



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 6, 2022 – 12:30 PM EDT

PDB ID : 8DC9
Title : RNA ligase RtcB from *Pyrococcus horikoshii* in complex with Mn²⁺ and GTP
Authors : Jacewicz, A.; Dantuluri, S.; Shuman, S.
Deposited on : 2022-06-16
Resolution : 2.47 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 : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.31.2
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.31.2

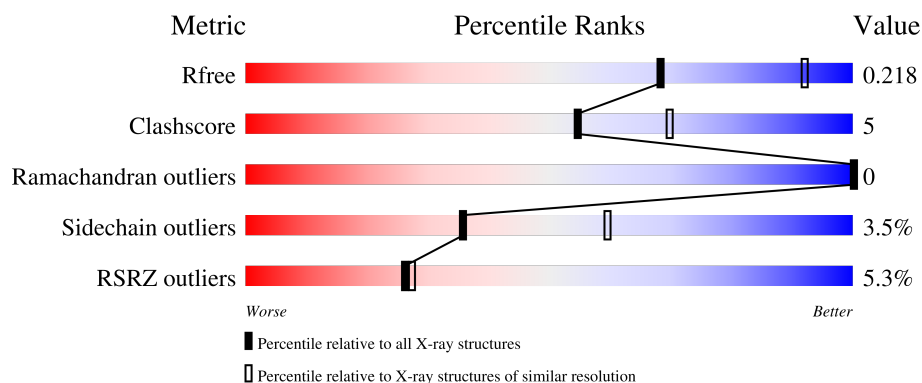
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.47 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}	130704	5857 (2.50-2.46)
Clashscore	141614	6594 (2.50-2.46)
Ramachandran outliers	138981	6469 (2.50-2.46)
Sidechain outliers	138945	6471 (2.50-2.46)
RSRZ outliers	127900	5738 (2.50-2.46)

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	501	<div> <div>5%</div> <div> <div></div> <div>80%</div> <div>16%</div> <div>.</div> </div> </div>
1	B	501	<div> <div>5%</div> <div> <div></div> <div>81%</div> <div>14%</div> <div>..</div> </div> </div>
2	C	2	<div> <div></div> <div>100%</div> </div>
2	D	2	<div> <div></div> <div>100%</div> </div>

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 7766 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 tRNA-splicing ligase RtcB.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	481	Total	C	N	O	S	0	0	0
			3767	2375	694	681	17			
1	B	481	Total	C	N	O	S	0	0	0
			3767	2375	694	681	17			

There are 40 discrepancies between the modelled and reference sequences:

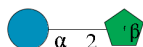
Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	initiating methionine	UNP O59245
A	-18	GLY	-	expression tag	UNP O59245
A	-17	SER	-	expression tag	UNP O59245
A	-16	SER	-	expression tag	UNP O59245
A	-15	HIS	-	expression tag	UNP O59245
A	-14	HIS	-	expression tag	UNP O59245
A	-13	HIS	-	expression tag	UNP O59245
A	-12	HIS	-	expression tag	UNP O59245
A	-11	HIS	-	expression tag	UNP O59245
A	-10	HIS	-	expression tag	UNP O59245
A	-9	SER	-	expression tag	UNP O59245
A	-8	SER	-	expression tag	UNP O59245
A	-7	GLY	-	expression tag	UNP O59245
A	-6	LEU	-	expression tag	UNP O59245
A	-5	VAL	-	expression tag	UNP O59245
A	-4	PRO	-	expression tag	UNP O59245
A	-3	ARG	-	expression tag	UNP O59245
A	-2	GLY	-	expression tag	UNP O59245
A	-1	SER	-	expression tag	UNP O59245
A	0	HIS	-	expression tag	UNP O59245
B	-19	MET	-	initiating methionine	UNP O59245
B	-18	GLY	-	expression tag	UNP O59245
B	-17	SER	-	expression tag	UNP O59245
B	-16	SER	-	expression tag	UNP O59245
B	-15	HIS	-	expression tag	UNP O59245

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-14	HIS	-	expression tag	UNP O59245
B	-13	HIS	-	expression tag	UNP O59245
B	-12	HIS	-	expression tag	UNP O59245
B	-11	HIS	-	expression tag	UNP O59245
B	-10	HIS	-	expression tag	UNP O59245
B	-9	SER	-	expression tag	UNP O59245
B	-8	SER	-	expression tag	UNP O59245
B	-7	GLY	-	expression tag	UNP O59245
B	-6	LEU	-	expression tag	UNP O59245
B	-5	VAL	-	expression tag	UNP O59245
B	-4	PRO	-	expression tag	UNP O59245
B	-3	ARG	-	expression tag	UNP O59245
B	-2	GLY	-	expression tag	UNP O59245
B	-1	SER	-	expression tag	UNP O59245
B	0	HIS	-	expression tag	UNP O59245

- Molecule 2 is an oligosaccharide called beta-D-fructofuranose-(2-1)-alpha-D-glucopyranose.



