W O R L D W I D E
wwPDB
PROTEIN DATA BANK

wwPDB X-ray Structure Validation Summary Report (i)

Dec 8, 2023 – 06:17 am GMT

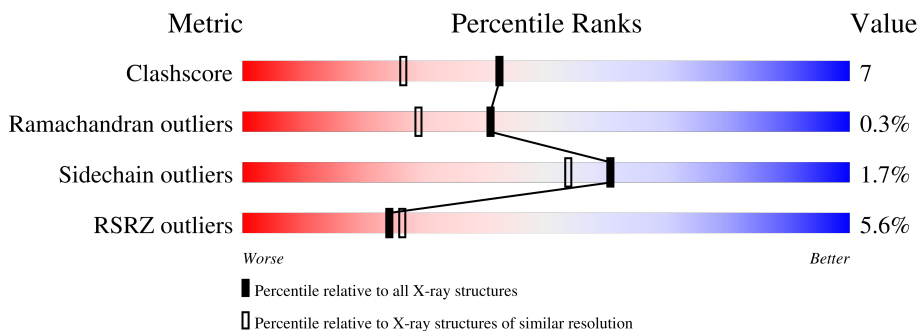|		
PDB ID	:	1OGG
Title	:	chitinase b from serratia marcescens mutant d142n in complex with inhibitor allosamidin
Authors	:	Vaaje-Kolstad, G.; Houston, D.R.; Rao, F.V.; Peter, M.G.; Synstad, B.; van Aalten, D.M.F.; Eijsink, V.G.H.
Deposited on	:	2003-04-30
Resolution	:	1.97 Å(reported)

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 1.97 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	Percentile Ranks	Value
Clashscore		7
Ramachandran outliers		0.3%
Sidechain outliers		1.7%
RSRZ outliers		5.6%

Worse *Better*

∎ Percentile relative to all X-ray structures

⧠ Percentile relative to X-ray structures of similar resolution

Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	141614	1014 (1.98-1.98)
Ramachandran outliers	138981	1006 (1.98-1.98)
Sidechain outliers	138945	1006 (1.98-1.98)
RSRZ outliers	127900	11410 (2.00-1.96)

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 <=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	499	7% / 83% / 15% / ·
1	B	499	4% / 88% / 11%
2	C	2	50% / 50%
2	D	2	50% / 50%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	GOL	B	1504	-	-	X	-

2 Entry composition ⓘ

There are 6 unique types of molecules in this entry. The entry contains 8811 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 CHITINASE B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	497	Total 3904	C 2497	N 661	O 732	S 14	23	1	1
1	B	497	Total 3919	C 2505	N 663	O 737	S 14	12	2	0

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	142	ASN	ASP	engineered mutation	UNP P11797
B	142	ASN	ASP	engineered mutation	UNP P11797
A	498	VAL	LEU	conflict	UNP P11797
B	498	VAL	LEU	conflict	UNP P11797

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

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

- Molecule 3 is ALLOSAMIZOLINE (three-letter code: AMI) (formula: $C_9H_{16}N_2O_4$).
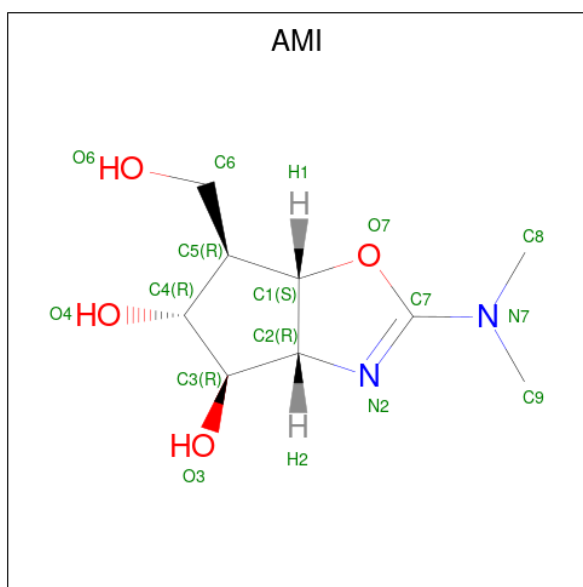
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			15	9	2	4		
3	B	1	Total	C	N	O	0	0
			15	9	2	4		

