Full wwPDB X-ray Structure Validation Report (i)

Jan 29, 2024 – 08:36 PM EST

PDB ID	:	1FEU
Title	:	CRYSTAL STRUCTURE OF RIBOSOMAL PROTEIN TL5, ONE OF THE CTC FAMILY PROTEINS, COMPLEXED WITH A FRAGMENT OF 5S RRNA.
Authors	:	Fedorov, R.V.; Meshcheryakov, V.A.; Gongadze, G.M.; Fomenkova, N.P.; Nevskaya, N.A.; Selmer, M.; Laurberg, M.; Kristensen, O.; Al-Karadaghi, S.; Liljas, A.; Garber, M.B.; Nikonov, S.V.
Deposited on	:	2000-07-23
Resolution	:	2.30 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:

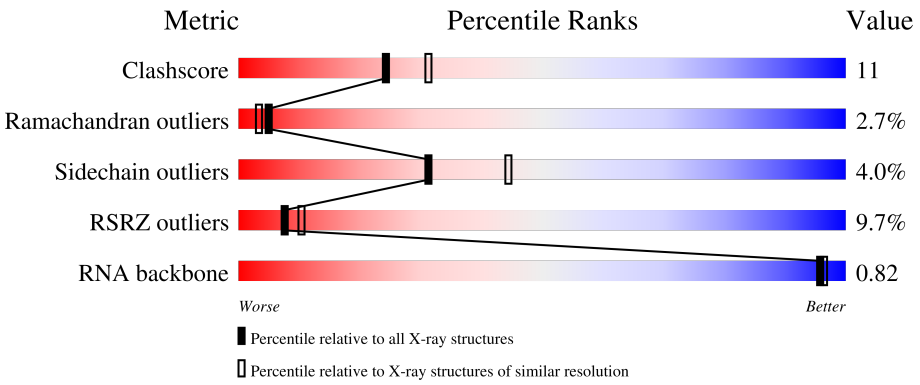
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

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 2.30 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))
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (2.30-2.30)
RNA backbone	3102	1090 (2.70-1.90)

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	B	19	79% 11% 11%
1	E	19	5% 63% 37%
2	C	21	81% 19%
2	F	21	86% 10% 5%

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Mol	Chain	Length	Quality of chain
3	A	206	12% 59% 27% • 10%
3	D	206	9% 68% 21% • 8%

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
4	MG	C	309	-	-	-	X
4	MG	D	310	-	-	-	X

2 Entry composition (i)

There are 6 unique types of molecules in this entry. The entry contains 5023 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 RNA chain called 19 NT FRAGMENT OF 5S RRNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
1	B	19	411	183	76	134	18	0	0	0
1	E	19	411	183	76	134	18	0	0	0

- Molecule 2 is a RNA chain called 21 NT FRAGMENT OF 5S RRNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
2	C	21	450	202	87	141	20	0	0	0
2	F	21	450	202	87	141	20	0	0	0

- Molecule 3 is a protein called 50S RIBOSOMAL PROTEIN L25.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
3	A	185	1476	942	262	269	3	0	0	0
3	D	189	1500	956	266	275	3	0	0	0

- Molecule 4 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Mg		
4	B	2	2	2	0	0
4	C	4	4	4	0	0
4	E	2	2	2	0	0
4	F	5	5	5	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Mg		
4	D	4	4	4	0	0

- Molecule 5 is CADMIUM ION (three-letter code: CD) (formula: Cd).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Cd		
5	A	6	6	6	0	0
			Total	Cd		
5	D	7	7	7	0	0

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
6	B	26	26	26	0	0
			Total	O		
6	C	37	37	37	0	0
			Total	O		
6	E	31	31	31	0	0
			Total	O		
6	F	40	40	40	0	0
			Total	O		
6	A	40	40	40	0	0
			Total	O		
6	D	121	121	121	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: 19 NT FRAGMENT OF 5S RRNA

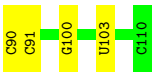
Chain B:

- Molecule 1: 19 NT FRAGMENT OF 5S RRNA

Chain E:

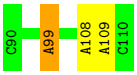
- Molecule 2: 21 NT FRAGMENT OF 5S RRNA

Chain C:



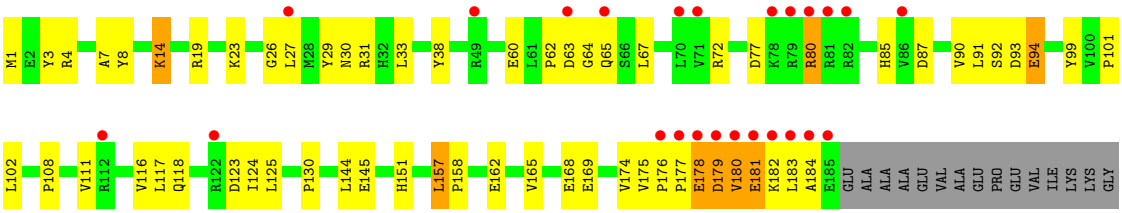
- Molecule 2: 21 NT FRAGMENT OF 5S RRNA

Chain F:



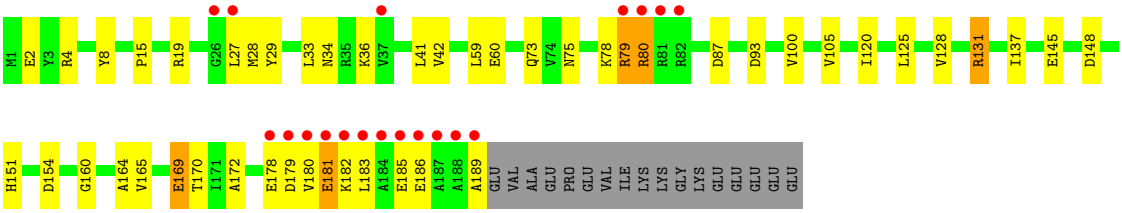
- Molecule 3: 50S RIBOSOMAL PROTEIN L25

Chain A:



• Molecule 3: 50S RIBOSOMAL PROTEIN L25

Chain D:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 31 1 2	Depositor
Cell constants a, b, c, α, β, γ	109.90Å 109.90Å 137.51Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	8.00 – 2.30 19.88 – 2.30	Depositor EDS
% Data completeness (in resolution range)	91.4 (8.00-2.30) 95.3 (19.88-2.30)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	3.87 (at 2.30Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.208 , 0.245 0.220 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å^2)	31.5	Xtriage
Anisotropy	0.125	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.33 , 43.8	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	0.039 for -h,-k,l	Xtriage
F_o,F_c correlation	0.93	EDS
Total number of atoms	5023	wwPDB-VP
Average B, all atoms (Å^2)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.90% 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: MG, CD

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	B	0.24	0/460	0.79	2/718 (0.3%)
1	E	0.25	0/460	0.66	0/718
2	C	0.23	0/504	0.81	5/785 (0.6%)
2	F	0.23	0/504	0.70	0/785
3	A	0.34	0/1508	0.60	0/2046
3	D	0.37	0/1532	0.65	0/2079
All	All	0.31	0/4968	0.68	7/7131 (0.1%)

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	81	G	P-O3'-C3'	6.88	127.96	119.70
1	B	81	G	OP1-P-O3'	6.17	118.77	105.20
2	C	103	U	P-O3'-C3'	-5.61	112.97	119.70
2	C	100	G	OP1-P-O3'	5.55	117.40	105.20
2	C	103	U	OP2-P-O3'	5.43	117.15	105.20
2	C	103	U	OP1-P-O3'	-5.25	93.66	105.20
2	C	100	G	P-O3'-C3'	5.21	125.95	119.70

There are no chirality outliers.

There are no planarity outliers.

