Full wwPDB X-ray Structure Validation Report ⓘ

Aug 28, 2023 – 05:09 PM EDT

PDB ID	:	3KO2
Title	:	I-MsoI re-designed for altered DNA cleavage specificity (-7C)
Authors	:	Taylor, G.K.; Stoddard, B.L.
Deposited on	:	2009-11-13
Resolution	:	2.90 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:

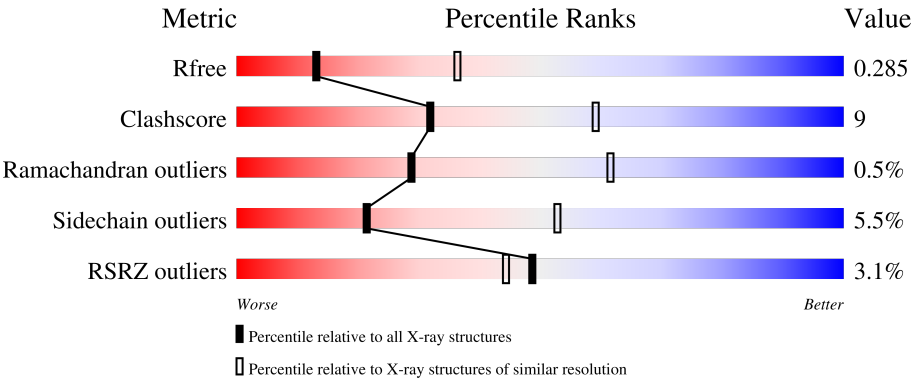
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.35
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35

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 2.90 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	1957 (2.90-2.90)
Clashscore	141614	2172 (2.90-2.90)
Ramachandran outliers	138981	2115 (2.90-2.90)
Sidechain outliers	138945	2117 (2.90-2.90)
RSRZ outliers	127900	1906 (2.90-2.90)

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	A	170	73% 21% • 5%
1	B	170	69% 22% • 5%
1	F	170	74% 19% • 5%
1	G	170	5% 69% 24% • 5%
2	C	24	8% 54% 42% •

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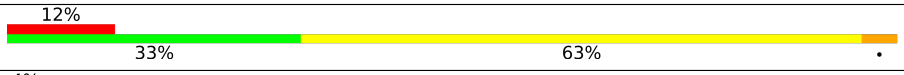
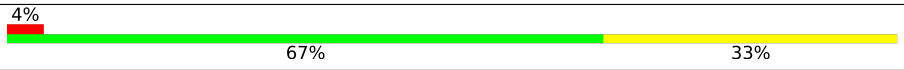
Mol	Chain	Length	Quality of chain
2	H	24	
3	D	24	
3	I	24	

2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 7232 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 Site-specific DNA endonuclease I-MsoI.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
1	A	161	Total	C	N	O	29	0	0
			1316	849	227	240			
1	B	161	Total	C	N	O	33	0	0
			1316	849	227	240			
1	F	161	Total	C	N	O	91	0	0
			1316	849	227	240			
1	G	161	Total	C	N	O	117	0	0
			1316	849	227	240			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	28	ARG	LYS	engineered mutation	UNP C0JWR6
A	43	GLU	SER	engineered mutation	UNP C0JWR6
A	70	THR	ASN	engineered mutation	UNP C0JWR6
A	85	TRP	ILE	engineered mutation	UNP C0JWR6
B	28	ARG	LYS	engineered mutation	UNP C0JWR6
B	43	GLU	SER	engineered mutation	UNP C0JWR6
B	70	THR	ASN	engineered mutation	UNP C0JWR6
B	85	TRP	ILE	engineered mutation	UNP C0JWR6
F	28	ARG	LYS	engineered mutation	UNP C0JWR6
F	43	GLU	SER	engineered mutation	UNP C0JWR6
F	70	THR	ASN	engineered mutation	UNP C0JWR6
F	85	TRP	ILE	engineered mutation	UNP C0JWR6
G	28	ARG	LYS	engineered mutation	UNP C0JWR6
G	43	GLU	SER	engineered mutation	UNP C0JWR6
G	70	THR	ASN	engineered mutation	UNP C0JWR6
G	85	TRP	ILE	engineered mutation	UNP C0JWR6