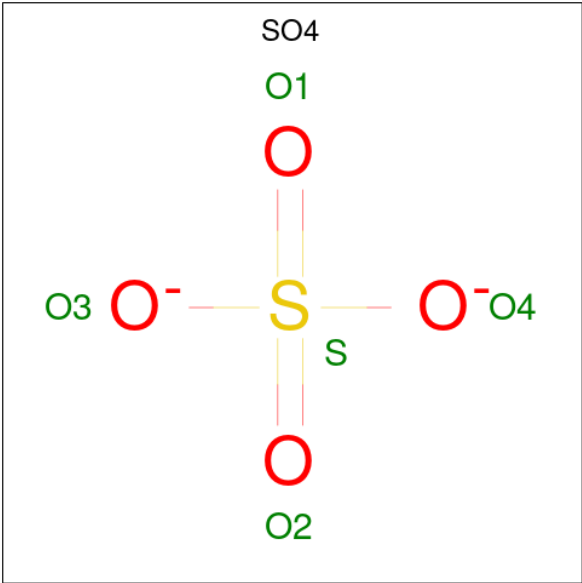
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
2	C	2	Total	C	O	0	0	0
			23	12	11			
2	D	2	Total	C	O	0	0	0
			23	12	11			

- Molecule 3 is GUANOSINE-5'-TRIPHOSPHATE (three-letter code: GTP) (formula: C₁₀H₁₆N₅O₁₄P₃) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total 32	C 10	N 5	O 14	P 3	0	0
3	B	1	Total 32	C 10	N 5	O 14	P 3	2	0

- Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



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

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

- Molecule 5 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	2	Total Mn 2 2	0	0
5	B	2	Total Mn 2 2	0	0

- Molecule 6 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	1	Total 1	Cl 1	0	0
6	B	2	Total 2	Cl 2	0	0

- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	9	Total 9	O 9	0	0
7	B	11	Total 11	O 11	0	0

GLC1
FRU2

- Molecule 2: beta-D-fructofuranose-(2-1)-alpha-D-glucopyranose

Chain D:

100%

GLC1
FRU2

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	80.74Å 137.28Å 150.16Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.39 – 2.47 49.39 – 2.47	Depositor EDS
% Data completeness (in resolution range)	100.0 (49.39-2.47) 100.0 (49.39-2.47)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.07 (at 2.48Å)	Xtriage
Refinement program	PHENIX 1.20.1_4487	Depositor
R, R_{free}	0.183 , 0.218 0.183 , 0.218	Depositor DCC
R_{free} test set	2979 reflections (4.91%)	wwPDB-VP
Wilson B-factor (Å ²)	59.1	Xtriage
Anisotropy	0.391	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 54.0	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	0.96	EDS
Total number of atoms	7766	wwPDB-VP
Average B, all atoms (Å ²)	61.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.15% 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: CL, FRU, GLC, GTP, SO4, MN

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	0.43	0/3844	0.65	0/5185
1	B	0.43	0/3844	0.66	0/5185
All	All	0.43	0/7688	0.65	0/10370

There are no bond length outliers.

There are no bond angle outliers.

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	3767	0	3794	44	0
1	B	3767	0	3794	35	0
2	C	23	0	21	1	0
2	D	23	0	21	1	0
3	A	32	0	12	4	0
3	B	32	0	12	0	0
4	A	55	0	0	3	0
4	B	40	0	0	1	0
5	A	2	0	0	0	0
5	B	2	0	0	0	0
6	A	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	B	2	0	0	0	0
7	A	9	0	0	1	0
7	B	11	0	0	1	0
All	All	7766	0	7654	80	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 5.

