



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 09:25 AM GMT

PDB ID : 3ID5  
Title : Crystal structure of Sulfolobus solfataricus C/D RNP assembled with Nop5, fibrillarin, L7Ae and a split half C/D RNA  
Authors : Ye, K.  
Deposited on : 2009-07-20  
Resolution : 4.01 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](mailto:validation@mail.wwpdb.org)  
A user guide is available at  
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>  
with specific help available everywhere you see the ⓘ symbol.

---

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

MolProbity : 4.02b-467  
Mogul : 1.7 (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

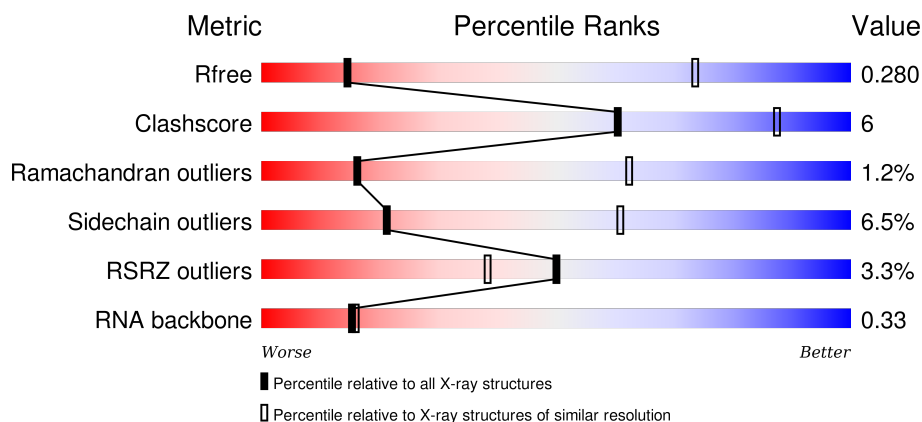
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 4.01 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	1015 (4.48-3.56)
Clashscore	102246	1054 (4.44-3.60)
Ramachandran outliers	100387	1006 (4.42-3.60)
Sidechain outliers	100360	1016 (4.46-3.58)
RSRZ outliers	91569	1018 (4.48-3.56)
RNA backbone	2183	1079 (5.04-2.80)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	388	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 13%, green 81%, grey 5%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> <span>%</span> <span>81%</span> <span>13%</span> <span>• •</span> </div> </div>
1	E	388	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 14%, green 80%, grey 4%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> <span>%</span> <span>80%</span> <span>14%</span> <span>• •</span> </div> </div>
2	B	232	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 22%, green 71%, grey 4%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> <span>%</span> <span>71%</span> <span>22%</span> <span>• •</span> </div> </div>
2	F	232	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, yellow 24%, green 69%, grey 7%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> <span>%</span> <span>69%</span> <span>24%</span> <span>• •</span> </div> </div>

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Length	Quality of chain
3	C	130	
3	G	130	
4	D	35	
4	H	35	

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	B	301	-	-	-	X
5	SAM	F	301	-	-	-	X

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 12730 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 Pre mRNA splicing protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	371	Total	C	N	O	S	0	0	0
			2931	1864	511	551	5			
1	E	371	Total	C	N	O	S	0	0	0
			2931	1864	511	551	5			

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	2	VAL	MET	ENGINEERED	UNP Q97ZH3
A	381	HIS	-	EXPRESSION TAG	UNP Q97ZH3
A	382	HIS	-	EXPRESSION TAG	UNP Q97ZH3
A	383	HIS	-	EXPRESSION TAG	UNP Q97ZH3
A	384	HIS	-	EXPRESSION TAG	UNP Q97ZH3
A	385	HIS	-	EXPRESSION TAG	UNP Q97ZH3
A	386	HIS	-	EXPRESSION TAG	UNP Q97ZH3
A	387	HIS	-	EXPRESSION TAG	UNP Q97ZH3
A	388	HIS	-	EXPRESSION TAG	UNP Q97ZH3
E	2	VAL	MET	ENGINEERED	UNP Q97ZH3
E	381	HIS	-	EXPRESSION TAG	UNP Q97ZH3
E	382	HIS	-	EXPRESSION TAG	UNP Q97ZH3
E	383	HIS	-	EXPRESSION TAG	UNP Q97ZH3
E	384	HIS	-	EXPRESSION TAG	UNP Q97ZH3
E	385	HIS	-	EXPRESSION TAG	UNP Q97ZH3
E	386	HIS	-	EXPRESSION TAG	UNP Q97ZH3
E	387	HIS	-	EXPRESSION TAG	UNP Q97ZH3
E	388	HIS	-	EXPRESSION TAG	UNP Q97ZH3

