



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 06:38 pm GMT

PDB ID : 2ADF
Title : Crystal Structure and Paratope Determination of 82D6A3, an Antithrombotic Antibody Directed Against the von Willebrand factor A3-Domain
Authors : Staelens, S.; Hadders, M.A.; Vauterin, S.; Platteau, C.; Vanhoorelbeke, K.; Huizinga, E.G.; Deckmyn, H.
Deposited on : 2005-07-20
Resolution : 1.90 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

<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	:	1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix)	:	1.9-1692
EDS	:	trunk28620
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
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)	:	recalc28949

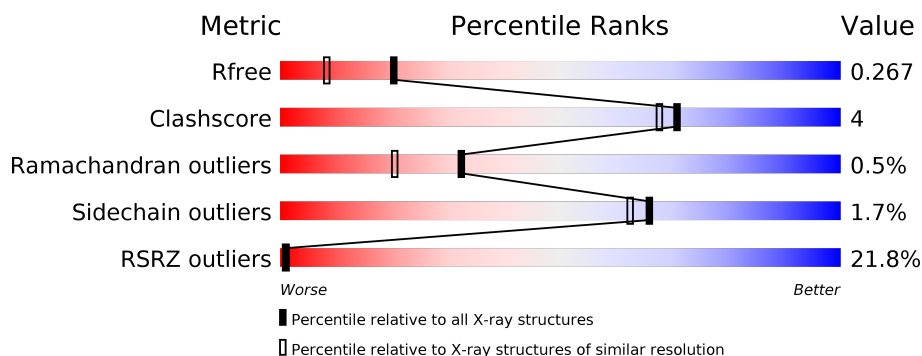
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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}	100719	5047 (1.90-1.90)
Clashscore	112137	5731 (1.90-1.90)
Ramachandran outliers	110173	5669 (1.90-1.90)
Sidechain outliers	110143	5670 (1.90-1.90)
RSRZ outliers	101464	5100 (1.90-1.90)

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 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	196	
2	H	218	
3	L	209	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	SO4	L	1003	-	-	-	X
6	GOL	H	1004	-	-	-	X

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 5165 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 Von Willebrand factor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	189	Total	C	N	O	S	0	0	0
			1414	894	242	272	6			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	916	GLY	-	EXPRESSION TAG	UNP P04275
A	917	SER	-	EXPRESSION TAG	UNP P04275
A	918	HIS	-	EXPRESSION TAG	UNP P04275
A	919	MET	-	EXPRESSION TAG	UNP P04275

- Molecule 2 is a protein called 82D6A3 IgG.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	218	Total	C	N	O	S	0	0	0
			1665	1056	272	330	7			

- Molecule 3 is a protein called 82D6A3 IgG.

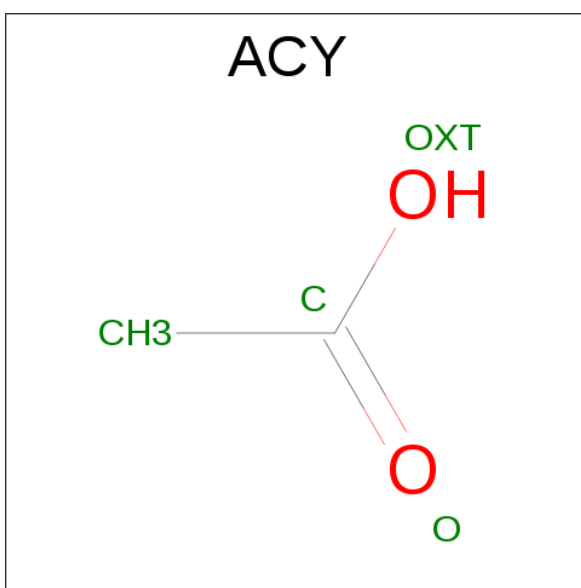
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	L	209	Total	C	N	O	S	0	0	0
			1624	1015	274	329	6			

- Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	H	1	Total	O	S	0	0
			5	4	1		
4	L	1	Total	O	S	0	0
			5	4	1		

- Molecule 5 is ACETIC ACID (three-letter code: ACY) (formula: $C_2H_4O_2$).



