



Full wwPDB X-ray Structure Validation Report ⓘ

May 26, 2020 – 11:52 pm BST

PDB ID : 5NWP
Title : Crystal Structure of the Lectin Domain From the F17-like Adhesin, UclD
Authors : Ruer, S.; Remaut, H.
Deposited on : 2017-05-08
Resolution : 1.05 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 following versions of software and data (see [references ⓘ](#)) 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 : 2.11
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) : 2.11

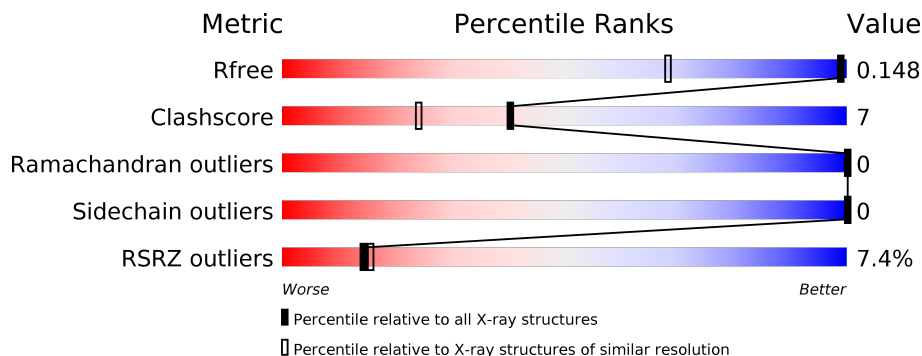
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.05 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	1202 (1.10-1.02)
Clashscore	141614	1252 (1.10-1.02)
Ramachandran outliers	138981	1204 (1.10-1.02)
Sidechain outliers	138945	1202 (1.10-1.02)
RSRZ outliers	127900	1178 (1.10-1.02)

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 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	203	
1	B	203	

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 6888 atoms, of which 3135 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 Adhesin.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S			
1	A	197	3115	986	1572	263	291	3	48	16	0
1	B	197	3106	985	1563	261	294	3	46	17	0

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	198	HIS	-	expression tag	UNP Q8GA71
A	199	HIS	-	expression tag	UNP Q8GA71
A	200	HIS	-	expression tag	UNP Q8GA71
A	201	HIS	-	expression tag	UNP Q8GA71
A	202	HIS	-	expression tag	UNP Q8GA71
A	203	HIS	-	expression tag	UNP Q8GA71
B	198	HIS	-	expression tag	UNP Q8GA71
B	199	HIS	-	expression tag	UNP Q8GA71
B	200	HIS	-	expression tag	UNP Q8GA71
B	201	HIS	-	expression tag	UNP Q8GA71
B	202	HIS	-	expression tag	UNP Q8GA71
B	203	HIS	-	expression tag	UNP Q8GA71

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		

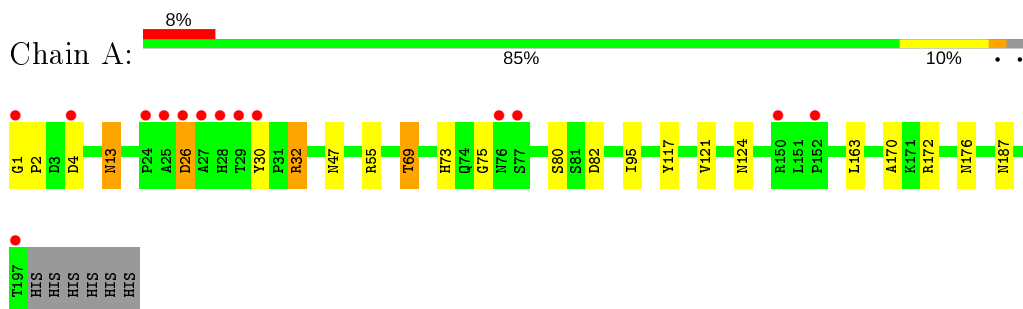
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	315	Total	O	0	5
			320	320		
3	B	331	Total	O	0	11
			342	342		

