



Full wwPDB X-ray Structure Validation Report

Feb 27, 2014 – 08:33 PM GMT

PDB ID : 2R2B
Title : Crystal structure of S25-2 Fab in complex with Kdo analogues
Authors : Brooks, C.L.; Evans, S.V.
Deposited on : 2007-08-24
Resolution : 1.60 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

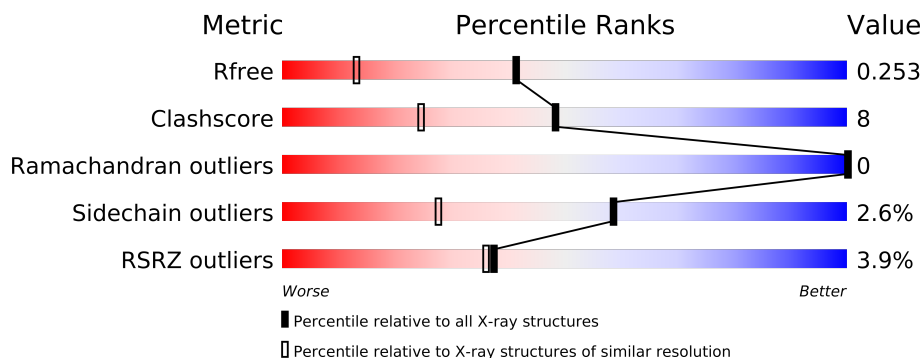
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 1.60 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}	66092	1872 (1.60-1.60)
Clashscore	79885	2199 (1.60-1.60)
Ramachandran outliers	78287	2126 (1.60-1.60)
Sidechain outliers	78261	2125 (1.60-1.60)
RSRZ outliers	66119	1872 (1.60-1.60)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	219	
2	B	222	

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 3698 atoms, of which 0 are hydrogen and 0 are deuterium.

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 Fab, antibody fragment (IgG1k), light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	217	Total	C	N	O	S	0	0	0
			1684	1047	286	343	8			

- Molecule 2 is a protein called Fab, antibody fragment (IgG1k), heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	214	Total	C	N	O	S	0	1	0
			1651	1051	272	321	7			

- Molecule 3 is a polymer of unknown type called SUGAR (3-MER).

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	3	Total	C	O	0	0
			49	27	22		

- Molecule 4 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total	Zn	0	0
			1	1		
4	A	1	Total	Zn	0	0
			1	1		

- Molecule 5 is water.

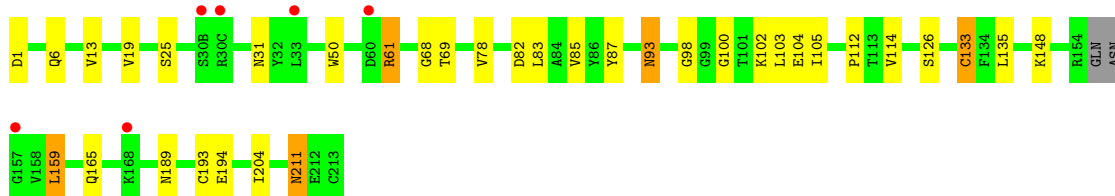
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	166	Total	O	0	0
			166	166		
5	B	146	Total	O	0	0
			146	146		

3 Residue-property plots

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.

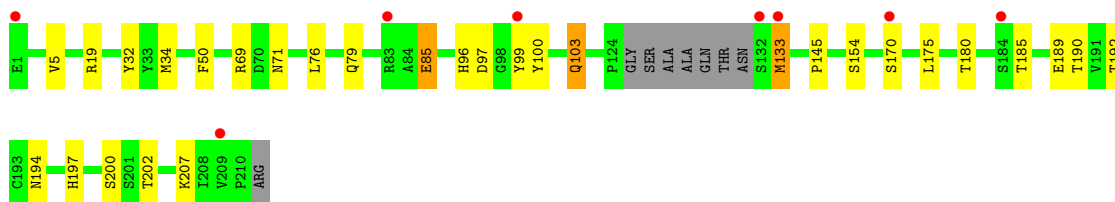
- Molecule 1: Fab, antibody fragment (IgG1k), light chain

Chain A: 



- Molecule 2: Fab, antibody fragment (IgG1k), heavy chain

Chain B: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	45.89Å 81.66Å 131.55Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 1.60 19.92 – 1.60	Depositor EDS
% Data completeness (in resolution range)	95.7 (20.00-1.60) 95.7 (19.92-1.60)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.54 (at 1.60Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.231 , 0.253 0.228 , 0.253	Depositor DCC
R_{free} test set	3212 reflections (5.35%)	DCC
Wilson B-factor (Å ²)	22.4	Xtriage
Anisotropy	0.134	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 33.8	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Outliers	0 of 63237 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	3698	wwPDB-VP
Average B, all atoms (Å ²)	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.27% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, KDO, KDA

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	0.60	1/1718 (0.1%)	0.71	2/2325 (0.1%)
2	B	0.59	0/1696	0.71	0/2315
All	All	0.59	1/3414 (0.0%)	0.71	2/4640 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	133	CYS	CB-SG	-5.33	1.73	1.81

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	61	ARG	NE-CZ-NH1	-6.68	116.96	120.30
1	A	159	LEU	CA-CB-CG	6.13	129.41	115.30

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1684	0	1634	32	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	1651	0	1607	21	0
3	B	49	0	37	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
5	A	166	0	0	9	0
5	B	146	0	0	1	0
All	All	3698	0	3278	52	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 8.