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

- Molecule 6 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	H	1	Total	C	O	0	0
			6	3	3		


- Molecule 7 is water.

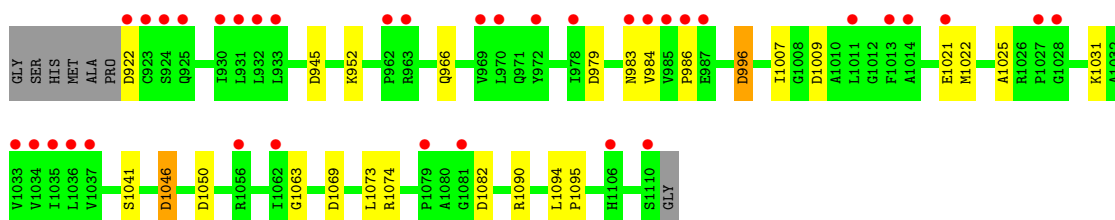
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	136	Total	O	0	0
			136	136		
7	H	155	Total	O	0	0
			155	155		
7	L	151	Total	O	0	0
			151	151		

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.

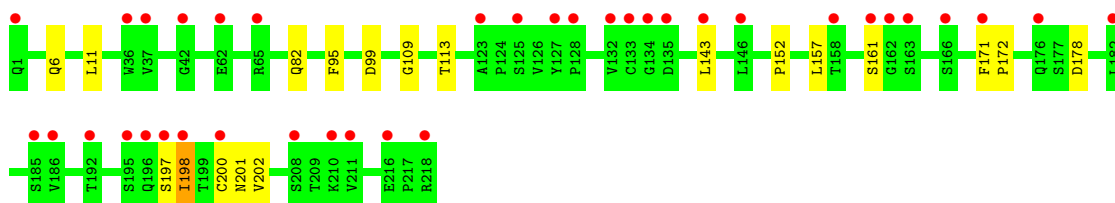
• Molecule 1: Von Willebrand factor

Chain A: 




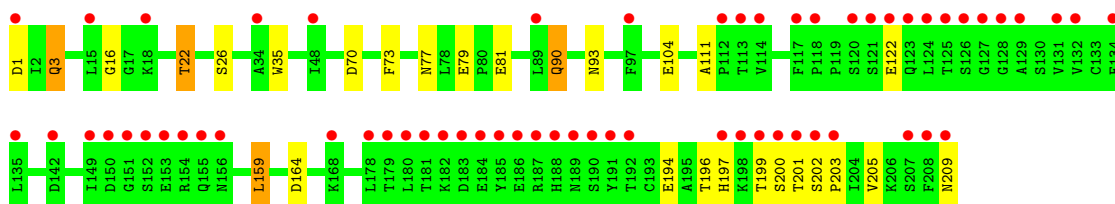
• Molecule 2: 82D6A3 IgG

Chain H: 



• Molecule 3: 82D6A3 IgG

Chain L: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	72.18Å 89.08Å 123.51Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	27.35 – 1.90 27.34 – 1.90	Depositor EDS
% Data completeness (in resolution range)	100.0 (27.35-1.90) 99.8 (27.34-1.90)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.25 (at 1.91Å)	Xtriage
Refinement program	REFMAC 5.2.0003	Depositor
R, R_{free}	0.191 , 0.220 0.246 , 0.267	Depositor DCC
R_{free} test set	3240 reflections (5.12%)	DCC
Wilson B-factor (Å ²)	22.9	Xtriage
Anisotropy	0.246	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 45.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	5165	wwPDB-VP
Average B, all atoms (Å ²)	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.23% 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: GOL, ACY, 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	0.48	0/1439	0.86	10/1957 (0.5%)
2	H	0.46	0/1709	0.74	2/2333 (0.1%)
3	L	0.60	2/1662 (0.1%)	0.81	3/2253 (0.1%)
All	All	0.52	2/4810 (0.0%)	0.80	15/6543 (0.2%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	L	209	ASN	C-OXT	13.29	1.48	1.23
3	L	122	GLU	CD-OE1	5.28	1.31	1.25

