



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 17, 2026 – 06:32 PM UTC

PDB ID : 6A8Q / pdb_00006a8q
Title : R207A mutant of highly active EfBSH
Authors : Ramasamy, S.; Yadav, Y.
Deposited on : 2018-07-10
Resolution : 2.00 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

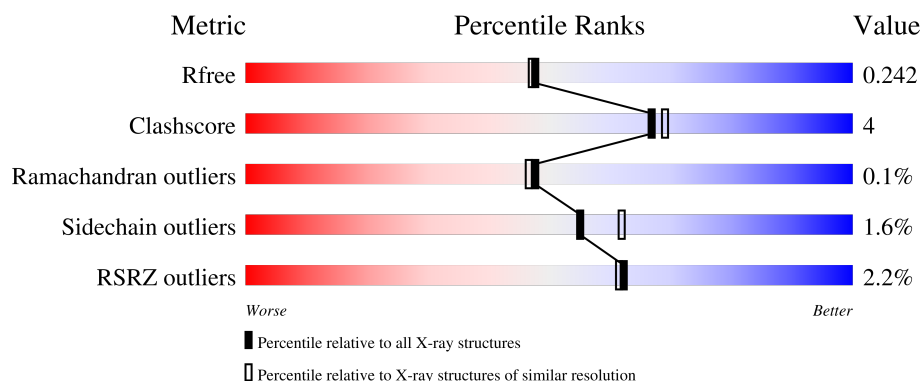
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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 $\leq 5\%$. 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	331	0% 92% 5%
1	B	331	2% 93% 5%
1	C	331	4% 89% 7%
1	D	331	2% 92% 5%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 10776 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 Bile salt hydrolase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	323	Total	C	N	O	S	0	2	0
			2600	1662	418	511	9			
1	B	323	Total	C	N	O	S	0	2	0
			2599	1662	417	511	9			
1	C	323	Total	C	N	O	S	0	2	0
			2599	1661	417	511	10			
1	D	323	Total	C	N	O	S	0	2	0
			2601	1663	417	512	9			

There are 36 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	207	ALA	ARG	engineered mutation	UNP C7CXJ5
A	325	LEU	-	expression tag	UNP C7CXJ5
A	326	GLU	-	expression tag	UNP C7CXJ5
A	327	HIS	-	expression tag	UNP C7CXJ5
A	328	HIS	-	expression tag	UNP C7CXJ5
A	329	HIS	-	expression tag	UNP C7CXJ5
A	330	HIS	-	expression tag	UNP C7CXJ5
A	331	HIS	-	expression tag	UNP C7CXJ5
A	332	HIS	-	expression tag	UNP C7CXJ5
B	207	ALA	ARG	engineered mutation	UNP C7CXJ5
B	325	LEU	-	expression tag	UNP C7CXJ5
B	326	GLU	-	expression tag	UNP C7CXJ5
B	327	HIS	-	expression tag	UNP C7CXJ5
B	328	HIS	-	expression tag	UNP C7CXJ5
B	329	HIS	-	expression tag	UNP C7CXJ5
B	330	HIS	-	expression tag	UNP C7CXJ5
B	331	HIS	-	expression tag	UNP C7CXJ5
B	332	HIS	-	expression tag	UNP C7CXJ5
C	207	ALA	ARG	engineered mutation	UNP C7CXJ5
C	325	LEU	-	expression tag	UNP C7CXJ5
C	326	GLU	-	expression tag	UNP C7CXJ5

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
C	327	HIS	-	expression tag	UNP C7CXJ5
C	328	HIS	-	expression tag	UNP C7CXJ5
C	329	HIS	-	expression tag	UNP C7CXJ5
C	330	HIS	-	expression tag	UNP C7CXJ5
C	331	HIS	-	expression tag	UNP C7CXJ5
C	332	HIS	-	expression tag	UNP C7CXJ5
D	207	ALA	ARG	engineered mutation	UNP C7CXJ5
D	325	LEU	-	expression tag	UNP C7CXJ5
D	326	GLU	-	expression tag	UNP C7CXJ5
D	327	HIS	-	expression tag	UNP C7CXJ5
D	328	HIS	-	expression tag	UNP C7CXJ5
D	329	HIS	-	expression tag	UNP C7CXJ5
D	330	HIS	-	expression tag	UNP C7CXJ5
D	331	HIS	-	expression tag	UNP C7CXJ5
D	332	HIS	-	expression tag	UNP C7CXJ5

- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	95	Total O 95 95	0	0
2	B	87	Total O 87 87	0	0
2	C	90	Total O 90 90	0	0
2	D	105	Total O 105 105	0	0

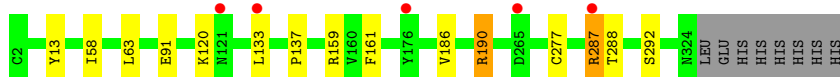
3 Residue-property plots [i](#)

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.