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	24	Total 491	C 233	N 94	O 141	P 23	0	0	0
2	H	24	Total 491	C 233	N 94	O 141	P 23	0	0	0

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	D	24	Total 487	C 232	N 89	O 143	P 23	0	0	0
3	I	24	Total 487	C 232	N 89	O 143	P 23	14	0	0

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

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	2	Total 2	Ca 2	0	0
4	C	1	Total 1	Ca 1	0	0
4	F	2	Total 2	Ca 2	0	0
4	H	1	Total 1	Ca 1	0	0

- Molecule 5 is water.

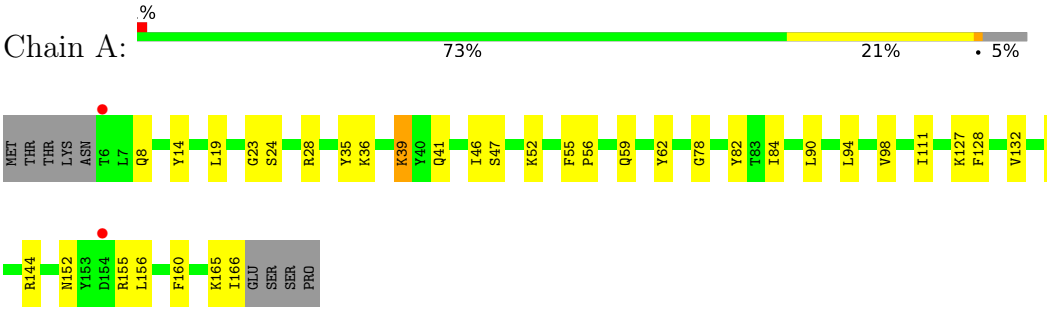
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total 1	O 1	0	0
5	B	3	Total 3	O 3	0	0
5	F	1	Total 1	O 1	0	0
5	G	1	Total 1	O 1	0	0

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.

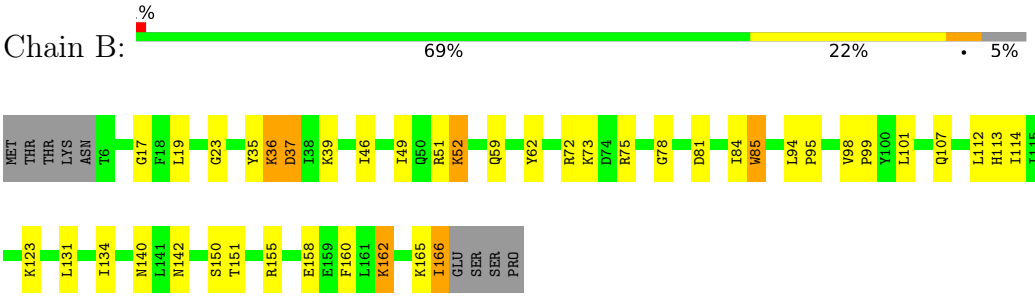
• Molecule 1: Site-specific DNA endonuclease I-MsoI

Chain A:



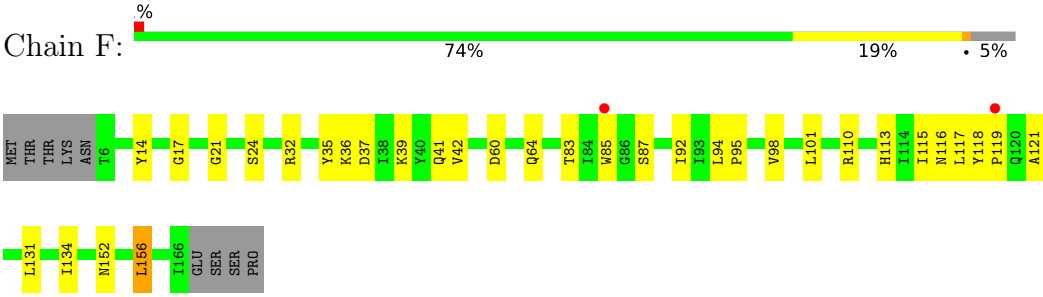
• Molecule 1: Site-specific DNA endonuclease I-MsoI