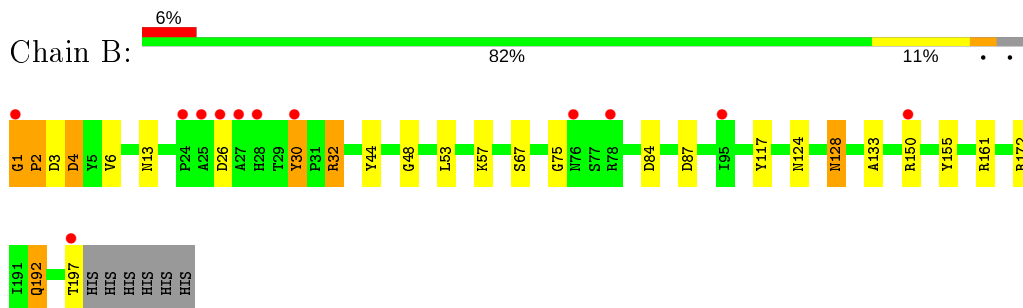
3 Residue-property plots [i](#)

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



- Molecule 1: Adhesin



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	30.09Å 91.87Å 64.93Å 90.00° 96.46° 90.00°	Depositor
Resolution (Å)	91.87 – 1.05 37.42 – 1.05	Depositor EDS
% Data completeness (in resolution range)	88.7 (91.87-1.05) 88.7 (37.42-1.05)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.76 (at 1.05Å)	Xtrriage
Refinement program	REFMAC 5.8.0124	Depositor
R, R_{free}	0.117 , 0.143 0.125 , 0.148	Depositor DCC
R_{free} test set	7242 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	7.7	Xtrriage
Anisotropy	0.096	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 50.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.98	EDS
Total number of atoms	6888	wwPDB-VP
Average B, all atoms (Å ²)	11.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.26% 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

Bond lengths and bond angles in the following residue types are not validated in this section: SO4

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	1.21	5/1614 (0.3%)	1.26	14/2204 (0.6%)
1	B	1.38	16/1621 (1.0%)	1.23	17/2213 (0.8%)
All	All	1.30	21/3235 (0.6%)	1.24	31/4417 (0.7%)

All (21) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	1	GLY	C-O	8.40	1.37	1.23
1	B	172	ARG	CA-CB	8.36	1.72	1.53
1	B	189	GLY	C-O	-7.51	1.11	1.23
1	B	192	GLN	CB-CG	-7.50	1.32	1.52
1	B	32[A]	ARG	CZ-NH1	-7.36	1.23	1.33
1	B	32[B]	ARG	CZ-NH1	-7.36	1.23	1.33
1	A	172	ARG	CA-CB	7.30	1.70	1.53
1	B	150	ARG	CZ-NH1	6.31	1.41	1.33
1	B	67[A]	SER	CA-CB	6.22	1.62	1.52
1	B	67[B]	SER	CA-CB	6.22	1.62	1.52
1	B	6	VAL	CA-CB	-6.07	1.42	1.54
1	B	2	PRO	N-CA	-6.01	1.37	1.47
1	B	44	TYR	CD1-CE1	5.79	1.48	1.39
1	B	128[A]	ASN	CB-CG	5.74	1.64	1.51
1	B	128[B]	ASN	CB-CG	5.74	1.64	1.51
1	B	6	VAL	CB-CG2	-5.62	1.41	1.52
1	A	13	ASN	CG-ND2	-5.37	1.19	1.32
1	B	44	TYR	CB-CG	-5.36	1.43	1.51
1	A	69[A]	THR	CB-CG2	5.32	1.70	1.52
1	A	69[B]	THR	CB-CG2	5.32	1.70	1.52
1	A	47	ASN	CG-OD1	5.00	1.34	1.24

