



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2026 – 05:28 PM UTC

PDB ID : 4HIJ / pdb_00004hij
Title : Anti-Streptococcus pneumoniae 23F Fab 023.102 with bound L-rhamnose-(1-2)-alpha-D-galactose-(3-O)-phosphate-2-glycerol
Authors : Bryson, S.; Risnes, L.; Damgupta, S.; Thomson, C.A.; Schrader, J.W.; Pai, E.F.
Deposited on : 2012-10-11
Resolution : 2.10 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

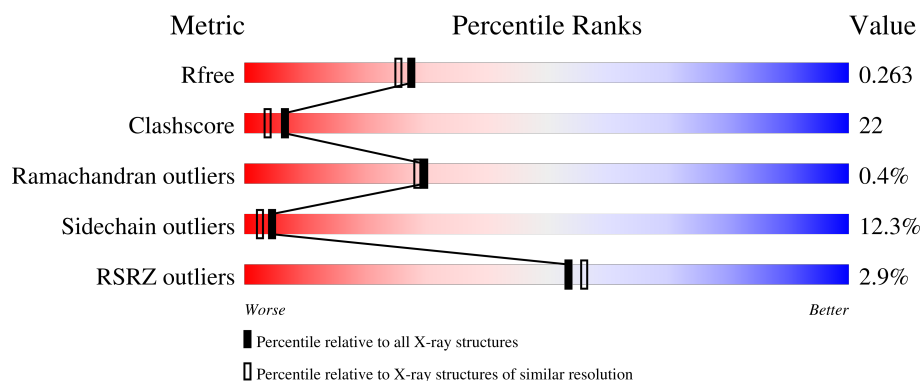
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

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	
1	C	219	
2	B	231	
2	D	231	
3	E	2	

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
3	F	2	50% 50%

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 6694 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.

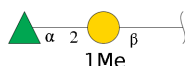
- Molecule 1 is a protein called Fab 023.102 light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	209	Total	C	N	O	S	0	0	0
			1614	1008	279	323	4			
1	C	212	Total	C	N	O	S	0	0	0
			1635	1021	283	327	4			

- Molecule 2 is a protein called Fab 023.102 heavy chain.

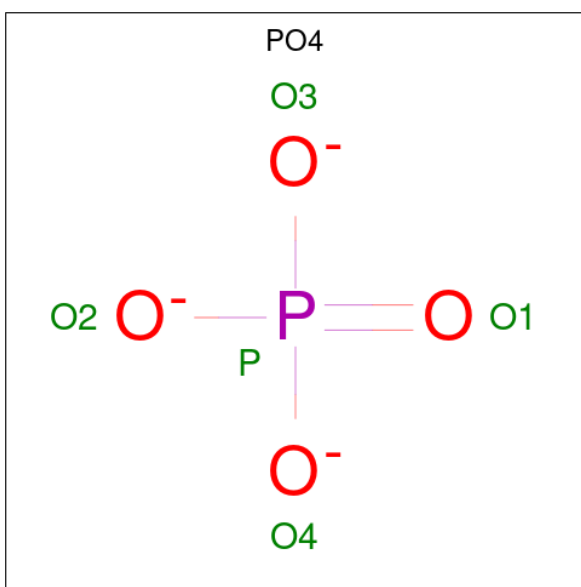
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	203	Total	C	N	O	S	0	0	0
			1534	969	261	298	6			
2	D	202	Total	C	N	O	S	0	0	0
			1528	966	260	296	6			

- Molecule 3 is an oligosaccharide called alpha-L-rhamnopyranose-(1-2)-methyl beta-D-galactopyranoside.



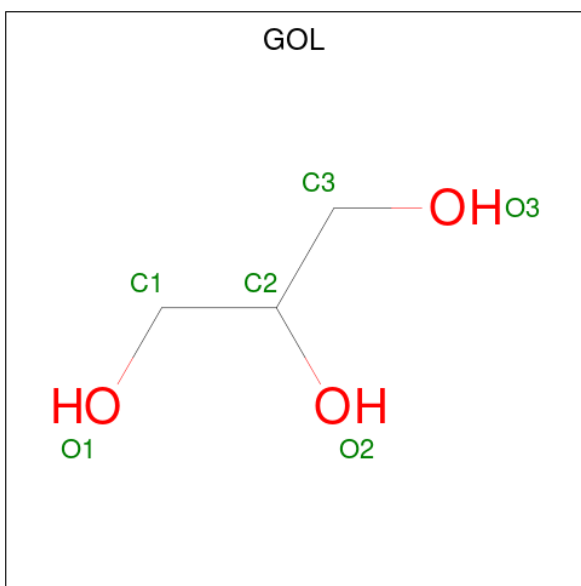
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
3	E	2	Total	C	O	0	0	0
			23	13	10			
3	F	2	Total	C	O	0	0	0
			23	13	10			

- Molecule 4 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	O	P	0	0
			4	3	1		
4	D	1	Total	O	P	0	0
			4	3	1		

- Molecule 5 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total	C	O	0	0
			5	3	2		
5	D	1	Total	C	O	0	0
			5	3	2		

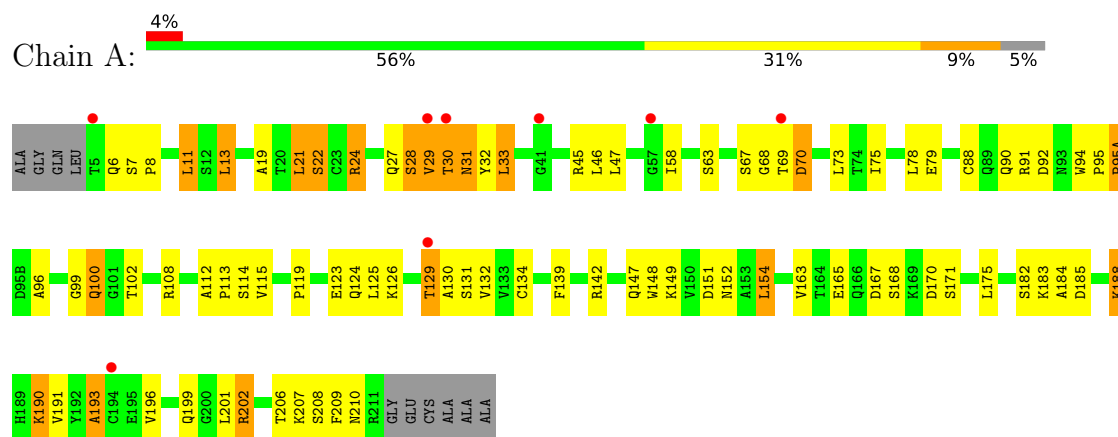
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	82	Total 82	O 82	0	0
6	B	64	Total 64	O 64	0	0
6	C	91	Total 91	O 91	0	0
6	D	82	Total 82	O 82	0	0

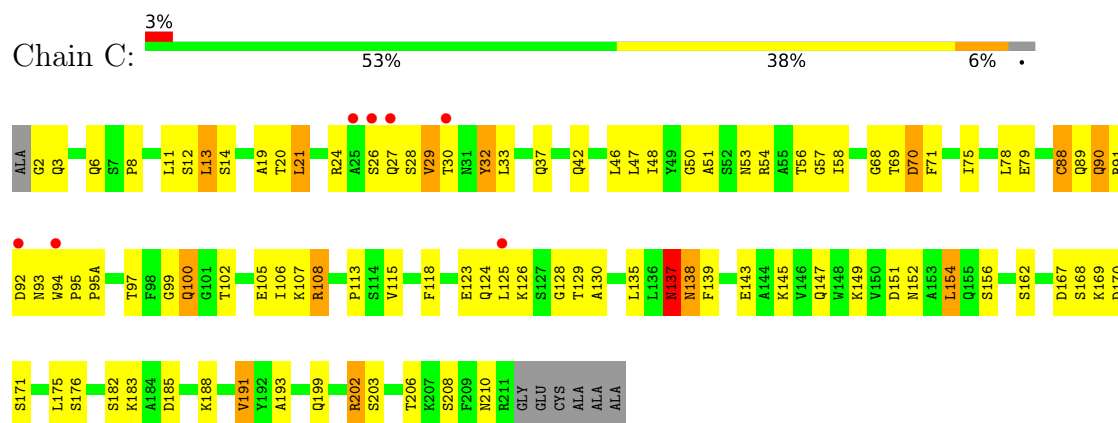
3 Residue-property plots [i](#)

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.