All (52) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:133:CYS:SG	5:A:377:HOH:O	2.06	1.12
1:A:25:SER:HB2	5:A:378:HOH:O	0.85	1.02
1:A:133:CYS:CB	5:A:377:HOH:O	2.16	0.94
2:B:69:ARG:HE	2:B:71:ASN:HD21	1.27	0.82
1:A:114:VAL:HG22	1:A:135:LEU:HD12	1.68	0.75
1:A:61:ARG:HD3	5:A:301:HOH:O	1.86	0.75
1:A:133:CYS:HB2	5:A:377:HOH:O	1.82	0.73
2:B:154:SER:H	2:B:194:ASN:HD21	1.33	0.73
1:A:189:ASN:HD21	1:A:211:ASN:H	1.37	0.72
1:A:105:ILE:H	1:A:165:GLN:HE22	1.44	0.66
2:B:69:ARG:NE	2:B:71:ASN:HD21	1.99	0.59
1:A:126:SER:HB3	5:A:343:HOH:O	2.02	0.59
2:B:145:PRO:O	2:B:197:HIS:HE1	1.87	0.57
2:B:96:HIS:HD2	2:B:100:TYR:O	1.87	0.57
2:B:5:VAL:HA	2:B:103:GLN:HE22	1.69	0.56
1:A:1:ASP:O	5:A:338:HOH:O	2.18	0.55
2:B:69:ARG:HE	2:B:71:ASN:ND2	2.02	0.55
1:A:83:LEU:HD13	1:A:165:GLN:HE21	1.70	0.55
2:B:190:THR:HG23	5:B:267:HOH:O	2.07	0.53
2:B:103:GLN:H	2:B:103:GLN:NE2	2.06	0.53
2:B:197:HIS:HD2	2:B:200:SER:OG	1.92	0.53
1:A:93:ASN:HD22	1:A:93:ASN:C	2.12	0.53
2:B:154:SER:H	2:B:194:ASN:ND2	2.04	0.53
1:A:31:ASN:ND2	1:A:68:GLY:H	2.07	0.52
2:B:192:THR:OG1	2:B:207[A]:LYS:HD3	2.11	0.50
1:A:61:ARG:NH1	1:A:82:ASP:OD2	2.36	0.50
1:A:193:CYS:CB	5:A:377:HOH:O	2.61	0.49
1:A:114:VAL:HG13	1:A:135:LEU:CD1	2.43	0.49
1:A:13:VAL:HG11	1:A:19:VAL:CG2	2.43	0.48

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:B:19:ARG:HH11	2:B:79:GLN:HE22	1.61	0.48
1:A:114:VAL:HG13	1:A:135:LEU:HD13	1.95	0.47
1:A:13:VAL:HG12	1:A:103:LEU:HD11	1.97	0.46
1:A:6:GLN:HE21	1:A:98:GLY:HA3	1.81	0.46
1:A:13:VAL:HG13	1:A:78:VAL:HG11	1.97	0.46
1:A:31:ASN:HD21	1:A:68:GLY:H	1.63	0.45
2:B:133:MET:HG3	2:B:180:THR:HG22	1.99	0.45
1:A:112:PRO:HG2	1:A:204:ILE:HD12	1.97	0.44
2:B:19:ARG:HH11	2:B:79:GLN:NE2	2.15	0.44
1:A:6:GLN:HE22	1:A:87:TYR:HA	1.82	0.44
1:A:193:CYS:HB2	5:A:377:HOH:O	2.17	0.43
1:A:85:VAL:HG22	1:A:102:LYS:HG2	2.01	0.42
1:A:25:SER:OG	1:A:69:THR:HA	2.19	0.42
2:B:185:THR:O	2:B:189:GLU:HB2	2.19	0.42
1:A:50:TRP:HH2	2:B:99:TYR:CE2	2.38	0.42
2:B:197:HIS:HB3	2:B:202:THR:HB	2.01	0.41
1:A:85:VAL:HG22	1:A:102:LYS:CG	2.50	0.41
1:A:6:GLN:NE2	1:A:100:GLY:H	2.19	0.41
1:A:148:LYS:HE2	1:A:194:GLU:OE1	2.20	0.41
2:B:32:TYR:CE1	2:B:97:ASP:HB3	2.55	0.41
2:B:34:MET:HB3	2:B:76:LEU:HD22	2.03	0.41
2:B:85:GLU:O	2:B:85:GLU:HG2	2.21	0.41
1:A:61:ARG:HH12	1:A:82:ASP:CG	2.22	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	213/219 (97%)	210 (99%)	3 (1%)	0	100	100
2	B	211/222 (95%)	204 (97%)	7 (3%)	0	100	100
All	All	424/441 (96%)	414 (98%)	10 (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	193/195 (99%)	189 (98%)	4 (2%)	66	37
2	B	186/190 (98%)	180 (97%)	6 (3%)	51	20
All	All	379/385 (98%)	369 (97%)	10 (3%)	59	28