Chain B:



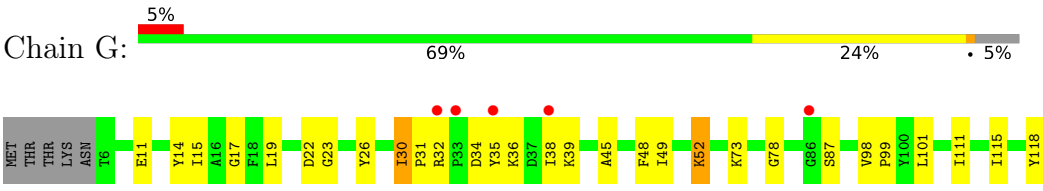
• Molecule 1: Site-specific DNA endonuclease I-MsoI

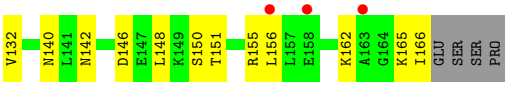
Chain F:



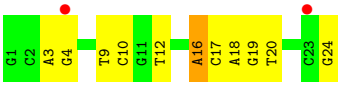
• Molecule 1: Site-specific DNA endonuclease I-MsoI

Chain G:

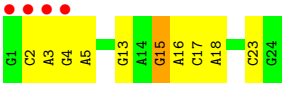




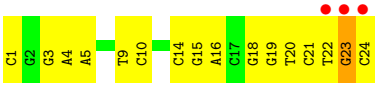
• Molecule 2: 5'-D(*GP*CP*AP*GP*AP*CP*CP*GP*TP*CP*GP*TP*GP*AP*GP*AP*CP*AP*GP*TP*TP*CP*CP*G)-3'



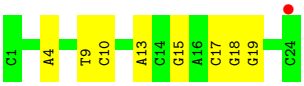
• Molecule 2: 5'-D(*GP*CP*AP*GP*AP*CP*CP*GP*TP*CP*GP*TP*GP*AP*GP*AP*CP*AP*GP*TP*TP*CP*CP*G)-3'



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



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



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α, β, γ	49.03Å 70.01Å 72.26Å 81.66° 70.00° 89.48°	Depositor
Resolution (Å)	33.56 – 2.90 33.56 – 2.90	Depositor EDS
% Data completeness (in resolution range)	95.9 (33.56-2.90) 93.7 (33.56-2.90)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	3.45 (at 2.90Å)	Xtriage
Refinement program	PHENIX 1.5_2	Depositor
R, R_{free}	0.248 , 0.286 0.245 , 0.285	Depositor DCC
R_{free} test set	1898 reflections (10.00%)	wwPDB-VP
Wilson B-factor (Å^2)	52.4	Xtriage
Anisotropy	0.586	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.29 , 27.7	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.30$	Xtriage
Estimated twinning fraction	0.298 for h,-k,h-l	Xtriage
F_o,F_c correlation	0.93	EDS
Total number of atoms	7232	wwPDB-VP
Average B, all atoms (Å^2)	60.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 28.27 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.8994e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

[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

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	A	0.43	0/1342	0.49	0/1812
1	B	0.30	0/1342	0.48	0/1812
1	F	0.28	0/1342	0.43	0/1812
1	G	0.28	0/1342	0.45	0/1812
2	C	0.56	0/551	1.19	3/849 (0.4%)
2	H	0.55	0/551	1.13	2/849 (0.2%)
3	D	0.56	0/545	1.26	3/839 (0.4%)
3	I	0.54	0/545	1.14	0/839
All	All	0.41	0/7560	0.77	8/10624 (0.1%)

There are no bond length outliers.

