



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 02:00 PM UTC

PDB ID : 1FIZ / pdb_00001fiz
Title : THREE DIMENSIONAL STRUCTURE OF BETA-ACROSIN FROM BOAR SPERMATOZOA
Authors : Tranter, R.; Read, J.A.; Jones, R.; Brady, R.L.
Deposited on : 2000-08-07
Resolution : 2.90 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

<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
Mogul	:	2022.3.0, CSD as543be (2022)
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

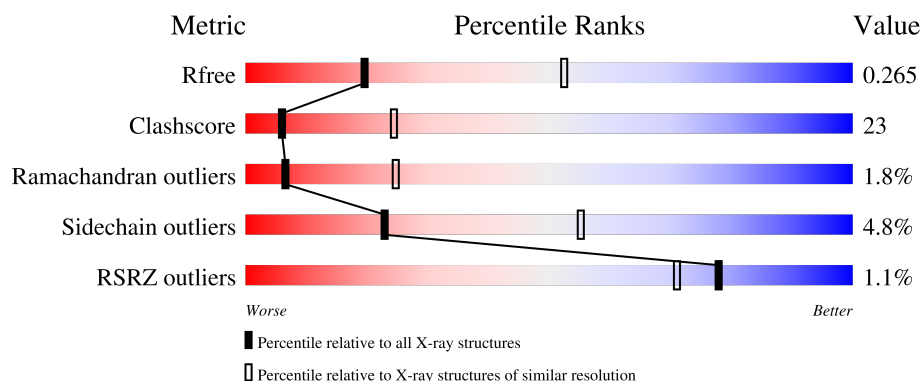
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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}	180053	2481 (2.90-2.90)
Clashscore	190562	2690 (2.90-2.90)
Ramachandran outliers	187476	2623 (2.90-2.90)
Sidechain outliers	187428	2625 (2.90-2.90)
RSRZ outliers	180081	2481 (2.90-2.90)

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 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	263	% 50% 38% 11%
2	L	23	4% 26% 22% 9% 43%
3	B	4	50% 50%

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	NAG	B	1	X	-	-	-
5	PBZ	A	308	-	X	-	-

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 2256 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.

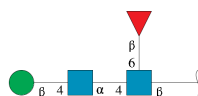
- Molecule 1 is a protein called BETA-ACROSIN HEAVY CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	263	Total	C	N	O	S	77	0	0
			2055	1317	367	356	15			

- Molecule 2 is a protein called BETA-ACROSIN LIGHT CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	L	13	Total	C	N	O	S	5	0	0
			97	58	20	17	2			

- Molecule 3 is an oligosaccharide called beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-alpha-D-glucopyranose-(1-4)-[beta-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose.



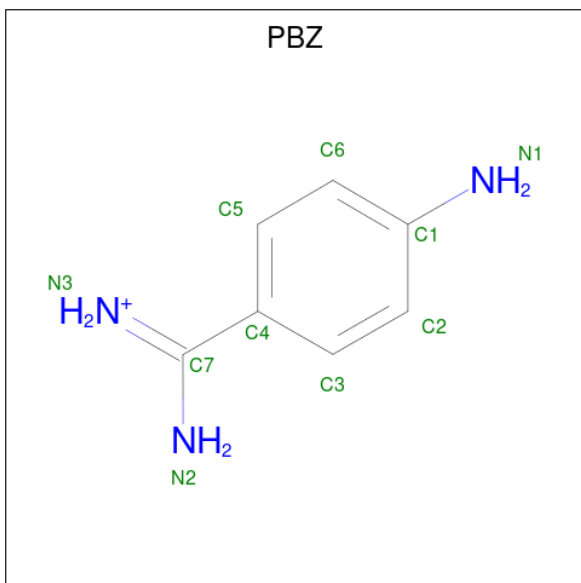
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	B	4	Total	C	N	O		0	0	0
			49	28	2	19				

- Molecule 4 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

- Molecule 5 is P-AMINO BENZAMIDINE (CCD ID: PBZ) (formula: $C_7H_{10}N_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	N	0	0
			10	7	3		