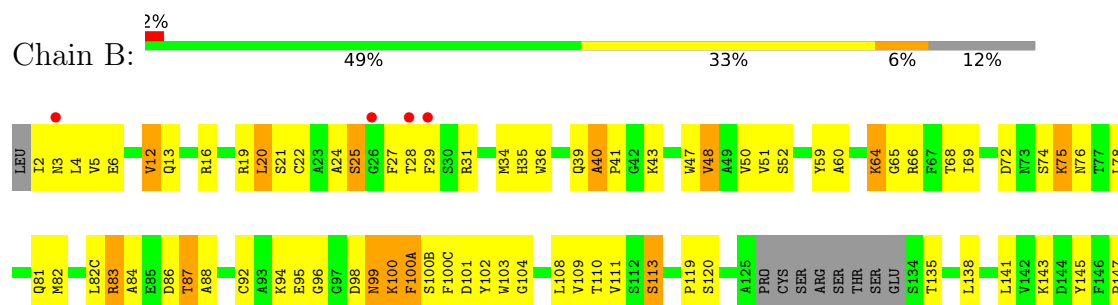
- Molecule 1: Fab 023.102 light chain

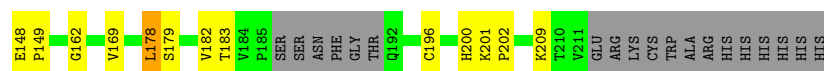


- Molecule 1: Fab 023.102 light chain

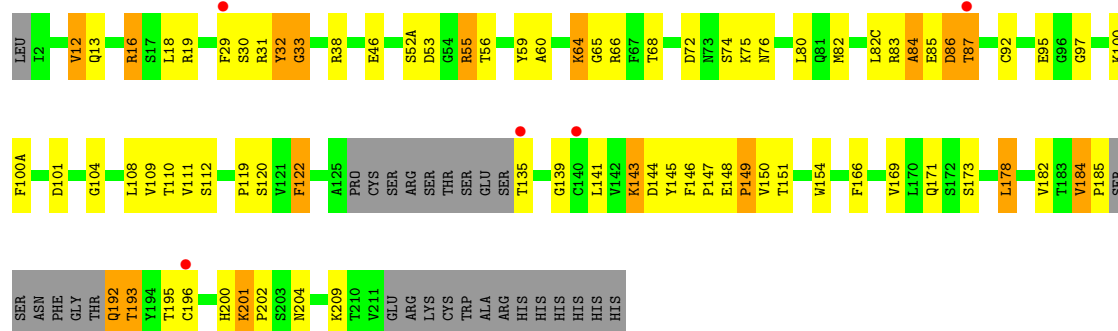


- Molecule 2: Fab 023.102 heavy chain





- Molecule 2: Fab 023.102 heavy chain



- Molecule 3: alpha-L-rhamnopyranose-(1-2)-methyl beta-D-galactopyranoside



- Molecule 3: alpha-L-rhamnopyranose-(1-2)-methyl beta-D-galactopyranoside



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	143.87Å 66.49Å 118.21Å 90.00° 111.58° 90.00°	Depositor
Resolution (Å)	17.00 – 2.10 17.00 – 2.10	Depositor EDS
% Data completeness (in resolution range)	(Not available) (17.00-2.10) 84.8 (17.00-2.10)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.45 (at 2.00Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.249 , 0.262 0.251 , 0.263	Depositor DCC
R_{free} test set	5391 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	33.4	Xtriage
Anisotropy	0.180	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 36.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	6694	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 41.05 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 2.5010e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: PO4, GOL, RAM, MBG

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	0.52	0/1651	1.18	19/2246 (0.8%)
1	C	0.50	0/1672	1.16	20/2274 (0.9%)
2	B	0.51	0/1569	1.17	16/2133 (0.8%)
2	D	0.51	0/1563	1.14	17/2125 (0.8%)
All	All	0.51	0/6455	1.16	72/8778 (0.8%)

There are no bond length outliers.

All (72) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	27	PHE	N-CA-C	-9.56	94.17	108.79
1	A	99	GLY	N-CA-C	-9.19	100.78	112.54
1	C	182	SER	N-CA-C	-8.12	99.59	110.55
1	A	95(A)	PRO	N-CA-C	7.92	124.14	113.84
1	C	58	ILE	CA-C-N	7.71	127.75	119.89
1	C	58	ILE	C-N-CA	7.71	127.75	119.89
2	B	143	LYS	N-CA-C	7.63	122.08	109.95
2	B	27	PHE	CA-C-N	-7.47	112.83	122.77
2	B	27	PHE	C-N-CA	-7.47	112.83	122.77
1	C	203	SER	N-CA-C	-7.45	101.51	108.22
1	A	31	ASN	N-CA-C	7.24	121.75	112.34
2	D	120	SER	N-CA-C	-7.16	98.07	109.59
2	D	32	TYR	N-CA-C	7.05	120.65	109.50
2	D	144	ASP	N-CA-C	7.04	119.82	111.02
1	A	79	GLU	CA-C-N	6.95	126.93	119.28
1	A	79	GLU	C-N-CA	6.95	126.93	119.28
2	B	99	ASN	N-CA-C	-6.94	96.02	110.80
1	C	79	GLU	N-CA-C	-6.86	101.41	110.40
2	B	100(A)	PHE	N-CA-C	-6.67	104.76	113.17
1	A	182	SER	N-CA-C	-6.66	100.63	110.48