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	B	411	0	206	2	0
1	E	411	0	206	8	0
2	C	450	0	230	1	0
2	F	450	0	230	2	0
3	A	1476	0	1506	46	1
3	D	1500	0	1527	41	0
4	B	2	0	0	0	0
4	C	4	0	0	0	0
4	D	4	0	0	0	0
4	E	2	0	0	0	0
4	F	5	0	0	0	0
5	A	6	0	0	0	0
5	D	7	0	0	0	0
6	A	40	0	0	1	0
6	B	26	0	0	0	0
6	C	37	0	0	0	0
6	D	121	0	0	5	0
6	E	31	0	0	0	0
6	F	40	0	0	1	0
All	All	5023	0	3905	95	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 (95) 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)
3:D:93:ASP:HB2	6:D:425:HOH:O	1.73	0.89
1:E:69:G:H3'	1:E:69:G:N3	1.91	0.85
3:D:4:ARG:HD3	3:D:60:GLU:OE1	1.78	0.83
1:E:75:G:H1'	3:D:27:LEU:HD22	1.59	0.83
3:A:30:ASN:HD22	3:A:90:VAL:HB	1.50	0.76
3:A:175:VAL:HB	3:A:180:VAL:HG21	1.73	0.71
3:A:108:PRO:HB2	3:A:111:VAL:HG23	1.72	0.71
3:D:131:ARG:HG2	3:D:131:ARG:HH11	1.56	0.70
3:A:4:ARG:HD3	3:A:60:GLU:OE2	1.92	0.70
3:D:79:ARG:O	3:D:80:ARG:HB2	1.92	0.70
3:A:179:ASP:HB2	3:A:183:LEU:HD22	1.74	0.69
3:D:28:MET:HE1	3:D:59:LEU:HD12	1.77	0.67
3:A:182:LYS:HE2	3:A:182:LYS:N	2.11	0.66
3:D:183:LEU:O	3:D:186:GLU:HG2	1.97	0.65
3:D:125:LEU:HB3	3:D:165:VAL:CG2	2.28	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:33:LEU:HD11	3:A:90:VAL:HG21	1.82	0.62
3:A:31:ARG:HD2	3:A:94:GLU:OE1	1.99	0.62
3:A:29:TYR:HA	3:A:33:LEU:HD12	1.83	0.61
3:A:77:ASP:OD1	3:A:80:ARG:HB2	2.01	0.60
3:D:28:MET:HE1	3:D:59:LEU:CD1	2.32	0.60
3:D:131:ARG:HG2	3:D:131:ARG:NH1	2.18	0.59
3:D:34:ASN:HB2	6:D:382:HOH:O	2.03	0.57
3:D:186:GLU:HA	3:D:189:ALA:HB3	1.87	0.57
3:A:182:LYS:HE2	3:A:182:LYS:CA	2.36	0.56
1:E:69:G:N3	1:E:69:G:C3'	2.67	0.56
3:D:78:LYS:HG2	6:D:379:HOH:O	2.06	0.55
3:D:79:ARG:O	3:D:80:ARG:CB	2.55	0.55
3:A:101:PRO:HA	3:A:123:ASP:OD1	2.07	0.55
3:A:30:ASN:ND2	3:A:90:VAL:HB	2.22	0.54
3:D:125:LEU:HB3	3:D:165:VAL:HG23	1.89	0.54
1:B:75:G:H21	3:A:85:HIS:CE1	2.25	0.54
3:D:165:VAL:CG1	3:D:169:GLU:HG3	2.38	0.54
1:E:74:U:O2	3:D:27:LEU:HD21	2.09	0.53
3:D:120:ILE:HG12	3:D:172:ALA:HA	1.91	0.52
3:A:116:VAL:HB	3:A:175:VAL:HG23	1.91	0.52
3:A:4:ARG:HD3	3:A:60:GLU:CD	2.30	0.52
2:F:99:A:H1'	6:F:346:HOH:O	2.10	0.51
3:D:125:LEU:HG	3:D:164:ALA:HB3	1.91	0.51
3:A:182:LYS:HE2	3:A:182:LYS:H	1.75	0.51
1:B:69:G:H5"	1:B:70:C:OP2	2.11	0.51
3:A:91:LEU:HD22	3:A:130:PRO:HG3	1.92	0.51
3:A:99:TYR:CE1	3:A:125:LEU:HD13	2.46	0.50
3:A:144:LEU:HB3	3:A:174:VAL:HG23	1.93	0.50
3:A:7:ALA:O	3:A:62:PRO:HD3	2.12	0.50
3:D:169:GLU:HA	3:D:169:GLU:OE1	2.12	0.50
3:D:75:ASN:HD22	3:D:75:ASN:N	2.09	0.49
3:A:157:LEU:HD23	3:A:162:GLU:HA	1.93	0.49
2:C:90:C:H2'	2:C:91:C:C6	2.48	0.49
3:A:176:PRO:O	3:A:180:VAL:HB	2.12	0.49
3:A:145:GLU:HB2	6:A:319:HOH:O	2.12	0.48
3:A:14:LYS:N	3:A:14:LYS:HD2	2.28	0.48
3:D:179:ASP:O	3:D:180:VAL:HB	2.13	0.48
1:E:79:G:H2'	1:E:80:U:O4'	2.14	0.48
3:A:124:ILE:HD11	3:A:165:VAL:HG21	1.95	0.48
3:A:8:TYR:HB2	3:A:38:TYR:CE2	2.49	0.47
3:A:182:LYS:HE2	3:A:182:LYS:HA	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:128:VAL:HG23	3:D:160:GLY:O	2.15	0.47
3:A:151:HIS:HD2	3:A:168:GLU:O	1.96	0.47
3:D:120:ILE:HG13	3:D:170:THR:HG22	1.97	0.47
1:E:75:G:H5"	3:D:36:LYS:HD3	1.97	0.46
3:D:29:TYR:HA	3:D:33:LEU:HD12	1.97	0.46
3:A:102:LEU:HD21	3:A:124:ILE:CG2	2.45	0.46
3:D:145:GLU:O	3:D:148:ASP:HB2	2.16	0.45
3:D:42:VAL:HG23	6:D:406:HOH:O	2.15	0.45
3:A:178:GLU:O	3:A:179:ASP:HB3	2.17	0.45
1:E:75:G:H4'	3:D:36:LYS:HG2	1.99	0.44
3:D:8:TYR:CD1	3:D:8:TYR:N	2.86	0.43
2:F:108:A:H2'	2:F:109:A:O4'	2.17	0.43
1:E:86:G:H2'	1:E:87:U:C6	2.54	0.43
3:D:181:GLU:O	3:D:185:GLU:N	2.43	0.43
3:A:108:PRO:HD2	3:A:111:VAL:HG21	2.01	0.42
3:A:180:VAL:HG22	3:A:180:VAL:O	2.18	0.42
3:D:151:HIS:O	3:D:154:ASP:HB2	2.20	0.42
3:A:62:PRO:C	3:A:64:GLY:H	2.23	0.42
3:A:92:SER:OG	3:A:93:ASP:N	2.50	0.42
3:A:179:ASP:HA	3:A:183:LEU:HD13	2.01	0.42
3:D:183:LEU:HD22	3:D:183:LEU:N	2.34	0.42
3:A:1:MET:HG2	3:A:3:TYR:OH	2.20	0.41
3:A:117:LEU:HD12	3:A:118:GLN:N	2.34	0.41
3:A:179:ASP:HB2	3:A:183:LEU:HB2	2.02	0.41
3:A:178:GLU:O	3:A:178:GLU:OE1	2.37	0.41
3:D:180:VAL:CG1	3:D:183:LEU:HB2	2.50	0.41
3:D:181:GLU:O	3:D:185:GLU:HG3	2.21	0.41
3:D:183:LEU:HA	3:D:186:GLU:OE2	2.21	0.41
3:A:181:GLU:O	3:A:182:LYS:HB2	2.21	0.41
3:D:100:VAL:HG11	3:D:137:ILE:HG13	2.02	0.41
3:A:108:PRO:HB3	3:A:174:VAL:HG21	2.02	0.41
3:A:65:GLN:OE1	3:A:67:LEU:HD21	2.20	0.41
3:D:15:PRO:O	3:D:19:ARG:HG3	2.21	0.41
3:D:182:LYS:HA	3:D:185:GLU:OE1	2.20	0.41
3:A:19:ARG:HA	3:A:23:LYS:O	2.20	0.40
3:D:41:LEU:HD22	6:D:348:HOH:O	2.21	0.40
3:A:26:GLY:HA2	3:A:85:HIS:CD2	2.57	0.40
3:A:157:LEU:HA	3:A:158:PRO:HD3	1.83	0.40
3:D:73:GLN:HB3	3:D:87:ASP:HB2	2.03	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)
3:A:80:ARG:NH2	3:A:80:ARG:NH2[6_555]	2.14	0.06