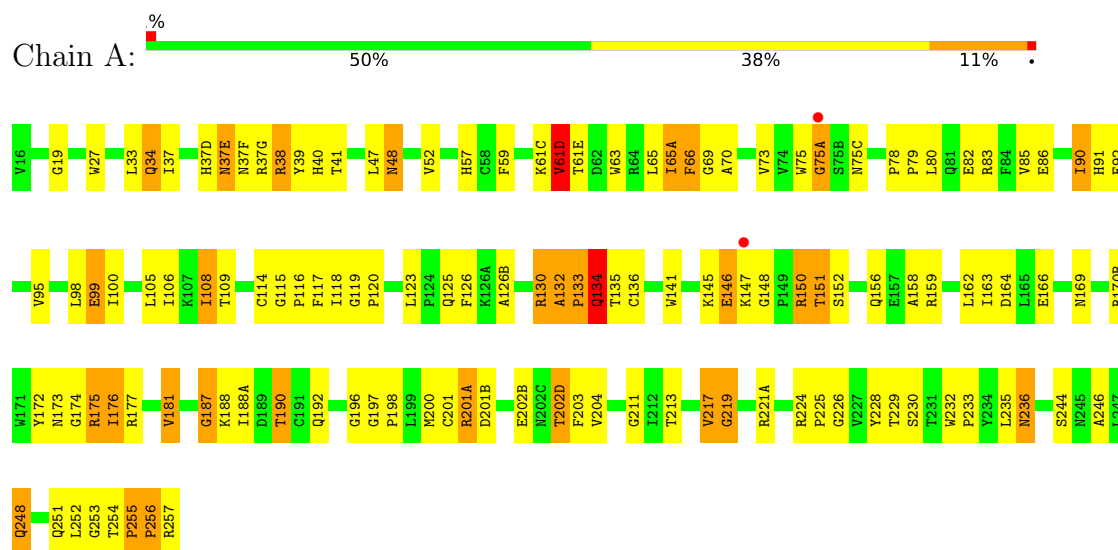
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	24	Total	O	0	0
			24	24		
6	L	1	Total	O	0	0
			1	1		

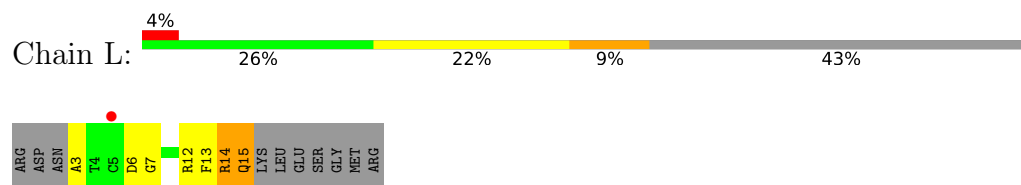
3 Residue-property plots

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: BETA-ACROSIN HEAVY CHAIN



• Molecule 2: BETA-ACROSIN LIGHT CHAIN



• Molecule 3: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-alpha-D-glucopyranose-(1-4)-[beta-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose



4 Data and refinement statistics

Property	Value	Source
Space group	P 43 3 2	Depositor
Cell constants a, b, c, α , β , γ	130.64Å 130.64Å 130.64Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.90 30.00 – 2.90	Depositor EDS
% Data completeness (in resolution range)	97.4 (30.00-2.90) 97.4 (30.00-2.90)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.54 (at 2.80Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.212 , 0.265 0.205 , 0.265	Depositor DCC
R_{free} test set	468 reflections (4.78%)	wwPDB-VP
Wilson B-factor (Å ²)	63.2	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 51.3	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	0.93	EDS
Total number of atoms	2256	wwPDB-VP
Average B, all atoms (Å ²)	49.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.46% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: NDG, PBZ, NAG, FUL, SO4, BMA

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	1.72	32/2118 (1.5%)	1.68	38/2886 (1.3%)
2	L	1.61	1/98 (1.0%)	1.58	1/130 (0.8%)
All	All	1.71	33/2216 (1.5%)	1.68	39/3016 (1.3%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