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	115	VAL	N-CA-C	6.54	118.45	108.71
1	C	14	SER	CA-C-N	6.48	126.82	119.83
1	C	14	SER	C-N-CA	6.48	126.82	119.83
2	D	60	ALA	N-CA-C	-6.38	102.10	110.53
2	D	122	PHE	CA-C-N	6.30	126.80	119.99
2	D	122	PHE	C-N-CA	6.30	126.80	119.99
2	D	84	ALA	N-CA-C	-6.26	104.51	111.71
1	A	79	GLU	N-CA-C	-6.24	100.84	110.58
1	C	99	GLY	N-CA-C	-6.15	103.10	112.51
2	B	40	ALA	CA-C-N	6.13	127.50	119.84
2	B	40	ALA	C-N-CA	6.13	127.50	119.84
1	C	137	ASN	N-CA-C	6.12	119.21	109.24
1	A	28	SER	N-CA-C	-6.05	105.20	113.30
2	B	148	GLU	N-CA-C	-5.99	101.41	110.39
1	A	115	VAL	N-CA-C	5.97	117.61	108.71
1	C	14	SER	N-CA-C	-5.86	102.72	110.40
2	B	92	CYS	N-CA-C	-5.83	100.78	110.17
1	A	196	VAL	N-CA-C	5.70	116.68	108.48
1	C	50	GLY	N-CA-C	-5.69	106.14	114.90
1	A	114	SER	N-CA-C	-5.62	100.13	109.46
2	B	104	GLY	N-CA-C	-5.55	103.92	112.85
2	D	143	LYS	N-CA-C	5.54	119.08	110.17
1	A	193	ALA	CA-C-N	-5.53	115.31	123.05
1	A	193	ALA	C-N-CA	-5.53	115.31	123.05
2	D	92	CYS	N-CA-C	-5.51	101.31	110.17
2	D	33	GLY	N-CA-C	-5.49	105.06	112.14
2	B	48	VAL	CB-CA-C	-5.46	107.07	111.71
1	C	32	TYR	N-CA-C	5.44	119.76	113.18
2	B	60	ALA	N-CA-C	-5.43	102.45	110.48
1	A	193	ALA	N-CA-C	5.42	118.02	108.75
2	D	173	SER	N-CA-C	-5.39	106.66	113.18
1	C	191	VAL	N-CA-C	5.37	115.84	107.99
1	C	88	CYS	N-CA-C	-5.35	101.56	110.17
2	D	193	THR	N-CA-C	5.32	118.73	110.17
1	A	22	SER	N-CA-C	5.29	118.36	109.95
1	C	27	GLN	N-CA-C	-5.27	101.17	109.50
2	B	120	SER	N-CA-C	-5.27	100.65	109.24
1	A	165	GLU	N-CA-C	-5.21	103.65	110.53
1	C	118	PHE	CA-C-N	-5.19	115.03	120.38
1	C	118	PHE	C-N-CA	-5.19	115.03	120.38
2	B	113	SER	N-CA-C	-5.19	107.12	112.93
1	A	33	LEU	N-CA-C	5.17	117.91	109.59

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	59	TYR	N-CA-C	5.16	117.52	109.52
2	D	148	GLU	N-CA-C	-5.13	102.81	110.20
2	D	195	THR	N-CA-C	5.10	117.72	108.69
1	A	190	LYS	N-CA-C	5.06	119.67	112.93
2	D	104	GLY	N-CA-C	-5.04	104.73	112.85
2	D	166	PHE	N-CA-C	5.03	117.02	110.08
2	D	86	ASP	N-CA-C	5.02	118.67	112.54
1	C	69	THR	N-CA-C	-5.01	107.01	113.02
1	C	57	GLY	N-CA-C	-5.01	108.41	114.92
1	A	88	CYS	N-CA-C	-5.00	102.38	110.14

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	1614	0	1562	66	0
1	C	1635	0	1584	79	0
2	B	1534	0	1490	68	0
2	D	1528	0	1485	70	0
3	E	23	0	22	2	0
3	F	23	0	22	0	0
4	B	4	0	0	1	0
4	D	4	0	0	0	0
5	B	5	0	5	1	0
5	D	5	0	5	0	0
6	A	82	0	0	3	0
6	B	64	0	0	1	0
6	C	91	0	0	4	0
6	D	82	0	0	7	0
All	All	6694	0	6175	270	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 22.

All (270) 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:B:28:THR:HB	2:B:31:ARG:NE	1.64	1.12
1:A:24:ARG:HG3	1:A:24:ARG:HH11	0.96	1.08
2:B:28:THR:HB	2:B:31:ARG:HE	0.94	1.06
2:B:64:LYS:HD2	2:B:65:GLY:N	1.74	1.01
1:A:24:ARG:HG3	1:A:24:ARG:NH1	1.76	0.94
1:A:202:ARG:HG3	1:A:202:ARG:HH11	1.33	0.93
1:C:2:GLY:HA2	1:C:97:THR:HG21	1.51	0.93
1:C:100:GLN:H	1:C:100:GLN:NE2	1.67	0.92
2:D:64:LYS:HD3	2:D:65:GLY:N	1.85	0.91
1:C:100:GLN:H	1:C:100:GLN:HE21	1.16	0.90
2:B:64:LYS:HD2	2:B:65:GLY:H	1.37	0.87
1:A:24:ARG:HH11	1:A:24:ARG:CG	1.88	0.83
2:B:84:ALA:O	2:B:87:THR:HG23	1.78	0.82
2:B:196:CYS:SG	2:B:209:LYS:HB3	2.19	0.81
2:B:119:PRO:HB3	2:B:145:TYR:HB3	1.62	0.80
2:D:192:GLN:HG3	2:D:193:THR:N	1.94	0.80
1:C:93:ASN:HD21	2:D:100:LYS:NZ	1.81	0.78
2:B:81:GLN:HE22	2:D:16:ARG:HH21	1.30	0.78
1:C:28:SER:HB2	1:C:68:GLY:O	1.83	0.77
1:C:125:LEU:HB3	1:C:183:LYS:HD2	1.65	0.76
1:A:13:LEU:HD23	1:A:78:LEU:HD11	1.67	0.76
1:C:147:GLN:HE21	1:C:149:LYS:HD2	1.51	0.75
1:C:113:PRO:HB3	1:C:139:PHE:HB3	1.68	0.74
1:C:93:ASN:ND2	2:D:100:LYS:NZ	2.36	0.74
2:D:84:ALA:O	2:D:87:THR:HG23	1.87	0.74
2:B:201:LYS:HE2	6:B:442:HOH:O	1.88	0.74
1:A:94:TRP:CE3	3:E:1:MBG:H71	2.23	0.73
2:D:66:ARG:NH2	2:D:86:ASP:OD2	2.22	0.73
1:C:100:GLN:HE21	1:C:100:GLN:N	1.86	0.73
2:D:64:LYS:HD3	2:D:65:GLY:H	1.51	0.72
1:A:113:PRO:HB3	1:A:139:PHE:HB3	1.71	0.72
1:C:147:GLN:NE2	1:C:149:LYS:HD2	2.05	0.72
1:C:21:LEU:HD23	1:C:21:LEU:N	2.05	0.71
1:C:169:LYS:HD2	1:C:169:LYS:C	2.16	0.71
1:A:28:SER:HB3	1:A:68:GLY:O	1.90	0.70
1:A:193:ALA:HB2	1:A:208:SER:HB3	1.74	0.70
2:D:55:ARG:HG3	2:D:55:ARG:HH11	1.59	0.68
2:B:20:LEU:HD13	2:B:82:MET:CE	2.23	0.68
1:C:70:ASP:OD1	1:C:70:ASP:N	2.27	0.67
2:D:154:TRP:CH2	2:D:196:CYS:SG	2.87	0.67
1:A:8:PRO:CG	1:A:11:LEU:HD13	2.25	0.67