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	190[A]	ARG	NE-CZ-NH2	-13.09	113.75	120.30
1	A	190[B]	ARG	NE-CZ-NH2	-13.09	113.75	120.30
1	B	44	TYR	CD1-CE1-CZ	-9.96	110.84	119.80
1	A	190[A]	ARG	NH1-CZ-NH2	8.95	129.25	119.40
1	A	190[B]	ARG	NH1-CZ-NH2	8.95	129.25	119.40
1	A	172	ARG	NE-CZ-NH1	8.88	124.74	120.30
1	B	4[A]	ASP	CB-CG-OD1	8.42	125.88	118.30
1	B	4[B]	ASP	CB-CG-OD1	8.42	125.88	118.30
1	A	82	ASP	CB-CG-OD2	7.90	125.41	118.30
1	B	30	TYR	CB-CG-CD1	7.49	125.50	121.00
1	B	161	ARG	NH1-CZ-NH2	-7.38	111.28	119.40
1	A	30	TYR	CB-CG-CD1	-7.25	116.65	121.00
1	B	3	ASP	CB-CG-OD2	-7.22	111.80	118.30
1	A	55	ARG	NE-CZ-NH2	-6.72	116.94	120.30
1	B	190	ARG	NE-CZ-NH2	-6.70	116.95	120.30
1	A	190[A]	ARG	NE-CZ-NH1	-6.68	116.96	120.30
1	A	190[B]	ARG	NE-CZ-NH1	-6.68	116.96	120.30
1	B	161	ARG	NE-CZ-NH1	6.59	123.59	120.30
1	B	150	ARG	NE-CZ-NH1	6.58	123.59	120.30
1	A	26	ASP	CB-CG-OD1	-6.37	112.56	118.30
1	B	197	THR	CA-C-O	-6.24	107.00	120.10
1	B	87	ASP	CB-CG-OD2	-6.11	112.80	118.30
1	B	161	ARG	NE-CZ-NH2	6.04	123.32	120.30
1	B	44	TYR	CZ-CE2-CD2	-6.02	114.38	119.80
1	A	172	ARG	NE-CZ-NH2	-5.84	117.38	120.30
1	B	84	ASP	CB-CG-OD1	5.77	123.50	118.30
1	A	32[A]	ARG	NE-CZ-NH1	5.09	122.85	120.30
1	A	32[B]	ARG	NE-CZ-NH1	5.09	122.85	120.30
1	B	172	ARG	NE-CZ-NH2	-5.06	117.77	120.30
1	B	48	GLY	CA-C-N	5.04	126.29	116.20
1	B	3	ASP	CB-CG-OD1	5.03	122.83	118.30

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	1543	1572	1577	24	2
1	B	1543	1563	1570	21	2
2	A	5	0	0	0	0
3	A	320	0	0	13	0
3	B	342	0	0	7	6
All	All	3753	3135	3147	45	6

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 (45) 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:117:TYR:H	1:A:124:ASN:HD21	1.18	0.86
1:B:117:TYR:H	1:B:124:ASN:HD21	1.18	0.86
1:B:26[A]:ASP:CG	3:B:303:HOH:O	2.19	0.80
1:B:26[A]:ASP:OD2	3:B:301:HOH:O	1.99	0.79
1:B:32[B]:ARG:HD2	3:B:545:HOH:O	1.84	0.77
1:A:1:GLY:N	1:A:2:PRO:CD	2.48	0.76
1:A:170:ALA:H	1:A:176:ASN:HD22	1.31	0.76
1:A:1:GLY:H3	1:A:2:PRO:CD	2.01	0.72
1:A:1:GLY:O	3:A:402:HOH:O	2.06	0.72
1:A:4[B]:ASP:OD1	3:A:402:HOH:O	2.10	0.70
1:A:69[B]:THR:HG22	1:A:80[B]:SER:OG	1.92	0.69
1:A:1:GLY:N	1:A:2:PRO:HD2	2.12	0.64
1:B:26[A]:ASP:OD1	3:B:303:HOH:O	2.13	0.63
1:A:32[A]:ARG:CD	3:A:401:HOH:O	2.48	0.60
1:B:53:LEU:HD23	1:B:133:ALA:N	2.20	0.56
1:B:13[A]:ASN:HD21	1:B:187:ASN:HD22	1.52	0.56
1:A:1:GLY:H2	1:A:2:PRO:HD2	1.71	0.56
1:B:1:GLY:O	1:B:4[B]:ASP:OD1	2.24	0.55
1:B:13[A]:ASN:ND2	1:B:187:ASN:HD22	2.07	0.53
1:A:1:GLY:H3	1:A:2:PRO:HD3	1.73	0.52
1:B:53:LEU:HD23	1:B:133:ALA:CA	2.40	0.52
1:A:121[B]:VAL:HG22	3:A:617:HOH:O	2.10	0.52
1:A:26:ASP:HB2	3:A:407:HOH:O	2.11	0.49
1:B:128[B]:ASN:OD1	3:B:304:HOH:O	2.19	0.49
1:B:32[B]:ARG:HD3	3:B:562:HOH:O	2.12	0.49
1:A:32[A]:ARG:HG2	3:A:401:HOH:O	2.12	0.49
1:B:53:LEU:CD2	1:B:133:ALA:N	2.76	0.49
1:A:170:ALA:HB1	3:A:532:HOH:O	2.13	0.48
1:A:73[A]:HIS:CD2	3:A:522:HOH:O	2.66	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1:GLY:N	1:B:2:PRO:HD2	2.28	0.47
1:A:26:ASP:CB	3:A:407:HOH:O	2.64	0.46
1:A:75:GLY:HA3	3:A:417:HOH:O	2.15	0.46
1:B:117:TYR:H	1:B:124:ASN:ND2	1.99	0.46
1:B:155:TYR:CD1	1:B:192:GLN:HG2	2.52	0.45
3:A:482[B]:HOH:O	1:B:57:LYS:NZ	2.49	0.45
1:A:13:ASN:HD22	1:A:187:ASN:HB2	1.82	0.44
1:A:69[B]:THR:HG22	1:A:80[B]:SER:HG	1.78	0.44
1:B:1:GLY:O	1:B:4[B]:ASP:CG	2.57	0.43
1:B:155:TYR:CE1	1:B:192:GLN:HG2	2.54	0.42
1:A:32[A]:ARG:CG	3:A:401:HOH:O	2.67	0.42
1:A:95[B]:ILE:HD11	1:A:163:LEU:HD22	2.00	0.42
1:B:75:GLY:C	3:B:308[B]:HOH:O	2.57	0.42
1:A:95[B]:ILE:HD11	1:A:163:LEU:CD2	2.51	0.41
1:A:1:GLY:HA2	3:A:403:HOH:O	2.20	0.40
1:B:1:GLY:N	1:B:2:PRO:CD	2.84	0.40