- Molecule 2 is a protein called Fibrillarin-like rRNA/tRNA 2'-O-methyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	223	Total	C	N	O	S	0	0	0
			1796	1156	302	334	4			

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	F	223	Total	C	N	O	S	0	0	0
			1796	1156	302	334	4			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	2	ALA	SER	ENGINEERED	UNP P58032
F	2	ALA	SER	ENGINEERED	UNP P58032

- Molecule 3 is a protein called 50S ribosomal protein L7Ae.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	120	Total	C	N	O	S	0	0	0
			911	578	153	178	2			
3	G	120	Total	C	N	O	S	0	0	0
			911	578	153	178	2			

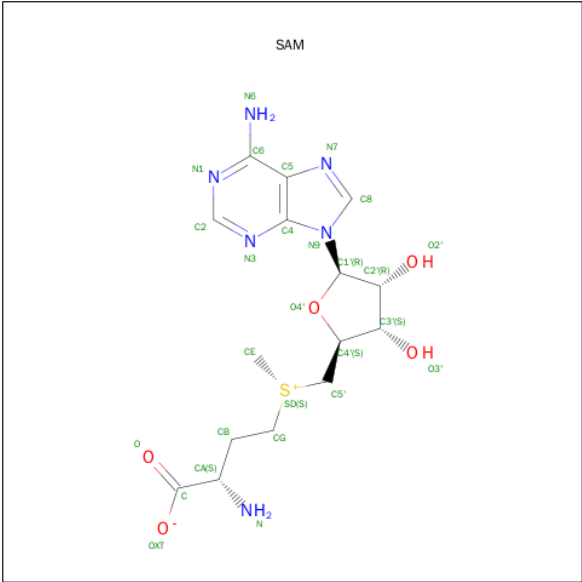
There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	1	MET	-	EXPRESSION TAG	UNP P55858
C	2	ASP	-	EXPRESSION TAG	UNP P55858
C	3	ALA	-	EXPRESSION TAG	UNP P55858
G	1	MET	-	EXPRESSION TAG	UNP P55858
G	2	ASP	-	EXPRESSION TAG	UNP P55858
G	3	ALA	-	EXPRESSION TAG	UNP P55858

- Molecule 4 is a RNA chain called half C/D RNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	34	Total	C	N	O	P	0	0	0
			700	309	121	236	34			
4	H	34	Total	C	N	O	P	0	0	0
			700	309	121	236	34			

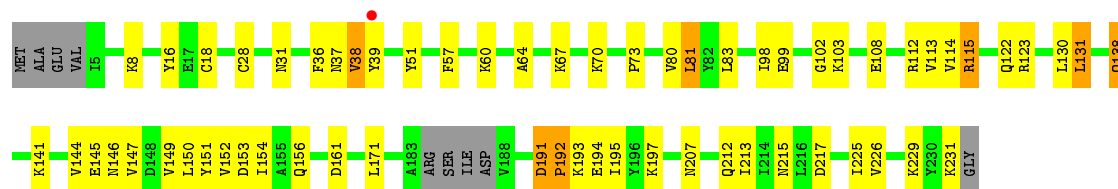
- Molecule 5 is S-ADENOSYLMETHIONINE (three-letter code: SAM) (formula: C<sub>15</sub>H<sub>22</sub>N<sub>6</sub>O<sub>5</sub>S).




