



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2026 – 06:12 PM UTC

PDB ID : 2D43 / pdb_00002d43
Title : Crystal structure of arabinofuranosidase complexed with arabinotriose
Authors : Miyanaga, A.; Koseki, T.; Miwa, Y.; Matsuzawa, H.; Wakagi, T.; Shoun, H.; Fushinobu, S.
Deposited on : 2005-10-07
Resolution : 2.80 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

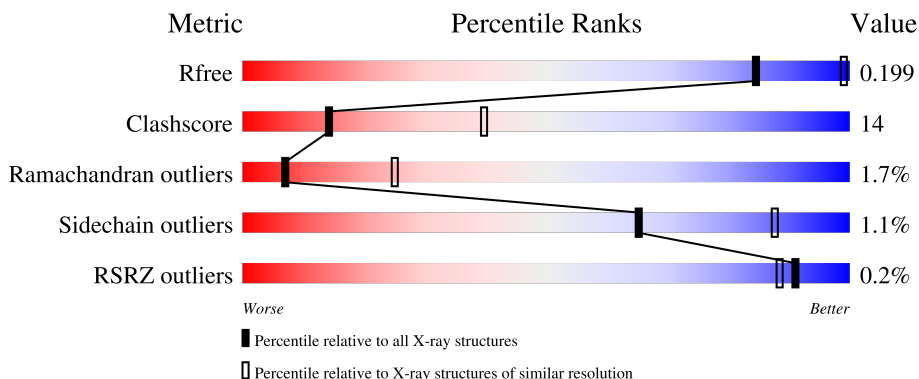
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION


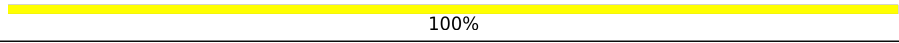
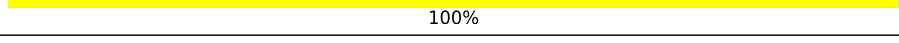
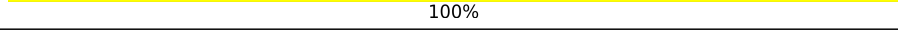
The reported resolution of this entry is 2.80 Å.

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	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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	482	 66% 33% .
2	B	2	 100%
3	C	2	 100%
3	D	2	 100%

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 3700 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 alpha-L-arabinofuranosidase B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	482	3575	2208	589	762	16	0	0	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	18	MET	-	initiating methionine	GB 21280333
A	221	ALA	GLU	engineered mutation	GB 21280333

- Molecule 2 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
2	B	2	28	16	2	10	0	0	0

- Molecule 3 is an oligosaccharide called alpha-L-arabinofuranose-(1-5)-alpha-L-arabinofuranose.



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
			Total	C	O			
3	C	2	19	10	9	0	0	0
3	D	2	19	10	9	0	0	0

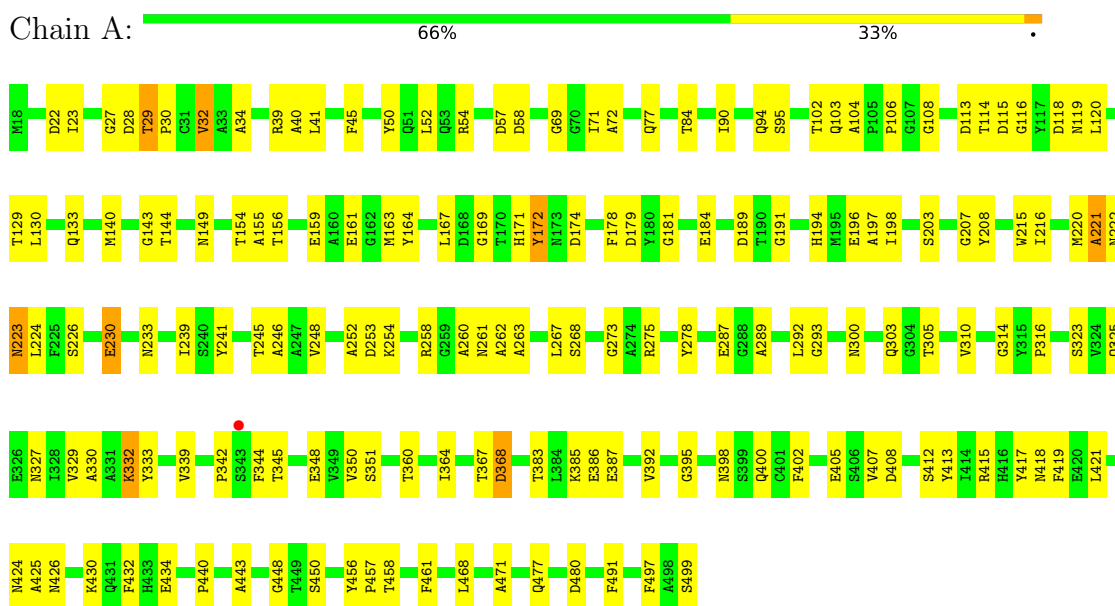
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	59	Total	O	0	0
			59	59		