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			6	3	3		
4	A	1	Total	C	O	0	0
			6	3	3		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total C O			0	0
			6	3	3		
4	B	1	Total C O			0	0
			6	3	3		
4	B	1	Total C O			0	0
			6	3	3		
4	B	1	Total C O			0	0
			6	3	3		

- Molecule 5 is SULFATE ION (three-letter code: SO4) (formula: O_4S).
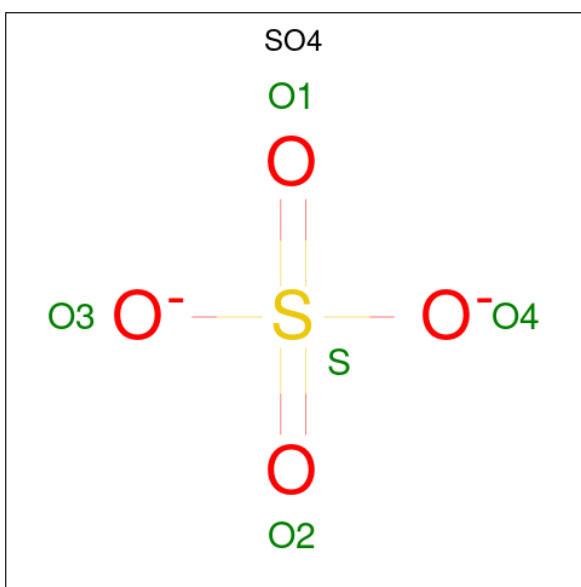
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total O S			0	0
			5	4	1		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	386	Total O		0	0
			386	386		
6	B	475	Total O		0	0
			475	475		

3 Residue-property plots ⓘ

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 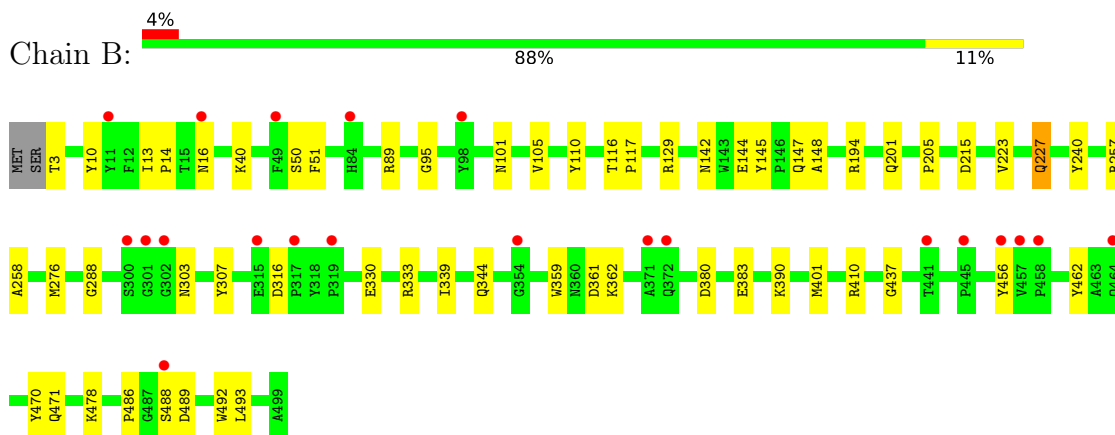or more. A red dot above a residue indicates a poor fit to the electron density (RSRZ > 2). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