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



Chain F: 




- Molecule 3: 50S ribosomal protein L7Ae

Chain C: 



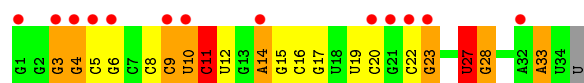
- Molecule 3: 50S ribosomal protein L7Ae

Chain G: 



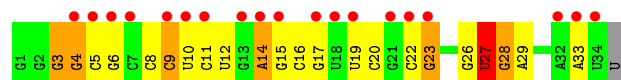
- Molecule 4: half C/D RNA

Chain D: 



- Molecule 4: half C/D RNA

Chain H: 





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	265.49 Å   265.49 Å   129.43 Å 90.00°   90.00°   90.00°	Depositor
Resolution (Å)	20.00 – 4.01 19.99 – 4.01	Depositor EDS
% Data completeness (in resolution range)	96.5 (20.00-4.01) 96.5 (19.99-4.01)	Depositor EDS
$R_{merge}$	0.14	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.22 (at 4.07 Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, $R_{free}$	0.285   ,   0.304 0.267   ,   0.280	Depositor DCC
$R_{free}$ test set	1888 reflections (5.28%)	DCC
Wilson B-factor (Å <sup>2</sup> )	135.9	Xtriage
Anisotropy	0.008	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.25 , 181.0	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	0 of 37665 reflections	Xtriage
$F_o, F_c$ correlation	0.90	EDS
Total number of atoms	12730	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	166.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.18% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  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	0.31	0/2977	0.47	0/4021
1	E	0.31	0/2977	0.47	0/4021
2	B	0.34	0/1829	0.50	0/2474
2	F	0.34	0/1829	0.50	0/2474
3	C	0.30	0/920	0.44	0/1239
3	G	0.30	0/920	0.44	0/1239
4	D	0.73	0/780	1.57	16/1214 (1.3%)
4	H	0.71	0/780	1.52	9/1214 (0.7%)
All	All	0.39	0/13012	0.72	25/17896 (0.1%)

There are no bond length outliers.

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	H	9	C	O5'-P-OP2	-21.97	84.33	110.70
4	D	9	C	O5'-P-OP1	-20.57	86.02	110.70
4	H	12	U	OP1-P-OP2	-17.74	93.00	119.60
4	D	12	U	OP1-P-OP2	-17.38	93.52	119.60
4	H	22	C	O3'-P-O5'	-10.76	83.56	104.00
4	D	22	C	O3'-P-O5'	-10.58	83.89	104.00
4	D	10	U	O3'-P-O5'	-10.27	84.48	104.00
4	D	22	C	OP2-P-O3'	-8.11	87.35	105.20
4	H	22	C	OP2-P-O3'	-7.80	88.03	105.20
4	D	14	A	P-O3'-C3'	7.72	128.96	119.70
4	H	14	A	P-O3'-C3'	7.65	128.88	119.70
4	H	22	C	OP1-P-O3'	-7.61	88.46	105.20
4	D	23	G	O5'-P-OP2	7.31	119.48	110.70
4	D	10	U	OP2-P-O3'	-7.22	89.31	105.20
4	H	8	C	O3'-P-O5'	7.22	117.72	104.00
4	D	8	C	O3'-P-O5'	7.10	117.49	104.00

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	H	27	U	P-O3'-C3'	6.89	127.97	119.70
4	D	11	C	OP1-P-OP2	6.74	129.71	119.60
4	D	9	C	P-O3'-C3'	6.66	127.69	119.70
4	D	27	U	P-O3'-C3'	6.44	127.43	119.70
4	D	10	U	OP1-P-O3'	-6.22	91.53	105.20
4	H	23	G	O5'-P-OP2	6.20	118.14	110.70
4	D	22	C	OP1-P-O3'	-6.04	91.92	105.20
4	D	11	C	O5'-P-OP2	5.89	117.77	110.70
4	D	33	A	P-O3'-C3'	5.07	125.79	119.70

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	2931	0	2951	34	0
1	E	2931	0	2951	38	0
2	B	1796	0	1828	27	0
2	F	1796	0	1828	41	0
3	C	911	0	961	10	0
3	G	911	0	961	9	0
4	D	700	0	351	6	0
4	H	700	0	351	5	0
5	B	27	0	22	0	0
5	F	27	0	22	0	0
All	All	12730	0	12226	140	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 6.

