Full wwPDB X-ray Structure Validation Report (i)

Feb 6, 2024 – 03:26 PM EST

PDB ID	:	2EZV
Title	:	Crystal structure of tetrameric restriction endonuclease SfiI bound to cognate DNA.
Authors	:	Aggarwal, A.K.; Vanamee, E.S.; Viadiu, H.
Deposited on	:	2005-11-10
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
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 (1)) were used in the production of this report:

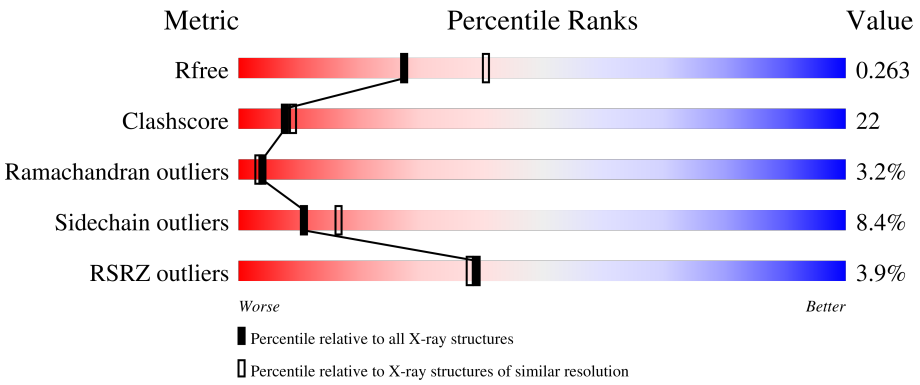
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation 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 2.40 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}	130704	3907 (2.40-2.40)
Clashscore	141614	4398 (2.40-2.40)
Ramachandran outliers	138981	4318 (2.40-2.40)
Sidechain outliers	138945	4319 (2.40-2.40)
RSRZ outliers	127900	3811 (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 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	F	21	5%; 10%; 62%; 29%
2	G	21	5%; 10%; 62%; 29%
3	A	269	2%; 65%; 27%; 7%
3	B	269	5%; 66%; 28%; 6%

2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 5029 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 DNA chain called 5'-D(*TP*GP*GP*CP*CP*AP*AP*CP*AP*AP*GP*GP*CP*CP*T)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
1	F	15	304	145	59	86	14	0	0	0

- Molecule 2 is a DNA chain called 5'-D(*AP*GP*GP*CP*CP*TP*TP*GP*TP*TP*GP*GP*CP*CP*A)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
2	G	15	305	146	55	90	14	0	0	0

- Molecule 3 is a protein called Type II restriction enzyme SfiI.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
3	A	267	2127	1351	372	399	5	0	0	0
3	B	269	2112	1342	366	399	5	0	0	0

- Molecule 4 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Ca		
4	A	1	1	1	0	0
4	B	1	1	1	0	0

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
5	F	8	8	8	0	0

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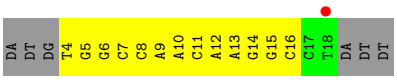
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
5	G	4	4	4	0	0
			Total	O		
5	A	92	92	92	0	0
			Total	O		
5	B	75	75	75	0	0

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.

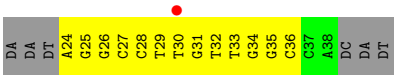
• Molecule 1: 5'-D(*TP*GP*GP*CP*CP*AP*AP*CP*AP*AP*GP*GP*CP*CP*T)-3'

Chain F: 5% 10% 62% 29%



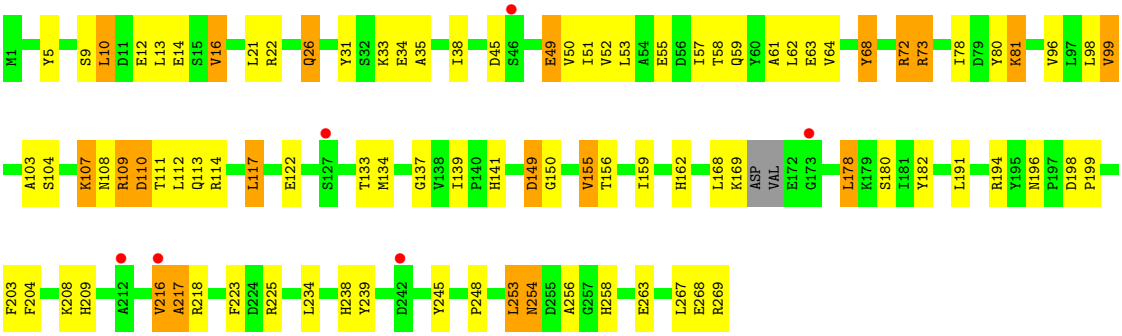
• Molecule 2: 5'-D(*AP*GP*GP*CP*CP*TP*TP*GP*TP*TP*GP*GP*CP*CP*A)-3'