- Molecule 1: CHITINASE B
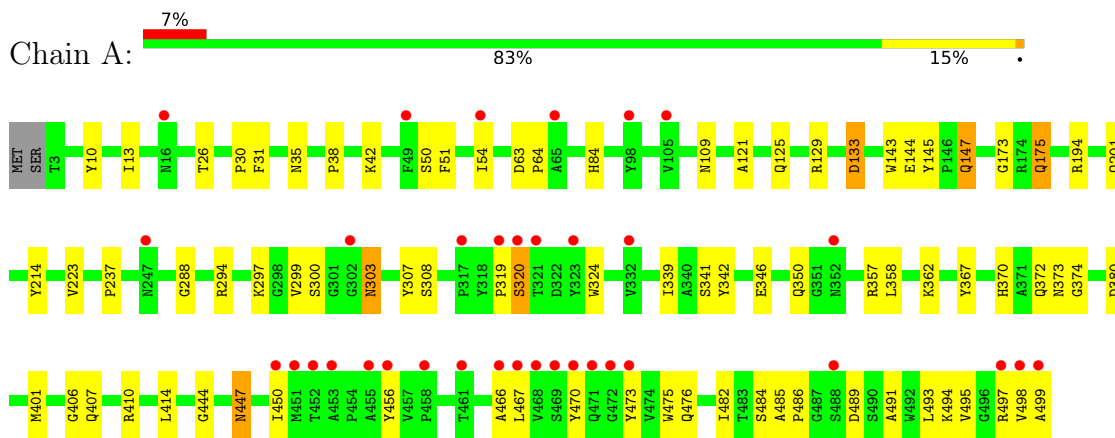
Chain A:

- Molecule 1: CHITINASE B

Chain B:

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

Chain C:

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

Chain D:

50%   50%

NAA1
NAA2

4 Data and refinement statistics ⓘ

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	55.87Å 104.74Å 186.73Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	24.76 – 1.97 25.00 – 1.97	Depositor EDS
% Data completeness (in resolution range)	95.1 (24.76-1.97) 95.2 (25.00-1.97)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.78 (at 1.98Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.198 , 0.238 0.183 , (Not available)	Depositor DCC
R_{free} test set	732 reflections (0.98%)	wwPDB-VP
Wilson B-factor (Å^2)	28.8	Xtriage
Anisotropy	0.391	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.35 , 56.7	EDS
L-test for twinning[2]	$< |L| > = 0.48$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	8811	wwPDB-VP
Average B, all atoms (Å^2)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.34% of the height of the origin peak. No significant pseudotranslation is detected.*

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[1]Intensities estimated from amplitudes.
[2]Theoretical values of $< |L| >$, $< L^2 >$ 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: GOL, AMI, SO4, NAA

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	0.54	0/4019	0.68	0/5478
1	B	57.27	2/4038 (0.0%)	1.12	6/5502 (0.1%)
All	All	40.55	2/8057 (0.0%)	0.92	6/10980 (0.1%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	361[A]	ASP	CA-CB	2573.25	58.15	1.53
1	B	361[B]	ASP	CA-CB	2573.25	58.15	1.53

The worst 5 of 6 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(º)	Ideal(º)
1	B	361[A]	ASP	CB-CA-C	-32.56	45.27	110.40
1	B	361[B]	ASP	CB-CA-C	-32.56	45.27	110.40
1	B	361[A]	ASP	CA-CB-CG	-27.11	53.76	113.40
1	B	361[B]	ASP	CA-CB-CG	-27.11	53.76	113.40
1	B	361[A]	ASP	N-CA-CB	-17.53	79.04	110.60

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	3904	0	3739	65	0
1	B	3919	0	3752	41	0
2	C	28	0	25	0	0
2	D	28	0	25	0	0
3	A	15	0	15	2	0
3	B	15	0	15	2	0
4	A	18	0	24	2	0
4	B	18	0	24	6	0
5	A	5	0	0	0	0
6	A	386	0	0	5	0
6	B	475	0	0	10	0
All	All	8811	0	7619	107	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 7.

The worst 5 of 107 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:444:GLY:H	1:A:447:ASN:HD21	1.08	0.95
1:A:294:ARG:NH1	1:A:339:ILE:HD11	1.84	0.91
1:A:175:GLN:H	1:A:175:GLN:HE21	1.20	0.88
1:B:330:GLU:HA	1:B:333[B]:ARG:NH2	1.89	0.85
1:B:89:ARG:HH12	4:B:1504:GOL:H11	1.44	0.81

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	496/499 (99%)	485 (98%)	9 (2%)	2 (0%)	34	22
1	B	497/499 (100%)	487 (98%)	9 (2%)	1 (0%)	47	38

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	993/998 (100%)	972 (98%)	18 (2%)	3 (0%)	41 29

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	498	VAL
1	B	148	ALA
1	A	320	SER

5.3.2   Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	403/405 (100%)	396 (98%)	7 (2%)	60 53
1	B	405/405 (100%)	398 (98%)	7 (2%)	60 53
All	All	808/810 (100%)	794 (98%)	14 (2%)	60 53