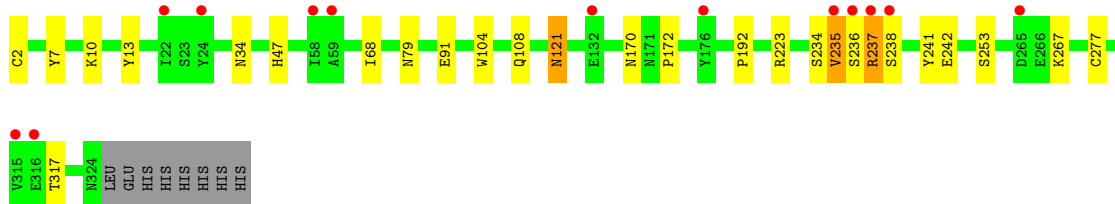
- Molecule 1: Bile salt hydrolase



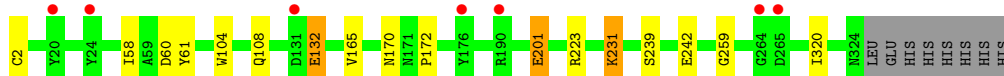
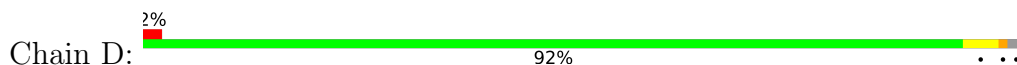
- Molecule 1: Bile salt hydrolase



- Molecule 1: Bile salt hydrolase



- Molecule 1: Bile salt hydrolase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	59.33Å 156.45Å 72.91Å 90.00° 98.62° 90.00°	Depositor
Resolution (Å)	39.01 – 2.00 39.01 – 2.00	Depositor EDS
% Data completeness (in resolution range)	94.1 (39.01-2.00) 94.1 (39.01-2.00)	Depositor EDS
R_{merge}	0.20	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.20 (at 2.00Å)	Xtriage
Refinement program	REFMAC 5.8.0222	Depositor
R, R_{free}	0.201 , 0.232 0.214 , 0.242	Depositor DCC
R_{free} test set	4095 reflections (4.63%)	wwPDB-VP
Wilson B-factor (Å ²)	13.4	Xtriage
Anisotropy	0.132	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 37.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	10776	wwPDB-VP
Average B, all atoms (Å ²)	16.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.26% 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, 0.333 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	0.90	2/2665 (0.1%)	0.96	2/3613 (0.1%)
1	B	0.89	1/2664 (0.0%)	0.96	1/3613 (0.0%)
1	C	0.85	0/2664	0.95	2/3612 (0.1%)
1	D	0.85	1/2666 (0.0%)	0.95	3/3616 (0.1%)
All	All	0.87	4/10659 (0.0%)	0.95	8/14454 (0.1%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	175	ASP	C-O	-8.58	1.14	1.24
1	D	60	ASP	C-O	-7.46	1.14	1.24
1	B	288	THR	C-N	-6.40	1.25	1.33
1	A	120	LYS	C-O	-5.17	1.17	1.24

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	121	ASN	CA-CB-CG	-7.88	104.72	112.60
1	D	58	ILE	O-C-N	7.07	129.79	122.09
1	C	317	THR	CA-CB-OG1	-6.63	99.65	109.60
1	D	201	GLU	CB-CG-CD	6.01	122.81	112.60
1	A	265	ASP	N-CA-C	5.85	119.26	111.30
1	A	201	GLU	CB-CG-CD	5.20	121.44	112.60
1	B	137	PRO	CB-CA-C	5.12	116.06	111.40
1	D	61	TYR	N-CA-C	5.12	117.37	110.31

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	2600	0	2514	8	0
1	B	2599	0	2512	16	0
1	C	2599	0	2510	43	0
1	D	2601	0	2511	14	0
2	A	95	0	0	2	0
2	B	87	0	0	0	0
2	C	90	0	0	1	0
2	D	105	0	0	1	0
All	All	10776	0	10047	75	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 4.

