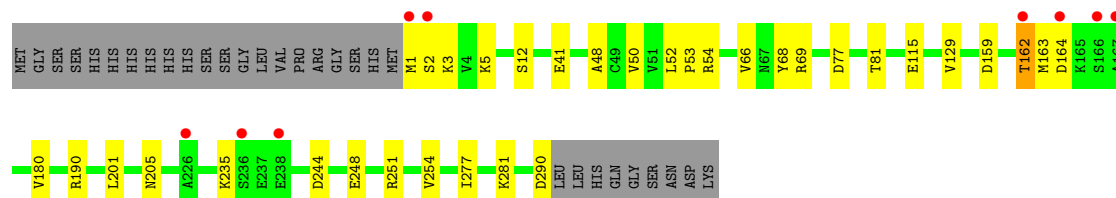
- Molecule 1: S-methyl-5'-thioadenosinephosphorylase

Chain A: 



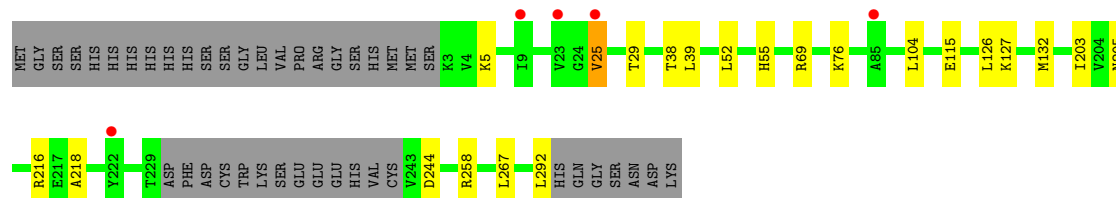
- Molecule 1: S-methyl-5'-thioadenosinephosphorylase

Chain B: 



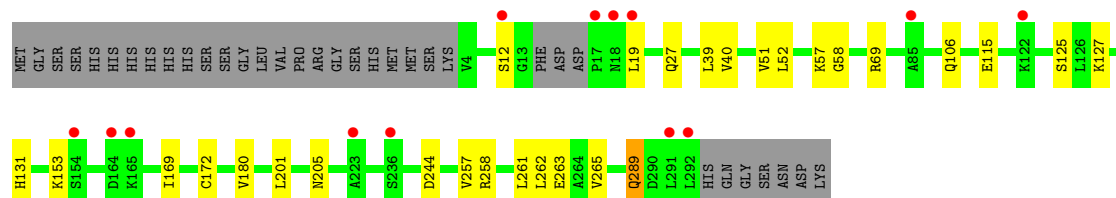
- Molecule 1: S-methyl-5'-thioadenosinephosphorylase

Chain C: 

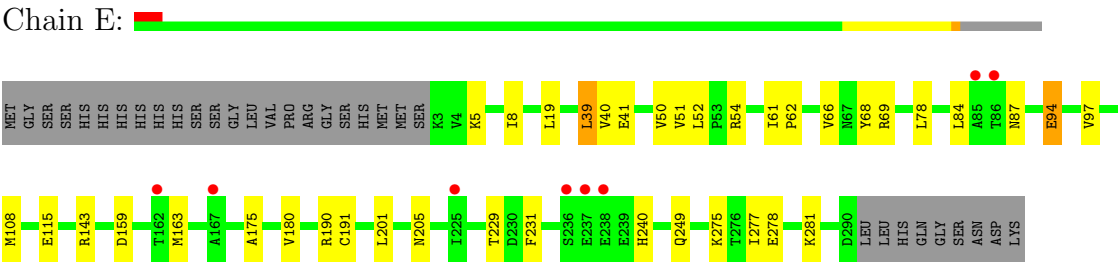


- Molecule 1: S-methyl-5'-thioadenosinephosphorylase

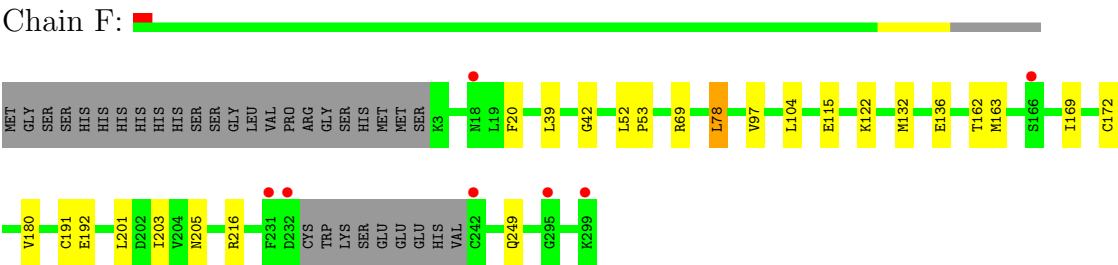
Chain D: 