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:202:ARG:HG3	1:A:202:ARG:NH1	2.05	0.67
1:C:93:ASN:ND2	2:D:100:LYS:HZ1	1.93	0.66
2:B:19:ARG:HD3	2:D:16:ARG:HD2	1.78	0.66
2:D:55:ARG:HG3	2:D:55:ARG:NH1	2.11	0.66
1:C:94:TRP:HA	1:C:95:PRO:C	2.22	0.65
2:D:139:GLY:HA2	2:D:154:TRP:CH2	2.32	0.65
2:B:40:ALA:HB1	2:B:43:LYS:HE2	1.78	0.65
1:C:13:LEU:HD23	1:C:78:LEU:HD11	1.78	0.65
1:A:24:ARG:NH1	1:A:70:ASP:OD2	2.30	0.64
2:D:38:ARG:NH2	2:D:46:GLU:OE1	2.30	0.64
2:D:119:PRO:HB3	2:D:145:TYR:HB3	1.79	0.64
1:C:93:ASN:HD21	2:D:100:LYS:HZ1	1.44	0.64
1:A:6:GLN:H	1:A:100:GLN:HE22	1.44	0.63
1:C:138:ASN:H	1:C:138:ASN:HD22	1.46	0.63
2:B:29:PHE:CD2	2:B:76:ASN:HA	2.34	0.63
1:A:151:ASP:O	1:A:152:ASN:HB2	1.97	0.62
1:C:202:ARG:HG2	6:C:319:HOH:O	1.97	0.62
2:B:20:LEU:HD13	2:B:82:MET:HE2	1.81	0.62
2:B:20:LEU:HD22	2:B:82:MET:HE2	1.80	0.62
2:D:29:PHE:CB	2:D:76:ASN:HD22	2.13	0.62
1:A:28:SER:CB	1:A:68:GLY:O	2.48	0.62
1:A:126:LYS:HG2	6:A:380:HOH:O	2.00	0.61
1:C:138:ASN:HD22	1:C:138:ASN:N	1.98	0.61
2:B:28:THR:OG1	2:B:31:ARG:NH2	2.30	0.61
1:A:175:LEU:C	1:A:175:LEU:HD23	2.26	0.61
2:D:87:THR:HB	2:D:110:THR:HA	1.82	0.61
2:B:28:THR:HB	2:B:31:ARG:CZ	2.29	0.61
1:C:156:SER:HB2	6:C:369:HOH:O	2.00	0.61
2:D:72:ASP:CG	2:D:75:LYS:HD3	2.26	0.61
2:D:12:VAL:HG13	2:D:16:ARG:HB2	1.81	0.61
1:C:11:LEU:HG	1:C:13:LEU:HD22	1.83	0.60
2:D:29:PHE:CD2	2:D:76:ASN:HA	2.37	0.60
4:B:303:PO4:P	5:B:304:GOL:O1	2.59	0.60
1:A:147:GLN:CD	1:A:154:LEU:CD2	2.75	0.59
2:B:13:GLN:NE2	2:B:113:SER:HA	2.18	0.59
2:D:192:GLN:HG3	2:D:193:THR:H	1.65	0.59
2:D:85:GLU:HB2	6:D:404:HOH:O	2.02	0.59
1:A:21:LEU:HD23	1:A:21:LEU:N	2.18	0.58
1:C:91:ARG:HD2	1:C:95(A):PRO:O	2.04	0.58
2:D:143:LYS:NZ	2:D:171:GLN:OE1	2.36	0.58
2:B:178:LEU:C	2:B:178:LEU:HD12	2.29	0.57

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:129:THR:HG22	1:C:130:ALA:N	2.19	0.57
2:D:209:LYS:HE2	6:D:464:HOH:O	2.04	0.57
2:D:184:VAL:HG22	2:D:184:VAL:O	2.05	0.57
1:A:125:LEU:HB3	1:A:183:LYS:HE3	1.87	0.57
1:C:143:GLU:H	1:C:143:GLU:CD	2.12	0.57
2:D:64:LYS:HD3	2:D:64:LYS:C	2.26	0.57
1:A:8:PRO:HG2	1:A:11:LEU:HD13	1.86	0.56
2:D:55:ARG:HH11	2:D:55:ARG:CG	2.18	0.56
1:C:93:ASN:HD21	2:D:100:LYS:HZ2	1.53	0.56
1:C:108:ARG:HD2	1:C:170:ASP:O	2.05	0.56
2:B:135:THR:HG23	2:B:135:THR:O	2.05	0.56
1:C:147:GLN:HG3	1:C:154:LEU:HD21	1.87	0.56
1:A:29:VAL:HG11	1:A:32:TYR:CE2	2.41	0.56
1:A:94:TRP:CE3	1:A:95(A):PRO:HD3	2.40	0.56
1:C:94:TRP:CE3	1:C:95(A):PRO:HD3	2.41	0.56
2:B:28:THR:CB	2:B:31:ARG:HH21	2.19	0.55
1:C:24:ARG:NH1	1:C:70:ASP:OD2	2.39	0.55
2:D:84:ALA:HA	2:D:111:VAL:HB	1.88	0.55
2:B:94:LYS:NZ	2:B:95:GLU:O	2.36	0.55
1:A:124:GLN:NE2	1:A:131:SER:H	2.03	0.55
1:C:33:LEU:HG	1:C:71:PHE:CD2	2.41	0.55
2:B:72:ASP:CG	2:B:75:LYS:HG3	2.32	0.55
1:A:19:ALA:HB3	1:A:75:ILE:HB	1.89	0.54
1:A:30:THR:HG23	1:A:31:ASN:H	1.73	0.54
1:C:11:LEU:CD2	1:C:21:LEU:HD22	2.37	0.54
1:C:48:ILE:HD13	1:C:54:ARG:HA	1.89	0.54
1:C:32:TYR:CZ	1:C:92:ASP:OD2	2.61	0.53
2:D:29:PHE:HB3	2:D:76:ASN:HD22	1.73	0.53
1:C:20:THR:C	1:C:21:LEU:HD23	2.33	0.53
1:C:30:THR:HB	1:C:51:ALA:HB3	1.90	0.53
1:C:169:LYS:HD2	1:C:169:LYS:O	2.08	0.53
2:B:82:MET:HB3	2:B:82(C):LEU:HD21	1.91	0.53
1:C:11:LEU:HD22	1:C:21:LEU:HD22	1.91	0.53
2:B:50:VAL:HG22	2:B:51:VAL:N	2.24	0.53
1:A:154:LEU:HD13	1:A:154:LEU:C	2.34	0.52
1:C:89:GLN:HG2	1:C:90:GLN:N	2.23	0.52
2:B:40:ALA:CB	2:B:43:LYS:HE2	2.39	0.52
1:C:138:ASN:N	1:C:138:ASN:ND2	2.56	0.52
2:D:13:GLN:O	2:D:16:ARG:HB2	2.09	0.52
1:C:135:LEU:HD11	1:C:137:ASN:HD22	1.74	0.52
2:D:135:THR:HG23	2:D:135:THR:O	2.09	0.52

