



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 15, 2026 – 11:48 AM UTC

PDB ID : 1Q9W / pdb_00001q9w
Title : S45-18 Fab pentasaccharide bisphosphate complex
Authors : Nguyen, H.P.; Seto, N.O.; MacKenzie, C.R.; Brade, L.; Kosma, P.; Brade, H.; Evans, S.V.
Deposited on : 2003-08-26
Resolution : 1.75 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

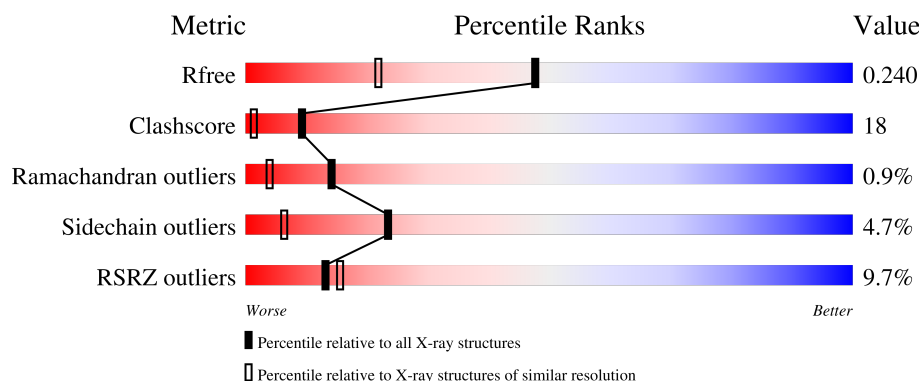
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.75 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	3183 (1.76-1.76)
Clashscore	190562	3299 (1.76-1.76)
Ramachandran outliers	187476	3274 (1.76-1.76)
Sidechain outliers	187428	3274 (1.76-1.76)
RSRZ outliers	180081	3183 (1.76-1.76)

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	219	8% 72% 25% .
1	C	219	7% 68% 26% 5%
2	B	226	13% 70% 27% .
2	D	226	10% 67% 29% .
3	E	5	40% 60%

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Mol	Chain	Length	Quality of chain
4	F	3	33% 67%

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	GP4	E	2	X	-	-	-

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 7744 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.

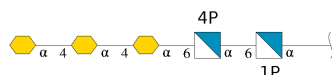
- Molecule 1 is a protein called S45-18 Fab (IgG1k) light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	219	Total	C	N	O	S	0	0	0
			1707	1066	289	344	8			
1	C	219	Total	C	N	O	S	0	0	0
			1707	1066	289	344	8			

- Molecule 2 is a protein called S45-18 Fab (IgG1k) heavy chain.

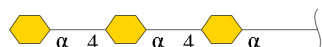
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	226	Total	C	N	O	S	0	0	0
			1718	1085	284	340	9			
2	D	226	Total	C	N	O	S	0	0	0
			1718	1085	284	340	9			

- Molecule 3 is an oligosaccharide called 3-deoxy-alpha-D-manno-oct-2-ulopyranosonic acid-(2-4)-3-deoxy-alpha-D-manno-oct-2-ulopyranosonic acid-(2-4)-3-deoxy-alpha-D-manno-oct-2-ulopyranosonic acid-(2-6)-2-amino-2-deoxy-4-O-phosphono-alpha-D-glucopyranose-(1-6)-2-amino-2-deoxy-1-O-phosphono-alpha-D-glucopyranose.



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	E	5	Total	C	N	O	P	0	0	0
			76	36	2	36	2			

- Molecule 4 is an oligosaccharide called 3-deoxy-alpha-D-manno-oct-2-ulopyranosonic acid-(2-4)-3-deoxy-alpha-D-manno-oct-2-ulopyranosonic acid-(2-4)-3-deoxy-alpha-D-manno-oct-2-ulopyranosonic acid.



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
4	F	3	Total	C	O	0	0	0
			46	24	22			

- Molecule 5 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	2	Total	Mg	0	0
			2	2		
5	B	1	Total	Mg	0	0
			1	1		
5	C	2	Total	Mg	0	0
			2	2		

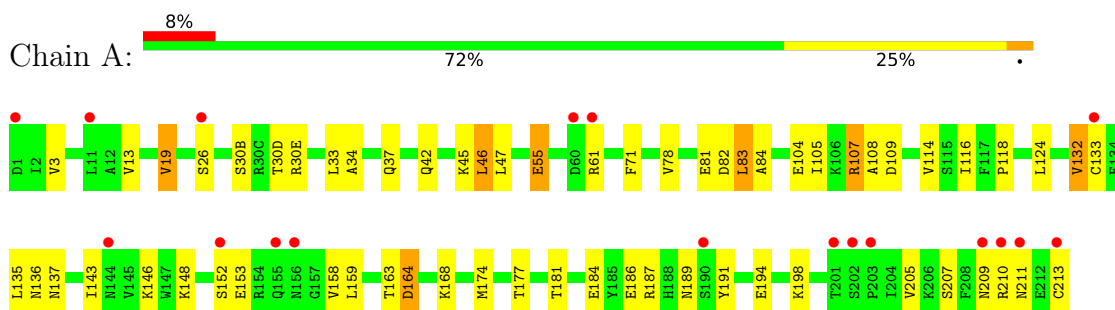
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	160	Total	O	0	0
			160	160		
6	B	220	Total	O	0	0
			220	220		
6	C	176	Total	O	0	0
			176	176		
6	D	211	Total	O	0	0
			211	211		

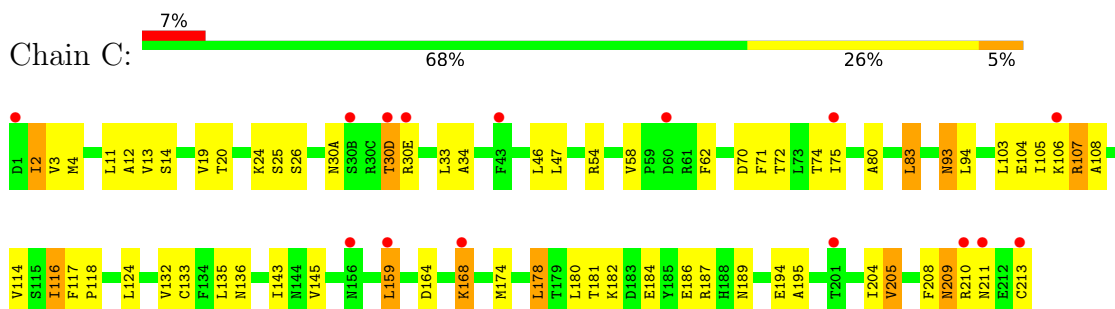
3 Residue-property plots [i](#)

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.