• Molecule 1: S-methyl-5'-thioadenosinephosphorylase



• Molecule 1: S-methyl-5'-thioadenosinephosphorylase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	81.27Å 82.55Å 150.30Å 90.00° 100.61° 90.00°	Depositor
Resolution (Å)	30.01 – 2.10 30.01 – 2.10	Depositor EDS
% Data completeness (in resolution range)	91.0 (30.01-2.10) 91.0 (30.01-2.10)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.12	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.89 (at 2.10Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.2_1309)	Depositor
R, R_{free}	0.192 , 0.241 0.191 , 0.242	Depositor DCC
R_{free} test set	5228 reflections (5.03%)	DCC
Wilson B-factor (Å ²)	24.9	Xtriage
Anisotropy	0.234	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 34.3	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	157 of 103947 reflections (0.151%)	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	13756	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 58.56 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 2.0086e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

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

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	0.39	0/2242	0.56	0/3041
1	B	0.37	0/2242	0.53	0/3043
1	C	0.35	0/2145	0.54	0/2908
1	D	0.37	0/2211	0.54	0/2999
1	E	0.37	0/2247	0.53	0/3048
1	F	0.38	0/2235	0.53	0/3029
All	All	0.37	0/13322	0.54	0/18068

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2197	0	2195	18	0
1	B	2197	0	2169	20	0
1	C	2105	0	2121	12	0
1	D	2168	0	2159	17	0
1	E	2202	0	2185	24	0
1	F	2193	0	2187	12	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	10	0	4	0	0
2	B	10	0	4	0	0
2	D	10	0	4	0	0
2	E	10	0	4	0	0
3	B	5	0	0	2	0
3	C	5	0	0	1	0
3	E	5	0	0	2	0
3	F	5	0	0	0	0
4	A	127	0	0	0	0
4	B	114	0	0	1	0
4	C	94	0	0	1	0
4	D	80	0	0	0	0
4	E	91	0	0	1	0
4	F	128	0	0	1	0
All	All	13756	0	13032	99	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 4.

All (99) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:54:ARG:NH1	3:B:302:SO4:O4	2.07	0.88
1:E:69:ARG:HD2	1:E:115:GLU:HB3	1.65	0.78
1:B:41:GLU:HG2	1:B:50:VAL:HG22	1.66	0.76
1:E:54:ARG:NH1	3:E:302:SO4:O1	2.18	0.76
1:F:191:CYS:SG	4:F:512:HOH:O	2.52	0.67
1:E:69:ARG:HD2	1:E:115:GLU:CB	2.25	0.67
1:E:40:VAL:HG13	1:E:51:VAL:HB	1.79	0.65
1:A:254:VAL:HG12	1:A:258:ARG:HE	1.62	0.65
1:B:3:LYS:HG3	1:B:81:THR:HG21	1.78	0.64
1:A:162:THR:HG23	1:A:163:MET:HG3	1.78	0.64
1:D:69:ARG:HD2	1:D:115:GLU:HB3	1.80	0.64
1:C:25:VAL:HG13	1:C:38:THR:HG21	1.80	0.63
1:A:180:VAL:HB	1:A:201:LEU:HD13	1.82	0.62
1:E:39:LEU:HD22	1:E:52:LEU:HD13	1.82	0.62
1:F:69:ARG:HD2	1:F:115:GLU:HB3	1.82	0.62
1:F:132:MET:HE2	1:F:216:ARG:HG2	1.83	0.60
1:B:69:ARG:HD2	1:B:115:GLU:CB	2.32	0.60
1:A:250:PHE:O	1:A:254:VAL:HG23	2.02	0.60
1:D:69:ARG:HD2	1:D:115:GLU:CB	2.33	0.59
1:C:132:MET:HE2	1:C:216:ARG:HG2	1.84	0.59
1:D:180:VAL:HB	1:D:201:LEU:HD13	1.85	0.57

