



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 28, 2018 – 07:13 AM EST

PDB ID : 2R23
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 : 1.65 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 : rb-20030736
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) : rb-20030736

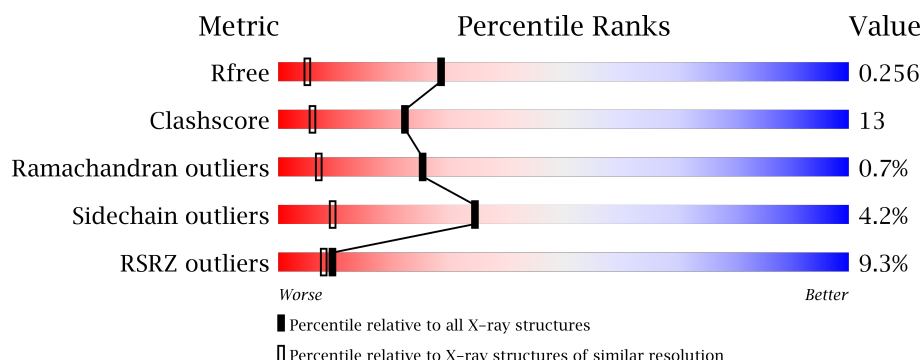
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.65 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	1368 (1.66-1.66)
Clashscore	112137	1468 (1.66-1.66)
Ramachandran outliers	110173	1438 (1.66-1.66)
Sidechain outliers	110143	1438 (1.66-1.66)
RSRZ outliers	101464	1371 (1.66-1.66)

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	219	
2	B	224	

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
3	MG	A	216	-	-	-	X

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 3809 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 Fab, antibody fragment (IgG1k), light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	217	Total	C	N	O	S	0	1	0
			1693	1053	288	344	8			

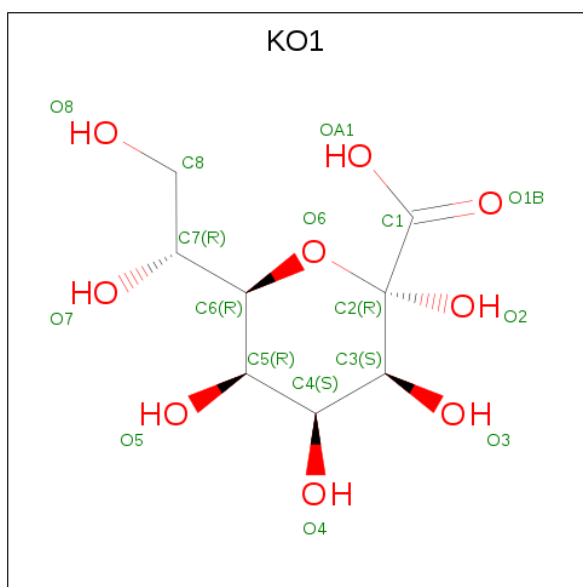
- Molecule 2 is a protein called Fab, antibody fragment (IgG1k), heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	222	Total	C	N	O	S	0	2	0
			1708	1079	284	337	8			

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

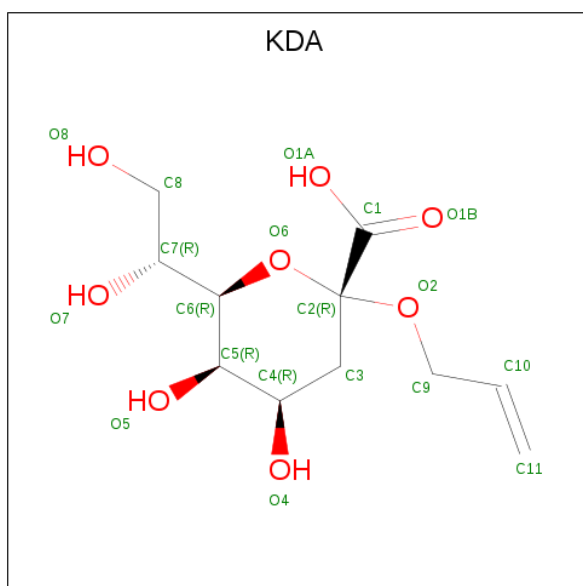
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Mg	0	0
			1	1		