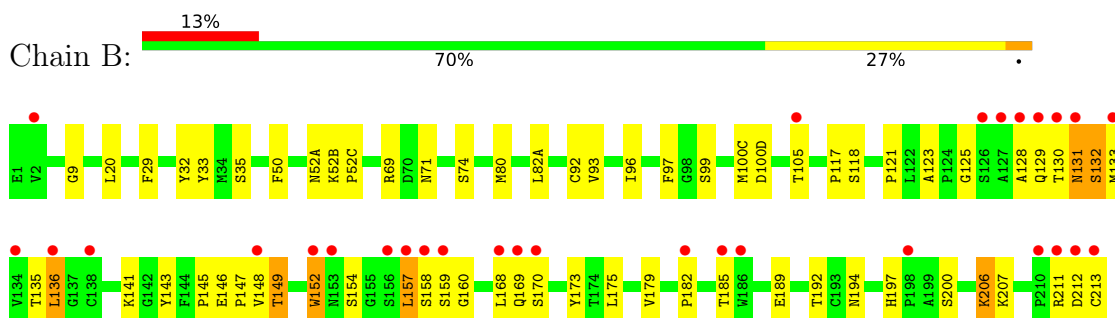
- Molecule 1: S45-18 Fab (IgG1k) light chain



- Molecule 1: S45-18 Fab (IgG1k) light chain

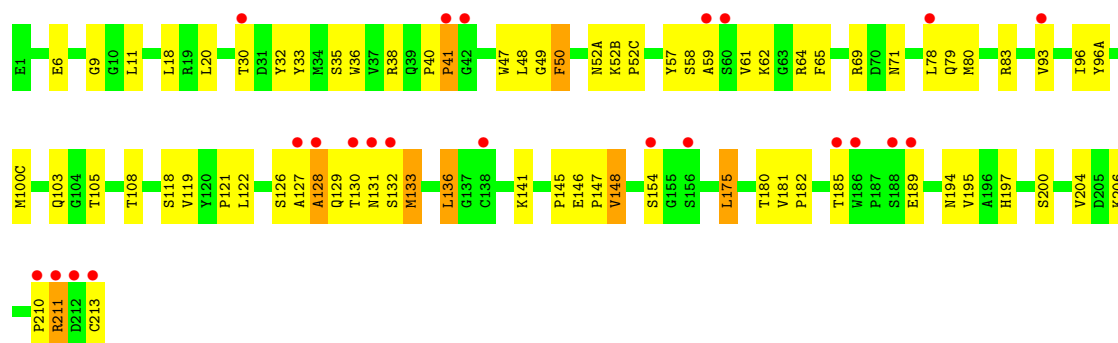


- Molecule 2: S45-18 Fab (IgG1k) heavy chain



- Molecule 2: S45-18 Fab (IgG1k) heavy chain





- Molecule 3: 3-deoxy-alpha-D-manno-oct-2-ulopyranosonic acid-(2-4)-3-deoxy-alpha-D-manno-oct-2-ulopyranosonic acid-(2-4)-3-deoxy-alpha-D-manno-oct-2-ulopyranosonic acid-(2-6)-2-amino-2-deoxy-4-O-phosphono-alpha-D-glucopyranose-(1-6)-2-amino-2-deoxy-1-O-phosphono-alpha-D-glucopyranose

Chain E: 40% 60%



- Molecule 4: 3-deoxy-alpha-D-manno-oct-2-ulopyranosonic acid-(2-4)-3-deoxy-alpha-D-manno-oct-2-ulopyranosonic acid-(2-4)-3-deoxy-alpha-D-manno-oct-2-ulopyranosonic acid

Chain F: 33% 67%



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	71.20Å 113.90Å 133.90Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.90 – 1.75 19.90 – 1.75	Depositor EDS
% Data completeness (in resolution range)	90.5 (19.90-1.75) 90.4 (19.90-1.75)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.33 (at 1.74Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.212 , 0.247 0.205 , 0.240	Depositor DCC
R_{free} test set	9994 reflections (9.42%)	wwPDB-VP
Wilson B-factor (Å ²)	23.4	Xtriage
Anisotropy	0.450	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 45.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7744	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 6.83% 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: KDO, GP4, GP1, MG

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	0.47	0/1744	0.97	5/2361 (0.2%)
1	C	0.49	0/1744	0.97	8/2361 (0.3%)
2	B	0.52	0/1763	1.06	12/2406 (0.5%)
2	D	0.52	0/1763	1.03	10/2406 (0.4%)
All	All	0.50	0/7014	1.01	35/9534 (0.4%)

There are no bond length outliers.

All (35) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	146	GLU	N-CA-C	-8.09	98.25	110.39
1	C	30(D)	THR	N-CA-C	-7.58	98.23	110.20
2	D	33	TYR	N-CA-C	-7.47	99.96	110.50
2	B	33	TYR	N-CA-C	-7.35	99.61	110.48
2	D	32	TYR	N-CA-C	7.27	121.30	109.59
2	D	118	SER	N-CA-C	-6.98	97.83	109.07
2	B	32	TYR	N-CA-C	6.93	120.75	109.59
1	C	136	ASN	N-CA-C	6.73	120.65	109.95
1	A	136	ASN	N-CA-C	6.57	120.39	109.95
2	B	99	SER	N-CA-C	6.36	118.01	111.14
2	D	141	LYS	N-CA-C	6.33	119.84	109.72
2	B	146	GLU	N-CA-C	-6.25	101.01	110.39
2	D	96	ILE	N-CA-C	6.14	117.15	110.21
1	A	164	ASP	N-CA-C	-6.08	102.69	110.53
1	C	209	ASN	N-CA-C	5.77	120.60	113.50
1	C	164	ASP	N-CA-C	-5.62	102.96	110.55
2	B	152	TRP	CA-C-N	-5.58	113.70	122.08
2	B	152	TRP	C-N-CA	-5.58	113.70	122.08
1	A	181	THR	N-CA-C	-5.58	102.27	110.52
1	C	204	ILE	N-CA-C	-5.52	99.78	108.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	205	VAL	N-CA-C	5.52	115.89	108.17
2	B	96	ILE	CB-CA-C	-5.39	105.42	111.45
2	B	118	SER	N-CA-C	-5.37	100.42	109.07
2	D	148	VAL	N-CA-C	-5.35	100.48	108.46
2	B	160	GLY	N-CA-C	-5.31	108.30	115.36
1	C	181	THR	N-CA-C	-5.30	103.39	110.55
2	B	141	LYS	N-CA-C	5.26	117.81	109.50
2	D	96	ILE	CB-CA-C	-5.21	105.61	111.45
1	A	137	ASN	N-CA-C	5.16	118.36	111.24
1	C	2	ILE	N-CA-C	5.14	116.31	109.37
2	D	40	PRO	N-CA-C	-5.13	105.54	110.47
2	B	92	CYS	N-CA-C	-5.12	100.97	109.46
2	B	100(D)	ASP	N-CA-C	5.08	119.18	112.89
1	A	132	VAL	N-CA-C	5.06	115.77	108.48
2	D	96(A)	TYR	N-CA-C	-5.00	107.86	114.31

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	1707	0	1653	64	0
1	C	1707	0	1653	77	0
2	B	1718	0	1671	43	0
2	D	1718	0	1671	70	0
3	E	76	0	34	3	0
4	F	46	0	35	1	0
5	A	2	0	0	0	0
5	B	1	0	0	0	0
5	C	2	0	0	0	0
6	A	160	0	0	3	0
6	B	220	0	0	1	0
6	C	176	0	0	3	0
6	D	211	0	0	4	0
All	All	7744	0	6717	248	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 18.