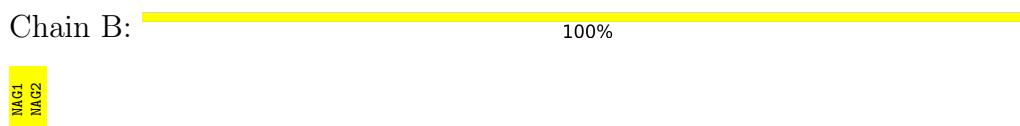
3 Residue-property plots [i](#)

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.

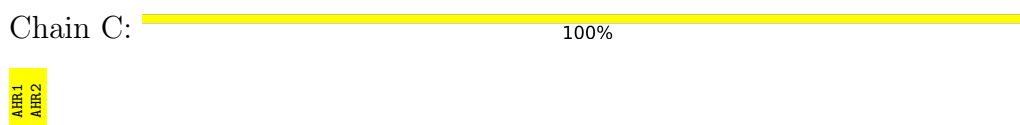
- Molecule 1: alpha-L-arabinofuranosidase B



- Molecule 2: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



- Molecule 3: alpha-L-arabinofuranose-(1-5)-alpha-L-arabinofuranose



- Molecule 3: alpha-L-arabinofuranose-(1-5)-alpha-L-arabinofuranose



AHR1
AHR2

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	39.48Å 98.75Å 144.47Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.72 – 2.80 46.72 – 2.80	Depositor EDS
% Data completeness (in resolution range)	(Not available) (46.72-2.80) 88.8 (46.72-2.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.30 (at 2.81Å)	Xtrriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.199 , 0.248 0.204 , 0.199	Depositor DCC
R_{free} test set	656 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	32.7	Xtrriage
Anisotropy	1.027	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 48.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.31$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	3700	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.04% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: NAG, AHR

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	0.42	0/3658	0.97	14/4990 (0.3%)

There are no bond length outliers.

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	32	VAL	N-CA-C	-9.64	103.74	113.47
1	A	29	THR	CA-C-N	6.93	126.92	119.78
1	A	29	THR	C-N-CA	6.93	126.92	119.78
1	A	191	GLY	N-CA-C	6.64	119.73	112.29
1	A	461	PHE	N-CA-C	-6.45	100.50	110.10
1	A	120	LEU	N-CA-C	-6.12	101.42	110.48
1	A	224	LEU	N-CA-C	-5.71	99.42	108.73
1	A	418	ASN	N-CA-C	-5.63	105.03	112.24
1	A	230	GLU	N-CA-C	-5.58	102.64	110.50
1	A	471	ALA	N-CA-C	5.33	118.79	110.32
1	A	39	ARG	N-CA-C	5.06	115.99	108.60
1	A	22	ASP	N-CA-C	-5.02	105.72	111.14
1	A	332	LYS	N-CA-C	5.01	117.29	111.02
1	A	108	GLY	N-CA-C	-5.00	108.11	114.16

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	3575	0	3250	98	0
2	B	28	0	25	0	0
3	C	19	0	0	0	0
3	D	19	0	0	0	0
4	A	59	0	0	1	0
All	All	3700	0	3275	98	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 14.