All (140) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:125:THR:HG21	2:B:130:LEU:HA	1.47	0.96

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:26:ILE:HD12	1:A:53:GLU:HB3	1.64	0.80
1:E:125:THR:HG21	2:F:130:LEU:HA	1.62	0.79
1:E:26:ILE:HD12	1:E:53:GLU:HB3	1.64	0.78
1:E:249:ASN:HB2	2:F:38:VAL:HG11	1.69	0.74
1:A:168:HIS:HB2	1:A:225:ILE:HD12	1.70	0.73
1:E:168:HIS:HB2	1:E:225:ILE:HD12	1.71	0.72
2:B:138:GLN:HE21	2:B:138:GLN:H	1.41	0.69
2:F:70:LYS:H	2:F:212:GLN:HE22	1.41	0.69
2:F:138:GLN:H	2:F:138:GLN:HE21	1.39	0.68
2:B:70:LYS:H	2:B:212:GLN:HE22	1.42	0.68
1:A:293:SER:H	3:C:40:ASN:HD21	1.45	0.65
2:B:80:VAL:HG12	2:B:149:VAL:HB	1.80	0.63
1:E:331:ARG:HA	1:E:334:ARG:HE	1.64	0.63
2:F:80:VAL:HG12	2:F:149:VAL:HB	1.81	0.63
4:H:3:G:H2'	4:H:4:G:H8	1.65	0.62
1:E:90:SER:HB2	2:F:141:LYS:NZ	2.14	0.62
1:E:370:ARG:NH2	4:H:29:A:OP1	2.34	0.61
1:A:125:THR:CG2	2:B:130:LEU:HA	2.26	0.61
4:D:3:G:H2'	4:D:4:G:H8	1.65	0.61
1:A:331:ARG:HA	1:A:334:ARG:HE	1.66	0.61
1:A:40:LEU:HD11	1:A:127:ARG:HG2	1.83	0.61
1:E:90:SER:HB2	2:F:141:LYS:HZ3	1.67	0.60
2:B:114:VAL:HG21	2:B:131:LEU:HD12	1.84	0.60
3:G:30:LYS:HE2	3:G:93:CYS:HA	1.84	0.60
3:C:30:LYS:HE2	3:C:93:CYS:HA	1.83	0.59
1:A:226:SER:HB3	2:F:112:ARG:HG3	1.85	0.59
2:B:28:CYS:SG	2:B:51:TYR:HB3	2.43	0.58
2:F:114:VAL:HG21	2:F:131:LEU:HD12	1.85	0.58
2:F:28:CYS:SG	2:F:51:TYR:HB3	2.44	0.58
1:E:40:LEU:HD11	1:E:127:ARG:HG2	1.84	0.58
2:B:22:ASP:HA	2:F:57:PHE:CZ	2.38	0.57
2:B:154:ILE:HG22	2:B:156:GLN:HG2	1.86	0.57
2:F:154:ILE:HG22	2:F:156:GLN:HG2	1.87	0.57
2:F:83:LEU:HD12	2:F:152:VAL:HG22	1.87	0.57
1:A:291:PRO:HG3	3:C:44:LYS:HG2	1.87	0.56
1:A:100:VAL:HA	1:A:103:ASP:HB2	1.87	0.56
3:C:42:THR:HG23	3:C:103:ALA:HB2	1.88	0.56
1:E:125:THR:CG2	2:F:130:LEU:HA	2.35	0.55
1:E:130:ARG:CZ	1:E:263:PRO:HG3	2.37	0.55
2:F:38:VAL:HG12	2:F:39:TYR:N	2.21	0.55
2:B:16:TYR:HB2	2:B:28:CYS:HB3	1.89	0.54