Chain G: 5% 10% 62% 29%



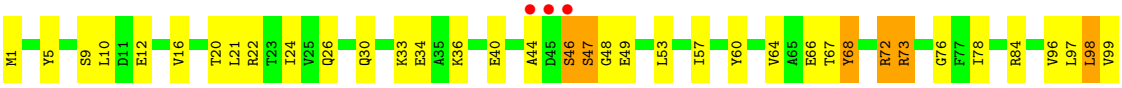
• Molecule 3: Type II restriction enzyme SfiI

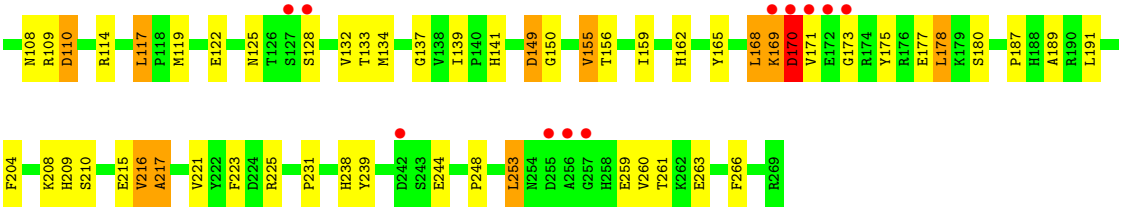
Chain A: 2% 65% 27% 7%



• Molecule 3: Type II restriction enzyme SfiI

Chain B: 5% 66% 28% 6%




N108 R109 D110 R114 L117 P118 M119 E122 N125 T126 S127 S128 V132 T133 M134 G137 V138 I139 P140 H141 D149 G150 V155 T156 I159 H162 Y165 L168 K169 D170 V171 E172 G173 R174 Y175 R176 E177 L178 K179 S180 P187 H188 A189 R190 L191
F204 K208 H209 S210 E215 V216 A217 V221 Y222 F223 D224 R225 P231 H238 Y239 D242 S243 E244 P248 L253 N254 D255 A256 G257 H258 E259 V260 T261 K262 E263 F266 H269

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α, β, γ	85.60Å 85.60Å 202.40Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	50.00 – 2.40 13.98 – 2.38	Depositor EDS
% Data completeness (in resolution range)	95.9 (50.00-2.40) 96.1 (13.98-2.38)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	5.14 (at 2.39Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.224 , 0.265 0.224 , 0.263	Depositor DCC
R_{free} test set	1694 reflections (4.87%)	wwPDB-VP
Wilson B-factor (Å^2)	26.0	Xtriage
Anisotropy	0.209	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.34 , 51.1	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.31$	Xtriage
Estimated twinning fraction	0.036 for -h,-k,l	Xtriage
F_o,F_c correlation	0.92	EDS
Total number of atoms	5029	wwPDB-VP
Average B, all atoms (Å^2)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.75% 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: CA

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	F	0.38	0/341	0.79	0/524
2	G	0.45	0/341	0.85	0/525
3	A	0.37	0/2176	0.67	0/2954
3	B	0.36	0/2161	0.65	1/2939 (0.0%)
All	All	0.37	0/5019	0.69	1/6942 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	B	168	LEU	CA-CB-CG	5.10	127.04	115.30

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	F	304	0	169	31	0
2	G	305	0	171	23	0
3	A	2127	0	2057	82	0
3	B	2112	0	2016	78	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	A	92	0	0	3	0
5	B	75	0	0	3	0
5	F	8	0	0	1	0
5	G	4	0	0	0	0
All	All	5029	0	4413	204	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 22.