All (33) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	90	ILE	CA-CB	-11.71	1.39	1.54
1	A	135	THR	CA-CB	10.17	1.66	1.53
1	A	217	VAL	C-N	-9.96	1.18	1.33
1	A	256	PRO	C-O	-7.16	1.16	1.23
1	A	164	ASP	CA-C	-7.10	1.44	1.52
2	L	15	GLN	CA-CB	-7.03	1.39	1.53
1	A	251	GLN	C-O	-6.95	1.15	1.24
1	A	108	ILE	CA-CB	-6.91	1.45	1.54
1	A	132	ALA	CA-CB	-6.89	1.42	1.53
1	A	163	ILE	CA-CB	-6.87	1.46	1.54
1	A	123	LEU	C-O	-6.44	1.18	1.24
1	A	201(A)	ARG	C-O	-6.33	1.17	1.23
1	A	27	TRP	C-O	-6.23	1.17	1.24
1	A	158	ALA	CA-CB	-6.07	1.43	1.53
1	A	120	PRO	C-O	-5.88	1.16	1.23
1	A	225	PRO	C-O	-5.77	1.17	1.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	151	THR	CA-C	-5.71	1.45	1.53
1	A	134	GLN	CA-C	5.54	1.59	1.52
1	A	156	GLN	CA-CB	-5.48	1.45	1.53
1	A	188(A)	ILE	CA-CB	-5.44	1.47	1.54
1	A	130	ARG	CA-CB	-5.42	1.46	1.53
1	A	57	HIS	C-N	-5.40	1.27	1.33
1	A	152	SER	C-O	-5.33	1.17	1.24
1	A	134	GLN	CB-CG	-5.32	1.36	1.52
1	A	226	GLY	C-O	-5.28	1.18	1.23
1	A	38	ARG	C-O	-5.27	1.17	1.24
1	A	150	ARG	N-CA	5.25	1.52	1.46
1	A	65(A)	ILE	CA-CB	-5.25	1.47	1.54
1	A	48	ASN	CA-C	-5.22	1.46	1.52
1	A	159	ARG	CA-CB	5.21	1.60	1.53
1	A	66	PHE	C-N	-5.19	1.26	1.33
1	A	248	GLN	C-O	-5.09	1.17	1.24
1	A	61(D)	VAL	CA-CB	-5.07	1.47	1.54

All (39) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	130	ARG	O-C-N	-16.23	104.47	123.13
1	A	34	GLN	O-C-N	10.05	136.25	123.19
1	A	34	GLN	CA-C-N	-9.79	110.60	122.93
1	A	34	GLN	C-N-CA	-9.79	110.60	122.93
1	A	126(B)	ALA	N-CA-C	8.38	120.04	111.07
1	A	217	VAL	CA-C-N	8.12	133.16	120.65
1	A	217	VAL	C-N-CA	8.12	133.16	120.65
1	A	133	PRO	N-CD-CG	-7.63	94.64	103.80
1	A	217	VAL	O-C-N	-7.23	114.16	122.62
1	A	236	ASN	N-CA-CB	-7.14	98.42	110.49
2	L	14	ARG	N-CA-C	-6.96	104.66	113.01
1	A	196	GLY	N-CA-C	-6.80	106.08	114.92
1	A	37(F)	ASN	N-CA-C	6.64	120.71	112.47
1	A	181	VAL	CB-CA-C	-6.64	100.40	111.29
1	A	202(D)	THR	N-CA-C	6.61	120.46	109.95
1	A	86	GLU	N-CA-C	-6.52	105.45	113.41
1	A	150	ARG	N-CA-C	6.08	118.82	110.24
1	A	176	ILE	N-CA-C	6.01	118.01	108.87
1	A	158	ALA	CA-C-N	-5.86	113.92	122.19
1	A	158	ALA	C-N-CA	-5.86	113.92	122.19
1	A	198	PRO	N-CA-C	5.79	121.43	111.77

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	197	GLY	CA-C-O	-5.72	116.91	122.46
1	A	83	ARG	N-CA-C	5.57	117.26	108.96
1	A	187	GLY	N-CA-C	5.57	121.38	111.02
1	A	164	ASP	N-CA-C	5.38	117.24	110.24
1	A	246	ALA	N-CA-C	-5.36	105.34	111.07
1	A	146	GLU	N-CA-C	-5.34	102.63	110.42
1	A	63	TRP	N-CA-C	5.32	118.25	109.85
1	A	162	LEU	N-CA-CB	5.30	117.84	109.83
1	A	164	ASP	CA-C-N	-5.28	112.79	120.29
1	A	164	ASP	C-N-CA	-5.28	112.79	120.29
1	A	57	HIS	CA-C-O	5.23	125.78	119.61
1	A	166	GLU	N-CA-CB	5.21	118.38	110.14
1	A	123	LEU	O-C-N	-5.12	117.46	121.14
1	A	219	GLY	N-CA-C	-5.08	103.38	112.22
1	A	34	GLN	N-CA-C	5.07	117.83	109.72
1	A	255	PRO	CA-C-N	5.05	124.81	119.76
1	A	255	PRO	C-N-CA	5.05	124.81	119.76
1	A	90	ILE	CA-C-O	-5.05	115.38	121.54

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	130	ARG	Mainchain

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	2055	0	2041	92	0
2	L	97	0	90	6	0
3	B	49	0	42	3	0
4	A	20	0	0	0	0
5	A	10	0	10	1	0
6	A	24	0	0	1	0
6	L	1	0	0	0	0
All	All	2256	0	2183	96	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 23.