All (248) 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)
2:D:154:SER:H	2:D:194:ASN:HD21	1.04	0.99
1:C:114:VAL:HG22	1:C:135:LEU:HD13	1.45	0.98
1:C:168:LYS:HZ3	1:C:168:LYS:HB2	1.39	0.88
2:D:105:THR:HG23	6:D:377:HOH:O	1.76	0.85
1:C:210:ARG:HA	1:C:213:CYS:SG	2.15	0.85
1:C:2:ILE:HG22	1:C:4:MET:CE	2.06	0.84
1:C:116:ILE:HD13	1:C:117:PHE:N	1.92	0.84
2:D:154:SER:H	2:D:194:ASN:ND2	1.76	0.82
1:A:194:GLU:HG2	1:A:205:VAL:HG12	1.61	0.81
2:B:69:ARG:HE	2:B:71:ASN:HD21	1.30	0.79
1:A:210:ARG:HA	1:A:213:CYS:SG	2.22	0.78
1:C:30(E):ARG:H	1:C:30(E):ARG:HD2	1.48	0.78
1:C:117:PHE:HB2	1:C:132:VAL:HG13	1.65	0.77
2:D:48:LEU:HD22	2:D:61:VAL:HG11	1.65	0.77
1:A:163:THR:HG23	6:A:269:HOH:O	1.83	0.77
2:D:119:VAL:HG21	2:D:204:VAL:HG11	1.67	0.77
1:C:2:ILE:HG22	1:C:4:MET:HE3	1.65	0.76
2:B:154:SER:H	2:B:194:ASN:HD21	1.34	0.76
2:B:168:LEU:HB2	2:B:173:TYR:CE1	2.20	0.75
1:C:30(A):ASN:O	1:C:30(D):THR:O	2.06	0.74
2:D:9:GLY:H	2:D:105:THR:HG21	1.52	0.73
1:C:189:ASN:ND2	1:C:211:ASN:HB3	2.04	0.72
1:A:159:LEU:HD21	2:B:169:GLN:NE2	2.05	0.72
1:C:132:VAL:HG11	2:D:122:LEU:HD13	1.71	0.72
2:B:9:GLY:HA3	2:B:105:THR:HG22	1.72	0.71
2:D:30:THR:OG1	2:D:52(B):LYS:HB2	1.90	0.71
1:C:3:VAL:C	1:C:4:MET:HE2	2.16	0.70
2:B:125:GLY:HA2	2:B:212:ASP:O	1.91	0.70
2:B:157:LEU:HD23	2:B:179:VAL:HG21	1.73	0.69
1:C:3:VAL:O	1:C:4:MET:HE2	1.91	0.69
2:B:136:LEU:HD22	2:B:136:LEU:N	2.08	0.68
1:C:213:CYS:SG	2:D:213:CYS:C	2.76	0.68
1:A:118:PRO:HG3	2:B:213:CYS:HB2	1.74	0.68
1:C:80:ALA:HA	1:C:105:ILE:HD13	1.76	0.68
2:B:9:GLY:H	2:B:105:THR:HG21	1.59	0.68
1:A:189:ASN:ND2	1:A:211:ASN:HB3	2.09	0.68
2:D:69:ARG:HE	2:D:71:ASN:HD21	1.41	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:83:LEU:HD23	1:A:105:ILE:HD13	1.78	0.66
1:A:13:VAL:HG11	1:A:19:VAL:HG22	1.78	0.66
1:A:184:GLU:HA	1:A:187:ARG:NH1	2.11	0.65
2:D:65:PHE:HD1	2:D:78:LEU:HD11	1.60	0.65
1:C:13:VAL:HG11	1:C:19:VAL:HG22	1.78	0.65
1:C:62:PHE:CD2	1:C:75:ILE:HD12	2.31	0.65
1:C:117:PHE:HB2	1:C:132:VAL:CG1	2.27	0.65
2:D:103:GLN:OE1	6:D:259:HOH:O	2.14	0.64
1:A:46:LEU:HD13	1:A:55:GLU:HG2	1.77	0.64
1:C:107:ARG:HD3	1:C:108:ALA:O	1.97	0.64
1:C:12:ALA:HB2	1:C:104:GLU:HG3	1.80	0.63
2:D:79:GLN:HG2	6:D:233:HOH:O	1.98	0.63
1:C:209:ASN:O	1:C:210:ARG:HB3	1.96	0.63
1:A:61:ARG:HD3	6:A:370:HOH:O	1.99	0.63
1:C:135:LEU:HD11	1:C:195:ALA:HB2	1.80	0.63
2:D:93:VAL:CG2	2:D:100(C):MET:HB3	2.28	0.63
1:C:168:LYS:HZ3	1:C:168:LYS:CB	2.12	0.62
1:A:163:THR:HG22	1:A:164:ASP:O	1.98	0.62
2:D:133:MET:HE3	2:D:180:THR:HG22	1.80	0.62
1:A:107:ARG:HD3	1:A:108:ALA:O	2.00	0.62
1:A:163:THR:HG22	1:A:164:ASP:N	2.15	0.62
2:D:35:SER:HB2	2:D:93:VAL:CG1	2.30	0.62
1:C:116:ILE:HG12	1:C:133:CYS:SG	2.40	0.62
2:D:93:VAL:HG21	2:D:100(C):MET:HB3	1.81	0.61
1:A:61:ARG:NH2	1:A:82:ASP:OD1	2.33	0.61
2:D:119:VAL:HG21	2:D:204:VAL:CG1	2.30	0.61
1:A:114:VAL:HG22	1:A:135:LEU:CD2	2.31	0.61
2:B:93:VAL:HG21	2:B:100(C):MET:HB3	1.83	0.61
1:A:146:LYS:HE2	1:A:153:GLU:OE2	2.01	0.61
2:B:152:TRP:O	2:B:192:THR:O	2.17	0.61
2:D:65:PHE:CD1	2:D:78:LEU:HD11	2.35	0.60
1:A:105:ILE:N	1:A:105:ILE:HD12	2.16	0.60
2:D:78:LEU:CD1	2:D:80:MET:HG2	2.32	0.60
2:B:69:ARG:NE	2:B:71:ASN:HD21	2.00	0.60
1:A:114:VAL:HG22	1:A:135:LEU:HD22	1.83	0.60
2:D:9:GLY:HA3	2:D:105:THR:HG22	1.84	0.60
2:D:57:TYR:HB2	2:D:62:LYS:HD2	1.85	0.59
1:C:11:LEU:CD1	1:C:103:LEU:HD13	2.33	0.59
2:D:64:ARG:HG2	2:D:64:ARG:HH11	1.68	0.59
2:B:145:PRO:O	2:B:197:HIS:HE1	1.86	0.58
1:A:118:PRO:HG3	2:B:213:CYS:CB	2.32	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:114:VAL:HG22	1:C:135:LEU:CD1	2.26	0.58
2:D:145:PRO:O	2:D:197:HIS:HE1	1.85	0.58
1:C:189:ASN:ND2	1:C:209:ASN:O	2.37	0.58
1:A:30(B):SER:HA	1:A:30(E):ARG:HH21	1.69	0.58
1:A:37:GLN:CB	1:A:47:LEU:HD11	2.33	0.58