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:180:VAL:HB	1:B:201:LEU:HD13	1.87	0.57
1:B:69:ARG:HD2	1:B:115:GLU:HB3	1.86	0.56
1:A:132:MET:HE2	1:A:216:ARG:HG2	1.88	0.54
1:A:251:ARG:NH2	1:C:292:LEU:HD11	2.23	0.54
1:F:69:ARG:HD2	1:F:115:GLU:CB	2.37	0.54
1:A:227:ILE:HD13	1:A:250:PHE:CZ	2.44	0.53
1:D:40:VAL:HG23	1:D:51:VAL:HB	1.92	0.52
1:B:5:LYS:NZ	1:B:41:GLU:OE2	2.39	0.52
1:E:190:ARG:HH22	1:E:240:HIS:CE1	2.28	0.52
1:A:22:LYS:HD2	1:A:25:VAL:HG23	1.92	0.51
1:C:258:ARG:NH2	4:C:475:HOH:O	2.31	0.51
1:A:247:LEU:HD13	1:C:292:LEU:HD22	1.91	0.51
1:E:277:ILE:HG22	1:E:281:LYS:HE2	1.92	0.51
1:C:69:ARG:HD2	1:C:115:GLU:CB	2.41	0.51
1:A:69:ARG:HD2	1:A:115:GLU:CB	2.41	0.51
1:B:244:ASP:OD2	4:B:496:HOH:O	2.19	0.51
1:C:76:LYS:HD2	1:C:218:ALA:HB1	1.93	0.50
1:E:275:LYS:HZ3	1:E:275:LYS:HA	1.76	0.50
1:B:129:VAL:HG12	1:B:290:ASP:HB3	1.94	0.50
1:D:39:LEU:HD13	1:D:52:LEU:HB2	1.92	0.50
1:A:4:VAL:HG21	1:A:269:GLY:HA2	1.93	0.49
1:E:5:LYS:NZ	1:E:41:GLU:OE2	2.37	0.49
1:A:146:LEU:HD13	1:A:225:ILE:HG12	1.95	0.49
1:E:143:ARG:NH2	1:E:175:ALA:O	2.46	0.48
1:B:162:THR:HG23	1:B:163:MET:HG3	1.94	0.48
1:E:94:GLU:HB2	4:E:417:HOH:O	2.12	0.48
1:F:39:LEU:HD12	1:F:78:LEU:HD22	1.94	0.48
1:B:12:SER:N	3:B:302:SO4:O3	2.46	0.48
1:E:159:ASP:O	1:E:163:MET:HG2	2.14	0.47
1:A:39:LEU:HG	1:A:52:LEU:HD13	1.95	0.47
1:A:254:VAL:CG1	1:A:258:ARG:HE	2.26	0.47
1:E:97:VAL:HG22	1:E:231:PHE:HZ	1.79	0.47
1:F:104:LEU:HD21	1:F:203:ILE:HD11	1.97	0.46
1:D:169:ILE:O	1:D:172:CYS:HB3	2.15	0.46
1:A:237:GLU:HA	1:A:240:HIS:CE1	2.51	0.46
1:C:104:LEU:HD21	1:C:203:ILE:HD11	1.96	0.46
1:B:69:ARG:HD2	1:B:115:GLU:HB2	1.98	0.46
1:C:55:HIS:NE2	3:C:301:SO4:O1	2.39	0.46
1:E:180:VAL:HB	1:E:201:LEU:HD13	1.98	0.45
1:E:108:MET:HG2	1:F:192:GLU:HG2	1.98	0.45
1:B:52:LEU:HD12	1:B:53:PRO:HD2	1.99	0.45
1:E:41:GLU:HG3	1:E:50:VAL:HG22	1.98	0.45