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:225:ILE:HD11	1:E:139:LEU:HD21	1.88	0.54
1:A:139:LEU:HD21	1:E:225:ILE:HD11	1.89	0.54
1:E:100:VAL:HA	1:E:103:ASP:HB2	1.88	0.54
2:B:83:LEU:HD12	2:B:152:VAL:HG22	1.89	0.54
2:F:16:TYR:HB2	2:F:28:CYS:HB3	1.89	0.53
1:A:235:MET:HE3	2:F:39:TYR:HB3	1.90	0.53
3:G:42:THR:HG23	3:G:103:ALA:HB2	1.89	0.53
1:A:235:MET:HE1	1:E:246:ILE:HA	1.90	0.52
1:A:11:ILE:HB	1:A:104:ILE:HD11	1.92	0.52
1:E:11:ILE:HB	1:E:104:ILE:HD11	1.93	0.51
3:G:30:LYS:CE	3:G:93:CYS:HA	2.41	0.51
1:E:9:HIS:HD2	1:E:11:ILE:HG12	1.75	0.51
1:E:126:ARG:NH2	2:F:122:GLN:HE21	2.09	0.51
2:B:192:PRO:HB3	2:B:195:ILE:HD12	1.93	0.51
1:E:293:SER:H	3:G:40:ASN:HD21	1.59	0.51
1:E:132:ALA:C	1:E:134:GLN:H	2.13	0.51
1:E:112:ASP:HA	1:E:115:ASN:HB2	1.94	0.50
1:A:246:ILE:HA	1:E:235:MET:HE1	1.94	0.50
2:B:145:GLU:HG2	2:B:146:ASN:N	2.27	0.50
1:E:41:ASN:HB3	1:E:46:ILE:HB	1.94	0.50
3:C:30:LYS:CE	3:C:93:CYS:HA	2.41	0.50
1:A:41:ASN:HB3	1:A:46:ILE:HB	1.94	0.50
1:A:9:HIS:HD2	1:A:11:ILE:HG12	1.77	0.49
1:E:129:LEU:HD11	2:F:115:ARG:HD2	1.94	0.49
2:F:145:GLU:HG2	2:F:146:ASN:N	2.27	0.49
1:A:112:ASP:HA	1:A:115:ASN:HB2	1.93	0.49
2:F:81:LEU:HB2	2:F:147:VAL:HG21	1.94	0.49
1:A:293:SER:H	3:C:40:ASN:ND2	2.11	0.48
1:A:296:GLN:NE2	4:D:11:C:H41	2.12	0.48
2:F:192:PRO:HB3	2:F:195:ILE:HD12	1.94	0.48
1:E:94:ARG:HH22	2:F:141:LYS:NZ	2.11	0.48
4:D:27:U:H4'	4:D:28:G:OP2	2.12	0.48
1:E:9:HIS:CE1	1:E:97:LEU:HD11	2.48	0.48
1:A:2:VAL:HB	1:A:61:GLN:HG2	1.95	0.48
2:B:213:ILE:HG12	2:B:226:VAL:HG22	1.96	0.48
3:G:27:ARG:HA	3:G:30:LYS:HD2	1.96	0.47
1:A:254:LEU:HD11	1:A:275:ALA:HB2	1.96	0.47
2:B:81:LEU:HB2	2:B:147:VAL:HG21	1.97	0.47
4:H:27:U:H4'	4:H:28:G:OP2	2.15	0.47
1:E:254:LEU:HD11	1:E:275:ALA:HB2	1.96	0.47
1:A:226:SER:HB3	2:F:112:ARG:HA	1.96	0.47

