Full wwPDB X-ray Structure Validation Report (i)

Feb 1, 2016 – 07:05 AM GMT

PDB ID	:	2ZIF
Title	:	Crystal Structure of TTHA0409, Putative DNA Modification Methylase from Thermus thermophilus HB8- Complexed with S-Adenosyl-L-Methionine
Authors	:	Morita, R.; Ishikawa, H.; Nakagawa, N.; Masui, R.; Yokoyama, S.; Kuramitsu, S.; RIKEN Structural Genomics/Proteomics Initiative (RSGI)
Deposited on	:	2008-02-15
Resolution	:	2.40 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
http://wwpdb.org/validation/2016/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

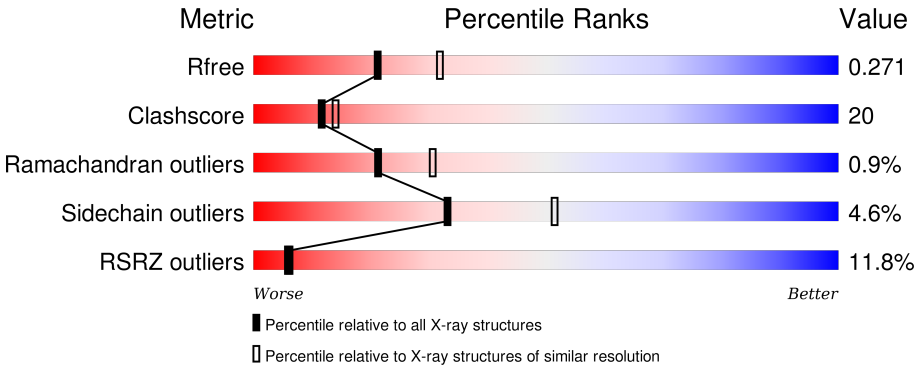
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk26865

1 Overall quality at a glance (i)

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

The reported resolution of this entry is 2.40 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}	91344	2919 (2.40-2.40)
Clashscore	102246	3407 (2.40-2.40)
Ramachandran outliers	100387	3351 (2.40-2.40)
Sidechain outliers	100360	3352 (2.40-2.40)
RSRZ outliers	91569	2928 (2.40-2.40)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=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	297	11% 51% 28% • 20%
1	B	297	8% 52% 26% • 18%

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
2	SAM	B	298	-	-	X	-

2 Entry composition

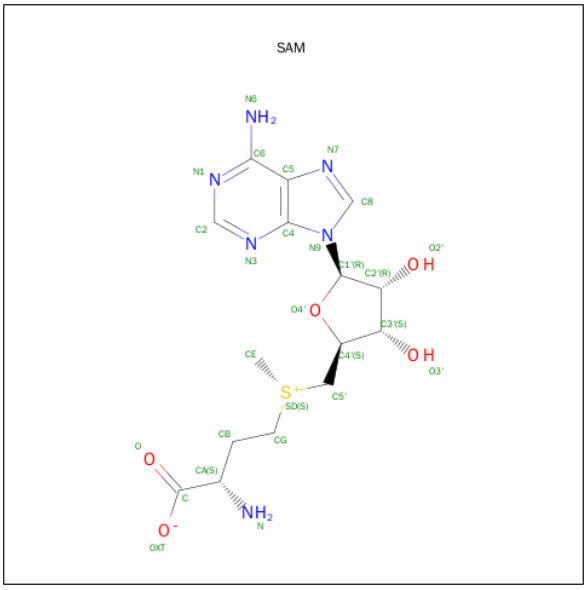
There are 3 unique types of molecules in this entry. The entry contains 4047 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Putative modification methylase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	239	Total	C	N	O	S	0	0	0
			1951	1265	352	331	3			
1	B	243	Total	C	N	O	S	0	0	0
			1981	1280	362	336	3			

- Molecule 2 is S-ADENOSYLMETHIONINE (three-letter code: SAM) (formula: $C_{15}H_{22}N_6O_5S$).



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

- Molecule 3 is water.