All (6) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:196:GLN:OE1	3:B:626:HOH:O[2_556]	0.46	1.74
1:B:30:TYR:O	3:B:623:HOH:O[1_655]	0.50	1.70
1:B:30:TYR:C	3:B:623:HOH:O[1_655]	1.26	0.94
1:A:196:GLN:CD	3:B:626:HOH:O[2_556]	1.42	0.78
3:B:486:HOH:O	3:B:550:HOH:O[1_455]	1.88	0.32
3:B:428:HOH:O	3:B:573:HOH:O[1_455]	1.89	0.31

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	211/203 (104%)	207 (98%)	4 (2%)	0	100	100
1	B	212/203 (104%)	207 (98%)	5 (2%)	0	100	100
All	All	423/406 (104%)	414 (98%)	9 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	171/161 (106%)	171 (100%)	0	100	100
1	B	172/161 (107%)	172 (100%)	0	100	100
All	All	343/322 (106%)	343 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	13	ASN
1	A	47	ASN
1	A	124	ASN
1	A	176	ASN
1	A	187	ASN
1	B	124	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](#)

1 ligand is modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	SO4	A	301	-	4,4,4	0.86	0	6,6,6	2.35	2 (33%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
2	A	301	SO4	O4-S-O3	4.37	127.72	109.06
2	A	301	SO4	O4-S-O1	-3.66	90.20	109.31

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	A	197/203 (97%)	1.11	17 (8%) 10 13	5, 8, 28, 54	0
1	B	197/203 (97%)	1.02	12 (6%) 21 20	5, 8, 19, 43	0
All	All	394/406 (97%)	1.06	29 (7%) 14 16	5, 8, 25, 54	0

All (29) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	27	ALA	12.6
1	A	25	ALA	9.2
1	B	25	ALA	5.6
1	A	26	ASP	5.2
1	B	28	HIS	5.0
1	A	28	HIS	4.9
1	B	78	ARG	4.8
1	B	76	ASN	4.5
1	A	24	PRO	4.4
1	A	76	ASN	4.4
1	A	150	ARG	3.5
1	B	27	ALA	3.5
1	A	30	TYR	3.5
1	B	1	GLY	3.2
1	B	26[A]	ASP	3.2
1	A	1	GLY	3.1
1	A	77	SER	2.9
1	A	29[A]	THR	2.8
1	B	95[A]	ILE	2.8
1	A	190[A]	ARG	2.6
1	A	152	PRO	2.5
1	B	197	THR	2.5
1	B	150	ARG	2.4
1	A	193	VAL	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	196	GLN	2.3
1	B	24	PRO	2.3
1	B	30	TYR	2.2
1	A	4[A]	ASP	2.1
1	A	197	THR	2.1

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](#)

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
2	SO4	A	301	5/5	0.94	0.20	6,10,12,14	5

6.5 Other polymers [i](#)

There are no such residues in this entry.