All (80) 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:B:336:GLU:OE2	1:B:343:ARG:HD2	1.91	0.71
4:A:512:SO4:O1	7:A:601:HOH:O	2.09	0.70
1:A:202:ASN:OD1	1:A:451:TYR:OH	2.10	0.69
1:B:217:GLU:OE2	1:B:217:GLU:N	2.24	0.68
4:B:504:SO4:O3	7:B:601:HOH:O	2.11	0.68
1:A:442:VAL:O	1:A:446:GLU:HG2	1.94	0.68
1:B:412:ARG:NH2	1:B:446:GLU:OE2	2.28	0.66
1:A:250:ARG:HG2	1:A:250:ARG:HH11	1.62	0.65
1:B:202:ASN:OD1	1:B:451:TYR:OH	2.12	0.64
1:A:141:TRP:O	1:A:144:ILE:HG13	1.97	0.64
1:B:6:LYS:HD2	1:B:16:PRO:HG3	1.79	0.63
1:A:257:ARG:NH2	4:A:508:SO4:O1	2.32	0.63
1:A:100:VAL:HG23	1:A:329:HIS:HE1	1.65	0.62
1:A:5:LEU:HD13	1:A:35:LEU:HD21	1.84	0.59
1:B:58:LYS:O	2:D:2:FRU:H11	2.04	0.58
1:A:381:MET:HB2	1:A:478:VAL:HG21	1.87	0.57
1:A:202:ASN:HB2	3:A:501:GTP:O1B	2.06	0.56
1:A:198:LEU:HA	1:A:235:THR:HG21	1.88	0.56
1:B:261:ILE:HD11	1:B:272:PRO:HG3	1.87	0.56
1:A:238:ARG:O	1:A:242:HIS:HD2	1.88	0.56
1:A:200:SER:OG	1:A:201:GLY:N	2.38	0.56
1:A:234:HIS:CE1	3:A:501:GTP:H5'	2.41	0.55
1:A:205:LEU:HD12	1:A:233:VAL:HG22	1.87	0.55
1:B:97:ASN:HD22	1:B:289:ALA:HB1	1.72	0.54
1:B:88:SER:HB2	1:B:270:SER:HB3	1.90	0.53
1:A:294:ALA:O	1:A:298:MET:HG3	2.09	0.52
1:A:100:VAL:HG23	1:A:329:HIS:CE1	2.43	0.52
1:A:441:ARG:O	1:A:445:GLU:HG3	2.10	0.52
1:A:203:HIS:CD2	1:A:235:THR:HA	2.44	0.52
1:A:66:GLY:HA2	1:A:74:ILE:O	2.11	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:66:GLY:HA2	1:B:74:ILE:O	2.11	0.51
1:A:442:VAL:HG11	1:A:481:GLY:O	2.11	0.50
1:A:23:ARG:NH1	4:A:510:SO4:O1	2.45	0.49
1:B:72:PHE:CG	1:B:91:GLY:HA2	2.48	0.49
1:A:409:VAL:HG13	1:A:410:LEU:HD12	1.94	0.49
1:A:58:LYS:HA	2:C:2:FRU:H11	1.95	0.48
1:B:424:ILE:HA	1:B:427:GLU:OE1	2.13	0.48
1:B:379:GLY:HA3	1:B:384:ALA:O	2.13	0.48
1:B:205:LEU:CD1	1:B:233:VAL:HG22	2.43	0.48
1:B:224:LEU:HA	1:B:228:GLN:OE1	2.13	0.48
1:A:144:ILE:CD1	1:A:192:ALA:HB2	2.44	0.47
1:B:264:PRO:HG2	1:B:268:LEU:HD12	1.96	0.47
1:A:161:GLY:HA3	1:A:297:GLN:HE21	1.81	0.46
1:B:5:LEU:HD13	1:B:35:LEU:HD21	1.96	0.46
1:A:365:PRO:HB2	1:A:367:LEU:HD23	1.97	0.45
1:A:480:LYS:HE2	3:A:501:GTP:O6	2.16	0.45
1:A:104:ARG:O	1:A:321:MET:HA	2.17	0.45
1:B:325:TYR:CE2	1:B:327:VAL:HG13	2.52	0.45
1:B:104:ARG:O	1:B:321:MET:HA	2.17	0.45
1:A:125:LYS:O	1:B:134:GLN:NE2	2.48	0.44
1:B:53:LEU:HB2	1:B:80:PHE:HE2	1.81	0.44
1:B:54:PRO:HG3	1:B:262:PRO:HG2	2.00	0.44
1:B:7:ARG:HD2	1:B:13:TRP:CH2	2.53	0.43
1:A:202:ASN:HB2	3:A:501:GTP:PB	2.59	0.43
1:A:225:PHE:CZ	1:A:228:GLN:HB2	2.53	0.43
1:B:215:ASP:OD2	1:B:468:LYS:HE3	2.18	0.43
1:B:282:PHE:CE2	1:B:286:LYS:HE3	2.54	0.43
1:B:480:LYS:HB2	1:B:480:LYS:HE2	1.69	0.43
1:A:282:PHE:CE2	1:A:286:LYS:HE3	2.53	0.43
1:A:308:GLN:HB2	1:A:315:PRO:HG3	2.01	0.43
1:B:11:ILE:HD13	1:B:464:ALA:HB2	2.01	0.43
1:B:119:LEU:HA	1:B:310:VAL:HG21	2.01	0.43
1:A:420:ARG:HE	1:A:421:GLY:H	1.65	0.43
1:B:91:GLY:O	1:B:349:HIS:HB3	2.18	0.42
1:B:117:LYS:HG2	1:B:434:TYR:CE1	2.55	0.42
1:B:376:LEU:O	1:B:377:ILE:HD13	2.20	0.42
1:A:119:LEU:HA	1:A:310:VAL:HG21	2.01	0.42
1:A:250:ARG:HG2	1:A:250:ARG:NH1	2.31	0.41
1:A:333:LYS:HE2	1:A:333:LYS:HB3	1.84	0.41
1:A:164:ARG:NH1	1:A:322:ASP:OD1	2.53	0.41
1:A:8:ILE:HD11	1:A:14:GLU:HB2	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:427:GLU:O	1:A:431:ARG:HG3	2.21	0.41
1:A:420:ARG:HG3	1:A:422:ASP:H	1.85	0.41
1:B:332:GLY:HA2	1:B:348:VAL:O	2.20	0.41
1:B:442:VAL:HG11	1:B:481:GLY:O	2.19	0.41
1:A:224:LEU:HA	1:A:228:GLN:OE1	2.21	0.41
1:A:433:ILE:HG12	1:A:475:PRO:O	2.21	0.40
1:B:205:LEU:HD12	1:B:205:LEU:HA	1.93	0.40
1:B:237:SER:OG	1:B:241:GLY:HA3	2.21	0.40
1:A:63:MET:HB3	1:A:64:PRO:HD2	2.04	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	479/501 (96%)	463 (97%)	16 (3%)	0	100	100
1	B	479/501 (96%)	463 (97%)	16 (3%)	0	100	100
All	All	958/1002 (96%)	926 (97%)	32 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	390/407 (96%)	377 (97%)	13 (3%)	38	61
1	B	390/407 (96%)	376 (96%)	14 (4%)	35	58
All	All	780/814 (96%)	753 (96%)	27 (4%)	36	59