All (98) 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:415:ARG:HH11	1:A:424:ASN:HD22	1.10	0.97
1:A:260:ALA:HB2	1:A:267:LEU:HA	1.62	0.79
1:A:119:ASN:O	1:A:144:THR:HG23	1.84	0.76
1:A:415:ARG:NH1	1:A:424:ASN:HD22	1.84	0.74
1:A:253:ASP:HA	1:A:275:ARG:HG2	1.71	0.70
1:A:386:GLU:HB3	1:A:408:ASP:HB3	1.72	0.70
1:A:34:ALA:HB1	1:A:41:LEU:HD12	1.77	0.67
1:A:90:ILE:HG12	1:A:102:THR:HG22	1.77	0.67
1:A:383:THR:O	1:A:387:GLU:HG3	1.95	0.66
1:A:230:GLU:HB2	1:A:278:TYR:CZ	2.30	0.66
1:A:40:ALA:HB3	1:A:69:GLY:HA2	1.81	0.63
1:A:23:ILE:HG21	1:A:323:SER:HB2	1.82	0.62
1:A:405:GLU:HB2	1:A:413:TYR:CE1	2.34	0.62
1:A:421:LEU:HD12	1:A:468:LEU:HB3	1.82	0.62
1:A:207:GLY:O	1:A:208:TYR:HB3	2.02	0.60
1:A:348:GLU:HG2	1:A:350:VAL:HG13	1.84	0.60
1:A:477:GLN:HB2	1:A:480:ASP:OD2	2.02	0.60
1:A:164:TYR:HA	1:A:246:ALA:O	2.02	0.59
1:A:50:TYR:HE1	1:A:52:LEU:HD21	1.67	0.58
1:A:220:MET:O	1:A:221:ALA:HB3	2.03	0.58
1:A:398:ASN:OD1	1:A:400:GLN:HB2	2.04	0.57
1:A:143:GLY:HA2	1:A:300:ASN:HD22	1.68	0.57
1:A:129:THR:HA	1:A:133:GLN:O	2.04	0.56
1:A:174:ASP:O	1:A:203:SER:HB2	2.06	0.56
1:A:178:PHE:CD1	1:A:293:GLY:HA2	2.42	0.54
1:A:386:GLU:HG2	1:A:407:VAL:HG12	1.90	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:339:VAL:HG13	1:A:440:PRO:HB2	1.87	0.54
1:A:106:PRO:HB3	1:A:116:GLY:O	2.08	0.54
1:A:181:GLY:HA3	1:A:197:ALA:HA	1.91	0.53
1:A:208:TYR:HB2	1:A:215:TRP:CZ3	2.44	0.53
1:A:169:GLY:HA3	1:A:241:TYR:O	2.10	0.51
1:A:239:ILE:HG23	1:A:239:ILE:O	2.10	0.51
1:A:254:LYS:HA	1:A:273:GLY:O	2.11	0.51
1:A:103:GLN:NE2	1:A:118:ASP:O	2.38	0.51
1:A:172:TYR:HA	1:A:303:GLN:O	2.12	0.50
1:A:364:ILE:HD11	1:A:491:PHE:CE1	2.46	0.50
1:A:54:ARG:HB2	1:A:84:THR:HG21	1.92	0.50
1:A:443:ALA:HA	1:A:450:SER:OG	2.12	0.49
1:A:184:GLU:HA	1:A:287:GLU:HB2	1.95	0.49
1:A:261:ASN:OD1	1:A:263:ALA:HB3	2.13	0.48
1:A:325:GLN:HE21	1:A:325:GLN:HA	1.78	0.48
1:A:314:GLY:O	1:A:316:PRO:HD3	2.14	0.48
1:A:40:ALA:CB	1:A:69:GLY:HA2	2.44	0.48
1:A:230:GLU:HB2	1:A:278:TYR:OH	2.12	0.48
1:A:41:LEU:N	1:A:94:GLN:OE1	2.33	0.47
1:A:325:GLN:HA	1:A:325:GLN:NE2	2.29	0.47
1:A:27:GLY:O	1:A:28:ASP:HB2	2.15	0.47
1:A:497:PHE:C	1:A:499:SER:H	2.22	0.47
1:A:161:GLU:OE2	1:A:196:GLU:OE1	2.33	0.47