- Molecule 4 is D-glycero-alpha-D-talo-oct-2-ulopyranosonic acid (three-letter code: KO1) (formula: C₈H₁₄O₉).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	C	O	0	0
			17	8	9		

- Molecule 5 is (3-DEOXY-D-MANNO-OCT-2-ULOSONIC ACID)-2-O-ALLYL (three-letter code: KDA) (formula: $C_{11}H_{18}O_8$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total	C	O	0	0
			18	11	7		

- Molecule 6 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	B	1	Total 1	Zn 1	0	0

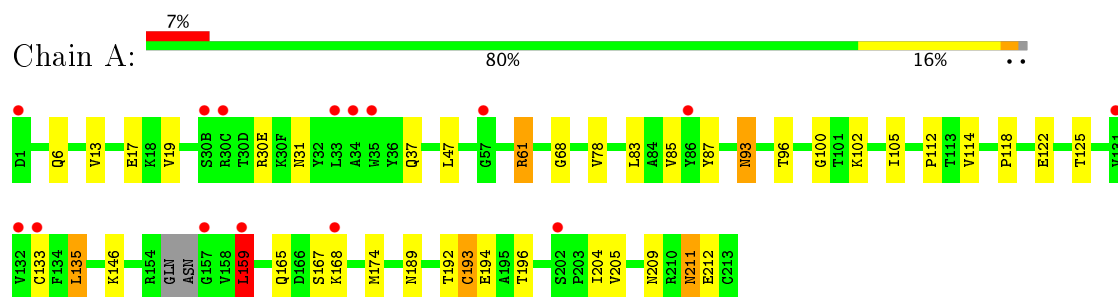
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	182	Total 182	O 182	0	0
7	B	189	Total 189	O 189	0	0

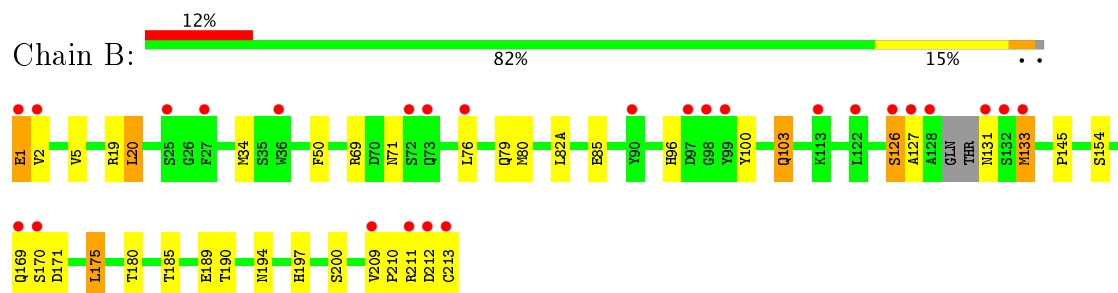
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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: Fab, antibody fragment (IgG1k), light chain



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



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	45.64Å 81.38Å 131.06Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 1.65 19.90 – 1.65	Depositor EDS
% Data completeness (in resolution range)	95.8 (20.00-1.65) 95.8 (19.90-1.65)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.74 (at 1.65Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.228 , 0.264 0.221 , 0.256	Depositor DCC
R_{free} test set	2889 reflections (5.33%)	DCC
Wilson B-factor (Å ²)	31.2	Xtriage
Anisotropy	0.068	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 39.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.52$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	3809	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.32% 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: KO1, ZN, MG, 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	0.78	1/1727 (0.1%)	0.75	1/2336 (0.0%)
2	B	0.77	0/1753	0.79	1/2392 (0.0%)
All	All	0.78	1/3480 (0.0%)	0.77	2/4728 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	193	CYS	CB-SG	-5.39	1.73	1.81

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	159	LEU	CA-CB-CG	8.26	134.29	115.30
2	B	175	LEU	CA-CB-CG	5.97	129.03	115.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	1693	0	1648	57	0
2	B	1708	0	1650	39	0
3	A	1	0	0	0	1

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	17	0	12	0	0
5	B	18	0	15	0	0
6	B	1	0	0	0	0
7	A	182	0	0	12	2
7	B	189	0	0	5	3
All	All	3809	0	3325	88	3

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 13.