All (96) 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:147:LYS:HG3	1:A:148:GLY:H	1.08	1.14
1:A:37:ILE:HD12	1:A:41:THR:HG21	1.46	0.97
1:A:147:LYS:HG3	1:A:148:GLY:N	1.83	0.92
1:A:169:ASN:OD1	1:A:174:GLY:N	2.20	0.73
1:A:59:PHE:HZ	1:A:106:ILE:HD11	1.54	0.71
1:A:61(D):VAL:HG13	1:A:85:VAL:O	1.90	0.71
1:A:37(G):ARG:HB3	1:A:39:TYR:CE1	2.26	0.70
1:A:114:CYS:SG	1:A:119:GLY:HA2	2.30	0.70
1:A:172:TYR:HE1	1:A:217:VAL:HG23	1.57	0.69
1:A:169:ASN:OD1	1:A:174:GLY:CA	2.40	0.68
1:A:98:LEU:O	1:A:99:GLU:HB2	1.94	0.67
1:A:95:VAL:HB	1:A:100:ILE:HB	1.77	0.65
1:A:169:ASN:OD1	1:A:174:GLY:HA2	1.94	0.65
1:A:117:PHE:C	1:A:118:ILE:HG13	2.20	0.64
1:A:145:LYS:O	1:A:146:GLU:C	2.37	0.64
1:A:190:THR:OG1	5:A:308:PBZ:N2	2.32	0.63
1:A:69:GLY:O	1:A:80:LEU:HD23	1.99	0.63
1:A:116:PRO:HA	2:L:14:ARG:HB2	1.81	0.63
1:A:253:GLY:O	1:A:255:PRO:HD3	1.99	0.62
1:A:75:TRP:O	1:A:75(A):GLY:C	2.42	0.62
1:A:52:VAL:CG2	1:A:108:ILE:HD11	2.30	0.62
1:A:66:PHE:CD1	1:A:66:PHE:N	2.68	0.60
1:A:172:TYR:CD2	1:A:176:ILE:HD11	2.37	0.60
1:A:147:LYS:CG	1:A:148:GLY:H	2.00	0.59
1:A:48:ASN:OD1	1:A:48:ASN:C	2.45	0.59
1:A:114:CYS:O	2:L:3:ALA:HB1	2.02	0.59
1:A:132:ALA:HB1	1:A:133:PRO:HA	1.86	0.58
1:A:146:GLU:OE1	1:A:219:GLY:HA3	2.03	0.58
1:A:90:ILE:HG22	1:A:91:HIS:N	2.15	0.58
1:A:37(G):ARG:HB3	1:A:39:TYR:HE1	1.68	0.57
1:A:38:ARG:O	1:A:75(A):GLY:HA2	2.05	0.56
1:A:177:ARG:HA	3:B:1:NAG:H83	1.87	0.56
1:A:115:GLY:O	2:L:14:ARG:HD3	2.05	0.56
1:A:117:PHE:O	1:A:118:ILE:HG13	2.06	0.56
1:A:201(A):ARG:HG2	1:A:203:PHE:CE2	2.41	0.55
1:A:52:VAL:HG23	1:A:108:ILE:HD11	1.88	0.55
1:A:66:PHE:CZ	1:A:108:ILE:HD13	2.42	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:70:ALA:HA	1:A:80:LEU:CD2	2.38	0.54
1:A:99:GLU:HA	1:A:99:GLU:OE1	2.10	0.52
1:A:134:GLN:OE1	1:A:201:CYS:HB3	2.09	0.52
1:A:253:GLY:C	1:A:255:PRO:HD3	2.35	0.52
1:A:61(C):LYS:O	1:A:61(E):THR:N	2.42	0.52
1:A:37(G):ARG:NH1	1:A:75(C):ASN:HD21	2.08	0.51
3:B:3:BMA:H62	3:B:3:BMA:O2	2.11	0.51
1:A:125:GLN:NE2	1:A:252:LEU:O	2.44	0.51
1:A:141:TRP:O	1:A:151:THR:HB	2.11	0.50
1:A:187:GLY:O	1:A:188:LYS:HB2	2.11	0.50
1:A:59:PHE:CD1	1:A:59:PHE:N	2.79	0.50
1:A:170(B):ARG:NH2	6:A:320:HOH:O	2.46	0.49
1:A:34:GLN:HG2	1:A:40:HIS:HA	1.93	0.49