All (75) 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:C:237:ARG:HD2	1:C:242:GLU:CG	1.51	1.39
1:C:237:ARG:CD	1:C:242:GLU:HB3	1.70	1.22
1:C:237:ARG:HD2	1:C:242:GLU:CB	1.76	1.16
1:C:237:ARG:HD2	1:C:242:GLU:HG2	1.12	1.07
1:C:237:ARG:CG	1:C:242:GLU:CB	2.33	1.07
1:C:237:ARG:CD	1:C:242:GLU:CG	2.33	1.07
1:C:237:ARG:CD	1:C:242:GLU:CB	2.31	1.05
1:C:237:ARG:HG3	1:C:242:GLU:OE1	1.58	1.03
1:C:237:ARG:HG3	1:C:242:GLU:CD	1.83	1.02
1:B:190:ARG:HH11	1:B:190:ARG:HG3	1.26	1.01
1:B:287:ARG:HD3	1:B:292:SER:O	1.68	0.94
1:C:236:SER:HA	1:C:237:ARG:HH11	1.36	0.90
1:C:237:ARG:CG	1:C:242:GLU:HB3	1.97	0.90
1:D:170:ASN:HD22	1:D:223:ARG:HH22	1.24	0.85
1:C:237:ARG:HD2	1:C:242:GLU:HB3	1.43	0.84
1:C:170:ASN:HD22	1:C:223:ARG:HH22	1.26	0.82
1:C:237:ARG:CD	1:C:242:GLU:HG2	2.03	0.81
1:D:165:VAL:O	1:D:231:LYS:HE3	1.83	0.79
1:C:237:ARG:HG2	1:C:242:GLU:CB	2.13	0.78

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:237:ARG:CG	1:C:242:GLU:HB2	2.13	0.78
1:C:237:ARG:O	1:C:238:SER:HB3	1.84	0.76
1:C:170:ASN:ND2	1:C:223:ARG:HH22	1.85	0.75
1:C:237:ARG:CG	1:C:242:GLU:CD	2.58	0.75
1:D:170:ASN:ND2	1:D:223:ARG:HH22	1.84	0.73
1:C:237:ARG:NE	1:C:242:GLU:HB3	2.06	0.68
1:C:237:ARG:HG3	1:C:242:GLU:CB	2.24	0.68
1:C:237:ARG:CG	1:C:242:GLU:CG	2.70	0.67
1:A:120:LYS:HA	1:A:120:LYS:HE2	1.80	0.64
1:C:236:SER:C	1:C:237:ARG:HE	2.06	0.64
1:C:237:ARG:CD	1:C:242:GLU:CD	2.71	0.63
1:C:237:ARG:HG2	1:C:242:GLU:HB3	1.73	0.63
1:B:287:ARG:HD3	1:B:292:SER:C	2.24	0.62
1:D:165:VAL:O	1:D:231:LYS:CE	2.46	0.62
1:C:237:ARG:HG2	1:C:242:GLU:HB2	1.78	0.62
1:C:79:ASN:ND2	2:C:402:HOH:O	2.34	0.60
1:C:237:ARG:HG3	1:C:242:GLU:CG	2.31	0.60
1:B:287:ARG:HD3	1:B:292:SER:HA	1.85	0.58
1:B:287:ARG:CD	1:B:292:SER:O	2.49	0.58
1:B:190:ARG:HG3	1:B:190:ARG:NH1	2.07	0.56
1:B:287:ARG:CD	1:B:292:SER:HA	2.35	0.56
1:B:58:ILE:C	1:B:58:ILE:HD12	2.31	0.56
1:C:2:CYS:N	1:C:170:ASN:HD21	2.05	0.55
1:B:190:ARG:HH11	1:B:190:ARG:CG	2.05	0.55
1:D:2:CYS:N	1:D:170:ASN:HD21	2.05	0.54
1:C:235:VAL:HG22	1:D:259:GLY:HA3	1.89	0.53
1:C:235:VAL:HG22	1:D:259:GLY:CA	2.39	0.52
1:C:237:ARG:HG3	1:C:242:GLU:HB2	1.87	0.51
1:A:104:TRP:CZ2	1:A:118:LEU:HD11	2.45	0.51
1:B:287:ARG:HD3	1:B:292:SER:CA	2.41	0.51
1:C:34:ASN:ND2	1:C:47:HIS:ND1	2.58	0.51
1:A:49:PHE:CE1	1:A:71:LYS:HE3	2.47	0.49
1:B:186:VAL:HG23	1:C:192:PRO:HB3	1.96	0.47
1:A:235:VAL:HG23	2:A:462:HOH:O	2.15	0.46
1:A:104:TRP:CH2	1:A:118:LEU:HG	2.50	0.46
1:A:239:SER:OG	1:A:242:GLU:OE1	2.28	0.46
1:A:79:ASN:ND2	2:A:402:HOH:O	2.48	0.45
1:C:253:SER:HA	2:D:476:HOH:O	2.16	0.45
1:C:241:TYR:CE1	1:D:320:ILE:HD12	2.52	0.44
1:D:201:GLU:O	1:D:201:GLU:HG2	2.17	0.44
1:B:159:ARG:HH11	1:B:161:PHE:HE2	1.65	0.44

