Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 06:04 PM UTC

PDB ID	:	9RGL / pdb_00009rgl
Title	:	Crystal structure of the SroF-Sar complex, repressor and antirepressor of phage phi3T.
Authors	:	Zamora-Caballero, S.; Marina, A.
Deposited on	:	2025-06-06
Resolution	:	3.05 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 ⓘ 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 ⓘ) were used in the production of this report:

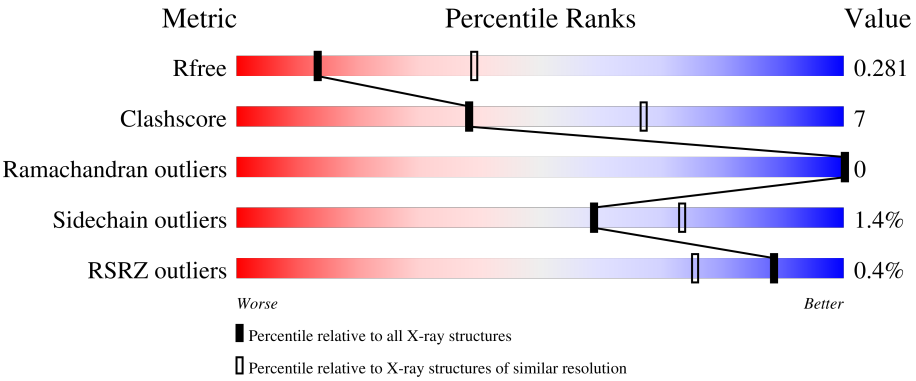
MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

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 3.05 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}	180053	2469 (3.10-3.02)
Clashscore	190562	2569 (3.10-3.02)
Ramachandran outliers	187476	2424 (3.10-3.02)
Sidechain outliers	187428	2423 (3.10-3.02)
RSRZ outliers	180081	2469 (3.10-3.02)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=5% The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	325	% 73% 16% 11%
1	G	325	76% 18% 5%
2	F	113	69% 19% • 12%
2	J	113	68% 19% • 12%

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 6615 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 Core-binding (CB) domain-containing protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	289	Total	C	N	O	S	0	0	0
			2419	1548	405	460	6			
1	G	308	Total	C	N	O	S	0	0	0
			2567	1640	430	489	8			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	expression tag	UNP A0A1P8CWX4
A	0	SER	-	expression tag	UNP A0A1P8CWX4
G	-1	GLY	-	expression tag	UNP A0A1P8CWX4
G	0	SER	-	expression tag	UNP A0A1P8CWX4

- Molecule 2 is a protein called Sar, phi3T_99, SPbeta arbitrium repressor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	F	100	Total	C	N	O	S	0	0	0
			806	507	130	166	3			
2	J	100	Total	C	N	O	S	0	0	0
			802	504	129	166	3			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
F	-1	GLY	-	expression tag	UNP A0A1P8CWX3
F	0	SER	-	expression tag	UNP A0A1P8CWX3
J	-1	GLY	-	expression tag	UNP A0A1P8CWX3
J	0	SER	-	expression tag	UNP A0A1P8CWX3

- Molecule 3 is water.

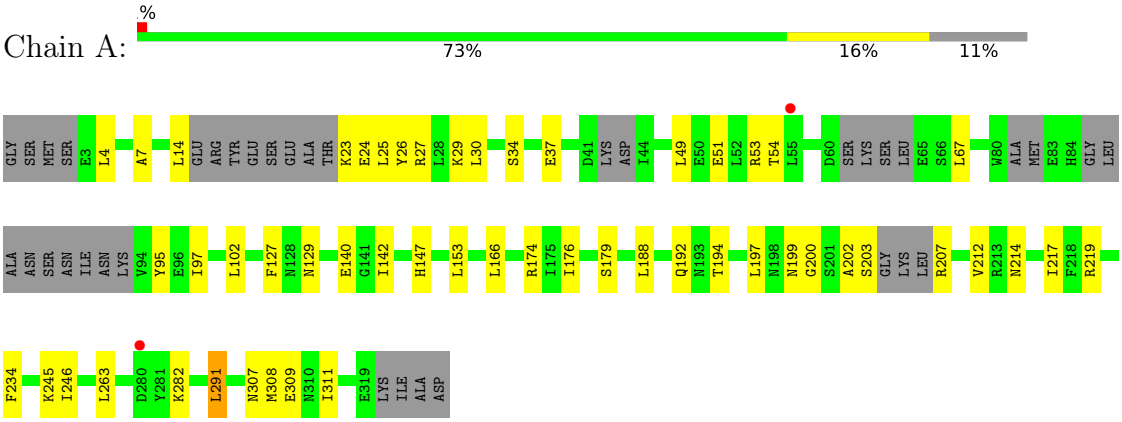
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	7	Total O 7 7	0	0
3	F	2	Total O 2 2	0	0
3	G	10	Total O 10 10	0	0
3	J	2	Total O 2 2	0	0

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.