5 of 14 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	227	GLN
1	B	240	TYR
1	B	410	ARG
1	B	339	ILE
1	B	380	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 18 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	273	GLN
1	B	394	GLN
1	B	344	GLN
1	B	35	ASN
1	B	227	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

4 monosaccharides are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#|Z| > 2$	Counts	RMSZ	$\#|Z| > 2$
2	NAA	C	1	3,2	14,14,15	1.01	0	17,19,21	0.92	1 (5%)
2	NAA	C	2	2	14,14,15	0.93	0	17,19,21	0.71	0
2	NAA	D	1	3,2	14,14,15	0.99	0	17,19,21	0.89	0
2	NAA	D	2	2	14,14,15	1.00	1 (7%)	17,19,21	0.82	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
2	NAA	C	1	3,2	-	0/6/23/26	0/1/1/1
2	NAA	C	2	2	-	0/6/23/26	0/1/1/1
2	NAA	D	1	3,2	-	0/6/23/26	0/1/1/1
2	NAA	D	2	2	-	0/6/23/26	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	2	NAA	C1-C2	2.31	1.55	1.52

All (1) bond angle outliers are listed below:

| Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	C	1	NAA	C1-C2-N2	-2.41	106.37	110.49

There are no chirality outliers.

There are no torsion outliers.

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.