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	59	TYR
1	A	110	LYS
1	A	129	SER
1	A	133	SER
1	A	153	LYS
1	A	176	GLU
1	A	266	ARG
1	A	285	MET
1	A	408	ARG
1	A	423	ARG
1	A	439	SER
1	A	440	MET
1	B	45	GLU
1	B	59	TYR
1	B	104	ARG
1	B	246	SER
1	B	256	ILE
1	B	257	ARG
1	B	327	VAL
1	B	346	VAL
1	B	362	GLU
1	B	366	ARG
1	B	409	VAL
1	B	411	SER
1	B	412	ARG
1	B	439	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	242	HIS
1	A	329	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 ⓘ

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	GLC	C	1	2	11,11,12	0.53	0	15,15,17	2.62	3 (20%)
2	FRU	C	2	2	11,12,12	0.78	0	10,18,18	0.83	0
2	GLC	D	1	2	11,11,12	0.51	0	15,15,17	1.76	1 (6%)
2	FRU	D	2	2	11,12,12	0.60	0	10,18,18	0.76	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	GLC	C	1	2	-	2/2/19/22	0/1/1/1
2	FRU	C	2	2	-	5/5/24/24	0/1/1/1
2	GLC	D	1	2	-	2/2/19/22	0/1/1/1
2	FRU	D	2	2	-	3/5/24/24	0/1/1/1

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	1	GLC	C1-O5-C5	8.08	123.14	112.19
2	D	1	GLC	C1-O5-C5	5.79	120.04	112.19
2	C	1	GLC	O5-C5-C6	4.76	114.67	107.20
2	C	1	GLC	C2-C3-C4	-2.31	106.89	110.89

There are no chirality outliers.

All (12) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	2	FRU	O1-C1-C2-C3
2	C	2	FRU	O1-C1-C2-O2
2	C	2	FRU	O1-C1-C2-O5
2	D	2	FRU	O1-C1-C2-C3
2	D	2	FRU	O1-C1-C2-O2
2	C	1	GLC	O5-C5-C6-O6
2	D	1	GLC	O5-C5-C6-O6
2	C	1	GLC	C4-C5-C6-O6
2	D	1	GLC	C4-C5-C6-O6
2	C	2	FRU	O5-C5-C6-O6
2	D	2	FRU	O1-C1-C2-O5
2	C	2	FRU	C4-C5-C6-O6