1:C:178:LEU:C	1:C:178:LEU:HD13	2.29	0.58
1:C:30(E):ARG:HD2	1:C:30(E):ARG:N	2.16	0.57
2:B:93:VAL:CG2	2:B:100(C):MET:HB3	2.33	0.57
1:C:3:VAL:HB	1:C:26:SER:OG	2.04	0.57
2:D:11:LEU:HD12	2:D:108:THR:O	2.05	0.56
1:A:158:VAL:C	1:A:159:LEU:HD12	2.30	0.56
2:B:123:ALA:HB3	2:B:212:ASP:HB2	1.86	0.56
1:C:20:THR:HG23	1:C:72:THR:CG2	2.36	0.56
1:C:213:CYS:SG	2:D:213:CYS:O	2.64	0.56
4:F:2:KDO:H5	4:F:3:KDO:O1A	2.05	0.56
1:C:74:THR:C	1:C:75:ILE:HD13	2.31	0.56
1:C:116:ILE:HD13	1:C:117:PHE:H	1.71	0.55
2:B:128:ALA:O	2:B:129:GLN:HB2	2.06	0.55
1:A:213:CYS:HB3	2:B:213:CYS:O	2.06	0.55
1:C:83:LEU:HD22	6:C:453:HOH:O	2.07	0.55
1:A:45:LYS:O	1:A:47:LEU:HD12	2.06	0.55
2:D:78:LEU:HD13	2:D:78:LEU:C	2.31	0.55
1:C:168:LYS:NZ	1:C:168:LYS:H	2.05	0.55
1:C:118:PRO:CG	2:D:213:CYS:HB3	2.37	0.54
1:C:24:LYS:HE2	1:C:70:ASP:OD1	2.06	0.54
1:A:33:LEU:HG	1:A:34:ALA:N	2.21	0.54
2:D:119:VAL:CG2	2:D:204:VAL:HG11	2.34	0.54
2:D:52(A):ASN:CG	2:D:52(C):PRO:HD2	2.33	0.54
2:B:212:ASP:O	2:B:213:CYS:HB3	2.06	0.54
1:A:158:VAL:O	1:A:159:LEU:HD12	2.08	0.53
2:B:52(B):LYS:HB3	2:B:52(C):PRO:HD3	1.89	0.53
2:D:69:ARG:NE	2:D:71:ASN:HD21	2.06	0.53
1:A:189:ASN:O	1:A:209:ASN:O	2.25	0.53
1:A:189:ASN:HD21	1:A:211:ASN:HB3	1.72	0.53
1:C:2:ILE:HG22	1:C:4:MET:HE1	1.89	0.53
2:D:35:SER:O	2:D:93:VAL:HG12	2.08	0.53
1:A:104:GLU:C	1:A:105:ILE:HD12	2.34	0.53
1:C:4:MET:HE1	1:C:25:SER:HB2	1.90	0.53
2:D:58:SER:O	2:D:62:LYS:HG2	2.09	0.53
2:D:195:VAL:HB	2:D:204:VAL:HG13	1.91	0.53
2:D:48:LEU:CD2	2:D:61:VAL:HG11	2.38	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:37:GLN:N	1:A:47:LEU:HD11	2.24	0.52
1:C:14:SER:N	1:C:106:LYS:HE2	2.23	0.52
2:D:132:SER:O	2:D:133:MET:HB2	2.08	0.52
2:D:175:LEU:HD12	2:D:175:LEU:C	2.34	0.52
1:C:13:VAL:C	1:C:106:LYS:HE2	2.34	0.52
1:C:194:GLU:HG2	1:C:205:VAL:HG12	1.90	0.52
1:A:135:LEU:HD13	1:A:143:ILE:HD13	1.91	0.52
1:C:143:ILE:HG23	1:C:174:MET:HE3	1.92	0.52
2:D:182:PRO:HG2	2:D:185:THR:OG1	2.10	0.52
1:C:189:ASN:HD21	1:C:211:ASN:HB3	1.72	0.52
2:D:103:GLN:CD	6:D:236:HOH:O	2.53	0.52
1:C:4:MET:HE1	1:C:25:SER:CB	2.40	0.52
1:C:178:LEU:CD1	1:C:180:LEU:HG	2.40	0.52
1:A:189:ASN:ND2	1:A:209:ASN:O	2.43	0.51
2:B:35:SER:HB2	2:B:93:VAL:CG1	2.40	0.51
2:B:135:THR:C	2:B:136:LEU:HD22	2.34	0.51
1:C:13:VAL:HA	1:C:106:LYS:HE2	1.91	0.51
1:A:33:LEU:HD22	1:A:71:PHE:CG	2.46	0.51
2:B:69:ARG:HE	2:B:71:ASN:ND2	2.05	0.51
1:C:143:ILE:CG2	1:C:174:MET:HE3	2.41	0.51
2:D:195:VAL:HB	2:D:204:VAL:CG1	2.39	0.51
1:C:118:PRO:HB3	1:C:208:PHE:CZ	2.45	0.51
1:A:148:LYS:HA	1:A:152:SER:O	2.10	0.51
2:D:197:HIS:HD2	2:D:200:SER:OG	1.93	0.50
2:D:210:PRO:O	2:D:211:ARG:HB3	2.12	0.50
1:C:159:LEU:HG	1:C:159:LEU:O	2.10	0.50
2:D:136:LEU:N	2:D:136:LEU:HD22	2.27	0.50
1:A:109:ASP:OD2	1:A:198:LYS:HE3	2.12	0.49
2:B:131:ASN:O	2:B:132:SER:HB2	2.11	0.49
3:E:4:KDO:H5	3:E:5:KDO:O1A	2.12	0.49
2:B:29:PHE:CD2	2:B:74:SER:HA	2.47	0.49
2:B:169:GLN:O	2:B:170:SER:HB2	2.11	0.49
1:C:12:ALA:CB	1:C:104:GLU:HG3	2.41	0.49
1:A:3:VAL:H	1:A:26:SER:HB2	1.76	0.49
1:A:19:VAL:CG2	1:A:78:VAL:HG21	2.43	0.49
1:A:83:LEU:HD13	1:A:83:LEU:C	2.38	0.49
1:A:163:THR:CG2	1:A:164:ASP:N	2.75	0.49
2:B:97:PHE:HB2	3:E:4:KDO:H32	1.95	0.48
1:C:189:ASN:HD21	1:C:211:ASN:CB	2.26	0.48
2:D:64:ARG:HB2	2:D:64:ARG:CZ	2.41	0.48
1:C:145:VAL:HG23	1:C:174:MET:HE1	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:80:MET:HB3	2:B:82(A):LEU:HD21	1.96	0.48
2:B:149:THR:HG22	6:B:436:HOH:O	2.13	0.48
2:B:117:PRO:HB3	2:B:143:TYR:HB3	1.94	0.48
1:C:93:ASN:HD22	1:C:93:ASN:C	2.20	0.48
2:D:59:ALA:HA	2:D:62:LYS:CG	2.43	0.47
2:B:154:SER:H	2:B:194:ASN:ND2	2.06	0.47
1:C:33:LEU:HD22	1:C:71:PHE:CG	2.49	0.47
1:C:83:LEU:HD21	6:C:390:HOH:O	2.14	0.47
1:C:13:VAL:CA	1:C:106:LYS:HE2	2.45	0.47