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
3	A	183/206 (89%)	162 (88%)	14 (8%)	7 (4%)	3 1
3	D	187/206 (91%)	170 (91%)	14 (8%)	3 (2%)	9 9
All	All	370/412 (90%)	332 (90%)	28 (8%)	10 (3%)	5 3

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	A	177	PRO
3	A	181	GLU
3	D	80	ARG
3	A	80	ARG
3	D	79	ARG
3	A	184	ALA
3	D	181	GLU
3	A	63	ASP
3	A	179	ASP
3	A	180	VAL

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	
3	A	163/179 (91%)	155 (95%)	8 (5%)	25	35
3	D	164/179 (92%)	159 (97%)	5 (3%)	41	57
All	All	327/358 (91%)	314 (96%)	13 (4%)	31	44

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	A	14	LYS
3	A	27	LEU
3	A	72	ARG
3	A	87	ASP
3	A	94	GLU
3	A	157	LEU
3	A	169	GLU
3	A	178	GLU
3	D	2	GLU
3	D	105	VAL
3	D	131	ARG
3	D	169	GLU
3	D	178	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
3	A	75	ASN
3	A	151	HIS
3	D	75	ASN

5.3.3 RNA (i)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	B	19/19 (100%)	1 (5%)	1 (5%)
1	E	18/19 (94%)	0	0
2	C	20/21 (95%)	0	0
2	F	20/21 (95%)	1 (5%)	0
All	All	77/80 (96%)	2 (2%)	1 (1%)

All (2) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	B	70	C
2	F	99	A

All (1) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
1	B	69	G

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)

Of 30 ligands modelled in this entry, 30 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