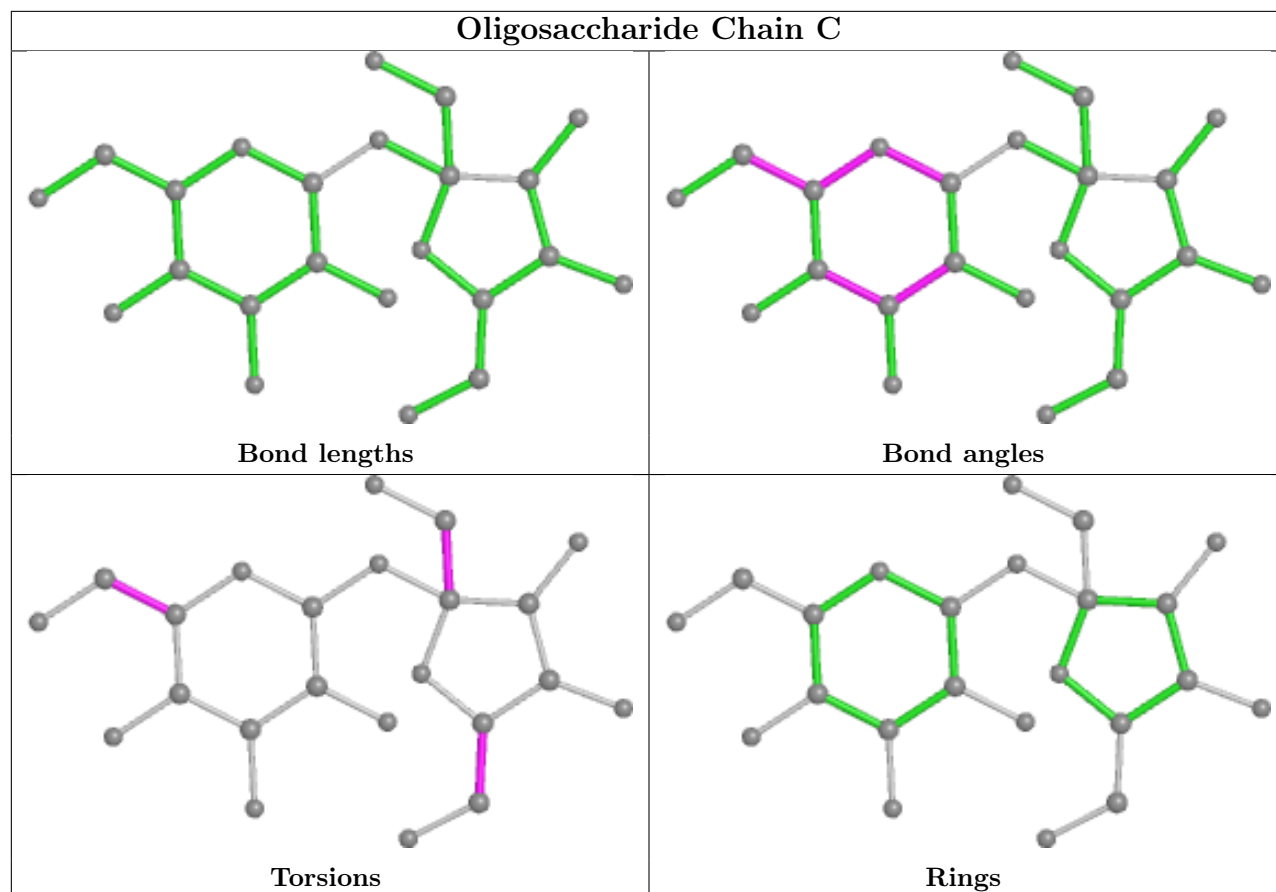
There are no ring outliers.

2 monomers are involved in 2 short contacts:

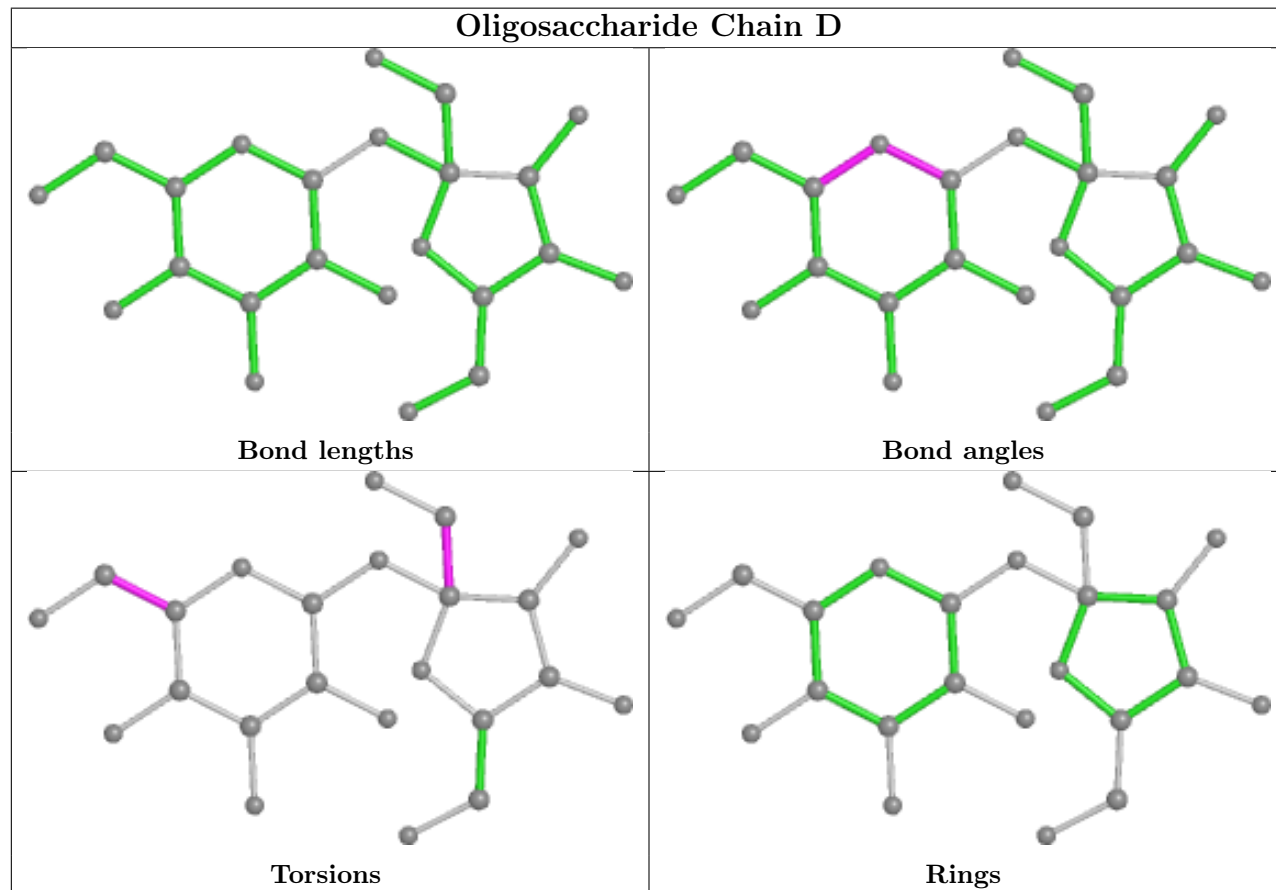
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	2	FRU	1	0
2	D	2	FRU	1	0