1:A:171:HIS:O	1:A:172:TYR:HB3	2.15	0.47
1:A:426:ASN:HA	1:A:432:PHE:CE2	2.49	0.47
1:A:45:PHE:CE2	1:A:94:GLN:HG2	2.50	0.46
1:A:57:ASP:O	1:A:58:ASP:HB2	2.14	0.46
1:A:330:ALA:O	1:A:332:LYS:HG3	2.15	0.46
1:A:385:LYS:HB3	1:A:497:PHE:HB3	1.98	0.46
1:A:171:HIS:HA	4:A:714:HOH:O	2.16	0.45
1:A:386:GLU:HG2	1:A:407:VAL:CG1	2.46	0.45
1:A:189:ASP:OD2	1:A:222:ASN:OD1	2.34	0.45
1:A:28:ASP:C	1:A:30:PRO:HD3	2.40	0.45
1:A:430:LYS:O	1:A:434:GLU:HG3	2.16	0.45
1:A:71:ILE:HG22	1:A:72:ALA:N	2.32	0.45
1:A:456:TYR:O	1:A:458:THR:N	2.49	0.45
1:A:412:SER:HA	1:A:425:ALA:HA	1.98	0.45
1:A:395:GLY:N	1:A:402:PHE:HA	2.32	0.45
1:A:178:PHE:CE1	1:A:293:GLY:HA2	2.53	0.44
1:A:448:GLY:HA3	1:A:491:PHE:O	2.17	0.44
1:A:258:ARG:HA	1:A:268:SER:O	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:156:THR:O	1:A:159:GLU:HG3	2.17	0.43
1:A:114:THR:HG22	1:A:115:ASP:OD2	2.18	0.43
1:A:220:MET:O	1:A:221:ALA:CB	2.66	0.43
1:A:216:ILE:HG13	1:A:239:ILE:HG21	1.99	0.43
1:A:345:THR:O	1:A:348:GLU:HB3	2.17	0.43
1:A:253:ASP:HA	1:A:275:ARG:CG	2.45	0.43
1:A:245:THR:OG1	1:A:262:ALA:HB2	2.18	0.43
1:A:415:ARG:HD3	1:A:415:ARG:C	2.43	0.43
1:A:327:ASN:O	1:A:330:ALA:HB3	2.17	0.43
1:A:104:ALA:HB3	1:A:118:ASP:OD2	2.19	0.43
1:A:32:VAL:HG21	1:A:154:THR:HA	2.01	0.42
1:A:41:LEU:HD13	1:A:310:VAL:CG1	2.49	0.42
1:A:179:ASP:HA	1:A:198:ILE:O	2.18	0.42
1:A:344:PHE:HB2	1:A:392:VAL:HG11	2.00	0.42
1:A:417:TYR:C	1:A:419:PHE:N	2.73	0.42
1:A:140:MET:CE	1:A:293:GLY:HA3	2.49	0.42
1:A:155:ALA:HB1	1:A:159:GLU:HB2	2.01	0.42
1:A:52:LEU:CD1	1:A:77:GLN:HG3	2.50	0.42
1:A:149:ASN:HA	1:A:289:ALA:HB1	2.01	0.42
1:A:163:MET:HG2	1:A:248:VAL:O	2.21	0.41
1:A:29:THR:N	1:A:30:PRO:HD3	2.35	0.41
1:A:194:HIS:ND1	1:A:287:GLU:OE2	2.49	0.41
1:A:325:GLN:O	1:A:329:VAL:HG23	2.20	0.41
1:A:164:TYR:CD2	1:A:310:VAL:HG22	2.56	0.41
1:A:367:THR:O	1:A:368:ASP:HB2	2.21	0.41
1:A:167:LEU:HA	1:A:305:THR:O	2.22	0.40
1:A:172:TYR:O	1:A:303:GLN:NE2	2.53	0.40
1:A:179:ASP:O	1:A:292:LEU:HA	2.20	0.40
1:A:226:SER:OG	1:A:233:ASN:HB2	2.21	0.40
1:A:130:LEU:HD11	1:A:333:TYR:CE2	2.57	0.40
1:A:344:PHE:HB3	1:A:348:GLU:OE1	2.22	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	480/482 (100%)	437 (91%)	35 (7%)	8 (2%)	7 25