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	979	ASP	CB-CG-OD2	7.34	124.91	118.30
3	L	159	LEU	CA-CB-CG	6.80	130.94	115.30
2	H	99	ASP	CB-CG-OD2	5.67	123.41	118.30
1	A	1009	ASP	CB-CG-OD2	5.66	123.40	118.30
1	A	1082	ASP	CB-CG-OD2	5.63	123.36	118.30
1	A	1050	ASP	CB-CG-OD2	5.54	123.28	118.30
1	A	1090	ARG	NE-CZ-NH2	-5.42	117.59	120.30
1	A	922	ASP	CB-CG-OD2	5.28	123.06	118.30
3	L	70	ASP	CB-CG-OD2	5.28	123.05	118.30
1	A	1046	ASP	CB-CG-OD2	5.24	123.02	118.30
1	A	996	ASP	CB-CG-OD2	5.20	122.98	118.30
3	L	164	ASP	CB-CG-OD2	5.08	122.88	118.30
1	A	1069	ASP	CB-CG-OD2	5.04	122.84	118.30
1	A	945	ASP	CB-CG-OD2	5.03	122.83	118.30
2	H	178	ASP	CB-CG-OD2	5.01	122.81	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	1414	0	1425	10	0
2	H	1665	0	1621	13	0
3	L	1624	0	1567	14	0
4	H	5	0	0	0	0
4	L	5	0	0	0	0
5	A	4	0	3	0	0
6	H	6	0	8	1	0
7	A	136	0	0	2	0
7	H	155	0	0	4	0
7	L	151	0	0	2	0
All	All	5165	0	4624	38	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 (38) 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)
3:L:200:SER:HB3	3:L:202:SER:O	1.63	0.99
2:H:161:SER:H	2:H:201:ASN:HD21	1.24	0.84
1:A:1021:GLU:HG2	7:A:1227:HOH:O	1.78	0.83
1:A:984:VAL:O	1:A:986:PRO:HD3	1.87	0.74
2:H:6:GLN:HE21	2:H:109:GLY:HA3	1.53	0.73
3:L:200:SER:CB	3:L:202:SER:O	2.37	0.72
3:L:22:THR:HG22	7:L:1049:HOH:O	1.93	0.66
2:H:161:SER:H	2:H:201:ASN:ND2	1.95	0.63
6:H:1004:GOL:H32	7:H:1066:HOH:O	1.99	0.62
2:H:157:LEU:HD11	2:H:200:CYS:HB2	1.82	0.61
3:L:16:GLY:HA2	3:L:77:ASN:HD22	1.65	0.61
1:A:952:LYS:NZ	1:A:996:ASP:OD1	2.34	0.58
2:H:157:LEU:HD13	2:H:202:VAL:HG22	1.84	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:L:200:SER:HB2	7:L:1024:HOH:O	2.04	0.57
2:H:113:THR:HG22	7:H:1014:HOH:O	2.02	0.57
3:L:199:THR:O	3:L:200:SER:HB2	2.06	0.55
3:L:79:GLU:OE1	3:L:81:GLU:OE2	2.27	0.53
1:A:966:GLN:HE21	1:A:983:ASN:H	1.57	0.52
3:L:202:SER:HB3	3:L:203:PRO:HD2	1.92	0.52
3:L:90:GLN:HE22	3:L:93:ASN:H	1.58	0.51
2:H:143:LEU:HD12	2:H:198:ILE:HG21	1.97	0.47
1:A:1007:ILE:HD12	1:A:1041:SER:HB2	1.97	0.46
3:L:35:TRP:CE2	3:L:73:PHE:HB2	2.51	0.45
2:H:6:GLN:HE22	2:H:95:PHE:HA	1.81	0.45
1:A:1094:LEU:HB3	1:A:1095:PRO:HD3	1.99	0.45
1:A:952:LYS:CE	1:A:996:ASP:OD1	2.64	0.45
3:L:194:GLU:HG2	3:L:205:VAL:HG22	2.00	0.44
3:L:111:ALA:HA	3:L:199:THR:HB	1.99	0.44
2:H:6:GLN:HE21	2:H:109:GLY:CA	2.28	0.42
3:L:196:THR:CG2	3:L:203:PRO:HB3	2.49	0.42
1:A:1063:GLY:HA3	1:A:1073:LEU:HD11	2.00	0.42
2:H:11:LEU:HD22	2:H:152:PRO:HG3	2.01	0.41
2:H:171:PHE:HA	2:H:172:PRO:HD3	1.96	0.41
3:L:3:GLN:HG2	3:L:26:SER:HB3	2.02	0.41
1:A:1074:ARG:NH2	7:A:1242:HOH:O	2.50	0.40
2:H:82:GLN:NE2	7:H:1158:HOH:O	2.55	0.40
1:A:1025:ALA:HB1	1:A:1031:LYS:HZ2	1.87	0.40
2:H:113:THR:HG21	7:H:1105:HOH:O	2.21	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	187/196 (95%)	180 (96%)	7 (4%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	H	216/218 (99%)	210 (97%)	4 (2%)	2 (1%)	20	8
3	L	207/209 (99%)	200 (97%)	6 (3%)	1 (0%)	32	20
All	All	610/623 (98%)	590 (97%)	17 (3%)	3 (0%)	32	20