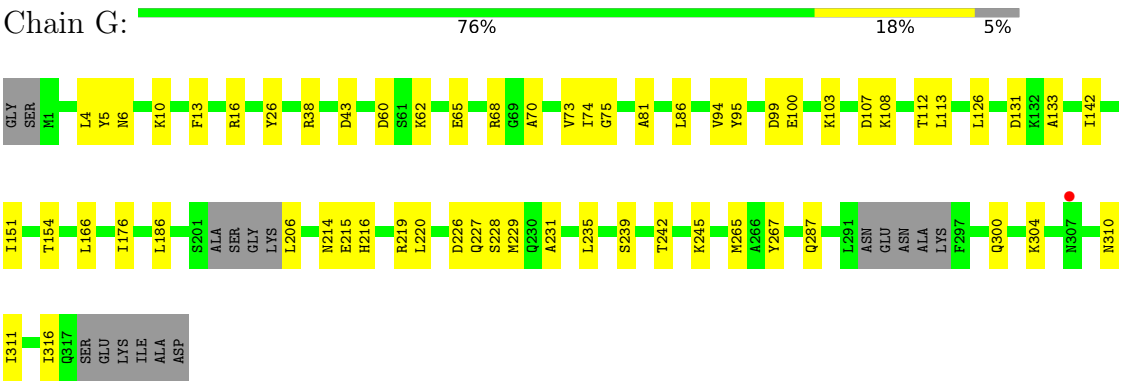
• Molecule 1: Core-binding (CB) domain-containing protein

Chain A:



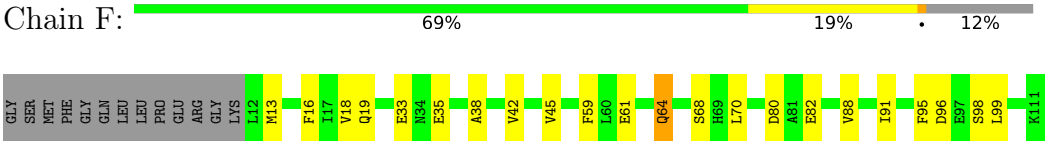
• Molecule 1: Core-binding (CB) domain-containing protein

Chain G:



• Molecule 2: Sar, phi3T_99, SPbeta arbitrium repressor

Chain F:



• Molecule 2: Sar, phi3T_99, SPbeta arbitrium repressor

Chain J:




GLY
SER
MET
PHE
GLY
GLN
LEU
LEU
PRO
GLU
ARG
GLY
K11
L12
M13
I17
V18
S23
R29
I30
E33
N34
E35
E36
Q37
A38
V42
K53
I57
E90
I91
D92
E93
M94
F95
S98
S105
F108
T109
A110
LYS

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	78.94Å 80.98Å 157.97Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	38.53 – 3.05 38.53 – 3.05	Depositor EDS
% Data completeness (in resolution range)	99.5 (38.53-3.05) 99.5 (38.53-3.05)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.64 (at 3.07Å)	Xtriage
Refinement program	PHENIX 1.21.1_5286	Depositor
R, R_{free}	0.240 , 0.282 0.239 , 0.281	Depositor DCC
R_{free} test set	984 reflections (4.26%)	wwPDB-VP
Wilson B-factor (Å^2)	86.5	Xtriage
Anisotropy	0.771	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.31 , 88.8	EDS
L-test for twinning[2]	$< \|L\| > = 0.47$, $< L^2 > = 0.29$	Xtriage
Estimated twinning fraction	0.017 for k,h,-l	Xtriage
F_o,F_c correlation	0.93	EDS
Total number of atoms	6615	wwPDB-VP
Average B, all atoms (Å^2)	127.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.67% 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

5.1 Standard geometry

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	0.10	0/2460	0.29	0/3297
1	G	0.10	0/2613	0.27	0/3507
2	F	0.09	0/817	0.27	0/1094
2	J	0.10	0/813	0.27	0/1090
All	All	0.10	0/6703	0.28	0/8988

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	2419	0	2378	32	0
1	G	2567	0	2541	34	0
2	F	806	0	778	18	0
2	J	802	0	767	19	0
3	A	7	0	0	1	0
3	F	2	0	0	0	0
3	G	10	0	0	1	0
3	J	2	0	0	0	0
All	All	6615	0	6464	90	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 7.