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	457	PRO
1	A	223	ASN
1	A	360	THR
1	A	221	ALA
1	A	252	ALA
1	A	172	TYR
1	A	342	PRO
1	A	368	ASP

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	376/376 (100%)	372 (99%)	4 (1%)	65 87

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	95	SER
1	A	113	ASP
1	A	223	ASN
1	A	351	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	97	ASN
1	A	222	ASN

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Mol	Chain	Res	Type
1	A	233	ASN
1	A	282	ASN
1	A	300	ASN
1	A	303	GLN
1	A	325	GLN
1	A	424	ASN
1	A	466	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

6 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
2	NAG	B	1	1,2	14,14,15	2.07	3 (21%)	17,19,21	1.49	3 (17%)
2	NAG	B	2	2	14,14,15	1.59	3 (21%)	17,19,21	2.01	9 (52%)
3	AHR	C	1	3	10,10,10	1.94	4 (40%)	13,14,14	0.83	1 (7%)
3	AHR	C	2	3	9,9,10	1.65	2 (22%)	11,12,14	1.61	1 (9%)
3	AHR	D	1	3	10,10,10	1.83	4 (40%)	13,14,14	0.69	1 (7%)
3	AHR	D	2	3	9,9,10	1.89	3 (33%)	11,12,14	0.82	1 (9%)

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
2	NAG	B	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	B	2	2	-	0/6/23/26	0/1/1/1
3	AHR	C	1	3	-	0/2/18/18	0/1/1/1
3	AHR	C	2	3	-	1/2/15/18	0/1/1/1
3	AHR	D	1	3	-	2/2/18/18	0/1/1/1
3	AHR	D	2	3	-	0/2/15/18	0/1/1/1

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1	NAG	C1-C2	6.29	1.60	1.52
2	B	2	NAG	C4-C5	4.01	1.61	1.53
3	D	1	AHR	C1-C2	3.39	1.56	1.52
3	D	2	AHR	O4-C4	-3.31	1.38	1.44
3	C	2	AHR	C1-C2	3.30	1.57	1.51
3	C	1	AHR	C1-C2	3.27	1.56	1.52
3	C	1	AHR	O4-C1	3.16	1.47	1.43
3	D	2	AHR	C1-C2	2.91	1.56	1.51
3	D	2	AHR	O2-C2	2.54	1.48	1.43
2	B	2	NAG	O4-C4	2.47	1.49	1.43
3	D	1	AHR	O4-C1	2.44	1.46	1.43
3	D	1	AHR	O2-C2	2.32	1.48	1.43
3	C	1	AHR	O1-C1	2.30	1.46	1.39
2	B	2	NAG	C3-C2	2.23	1.57	1.52
3	D	1	AHR	O1-C1	2.13	1.46	1.39
2	B	1	NAG	C4-C3	2.13	1.57	1.52
3	C	1	AHR	O2-C2	2.10	1.48	1.43
3	C	2	AHR	C5-C4	2.02	1.58	1.51
2	B	1	NAG	O5-C1	2.01	1.47	1.43

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	2	AHR	O4-C4-C3	-4.65	100.34	104.63
2	B	1	NAG	C1-O5-C5	3.49	116.86	112.19
2	B	2	NAG	C6-C5-C4	3.36	121.26	113.02
2	B	2	NAG	O4-C4-C3	-3.19	102.86	110.38
2	B	1	NAG	C2-N2-C7	3.05	126.99	122.90
2	B	2	NAG	O6-C6-C5	-2.69	102.19	111.33
2	B	2	NAG	O5-C5-C6	-2.49	102.82	107.66
2	B	2	NAG	C1-O5-C5	2.43	115.45	112.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	2	NAG	C8-C7-N2	-2.16	112.53	116.12
3	D	2	AHR	O4-C4-C3	-2.14	102.65	104.63
3	D	1	AHR	C2-C3-C4	2.12	106.71	102.61
2	B	2	NAG	O4-C4-C5	2.12	114.53	109.32
3	C	1	AHR	C2-C3-C4	2.11	106.68	102.61
2	B	2	NAG	O3-C3-C4	-2.10	105.43	110.38
2	B	2	NAG	C1-C2-N2	-2.07	107.17	110.43
2	B	1	NAG	O5-C1-C2	-2.01	108.18	111.29

There are no chirality outliers.