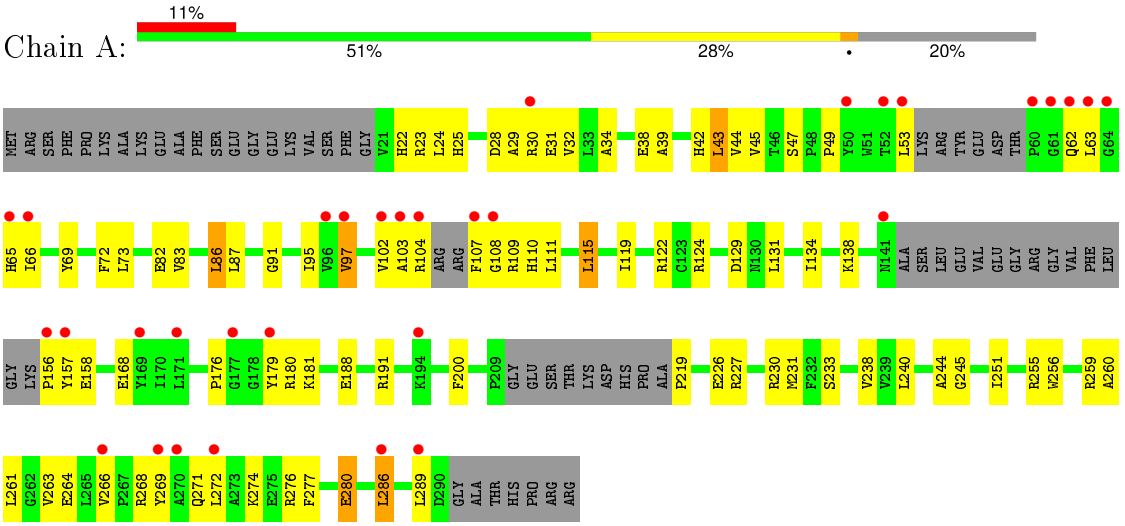
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	35	Total O 35 35	0	0
3	B	26	Total O 26 26	0	0

3 Residue-property plots (i)

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

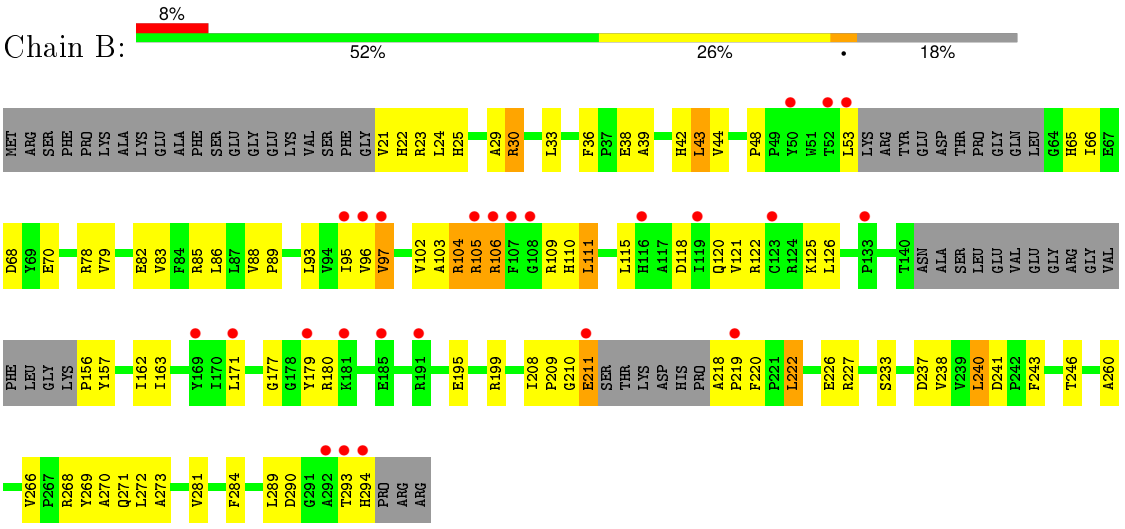
- Molecule 1: Putative modification methylase

Chain A:



- Molecule 1: Putative modification methylase

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	63.90Å 58.64Å 81.22Å 90.00° 106.06° 90.00°	Depositor
Resolution (Å)	40.67 – 2.40 40.66 – 2.40	Depositor EDS
% Data completeness (in resolution range)	93.5 (40.67-2.40) 93.3 (40.66-2.40)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	0.03	Depositor
$< I/\sigma(I) >$ [1]	4.10 (at 2.39Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.217 , 0.273 0.216 , 0.271	Depositor DCC
R_{free} test set	2105 reflections (9.86%)	DCC
Wilson B-factor (Å^2)	42.0	Xtriage
Anisotropy	0.341	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.35 , 49.9	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Outliers	2 of 22664 reflections (0.009%)	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	4047	wwPDB-VP
Average B, all atoms (Å^2)	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.51% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

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

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	0.38	0/2004	0.63	0/2714
1	B	0.36	0/2035	0.63	0/2757
All	All	0.37	0/4039	0.63	0/5471

There are no bond length outliers.

There are no bond angle outliers.

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	1951	0	1955	78	0
1	B	1981	0	1981	84	0
2	A	27	0	22	3	0
2	B	27	0	22	11	0
3	A	35	0	0	2	0
3	B	26	0	0	2	0
All	All	4047	0	3980	162	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 20.