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:38:VAL:HG12	2:B:39:TYR:N	2.30	0.47
3:C:27:ARG:HA	3:C:30:LYS:HD2	1.96	0.47
2:B:193:LYS:HD3	2:B:194:GLU:HG2	1.98	0.46
1:E:269:VAL:HB	1:E:273:LEU:HD23	1.98	0.46
2:F:193:LYS:HD3	2:F:194:GLU:HG2	1.98	0.46
1:E:245:ASN:HB3	2:F:38:VAL:HG13	1.96	0.46
2:F:213:ILE:HG12	2:F:226:VAL:HG22	1.96	0.46
1:E:2:VAL:HB	1:E:61:GLN:HG2	1.98	0.45
1:A:9:HIS:CE1	1:A:97:LEU:HD11	2.51	0.45
3:G:37:LYS:HG3	4:H:26:G:C8	2.52	0.45
1:A:168:HIS:NE2	1:A:192:ARG:HG3	2.32	0.45
2:B:18:CYS:SG	2:B:28:CYS:SG	3.15	0.45
2:F:36:PHE:HB3	2:F:37:ASN:H	1.56	0.45
2:B:112:ARG:HG3	1:E:226:SER:HB3	2.00	0.44
1:E:245:ASN:CB	2:F:38:VAL:HG13	2.47	0.44
2:B:191:ASP:N	2:B:192:PRO:CD	2.80	0.44
1:A:269:VAL:HB	1:A:273:LEU:HD23	1.99	0.44
2:F:191:ASP:N	2:F:192:PRO:CD	2.81	0.44
2:F:98:ILE:HG23	2:F:102:GLY:HA3	1.99	0.44
2:F:207:ASN:N	2:F:207:ASN:OD1	2.50	0.44
3:G:35:ILE:HG22	3:G:104:ILE:HA	2.00	0.43
1:E:132:ALA:C	1:E:134:GLN:N	2.70	0.43
4:D:16:C:H2'	4:D:17:G:H8	1.84	0.43
1:E:168:HIS:NE2	1:E:192:ARG:HG3	2.33	0.43
1:E:9:HIS:CD2	1:E:11:ILE:HG12	2.54	0.43
4:D:3:G:H2'	4:D:4:G:C8	2.51	0.42
2:B:207:ASN:OD1	2:B:207:ASN:N	2.51	0.42
2:F:70:LYS:N	2:F:212:GLN:HE22	2.13	0.42
1:A:370:ARG:HH22	4:D:28:G:H5''	1.84	0.42
3:C:106:GLU:HA	3:C:107:PRO:HD3	1.91	0.42
1:A:226:SER:HB3	2:F:112:ARG:CG	2.48	0.42
4:H:16:C:H2'	4:H:17:G:H8	1.84	0.42
1:A:132:ALA:C	1:A:134:GLN:H	2.23	0.42
3:C:26:VAL:HG12	3:C:30:LYS:HE3	2.01	0.42
2:F:108:GLU:HG3	2:F:114:VAL:HG23	2.01	0.42
2:B:110:SER:HA	2:B:111:PRO:HD3	1.90	0.42
3:C:35:ILE:HG22	3:C:104:ILE:HA	2.01	0.42
3:G:26:VAL:HG12	3:G:30:LYS:HE3	2.00	0.42
2:F:73:PRO:HB2	2:F:149:VAL:CG2	2.49	0.41
2:F:18:CYS:SG	2:F:28:CYS:SG	3.17	0.41
2:B:98:ILE:HG23	2:B:102:GLY:HA3	2.01	0.41

*Continued on next page...*

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:231:SER:O	1:A:235:MET:HG3	2.20	0.41
2:B:101:ASN:HA	2:B:101:ASN:HD22	1.69	0.41
1:E:316:LYS:O	1:E:318:GLY:N	2.53	0.41
1:E:231:SER:O	1:E:235:MET:HG3	2.20	0.41
1:A:71:VAL:HA	1:A:81:VAL:HG11	2.02	0.41
1:A:84:GLU:HA	1:A:85:PRO:HD3	1.91	0.41
2:F:98:ILE:HG22	2:F:99:GLU:O	2.20	0.41
1:E:127:ARG:HE	1:E:127:ARG:HB2	1.70	0.41
2:B:207:ASN:O	2:B:231:LYS:HG2	2.20	0.41
2:B:98:ILE:HG22	2:B:99:GLU:O	2.21	0.41
2:F:151:TYR:CE2	2:F:153:ASP:HB2	2.55	0.41
2:F:64:ALA:HB2	2:F:225:ILE:HD12	2.03	0.41
1:A:9:HIS:CD2	1:A:11:ILE:HG12	2.56	0.40
2:F:207:ASN:O	2:F:231:LYS:HG2	2.19	0.40
2:B:124:ARG:HA	2:B:125:PRO:HD3	1.85	0.40
3:G:81:VAL:HG22	3:G:121:VAL:HG11	2.03	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	367/388 (95%)	341 (93%)	22 (6%)	4 (1%)	17	64
1	E	367/388 (95%)	342 (93%)	20 (5%)	5 (1%)	14	59
2	B	219/232 (94%)	196 (90%)	19 (9%)	4 (2%)	11	55
2	F	219/232 (94%)	197 (90%)	18 (8%)	4 (2%)	11	55
3	C	118/130 (91%)	117 (99%)	1 (1%)	0	100	100
3	G	118/130 (91%)	117 (99%)	1 (1%)	0	100	100
All	All	1408/1500 (94%)	1310 (93%)	81 (6%)	17 (1%)	16	62

