



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 10:21 PM GMT

PDB ID : 1T82
Title : Crystal Structure of the putative thioesterase from *Shewanella oneidensis*, Northeast Structural Genomics Target SoR51
Authors : Forouhar, F.; Lee, I.; Vorobiev, S.M.; Xiao, R.; Acton, T.B.; Montelione, G.T.; Hunt, J.F.; Tong, L.; Northeast Structural Genomics Consortium (NESG)
Deposited on : 2004-05-11
Resolution : 1.70 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

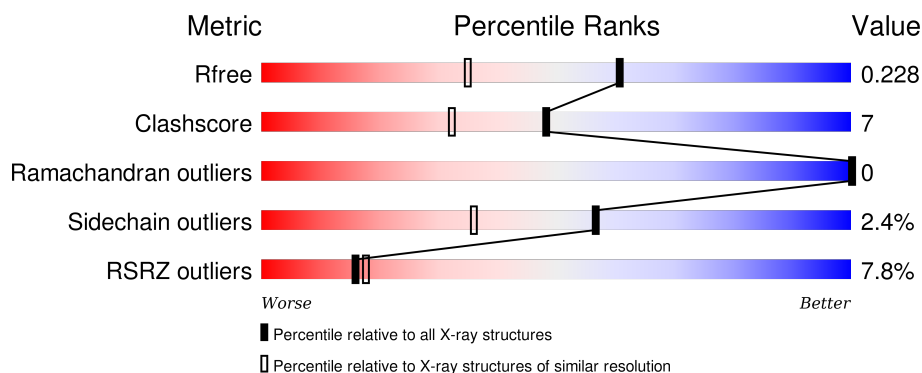
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.70 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}	91344	3190 (1.70-1.70)
Clashscore	102246	3585 (1.70-1.70)
Ramachandran outliers	100387	3527 (1.70-1.70)
Sidechain outliers	100360	3527 (1.70-1.70)
RSRZ outliers	91569	3200 (1.70-1.70)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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 $\leq 5\%$. 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	155	<div> <div>7%</div> <div>82% 10% 8%</div> </div>
1	B	155	<div> <div>5%</div> <div>74% 19% • 5%</div> </div>
1	C	155	<div> <div>5%</div> <div>79% 14% 7%</div> </div>
1	D	155	<div> <div>11%</div> <div>70% 16% • 12%</div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4931 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 hypothetical acetyltransferase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	143	Total	C	N	O	S	Se	0	0	0
			1125	724	191	203	2	5			
1	B	147	Total	C	N	O	S	Se	0	0	0
			1165	748	203	207	2	5			
1	C	144	Total	C	N	O	S	Se	0	0	0
			1134	730	193	204	2	5			
1	D	136	Total	C	N	O	S	Se	0	0	0
			1073	692	183	191	2	5			

There are 64 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-9	MSE	-	EXPRESSION TAG	UNP Q8E989
A	-8	GLY	-	EXPRESSION TAG	UNP Q8E989
A	-7	HIS	-	EXPRESSION TAG	UNP Q8E989
A	-6	HIS	-	EXPRESSION TAG	UNP Q8E989
A	-5	HIS	-	EXPRESSION TAG	UNP Q8E989
A	-4	HIS	-	EXPRESSION TAG	UNP Q8E989
A	-3	HIS	-	EXPRESSION TAG	UNP Q8E989
A	-2	HIS	-	EXPRESSION TAG	UNP Q8E989
A	-1	SER	-	EXPRESSION TAG	UNP Q8E989
A	0	HIS	-	EXPRESSION TAG	UNP Q8E989
A	1	MSE	MET	MODIFIED RESIDUE	UNP Q8E989
A	22	MSE	MET	MODIFIED RESIDUE	UNP Q8E989
A	50	MSE	MET	MODIFIED RESIDUE	UNP Q8E989
A	59	MSE	MET	MODIFIED RESIDUE	UNP Q8E989
A	66	MSE	MET	MODIFIED RESIDUE	UNP Q8E989
A	145	MSE	MET	MODIFIED RESIDUE	UNP Q8E989
B	-10	MSE	-	EXPRESSION TAG	UNP Q8E989
B	-9	GLY	-	EXPRESSION TAG	UNP Q8E989
B	-8	HIS	-	EXPRESSION TAG	UNP Q8E989
B	-7	HIS	-	EXPRESSION TAG	UNP Q8E989
B	-6	HIS	-	EXPRESSION TAG	UNP Q8E989