Oligosaccharide Chain C

Bond lengths

Bond angles

Torsions

Rings

Oligosaccharide Chain D

Bond lengths | Bond angles

Torsions | Rings
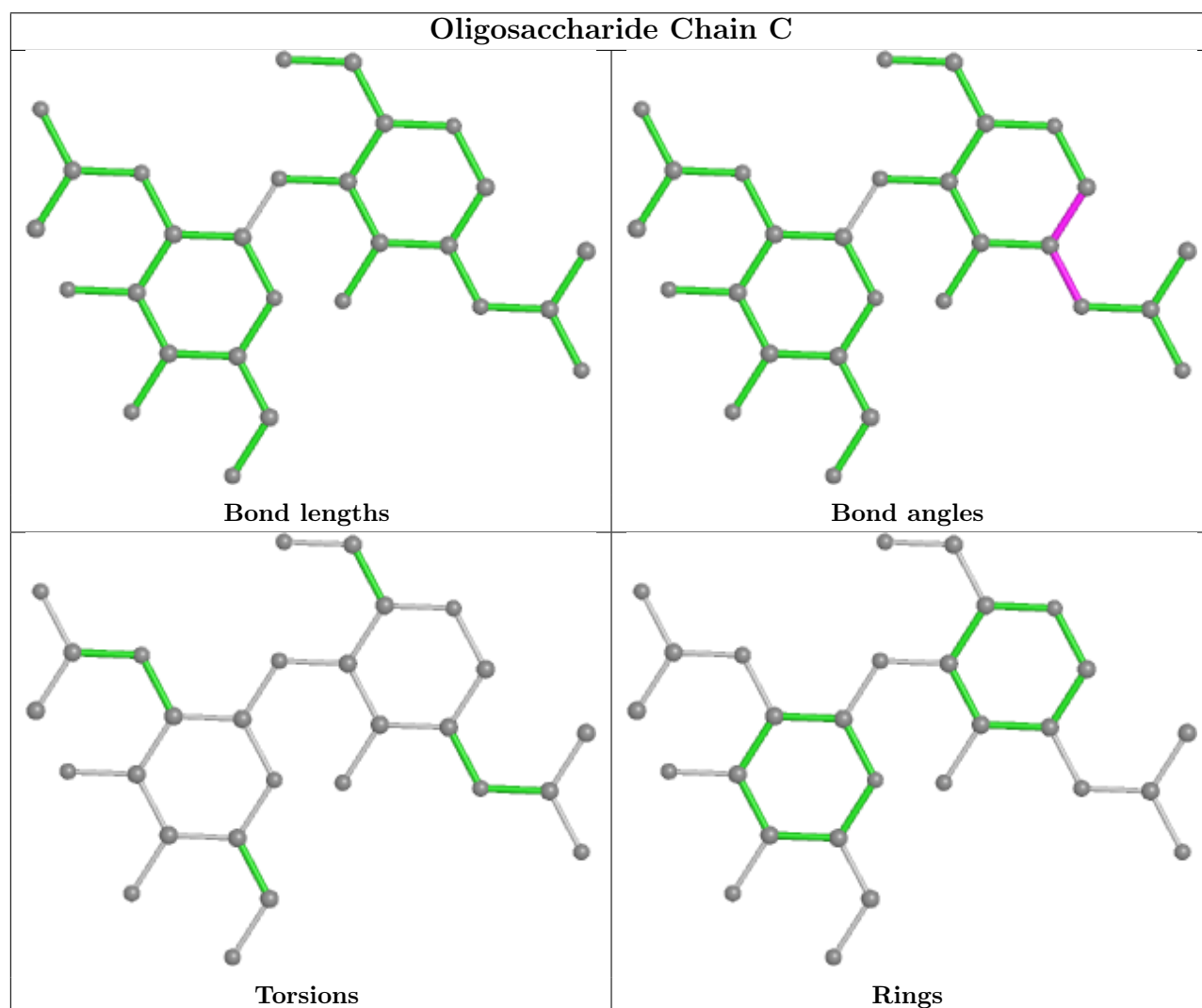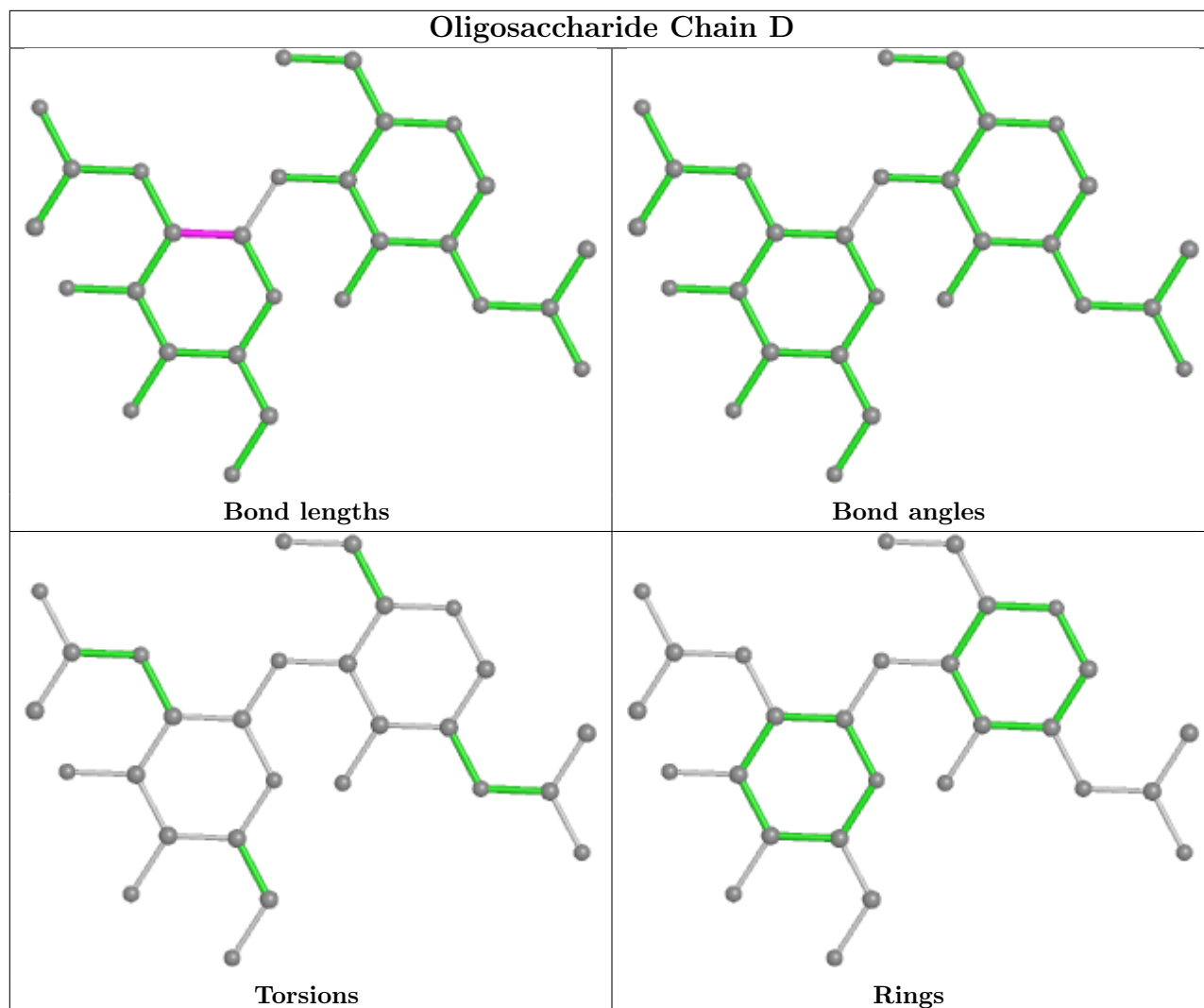
5.6 Ligand geometry ⓘ