All (17) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	315	PRO
1	A	316	LYS
1	A	317	HIS
1	E	315	PRO
1	E	316	LYS
1	E	317	HIS
1	A	85	PRO
1	E	85	PRO
2	B	38	VAL
1	E	133	ALA
2	F	38	VAL
2	B	192	PRO
2	F	192	PRO
2	B	191	ASP
2	F	191	ASP
2	B	144	VAL
2	F	144	VAL

### 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	307/330 (93%)	287 (94%)	20 (6%)	21	61
1	E	307/330 (93%)	288 (94%)	19 (6%)	23	62
2	B	198/205 (97%)	179 (90%)	19 (10%)	10	44
2	F	198/205 (97%)	180 (91%)	18 (9%)	12	47
3	C	98/107 (92%)	97 (99%)	1 (1%)	82	92
3	G	98/107 (92%)	97 (99%)	1 (1%)	82	92
All	All	1206/1284 (94%)	1128 (94%)	78 (6%)	21	61

All (78) residues with a non-rotameric sidechain are listed below:



Mol	Chain	Res	Type
1	A	24	ASP
1	A	26	ILE
1	A	53	GLU
1	A	64	VAL
1	A	77	LEU
1	A	97	LEU
1	A	117	LEU
1	A	127	ARG
1	A	152	LYS
1	A	169	PHE
1	A	177	GLU
1	A	196	THR
1	A	225	ILE
1	A	226	SER
1	A	244	TYR
1	A	319	ILE
1	A	321	PHE
1	A	341	LEU
1	A	362	GLN
1	A	370	ARG
2	B	8	LYS
2	B	26	ARG
2	B	31	ASN
2	B	60	LYS
2	B	67	LYS
2	B	81	LEU
2	B	103	LYS
2	B	113	VAL
2	B	115	ARG
2	B	123	ARG
2	B	131	LEU
2	B	138	GLN
2	B	150	LEU
2	B	161	ASP
2	B	171	LEU
2	B	197	LYS
2	B	215	ASN
2	B	217	ASP
2	B	229	LYS
3	C	87	LYS
1	E	24	ASP
1	E	53	GLU
1	E	64	VAL

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type
1	E	77	LEU
1	E	97	LEU
1	E	117	LEU
1	E	127	ARG
1	E	152	LYS
1	E	169	PHE
1	E	177	GLU
1	E	196	THR
1	E	225	ILE
1	E	226	SER
1	E	244	TYR
1	E	319	ILE
1	E	321	PHE
1	E	341	LEU
1	E	362	GLN
1	E	370	ARG
2	F	8	LYS
2	F	31	ASN
2	F	60	LYS
2	F	67	LYS
2	F	81	LEU
2	F	103	LYS
2	F	113	VAL
2	F	115	ARG
2	F	123	ARG
2	F	131	LEU
2	F	138	GLN
2	F	150	LEU
2	F	161	ASP
2	F	171	LEU
2	F	197	LYS
2	F	215	ASN
2	F	217	ASP
2	F	229	LYS
3	G	87	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (24) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	9	HIS
1	A	59	ASN
1	A	251	ASN

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type
1	A	296	GLN
2	B	31	ASN
2	B	101	ASN
2	B	138	GLN
2	B	174	ASN
2	B	212	GLN
2	B	215	ASN
3	C	40	ASN
1	E	9	HIS
1	E	59	ASN
1	E	249	ASN
1	E	251	ASN
1	E	296	GLN
2	F	31	ASN
2	F	101	ASN
2	F	122	GLN
2	F	138	GLN
2	F	174	ASN
2	F	212	GLN
2	F	215	ASN
3	G	40	ASN

### 5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
4	D	32/35 (91%)	15 (46%)	5 (15%)
4	H	32/35 (91%)	15 (46%)	4 (12%)
All	All	64/70 (91%)	30 (46%)	9 (14%)

All (30) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
4	D	3	G
4	D	4	G
4	D	5	C
4	D	6	G
4	D	9	C
4	D	10	U
4	D	11	C
4	D	14	A
4	D	15	G

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type
4	D	19	U
4	D	20	C
4	D	23	G
4	D	27	U
4	D	28	G
4	D	33	A
4	H	3	G
4	H	4	G
4	H	5	C
4	H	6	G
4	H	9	C
4	H	10	U
4	H	11	C
4	H	14	A
4	H	15	G
4	H	19	U
4	H	20	C
4	H	23	G
4	H	27	U
4	H	28	G
4	H	33	A