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
B	-5	HIS	-	EXPRESSION TAG	UNP Q8E989
B	-4	HIS	-	EXPRESSION TAG	UNP Q8E989
B	-3	HIS	-	EXPRESSION TAG	UNP Q8E989
B	-3A	SER	-	EXPRESSION TAG	UNP Q8E989
B	-2	HIS	-	EXPRESSION TAG	UNP Q8E989
B	1	MSE	MET	MODIFIED RESIDUE	UNP Q8E989
B	22	MSE	MET	MODIFIED RESIDUE	UNP Q8E989
B	50	MSE	MET	MODIFIED RESIDUE	UNP Q8E989
B	59	MSE	MET	MODIFIED RESIDUE	UNP Q8E989
B	66	MSE	MET	MODIFIED RESIDUE	UNP Q8E989
B	145	MSE	MET	MODIFIED RESIDUE	UNP Q8E989
C	-9	MSE	-	EXPRESSION TAG	UNP Q8E989
C	-8	GLY	-	EXPRESSION TAG	UNP Q8E989
C	-7	HIS	-	EXPRESSION TAG	UNP Q8E989
C	-6	HIS	-	EXPRESSION TAG	UNP Q8E989
C	-5	HIS	-	EXPRESSION TAG	UNP Q8E989
C	-4	HIS	-	EXPRESSION TAG	UNP Q8E989
C	-3	HIS	-	EXPRESSION TAG	UNP Q8E989
C	-2	HIS	-	EXPRESSION TAG	UNP Q8E989
C	-1	SER	-	EXPRESSION TAG	UNP Q8E989
C	0	HIS	-	EXPRESSION TAG	UNP Q8E989
C	1	MSE	MET	MODIFIED RESIDUE	UNP Q8E989
C	22	MSE	MET	MODIFIED RESIDUE	UNP Q8E989
C	50	MSE	MET	MODIFIED RESIDUE	UNP Q8E989
C	59	MSE	MET	MODIFIED RESIDUE	UNP Q8E989
C	66	MSE	MET	MODIFIED RESIDUE	UNP Q8E989
C	145	MSE	MET	MODIFIED RESIDUE	UNP Q8E989
D	-9	MSE	-	EXPRESSION TAG	UNP Q8E989
D	-8	GLY	-	EXPRESSION TAG	UNP Q8E989
D	-7	HIS	-	EXPRESSION TAG	UNP Q8E989
D	-6	HIS	-	EXPRESSION TAG	UNP Q8E989
D	-5	HIS	-	EXPRESSION TAG	UNP Q8E989
D	-4	HIS	-	EXPRESSION TAG	UNP Q8E989
D	-3	HIS	-	EXPRESSION TAG	UNP Q8E989
D	-2	HIS	-	EXPRESSION TAG	UNP Q8E989
D	-1	SER	-	EXPRESSION TAG	UNP Q8E989
D	0	HIS	-	EXPRESSION TAG	UNP Q8E989
D	1	MSE	MET	MODIFIED RESIDUE	UNP Q8E989
D	22	MSE	MET	MODIFIED RESIDUE	UNP Q8E989
D	50	MSE	MET	MODIFIED RESIDUE	UNP Q8E989
D	59	MSE	MET	MODIFIED RESIDUE	UNP Q8E989
D	66	MSE	MET	MODIFIED RESIDUE	UNP Q8E989

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
D	145	MSE	MET	MODIFIED RESIDUE	UNP Q8E989

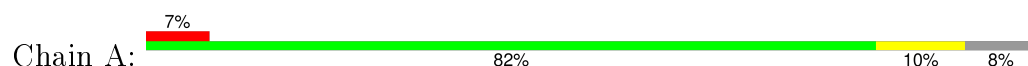
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	84	Total 84	O 84	0	0
2	B	112	Total 112	O 112	0	0
2	C	128	Total 128	O 128	0	0
2	D	110	Total 110	O 110	0	0