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	H	197	SER
3	L	197	HIS
2	H	198	ILE

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	156/160 (98%)	154 (99%)	2 (1%)	73	72
2	H	187/187 (100%)	187 (100%)	0	100	100
3	L	185/185 (100%)	178 (96%)	7 (4%)	38	27
All	All	528/532 (99%)	519 (98%)	9 (2%)	66	62

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1022	MET
1	A	1046	ASP
3	L	1	ASP
3	L	3	GLN
3	L	22	THR
3	L	90	GLN
3	L	104	GLU
3	L	159	LEU
3	L	201	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such

sidechains are listed below:

Mol	Chain	Res	Type
1	A	959	ASN
1	A	966	GLN
2	H	6	GLN
2	H	169	HIS
2	H	201	ASN
3	L	77	ASN
3	L	90	GLN
3	L	137	ASN
3	L	160	ASN
3	L	189	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](#)

4 ligands are 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	$\# Z > 2$
5	ACY	A	1112	-	1,3,3	1.41	0	0,3,3	0.00	-
4	SO4	H	1002	-	4,4,4	0.13	0	6,6,6	0.19	0
6	GOL	H	1004	-	5,5,5	0.36	0	5,5,5	0.30	0
4	SO4	L	1003	-	4,4,4	0.21	0	6,6,6	0.25	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	ACY	A	1112	-	-	0/0/0/0	0/0/0/0
4	SO4	H	1002	-	-	0/0/0/0	0/0/0/0
6	GOL	H	1004	-	-	0/4/4/4	0/0/0/0
4	SO4	L	1003	-	-	0/0/0/0	0/0/0/0

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.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	H	1004	GOL	1	0

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	189/196 (96%)	0.91	36 (19%)  	10, 16, 30, 38	0
2	H	218/218 (100%)	1.03	37 (16%)  	10, 22, 34, 35	0
3	L	209/209 (100%)	1.58	61 (29%)  	13, 17, 25, 43	0
All	All	616/623 (98%)	1.18	134 (21%)  	10, 17, 33, 43	0

All (134) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	L	201	THR	12.4
1	A	985	VAL	9.4
3	L	199	THR	8.1
1	A	984	VAL	7.9
3	L	181	THR	6.6
3	L	128	GLY	6.4
3	L	185	TYR	6.2
3	L	124	LEU	6.2
3	L	191	TYR	5.9
3	L	125	THR	5.9
3	L	180	LEU	5.6
2	H	133	CYS	5.4
3	L	183	ASP	5.3
3	L	190	SER	5.3
3	L	198	LYS	5.3
3	L	149	ILE	5.0
3	L	202	SER	5.0
3	L	155	GLN	4.9
2	H	197	SER	4.9
3	L	200	SER	4.7
2	H	163	SER	4.6
3	L	121	SER	4.6
3	L	189	ASN	4.5