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:13:GLN:HA	2:D:112:SER:O	2.10	0.52
2:D:30:SER:O	2:D:52(A):SER:HB2	2.10	0.52
2:B:12:VAL:HG13	2:B:16:ARG:HB2	1.92	0.52
1:A:112:ALA:HB1	1:A:201:LEU:HD23	1.91	0.52
2:B:19:ARG:CD	2:D:16:ARG:HD2	2.40	0.52
1:A:94:TRP:HA	1:A:95:PRO:C	2.33	0.52
1:C:93:ASN:ND2	2:D:100:LYS:HZ2	2.07	0.51
1:A:47:LEU:HA	1:A:58:ILE:HG13	1.91	0.51
1:C:21:LEU:N	1:C:21:LEU:CD2	2.72	0.51
1:A:75:ILE:HG21	1:A:78:LEU:HD23	1.93	0.51
1:A:142:ARG:CZ	1:A:163:VAL:HG21	2.41	0.51
1:C:105:GLU:HG2	1:C:106:ILE:N	2.24	0.51
1:C:129:THR:HG22	1:C:130:ALA:H	1.76	0.51
2:B:64:LYS:HD2	2:B:64:LYS:C	2.21	0.51
1:A:8:PRO:O	1:A:102:THR:HG23	2.11	0.51
2:B:6:GLU:HA	2:B:21:SER:O	2.11	0.51
1:A:7:SER:HA	1:A:8:PRO:C	2.36	0.51
2:B:41:PRO:O	2:B:43:LYS:NZ	2.42	0.50
1:C:100:GLN:NE2	1:C:100:GLN:N	2.48	0.50
2:D:53:ASP:OD1	2:D:55:ARG:HG2	2.12	0.50
2:D:19:ARG:HD3	6:D:479:HOH:O	2.11	0.50
2:B:39:GLN:C	2:B:88:ALA:HB1	2.37	0.50
1:C:124:GLN:HA	2:D:122:PHE:CE1	2.47	0.50
2:D:139:GLY:HA2	2:D:154:TRP:HH2	1.74	0.50
2:B:34:MET:HB3	2:B:78:LEU:HD22	1.94	0.50
2:B:36:TRP:HD1	2:B:69:ILE:HD12	1.77	0.50
2:D:16:ARG:HD3	6:D:466:HOH:O	2.11	0.50
1:C:191:VAL:HG22	1:C:210:ASN:OD1	2.11	0.49
2:D:145:TYR:CE2	2:D:150:VAL:HG13	2.47	0.49
2:D:178:LEU:C	2:D:178:LEU:HD12	2.38	0.49
1:A:184:ALA:O	1:A:188:LYS:HG2	2.12	0.49
1:C:48:ILE:HA	1:C:53:ASN:O	2.12	0.49
1:A:29:VAL:HB	1:A:32:TYR:CD2	2.47	0.49
1:A:24:ARG:CG	1:A:69:THR:HG23	2.43	0.48
1:A:45:ARG:HD2	6:A:320:HOH:O	2.12	0.48
2:B:87:THR:HG22	2:B:111:VAL:H	1.79	0.48
1:C:29:VAL:HG11	1:C:32:TYR:CE2	2.48	0.48
2:D:154:TRP:CZ2	2:D:196:CYS:SG	3.06	0.48
2:B:87:THR:HA	2:B:109:VAL:O	2.14	0.48
1:C:167:ASP:O	1:C:171:SER:HA	2.14	0.48
2:D:82:MET:HE1	2:D:109:VAL:HG21	1.95	0.48

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:147:GLN:NE2	1:A:154:LEU:HD21	2.29	0.47
2:D:95:GLU:HG2	2:D:100:LYS:O	2.14	0.47
1:A:167:ASP:HB3	1:A:170:ASP:OD1	2.14	0.47
1:C:151:ASP:O	1:C:152:ASN:HB2	2.15	0.47
1:A:119:PRO:HG3	1:A:209:PHE:CD2	2.50	0.47
2:B:95:GLU:HG2	2:B:100:LYS:O	2.15	0.47
1:A:24:ARG:HG2	1:A:69:THR:HG23	1.97	0.46
2:B:3:ASN:N	2:B:25:SER:O	2.47	0.46
2:D:33:GLY:O	2:D:95:GLU:HB2	2.15	0.46
1:A:30:THR:HG23	1:A:31:ASN:N	2.30	0.46
2:B:96:GLY:HA3	2:B:99:ASN:O	2.16	0.46
2:D:59:TYR:CD1	2:D:64:LYS:HE2	2.50	0.46
2:B:66:ARG:NH2	2:B:86:ASP:OD2	2.48	0.46
2:D:151:THR:HG23	6:D:470:HOH:O	2.15	0.46
1:A:21:LEU:HG	1:A:73:LEU:HB3	1.97	0.46
2:B:20:LEU:CD1	2:B:82:MET:CE	2.94	0.46
1:C:123:GLU:O	1:C:126:LYS:HB2	2.16	0.46
1:C:3:GLN:HE21	1:C:26:SER:HB3	1.81	0.46
1:A:28:SER:O	1:A:30:THR:N	2.48	0.46
1:C:91:ARG:HG2	1:C:92:ASP:N	2.31	0.46
2:B:12:VAL:HG13	2:B:13:GLN:O	2.15	0.45
2:D:66:ARG:HH22	2:D:86:ASP:CG	2.23	0.45
1:C:91:ARG:HG2	1:C:92:ASP:H	1.81	0.45
2:B:35:HIS:ND1	2:B:50:VAL:HB	2.32	0.45
1:C:169:LYS:C	1:C:169:LYS:CD	2.83	0.45
2:D:12:VAL:HG13	2:D:16:ARG:CB	2.45	0.45
2:D:32:TYR:HE2	2:D:97:GLY:HA2	1.82	0.45
2:D:84:ALA:O	2:D:87:THR:CG2	2.62	0.45
1:A:147:GLN:CD	1:A:154:LEU:HD21	2.41	0.45
1:A:123:GLU:OE1	1:A:123:GLU:N	2.50	0.45
1:A:28:SER:O	1:A:29:VAL:C	2.60	0.44
2:D:100:LYS:HE3	2:D:100(A):PHE:CE1	2.52	0.44
2:D:83:ARG:O	2:D:111:VAL:HG21	2.18	0.44
1:A:193:ALA:CB	1:A:208:SER:HB3	2.45	0.44
2:B:162:GLY:O	2:B:182:VAL:HA	2.18	0.44
1:A:24:ARG:HA	1:A:69:THR:O	2.17	0.44
2:B:87:THR:HB	2:B:110:THR:HA	1.99	0.44
1:C:28:SER:HB2	1:C:68:GLY:C	2.41	0.44
2:D:29:PHE:HB2	2:D:76:ASN:HD22	1.81	0.44
2:B:102:TYR:C	2:B:103:TRP:CD1	2.96	0.44
1:A:167:ASP:O	1:A:171:SER:HA	2.18	0.44