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:E:8:ILE:HD13	1:E:84:LEU:HB2	1.99	0.45
1:F:162:THR:HG23	1:F:163:MET:HG3	1.97	0.45
1:C:69:ARG:HD2	1:C:115:GLU:HB3	1.98	0.45
1:D:12:SER:O	1:D:258:ARG:NH1	2.50	0.45
1:E:275:LYS:NZ	1:E:278:GLU:OE1	2.47	0.44
1:B:5:LYS:HE2	1:B:48:ALA:HB1	1.98	0.43
1:E:61:ILE:HA	1:E:62:PRO:HD3	1.91	0.43
1:E:66:VAL:HB	1:E:68:TYR:CE1	2.53	0.43
1:A:141:ARG:HH22	1:A:273:TRP:HA	1.82	0.43
1:C:127:LYS:HD2	1:C:127:LYS:N	2.33	0.43
1:E:229:THR:HA	1:E:249:GLN:HG2	2.00	0.43
1:B:190:ARG:NH2	1:B:235:LYS:O	2.50	0.43
1:D:57:LYS:HA	1:D:58:GLY:HA2	1.71	0.43
1:D:40:VAL:CG2	1:D:51:VAL:HB	2.49	0.43
1:D:153:LYS:HE3	1:D:263:GLU:HB2	2.00	0.42
1:B:66:VAL:HB	1:B:68:TYR:CE1	2.54	0.42
1:D:106:GLN:HB3	1:E:191:CYS:HB2	2.01	0.42
1:D:261:LEU:O	1:D:265:VAL:HG13	2.19	0.42
1:D:262:LEU:O	1:D:265:VAL:HG22	2.19	0.42
1:A:69:ARG:HD2	1:A:115:GLU:HB3	2.01	0.42
1:B:159:ASP:HB3	1:B:162:THR:HG22	2.00	0.41
1:B:3:LYS:HD2	1:B:277:ILE:HD13	2.02	0.41
1:D:19:LEU:HD23	1:D:262:LEU:HD13	2.00	0.41
1:F:180:VAL:HB	1:F:201:LEU:HD13	2.02	0.41
1:F:169:ILE:O	1:F:172:CYS:HB3	2.20	0.41
1:A:169:ILE:O	1:A:172:CYS:HB3	2.20	0.41
1:F:20:PHE:CE1	1:F:42:GLY:HA3	2.56	0.41
1:C:39:LEU:HD22	1:C:52:LEU:HB2	2.03	0.41
1:D:69:ARG:NH2	1:D:125:SER:OG	2.54	0.40
1:E:39:LEU:HD12	1:E:78:LEU:HD12	2.02	0.40
1:E:87:ASN:HA	3:E:302:SO4:O1	2.22	0.40
1:F:52:LEU:HD12	1:F:53:PRO:HD2	2.03	0.40
1:B:248:GLU:HA	1:B:251:ARG:NH1	2.36	0.40
1:B:1:MET:HA	1:B:2:SER:HA	1.61	0.40
1:D:127:LYS:HA	1:D:127:LYS:HD3	1.83	0.40
1:D:131:HIS:HD2	1:D:289:GLN:HG3	1.86	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	287/320 (90%)	283 (99%)	4 (1%)	0	100	100
1	B	288/320 (90%)	285 (99%)	3 (1%)	0	100	100
1	C	273/320 (85%)	272 (100%)	1 (0%)	0	100	100
1	D	282/320 (88%)	274 (97%)	8 (3%)	0	100	100
1	E	286/320 (89%)	281 (98%)	5 (2%)	0	100	100
1	F	284/320 (89%)	282 (99%)	2 (1%)	0	100	100
All	All	1700/1920 (88%)	1677 (99%)	23 (1%)	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	238/274 (87%)	232 (98%)	6 (2%)	60	63
1	B	238/274 (87%)	232 (98%)	6 (2%)	60	63
1	C	230/274 (84%)	223 (97%)	7 (3%)	53	55
1	D	234/274 (85%)	229 (98%)	5 (2%)	66	70
1	E	240/274 (88%)	236 (98%)	4 (2%)	73	78
1	F	239/274 (87%)	233 (98%)	6 (2%)	60	63
All	All	1419/1644 (86%)	1385 (98%)	34 (2%)	61	65