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	D	16	DA	O4'-C1'-N9	8.44	113.91	108.00
2	C	16	DA	O4'-C1'-N9	8.24	113.77	108.00
3	D	23	DG	O4'-C1'-N9	6.22	112.35	108.00
2	H	23	DC	O4'-C1'-N1	5.76	112.03	108.00
2	C	16	DA	C1'-O4'-C4'	-5.63	104.47	110.10
3	D	14	DC	O4'-C1'-C2'	-5.51	101.49	105.90
2	C	12	DT	N3-C4-O4	5.43	123.16	119.90
2	H	15	DG	P-O3'-C3'	-5.17	113.50	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	A	1316	0	1355	22	0
1	B	1316	0	1355	31	0
1	F	1316	0	1355	23	0
1	G	1316	0	1355	23	0
2	C	491	0	270	8	0
2	H	491	0	270	7	0
3	D	487	0	271	14	0
3	I	487	0	271	7	0
4	A	2	0	0	0	0
4	C	1	0	0	0	0
4	F	2	0	0	0	0
4	H	1	0	0	0	0
5	A	1	0	0	0	0
5	B	3	0	0	0	0
5	F	1	0	0	0	0
5	G	1	0	0	0	0
All	All	7232	0	6502	116	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 9.

All (116) 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:36:LYS:HG3	1:F:37:ASP:H	1.42	0.84
3:D:9:DT:H2"	3:D:10:DC:C5	2.15	0.81
1:G:140:ASN:HD21	1:G:150:SER:H	1.29	0.81
1:B:140:ASN:HD21	1:B:150:SER:H	1.36	0.72
1:A:90:LEU:HB3	1:A:94:LEU:HD12	1.73	0.71
1:A:46:ILE:HD11	1:A:84:ILE:HD12	1.74	0.69
2:C:9:DT:H2"	2:C:10:DC:C5	2.28	0.68
1:A:136:ASP:OD1	1:A:152:ASN:HB2	1.94	0.68
1:B:98:VAL:HA	1:B:101:LEU:HD12	1.74	0.66
1:B:36:LYS:HG2	1:B:37:ASP:N	2.10	0.66
1:A:19:LEU:HD12	1:A:23:GLY:HA3	1.79	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:32:ARG:HH12	3:I:4:DA:H62	1.44	0.64
3:D:18:DG:H2"	3:D:19:DG:O5'	1.98	0.63
1:B:151:THR:HG23	1:B:155:ARG:HE	1.63	0.63
2:H:17:DC:H2'	2:H:18:DA:C8	2.33	0.63
1:G:19:LEU:HD12	1:G:23:GLY:HA3	1.80	0.63
1:G:151:THR:HG23	1:G:155:ARG:HE	1.64	0.62
1:F:98:VAL:HA	1:F:101:LEU:HD12	1.82	0.62
1:B:35:TYR:HB2	1:B:39:LYS:HA	1.82	0.61
1:F:36:LYS:HG3	1:F:37:ASP:N	2.13	0.61
1:A:39:LYS:NZ	1:A:39:LYS:HB3	2.16	0.60
1:A:35:TYR:HB2	1:A:39:LYS:HA	1.84	0.59
1:B:36:LYS:HG2	1:B:37:ASP:H	1.67	0.59
1:B:142:ASN:HB3	2:C:16:DA:H5"	1.86	0.58
1:B:72:ARG:NH2	1:B:75:ARG:HG2	2.17	0.58
2:H:13:DG:N2	3:I:13:DA:C2	2.71	0.58
1:B:59:GLN:O	1:B:62:TYR:HB3	2.05	0.57