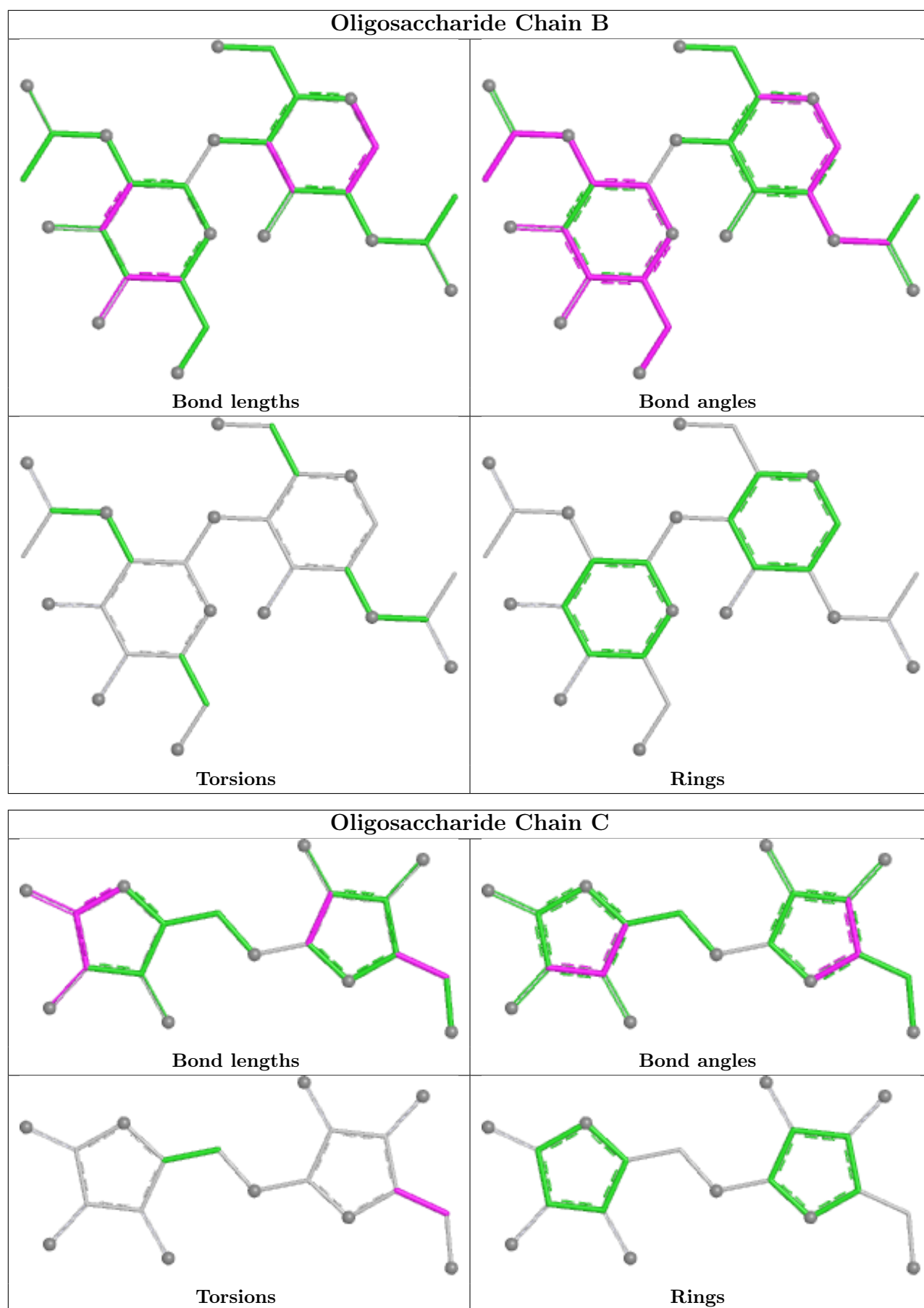
All (3) torsion outliers are listed below:

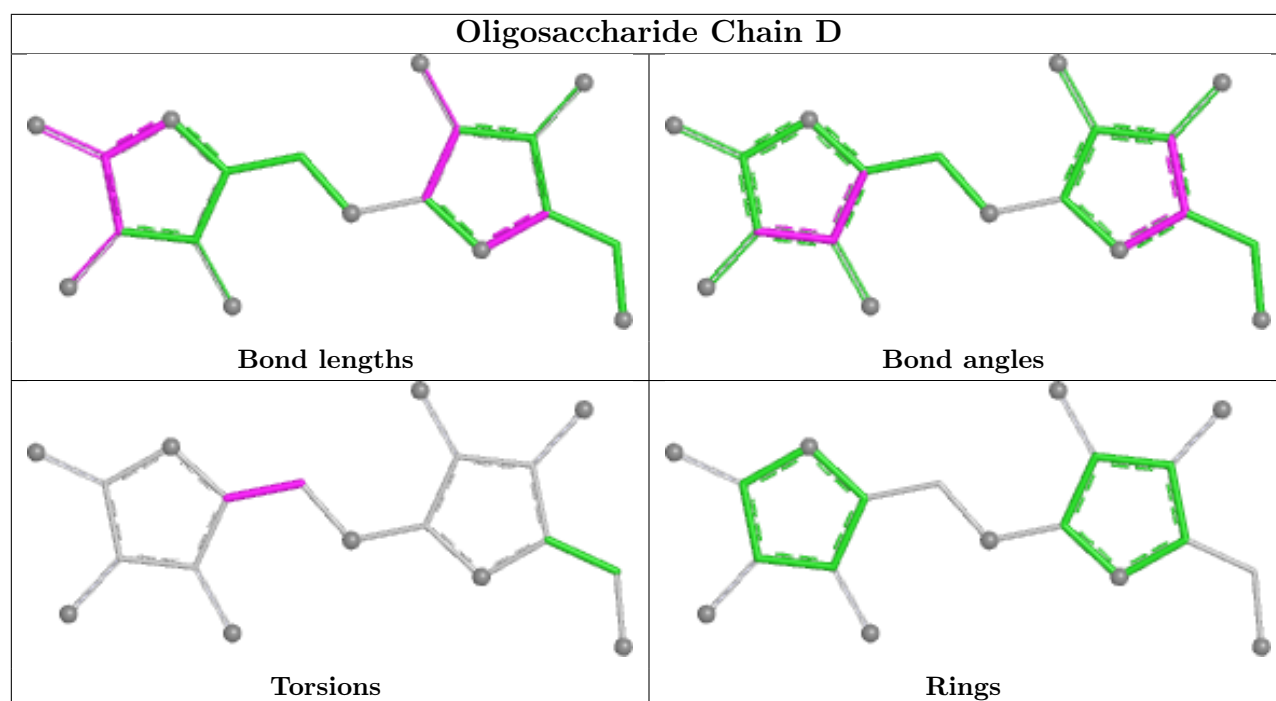
Mol	Chain	Res	Type	Atoms
3	C	2	AHR	C3-C4-C5-O5
3	D	1	AHR	C3-C4-C5-O5
3	D	1	AHR	O4-C4-C5-O5

There are no ring outliers.

No monomer is involved in short contacts.

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](#)

There are no ligands in this entry.

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 [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	482/482 (100%)	-0.43	1 (0%) 91 88	19, 30, 39, 59	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	343	SER	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](#)

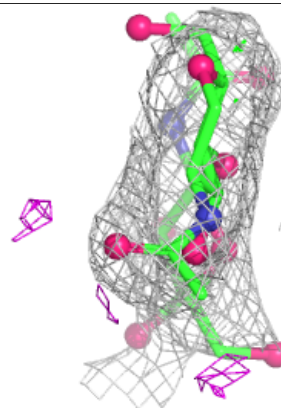
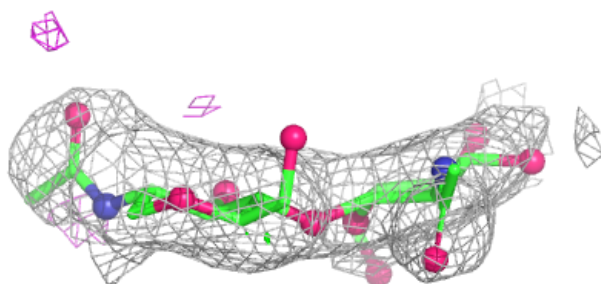
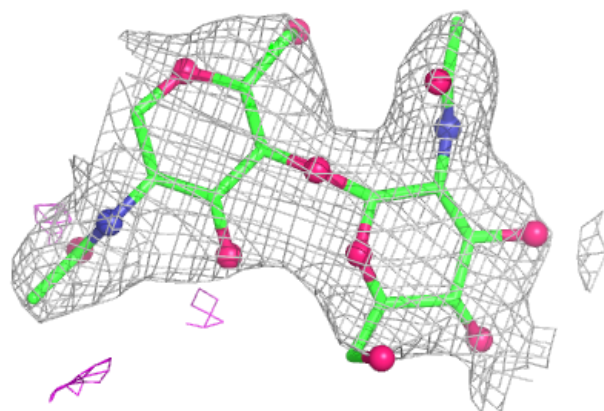
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q < 0.9’ lists 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	AHR	C	1	10/10	0.78	0.15	43,55,58,63	0
3	AHR	D	1	10/10	0.79	0.14	45,55,59,61	0
2	NAG	B	1	14/15	0.85	0.12	38,41,45,48	0
2	NAG	B	2	14/15	0.90	0.11	47,52,55,57	0
3	AHR	C	2	9/10	0.95	0.07	30,33,36,37	0
3	AHR	D	2	9/10	0.96	0.09	36,37,39,42	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.