All (34) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	15	ASP
1	A	21	LYS
1	A	29	THR
1	A	205	ASN
1	A	220	LEU
1	A	244	ASP
1	B	77	ASP
1	B	162	THR
1	B	164	ASP
1	B	205	ASN
1	B	254	VAL
1	B	281	LYS
1	C	5	LYS
1	C	25	VAL
1	C	29	THR
1	C	126	LEU
1	C	205	ASN
1	C	244	ASP
1	C	267	LEU
1	D	27	GLN
1	D	205	ASN
1	D	244	ASP
1	D	257	VAL
1	D	289	GLN
1	E	19	LEU
1	E	39	LEU
1	E	94	GLU
1	E	205	ASN
1	F	78	LEU
1	F	97	VAL
1	F	122	LYS
1	F	136	GLU
1	F	205	ASN
1	F	249	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA ⓘ

There are no RNA chains in this entry.

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

Mogul failed to run properly - this section will therefore be empty.

5.5 Carbohydrates ⓘ

Mogul failed to run properly - this section will therefore be empty.

5.6 Ligand geometry ⓘ

Mogul failed to run properly - this section will therefore be empty.

5.7 Other polymers ⓘ

Mogul failed to run properly - this section will therefore be empty.

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	289/320 (90%)	-0.06	9 (3%)	47	52	11, 24, 49, 68	0
1	B	290/320 (90%)	0.13	9 (3%)	47	52	15, 27, 54, 88	0
1	C	277/320 (86%)	0.01	5 (1%)	65	69	16, 31, 56, 70	0
1	D	286/320 (89%)	0.11	13 (4%)	32	35	16, 30, 50, 75	0
1	E	288/320 (90%)	0.00	8 (2%)	50	55	14, 28, 55, 83	0
1	F	288/320 (90%)	-0.09	7 (2%)	56	61	14, 25, 48, 70	0
All	All	1718/1920 (89%)	0.02	51 (2%)	48	53	11, 28, 53, 88	0

All (51) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	2	SER	6.2
1	B	1	MET	4.9
1	D	292	LEU	4.3
1	C	25	VAL	4.3
1	D	19	LEU	3.7
1	B	238	GLU	3.6
1	E	237	GLU	3.4
1	E	238	GLU	3.4
1	D	18	ASN	3.3
1	B	166	SER	3.2
1	E	162	THR	3.2
1	B	162	THR	3.1
1	C	23	VAL	3.1
1	F	231	PHE	3.1
1	A	57	LYS	3.1
1	E	225	ILE	3.0
1	B	226	ALA	3.0
1	B	236	SER	3.0
1	A	237	GLU	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	16	ASP	2.9
1	F	299	LYS	2.9
1	B	164	ASP	2.8
1	C	9	ILE	2.8
1	A	17	PRO	2.7
1	F	242	CYS	2.7
1	F	232	ASP	2.5
1	E	167	ALA	2.5
1	E	85	ALA	2.4
1	D	17	PRO	2.4
1	A	18	ASN	2.4
1	D	236	SER	2.4
1	E	86	THR	2.4
1	A	223	ALA	2.4
1	B	167	ALA	2.3
1	F	18	ASN	2.3
1	D	85	ALA	2.3
1	C	222	TYR	2.3
1	F	295	GLY	2.3
1	D	154	SER	2.2
1	D	165	LYS	2.2
1	D	223	ALA	2.2
1	D	12	SER	2.2
1	C	85	ALA	2.2
1	A	20	PHE	2.1
1	A	13	GLY	2.1
1	D	164	ASP	2.1
1	D	122	LYS	2.1
1	D	291	LEU	2.1
1	A	278	GLU	2.0
1	E	236	SER	2.0
1	F	166	SER	2.0

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

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(Å ²)	Q<0.9
3	SO4	E	302	5/5	0.18	0.36	40,46,54,55	0
3	SO4	B	302	5/5	0.18	0.22	40,50,54,56	0
2	ADE	B	301	10/10	0.13	0.20	27,28,29,30	0
2	ADE	A	301	10/10	0.10	-0.46	14,16,18,18	0
2	ADE	E	301	10/10	0.11	-0.66	22,24,27,28	0
2	ADE	D	301	10/10	0.08	-1.61	19,25,28,30	0
3	SO4	C	301	5/5	0.09	-1.65	27,34,38,38	0
3	SO4	F	301	5/5	0.07	-1.65	21,24,27,27	0

6.5 Other polymers

There are no such residues in this entry.