All (162) 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:298:SAM:CB	2:B:298:SAM:CG	1.78	1.56
1:B:219:PRO:HD2	2:B:298:SAM:HB2	1.49	0.91
1:B:42:HIS:HD2	1:B:238:VAL:H	1.08	0.89
1:A:42:HIS:HD2	1:A:238:VAL:H	1.18	0.86
1:B:22:HIS:HD2	1:B:260:ALA:H	1.24	0.86
1:B:44:VAL:HG22	1:B:240:LEU:HB3	1.59	0.83
1:B:103:ALA:HB2	1:B:110:HIS:ND1	1.95	0.81
1:B:96:VAL:HG21	1:B:220:PHE:HD2	1.49	0.75
1:B:243:PHE:O	2:B:298:SAM:HG2	1.86	0.74
1:A:62:GLN:HG2	1:A:63:LEU:HD12	1.68	0.74
1:A:22:HIS:HD2	1:A:260:ALA:H	1.35	0.74
1:B:293:THR:HG22	1:B:294:HIS:H	1.53	0.73
1:A:95:ILE:HG22	1:A:97:VAL:HG22	1.69	0.73
1:B:96:VAL:HG21	1:B:220:PHE:CD2	2.24	0.71
1:A:34:ALA:HA	1:A:82:GLU:OE2	1.90	0.71
1:B:219:PRO:CD	2:B:298:SAM:HB2	2.19	0.71
1:B:42:HIS:CD2	1:B:238:VAL:H	2.00	0.70
1:B:30:ARG:NH1	1:B:78:ARG:HH21	1.91	0.69
1:A:83:VAL:HA	1:A:86:LEU:HD11	1.74	0.68
1:A:43:LEU:HD23	1:A:44:VAL:N	2.08	0.68
1:A:271:GLN:HE22	1:A:274:LYS:HE3	1.60	0.67
1:A:30:ARG:HH11	1:A:30:ARG:HG3	1.60	0.67
1:B:21:VAL:HG11	1:B:23:ARG:HH21	1.59	0.67
1:B:43:LEU:HD12	1:B:233:SER:HB3	1.77	0.66
1:A:42:HIS:CD2	1:A:238:VAL:H	2.08	0.65
1:A:179:TYR:HE1	1:A:181:LYS:HZ2	1.45	0.65
1:A:49:PRO:HA	2:A:298:SAM:H5'1	1.79	0.64
1:B:180:ARG:HH22	1:B:237:ASP:CG	2.00	0.64
1:B:42:HIS:HB3	1:B:180:ARG:HH21	1.63	0.64
1:A:268:ARG:O	1:A:272:LEU:HG	1.98	0.63
1:B:96:VAL:HG23	1:B:96:VAL:O	1.99	0.62
1:B:66:ILE:O	1:B:70:GLU:HG2	1.99	0.62
1:A:188:GLU:HA	1:A:191:ARG:HH11	1.64	0.62
1:A:111:LEU:HD11	3:A:324:HOH:O	1.99	0.62
1:A:219:PRO:HA	3:A:323:HOH:O	2.00	0.61
1:A:43:LEU:HD12	1:A:233:SER:HB3	1.83	0.60
1:A:240:LEU:HA	1:A:261:LEU:O	2.01	0.60
1:A:62:GLN:HG2	1:A:63:LEU:CD1	2.32	0.60
1:B:42:HIS:HB3	1:B:180:ARG:NH2	2.16	0.60
2:B:298:SAM:CA	2:B:298:SAM:CG	2.77	0.59
1:B:243:PHE:HB3	2:B:298:SAM:H5'2	1.84	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:179:TYR:HE1	1:A:181:LYS:NZ	2.00	0.59
1:A:103:ALA:HA	1:A:110:HIS:HA	1.85	0.59
2:B:298:SAM:CB	2:B:298:SAM:SD	2.91	0.58
1:B:268:ARG:O	1:B:272:LEU:HG	2.03	0.58
1:B:240:LEU:HD23	1:B:241:ASP:N	2.18	0.58
1:B:30:ARG:HG3	1:B:30:ARG:HH11	1.69	0.58
1:B:22:HIS:CD2	1:B:260:ALA:H	2.15	0.58
1:B:38:GLU:HG2	1:B:39:ALA:N	2.19	0.58
1:A:23:ARG:HH11	1:A:25:HIS:CE1	2.21	0.58
1:B:25:HIS:CE1	1:B:289:LEU:HD12	2.40	0.57
1:A:226:GLU:OE1	1:A:255:ARG:HD2	2.04	0.57
1:B:195:GLU:HG2	1:B:199:ARG:HH22	1.69	0.57
1:A:43:LEU:HD23	1:A:44:VAL:H	1.70	0.57
1:B:156:PRO:HG2	1:B:157:TYR:H	1.70	0.56
1:B:103:ALA:HB1	1:B:110:HIS:HB2	1.88	0.56
1:B:82:GLU:O	1:B:86:LEU:HD13	2.05	0.56
1:A:227:ARG:O	1:A:231:MET:HG3	2.06	0.56
1:B:79:VAL:O	1:B:83:VAL:HG23	2.05	0.56
1:A:83:VAL:HA	1:A:86:LEU:CD1	2.36	0.56
1:B:122:ARG:O	1:B:126:LEU:HG	2.07	0.55
1:B:53:LEU:HD21	1:B:102:VAL:HG11	1.88	0.54
1:A:276:ARG:NH2	1:A:280:GLU:OE2	2.40	0.54
1:A:124:ARG:NH2	1:B:118:ASP:OD2	2.41	0.54
1:B:30:ARG:HH12	1:B:78:ARG:HH21	1.56	0.54
1:A:43:LEU:CD2	1:A:44:VAL:N	2.69	0.54
1:A:29:ALA:HA	1:A:263:VAL:HG11	1.89	0.54
1:B:156:PRO:O	1:B:157:TYR:HB2	2.08	0.54
1:A:30:ARG:HD2	1:A:30:ARG:O	2.07	0.54
1:A:134:ILE:HD11	1:B:163:ILE:CD1	2.38	0.54
1:A:109:ARG:HG3	1:A:109:ARG:O	2.08	0.54
1:A:25:HIS:CE1	1:A:289:LEU:HD11	2.43	0.54
1:A:73:LEU:HB3	1:A:122:ARG:NE	2.23	0.54
1:A:82:GLU:O	1:A:86:LEU:HG	2.07	0.53
1:A:188:GLU:HA	1:A:191:ARG:NH1	2.23	0.53
1:A:24:LEU:HB2	1:A:286:LEU:HD21	1.89	0.53
1:A:104:ARG:HA	1:A:108:GLY:O	2.08	0.53
1:B:65:HIS:NE2	1:B:106:ARG:HG2	2.25	0.52
1:A:43:LEU:HD22	1:A:45:VAL:HG23	1.91	0.52
1:B:102:VAL:HG13	1:B:102:VAL:O	2.09	0.52
1:A:102:VAL:O	1:A:102:VAL:HG23	2.09	0.52