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	93	ASN
1	A	104	GLU
1	A	159	LEU
1	A	211	ASN
2	B	50	PHE
2	B	85	GLU
2	B	103	GLN
2	B	133	MET
2	B	170	SER
2	B	175	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	6	GLN
1	A	31	ASN
1	A	42	GLN
1	A	93	ASN
1	A	136	ASN
1	A	160	ASN
1	A	165	GLN
1	A	189	ASN
1	A	211	ASN
2	B	71	ASN
2	B	79	GLN
2	B	96	HIS
2	B	103	GLN
2	B	169	GLN

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Mol	Chain	Res	Type
2	B	194	ASN
2	B	197	HIS

5.3.3 RNA ⓘ

There are no RNA chains in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

3 carbohydrates 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$
3	KDA	B	214	3	18,18,19	2.20	5 (27%)	21,25,27	2.95	7 (33%)
3	KDO	B	215	3	15,15,16	2.62	5 (33%)	18,22,24	1.86	4 (22%)
3	KDO	B	216	3	16,16,16	1.76	3 (18%)	24,24,24	1.38	4 (16%)

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
3	KDA	B	214	3	-	0/17/32/35	0/1/1/1
3	KDO	B	215	3	-	0/12/27/30	0/1/1/1
3	KDO	B	216	3	-	0/12/30/30	0/1/1/1

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	214	KDA	C3-C2	6.18	1.56	1.52
3	B	215	KDO	O2-C2	5.63	1.47	1.39
3	B	215	KDO	C3-C2	5.32	1.56	1.52
3	B	215	KDO	O6-C2	5.02	1.47	1.42
3	B	214	KDA	O2-C2	4.12	1.46	1.41
3	B	216	KDO	O2-C2	4.10	1.45	1.39
3	B	216	KDO	O6-C2	3.81	1.46	1.42
3	B	214	KDA	O6-C2	3.19	1.47	1.42
3	B	216	KDO	C3-C2	2.55	1.55	1.51
3	B	214	KDA	O2-C9	-2.51	1.38	1.43
3	B	215	KDO	O6-C6	2.14	1.47	1.44
3	B	215	KDO	C6-C5	2.13	1.54	1.51
3	B	214	KDA	C6-C5	2.06	1.54	1.51

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	214	KDA	C9-O2-C2	9.40	128.03	115.49
3	B	214	KDA	O2-C9-C10	5.65	128.32	109.89
3	B	214	KDA	C2-O6-C6	5.10	122.93	113.99
3	B	215	KDO	C2-O6-C6	4.32	120.51	113.93
3	B	215	KDO	O6-C6-C7	4.09	114.01	106.69
3	B	214	KDA	O6-C6-C7	3.67	113.26	106.69
3	B	216	KDO	O1A-C1-C2	-3.20	119.04	123.48
3	B	215	KDO	O1A-C1-C2	-3.08	119.20	123.48
3	B	216	KDO	C3-C2-C1	-2.93	108.66	113.58
3	B	215	KDO	C3-C2-C1	-2.83	107.21	112.91
3	B	216	KDO	O2-C2-C1	2.72	114.48	109.86
3	B	214	KDA	O1A-C1-C2	2.63	118.57	113.33
3	B	216	KDO	O6-C6-C7	2.51	111.17	106.69
3	B	214	KDA	O6-C2-C3	2.25	114.56	111.16
3	B	214	KDA	O1B-C1-C2	-2.04	117.09	124.03

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.6 Ligand geometry ⓘ

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

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.

5.7 Other polymers ⓘ

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	217/219 (99%)	0.09	6 (2%)	50	49	13, 19, 27, 36	0
2	B	214/222 (96%)	0.13	8 (3%)	39	38	13, 18, 29, 35	0
All	All	431/441 (97%)	0.11	14 (3%)	37	44	13, 18, 29, 36	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	99	TYR	5.2
2	B	132	SER	2.9
1	A	30(C)	ARG	2.8
2	B	170	SER	2.8
1	A	168	LYS	2.6
2	B	184	SER	2.6
2	B	209	VAL	2.5
2	B	1	GLU	2.5
1	A	60	ASP	2.4
2	B	133	MET	2.3
1	A	30(B)	SER	2.3
1	A	33	LEU	2.2
1	A	157	GLY	2.0
2	B	83	ARG	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

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	RSR	LLDF	B-factors(Å ²)	Q<0.9
3	KDO	B	216	16/16	0.09	-	15,17,25,30	0
3	KDO	B	215	15/16	0.23	-	27,29,35,36	7
3	KDA	B	214	18/19	0.64	-	37,39,40,40	18

6.4 Ligands

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	RSR	LLDF	B-factors(Å ²)	Q<0.9
4	ZN	A	214	1/1	0.02	-	18,18,18,18	0
4	ZN	B	217	1/1	0.26	-	47,47,47,47	0

6.5 Other polymers

There are no such residues in this entry.