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 (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	B	19/19 (100%)	-0.04	0 100 100	21, 30, 46, 56	0
1	E	19/19 (100%)	0.14	1 (5%) 26 33	19, 23, 55, 69	0
2	C	21/21 (100%)	-0.25	0 100 100	19, 32, 46, 48	0
2	F	21/21 (100%)	-0.01	0 100 100	18, 26, 44, 47	0
3	A	185/206 (89%)	0.74	24 (12%) 3 4	25, 41, 68, 70	0
3	D	189/206 (91%)	0.56	19 (10%) 7 9	16, 28, 69, 70	0
All	All	454/492 (92%)	0.53	44 (9%) 7 10	16, 34, 69, 70	0

All (44) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	D	183	LEU	12.4
3	D	187	ALA	11.5
3	D	189	ALA	10.5
3	A	184	ALA	10.4
3	D	188	ALA	10.1
3	A	183	LEU	9.1
3	D	184	ALA	9.0
3	D	180	VAL	8.2
3	D	181	GLU	7.8
3	A	180	VAL	7.3
1	E	69	G	7.3
3	D	179	ASP	7.2
3	D	186	GLU	7.1
3	A	182	LYS	7.0
3	D	185	GLU	6.6
3	A	185	GLU	6.4
3	A	81	ARG	6.4
3	A	79	ARG	5.8
3	A	181	GLU	5.6

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Mol	Chain	Res	Type	RSRZ
3	D	182	LYS	5.6
3	D	79	ARG	4.4
3	A	178	GLU	4.3
3	A	112	ARG	3.8
3	D	178	GLU	3.7
3	A	78	LYS	3.7
3	A	122	ARG	3.6
3	D	81	ARG	3.3
3	A	179	ASP	3.3
3	D	80	ARG	3.3
3	A	80	ARG	3.3
3	A	82	ARG	3.2
3	A	177	PRO	3.2
3	A	27	LEU	2.7
3	D	27	LEU	2.6
3	D	82	ARG	2.5
3	A	70	LEU	2.5
3	A	63	ASP	2.4
3	A	86	VAL	2.4
3	A	49	ARG	2.3
3	D	37	VAL	2.3
3	A	176	PRO	2.2
3	D	26	GLY	2.2
3	A	65	GLN	2.1
3	A	71	VAL	2.1

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
4	MG	C	309	1/1	0.36	0.67	40,40,40,40	1
4	MG	C	318	1/1	0.73	0.15	43,43,43,43	0
4	MG	D	310	1/1	0.73	0.70	47,47,47,47	1
4	MG	C	314	1/1	0.86	0.05	32,32,32,32	0
4	MG	F	317	1/1	0.88	0.12	52,52,52,52	0
4	MG	F	323	1/1	0.89	0.08	38,38,38,38	0
4	MG	D	319	1/1	0.89	0.23	48,48,48,48	0
4	MG	D	322	1/1	0.89	0.19	43,43,43,43	0
4	MG	F	324	1/1	0.90	0.07	32,32,32,32	0
4	MG	F	316	1/1	0.92	0.19	31,31,31,31	0
4	MG	E	315	1/1	0.95	0.19	43,43,43,43	0
4	MG	E	321	1/1	0.95	0.09	38,38,38,38	0
4	MG	F	325	1/1	0.95	0.18	46,46,46,46	0
5	CD	D	307	1/1	0.96	0.04	46,46,46,46	1
4	MG	B	311	1/1	0.97	0.12	42,42,42,42	0
5	CD	A	299	1/1	0.97	0.04	47,47,47,47	1
4	MG	B	313	1/1	0.97	0.10	26,26,26,26	0
4	MG	D	320	1/1	0.98	0.05	28,28,28,28	0
5	CD	A	302	1/1	0.98	0.07	41,41,41,41	1
4	MG	C	312	1/1	0.98	0.15	33,33,33,33	0
5	CD	A	296	1/1	0.99	0.09	47,47,47,47	0
5	CD	A	301	1/1	0.99	0.09	43,43,43,43	1
5	CD	A	297	1/1	0.99	0.05	40,40,40,40	0
5	CD	D	300	1/1	0.99	0.04	42,42,42,42	1
5	CD	D	304	1/1	0.99	0.12	28,28,28,28	0
5	CD	D	305	1/1	0.99	0.07	40,40,40,40	0
5	CD	D	306	1/1	0.99	0.11	30,30,30,30	1
5	CD	A	298	1/1	0.99	0.03	53,53,53,53	0
5	CD	D	308	1/1	0.99	0.12	35,35,35,35	1
5	CD	D	303	1/1	1.00	0.09	33,33,33,33	0

6.5 Other polymers (i)

There are no such residues in this entry.