All (88) 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:125:THR:HB	7:A:361:HOH:O	1.36	1.23
2:B:20:LEU:HG	7:B:483:HOH:O	0.99	1.17
1:A:133:CYS:SG	1:A:193:CYS:HB2	1.92	1.09
1:A:159:LEU:HD22	2:B:169:GLN:NE2	1.77	0.99
1:A:61:ARG:HH11	1:A:61:ARG:HG3	1.27	0.97
1:A:133:CYS:SG	1:A:193:CYS:CB	2.54	0.95
1:A:159:LEU:CD2	2:B:169:GLN:NE2	2.30	0.93
2:B:20:LEU:HD11	2:B:80:MET:CE	1.99	0.93
1:A:85:VAL:HG22	1:A:102[B]:LYS:HG3	1.54	0.90
2:B:20:LEU:HD11	2:B:80:MET:HE1	1.54	0.88
1:A:19:VAL:HG23	1:A:78:VAL:HG11	1.57	0.86
1:A:159:LEU:HD21	7:B:463:HOH:O	1.76	0.86
2:B:154:SER:H	2:B:194:ASN:HD21	1.23	0.84
2:B:20:LEU:HD12	2:B:20:LEU:N	1.95	0.81
2:B:69:ARG:HE	2:B:71:ASN:HD21	1.27	0.80
1:A:196:THR:HG23	7:A:261:HOH:O	1.80	0.78
1:A:61:ARG:NH1	1:A:61:ARG:HG3	1.93	0.75
1:A:159:LEU:HB3	7:A:301:HOH:O	1.85	0.75
1:A:105:ILE:H	1:A:165:GLN:HE22	1.31	0.75
1:A:167:SER:OG	7:A:341:HOH:O	2.06	0.73
7:A:294:HOH:O	2:B:209:VAL:HG11	1.90	0.70
2:B:20:LEU:HD11	2:B:80:MET:HE2	1.74	0.68
1:A:159:LEU:HD22	2:B:169:GLN:CD	2.12	0.68
1:A:159:LEU:HD22	2:B:169:GLN:HE22	1.59	0.67
2:B:96:HIS:HD2	2:B:100:TYR:O	1.77	0.67
1:A:19:VAL:CG2	1:A:78:VAL:HG11	2.25	0.66
1:A:114:VAL:HG13	1:A:135:LEU:CD1	2.25	0.66
1:A:159:LEU:HD21	2:B:169:GLN:NE2	2.11	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:190:THR:HG23	7:B:395:HOH:O	1.97	0.65
1:A:204:ILE:HD11	7:A:228:HOH:O	1.96	0.65
1:A:133:CYS:HG	1:A:193:CYS:HB2	1.63	0.63
2:B:20:LEU:HD12	2:B:20:LEU:H	1.63	0.62
2:B:69:ARG:NE	2:B:71:ASN:HD21	1.96	0.61
2:B:154:SER:H	2:B:194:ASN:ND2	1.98	0.60
1:A:159:LEU:CD2	2:B:169:GLN:HE22	2.10	0.59
1:A:209:ASN:HB3	1:A:211:ASN:ND2	2.17	0.59
1:A:189:ASN:HD21	1:A:211:ASN:H	1.50	0.59
2:B:5:VAL:HA	2:B:103:GLN:HE22	1.69	0.58
2:B:197:HIS:HD2	2:B:200:SER:OG	1.86	0.58
2:B:69:ARG:HE	2:B:71:ASN:ND2	2.00	0.57
1:A:159:LEU:HD12	7:A:301:HOH:O	2.06	0.56
1:A:133:CYS:CB	1:A:193:CYS:SG	2.93	0.56
1:A:6:GLN:HE22	1:A:87:TYR:HA	1.71	0.55