1:A:138:LYS:HD2	1:A:168:GLU:HG3	1.92	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:42:HIS:HD2	1:B:238:VAL:N	1.92	0.51
1:B:222:LEU:O	1:B:226:GLU:HG3	2.09	0.51
1:B:290:ASP:HB2	3:B:324:HOH:O	2.09	0.51
1:A:97:VAL:HG11	1:A:115:LEU:HD13	1.93	0.51
1:A:251:ILE:HD11	1:A:276:ARG:HG2	1.93	0.51
1:A:28:ASP:OD2	1:A:30:ARG:HB3	2.09	0.51
1:A:191:ARG:HG3	1:A:191:ARG:HH11	1.77	0.50
1:A:240:LEU:HD12	1:A:261:LEU:O	2.12	0.50
1:B:29:ALA:O	1:B:33:LEU:HG	2.11	0.49
1:A:180:ARG:HH21	1:A:180:ARG:HG2	1.76	0.49
1:B:89:PRO:HB3	1:B:177:GLY:HA3	1.94	0.49
1:A:277:PHE:CG	1:A:286:LEU:HB2	2.46	0.49
1:B:48:PRO:CG	1:B:97:VAL:HG13	2.41	0.49
1:A:97:VAL:CG1	1:A:115:LEU:HD13	2.42	0.49
1:A:53:LEU:HD12	1:A:53:LEU:N	2.28	0.49
1:B:83:VAL:HG11	1:B:93:LEU:HD13	1.96	0.48
2:B:298:SAM:HG1	3:B:311:HOH:O	2.12	0.48
1:B:39:ALA:O	1:B:88:VAL:HG12	2.13	0.48
1:B:22:HIS:HB3	1:B:260:ALA:HB3	1.96	0.47
1:B:210:GLY:O	1:B:211:GLU:HB2	2.15	0.47
1:B:95:ILE:HG22	1:B:97:VAL:HG22	1.97	0.47
1:A:115:LEU:HD22	1:A:119:ILE:HD11	1.96	0.47
1:B:44:VAL:HG21	1:B:83:VAL:HG13	1.97	0.47
1:A:129:ASP:OD2	1:A:176:PRO:HB3	2.15	0.47
1:B:240:LEU:HD23	1:B:241:ASP:H	1.77	0.47
1:A:30:ARG:NH1	1:A:30:ARG:HG3	2.24	0.47
1:B:102:VAL:HG22	1:B:105:ARG:HB2	1.97	0.46
1:B:120:GLN:HG3	1:B:171:LEU:HD22	1.98	0.46
1:A:65:HIS:ND1	1:A:66:ILE:N	2.64	0.46
1:B:53:LEU:HD12	1:B:53:LEU:O	2.16	0.45
1:B:104:ARG:HH11	1:B:104:ARG:HG3	1.81	0.45
1:B:281:VAL:HG12	1:B:284:PHE:HB2	1.99	0.45
1:B:23:ARG:NH1	1:B:289:LEU:HD11	2.31	0.45
1:A:277:PHE:CD2	1:A:286:LEU:HB2	2.52	0.45
1:A:53:LEU:HD23	1:A:107:PHE:HZ	1.82	0.45
1:A:28:ASP:O	1:A:32:VAL:HG23	2.16	0.44
1:B:111:LEU:O	1:B:111:LEU:HD23	2.18	0.44
1:A:87:LEU:HG	1:A:91:GLY:HA3	1.98	0.44
1:A:65:HIS:ND1	1:A:66:ILE:HG23	2.32	0.44
1:A:271:GLN:NE2	1:A:274:LYS:HE3	2.29	0.44
1:A:230:ARG:HD2	1:A:256:TRP:CE2	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:30:ARG:HG3	1:B:30:ARG:NH1	2.32	0.44
1:B:36:PHE:O	1:B:85:ARG:NH1	2.47	0.43
1:A:69:TYR:O	1:A:72:PHE:HB3	2.17	0.43
1:A:138:LYS:HD2	1:A:168:GLU:CG	2.48	0.43
1:B:110:HIS:CD2	1:B:162:ILE:HD12	2.53	0.43
1:A:156:PRO:O	1:A:158:GLU:N	2.49	0.43
1:B:208:ILE:HD11	1:B:227:ARG:CZ	2.49	0.43
1:A:245:GLY:HA3	2:A:298:SAM:O	2.19	0.43
1:B:246:THR:N	2:B:298:SAM:O	2.44	0.42
1:B:38:GLU:HG3	1:B:85:ARG:O	2.19	0.42
1:A:129:ASP:OD2	1:B:109:ARG:HD2	2.19	0.42
1:A:31:GLU:O	1:A:34:ALA:HB3	2.20	0.42
1:B:121:VAL:HG12	1:B:125:LYS:HZ3	1.84	0.42
1:B:121:VAL:HG12	1:B:125:LYS:NZ	2.35	0.42
1:A:180:ARG:NH2	1:A:180:ARG:HG2	2.33	0.42
1:B:65:HIS:HA	1:B:68:ASP:OD2	2.19	0.42
1:B:109:ARG:NH1	1:B:111:LEU:HD13	2.35	0.42
1:B:30:ARG:HD2	1:B:30:ARG:O	2.19	0.42
1:A:28:ASP:OD1	2:A:298:SAM:N6	2.48	0.42
1:B:23:ARG:HH11	1:B:289:LEU:HD11	1.84	0.41
1:B:96:VAL:O	1:B:96:VAL:CG2	2.67	0.41
1:B:270:ALA:O	1:B:273:ALA:HB3	2.19	0.41
1:B:48:PRO:HD3	1:B:95:ILE:HG23	2.02	0.41
1:A:38:GLU:HG2	1:A:39:ALA:N	2.36	0.41
1:B:82:GLU:OE1	1:B:82:GLU:HA	2.20	0.41
1:B:243:PHE:HB3	2:B:298:SAM:C4'	2.50	0.41
1:A:131:LEU:HD22	1:B:110:HIS:CD2	2.56	0.41
1:A:244:ALA:HB3	1:A:264:GLU:HB2	2.03	0.41
1:A:24:LEU:CB	1:A:286:LEU:HD21	2.50	0.41
1:B:65:HIS:CD2	1:B:106:ARG:HG2	2.56	0.41
1:B:218:ALA:HA	1:B:219:PRO:HD2	1.83	0.40
1:B:243:PHE:HB3	2:B:298:SAM:C5'	2.52	0.40
1:B:103:ALA:C	1:B:105:ARG:H	2.23	0.40
1:B:208:ILE:HA	1:B:209:PRO:HD2	1.94	0.40
1:A:102:VAL:HG22	1:A:111:LEU:O	2.20	0.40
1:A:266:VAL:HB	1:A:269:TYR:CD2	2.56	0.40
1:B:266:VAL:HB	1:B:269:TYR:CD1	2.56	0.40
1:A:200:PHE:N	1:A:200:PHE:CD1	2.89	0.40
1:A:63:LEU:HG	1:A:66:ILE:HD11	2.03	0.40