1:B:19:LEU:HD23	1:B:101:LEU:HD21	1.85	0.57
1:B:140:ASN:ND2	1:B:150:SER:H	2.03	0.56
1:G:142:ASN:HB3	2:H:16:DA:H5"	1.86	0.56
1:A:52:LYS:HB2	1:A:78:GLY:O	2.05	0.56
1:G:32:ARG:HH12	3:I:4:DA:N6	2.02	0.55
3:I:18:DG:H2"	3:I:19:DG:H5"	1.89	0.55
1:F:113:HIS:NE2	1:F:134:ILE:HD13	2.22	0.54
1:F:152:ASN:O	1:F:156:LEU:HB2	2.08	0.54
2:C:19:DG:H2"	2:C:20:DT:H5'	1.88	0.54
3:D:20:DT:H2"	3:D:21:DC:H5'	1.89	0.54
1:A:28:ARG:NH1	3:D:18:DG:N7	2.56	0.53
1:G:98:VAL:N	1:G:99:PRO:HD2	2.22	0.53
1:B:160:PHE:HB3	1:B:166:ILE:HG22	1.90	0.53
3:D:18:DG:H2"	3:D:19:DG:C5'	2.37	0.53
1:F:118:TYR:N	1:F:119:PRO:CD	2.72	0.53
1:G:19:LEU:HD23	1:G:101:LEU:HD21	1.91	0.53
1:F:113:HIS:HA	1:F:116:ASN:HD22	1.74	0.53
1:G:111:ILE:O	1:G:115:ILE:HG13	2.09	0.52
1:A:39:LYS:HB3	1:A:39:LYS:HZ2	1.74	0.52
1:A:144:ARG:HG2	1:B:51:ARG:NH1	2.25	0.52
1:F:32:ARG:HH21	1:F:41:GLN:HE22	1.56	0.51
3:D:3:DG:H2"	3:D:4:DA:C8	2.47	0.50
1:B:52:LYS:HD2	1:B:78:GLY:HA2	1.94	0.50
3:D:22:DT:H2'	3:D:23:DG:C8	2.47	0.50
1:F:83:THR:HG21	1:F:85:TRP:HE1	1.76	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:158:GLU:O	1:B:162:LYS:HB2	2.13	0.49
2:C:24:DG:N2	3:D:1:DC:H42	2.11	0.49
2:C:24:DG:H22	3:D:1:DC:H42	1.58	0.49
1:F:14:TYR:CE2	1:G:17:GLY:HA2	2.48	0.49
1:A:144:ARG:HG2	1:B:51:ARG:CZ	2.42	0.49
1:B:113:HIS:ND1	1:B:134:ILE:HD13	2.27	0.49
1:G:87:SER:HB2	1:G:115:ILE:HG23	1.95	0.49
3:I:17:DC:H2'	3:I:18:DG:C8	2.47	0.49
1:B:72:ARG:HB3	1:B:81:ASP:OD2	2.13	0.48
1:F:92:ILE:O	1:F:95:PRO:HD2	2.13	0.48
1:F:121:ALA:HA	1:F:127:LYS:HB3	1.96	0.48
1:G:35:TYR:HB3	1:G:39:LYS:HA	1.95	0.48
1:A:55:PHE:HB3	1:A:56:PRO:HD3	1.96	0.48
2:C:10:DC:H42	3:D:15:DG:H1	1.62	0.47
2:H:4:DG:H2"	2:H:5:DA:H8	1.78	0.47
2:C:17:DC:H2'	2:C:18:DA:C8	2.50	0.47
1:G:132:VAL:HG11	1:G:156:LEU:HD12	1.97	0.47
1:F:35:TYR:HB2	1:F:39:LYS:HA	1.97	0.47
1:B:114:ILE:HG23	1:B:131:LEU:HD21	1.96	0.47
1:F:17:GLY:HA2	1:G:14:TYR:CE2	2.50	0.47
1:F:110:ARG:NE	1:F:110:ARG:HA	2.30	0.46
1:B:107:GLN:OE1	1:B:107:GLN:N	2.48	0.46
1:A:160:PHE:CD2	1:A:165:LYS:HD2	2.51	0.46
3:D:18:DG:H2"	3:D:19:DG:H5"	1.98	0.46