All (9) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
4	D	9	C
4	D	10	U
4	D	14	A
4	D	20	C
4	D	27	U
4	H	9	C
4	H	14	A
4	H	20	C
4	H	27	U

## 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	$\# Z  > 2$
5	SAM	B	301	-	21,29,29	1.20	2 (9%)	17,42,42	2.96	2 (11%)
5	SAM	F	301	-	21,29,29	1.19	2 (9%)	17,42,42	2.97	2 (11%)

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	B	301	-	-	0/8/33/33	0/3/3/3
5	SAM	F	301	-	-	0/8/33/33	0/3/3/3

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	F	301	SAM	C2-N1	2.53	1.38	1.33
5	B	301	SAM	C2-N1	2.63	1.38	1.33
5	F	301	SAM	C2-N3	3.85	1.39	1.32
5	B	301	SAM	C2-N3	3.87	1.39	1.32

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	301	SAM	N3-C2-N1	-11.50	120.09	128.89
5	F	301	SAM	N3-C2-N1	-11.46	120.12	128.89
5	B	301	SAM	O4'-C1'-N9	2.59	113.51	108.10

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	F	301	SAM	O4'-C1'-N9	2.68	113.70	108.10

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

## 6 Fit of model and data ⓘ

### 6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	371/388 (95%)	-0.28	4 (1%)	82	75	63, 206, 210, 251	0
1	E	371/388 (95%)	-0.24	5 (1%)	79	70	63, 206, 210, 251	0
2	B	223/232 (96%)	-0.24	2 (0%)	85	80	90, 104, 180, 211	1 (0%)
2	F	223/232 (96%)	-0.24	1 (0%)	93	90	90, 104, 180, 211	1 (0%)
3	C	120/130 (92%)	0.13	4 (3%)	50	38	207, 209, 210, 210	0
3	G	120/130 (92%)	0.06	1 (0%)	87	82	207, 209, 210, 210	0
4	D	34/35 (97%)	2.00	13 (38%)	0	1	207, 211, 281, 283	0
4	H	34/35 (97%)	2.52	19 (55%)	0	1	207, 211, 281, 283	0
All	All	1496/1570 (95%)	-0.08	49 (3%)	50	38	63, 205, 210, 283	2 (0%)

All (49) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
4	H	9	C	5.4
4	D	23	G	5.2
4	H	22	C	4.6
4	D	5	C	4.5
4	H	33	A	4.3
4	H	5	C	4.1
4	H	23	G	4.1
4	D	21	G	4.1
4	H	34	U	4.1
4	H	4	G	4.0
1	E	208	GLN	4.0
3	C	17	ASP	3.8
4	H	21	G	3.7
4	H	18	U	3.6
4	H	6	G	3.6
3	C	13	GLU	3.6

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	RSRZ
4	D	4	G	3.6
4	H	10	U	3.3
4	H	11	C	3.2
4	D	1	G	3.1
4	D	10	U	3.1
4	D	9	C	3.1
1	A	317	HIS	3.0
1	A	332	TRP	3.0
4	H	32	A	2.8
4	D	22	C	2.8
4	D	3	G	2.7
1	A	376	GLU	2.7
4	H	19	U	2.6
4	H	14	A	2.6
4	D	32	A	2.5
4	H	15	G	2.5
3	C	109	GLU	2.4
4	D	6	G	2.4
4	H	17	G	2.3
1	E	317	HIS	2.3
2	B	19	GLU	2.3
3	C	96	GLN	2.3
2	B	41	GLU	2.3
4	H	13	G	2.3
1	E	332	TRP	2.3
1	E	328	THR	2.2
2	F	39	TYR	2.2
4	H	7	C	2.2
1	A	372	ASP	2.1
4	D	20	C	2.1
1	E	207	GLU	2.1
4	D	14	A	2.1
3	G	15	PRO	2.1

## 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 [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. 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<sup>th</sup> 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( $\text{\AA}^2$ )	Q<0.9
5	SAM	F	301	27/27	0.66	0.53	3.34	145,145,146,147	0
5	SAM	B	301	27/27	0.76	0.48	2.39	145,145,146,147	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.