9 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	SO4	A	1505	-	4,4,4	0.30	0	6,6,6	0.12	0
4	GOL	A	1503	-	5,5,5	0.48	0	5,5,5	0.12	0
3	AMI	A	1501	2	13,16,16	0.98	1 (7%)	16,24,24	2.09	4 (25%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	#|Z| > 2	Counts	RMSZ	#|Z| > 2
4	GOL	B	1506	-	5,5,5	0.48	0	5,5,5	0.12	0
4	GOL	B	1504	-	5,5,5	0.41	0	5,5,5	0.25	0
3	AMI	B	1502	2	13,16,16	0.85	0	16,24,24	2.15	3 (18%)
4	GOL	B	1503	-	5,5,5	0.37	0	5,5,5	0.28	0
4	GOL	A	1502	-	5,5,5	0.36	0	5,5,5	0.21	0
4	GOL	A	1504	-	5,5,5	0.46	0	5,5,5	0.24	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
4	GOL	A	1503	-	-	0/4/4/4	-
3	AMI	A	1501	2	-	0/6/34/34	0/2/2/2
4	GOL	B	1506	-	-	0/4/4/4	-
4	GOL	B	1504	-	-	0/4/4/4	-
3	AMI	B	1502	2	-	0/6/34/34	0/2/2/2
4	GOL	B	1503	-	-	0/4/4/4	-
4	GOL	A	1502	-	-	0/4/4/4	-
4	GOL	A	1504	-	-	0/4/4/4	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1501	AMI	O7-C7	2.10	1.37	1.34

The worst 5 of 7 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	A	1501	AMI	O7-C7-N2	-5.79	114.73	117.80
3	B	1502	AMI	O7-C7-N2	-5.50	114.88	117.80
3	B	1502	AMI	C5-C4-C3	-4.52	97.99	103.81
3	A	1501	AMI	C5-C4-C3	-3.56	99.23	103.81
3	B	1502	AMI	C1-C2-N2	-2.71	101.96	104.44

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

6 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	1503	GOL	1	0
3	A	1501	AMI	2	0
4	B	1504	GOL	5	0
3	B	1502	AMI	2	0
4	B	1503	GOL	1	0
4	A	1504	GOL	1	0

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(Å²)	Q<0.9
1	A	497/499 (99%)	0.37	35 (7%)	16	18	17, 32, 56, 71	7 (1%)
1	B	497/499 (99%)	0.19	21 (4%)	36	38	18, 28, 48, 58	6 (1%)
All	All	994/998 (99%)	0.28	56 (5%)	24	26	17, 30, 52, 71	13 (1%)

The worst 5 of 56 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	499	ALA	20.0
1	A	498	VAL	6.6
1	A	450	ILE	6.4
1	A	452	THR	4.4
1	B	445	PRO	3.9