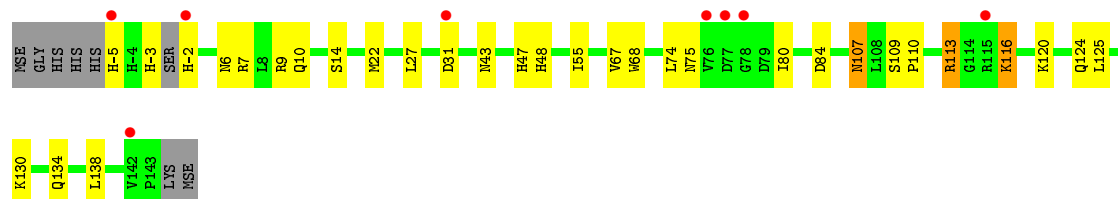
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

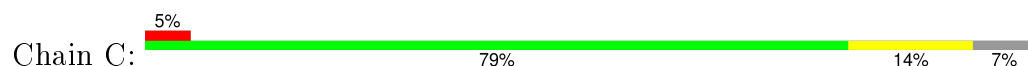
- Molecule 1: hypothetical acetyltransferase



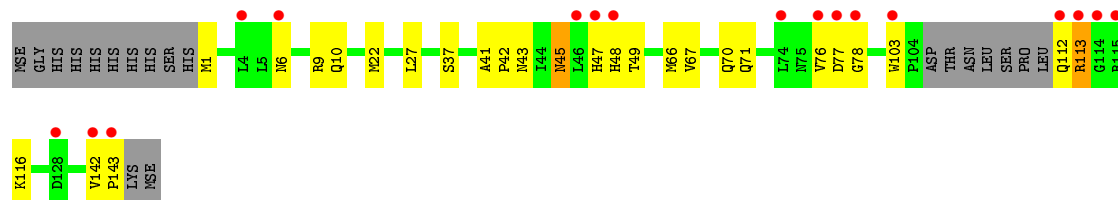
- Molecule 1: hypothetical acetyltransferase



- Molecule 1: hypothetical acetyltransferase



- Molecule 1: hypothetical acetyltransferase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	54.98Å 49.67Å 99.67Å 90.00° 91.09° 90.00°	Depositor
Resolution (Å)	29.81 – 1.70 29.81 – 1.69	Depositor EDS
% Data completeness (in resolution range)	92.9 (29.81-1.70) 96.8 (29.81-1.69)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	8.34 (at 1.70Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.194 , 0.222 0.202 , 0.228	Depositor DCC
R_{free} test set	5603 reflections (9.68%)	DCC
Wilson B-factor (Å ²)	12.0	Xtriage
Anisotropy	0.160	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 48.9	EDS
Estimated twinning fraction	0.014 for h,-k,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	1 of 112695 reflections (0.001%)	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	4931	wwPDB-VP
Average B, all atoms (Å ²)	15.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 15.56% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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	0.31	0/1149	0.55	0/1561
1	B	0.29	0/1192	0.54	0/1618
1	C	0.29	0/1157	0.54	0/1569
1	D	0.30	0/1095	0.54	0/1484
All	All	0.30	0/4593	0.54	0/6232