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

Oligosaccharide Chain C



Oligosaccharide Chain D



5.6 Ligand geometry

Of 28 ligands modelled in this entry, 7 are monoatomic - leaving 21 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	SO4	B	504	-	4,4,4	0.13	0	6,6,6	0.11	0
4	SO4	A	505	-	4,4,4	0.19	0	6,6,6	0.12	0
4	SO4	A	511	-	4,4,4	0.12	0	6,6,6	0.19	0
4	SO4	B	502	-	4,4,4	0.17	0	6,6,6	0.49	0
4	SO4	A	503	-	4,4,4	0.18	0	6,6,6	0.24	0
4	SO4	A	506	-	4,4,4	0.12	0	6,6,6	0.17	0
4	SO4	A	509	-	4,4,4	0.14	0	6,6,6	0.15	0
4	SO4	B	507	-	4,4,4	0.10	0	6,6,6	0.30	0
4	SO4	A	508	-	4,4,4	0.10	0	6,6,6	0.20	0
4	SO4	B	508	-	4,4,4	0.20	0	6,6,6	0.24	0
4	SO4	B	509	-	4,4,4	0.17	0	6,6,6	0.16	0
3	GTP	B	501	5	26,34,34	1.14	2 (7%)	32,54,54	1.83	6 (18%)
4	SO4	A	504	-	4,4,4	0.15	0	6,6,6	0.09	0
3	GTP	A	501	5	26,34,34	1.15	2 (7%)	32,54,54	1.43	5 (15%)
4	SO4	A	510	-	4,4,4	0.15	0	6,6,6	0.15	0
4	SO4	B	505	-	4,4,4	0.16	0	6,6,6	0.14	0
4	SO4	B	503	-	4,4,4	0.23	0	6,6,6	0.58	0
4	SO4	A	502	-	4,4,4	0.15	0	6,6,6	0.48	0
4	SO4	B	506	-	4,4,4	0.13	0	6,6,6	0.18	0
4	SO4	A	512	-	4,4,4	0.13	0	6,6,6	0.28	0
4	SO4	A	507	-	4,4,4	0.22	0	6,6,6	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
3	GTP	B	501	5	-	4/18/38/38	0/3/3/3
3	GTP	A	501	5	-	8/18/38/38	0/3/3/3

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	501	GTP	C5-C6	-3.83	1.39	1.47
3	B	501	GTP	C5-C6	-3.80	1.39	1.47
3	B	501	GTP	C2-N3	2.30	1.38	1.33
3	A	501	GTP	C2-N3	2.22	1.38	1.33

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	501	GTP	PB-O3B-PG	-6.29	111.24	132.83
3	A	501	GTP	PB-O3B-PG	-3.50	120.83	132.83
3	B	501	GTP	C8-N7-C5	3.26	109.20	102.99
3	A	501	GTP	C8-N7-C5	3.20	109.09	102.99
3	B	501	GTP	C3'-C2'-C1'	3.20	105.80	100.98
3	B	501	GTP	C5-C6-N1	3.15	119.52	113.95
3	A	501	GTP	C5-C6-N1	3.01	119.27	113.95
3	B	501	GTP	C2-N1-C6	-2.73	120.06	125.10
3	B	501	GTP	PA-O3A-PB	-2.61	123.86	132.83
3	A	501	GTP	C2-N1-C6	-2.43	120.63	125.10
3	A	501	GTP	C3'-C2'-C1'	2.27	104.40	100.98

There are no chirality outliers.