All (204) 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:F:15:DG:H2"	1:F:16:DC:H5"	1.26	1.15
1:F:12:DA:H2"	1:F:13:DA:H5'	1.21	1.11
3:A:149:ASP:CG	3:A:150:GLY:H	1.51	1.10
3:A:22:ARG:HH12	3:A:26:GLN:HG2	1.18	1.04
3:B:22:ARG:HH12	3:B:26:GLN:HG2	1.17	1.02
2:G:25:DG:H2"	2:G:26:DG:H5'	1.41	1.01
3:B:149:ASP:CG	3:B:150:GLY:H	1.60	0.99
3:A:80:TYR:HB3	3:A:99:VAL:HG22	1.46	0.97
1:F:10:DA:H2"	1:F:11:DC:H5"	1.47	0.96
1:F:15:DG:C2'	1:F:16:DC:H5"	1.98	0.94
3:A:72:ARG:HH11	3:A:72:ARG:HG3	1.35	0.91
1:F:10:DA:C2'	1:F:11:DC:H5"	2.02	0.90
1:F:5:DG:H2"	1:F:6:DG:H5'	1.54	0.89
3:A:45:ASP:H	3:A:49:GLU:HG3	1.36	0.89
3:A:149:ASP:CG	3:A:150:GLY:N	2.27	0.88
1:F:15:DG:H2"	1:F:16:DC:C5'	2.02	0.88
1:F:4:DT:H2"	1:F:5:DG:H5'	1.56	0.87
2:G:35:DG:H2"	2:G:36:DC:C5'	2.05	0.86
1:F:13:DA:H2"	1:F:14:DG:H5"	1.59	0.84
3:B:22:ARG:NH1	3:B:26:GLN:HG2	1.92	0.84
1:F:12:DA:H2"	1:F:13:DA:C5'	2.09	0.82
3:B:149:ASP:CG	3:B:150:GLY:N	2.31	0.82
3:A:149:ASP:OD1	3:A:150:GLY:N	2.13	0.80
3:A:114:ARG:HA	3:A:134:MET:HE1	1.62	0.80
3:B:24:ILE:HD12	3:B:159:ILE:HD13	1.62	0.79
3:A:209:HIS:CD2	3:A:217:ALA:HA	2.20	0.77
1:F:12:DA:C2'	1:F:13:DA:H5'	2.10	0.76
3:A:254:ASN:HB3	3:A:256:ALA:H	1.49	0.76
1:F:5:DG:C2'	1:F:6:DG:H5'	2.15	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:G:31:DG:H2"	2:G:32:DT:OP2	1.85	0.75
2:G:25:DG:C2'	2:G:26:DG:H5'	2.15	0.75
1:F:10:DA:H2"	1:F:11:DC:C5'	2.17	0.74
2:G:35:DG:H2"	2:G:36:DC:H5'	1.68	0.74
3:A:254:ASN:ND2	3:A:258:HIS:HB3	2.03	0.74
2:G:35:DG:H2"	2:G:36:DC:H5"	1.69	0.73
1:F:4:DT:H2"	1:F:5:DG:C5'	2.17	0.73
3:A:22:ARG:HH12	3:A:26:GLN:CG	1.99	0.73
3:B:109:ARG:N	5:B:285:HOH:O	2.21	0.73
3:B:53:LEU:O	3:B:57:ILE:HG12	1.88	0.73
2:G:28:DC:H2"	2:G:29:DT:OP2	1.89	0.72
2:G:35:DG:C2'	2:G:36:DC:H5"	2.19	0.72
3:A:156:THR:HG23	5:A:308:HOH:O	1.89	0.72
3:B:204:PHE:O	3:B:225:ARG:NH1	2.20	0.72
3:A:114:ARG:HA	3:A:134:MET:CE	2.20	0.71
2:G:34:DG:H5'	3:A:103:ALA:O	1.90	0.70
3:A:22:ARG:NH1	3:A:26:GLN:HG2	2.01	0.70
3:B:169:LYS:HA	3:B:175:TYR:O	1.93	0.69
3:B:44:ALA:HA	3:B:49:GLU:HG3	1.73	0.69
3:B:169:LYS:O	3:B:170:ASP:O	2.11	0.68
3:B:96:VAL:HG12	3:B:98:LEU:HD13	1.77	0.66
1:F:7:DC:H1'	1:F:8:DC:H5"	1.78	0.66
1:F:10:DA:H1'	1:F:11:DC:H5"	1.77	0.66