1:A:65:LEU:O	1:A:82:GLU:HA	2.13	0.49
1:A:181:VAL:HG23	1:A:230:SER:HB2	1.94	0.49
2:L:12:ARG:O	2:L:13:PHE:C	2.54	0.49
2:L:13:PHE:C	2:L:15:GLN:N	2.66	0.48
2:L:6:ASP:CG	2:L:7:GLY:H	2.21	0.48
1:A:92:GLU:H	1:A:92:GLU:CD	2.22	0.48
1:A:78:PRO:HA	1:A:79:PRO:C	2.38	0.48
1:A:146:GLU:OE1	1:A:221(A):ARG:NE	2.36	0.47
1:A:147:LYS:CG	1:A:148:GLY:N	2.66	0.46
1:A:202(D):THR:HG22	1:A:203:PHE:N	2.30	0.46
1:A:37(G):ARG:NH1	1:A:75(C):ASN:ND2	2.64	0.46
1:A:201(B):ASP:HB2	1:A:202(B):GLU:HG3	1.98	0.45
1:A:73:VAL:HG23	1:A:141:TRP:NE1	2.32	0.45
1:A:172:TYR:O	1:A:173:ASN:C	2.60	0.45
1:A:37(E):ASN:ND2	1:A:37(G):ARG:HB2	2.32	0.45
1:A:33:LEU:CD2	1:A:65:LEU:HD22	2.47	0.45
1:A:213:THR:HA	1:A:228:TYR:CD2	2.52	0.44
1:A:136:CYS:HA	1:A:200:MET:O	2.18	0.44
1:A:201(A):ARG:HG2	1:A:203:PHE:CD2	2.52	0.44
1:A:69:GLY:HA3	1:A:118:ILE:HD11	2.00	0.43
1:A:52:VAL:HG21	1:A:108:ILE:HD11	2.00	0.43
1:A:61(D):VAL:CG1	1:A:85:VAL:O	2.63	0.43
1:A:47:LEU:HA	1:A:47:LEU:HD23	1.65	0.43
1:A:230:SER:O	1:A:233:PRO:HD2	2.19	0.43
1:A:211:GLY:HA2	1:A:229:THR:O	2.18	0.43
1:A:187:GLY:O	1:A:188:LYS:CB	2.62	0.42
1:A:200:MET:HE3	1:A:200:MET:HB2	1.87	0.42
1:A:37(D):HIS:O	1:A:37(E):ASN:HB3	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:105:LEU:HD23	1:A:105:LEU:HA	1.88	0.42
1:A:172:TYR:HB2	1:A:176:ILE:CD1	2.50	0.41
1:A:177:ARG:CA	3:B:1:NAG:H83	2.50	0.41
1:A:59:PHE:CZ	1:A:106:ILE:HD11	2.45	0.41
1:A:65:LEU:O	1:A:65(A):ILE:HD13	2.20	0.41
1:A:146:GLU:O	1:A:147:LYS:HB3	2.20	0.41
1:A:33:LEU:HD22	1:A:65:LEU:HD22	2.02	0.41
1:A:172:TYR:O	1:A:175:ARG:HD3	2.20	0.41
1:A:200:MET:HA	1:A:204:VAL:O	2.20	0.41
1:A:109:THR:O	1:A:109:THR:HG23	2.20	0.41
1:A:255:PRO:CB	1:A:256:PRO:CD	2.98	0.41
1:A:66:PHE:CZ	1:A:108:ILE:CD1	3.03	0.41
1:A:125:GLN:O	1:A:126:PHE:C	2.62	0.41
1:A:66:PHE:HZ	1:A:108:ILE:CD1	2.34	0.40
1:A:232:TRP:O	1:A:235:LEU:HG	2.21	0.40
1:A:61(C):LYS:O	1:A:61(D):VAL:C	2.63	0.40
1:A:66:PHE:HZ	1:A:108:ILE:HD13	1.86	0.40
1:A:33:LEU:HD22	1:A:65:LEU:CD2	2.52	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	261/263 (99%)	227 (87%)	29 (11%)	5 (2%)	6	23
2	L	11/23 (48%)	9 (82%)	2 (18%)	0	100	100
All	All	272/286 (95%)	236 (87%)	31 (11%)	5 (2%)	6	25