1:C:19:VAL:HG12	1:C:20:THR:N	2.29	0.47
1:C:182:LYS:O	1:C:186:GLU:HG3	2.14	0.47
2:D:61:VAL:HB	2:D:65:PHE:CD2	2.49	0.47
2:D:130:THR:HG22	2:D:131:ASN:N	2.30	0.47
1:A:30(D):THR:O	1:A:30(E):ARG:HB2	2.14	0.47
1:C:83:LEU:HD13	1:C:83:LEU:C	2.40	0.47
1:C:168:LYS:NZ	1:C:168:LYS:N	2.63	0.46
2:D:36:TRP:CE2	2:D:78:LEU:HB2	2.50	0.46
2:B:52(A):ASN:CG	2:B:52(C):PRO:HD2	2.40	0.46
2:B:185:THR:O	2:B:189:GLU:HB2	2.15	0.46
1:C:20:THR:HG23	1:C:72:THR:HG23	1.97	0.46
2:D:69:ARG:HE	2:D:71:ASN:ND2	2.11	0.46
1:A:47:LEU:HD12	1:A:47:LEU:N	2.31	0.46
2:B:175:LEU:C	2:B:175:LEU:HD12	2.40	0.46
2:D:6:GLU:HB2	2:D:105:THR:OG1	2.16	0.46
2:B:197:HIS:HD2	2:B:200:SER:OG	1.99	0.46
2:D:121:PRO:HD3	2:D:206:LYS:HG2	1.98	0.46
1:A:81:GLU:CD	1:A:81:GLU:H	2.24	0.45
2:D:185:THR:O	2:D:189:GLU:N	2.36	0.45
1:C:135:LEU:HD11	1:C:195:ALA:CB	2.47	0.45
2:B:136:LEU:N	2:B:136:LEU:CD2	2.78	0.45
1:C:2:ILE:CG2	1:C:4:MET:CE	2.89	0.45
1:A:118:PRO:CG	2:B:213:CYS:HB2	2.44	0.45
1:A:209:ASN:O	1:A:210:ARG:CB	2.62	0.45
1:A:116:ILE:HG13	1:A:133:CYS:SG	2.57	0.45
1:C:54:ARG:HG2	1:C:58:VAL:HB	1.98	0.45
1:A:19:VAL:HG23	1:A:78:VAL:HG21	1.99	0.45
2:D:64:ARG:NH2	2:D:83:ARG:HH21	2.15	0.44
1:A:3:VAL:HB	1:A:26:SER:OG	2.17	0.44
2:D:38:ARG:CD	2:D:48:LEU:HD21	2.48	0.44
1:C:189:ASN:O	1:C:209:ASN:O	2.36	0.44
2:D:35:SER:HB2	2:D:93:VAL:HG13	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:118:PRO:HB3	1:C:208:PHE:CE2	2.53	0.44
1:A:83:LEU:O	1:A:84:ALA:HB2	2.18	0.44
1:C:2:ILE:CG2	1:C:4:MET:HE3	2.43	0.44
2:D:47:TRP:CH2	2:D:49:GLY:HA2	2.53	0.44
2:B:133:MET:SD	2:B:182:PRO:HA	2.58	0.43
1:C:104:GLU:CD	6:C:455:HOH:O	2.62	0.43
2:D:64:ARG:HH11	2:D:64:ARG:CG	2.29	0.43
1:A:19:VAL:HG23	1:A:78:VAL:CG2	2.49	0.43
2:B:182:PRO:HG2	2:B:185:THR:OG1	2.18	0.43
2:D:59:ALA:HA	2:D:62:LYS:HG3	2.01	0.43
1:A:37:GLN:N	1:A:47:LEU:CD1	2.81	0.43
1:C:186:GLU:HA	1:C:210:ARG:NH2	2.34	0.43
1:A:105:ILE:N	1:A:105:ILE:CD1	2.82	0.42
2:D:127:ALA:O	2:D:128:ALA:C	2.62	0.42
1:A:42:GLN:HG3	6:A:228:HOH:O	2.19	0.42
2:B:131:ASN:O	2:B:132:SER:CB	2.67	0.42
2:D:126:SER:N	2:D:213:CYS:OXT	2.52	0.42
2:D:154:SER:N	2:D:194:ASN:ND2	2.57	0.42
1:C:11:LEU:CD1	1:C:103:LEU:CD1	2.97	0.42
1:C:116:ILE:HD13	1:C:116:ILE:C	2.43	0.42
2:D:50:PHE:CD1	2:D:50:PHE:C	2.97	0.42
2:D:78:LEU:HD12	2:D:80:MET:HG2	2.00	0.42
2:D:189:GLU:OE2	2:D:189:GLU:HA	2.20	0.42
1:C:168:LYS:H	1:C:168:LYS:HZ2	1.66	0.42
2:D:47:TRP:CZ2	2:D:49:GLY:HA2	2.54	0.42
1:A:132:VAL:HG22	1:A:177:THR:HG23	2.00	0.41
1:C:33:LEU:HG	1:C:34:ALA:N	2.34	0.41
2:D:210:PRO:O	2:D:211:ARG:CB	2.68	0.41
1:A:83:LEU:HB2	1:A:105:ILE:HD13	2.01	0.41
2:B:157:LEU:O	2:B:158:SER:C	2.62	0.41
1:A:61:ARG:HH21	1:A:82:ASP:CG	2.28	0.41
1:A:168:LYS:HB3	1:A:168:LYS:NZ	2.35	0.41
1:A:191:TYR:O	1:A:207:SER:HB2	2.19	0.41
1:C:184:GLU:OE2	1:C:187:ARG:CZ	2.69	0.41
3:E:3:KDO:H5	3:E:4:KDO:C1	2.50	0.41
1:C:178:LEU:HD11	1:C:180:LEU:HG	2.02	0.41
2:D:30:THR:HG1	2:D:52(B):LYS:HB2	1.86	0.41
2:D:52(B):LYS:HB3	2:D:52(C):PRO:HD3	2.02	0.41
1:A:46:LEU:HD13	1:A:55:GLU:CG	2.46	0.40
1:A:146:LYS:NZ	1:A:153:GLU:HB2	2.35	0.40
1:A:189:ASN:HD21	1:A:211:ASN:CB	2.34	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:37:GLN:HB2	1:A:47:LEU:HD11	2.02	0.40
1:A:186:GLU:O	1:A:210:ARG:NH2	2.54	0.40
2:D:181:VAL:HB	2:D:182:PRO:HD2	2.03	0.40
1:A:135:LEU:HD13	1:A:143:ILE:CD1	2.51	0.40
2:B:121:PRO:HB3	2:B:206:LYS:HG2	2.03	0.40
2:D:64:ARG:NH2	2:D:83:ARG:NH2	2.70	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	217/219 (99%)	209 (96%)	8 (4%)	0	100	100
1	C	217/219 (99%)	205 (94%)	12 (6%)	0	100	100
2	B	224/226 (99%)	208 (93%)	12 (5%)	4 (2%)	6	1
2	D	224/226 (99%)	214 (96%)	6 (3%)	4 (2%)	6	1
All	All	882/890 (99%)	836 (95%)	38 (4%)	8 (1%)	14	4