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:37:GLN:HB2	1:C:47:LEU:HD11	2.00	0.43
1:C:46:LEU:HD22	2:D:101:ASP:HA	2.00	0.43
1:C:89:GLN:HG2	1:C:90:GLN:H	1.82	0.43
2:B:28:THR:CB	2:B:31:ARG:NH2	2.80	0.43
2:D:146:PHE:HA	2:D:147:PRO:HA	1.82	0.43
1:A:46:LEU:HD22	2:B:101:ASP:HA	2.00	0.43
1:A:95:PRO:HG2	6:A:377:HOH:O	2.18	0.43
2:B:147:PRO:HD2	2:B:202:PRO:CB	2.48	0.43
2:D:12:VAL:O	2:D:111:VAL:HA	2.19	0.43
2:B:4:LEU:CD2	2:B:24:ALA:HB2	2.49	0.43
1:A:113:PRO:CB	1:A:139:PHE:HB3	2.46	0.43
2:D:75:LYS:HE2	6:D:481:HOH:O	2.17	0.43
1:C:6:GLN:CG	1:C:88:CYS:SG	3.06	0.43
2:D:201:LYS:N	2:D:202:PRO:CD	2.82	0.43
1:A:193:ALA:HB2	1:A:208:SER:CB	2.46	0.42
1:A:67:SER:O	1:A:70:ASP:OD1	2.37	0.42
1:A:112:ALA:HB1	1:A:201:LEU:CD2	2.49	0.42
2:B:66:ARG:NH2	2:B:83:ARG:HG3	2.34	0.42
2:B:81:GLN:HE22	2:D:16:ARG:NH2	2.08	0.42
2:B:82:MET:HE1	2:B:109:VAL:HG21	2.00	0.42
1:C:175:LEU:C	1:C:175:LEU:HD23	2.45	0.42
1:A:96:ALA:HB3	2:B:47:TRP:CD1	2.55	0.42
1:A:129:THR:HG22	1:A:130:ALA:H	1.84	0.42
2:B:83:ARG:HG3	2:B:86:ASP:OD2	2.20	0.42
1:C:94:TRP:CA	1:C:95:PRO:C	2.92	0.42
2:B:36:TRP:O	2:B:48:VAL:HB	2.18	0.42
2:D:18:LEU:HD12	2:D:109:VAL:HG13	2.02	0.42
1:C:32:TYR:CE2	1:C:92:ASP:OD2	2.72	0.42
1:A:91:ARG:HD2	1:A:95(A):PRO:O	2.20	0.42
2:D:80:LEU:HD12	2:D:80:LEU:HA	1.88	0.42
2:B:200:HIS:CD2	2:B:202:PRO:HD2	2.54	0.42
1:C:113:PRO:CB	1:C:139:PHE:HB3	2.46	0.41
1:C:124:GLN:HG2	1:C:129:THR:O	2.20	0.41
1:A:207:LYS:HD3	1:A:207:LYS:HA	1.83	0.41
2:B:12:VAL:CG1	2:B:16:ARG:HB2	2.50	0.41
1:C:175:LEU:HD23	1:C:176:SER:N	2.35	0.41
1:A:134:CYS:HB2	1:A:148:TRP:CZ2	2.55	0.41
2:B:12:VAL:HG13	2:B:13:GLN:N	2.35	0.41
1:C:92:ASP:HB3	6:C:323:HOH:O	2.19	0.41
2:B:95:GLU:HB3	2:B:100:LYS:HB2	2.02	0.41
2:D:87:THR:HA	2:D:109:VAL:O	2.21	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:191:VAL:HG22	1:A:210:ASN:OD1	2.21	0.41
2:B:20:LEU:HD22	2:B:82:MET:CE	2.47	0.41
1:C:19:ALA:HB3	1:C:75:ILE:HB	2.03	0.41
2:B:34:MET:O	2:B:50:VAL:HG23	2.20	0.41
1:C:199:GLN:HB2	6:C:302:HOH:O	2.19	0.41
2:B:52:SER:HA	3:E:2:RAM:H62	2.02	0.41
2:B:84:ALA:O	2:B:87:THR:CG2	2.61	0.41
1:C:8:PRO:O	1:C:102:THR:HG23	2.20	0.41
2:B:50:VAL:CG2	2:B:51:VAL:N	2.84	0.41
1:C:12:SER:O	1:C:107:LYS:HE2	2.21	0.41
1:C:105:GLU:HG2	1:C:106:ILE:H	1.85	0.41
1:C:125:LEU:O	1:C:126:LYS:C	2.63	0.41
1:C:125:LEU:O	1:C:128:GLY:N	2.50	0.41
1:C:193:ALA:HB2	1:C:208:SER:HB3	2.03	0.41
2:D:82:MET:HE3	2:D:82(C):LEU:HD21	2.02	0.41
2:D:200:HIS:CD2	2:D:202:PRO:HD2	2.56	0.41
1:A:147:GLN:OE1	1:A:154:LEU:CD2	2.69	0.41
2:B:4:LEU:HD23	2:B:24:ALA:HB2	2.03	0.41
2:B:5:VAL:O	2:B:22:CYS:HA	2.21	0.41
1:C:143:GLU:CD	1:C:143:GLU:N	2.79	0.40
1:A:202:ARG:HA	1:A:202:ARG:HD2	1.96	0.40
1:A:147:GLN:OE1	1:A:154:LEU:HD22	2.22	0.40
2:B:34:MET:HB2	2:B:78:LEU:HD13	2.03	0.40
2:D:72:ASP:HA	6:D:457:HOH:O	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	207/219 (94%)	198 (96%)	8 (4%)	1 (0%)	24 22

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	210/219 (96%)	203 (97%)	6 (3%)	1 (0%)	24	22
2	B	197/231 (85%)	188 (95%)	9 (5%)	0	100	100
2	D	196/231 (85%)	188 (96%)	7 (4%)	1 (0%)	24	22
All	All	810/900 (90%)	777 (96%)	30 (4%)	3 (0%)	30	28

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	138	ASN
1	A	29	VAL
2	D	149	PRO

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	180/184 (98%)	155 (86%)	25 (14%)	3	2
1	C	182/184 (99%)	164 (90%)	18 (10%)	7	5
2	B	170/196 (87%)	147 (86%)	23 (14%)	4	2
2	D	169/196 (86%)	149 (88%)	20 (12%)	5	3
All	All	701/760 (92%)	615 (88%)	86 (12%)	4	2

All (86) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	LEU
1	A	13	LEU
1	A	21	LEU
1	A	22	SER
1	A	24	ARG
1	A	27	GLN
1	A	30	THR
1	A	33	LEU

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	63	SER
1	A	70	ASP
1	A	90	GLN
1	A	92	ASP
1	A	100	GLN
1	A	108	ARG
1	A	129	THR
1	A	132	VAL
1	A	149	LYS
1	A	154	LEU
1	A	168	SER
1	A	185	ASP
1	A	188	LYS
1	A	190	LYS
1	A	199	GLN
1	A	202	ARG
1	A	206	THR
2	B	2	ILE
2	B	12	VAL
2	B	20	LEU
2	B	25	SER
2	B	64	LYS
2	B	68	THR
2	B	74	SER
2	B	75	LYS
2	B	83	ARG
2	B	87	THR
2	B	98	ASP
2	B	100	LYS
2	B	100(A)	PHE
2	B	100(B)	SER
2	B	100(C)	PHE
2	B	108	LEU
2	B	138	LEU
2	B	141	LEU
2	B	149	PRO
2	B	169	VAL
2	B	178	LEU
2	B	179	SER
2	B	183	THR
1	C	13	LEU
1	C	21	LEU

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	C	29	VAL
1	C	42	GLN
1	C	56	THR
1	C	70	ASP
1	C	90	GLN
1	C	100	GLN
1	C	108	ARG
1	C	137	ASN
1	C	145	LYS
1	C	154	LEU
1	C	162	SER
1	C	168	SER
1	C	185	ASP
1	C	188	LYS
1	C	202	ARG
1	C	206	THR
2	D	12	VAL
2	D	16	ARG
2	D	31	ARG
2	D	55	ARG
2	D	56	THR
2	D	64	LYS
2	D	68	THR
2	D	74	SER
2	D	87	THR
2	D	108	LEU
2	D	141	LEU
2	D	149	PRO
2	D	169	VAL
2	D	178	LEU
2	D	182	VAL
2	D	184	VAL
2	D	185	PRO
2	D	192	GLN
2	D	201	LYS
2	D	204	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	38	GLN
1	A	42	GLN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	100	GLN
1	A	124	GLN
1	A	137	ASN
1	A	138	ASN
2	B	39	GLN
2	B	76	ASN
2	B	81	GLN
2	B	164	HIS
1	C	93	ASN
1	C	100	GLN
1	C	124	GLN
1	C	137	ASN
1	C	138	ASN
1	C	147	GLN
1	C	160	GLN
1	C	189	HIS
1	C	199	GLN
2	D	3	ASN
2	D	76	ASN

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	MBG	E	1	4,3	13,13,13	0.72	1 (7%)	18,18,18	1.31	4 (22%)
3	RAM	E	2	3	10,10,11	0.41	0	14,14,16	0.70	0
3	MBG	F	1	4,3	13,13,13	0.73	1 (7%)	18,18,18	0.94	1 (5%)
3	RAM	F	2	3	10,10,11	0.41	0	14,14,16	0.68	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
3	MBG	E	1	4,3	-	0/4/24/24	0/1/1/1
3	RAM	E	2	3	-	-	0/1/1/1
3	MBG	F	1	4,3	-	1/4/24/24	0/1/1/1
3	RAM	F	2	3	-	-	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	F	1	MBG	O1-C1	2.24	1.44	1.40
3	E	1	MBG	O1-C1	2.20	1.43	1.40

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	E	1	MBG	C7-O1-C1	-3.17	108.44	113.26
3	F	1	MBG	C7-O1-C1	-2.50	109.46	113.26
3	E	1	MBG	O1-C1-C2	2.40	110.90	108.14
3	E	1	MBG	C1-O5-C5	-2.39	109.05	113.72
3	E	1	MBG	O5-C5-C6	2.06	111.54	106.44

There are no chirality outliers.