There are no symmetry-related clashes.

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	229/297 (77%)	217 (95%)	11 (5%)	1 (0%)	39 56
1	B	235/297 (79%)	220 (94%)	12 (5%)	3 (1%)	15 21
All	All	464/594 (78%)	437 (94%)	23 (5%)	4 (1%)	21 30

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	105	ARG
1	A	157	TYR
1	B	106	ARG
1	B	179	TYR

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	205/251 (82%)	197 (96%)	8 (4%)	39 59
1	B	206/251 (82%)	195 (95%)	11 (5%)	28 44
All	All	411/502 (82%)	392 (95%)	19 (5%)	33 51

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	43	LEU
1	A	47	SER

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Mol	Chain	Res	Type
1	A	86	LEU
1	A	97	VAL
1	A	115	LEU
1	A	259	ARG
1	A	280	GLU
1	A	286	LEU
1	B	24	LEU
1	B	30	ARG
1	B	43	LEU
1	B	97	VAL
1	B	104	ARG
1	B	111	LEU
1	B	115	LEU
1	B	211	GLU
1	B	222	LEU
1	B	240	LEU
1	B	271	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	22	HIS
1	A	42	HIS
1	A	271	GLN
1	B	22	HIS
1	B	25	HIS
1	B	42	HIS
1	B	184	GLN
1	B	271	GLN

5.3.3 RNA

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates (i)

There are no carbohydrates in this entry.