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	37(E)	ASN
1	A	61(D)	VAL
1	A	75(A)	GLY
1	A	236	ASN
1	A	19	GLY

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	220/220 (100%)	209 (95%)	11 (5%)	22	53
2	L	10/19 (53%)	10 (100%)	0	100	100
All	All	230/239 (96%)	219 (95%)	11 (5%)	23	55

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	99	GLU
1	A	134	GLN
1	A	150	ARG
1	A	175	ARG
1	A	190	THR
1	A	192	GLN
1	A	224	ARG
1	A	244	SER
1	A	248	GLN
1	A	254	THR
1	A	257	ARG

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

Mol	Chain	Res	Type
1	A	71	ASN
1	A	75(C)	ASN
1	A	125	GLN

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Mol	Chain	Res	Type
1	A	192	GLN
1	A	245	ASN
1	A	248	GLN

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 ⓘ

4 monosaccharides 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	NAG	B	1	1,3	14,14,15	1.80	3 (21%)	17,19,21	3.18	4 (23%)
3	NDG	B	2	3	14,14,15	0.70	0	17,19,21	1.60	5 (29%)
3	BMA	B	3	3	11,11,12	0.54	0	15,15,17	2.25	3 (20%)
3	FUL	B	4	3	10,10,11	0.80	0	14,14,16	2.15	6 (42%)

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	NAG	B	1	1,3	1/1/5/7	2/6/23/26	0/1/1/1
3	NDG	B	2	3	-	2/6/23/26	0/1/1/1
3	BMA	B	3	3	-	1/2/19/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	FUL	B	4	3	-	-	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	1	NAG	O5-C1	-5.23	1.34	1.43
3	B	1	NAG	O3-C3	-2.04	1.37	1.43
3	B	1	NAG	C1-C2	2.01	1.55	1.52

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	1	NAG	O5-C1-C2	-9.35	96.83	111.29
3	B	1	NAG	C3-C4-C5	-7.04	97.47	110.23
3	B	3	BMA	C1-C2-C3	6.22	118.71	109.64
3	B	3	BMA	C1-O5-C5	4.47	118.17	112.19
3	B	1	NAG	O5-C5-C4	4.19	121.02	110.83
3	B	4	FUL	C1-C2-C3	4.08	115.59	109.64
3	B	2	NDG	C4-C3-C2	-3.27	106.23	111.02
3	B	4	FUL	O2-C2-C3	-3.16	103.60	110.15
3	B	2	NDG	O5-C5-C6	3.11	113.71	107.66
3	B	4	FUL	C1-O5-C5	3.03	120.12	112.97
3	B	2	NDG	O4-C4-C3	2.89	117.20	110.38
3	B	4	FUL	O3-C3-C4	-2.75	103.89	110.38
3	B	4	FUL	O5-C1-C2	2.57	116.92	110.79
3	B	1	NAG	O4-C4-C3	-2.47	104.54	110.38
3	B	3	BMA	C3-C4-C5	2.45	114.67	110.23
3	B	4	FUL	C3-C4-C5	2.43	113.51	109.81
3	B	2	NDG	O7-C7-C8	-2.12	118.27	122.05
3	B	2	NDG	C8-C7-N2	2.08	119.57	116.12

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	B	1	NAG	C1

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	1	NAG	C4-C5-C6-O6
3	B	2	NDG	O7-C7-N2-C2