2:G:32:DT:H2"	2:G:33:DT:H5'	1.78	0.65
3:B:72:ARG:HH11	3:B:72:ARG:HG3	1.61	0.65
3:A:13:LEU:O	3:A:16:VAL:HG13	1.98	0.64
3:A:254:ASN:ND2	5:A:289:HOH:O	2.30	0.64
3:A:72:ARG:HD3	5:A:357:HOH:O	1.97	0.63
3:B:209:HIS:CG	3:B:217:ALA:HA	2.32	0.63
1:F:13:DA:C2'	1:F:14:DG:H5"	2.27	0.63
3:B:96:VAL:HG12	3:B:98:LEU:CD1	2.29	0.63
1:F:10:DA:C1'	1:F:11:DC:H5"	2.29	0.63
3:B:5:TYR:OH	3:B:141:HIS:HD2	1.82	0.62
3:B:117:LEU:HD12	3:B:134:MET:CE	2.30	0.62
3:B:209:HIS:CD2	3:B:217:ALA:HA	2.35	0.62
3:A:216:VAL:O	3:A:217:ALA:CB	2.47	0.61
3:A:35:ALA:HA	3:A:57:ILE:HD12	1.81	0.61
3:A:178:LEU:HD13	3:A:239:TYR:HE1	1.64	0.61
3:A:108:ASN:O	3:A:109:ARG:O	2.18	0.61
3:A:5:TYR:OH	3:A:141:HIS:HD2	1.83	0.61
3:A:10:LEU:HG	3:A:191:LEU:HD23	1.82	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:59:GLN:O	3:A:63:GLU:HG2	2.01	0.60
3:B:109:ARG:HA	3:B:221:VAL:O	2.01	0.60
1:F:5:DG:H1'	1:F:6:DG:H5'	1.83	0.59
1:F:13:DA:N7	5:F:22:HOH:O	2.32	0.59
3:A:38:ILE:HB	3:A:57:ILE:HD11	1.83	0.59
2:G:35:DG:H1'	2:G:36:DC:H5"	1.84	0.59
3:B:216:VAL:O	3:B:217:ALA:HB3	2.01	0.59
3:B:97:LEU:HD11	3:B:159:ILE:HD12	1.85	0.59
3:A:53:LEU:O	3:A:57:ILE:HG12	2.04	0.57
2:G:29:DT:H2"	2:G:30:DT:H5'	1.87	0.57
3:A:107:LYS:HG2	3:A:182:TYR:OH	2.05	0.57
1:F:7:DC:C2'	1:F:8:DC:H5"	2.35	0.56
3:A:52:VAL:HG21	3:B:76:GLY:O	2.04	0.56
3:A:268:GLU:O	3:A:269:ARG:CB	2.53	0.56
3:A:122:GLU:H	3:A:196:ASN:HD21	1.52	0.56
3:A:9:SER:OG	3:A:12:GLU:HG3	2.05	0.55
3:B:99:VAL:HA	3:B:159:ILE:O	2.06	0.55
3:B:114:ARG:O	3:B:137:GLY:HA3	2.06	0.55
3:B:122:GLU:HA	3:B:132:VAL:O	2.06	0.55
3:A:72:ARG:HG3	3:A:72:ARG:NH1	2.09	0.55
3:B:117:LEU:HD12	3:B:134:MET:HE3	1.88	0.55
3:B:5:TYR:HB3	3:B:155:VAL:HG22	1.89	0.55
3:A:111:THR:C	3:A:112:LEU:HD12	2.27	0.54
3:B:72:ARG:HG3	3:B:72:ARG:NH1	2.20	0.54
1:F:7:DC:H2"	1:F:8:DC:C5'	2.37	0.54
3:B:78:ILE:CD1	3:B:98:LEU:HG	2.37	0.54
3:B:180:SER:OG	3:B:238:HIS:HD2	1.90	0.54
1:F:7:DC:H2"	1:F:8:DC:H5"	1.89	0.54
3:A:73:ARG:HD2	3:B:34:GLU:OE1	2.08	0.54
3:A:78:ILE:O	3:A:81:LYS:HE2	2.07	0.54
3:B:33:LYS:HE2	3:B:244:GLU:OE1	2.08	0.54
3:A:178:LEU:HD13	3:A:239:TYR:CE1	2.42	0.53
3:A:33:LYS:HD2	3:A:245:TYR:CE1	2.44	0.52
3:B:40:GLU:OE2	3:B:177:GLU:HA	2.08	0.52
3:B:78:ILE:HD13	3:B:98:LEU:HG	1.91	0.52