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	132	SER
2	B	211	ARG
2	D	41	PRO
2	D	128	ALA
2	D	211	ARG
2	B	130	THR
2	B	159	SER
2	D	133	MET

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	194/194 (100%)	187 (96%)	7 (4%)	31	11
1	C	194/194 (100%)	183 (94%)	11 (6%)	18	4
2	B	196/196 (100%)	186 (95%)	10 (5%)	21	5
2	D	196/196 (100%)	187 (95%)	9 (5%)	24	6
All	All	780/780 (100%)	743 (95%)	37 (5%)	23	6

All (37) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	19	VAL
1	A	46	LEU
1	A	55	GLU
1	A	83	LEU
1	A	107	ARG
1	A	124	LEU
1	A	174	MET
2	B	20	LEU
2	B	50	PHE
2	B	131	ASN
2	B	136	LEU
2	B	147	PRO
2	B	148	VAL
2	B	149	THR
2	B	157	LEU
2	B	206	LYS
2	B	207	LYS
1	C	46	LEU
1	C	47	LEU
1	C	83	LEU
1	C	93	ASN
1	C	94	LEU
1	C	107	ARG
1	C	116	ILE

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Mol	Chain	Res	Type
1	C	124	LEU
1	C	159	LEU
1	C	168	LYS
1	C	178	LEU
2	D	18	LEU
2	D	20	LEU
2	D	41	PRO
2	D	50	PHE
2	D	129	GLN
2	D	136	LEU
2	D	147	PRO
2	D	148	VAL
2	D	175	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	137	ASN
1	A	155	GLN
1	A	189	ASN
1	A	209	ASN
2	B	71	ASN
2	B	79	GLN
2	B	81	ASN
2	B	162	HIS
2	B	169	GLN
2	B	194	ASN
2	B	197	HIS
1	C	27	GLN
1	C	38	GLN
1	C	42	GLN
1	C	79	GLN
1	C	93	ASN
1	C	189	ASN
1	C	209	ASN
2	D	13	GLN
2	D	39	GLN
2	D	71	ASN
2	D	194	ASN
2	D	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 ⓘ

8 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	GP1	E	1	3	15,16,16	1.70	2 (13%)	24,24,24	1.02	2 (8%)
3	GP4	E	2	3	15,15,16	1.51	1 (6%)	20,22,24	1.20	1 (5%)
3	KDO	E	3	3	15,15,16	2.02	4 (26%)	17,21,24	1.54	4 (23%)
3	KDO	E	4	3	15,15,16	1.86	4 (26%)	17,21,24	1.59	4 (23%)
3	KDO	E	5	3	15,15,16	1.64	3 (20%)	17,21,24	1.35	3 (17%)
4	KDO	F	1	4	16,16,16	1.90	4 (25%)	17,24,24	1.48	3 (17%)
4	KDO	F	2	4	15,15,16	1.92	4 (26%)	17,21,24	1.73	5 (29%)
4	KDO	F	3	4	15,15,16	1.74	4 (26%)	17,21,24	1.39	3 (17%)

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	GP1	E	1	3	-	0/6/27/27	0/1/1/1
3	GP4	E	2	3	1/1/5/6	1/7/24/27	0/1/1/1
3	KDO	E	3	3	-	0/10/26/30	0/1/1/1
3	KDO	E	4	3	-	0/10/26/30	0/1/1/1
3	KDO	E	5	3	-	0/10/26/30	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	KDO	F	1	4	-	1/12/30/30	0/1/1/1
4	KDO	F	2	4	-	2/10/26/30	0/1/1/1
4	KDO	F	3	4	-	0/10/26/30	0/1/1/1

All (26) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	F	2	KDO	C4-C5	5.11	1.60	1.52
3	E	3	KDO	C4-C5	4.88	1.60	1.52
3	E	4	KDO	C4-C5	4.81	1.59	1.52
4	F	3	KDO	C4-C5	4.32	1.59	1.52
3	E	1	GP1	P4B-O1	-4.25	1.52	1.59
3	E	5	KDO	C4-C5	4.04	1.58	1.52
4	F	1	KDO	O6-C2	3.91	1.47	1.43
3	E	3	KDO	O6-C6	3.84	1.49	1.44
3	E	2	GP4	P4A-O4	-3.81	1.52	1.59
3	E	3	KDO	O6-C2	3.38	1.50	1.43
4	F	1	KDO	O2-C2	3.34	1.44	1.39
3	E	1	GP1	C3-C2	3.19	1.57	1.53
4	F	1	KDO	C4-C5	3.08	1.57	1.52
3	E	4	KDO	O6-C2	2.89	1.49	1.43
4	F	3	KDO	O6-C2	2.83	1.49	1.43
4	F	1	KDO	O6-C6	2.73	1.48	1.44
3	E	5	KDO	O6-C2	2.71	1.48	1.43
4	F	2	KDO	O6-C2	2.70	1.48	1.43
4	F	2	KDO	O6-C6	2.63	1.48	1.44
4	F	3	KDO	O6-C6	2.62	1.48	1.44
3	E	4	KDO	O6-C6	2.19	1.47	1.44
3	E	5	KDO	O6-C6	2.13	1.47	1.44
4	F	3	KDO	O1B-C1	-2.11	1.23	1.30
4	F	2	KDO	O5-C5	2.10	1.48	1.43
3	E	3	KDO	O1B-C1	-2.03	1.24	1.30
3	E	4	KDO	O5-C5	2.00	1.47	1.43

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	F	1	KDO	C3-C4-C5	-3.60	107.34	110.84
4	F	2	KDO	O6-C2-C3	3.52	115.29	110.56
3	E	2	GP4	C1-O5-C5	3.49	116.86	112.19
3	E	1	GP1	P4B-O1-C1	-3.36	114.52	123.54
3	E	4	KDO	O6-C2-C3	3.23	114.90	110.56

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	E	5	KDO	O1B-C1-O1A	-3.00	117.28	124.08
4	F	2	KDO	O6-C2-C1	-3.00	102.06	107.72
4	F	1	KDO	O1A-C1-C2	-2.96	118.91	123.85
4	F	3	KDO	O1B-C1-O1A	-2.62	118.14	124.08
4	F	2	KDO	C3-C4-C5	-2.54	107.00	110.67
3	E	3	KDO	O1B-C1-O1A	-2.53	118.35	124.08
4	F	1	KDO	C3-C2-C1	-2.52	108.17	112.84
3	E	4	KDO	O6-C2-C1	-2.40	103.19	107.72
3	E	3	KDO	O6-C2-C3	2.40	113.78	110.56
3	E	4	KDO	O1B-C1-O1A	-2.38	118.67	124.08
4	F	3	KDO	O6-C2-C1	-2.36	103.27	107.72
3	E	3	KDO	O1B-C1-C2	2.31	118.72	112.71
3	E	5	KDO	O6-C2-C1	-2.27	103.44	107.72
4	F	3	KDO	O1B-C1-C2	2.25	118.56	112.71
3	E	1	GP1	C4-C3-C2	-2.24	107.28	110.99
3	E	3	KDO	O6-C2-C1	-2.22	103.52	107.72
3	E	4	KDO	C3-C4-C5	-2.22	107.47	110.67
4	F	2	KDO	O1B-C1-C2	2.19	118.41	112.71
3	E	5	KDO	C3-C4-C5	-2.15	107.56	110.67
4	F	2	KDO	O1B-C1-O1A	-2.05	119.42	124.08

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	E	2	GP4	C1

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	F	2	KDO	C6-C7-C8-O8
4	F	2	KDO	O7-C7-C8-O8
3	E	2	GP4	C4-O4-P4A-O7A
4	F	1	KDO	O1B-C1-C2-O2

There are no ring outliers.

5 monomers are involved in 4 short contacts:

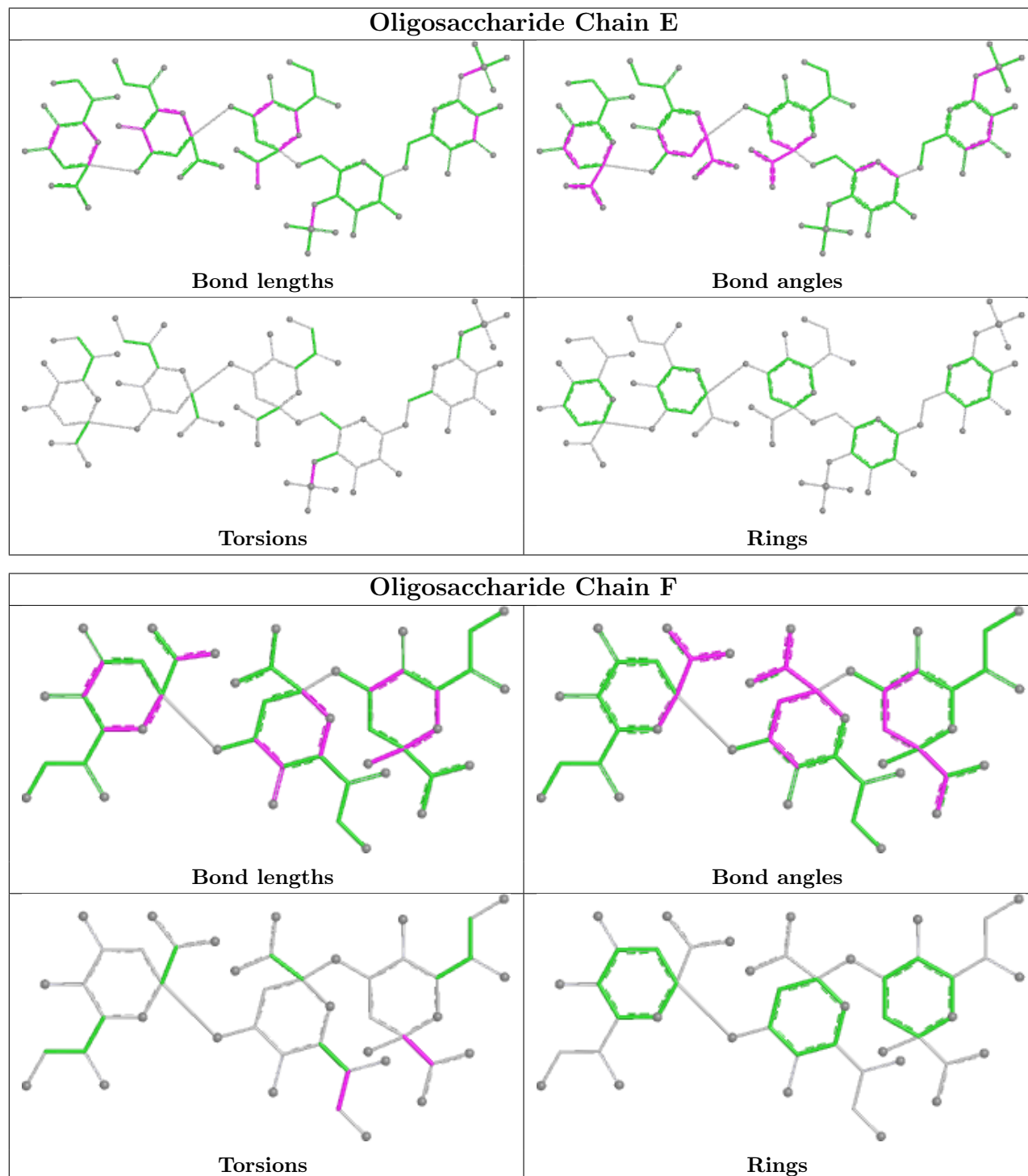
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	E	3	KDO	1	0
3	E	4	KDO	3	0
4	F	2	KDO	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	F	3	KDO	1	0
3	E	5	KDO	1	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

Of 5 ligands modelled in this entry, 5 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.

No monomer is involved in short contacts.

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	219/219 (100%)	0.50	18 (8%)	17 21	16, 29, 48, 70	0
1	C	219/219 (100%)	0.52	15 (6%)	23 27	18, 29, 42, 55	0
2	B	226/226 (100%)	0.55	30 (13%)	7 8	17, 27, 55, 77	0
2	D	226/226 (100%)	0.52	23 (10%)	12 14	17, 25, 50, 70	0
All	All	890/890 (100%)	0.52	86 (9%)	13 15	16, 28, 50, 77	0

All (86) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	213	CYS	8.8
2	B	213	CYS	5.9
1	C	159	LEU	5.1
2	D	138	CYS	4.9
1	A	213	CYS	4.4
2	B	128	ALA	4.2
2	B	212	ASP	4.2
2	D	211	ARG	4.1
2	D	131	ASN	4.0
2	B	129	GLN	4.0
2	D	130	THR	3.9
2	B	158	SER	3.9
2	D	132	SER	3.8
2	B	130	THR	3.7
2	D	128	ALA	3.7
2	D	212	ASP	3.5
2	B	131	ASN	3.4
2	B	157	LEU	3.4
1	A	211	ASN	3.4
2	B	170	SER	3.3
1	C	30(D)	THR	3.3

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Mol	Chain	Res	Type	RSRZ
2	B	105	THR	3.2
1	C	213	CYS	3.1
1	A	26	SER	3.0
1	A	60	ASP	3.0
2	B	138	CYS	3.0
1	A	209	ASN	2.9
2	B	211	ARG	2.9
1	C	30(B)	SER	2.8
2	D	189	GLU	2.8
1	C	156	ASN	2.7
2	B	152	TRP	2.7
2	B	168	LEU	2.7
2	B	156	SER	2.7
2	B	169	GLN	2.7
1	A	11	LEU	2.7
1	A	201	THR	2.7
2	B	127	ALA	2.7
1	A	190	SER	2.6
2	D	30	THR	2.6
2	D	41	PRO	2.6
2	B	126	SER	2.6
1	A	155	GLN	2.5
2	D	156	SER	2.5
2	D	42	GLY	2.5
1	C	60	ASP	2.5
2	D	185	THR	2.5
2	B	159	SER	2.5
2	D	154	SER	2.4
1	C	1	ASP	2.4
1	C	201	THR	2.4
2	B	134	VAL	2.4
1	A	156	ASN	2.3
1	A	144	ASN	2.3
1	A	152	SER	2.3
2	D	210	PRO	2.3
2	B	185	THR	2.3
2	D	127	ALA	2.3
2	B	136	LEU	2.3
2	B	186	TRP	2.3
2	D	60	SER	2.3
2	D	186	TRP	2.3
2	B	2	VAL	2.3

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Mol	Chain	Res	Type	RSRZ
2	B	182	PRO	2.3
2	D	59	ALA	2.3
2	B	198	PRO	2.2
1	C	75	ILE	2.2
1	C	211	ASN	2.2
2	B	148	VAL	2.2
1	C	43	PHE	2.2
1	C	106	LYS	2.2
2	D	93	VAL	2.1
2	D	78	LEU	2.1
1	A	202	SER	2.1
1	A	61	ARG	2.1
1	A	210	ARG	2.1
2	B	133	MET	2.1
1	A	203	PRO	2.1
2	D	188	SER	2.1
1	C	210	ARG	2.1
2	B	153	ASN	2.0
2	B	210	PRO	2.0
1	C	30(E)	ARG	2.0
1	A	1	ASP	2.0
1	A	133	CYS	2.0
1	C	168	LYS	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.

