Full wwPDB X-ray Structure Validation Report (i)

Jan 17, 2023 – 05:25 AM EST

PDB ID	:	1VL5
Title	:	CRYSTAL STRUCTURE OF A PUTATIVE METHYLTRANSFERASE (BH2331) FROM BACILLUS HALODURANS C-125 AT 1.95 A RESOLUTION
Authors	:	Joint Center for Structural Genomics (JCSG)
Deposited on	:	2004-07-09
Resolution	:	1.95 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:

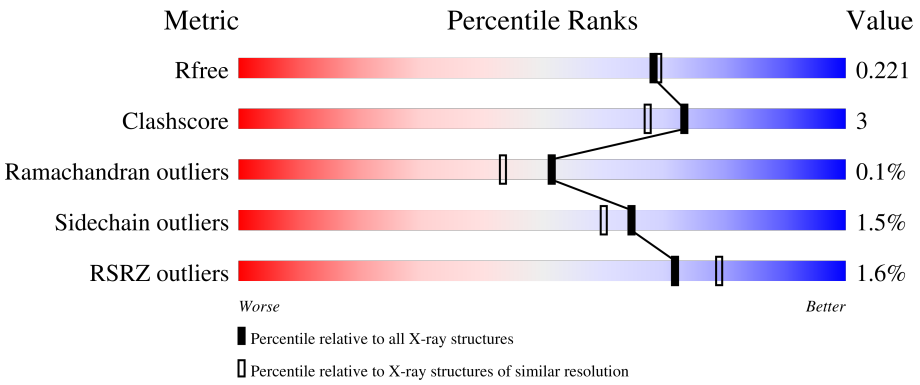
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.31.2
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.31.2

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 1.95 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	2580 (1.96-1.96)
Clashscore	141614	2705 (1.96-1.96)
Ramachandran outliers	138981	2678 (1.96-1.96)
Sidechain outliers	138945	2678 (1.96-1.96)
RSRZ outliers	127900	2539 (1.96-1.96)

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	260	82% 7% 11%
1	B	260	85% • 11%
1	C	260	2% 74% 8% 18%
1	D	260	3% 80% 8% • 11%

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 7836 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 unknown conserved protein BH2331.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	S	Se			
1	A	231	1846	1184	311	342	3	6	0	1	0
1	B	231	1840	1183	308	340	3	6	0	1	0
1	C	213	1697	1094	284	311	3	5	0	0	0
1	D	231	1819	1169	306	336	3	5	0	0	0

There are 72 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	6	MSE	-	expression tag	UNP Q9KAF6
A	7	GLY	-	expression tag	UNP Q9KAF6
A	8	SER	-	expression tag	UNP Q9KAF6
A	9	ASP	-	expression tag	UNP Q9KAF6
A	10	LYS	-	expression tag	UNP Q9KAF6
A	11	ILE	-	expression tag	UNP Q9KAF6
A	12	HIS	-	expression tag	UNP Q9KAF6
A	13	HIS	-	expression tag	UNP Q9KAF6
A	14	HIS	-	expression tag	UNP Q9KAF6
A	15	HIS	-	expression tag	UNP Q9KAF6
A	16	HIS	-	expression tag	UNP Q9KAF6
A	17	HIS	-	expression tag	UNP Q9KAF6
A	18	MSE	-	expression tag	UNP Q9KAF6
A	35	MSE	MET	modified residue	UNP Q9KAF6
A	102	MSE	LEU	SEE REMARK 999	UNP Q9KAF6
A	181	MSE	MET	modified residue	UNP Q9KAF6
A	208	MSE	MET	modified residue	UNP Q9KAF6
A	252	MSE	MET	modified residue	UNP Q9KAF6
B	6	MSE	-	expression tag	UNP Q9KAF6
B	7	GLY	-	expression tag	UNP Q9KAF6
B	8	SER	-	expression tag	UNP Q9KAF6