All (1) torsion outliers are listed below:

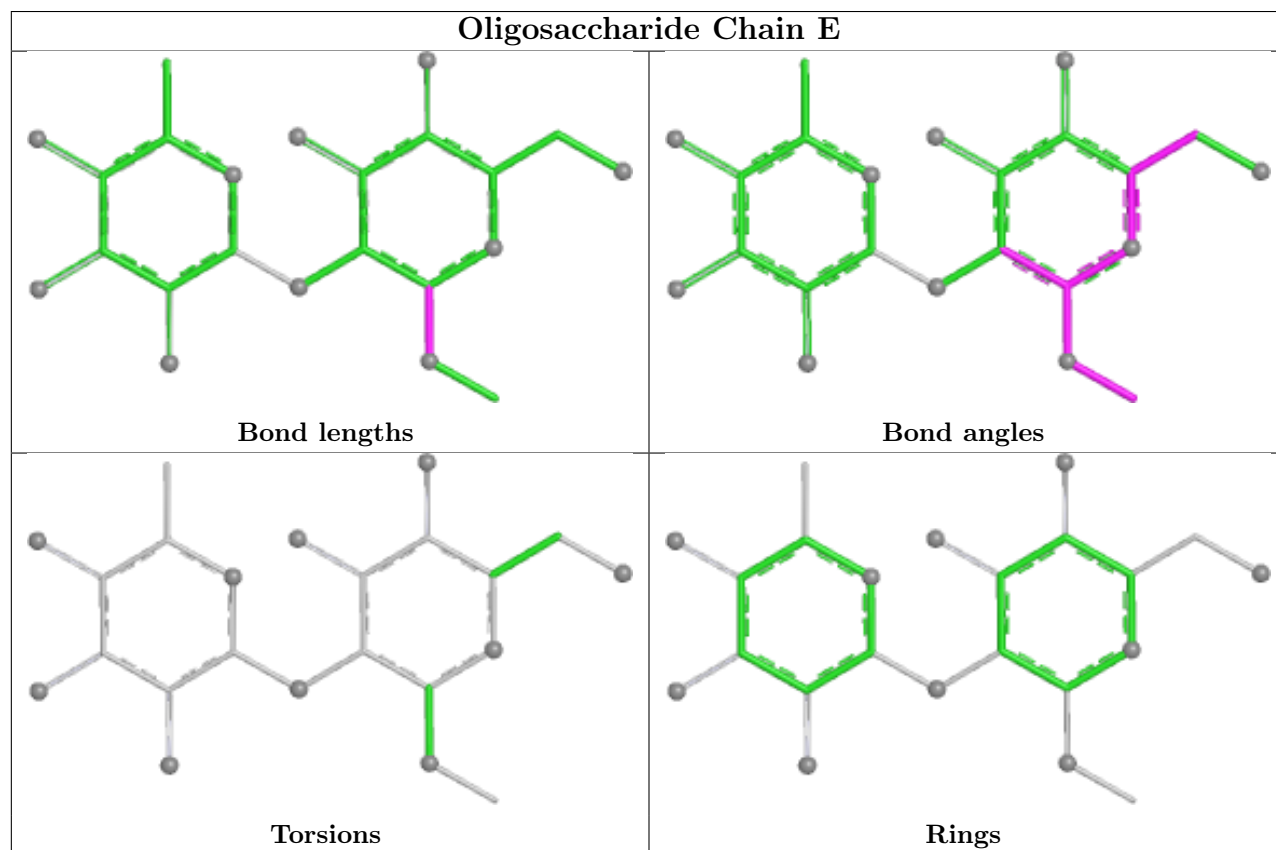
Mol	Chain	Res	Type	Atoms
3	F	1	MBG	O5-C5-C6-O6

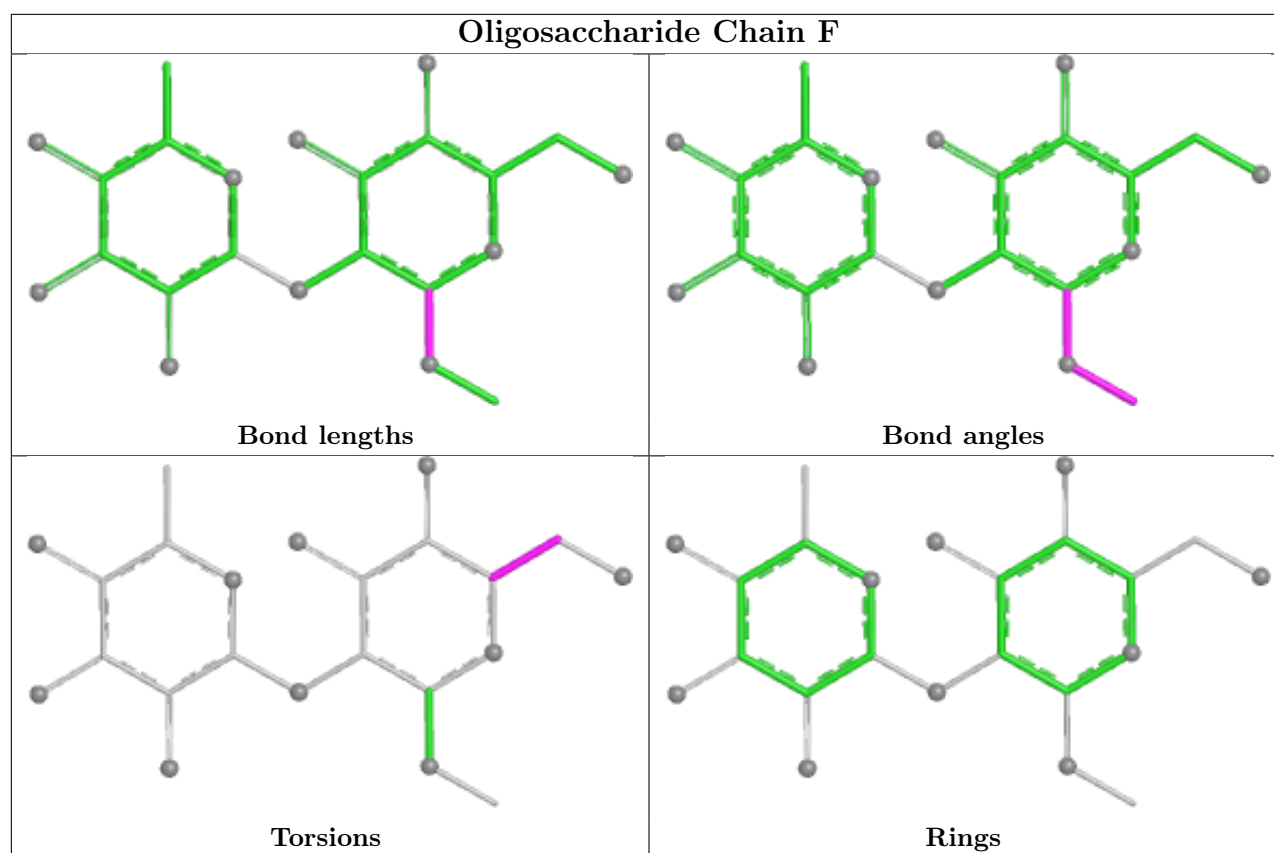
There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	E	1	MBG	1	0
3	E	2	RAM	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 [i](#)