All (90) 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:29:LYS:HD3	1:A:309:GLU:HG3	1.68	0.74
2:F:42:VAL:HG21	2:J:18:VAL:HG11	1.74	0.68
1:A:129:ASN:HD22	1:A:188:LEU:HB3	1.61	0.65
2:J:30:ILE:HD13	2:J:42:VAL:HG13	1.79	0.64
1:G:38:ARG:NH1	2:J:37:GLN:OE1	2.32	0.63
1:A:153:LEU:HG	1:A:217:ILE:HD12	1.83	0.60
1:G:13:PHE:O	1:G:16:ARG:NH1	2.35	0.60
1:A:282:LYS:HD2	1:A:291:LEU:HD11	1.85	0.58
1:A:202:ALA:HB1	1:A:207:ARG:HB3	1.85	0.58
1:A:245:LYS:HD3	2:J:98:SER:HA	1.86	0.57
1:A:25:LEU:HD13	1:A:308:MET:HE2	1.86	0.57
1:G:6:ASN:ND2	1:G:43:ASP:OD2	2.37	0.57
1:A:147:HIS:NE2	3:A:401:HOH:O	2.32	0.56
1:A:203:SER:HB2	1:A:207:ARG:HH21	1.71	0.56
1:G:176:ILE:HD11	1:G:287:GLN:HA	1.88	0.55
2:F:61:GLU:HG2	2:F:64:GLN:HB2	1.88	0.55
1:G:227:GLN:HE21	1:G:228:SER:H	1.55	0.55
2:F:18:VAL:HG11	2:J:42:VAL:HG21	1.87	0.54
1:A:25:LEU:HD22	1:A:308:MET:HG3	1.91	0.53
2:F:35:GLU:HB3	2:F:88:VAL:HG11	1.90	0.53
1:G:62:LYS:NZ	1:G:112:THR:OG1	2.38	0.52
1:G:65:GLU:HG2	1:G:310:ASN:HB3	1.92	0.52
1:G:4:LEU:HD21	1:G:10:LYS:HD2	1.91	0.52
1:A:142:ILE:HG12	1:A:166:LEU:HD11	1.91	0.51
2:F:16:PHE:CE2	2:J:90:GLU:HG2	2.46	0.51
1:G:214:ASN:H	1:G:219:ARG:HD2	1.75	0.51
1:A:4:LEU:HD23	1:A:7:ALA:HA	1.93	0.51
1:A:23:LYS:HB2	1:A:26:TYR:CE2	2.45	0.51
1:G:142:ILE:HG12	1:G:166:LEU:HD11	1.94	0.50
1:G:239:SER:O	1:G:242:THR:OG1	2.29	0.50
1:G:74:ILE:HG22	1:G:94:VAL:HG21	1.94	0.49
2:F:59:PHE:HE2	2:J:53:LYS:HE2	1.78	0.49
1:G:126:LEU:HD13	1:G:131:ASP:HB3	1.94	0.49
2:J:30:ILE:HD11	2:J:38:ALA:HB1	1.94	0.48
1:A:174:ARG:HH11	1:A:176:ILE:HD12	1.78	0.48
2:F:59:PHE:HD1	2:J:57:ILE:HG22	1.78	0.48
1:G:151:ILE:HG23	1:G:231:ALA:HB3	1.96	0.48
1:A:127:PHE:CD2	1:A:200:GLY:HA3	2.47	0.48
1:G:107:ASP:O	1:G:108:LYS:HG2	2.14	0.48
1:A:140:GLU:OE2	1:A:179:SER:OG	2.29	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:J:30:ILE:HG23	2:J:108:PHE:HD2	1.79	0.48
2:F:82:GLU:HB2	2:J:23:SER:HB2	1.96	0.48
1:G:60:ASP:OD1	1:G:60:ASP:N	2.46	0.48
1:G:300:GLN:NE2	3:G:402:HOH:O	2.45	0.47
1:A:34:SER:HA	1:A:37:GLU:HG2	1.97	0.47
1:G:68:ARG:NH1	1:G:99:ASP:OD2	2.46	0.47
1:A:192:GLN:NE2	1:A:194:THR:O	2.42	0.47
2:F:91:ILE:HD11	2:J:17:ILE:HG23	1.96	0.47
2:J:91:ILE:O	2:J:95:PHE:HB2	2.15	0.47
1:G:100:GLU:HA	1:G:103:LYS:HE3	1.97	0.46
1:A:214:ASN:HD22	1:A:219:ARG:HG2	1.80	0.46
1:G:133:ALA:HB1	1:G:186:LEU:HD12	1.96	0.46
1:G:154:THR:HA	1:G:216:HIS:HA	1.98	0.46
2:F:33:GLU:HB2	2:J:13:MET:HE1	1.98	0.45
2:F:42:VAL:HA	2:F:45:VAL:HG12	1.99	0.45
2:F:13:MET:HG2	2:J:33:GLU:HB3	1.97	0.45
1:A:307:ASN:CG	1:A:308:MET:H	2.25	0.45
1:G:81:ALA:HB1	1:G:86:LEU:HB2	1.99	0.45
1:A:212:VAL:O	1:A:219:ARG:HD2	2.17	0.45
1:A:24:GLU:HA	1:A:27:ARG:HB3	1.98	0.44
2:J:29:ARG:NE	2:J:105:SER:O	2.50	0.44
2:J:92:ASP:OD1	2:J:94:MET:N	2.48	0.44
1:A:67:LEU:HB3	1:A:102:LEU:HD22	1.99	0.44
1:A:95:TYR:O	1:A:97:ILE:N	2.51	0.44