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Chain	Residue	Modelled	Actual	Comment	Reference
B	9	ASP	-	expression tag	UNP Q9KAF6
B	10	LYS	-	expression tag	UNP Q9KAF6
B	11	ILE	-	expression tag	UNP Q9KAF6
B	12	HIS	-	expression tag	UNP Q9KAF6
B	13	HIS	-	expression tag	UNP Q9KAF6
B	14	HIS	-	expression tag	UNP Q9KAF6
B	15	HIS	-	expression tag	UNP Q9KAF6
B	16	HIS	-	expression tag	UNP Q9KAF6
B	17	HIS	-	expression tag	UNP Q9KAF6
B	18	MSE	-	expression tag	UNP Q9KAF6
B	35	MSE	MET	modified residue	UNP Q9KAF6
B	102	MSE	LEU	SEE REMARK 999	UNP Q9KAF6
B	181	MSE	MET	modified residue	UNP Q9KAF6
B	208	MSE	MET	modified residue	UNP Q9KAF6
B	252	MSE	MET	modified residue	UNP Q9KAF6
C	6	MSE	-	expression tag	UNP Q9KAF6
C	7	GLY	-	expression tag	UNP Q9KAF6
C	8	SER	-	expression tag	UNP Q9KAF6
C	9	ASP	-	expression tag	UNP Q9KAF6
C	10	LYS	-	expression tag	UNP Q9KAF6
C	11	ILE	-	expression tag	UNP Q9KAF6
C	12	HIS	-	expression tag	UNP Q9KAF6
C	13	HIS	-	expression tag	UNP Q9KAF6
C	14	HIS	-	expression tag	UNP Q9KAF6
C	15	HIS	-	expression tag	UNP Q9KAF6
C	16	HIS	-	expression tag	UNP Q9KAF6
C	17	HIS	-	expression tag	UNP Q9KAF6
C	18	MSE	-	expression tag	UNP Q9KAF6
C	35	MSE	MET	modified residue	UNP Q9KAF6
C	102	MSE	LEU	SEE REMARK 999	UNP Q9KAF6
C	181	MSE	MET	modified residue	UNP Q9KAF6
C	208	MSE	MET	modified residue	UNP Q9KAF6
C	252	MSE	MET	modified residue	UNP Q9KAF6
D	6	MSE	-	expression tag	UNP Q9KAF6
D	7	GLY	-	expression tag	UNP Q9KAF6
D	8	SER	-	expression tag	UNP Q9KAF6
D	9	ASP	-	expression tag	UNP Q9KAF6
D	10	LYS	-	expression tag	UNP Q9KAF6
D	11	ILE	-	expression tag	UNP Q9KAF6
D	12	HIS	-	expression tag	UNP Q9KAF6
D	13	HIS	-	expression tag	UNP Q9KAF6
D	14	HIS	-	expression tag	UNP Q9KAF6

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Chain	Residue	Modelled	Actual	Comment	Reference
D	15	HIS	-	expression tag	UNP Q9KAF6
D	16	HIS	-	expression tag	UNP Q9KAF6
D	17	HIS	-	expression tag	UNP Q9KAF6
D	18	MSE	-	expression tag	UNP Q9KAF6
D	35	MSE	MET	modified residue	UNP Q9KAF6
D	102	MSE	LEU	SEE REMARK 999	UNP Q9KAF6
D	181	MSE	MET	modified residue	UNP Q9KAF6
D	208	MSE	MET	modified residue	UNP Q9KAF6
D	252	MSE	MET	modified residue	UNP Q9KAF6

- Molecule 2 is water.

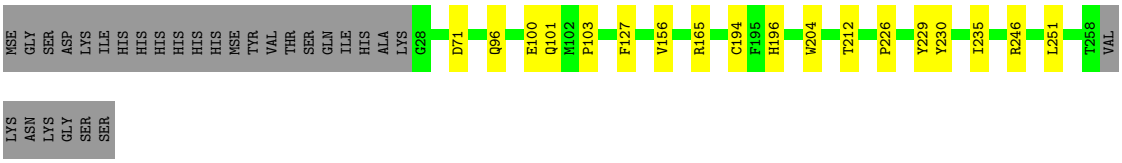
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
2	A	199	199	199	0	0
			Total	O		
2	B	166	166	166	0	0
			Total	O		
2	C	148	148	148	0	0
			Total	O		
2	D	121	121	121	0	0

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: unknown conserved protein BH2331

Chain A:

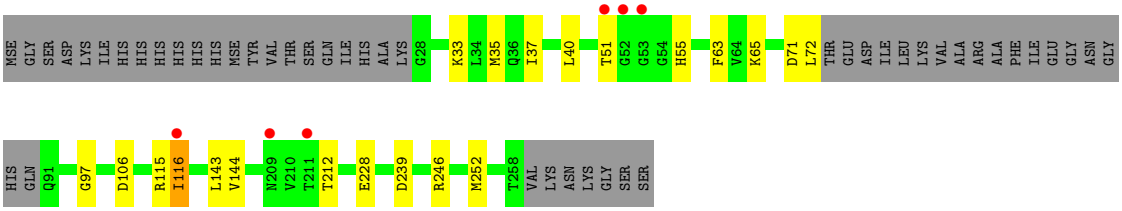


• Molecule 1: unknown conserved protein BH2331

Chain B:

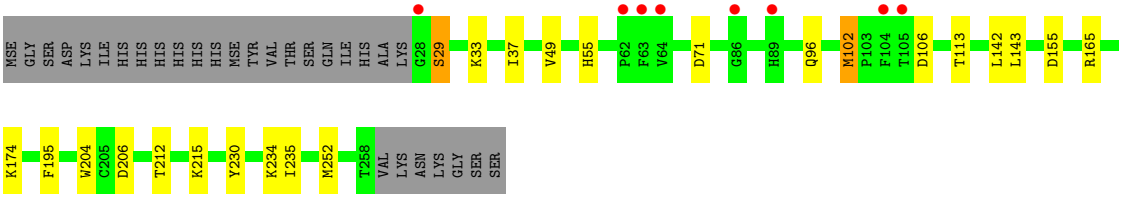
• Molecule 1: unknown conserved protein BH2331

Chain C:



• Molecule 1: unknown conserved protein BH2331

Chain D:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	64.01Å 87.47Å 105.24Å 90.00° 100.57° 90.00°	Depositor
Resolution (Å)	49.83 – 1.95 49.85 – 1.85	Depositor EDS
% Data completeness (in resolution range)	100.0 (49.83-1.95) 92.9 (49.85-1.85)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
$< I/\sigma(I) >$ [1]	1.58 (at 1.84Å)	Xtriage
Refinement program	REFMAC 5.2.0001	Depositor
R, R_{free}	0.177 , 0.209 0.189 , 0.221	Depositor DCC
R_{free} test set	4076 reflections (4.50%)	wwPDB-VP
Wilson B-factor (Å^2)	21.2	Xtriage
Anisotropy	0.573	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.35 , 46.3	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	7836	wwPDB-VP
Average B, all atoms (Å^2)	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.24% 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)

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	0.89	1/1884 (0.1%)	0.87	3/2534 (0.1%)
1	B	0.82	0/1880	0.79	0/2532
1	C	0.85	1/1734 (0.1%)	0.86	4/2331 (0.2%)
1	D	0.77	0/1859	0.78	2/2505 (0.1%)
All	All	0.83	2/7357 (0.0%)	0.83	9/9902 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand.A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	C	0	1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	212	THR	C-O	5.34	1.33	1.23
1	A	194	CYS	CB-SG	-5.19	1.73	1.81

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	246	ARG	NE-CZ-NH1	13.41	127.01	120.30
1	C	246	ARG	NE-CZ-NH2	-13.19	113.70	120.30
1	C	246	ARG	NE-CZ-NH1	12.88	126.74	120.30
1	A	246	ARG	NE-CZ-NH2	-12.48	114.06	120.30
1	C	106	ASP	CB-CG-OD1	5.85	123.57	118.30
1	D	106	ASP	CB-CG-OD1	5.30	123.08	118.30
1	A	251	LEU	CA-CB-CG	5.29	127.47	115.30
1	C	239	ASP	CB-CG-OD1	5.16	122.94	118.30
1	D	102	MSE	CG-SE-CE	-5.14	87.60	98.90

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	71	ASP	Peptide

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	1846	0	1757	10	0
1	B	1840	0	1746	15	0
1	C	1697	0	1622	10	0
1	D	1819	0	1718	12	0
2	A	199	0	0	1	0
2	B	166	0	0	0	0
2	C	148	0	0	2	0
2	D	121	0	0	0	0
All	All	7836	0	6843	46	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 3.