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	1125	0	1135	9	0
1	B	1165	0	1165	20	0
1	C	1134	0	1145	12	0
1	D	1073	0	1083	25	0
2	A	84	0	0	1	0
2	B	112	0	0	3	0
2	C	128	0	0	1	0
2	D	110	0	0	1	0
All	All	4931	0	4528	66	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 (66) 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:118:LYS:HG2	1:A:140:VAL:HG22	1.72	0.71
1:A:88:ARG:HH21	1:A:134:GLN:HG2	1.55	0.71
1:D:66:MSE:HE2	1:D:70:GLN:HG2	1.77	0.67
1:B:-5:HIS:HB2	1:B:6:ASN:HD21	1.59	0.66
1:D:45:ASN:HD22	1:D:45:ASN:C	1.98	0.64
1:C:41:ALA:HB3	1:C:42:PRO:HD3	1.80	0.63
1:B:107:ASN:ND2	1:B:109:SER:H	1.97	0.62
1:D:76:VAL:HG11	1:D:112:GLN:HB2	1.82	0.62
1:D:41:ALA:HB3	1:D:42:PRO:HD3	1.83	0.60
1:D:76:VAL:HG11	1:D:113:ARG:H	1.69	0.56
1:B:22:MSE:HG2	1:B:43:ASN:HD21	1.70	0.56
1:D:45:ASN:HD21	1:D:49:THR:H	1.53	0.56
1:D:67:VAL:O	1:D:71:GLN:HG3	2.06	0.54
1:C:40:LEU:HG	1:C:44:ILE:HG22	1.89	0.54
1:D:70:GLN:HG3	1:D:103:TRP:CE3	2.42	0.54
1:D:71:GLN:HB3	1:D:76:VAL:O	2.11	0.51
1:B:9:ARG:HH21	1:B:9:ARG:HG3	1.74	0.51
1:B:130:LYS:HE2	2:B:181:HOH:O	2.11	0.51
1:B:120:LYS:HG2	2:B:250:HOH:O	2.11	0.51
1:D:116:LYS:HG3	1:D:142:VAL:HG22	1.95	0.49
1:C:144:LYS:HB3	1:C:144:LYS:NZ	2.27	0.49
1:D:1:MSE:HB2	2:D:191:HOH:O	2.11	0.49
1:A:102:ARG:HD2	2:A:216:HOH:O	2.13	0.49
1:D:1:MSE:N	1:D:1:MSE:SE	2.96	0.49
1:D:6:ASN:ND2	1:D:9:ARG:HH22	2.10	0.48
1:B:47:HIS:O	1:B:48:HIS:HB2	2.13	0.48
1:C:109:SER:N	1:C:110:PRO:CD	2.77	0.47
1:C:68:TRP:O	1:C:72:GLN:HG2	2.14	0.47
1:A:27:LEU:HD11	1:A:37:SER:HB2	1.96	0.47
1:D:45:ASN:ND2	1:D:49:THR:H	2.13	0.47
1:C:3:GLU:HG2	2:C:232:HOH:O	2.15	0.46
1:A:9:ARG:HG2	1:A:9:ARG:HH11	1.80	0.46
1:C:55:ILE:HG23	1:C:125:LEU:HD13	1.98	0.46
1:B:113:ARG:HG2	1:B:113:ARG:HH21	1.80	0.45
1:A:109:SER:OG	1:A:110:PRO:HD3	2.16	0.45
1:A:70:GLN:HG3	1:A:103:TRP:CE3	2.52	0.45
1:B:10:GLN:NE2	1:B:14:SER:HB3	2.31	0.44
1:D:6:ASN:O	1:D:10:GLN:HG3	2.18	0.44
1:D:27:LEU:HD11	1:D:37:SER:HB2	1.99	0.44
1:A:55:ILE:O	1:A:58:ILE:HG22	2.18	0.44

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:45:ASN:C	1:D:45:ASN:ND2	2.69	0.44
1:D:78:GLY:HA2	1:D:143:PRO:HA	2.00	0.43
1:D:113:ARG:HA	1:D:113:ARG:NE	2.33	0.43
1:B:74:LEU:HD11	2:B:198:HOH:O	2.18	0.43
1:B:109:SER:OG	1:B:110:PRO:HD3	2.17	0.43
1:B:3:HIS:HB3	1:B:27:LEU:O	2.19	0.43
1:C:47:HIS:O	1:C:48:HIS:HB2	2.19	0.43
1:A:109:SER:N	1:A:110:PRO:CD	2.81	0.43
1:B:67:VAL:HB	1:B:80:ILE:HD13	2.00	0.43
1:C:70:GLN:HG3	1:C:103:TRP:CD2	2.53	0.43
1:B:109:SER:N	1:B:110:PRO:CD	2.81	0.43
1:D:76:VAL:HG21	1:D:112:GLN:CB	2.48	0.43
1:D:70:GLN:HG3	1:D:103:TRP:CD2	2.54	0.42
1:C:70:GLN:HG3	1:C:103:TRP:CE2	2.55	0.42
1:D:22:MSE:HG2	1:D:43:ASN:OD1	2.19	0.42
1:C:36:VAL:HG12	1:C:37:SER:N	2.35	0.42
1:B:7:ARG:HG2	1:B:68:TRP:HH2	1.84	0.42
1:B:55:ILE:HG23	1:B:125:LEU:HD13	2.01	0.41
1:B:124:GLN:HG2	1:B:134:GLN:HE22	1.86	0.41
1:D:47:HIS:O	1:D:48:HIS:HB2	2.20	0.41
1:B:116:LYS:HB3	1:B:116:LYS:NZ	2.36	0.41
1:D:76:VAL:HG21	1:D:112:GLN:HB2	2.02	0.41
1:B:107:ASN:HD21	1:B:109:SER:CB	2.34	0.40
1:D:76:VAL:HG21	1:D:112:GLN:HG3	2.02	0.40
1:C:104:PRO:HG2	1:C:106:THR:OG1	2.22	0.40
1:B:84:ASP:OD1	1:B:138:LEU:HB3	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	141/155 (91%)	139 (99%)	2 (1%)	0	100	100
1	B	143/155 (92%)	140 (98%)	3 (2%)	0	100	100
1	C	142/155 (92%)	138 (97%)	4 (3%)	0	100	100
1	D	132/155 (85%)	128 (97%)	4 (3%)	0	100	100
All	All	558/620 (90%)	545 (98%)	13 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