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Mol	Chain	Res	Type	RSRZ
3	L	187	ARG	4.3
3	L	152	SER	4.3
3	L	156	ASN	4.3
3	L	118	PRO	4.2
3	L	168	LYS	4.2
3	L	203	PRO	4.2
3	L	1	ASP	4.1
1	A	922	ASP	4.1
2	H	127	TYR	4.0
2	H	134	GLY	4.0
2	H	162	GLY	4.0
2	H	195	SER	4.0
2	H	210	LYS	3.9
2	H	146	LEU	3.9
3	L	208	PHE	3.8
3	L	209	ASN	3.8
3	L	184	GLU	3.7
1	A	923	CYS	3.7
3	L	123	GLN	3.6
3	L	188	HIS	3.6
3	L	154	ARG	3.6
1	A	1021	GLU	3.6
3	L	151	GLY	3.5
3	L	182	LYS	3.5
3	L	129	ALA	3.5
2	H	123	ALA	3.5
3	L	122	GLU	3.4
1	A	1110	SER	3.4
2	H	208	SER	3.3
1	A	932	LEU	3.3
3	L	135	LEU	3.3
1	A	986	PRO	3.3
3	L	179	THR	3.2
1	A	970	LEU	3.2
2	H	128	PRO	3.2
2	H	1	GLN	3.2
1	A	1014	ALA	3.2
3	L	126	SER	3.1
1	A	925	GLN	3.1
1	A	1033	VAL	3.1
3	L	207	SER	3.1
3	L	89	LEU	3.1

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Mol	Chain	Res	Type	RSRZ
3	L	186	GLU	3.0
2	H	166	SER	3.0
1	A	1013	PHE	3.0
3	L	192	THR	2.9
3	L	150	ASP	2.9
1	A	1034	VAL	2.8
1	A	1036	LEU	2.8
3	L	178	LEU	2.8
3	L	197	HIS	2.8
3	L	120	SER	2.8
2	H	186	VAL	2.8
1	A	1079	PRO	2.7
1	A	931	LEU	2.7
1	A	983	ASN	2.7
1	A	1035	ILE	2.7
1	A	1106	HIS	2.7
3	L	153	GLU	2.7
2	H	198	ILE	2.6
2	H	200	CYS	2.6
1	A	1081	GLY	2.6
3	L	131	VAL	2.6
2	H	36	TRP	2.6
1	A	963	ARG	2.5
2	H	211	VAL	2.5
2	H	161	SER	2.5
1	A	1027	PRO	2.5
3	L	132	VAL	2.5
2	H	196	GLN	2.5
1	A	1056	ARG	2.5
1	A	924	SER	2.4
1	A	1011	LEU	2.4
3	L	127	GLY	2.4
3	L	18	LYS	2.4
1	A	987	GLU	2.4
1	A	930	ILE	2.4
1	A	1028	GLY	2.4
1	A	969	VAL	2.3
3	L	97	PHE	2.3
1	A	962	PRO	2.3
2	H	135	ASP	2.3
2	H	218	ARG	2.3
3	L	34	ALA	2.3

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Mol	Chain	Res	Type	RSRZ
2	H	143	LEU	2.3
3	L	15	LEU	2.3
1	A	978	ILE	2.2
2	H	185	SER	2.2
3	L	142	ASP	2.2
2	H	37	VAL	2.2
3	L	114	VAL	2.2
2	H	216	GLU	2.2
1	A	1062	ILE	2.2
3	L	48	ILE	2.2
2	H	182	LEU	2.2
1	A	972	TYR	2.2
3	L	134	PHE	2.2
1	A	1037	VAL	2.2
3	L	112	PRO	2.1
2	H	192	THR	2.1
2	H	171	PHE	2.1
3	L	117	PHE	2.1
2	H	176	GLN	2.1
2	H	42	GLY	2.1
2	H	132	VAL	2.1
2	H	125	SER	2.1
2	H	62	GLU	2.1
2	H	65	ARG	2.1
2	H	158	THR	2.1
1	A	933	LEU	2.1
3	L	113	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 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains.

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	LLDF	B-factors(\AA^2)	Q<0.9
4	SO4	L	1003	5/5	0.95	0.19	4.59	38,39,40,40	0
6	GOL	H	1004	6/6	0.85	0.20	2.73	41,42,43,43	0
4	SO4	H	1002	5/5	0.97	0.18	0.96	29,30,33,33	0
5	ACY	A	1112	4/4	0.92	0.11	-0.86	35,35,35,35	0

6.5 Other polymers [i](#)

There are no such residues in this entry.