All (46) 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:100:GLU:O	1:A:101:GLN:N	1.86	1.08
1:A:100:GLU:O	1:A:100:GLU:CA	2.11	0.99
1:A:100:GLU:CA	1:A:101:GLN:N	2.34	0.91
1:B:102[B]:MSE:HE2	1:B:130:GLU:HB3	1.63	0.80
1:B:102[B]:MSE:HE3	1:B:134:VAL:HG11	1.66	0.76
1:A:156:VAL:HG12	1:B:156:VAL:HG12	1.72	0.71
1:B:102[B]:MSE:HE1	1:B:134:VAL:CG2	2.27	0.64
1:B:102[B]:MSE:HE1	1:B:134:VAL:HG22	1.80	0.64
1:B:102[B]:MSE:HE3	1:B:134:VAL:CG1	2.28	0.62
1:B:143:LEU:HD23	1:B:252:MSE:CE	2.30	0.62
1:B:143:LEU:HD23	1:B:252:MSE:HE3	1.84	0.59
1:D:29:SER:HB2	1:D:195:PHE:CD2	2.39	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:72:LEU:HG	1:C:97:GLY:O	2.05	0.56
1:B:102[B]:MSE:CE	1:B:134:VAL:CG1	2.84	0.55
1:D:33:LYS:O	1:D:37:ILE:HG12	2.06	0.54
1:C:33:LYS:HE3	1:C:37:ILE:HD11	1.91	0.52
1:D:143:LEU:HD23	1:D:252:MSE:CE	2.40	0.51
1:A:165:ARG:HD2	1:A:204:TRP:CZ2	2.48	0.49
1:C:116:ILE:O	1:C:116:ILE:HG22	2.14	0.48
1:D:165:ARG:HD2	1:D:204:TRP:CZ2	2.50	0.47
1:A:230:TYR:HA	1:A:235:ILE:HD12	1.97	0.47
1:A:226:PRO:HD2	1:A:229:TYR:CD1	2.50	0.46
1:C:228:GLU:CG	2:C:349:HOH:O	2.63	0.46
1:D:143:LEU:HD23	1:D:252:MSE:HE3	1.98	0.46
1:C:35:MSE:HE1	1:C:63:PHE:CD2	2.50	0.46
1:A:71:ASP:O	1:A:96:GLN:HA	2.17	0.45
1:B:143:LEU:HD23	1:B:252:MSE:HE1	1.98	0.45
1:C:115:ARG:HG2	1:C:116:ILE:HG12	1.98	0.45
1:D:155:ASP:OD2	1:D:174:LYS:HE3	2.16	0.45
1:C:143:LEU:HD22	1:C:252:MSE:CE	2.47	0.44
1:C:55:HIS:HB2	2:C:407:HOH:O	2.17	0.44
1:D:71:ASP:O	1:D:96:GLN:HA	2.18	0.43
1:B:192:LEU:HD13	1:B:252:MSE:CG	2.50	0.42
1:C:115:ARG:HG3	1:C:144:VAL:HB	2.00	0.42
1:D:49:VAL:HG22	1:D:102:MSE:HE1	2.00	0.42
1:D:230:TYR:HA	1:D:235:ILE:HD12	2.02	0.42
1:B:192:LEU:HD13	1:B:252:MSE:HG2	2.01	0.41
1:B:102[B]:MSE:CE	1:B:134:VAL:HG11	2.42	0.41
1:B:165:ARG:HD2	1:B:204:TRP:CZ2	2.56	0.41
1:A:196:HIS:HD2	2:A:438:HOH:O	2.03	0.41
1:D:206:ASP:OD1	1:D:215:LYS:NZ	2.41	0.41
1:D:234:LYS:HD3	1:D:234:LYS:HA	1.89	0.41
1:C:35:MSE:HE3	1:C:40:LEU:HD12	2.03	0.41
1:B:102[B]:MSE:CE	1:B:134:VAL:HG13	2.50	0.40
1:A:100:GLU:O	1:A:101:GLN:CA	2.64	0.40
1:D:113:THR:HG22	1:D:142:LEU:HB2	2.04	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	228/260 (88%)	222 (97%)	6 (3%)	0	100	100
1	B	230/260 (88%)	225 (98%)	5 (2%)	0	100	100
1	C	209/260 (80%)	205 (98%)	3 (1%)	1 (0%)	29	17
1	D	229/260 (88%)	223 (97%)	6 (3%)	0	100	100
All	All	896/1040 (86%)	875 (98%)	20 (2%)	1 (0%)	51	43

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	116	ILE

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	188/217 (87%)	185 (98%)	3 (2%)	62	58
1	B	185/217 (85%)	181 (98%)	4 (2%)	52	44
1	C	173/217 (80%)	171 (99%)	2 (1%)	71	68
1	D	182/217 (84%)	179 (98%)	3 (2%)	62	58
All	All	728/868 (84%)	716 (98%)	12 (2%)	65	58

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	103	PRO
1	A	127	PHE
1	A	212	THR
1	B	55	HIS
1	B	102[A]	MSE
1	B	102[B]	MSE
1	B	212	THR
1	C	51	THR
1	C	65	LYS
1	D	29	SER
1	D	55	HIS
1	D	212	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	D	87	ASN

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)

There are no ligands in this entry.

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	A	226/260 (86%)	-0.06	0 100 100	24, 30, 39, 46	0
1	B	226/260 (86%)	-0.27	0 100 100	22, 30, 38, 47	0
1	C	208/260 (80%)	-0.09	6 (2%) 51 60	22, 31, 43, 59	0
1	D	226/260 (86%)	-0.00	8 (3%) 44 53	22, 31, 40, 53	0
All	All	886/1040 (85%)	-0.11	14 (1%) 72 79	22, 30, 40, 59	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	86	GLY	4.2
1	D	89	HIS	3.3
1	C	51	THR	3.0
1	D	104	PHE	2.9
1	C	53	GLY	2.9
1	C	116	ILE	2.8
1	C	209	ASN	2.7
1	D	28	GLY	2.3
1	D	63	PHE	2.3
1	C	52	GLY	2.2
1	D	64	VAL	2.2
1	D	62	PRO	2.2
1	D	105	THR	2.2
1	C	211	THR	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)

There are no ligands in this entry.

6.5 Other polymers (i)

There are no such residues in this entry.