2:G:24:DA:H1'	2:G:25:DG:H5"	1.91	0.52
3:B:33:LYS:HE2	3:B:244:GLU:CD	2.30	0.52
1:F:5:DG:C1'	1:F:6:DG:H5'	2.40	0.51
3:B:134:MET:HA	5:B:338:HOH:O	2.09	0.51
1:F:14:DG:H2"	1:F:15:DG:C8	2.45	0.51
3:B:5:TYR:HB3	3:B:155:VAL:CG2	2.40	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:253:LEU:HD23	3:B:253:LEU:O	2.10	0.51
1:F:9:DA:H2"	1:F:10:DA:H5'	1.93	0.51
3:A:113:GLN:NE2	3:A:218:ARG:HD3	2.26	0.51
3:B:110:ASP:OD2	3:B:162:HIS:HD2	1.94	0.51
3:B:24:ILE:CD1	3:B:159:ILE:HD13	2.38	0.50
3:A:122:GLU:N	3:A:196:ASN:HD21	2.09	0.50
3:B:22:ARG:HH12	3:B:26:GLN:CG	2.06	0.50
3:B:175:TYR:CD1	3:B:175:TYR:N	2.79	0.50
3:A:168:LEU:O	3:A:169:LYS:HD3	2.11	0.50
2:G:24:DA:H2"	2:G:25:DG:H5'	1.94	0.50
3:A:113:GLN:HE22	3:A:218:ARG:HD3	1.77	0.49
2:G:35:DG:C1'	2:G:36:DC:H5"	2.41	0.49
2:G:26:DG:H1'	2:G:27:DC:C5	2.47	0.49
3:B:30:GLN:HG2	3:B:266:PHE:HB2	1.94	0.49
3:B:139:ILE:HD12	3:B:141:HIS:O	2.13	0.49
3:B:248:PRO:O	3:B:263:GLU:HA	2.12	0.49
1:F:14:DG:N7	3:B:208:LYS:HD3	2.28	0.49
3:A:254:ASN:HB2	3:A:258:HIS:H	1.77	0.49
3:A:99:VAL:HA	3:A:159:ILE:O	2.13	0.48
3:A:253:LEU:N	3:A:253:LEU:HD23	2.29	0.48
3:A:254:ASN:HD22	3:A:258:HIS:HB3	1.78	0.48
3:A:122:GLU:OE2	3:A:199:PRO:HD3	2.15	0.47
3:B:125:ASN:HB3	3:B:128:SER:OG	2.15	0.47
1:F:11:DC:H2"	1:F:12:DA:C8	2.49	0.47
3:A:168:LEU:O	3:A:169:LYS:HG2	2.14	0.47
3:B:117:LEU:HD22	3:B:119:MET:O	2.15	0.47
3:A:117:LEU:HD12	3:A:134:MET:CE	2.44	0.47
2:G:24:DA:H2"	2:G:25:DG:C5'	2.45	0.47
3:A:58:THR:O	3:A:62:LEU:HG	2.14	0.47
3:B:178:LEU:HD13	3:B:239:TYR:CE1	2.50	0.47
3:A:51:ILE:O	3:A:55:GLU:HG3	2.15	0.46
3:B:9:SER:OG	3:B:12:GLU:HG3	2.14	0.46
1:F:7:DC:C1'	1:F:8:DC:H5"	2.44	0.46
3:A:73:ARG:HB3	3:B:60:TYR:CE1	2.50	0.46
3:A:139:ILE:O	3:A:156:THR:HG21	2.15	0.46
3:B:36:LYS:HD2	3:B:239:TYR:CD2	2.51	0.46
3:A:209:HIS:CG	3:A:217:ALA:HA	2.49	0.46
3:B:66:GLU:OE2	3:B:84:ARG:HD3	2.16	0.46
3:A:117:LEU:HD12	3:A:134:MET:HE3	1.97	0.46
3:B:108:ASN:HA	3:B:223:PHE:CD1	2.51	0.46
3:B:165:TYR:HB2	3:B:177:GLU:O	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:169:LYS:CA	3:B:175:TYR:O	2.62	0.46
3:A:96:VAL:O	3:A:98:LEU:HD22	2.16	0.45
3:B:10:LEU:HG	3:B:191:LEU:HD23	1.98	0.45
3:B:73:ARG:NH1	5:B:336:HOH:O	2.49	0.45