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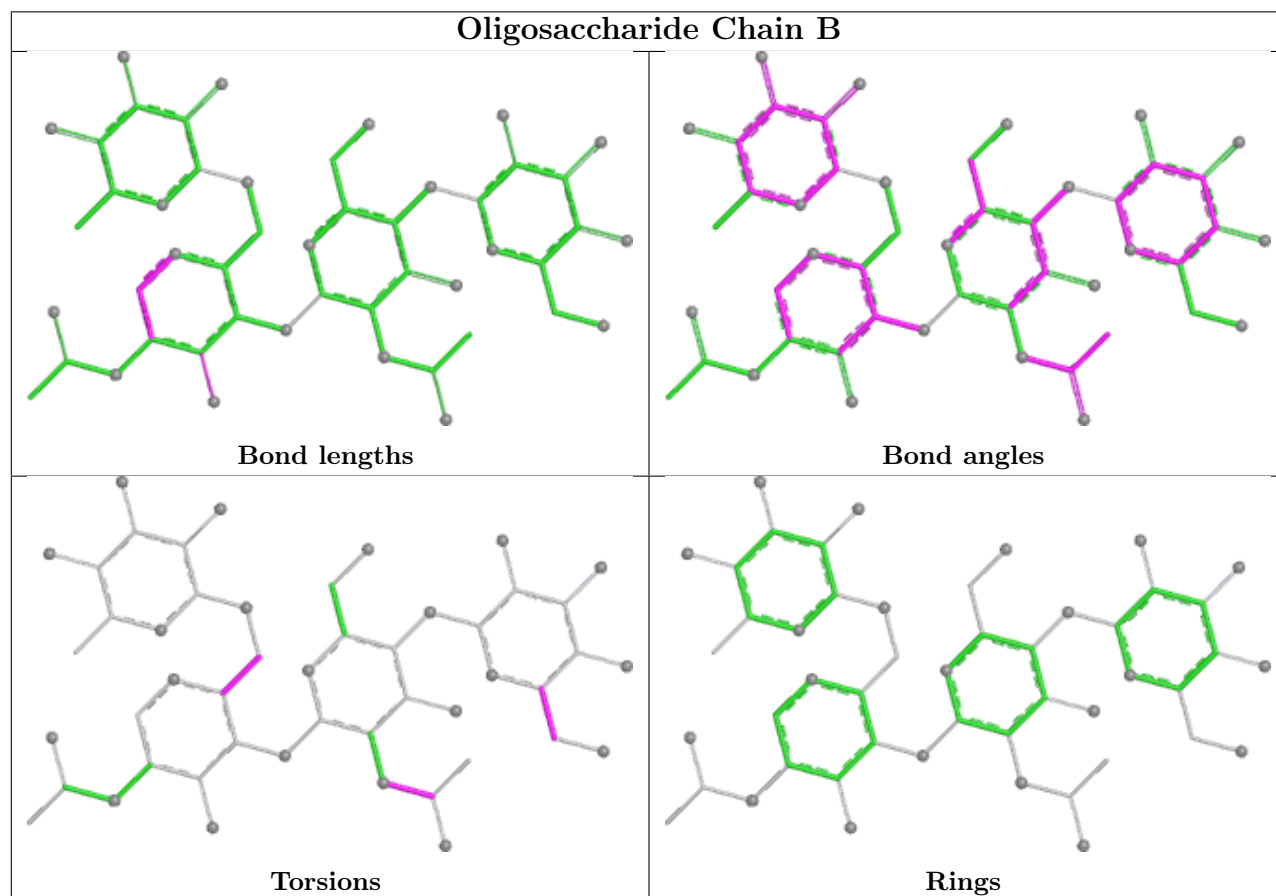
Mol	Chain	Res	Type	Atoms
3	B	1	NAG	O5-C5-C6-O6
3	B	2	NDG	C8-C7-N2-C2
3	B	3	BMA	O5-C5-C6-O6

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	3	BMA	1	0
3	B	1	NAG	2	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.



5.6 Ligand geometry [i](#)

5 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$
4	SO4	A	306	-	4,4,4	0.45	0	6,6,6	1.16	1 (16%)
4	SO4	A	304	-	4,4,4	0.23	0	6,6,6	0.35	0
5	PBZ	A	308	-	10,10,10	1.72	3 (30%)	9,13,13	3.41	5 (55%)
4	SO4	A	305	-	4,4,4	0.72	0	6,6,6	0.43	0
4	SO4	A	307	-	4,4,4	0.28	0	6,6,6	0.70	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	PBZ	A	308	-	-	4/4/4/4	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	308	PBZ	C6-C5	2.78	1.43	1.38
5	A	308	PBZ	C4-C7	-2.49	1.42	1.47
5	A	308	PBZ	C5-C4	-2.14	1.36	1.39

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	308	PBZ	C2-C3-C4	5.96	127.17	120.80
5	A	308	PBZ	C6-C5-C4	-4.87	115.60	120.80
5	A	308	PBZ	C5-C6-C1	4.06	125.62	120.66
5	A	308	PBZ	C3-C2-C1	-3.63	116.23	120.66
5	A	308	PBZ	C4-C7-N2	-3.08	113.34	118.01
4	A	306	SO4	O3-S-O2	2.16	120.83	109.56

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	308	PBZ	C3-C4-C7-N2
5	A	308	PBZ	C5-C4-C7-N2
5	A	308	PBZ	C3-C4-C7-N3
5	A	308	PBZ	C5-C4-C7-N3

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	308	PBZ	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	A	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	A	217:VAL	C	219:GLY	N	1.18

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	263/263 (100%)	-0.23	2 (0%) 82 77	16, 44, 69, 77	15 (5%)
2	L	13/23 (56%)	0.82	1 (7%) 19 16	31, 60, 92, 93	1 (7%)
All	All	276/286 (96%)	-0.18	3 (1%) 78 71	16, 44, 73, 93	16 (5%)

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	147	LYS	2.6
2	L	5	CYS	2.4
1	A	75(A)	GLY	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