5.6 Ligand geometry (i)

2 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	$\#\|Z\| > 2$
2	SAM	A	298	-	21,29,29	2.63	1 (4%)	17,42,42	2.53	3 (17%)
2	SAM	B	298	-	21,29,29	4.55	1 (4%)	17,42,42	2.28	3 (17%)

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	SAM	A	298	-	-	0/8/33/33	0/3/3/3
2	SAM	B	298	-	-	0/8/33/33	0/3/3/3

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	298	SAM	CG-CB	11.60	1.67	1.52
2	B	298	SAM	CG-CB	20.62	1.78	1.52

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	298	SAM	N3-C2-N1	-7.60	123.08	128.89
2	A	298	SAM	N3-C2-N1	-7.56	123.10	128.89
2	A	298	SAM	C1'-N9-C4	-5.68	118.37	126.94
2	B	298	SAM	C1'-N9-C4	-3.33	121.91	126.94
2	B	298	SAM	C4-C5-N7	-2.76	106.94	109.48

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	298	SAM	C4-C5-N7	-2.75	106.95	109.48

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 14 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	298	SAM	3	0
2	B	298	SAM	11	0

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, 95^{th} 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	239/297 (80%)	0.72	32 (13%)	4	4	19, 40, 74, 104	0
1	B	243/297 (81%)	0.55	25 (10%)	9	8	22, 43, 75, 96	0
All	All	482/594 (81%)	0.64	57 (11%)	6	6	19, 42, 75, 104	0

All (57) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	61	GLY	7.7
1	A	60	PRO	6.8
1	B	107	PHE	5.7
1	A	63	LEU	5.3
1	A	62	GLN	5.1
1	A	107	PHE	5.0
1	B	294	HIS	5.0
1	A	53	LEU	4.7
1	B	106	ARG	4.5
1	B	179	TYR	4.2
1	A	179	TYR	4.0
1	A	65	HIS	3.7
1	B	50	TYR	3.7
1	A	102	VAL	3.5
1	A	66	ILE	3.4
1	B	96	VAL	3.4
1	A	50	TYR	3.4
1	A	270	ALA	3.3
1	A	269	TYR	3.2
1	A	103	ALA	3.1
1	A	64	GLY	3.0
1	B	169	TYR	3.0
1	B	171	LEU	2.9
1	B	95	ILE	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	104	ARG	2.9
1	B	108	GLY	2.9
1	A	97	VAL	2.7
1	A	156	PRO	2.6
1	B	292	ALA	2.6
1	B	53	LEU	2.6
1	A	141	ASN	2.6
1	B	293	THR	2.6
1	A	52	THR	2.5
1	B	119	ILE	2.5
1	B	219	PRO	2.5
1	A	194	LYS	2.5
1	A	169	TYR	2.4
1	B	123	CYS	2.4
1	A	286	LEU	2.4
1	A	157	TYR	2.4
1	B	191	ARG	2.4
1	B	181	LYS	2.4
1	A	30	ARG	2.3
1	B	211	GLU	2.3
1	B	97	VAL	2.3
1	B	105	ARG	2.3
1	A	266	VAL	2.3
1	A	108	GLY	2.2
1	A	171	LEU	2.2
1	A	96	VAL	2.2
1	A	272	LEU	2.1
1	B	52	THR	2.1
1	B	185	GLU	2.1
1	A	289	LEU	2.1
1	B	116	HIS	2.1
1	B	133	PRO	2.1
1	A	177	GLY	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 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, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å^2)	Q<0.9
2	SAM	B	298	27/27	0.85	0.24	0.97	62,66,69,70	0
2	SAM	A	298	27/27	0.87	0.24	0.70	64,65,76,77	0

6.5 Other polymers

There are no such residues in this entry.