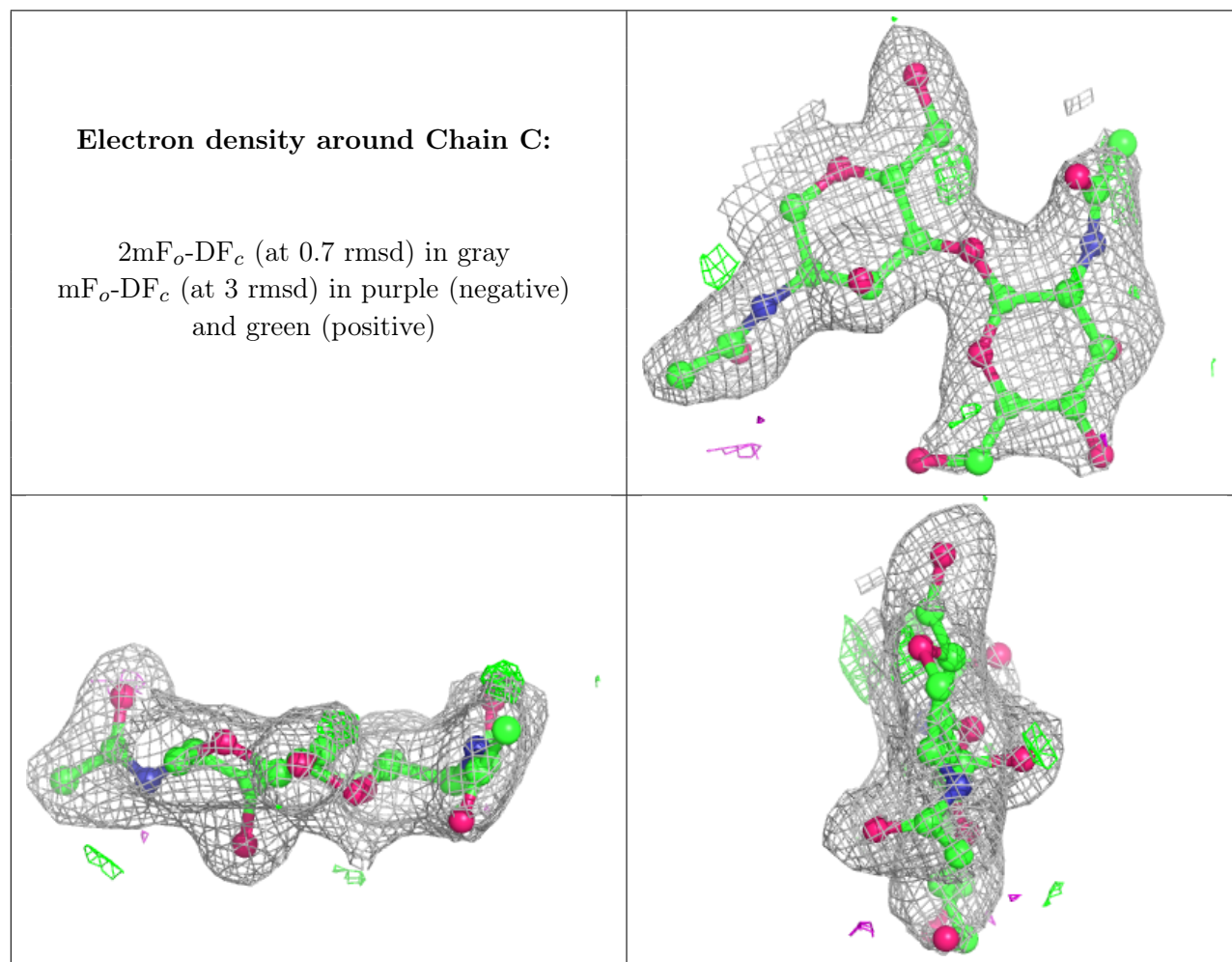## 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