1:A:135:LEU:HD23	1:A:174:MET:HG2	1.89	0.55
1:A:61:ARG:HH11	1:A:61:ARG:CG	2.07	0.55
2:B:145:PRO:O	2:B:197:HIS:HE1	1.91	0.54
1:A:13:VAL:HG11	1:A:19:VAL:CG2	2.39	0.53
1:A:209:ASN:HB3	1:A:211:ASN:HD22	1.72	0.53
2:B:212:ASP:C	2:B:213:CYS:SG	2.88	0.52
1:A:159:LEU:HD23	7:A:229:HOH:O	2.10	0.52
1:A:96:THR:CG2	7:A:241:HOH:O	2.58	0.52
1:A:83:LEU:HD13	1:A:165:GLN:HE21	1.75	0.51
1:A:13:VAL:CG2	1:A:17:GLU:HB2	2.40	0.51
1:A:114:VAL:HG22	1:A:135:LEU:HD12	1.93	0.50
1:A:194:GLU:HG2	1:A:205:VAL:HG22	1.94	0.49
2:B:133:MET:HG3	2:B:180:THR:HG22	1.95	0.49
2:B:19:ARG:HH11	2:B:79:GLN:HE22	1.59	0.49
2:B:34:MET:HB3	2:B:76:LEU:HD22	1.95	0.49
2:B:103:GLN:H	2:B:103:GLN:HE21	1.59	0.49
2:B:103:GLN:H	2:B:103:GLN:NE2	2.11	0.48
1:A:114:VAL:HG13	1:A:135:LEU:HD12	1.94	0.48
2:B:185:THR:O	2:B:189:GLU:HB2	2.13	0.48
1:A:133:CYS:HB3	1:A:193:CYS:SG	2.54	0.48
2:B:20:LEU:CD1	2:B:20:LEU:N	2.67	0.47
1:A:13:VAL:HG22	1:A:17:GLU:HB2	1.96	0.47
2:B:210:PRO:HB2	2:B:213:CYS:SG	2.53	0.47
2:B:1:GLU:HG2	2:B:2:VAL:N	2.30	0.47
1:A:30(E):ARG:HD3	7:A:331:HOH:O	2.15	0.47
1:A:31:ASN:ND2	1:A:68:GLY:H	2.15	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:93:ASN:HD22	1:A:93:ASN:C	2.20	0.45
1:A:168:LYS:HE2	7:A:300:HOH:O	2.16	0.45
1:A:31:ASN:HD21	1:A:68:GLY:H	1.65	0.44
1:A:192:THR:HG22	1:A:194:GLU:HG3	1.99	0.43
1:A:13:VAL:HG13	1:A:78:VAL:HG21	2.00	0.43
1:A:37:GLN:HB2	1:A:47:LEU:HD11	1.99	0.43
1:A:212:GLU:O	2:B:126:SER:HB2	2.19	0.42
2:B:19:ARG:HH11	2:B:79:GLN:NE2	2.18	0.42
2:B:96:HIS:CD2	2:B:100:TYR:O	2.67	0.42
1:A:6:GLN:NE2	1:A:100:GLY:H	2.18	0.42
1:A:112:PRO:HG2	1:A:204:ILE:HD13	2.01	0.41
1:A:61:ARG:NH1	1:A:61:ARG:CG	2.67	0.41
2:B:211:ARG:NH2	7:B:394:HOH:O	2.54	0.41
1:A:105:ILE:H	1:A:165:GLN:NE2	2.07	0.41
2:B:20:LEU:H	2:B:20:LEU:CD1	2.29	0.41
1:A:96:THR:HG22	7:A:241:HOH:O	2.20	0.40
1:A:118:PRO:HG2	2:B:211:ARG:HH21	1.85	0.40
1:A:122:GLU:HG3	7:B:363:HOH:O	2.20	0.40
1:A:17:GLU:O	1:A:78:VAL:HG22	2.21	0.40