3:B:119:MET:HB3	3:B:189:ALA:HA	1.99	0.45
3:A:81:LYS:N	3:A:81:LYS:HD2	2.32	0.45
3:A:110:ASP:OD2	3:A:162:HIS:HD2	1.99	0.45
3:A:168:LEU:O	3:A:169:LYS:CD	2.65	0.45
3:A:168:LEU:O	3:A:169:LYS:CG	2.65	0.45
3:A:117:LEU:HG	3:A:134:MET:HE2	1.99	0.44
3:A:31:TYR:CD2	3:A:61:ALA:HB2	2.52	0.44
3:B:46:SER:O	3:B:47:SER:O	2.36	0.44
3:B:117:LEU:HD12	3:B:134:MET:HE2	2.00	0.44
3:A:248:PRO:O	3:A:263:GLU:HA	2.17	0.44
2:G:27:DC:H1'	2:G:28:DC:H5'	2.00	0.43
3:B:119:MET:HE3	3:B:187:PRO:O	2.19	0.43
3:B:208:LYS:O	3:B:208:LYS:HG2	2.18	0.43
1:F:11:DC:H2"	1:F:12:DA:O5'	2.17	0.43
1:F:11:DC:H2"	1:F:12:DA:H8	1.83	0.43
3:A:198:ASP:HB2	3:A:199:PRO:HD2	2.01	0.43
3:A:216:VAL:O	3:A:217:ALA:HB2	2.19	0.43
3:A:5:TYR:HB3	3:A:155:VAL:HG22	2.00	0.43
3:A:114:ARG:O	3:A:137:GLY:HA3	2.19	0.43
3:B:231:PRO:HB3	3:B:259:GLU:HG3	2.00	0.43
3:A:180:SER:HB3	3:A:238:HIS:HD2	1.84	0.43
3:B:210:SER:N	3:B:215:GLU:OE2	2.44	0.42
3:B:67:THR:O	3:B:68:TYR:C	2.57	0.42
3:A:34:GLU:OE1	3:B:73:ARG:NH1	2.53	0.42
3:B:216:VAL:O	3:B:217:ALA:CB	2.66	0.42
2:G:34:DG:C5	3:A:208:LYS:HD3	2.55	0.42
3:A:14:GLU:OE1	3:A:194:ARG:NH1	2.50	0.42
3:A:204:PHE:O	3:A:225:ARG:NH2	2.53	0.42
3:A:73:ARG:HG2	3:B:60:TYR:CG	2.55	0.41
3:B:216:VAL:O	3:B:216:VAL:HG12	2.19	0.41
3:B:260:VAL:HG12	3:B:261:THR:N	2.34	0.41
3:A:104:SER:HB3	3:A:162:HIS:CE1	2.56	0.41
3:A:223:PHE:N	3:A:223:PHE:CD2	2.89	0.41
3:A:204:PHE:O	3:A:225:ARG:NH1	2.54	0.41
3:B:24:ILE:HD12	3:B:159:ILE:CD1	2.40	0.41
3:B:44:ALA:O	3:B:49:GLU:HG3	2.20	0.41
2:G:35:DG:C2'	2:G:36:DC:C5'	2.82	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:G:31:DG:H5"	3:B:47:SER:CB	2.50	0.41
2:G:32:DT:H2"	2:G:33:DT:C5'	2.49	0.41
3:B:20:THR:HG22	3:B:159:ILE:HD11	2.03	0.41
3:B:44:ALA:O	3:B:49:GLU:CG	2.69	0.41
3:A:208:LYS:O	3:A:208:LYS:HG2	2.20	0.41
2:G:35:DG:N7	3:A:208:LYS:NZ	2.68	0.40
3:A:49:GLU:O	3:A:52:VAL:HG12	2.22	0.40
3:B:169:LYS:HD2	3:B:175:TYR:O	2.20	0.40
3:A:50:VAL:HA	3:A:53:LEU:HD12	2.04	0.40
3:A:180:SER:HB3	3:A:238:HIS:CD2	2.57	0.40
3:B:170:ASP:OD1	3:B:170:ASP:N	2.54	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
3	A	263/269 (98%)	251 (95%)	6 (2%)	6 (2%)	6 7
3	B	267/269 (99%)	244 (91%)	12 (4%)	11 (4%)	3 2
All	All	530/538 (98%)	495 (93%)	18 (3%)	17 (3%)	4 3