3:I:9:DT:H2"	3:I:10:DC:C5	2.51	0.45
1:F:110:ARG:HA	1:F:110:ARG:HE	1.82	0.45
1:A:128:PHE:O	1:A:132:VAL:HG23	2.17	0.45
1:B:85:TRP:HZ3	3:D:5:DA:OP2	2.00	0.44
1:B:46:ILE:O	1:B:46:ILE:HG13	2.17	0.44
1:B:94:LEU:HB2	1:B:95:PRO:HD3	1.99	0.44
1:B:49:ILE:HD12	1:B:49:ILE:N	2.32	0.44
1:F:42:VAL:HG21	1:F:131:LEU:HD22	2.00	0.44
1:F:94:LEU:O	1:F:98:VAL:HG23	2.18	0.44
1:B:72:ARG:CZ	1:B:75:ARG:HG2	2.47	0.43
1:A:24:SER:HB3	1:A:47:SER:OG	2.17	0.43
1:G:118:TYR:N	1:G:119:PRO:CD	2.81	0.43
1:G:140:ASN:ND2	1:G:150:SER:H	2.08	0.43
1:G:26:TYR:CZ	1:G:45:ALA:HB3	2.54	0.43
1:A:35:TYR:CB	1:A:39:LYS:HA	2.48	0.43
1:F:21:GLY:HA2	1:G:22:ASP:OD1	2.19	0.43
1:A:14:TYR:CE2	1:B:17:GLY:HA2	2.53	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:46:ILE:HD11	1:B:84:ILE:HD12	2.01	0.42
1:F:60:ASP:O	1:F:64:GLN:HG3	2.19	0.42
3:D:23:DG:H2"	3:D:24:DC:OP2	2.19	0.42
1:F:36:LYS:HG2	2:H:2:DC:OP2	2.19	0.42
1:A:94:LEU:O	1:A:98:VAL:HG23	2.20	0.42
1:G:30:ILE:HA	1:G:31:PRO:HD3	1.74	0.42
1:A:62:TYR:HD1	1:A:82:TYR:CE1	2.38	0.42
1:B:72:ARG:NH2	1:B:81:ASP:OD1	2.53	0.42
2:C:3:DA:H2"	2:C:4:DG:H5"	2.02	0.41
1:B:98:VAL:N	1:B:99:PRO:CD	2.83	0.41
1:G:52:LYS:HB3	1:G:78:GLY:O	2.20	0.41
1:A:55:PHE:CE2	1:A:59:GLN:HG3	2.56	0.41
1:F:24:SER:HB2	3:I:15:DG:H3'	2.03	0.41
2:H:2:DC:H2"	2:H:3:DA:O4'	2.21	0.41
1:B:19:LEU:HD12	1:B:23:GLY:HA3	2.03	0.41
1:G:11:GLU:O	1:G:15:ILE:HG13	2.20	0.41
1:G:49:ILE:HD13	2:H:15:DG:C8	2.55	0.41
1:F:87:SER:HB2	1:F:115:ILE:HG23	2.03	0.41
1:G:48:PHE:C	1:G:49:ILE:HG13	2.40	0.41
1:A:165:LYS:O	1:A:166:ILE:HG23	2.21	0.40
3:D:19:DG:H5"	3:D:19:DG:H8	1.86	0.40
1:B:118:TYR:N	1:B:119:PRO:CD	2.84	0.40
1:A:135:VAL:O	1:A:139:GLN:HG3	2.22	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	159/170 (94%)	152 (96%)	6 (4%)	1 (1%)	25 58
1	B	159/170 (94%)	149 (94%)	10 (6%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	F	159/170 (94%)	151 (95%)	8 (5%)	0	100 100
1	G	159/170 (94%)	148 (93%)	9 (6%)	2 (1%)	12 37
All	All	636/680 (94%)	600 (94%)	33 (5%)	3 (0%)	29 61