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:120:LYS:HA	1:B:120:LYS:CE	2.47	0.43
1:B:63:LEU:HD13	1:B:133:LEU:HD21	1.99	0.43
1:D:132:GLU:H	1:D:132:GLU:HG2	1.64	0.43
1:C:7:TYR:OH	1:C:236:SER:HB3	2.18	0.43
1:C:236:SER:HA	1:C:237:ARG:NH1	2.18	0.43
1:D:239:SER:OG	1:D:242:GLU:OE1	2.31	0.42
1:D:104:TRP:O	1:D:108:GLN:HG2	2.19	0.42
1:C:241:TYR:CE1	1:D:320:ILE:CD1	3.03	0.41
1:A:13:TYR:HA	1:A:277:CYS:O	2.20	0.41
1:C:13:TYR:HA	1:C:277:CYS:O	2.21	0.41
1:C:68:ILE:HD13	1:C:277:CYS:HB3	2.03	0.41
1:B:58:ILE:HD11	1:B:133:LEU:HD22	2.04	0.40
1:B:13:TYR:HA	1:B:277:CYS:O	2.21	0.40
1:C:104:TRP:O	1:C:108:GLN:HG2	2.21	0.40
1:C:241:TYR:CZ	1:D:320:ILE:HD12	2.57	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	323/331 (98%)	309 (96%)	14 (4%)	0	100	100
1	B	323/331 (98%)	313 (97%)	10 (3%)	0	100	100
1	C	323/331 (98%)	310 (96%)	12 (4%)	1 (0%)	36	35
1	D	323/331 (98%)	313 (97%)	10 (3%)	0	100	100
All	All	1292/1324 (98%)	1245 (96%)	46 (4%)	1 (0%)	48	46

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	10	LYS

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	291/297 (98%)	286 (98%)	5 (2%)	53	60
1	B	291/297 (98%)	288 (99%)	3 (1%)	68	75
1	C	291/297 (98%)	284 (98%)	7 (2%)	43	47
1	D	291/297 (98%)	288 (99%)	3 (1%)	68	75
All	All	1164/1188 (98%)	1146 (98%)	18 (2%)	55	64

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	91	GLU
1	A	120	LYS
1	A	201	GLU
1	A	267	LYS
1	A	308	SER
1	B	91	GLU
1	B	190	ARG
1	B	287	ARG
1	C	91	GLU
1	C	121	ASN
1	C	172	PRO
1	C	234	SER
1	C	235	VAL
1	C	237	ARG
1	C	267	LYS
1	D	132	GLU
1	D	172	PRO
1	D	231	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	79	ASN
1	A	171	ASN
1	A	180	ASN
1	A	183	ASN
1	A	300	ASN
1	B	79	ASN
1	B	171	ASN
1	B	293	GLN
1	B	300	ASN
1	C	34	ASN
1	C	79	ASN
1	C	170	ASN
1	C	171	ASN
1	C	183	ASN
1	C	199	GLN
1	D	170	ASN
1	D	171	ASN
1	D	180	ASN
1	D	183	ASN
1	D	199	GLN
1	D	293	GLN
1	D	300	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 oligosaccharides 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

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	323/331 (97%)	-0.08	4 (1%) 76 76	5, 13, 30, 43	3 (0%)
1	B	323/331 (97%)	0.01	5 (1%) 72 71	5, 14, 30, 49	3 (0%)
1	C	323/331 (97%)	0.09	13 (4%) 42 41	4, 14, 36, 69	3 (0%)
1	D	323/331 (97%)	-0.09	7 (2%) 62 61	6, 13, 30, 52	3 (0%)
All	All	1292/1324 (97%)	-0.02	29 (2%) 62 61	4, 14, 32, 69	12 (0%)

All (29) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	236	SER	8.0
1	C	237	ARG	7.2
1	C	24	TYR	6.5
1	D	176	TYR	6.4
1	B	176	TYR	6.2
1	C	176	TYR	6.2
1	A	176	TYR	5.9
1	C	315	VAL	4.7
1	C	238	SER	4.4
1	B	287	ARG	4.1
1	C	58	ILE	3.9
1	B	121	ASN	3.3
1	B	265	ASP	3.2
1	C	22	ILE	2.9
1	C	59	ALA	2.9
1	C	316	GLU	2.8
1	D	131	ASP	2.8
1	D	20	TYR	2.7
1	C	235	VAL	2.5
1	B	133	LEU	2.5
1	C	132	GLU	2.4

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	D	190	ARG	2.4
1	D	265	ASP	2.4
1	A	58	ILE	2.3
1	D	24	TYR	2.3
1	A	22	ILE	2.2
1	C	265	ASP	2.2
1	D	264	GLY	2.1
1	A	132	GLU	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 oligosaccharides 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.