All (12) torsion outliers are listed below:

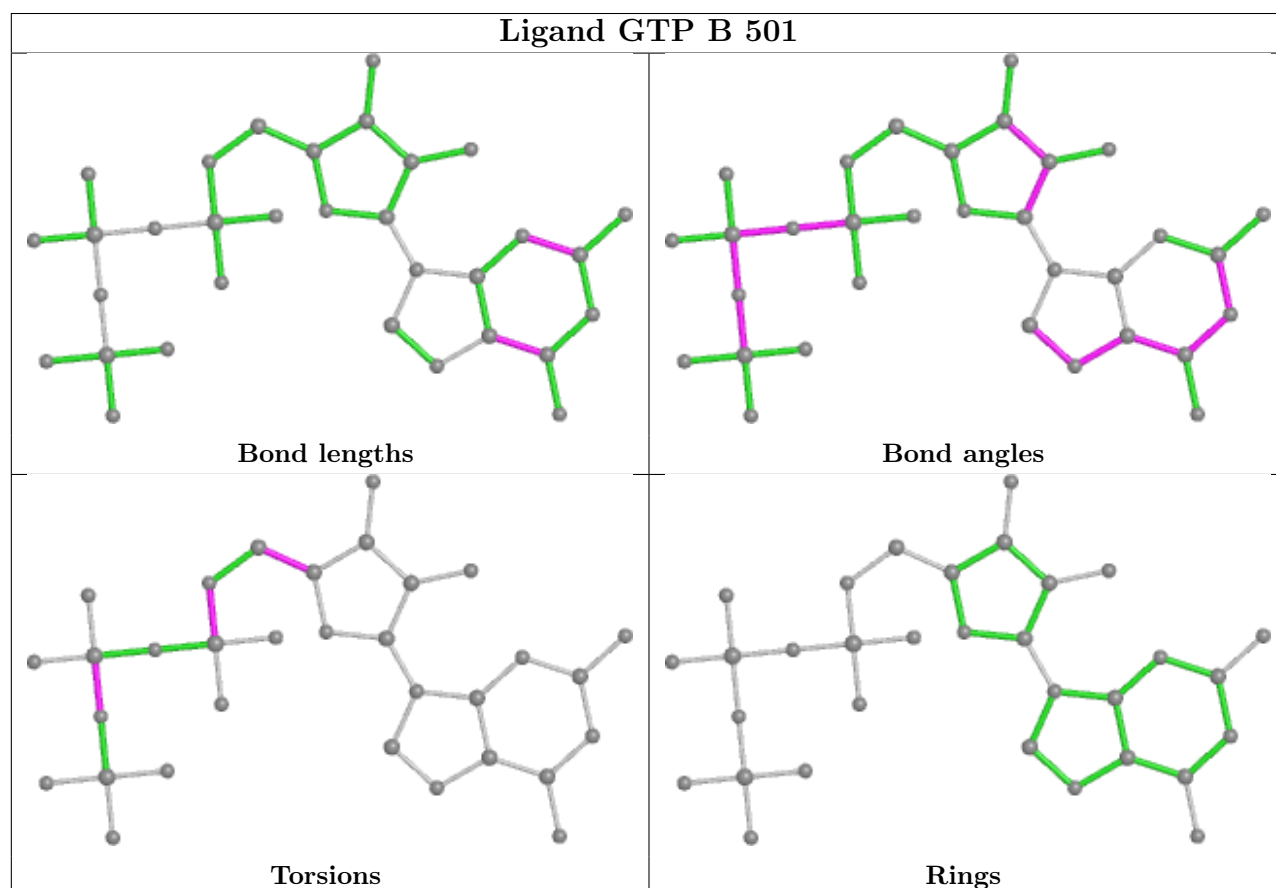
Mol	Chain	Res	Type	Atoms
3	A	501	GTP	C5'-O5'-PA-O3A
3	B	501	GTP	O4'-C4'-C5'-O5'
3	B	501	GTP	C3'-C4'-C5'-O5'
3	A	501	GTP	O4'-C4'-C5'-O5'
3	A	501	GTP	C3'-C4'-C5'-O5'
3	A	501	GTP	PB-O3A-PA-O1A
3	A	501	GTP	PG-O3B-PB-O1B
3	A	501	GTP	C5'-O5'-PA-O1A
3	A	501	GTP	PB-O3A-PA-O2A
3	B	501	GTP	PG-O3B-PB-O1B
3	A	501	GTP	PG-O3B-PB-O2B
3	B	501	GTP	C5'-O5'-PA-O1A