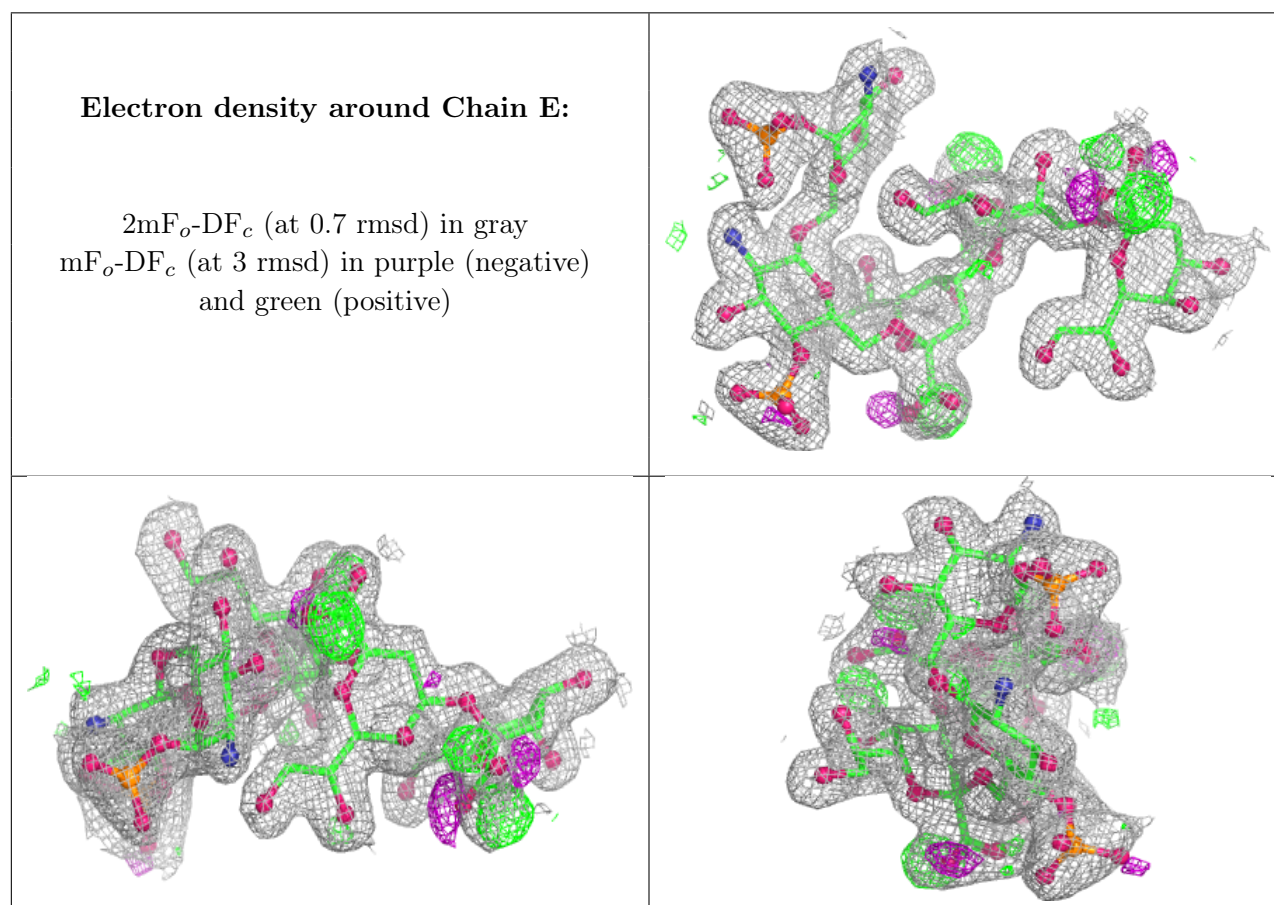
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	KDO	F	1	16/16	0.77	0.17	40,49,52,53	0
4	KDO	F	2	15/16	0.82	0.14	28,35,42,45	0
4	KDO	F	3	15/16	0.84	0.12	25,27,35,39	0
3	KDO	E	5	15/16	0.89	0.10	17,20,29,37	0
3	GP4	E	2	15/16	0.89	0.10	27,32,47,47	0

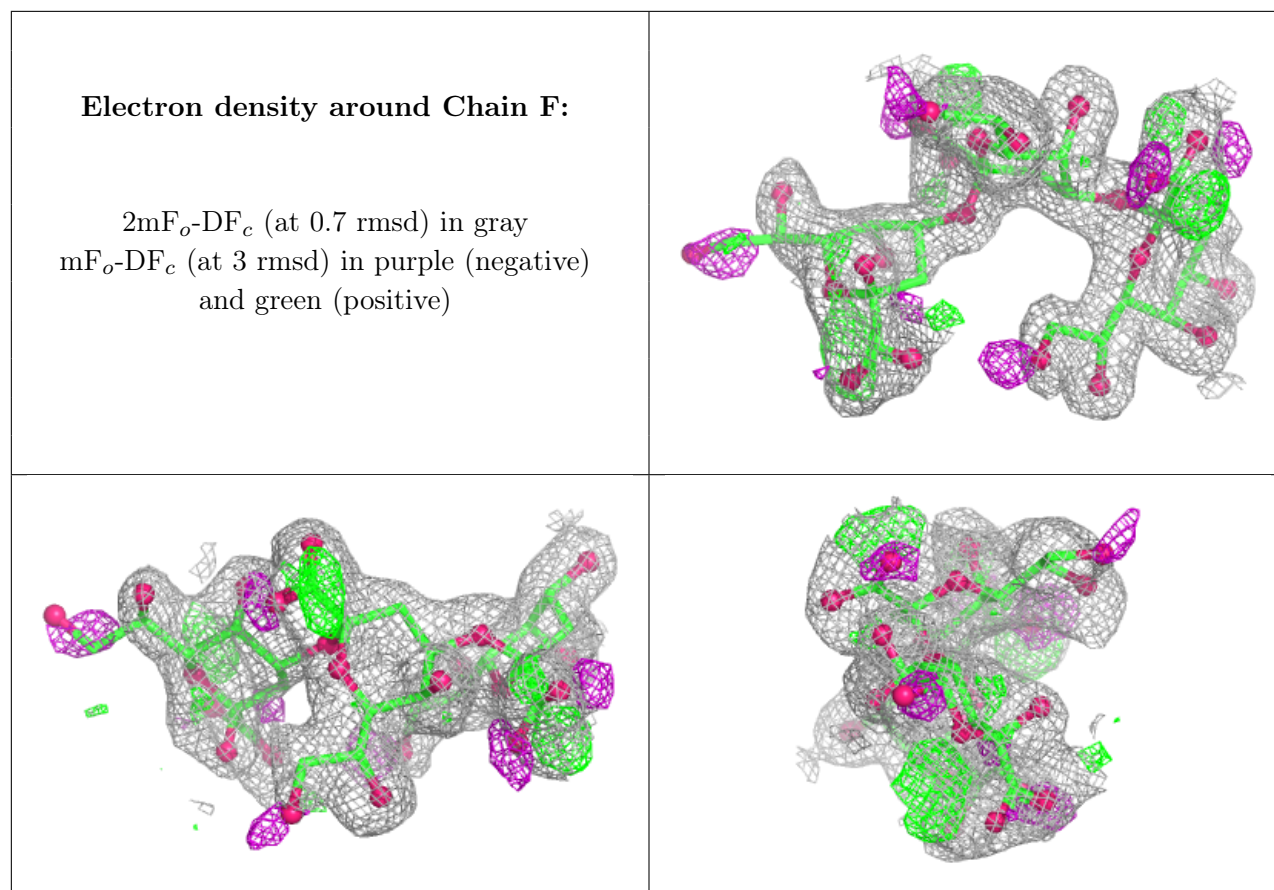
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	KDO	E	3	15/16	0.89	0.10	22,25,32,35	0
3	KDO	E	4	15/16	0.89	0.10	20,24,29,32	0
3	GP1	E	1	16/16	0.93	0.09	27,29,36,38	0

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
5	MG	A	215	1/1	0.97	0.06	26,26,26,26	0
5	MG	B	306	1/1	0.97	0.06	23,23,23,23	0
5	MG	C	314	1/1	0.99	0.03	24,24,24,24	0
5	MG	C	315	1/1	0.99	0.09	32,32,32,32	0
5	MG	A	214	1/1	1.00	0.02	16,16,16,16	0

6.5 Other polymers [i](#)

There are no such residues in this entry.