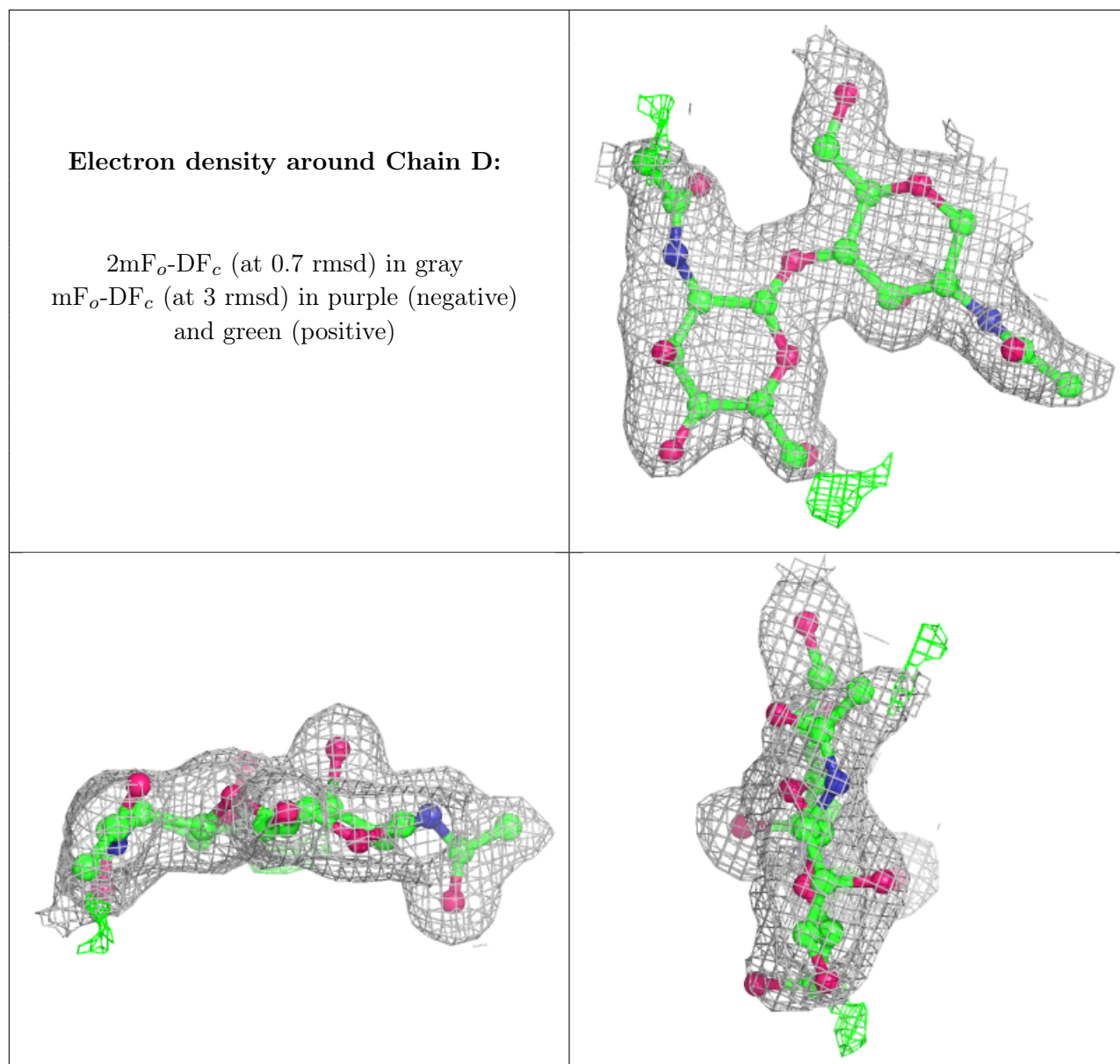There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

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(Å²)	Q<0.9
2	NAA	C	2	14/15	0.88	0.21	44,50,54,54	0
2	NAA	D	2	14/15	0.88	0.18	42,48,51,53	0
2	NAA	C	1	14/15	0.96	0.08	26,29,32,39	0
2	NAA	D	1	14/15	0.97	0.08	20,26,28,32	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 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:

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

6.4 Ligands ⓘ

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	GOL	A	1503	6/6	0.80	0.21	57,60,61,62	0
4	GOL	B	1504	6/6	0.82	0.23	51,53,53,54	0
4	GOL	A	1504	6/6	0.88	0.17	52,54,55,55	0
4	GOL	A	1502	6/6	0.89	0.17	30,38,41,42	0
4	GOL	B	1503	6/6	0.93	0.16	41,44,46,46	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
4	GOL	B	1506	6/6	0.93	0.12	33,36,37,38	0
5	SO4	A	1505	5/5	0.95	0.16	86,87,87,87	0
3	AMI	B	1502	15/15	0.97	0.10	18,21,25,26	0
3	AMI	A	1501	15/15	0.97	0.09	22,24,27,30	0

6.5 Other polymers ⓘ

There are no such residues in this entry.