All (3) 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)
3:A:216:MG:MG	7:B:379:HOH:O[2_454]	0.79	1.41
7:A:269:HOH:O	7:B:410:HOH:O[3_555]	2.16	0.04
7:A:313:HOH:O	7:B:548:HOH:O[3_555]	2.17	0.03

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	214/219 (98%)	209 (98%)	5 (2%)	0	100	100
2	B	220/224 (98%)	212 (96%)	5 (2%)	3 (1%)	13	1
All	All	434/443 (98%)	421 (97%)	10 (2%)	3 (1%)	25	7

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	170	SER
2	B	127	ALA
2	B	126	SER

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	194/195 (100%)	188 (97%)	6 (3%)	45	15
2	B	192/192 (100%)	182 (95%)	10 (5%)	27	6
All	All	386/387 (100%)	370 (96%)	16 (4%)	34	10

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	61	ARG
1	A	93	ASN
1	A	135	LEU
1	A	146	LYS
1	A	159	LEU
1	A	211	ASN
2	B	1	GLU
2	B	20	LEU
2	B	50	PHE
2	B	82(A)	LEU
2	B	85	GLU
2	B	103	GLN
2	B	131	ASN

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Mol	Chain	Res	Type
2	B	133	MET
2	B	171	ASP
2	B	175	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) 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	13	GLN
2	B	71	ASN
2	B	79	GLN
2	B	96	HIS
2	B	103	GLN
2	B	169	GLN
2	B	194	ASN
2	B	197	HIS

5.3.3 RNA ⓘ

There are no RNA molecules 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 ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry

Of 4 ligands modelled in this entry, 2 are monoatomic - leaving 2 for Mogul analysis.

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
4	KO1	B	214	5	12,17,17	1.90	2 (16%)	13,26,26	1.92	3 (23%)
5	KDA	B	215	4	15,18,19	2.06	5 (33%)	16,25,27	2.34	3 (18%)

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
4	KO1	B	214	5	-	0/6/35/35	0/1/1/1
5	KDA	B	215	4	-	0/11/32/35	0/1/1/1

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	215	KDA	C4-C5	2.11	1.55	1.52
5	B	215	KDA	O6-C2	2.67	1.46	1.42
5	B	215	KDA	O2-C2	2.69	1.45	1.40
5	B	215	KDA	O6-C6	3.07	1.49	1.44
4	B	214	KO1	O2-C2	3.67	1.47	1.40
5	B	215	KDA	C3-C2	4.00	1.56	1.51
4	B	214	KO1	O6-C6	4.51	1.51	1.44

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	214	KO1	O8-C8-C7	-4.82	100.47	111.11
4	B	214	KO1	O7-C7-C6	-2.14	103.97	109.12
4	B	214	KO1	O7-C7-C8	2.78	115.63	109.21
5	B	215	KDA	O2-C2-C3	2.86	114.75	107.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	215	KDA	O6-C6-C7	3.67	113.12	106.65
5	B	215	KDA	C2-O6-C6	6.68	124.86	113.91

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 ⓘ

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%)	0.51	15 (6%) 18 16	20, 28, 38, 49	0
2	B	222/224 (99%)	0.70	26 (11%) 5 4	20, 28, 43, 57	0
All	All	439/443 (99%)	0.61	41 (9%) 9 8	20, 28, 40, 57	0

All (41) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	213	CYS	7.6
2	B	99	TYR	6.6
2	B	126	SER	5.6
2	B	128	ALA	5.6
2	B	131	ASN	5.0
2	B	127	ALA	4.6
2	B	132	SER	4.3
1	A	33	LEU	3.9
1	A	1	ASP	3.9
2	B	98	GLY	3.7
2	B	25	SER	3.7
2	B	2	VAL	3.7
2	B	170	SER	3.3
1	A	34	ALA	3.3
1	A	159	LEU	3.0
2	B	212	ASP	3.0
1	A	57	GLY	2.9
2	B	211	ARG	2.8
2	B	97	ASP	2.8
2	B	1	GLU	2.7
2	B	73	GLN	2.7
1	A	157	GLY	2.7
2	B	133	MET	2.6
1	A	30(B)	SER	2.5

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Mol	Chain	Res	Type	RSRZ
2	B	76	LEU	2.5
1	A	133	CYS	2.5
1	A	30(C)	ARG	2.5
2	B	90	TYR	2.5
1	A	202	SER	2.4
1	A	132	VAL	2.4
2	B	72[A]	SER	2.4
2	B	27	PHE	2.3
1	A	131	VAL	2.3
1	A	86	TYR	2.3
2	B	169	GLN	2.2
1	A	35	TRP	2.2
2	B	113	LYS	2.2
2	B	122	LEU	2.1
2	B	209	VAL	2.0
1	A	168	LYS	2.0
2	B	36	TRP	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(Å ²)	Q<0.9
3	MG	A	216	1/1	0.94	0.37	7.15	49,49,49,49	0
5	KDA	B	215	18/19	0.81	0.30	1.96	39,43,47,47	7
4	KO1	B	214	17/17	0.91	0.10	-0.43	28,30,38,43	0
6	ZN	B	217	1/1	0.98	0.05	-3.09	31,31,31,31	0

6.5 Other polymers

There are no such residues in this entry.