All (17) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	A	109	ARG
3	A	217	ALA
3	A	254	ASN
3	B	47	SER
3	B	169	LYS
3	B	170	ASP
3	B	171	VAL

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Mol	Chain	Res	Type
3	B	173	GLY
3	B	216	VAL
3	A	68	TYR
3	A	216	VAL
3	B	68	TYR
3	B	149	ASP
3	A	149	ASP
3	B	46	SER
3	B	48	GLY
3	B	217	ALA

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	223/237 (94%)	202 (91%)	21 (9%)	8 13
3	B	217/237 (92%)	201 (93%)	16 (7%)	13 22
All	All	440/474 (93%)	403 (92%)	37 (8%)	11 16

All (37) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	A	10	LEU
3	A	16	VAL
3	A	21	LEU
3	A	26	GLN
3	A	49	GLU
3	A	64	VAL
3	A	68	TYR
3	A	72	ARG
3	A	73	ARG
3	A	81	LYS
3	A	99	VAL
3	A	107	LYS
3	A	110	ASP

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Mol	Chain	Res	Type
3	A	117	LEU
3	A	133	THR
3	A	155	VAL
3	A	178	LEU
3	A	203	PHE
3	A	234	LEU
3	A	253	LEU
3	A	267	LEU
3	B	1	MET
3	B	16	VAL
3	B	21	LEU
3	B	64	VAL
3	B	72	ARG
3	B	73	ARG
3	B	98	LEU
3	B	110	ASP
3	B	117	LEU
3	B	133	THR
3	B	155	VAL
3	B	156	THR
3	B	168	LEU
3	B	170	ASP
3	B	178	LEU
3	B	253	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
3	A	26	GLN
3	A	29	GLN
3	A	59	GLN
3	A	113	GLN
3	A	141	HIS
3	A	162	HIS
3	A	196	ASN
3	A	238	HIS
3	A	247	GLN
3	A	254	ASN
3	B	29	GLN
3	B	59	GLN
3	B	141	HIS
3	B	162	HIS

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Mol	Chain	Res	Type
3	B	238	HIS
3	B	247	GLN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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 2 ligands modelled in this entry, 2 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	F	15/21 (71%)	0.67	1 (6%)	17	16	32, 50, 67, 81	0
2	G	15/21 (71%)	0.73	1 (6%)	17	16	31, 49, 68, 70	0
3	A	267/269 (99%)	-0.46	6 (2%)	62	60	10, 20, 44, 60	0
3	B	269/269 (100%)	-0.24	14 (5%)	27	26	11, 27, 50, 67	0
All	All	566/580 (97%)	-0.30	22 (3%)	39	38	10, 24, 53, 81	0

All (22) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	B	45	ASP	5.4
3	A	46	SER	4.9
3	B	171	VAL	3.9
3	B	173	GLY	3.8
3	B	46	SER	3.7
3	B	170	ASP	3.6
3	B	44	ALA	3.1
3	A	242	ASP	3.0
3	A	212	ALA	3.0
3	B	169	LYS	2.8
1	F	18	DT	2.6
3	A	216	VAL	2.5
3	A	173	GLY	2.3
3	B	257	GLY	2.3
3	B	256	ALA	2.3
3	B	172	GLU	2.3
3	B	255	ASP	2.2
2	G	30	DT	2.2
3	B	128	SER	2.1
3	B	127	SER	2.1
3	A	127	SER	2.0
3	B	242	ASP	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 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	CA	A	270	1/1	0.87	0.06	36,36,36,36	0
4	CA	B	270	1/1	0.93	0.06	61,61,61,61	0

6.5 Other polymers (i)

There are no such residues in this entry.