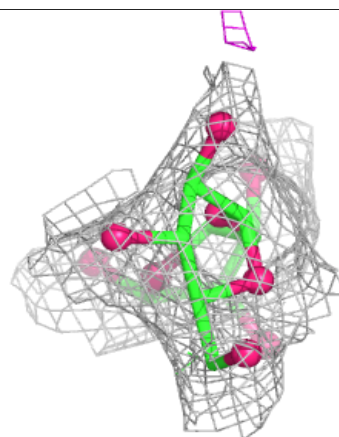
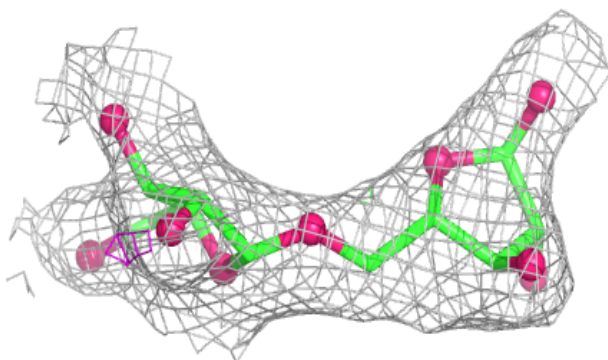
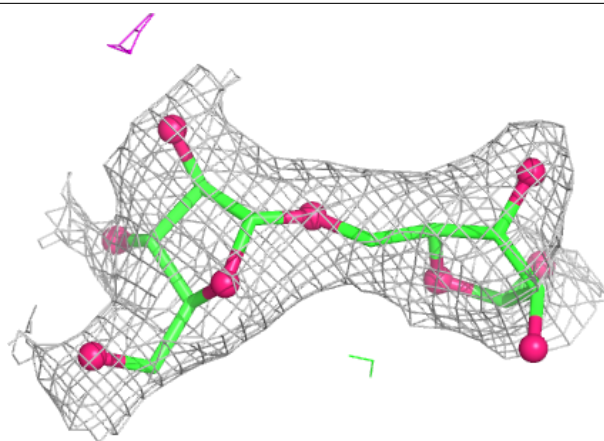
Electron density around Chain B:

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

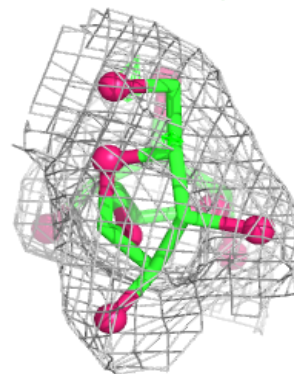
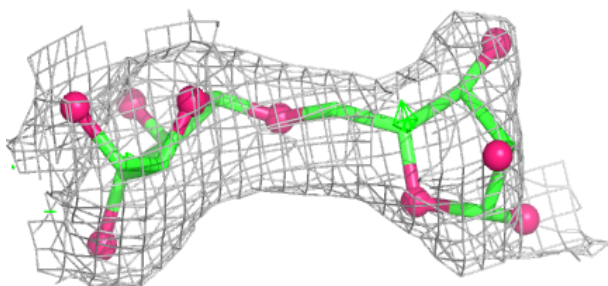
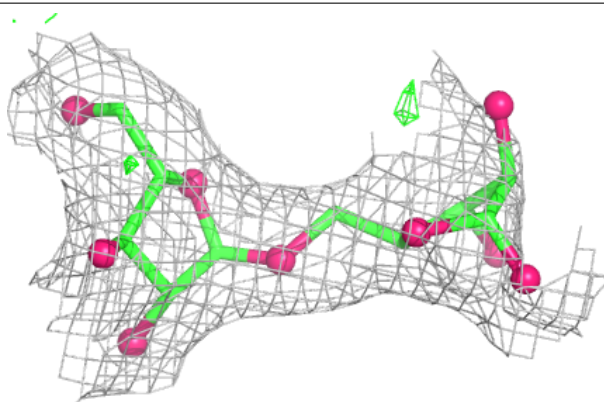


Electron density around Chain C:

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

**Electron density around Chain D:**

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



6.4 Ligands

There are no ligands in this entry.

6.5 Other polymers

There are no such residues in this entry.