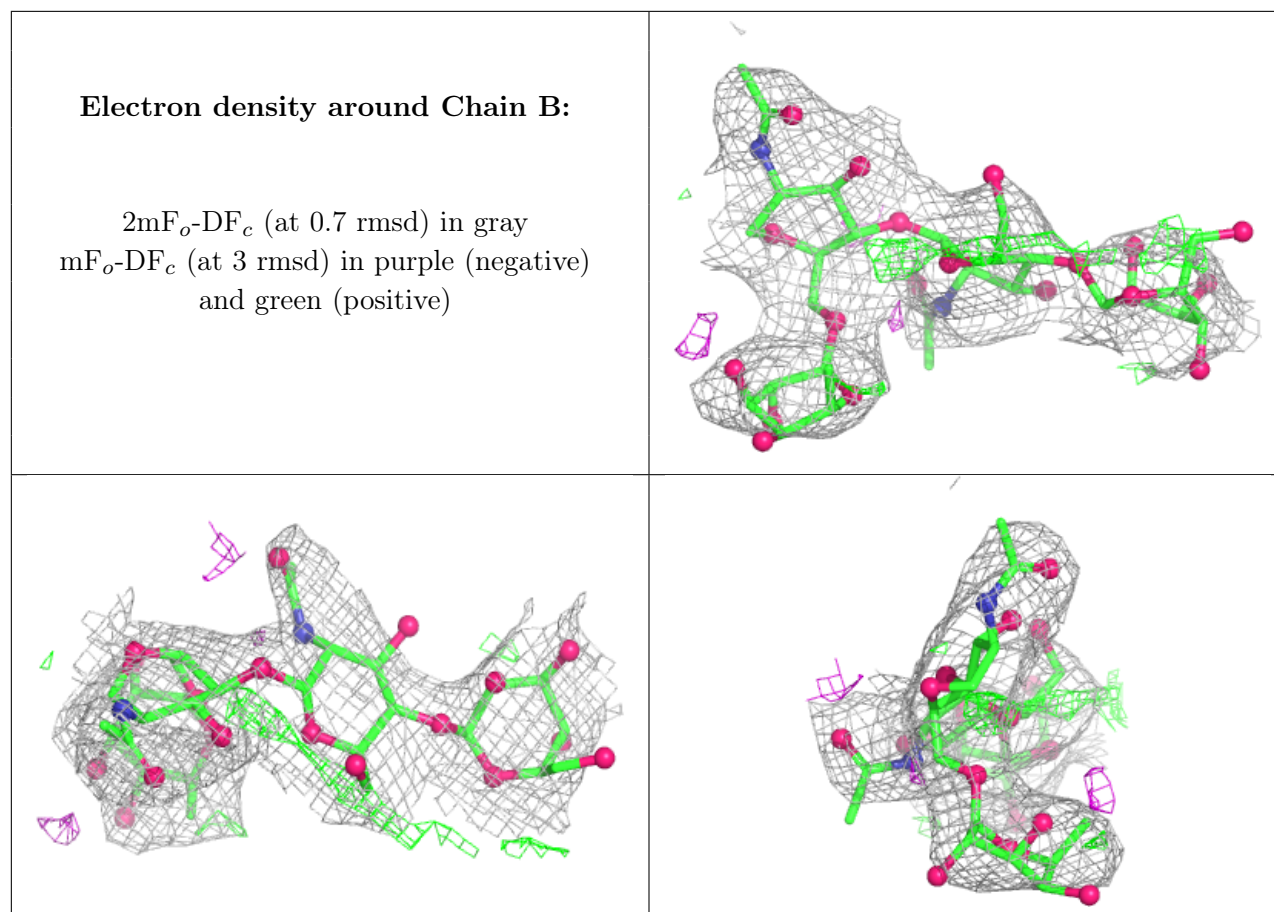
There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

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
3	NAG	B	1	14/15	-	-	54,65,79,82	0
3	NDG	B	2	14/15	-	-	93,103,108,116	0
3	BMA	B	3	11/12	-	-	121,123,126,127	0
3	FUL	B	4	10/11	-	-	84,88,91,92	0

The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.



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($\text{\AA}^2$)	Q<0.9
4	SO4	A	306	5/5	0.82	0.14	62,62,67,67	0
4	SO4	A	307	5/5	0.95	0.09	68,68,71,75	0
5	PBZ	A	308	10/10	0.96	0.11	38,40,42,45	0
4	SO4	A	305	5/5	0.97	0.06	45,49,50,55	0
4	SO4	A	304	5/5	0.98	0.07	37,39,40,41	0

6.5 Other polymers [i](#)

There are no such residues in this entry.