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	125/129 (97%)	123 (98%)	2 (2%)	70	54
1	B	129/129 (100%)	123 (95%)	6 (5%)	32	12
1	C	126/129 (98%)	125 (99%)	1 (1%)	86	79
1	D	118/129 (92%)	115 (98%)	3 (2%)	55	34
All	All	498/516 (96%)	486 (98%)	12 (2%)	57	36

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

Mol	Chain	Res	Type
1	A	1	MSE
1	A	115	ARG
1	B	-2	HIS
1	B	31	ASP
1	B	75	ASN
1	B	107	ASN
1	B	113	ARG
1	B	116	LYS
1	C	136	ASP
1	D	45	ASN
1	D	77	ASP
1	D	113	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	134	GLN
1	B	-3	HIS
1	B	6	ASN
1	B	10	GLN
1	B	43	ASN
1	B	71	GLN
1	B	107	ASN
1	B	124	GLN
1	B	134	GLN
1	C	6	ASN
1	C	43	ASN
1	C	71	GLN
1	C	134	GLN
1	D	6	ASN
1	D	45	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 carbohydrates 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

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	138/155 (89%)	0.51	11 (7%) 15 17	4, 11, 26, 32	0
1	B	142/155 (91%)	0.51	8 (5%) 28 30	5, 12, 24, 30	0
1	C	139/155 (89%)	0.43	7 (5%) 32 35	5, 14, 24, 30	0
1	D	131/155 (84%)	0.80	17 (12%) 5 5	5, 14, 28, 31	0
All	All	550/620 (88%)	0.56	43 (7%) 16 18	4, 13, 26, 32	0

All (43) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	-2	HIS	7.5
1	D	113	ARG	7.4
1	D	77	ASP	5.5
1	A	113	ARG	4.9
1	A	115	ARG	4.6
1	B	77	ASP	4.5
1	A	114	GLY	4.2
1	D	76	VAL	4.1
1	D	115	ARG	3.9
1	D	143	PRO	3.9
1	C	105	ASP	3.7
1	D	48	HIS	3.7
1	C	107	ASN	3.7
1	D	142	VAL	3.6
1	D	46	LEU	3.4
1	D	114	GLY	3.3
1	B	-5	HIS	3.0
1	C	106	THR	3.0
1	A	105	ASP	2.9
1	D	47	HIS	2.9
1	D	128	ASP	2.8

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	B	142	VAL	2.8
1	D	112	GLN	2.8
1	C	2	ASP	2.8
1	A	106	THR	2.7
1	C	136	ASP	2.6
1	D	74	LEU	2.6
1	A	2	ASP	2.6
1	D	78	GLY	2.6
1	A	88	ARG	2.5
1	B	115	ARG	2.5
1	A	3	GLU	2.5
1	C	144	LYS	2.4
1	B	78	GLY	2.3
1	B	31	ASP	2.3
1	A	143	PRO	2.2
1	C	86	HIS	2.2
1	D	4	LEU	2.2
1	D	103	TRP	2.1
1	B	76	VAL	2.1
1	D	6	ASN	2.1
1	A	111	LEU	2.0
1	A	10	GLN	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 carbohydrates 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.