1:G:70:ALA:O	1:G:74:ILE:HG13	2.18	0.44
1:A:263:LEU:HD11	1:A:311:ILE:HA	1.99	0.43
1:A:51:GLU:HA	1:A:54:THR:HG22	2.00	0.43
1:G:113:LEU:HD13	1:G:311:ILE:HG21	1.99	0.43
1:G:227:GLN:HG3	1:G:229:MET:H	1.82	0.43
1:G:74:ILE:HB	1:G:94:VAL:HG11	2.00	0.43
1:G:26:TYR:CE2	1:G:73:VAL:HG22	2.54	0.43
1:A:14:LEU:HG	1:A:27:ARG:NE	2.33	0.42
2:J:92:ASP:OD1	2:J:93:GLU:N	2.52	0.42
1:G:75:GLY:HA3	1:G:95:TYR:CZ	2.54	0.42
1:G:176:ILE:HD12	1:G:176:ILE:HA	1.88	0.42
2:F:38:ALA:O	2:F:42:VAL:HG13	2.20	0.42
1:A:95:TYR:O	1:A:97:ILE:HG12	2.20	0.42
2:F:59:PHE:HB2	2:F:70:LEU:HD11	2.02	0.42
1:A:197:LEU:HD22	1:A:207:ARG:HD2	2.01	0.42
1:G:265:MET:HE1	1:G:287:GLN:HG3	2.01	0.42
2:F:13:MET:HE3	1:G:5:TYR:HE1	1.85	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:96:ASP:C	2:F:98:SER:H	2.28	0.41
2:F:95:PHE:HE2	2:F:99:LEU:HB2	1.85	0.41
1:A:234:PHE:HB3	2:F:19:GLN:OE1	2.21	0.41
1:A:49:LEU:O	1:A:53:ARG:HG3	2.21	0.40
1:G:267:TYR:HE1	1:G:316:ILE:HG22	1.86	0.40
1:A:199:ASN:HB2	1:A:246:ILE:HD13	2.02	0.40
1:G:300:GLN:O	1:G:304:LYS:HG2	2.21	0.40
1:G:220:LEU:HD23	1:G:235:LEU:HD13	2.03	0.40
2:J:34:ASN:OD1	2:J:35:GLU:N	2.53	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	275/325 (85%)	265 (96%)	10 (4%)	0	100	100
1	G	302/325 (93%)	287 (95%)	15 (5%)	0	100	100
2	F	98/113 (87%)	95 (97%)	3 (3%)	0	100	100
2	J	98/113 (87%)	94 (96%)	4 (4%)	0	100	100
All	All	773/876 (88%)	741 (96%)	32 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	265/296 (90%)	263 (99%)	2 (1%)	73	79
1	G	282/296 (95%)	278 (99%)	4 (1%)	59	73
2	F	89/99 (90%)	86 (97%)	3 (3%)	32	59
2	J	88/99 (89%)	87 (99%)	1 (1%)	65	76
All	All	724/790 (92%)	714 (99%)	10 (1%)	59	73

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	30	LEU
1	A	291	LEU
2	F	64	GLN
2	F	68	SER
2	F	80	ASP
1	G	206	LEU
1	G	215	GLU
1	G	226	ASP
1	G	245	LYS
2	J	57	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	6	ASN
1	A	230	GLN
2	F	34	ASN
2	F	69	HIS
1	G	116	ASN
1	G	156	ASN
1	G	214	ASN
1	G	216	HIS
2	J	69	HIS

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

There are no oligosaccharides in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

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, 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	289/325 (88%)	-0.22	2 (0%) 84 66	69, 125, 224, 343	0
1	G	308/325 (94%)	-0.26	1 (0%) 90 80	77, 124, 193, 230	0
2	F	100/113 (88%)	-0.31	0 100 100	49, 102, 169, 206	0
2	J	100/113 (88%)	-0.34	0 100 100	66, 96, 161, 222	0
All	All	797/876 (90%)	-0.26	3 (0%) 88 75	49, 116, 204, 343	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	280	ASP	2.5
1	G	307	ASN	2.3
1	A	55	LEU	2.3

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 oligosaccharides in this entry.

6.4 Ligands

There are no ligands in this entry.

6.5 Other polymers

There are no such residues in this entry.