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	G	36	LYS
1	G	38	ILE
1	A	36	LYS

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	142/151 (94%)	135 (95%)	7 (5%)	25 57
1	B	142/151 (94%)	132 (93%)	10 (7%)	15 41
1	F	142/151 (94%)	138 (97%)	4 (3%)	43 76
1	G	142/151 (94%)	132 (93%)	10 (7%)	15 41
All	All	568/604 (94%)	537 (94%)	31 (6%)	21 53

All (31) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	8	GLN
1	A	39	LYS
1	A	41	GLN
1	A	111	ILE
1	A	127	LYS
1	A	155	ARG
1	A	156	LEU
1	B	36	LYS
1	B	37	ASP

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Mol	Chain	Res	Type
1	B	52	LYS
1	B	73	LYS
1	B	85	TRP
1	B	112	LEU
1	B	123	LYS
1	B	162	LYS
1	B	165	LYS
1	B	166	ILE
1	F	117	LEU
1	F	123	LYS
1	F	127	LYS
1	F	156	LEU
1	G	30	ILE
1	G	34	ASP
1	G	52	LYS
1	G	73	LYS
1	G	123	LYS
1	G	146	ASP
1	G	148	LEU
1	G	162	LYS
1	G	165	LYS
1	G	166	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	109	ASN
1	B	140	ASN
1	F	41	GLN
1	F	116	ASN
1	G	50	GLN
1	G	109	ASN
1	G	140	ASN

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 monosaccharides in this entry.

5.6 Ligand geometry

Of 6 ligands modelled in this entry, 6 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

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	161/170 (94%)	-0.26	2 (1%)	79	79	16, 37, 65, 115	9 (5%)
1	B	161/170 (94%)	-0.24	1 (0%)	89	89	13, 37, 65, 83	13 (8%)
1	F	160/170 (94%)	0.25	2 (1%)	77	77	35, 64, 100, 138	26 (16%)
1	G	156/170 (91%)	0.30	8 (5%)	28	24	25, 59, 107, 131	21 (13%)
2	C	24/24 (100%)	0.14	2 (8%)	11	8	22, 62, 108, 136	0
2	H	24/24 (100%)	0.46	4 (16%)	1	1	40, 84, 165, 181	0
3	D	24/24 (100%)	0.05	3 (12%)	3	3	23, 57, 106, 135	0
3	I	24/24 (100%)	0.17	1 (4%)	36	32	38, 82, 119, 173	1 (4%)
All	All	734/776 (94%)	0.03	23 (3%)	49	44	13, 51, 104, 181	70 (9%)

All (23) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	G	35	TYR	5.0
1	G	32	ARG	4.4
2	H	4	DG	4.0
3	I	24	DC	3.6
1	F	85	TRP	3.5
1	G	158	GLU	3.3
1	G	156	LEU	3.3
1	G	163	ALA	2.8
2	C	23	DC	2.8
2	H	1	DG	2.8
2	C	4	DG	2.7
3	D	24	DC	2.6
1	A	6	THR	2.6
1	G	33	PRO	2.6
1	G	38	ILE	2.5
1	F	119	PRO	2.2

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Mol	Chain	Res	Type	RSRZ
2	H	2	DC	2.1
2	H	3	DA	2.1
1	B	116	ASN	2.1
3	D	22	DT	2.1
1	A	154	ASP	2.0
3	D	23	DG	2.0
1	G	86	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 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, 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	B-factors(Å^2)	Q<0.9
4	CA	C	25	1/1	0.59	0.26	132,132,132,132	0
4	CA	F	171	1/1	0.73	0.28	80,80,80,80	0
4	CA	H	25	1/1	0.81	0.10	88,88,88,88	0
4	CA	A	172	1/1	0.84	0.20	46,46,46,46	0
4	CA	A	171	1/1	0.88	0.15	56,56,56,56	0
4	CA	F	172	1/1	0.92	0.15	64,64,64,64	0

6.5 Other polymers (i)

There are no such residues in this entry.