4 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
5	GOL	D	304	4	4,4,5	1.86	1 (25%)	3,3,5	0.90	0
4	PO4	B	303	5,3	0,3,4	-	-	0,3,6	-	-
5	GOL	B	304	4	4,4,5	1.86	1 (25%)	3,3,5	1.06	0
4	PO4	D	303	5,3	0,3,4	-	-	0,3,6	-	-

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	GOL	D	304	4	-	1/2/2/4	-
5	GOL	B	304	4	-	1/2/2/4	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	D	304	GOL	O1-C1	-3.66	1.23	1.42
5	B	304	GOL	O1-C1	-3.66	1.23	1.42

There are no bond angle outliers.

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	B	304	GOL	C1-C2-C3-O3
5	D	304	GOL	O1-C1-C2-C3

There are no ring outliers.

2 monomers are involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	303	PO4	1	0
5	B	304	GOL	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	209/219 (95%)	0.46	8 (3%) 44 46	25, 40, 61, 66	0
1	C	212/219 (96%)	0.36	7 (3%) 49 51	25, 38, 58, 67	0
2	B	203/231 (87%)	0.42	4 (1%) 65 67	24, 40, 60, 66	0
2	D	202/231 (87%)	0.33	5 (2%) 58 61	26, 38, 51, 60	0
All	All	826/900 (91%)	0.39	24 (2%) 53 56	24, 39, 58, 67	0

All (24) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	29	VAL	4.4
1	C	25	ALA	4.4
2	D	196	CYS	4.1
1	A	129	THR	3.1
1	A	69	THR	3.0
2	B	3	ASN	2.7
2	B	28	THR	2.6
2	D	140	CYS	2.6
1	C	125	LEU	2.6
2	D	29	PHE	2.6
1	A	5	THR	2.6
1	A	30	THR	2.6
2	B	29	PHE	2.5
1	A	194	CYS	2.5
1	C	27	GLN	2.4
1	A	57	GLY	2.4
2	D	87	THR	2.4
1	C	94	TRP	2.3
1	C	26	SER	2.1
1	C	92	ASP	2.1
2	D	135	THR	2.1

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	C	30	THR	2.0
1	A	41	GLY	2.0
2	B	26	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](#)

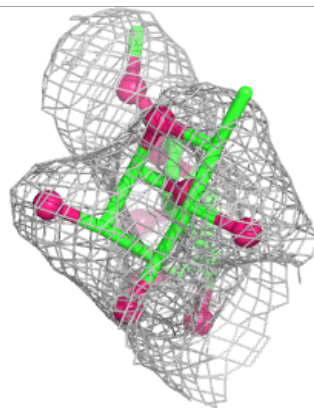
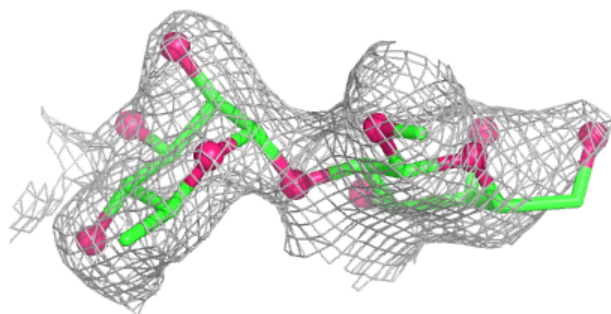
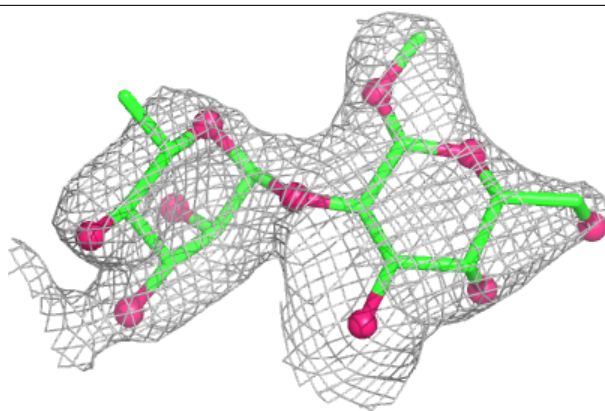
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	MBG	F	1	13/13	0.68	0.19	61,69,71,72	0
3	MBG	E	1	13/13	0.80	0.13	68,73,75,76	0
3	RAM	E	2	10/11	0.90	0.11	59,61,62,64	0
3	RAM	F	2	10/11	0.91	0.11	44,48,51,54	0

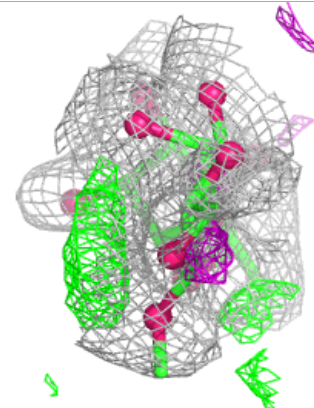
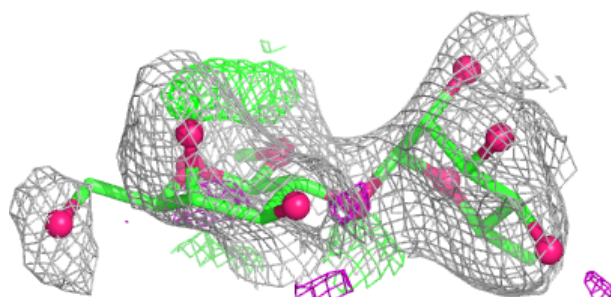
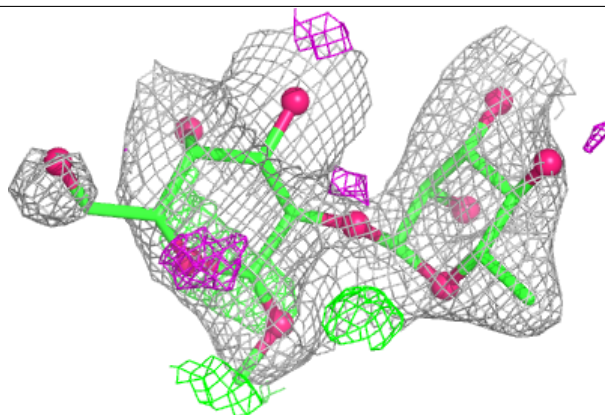
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around Chain E:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

**Electron density around Chain F:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



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(Å ²)	Q<0.9
5	GOL	B	304	5/6	0.64	0.19	83,83,84,85	0
5	GOL	D	304	5/6	0.76	0.25	77,77,78,80	0
4	PO4	D	303	4/5	0.84	0.17	74,74,74,75	0
4	PO4	B	303	4/5	0.85	0.14	78,78,79,81	0

6.5 Other polymers [i](#)

There are no such residues in this entry.