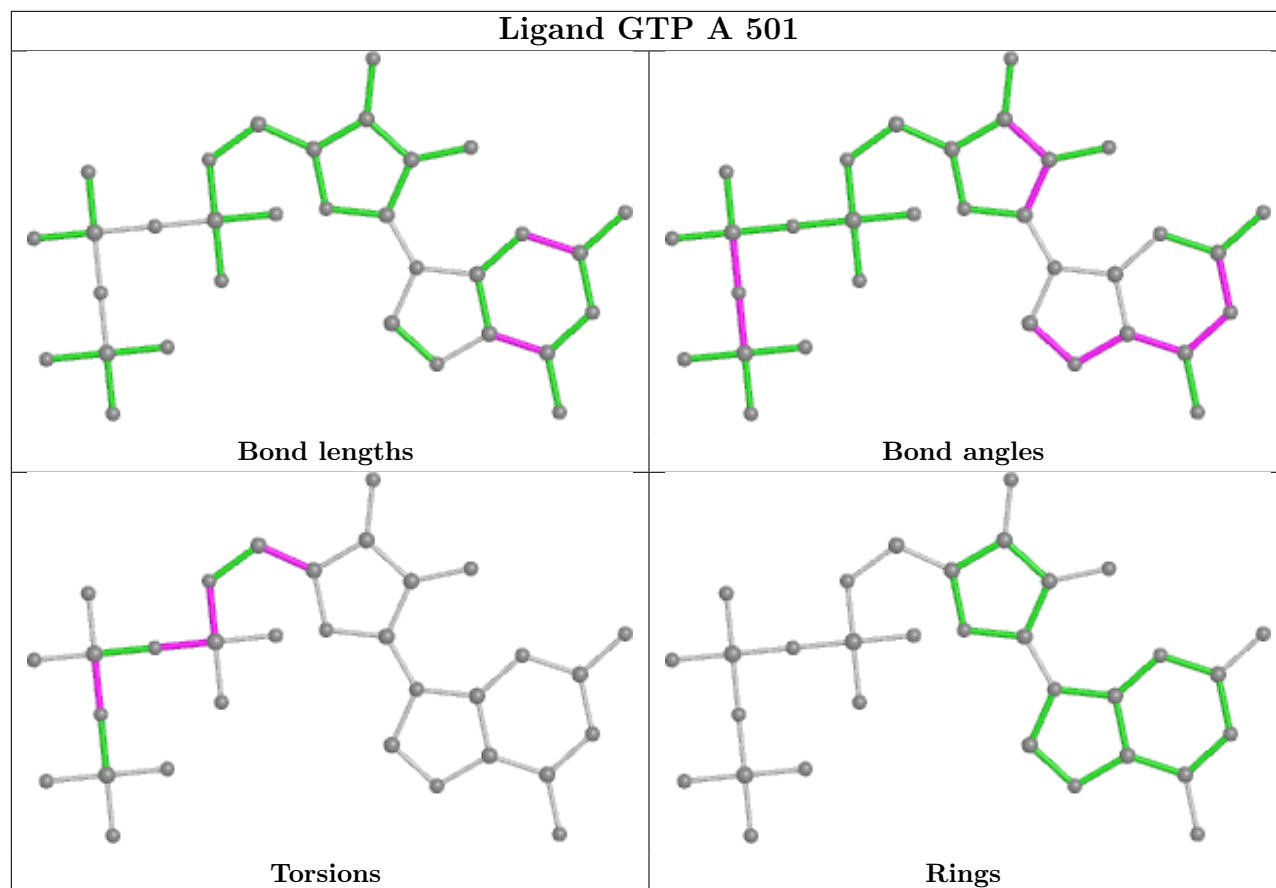
There are no ring outliers.

5 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	504	SO4	1	0
4	A	508	SO4	1	0
3	A	501	GTP	4	0
4	A	510	SO4	1	0
4	A	512	SO4	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 all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





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	481/501 (96%)	0.17	25 (5%)	27 28	43, 55, 93, 123	0
1	B	481/501 (96%)	0.20	26 (5%)	25 26	43, 57, 88, 117	0
All	All	962/1002 (96%)	0.19	51 (5%)	26 27	43, 56, 90, 123	0

All (51) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	141	TRP	4.9
1	A	419	TYR	4.9
1	A	1	MET	4.7
1	B	138	LYS	4.4
1	B	423	ARG	4.3
1	B	142	THR	4.0
1	A	413	LYS	3.7
1	A	343	ARG	3.6
1	B	263	TRP	3.6
1	A	431	ARG	3.6
1	A	423	ARG	3.5
1	B	419	TYR	3.4
1	B	430	ASN	3.3
1	B	417	ARG	3.3
1	A	430	ASN	3.2
1	B	2	VAL	3.2
1	B	1	MET	3.2
1	B	256	ILE	3.1
1	B	343	ARG	2.8
1	B	429	LEU	2.8
1	B	426	GLN	2.7
1	A	481	GLY	2.7
1	B	140	HIS	2.7
1	A	136	ARG	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	431	ARG	2.6
1	A	415	ALA	2.6
1	A	342	LYS	2.6
1	A	426	GLN	2.5
1	A	410	LEU	2.5
1	B	157	ASP	2.4
1	A	263	TRP	2.4
1	A	422	ASP	2.4
1	B	410	LEU	2.4
1	A	340	ASP	2.3
1	B	418	GLN	2.3
1	A	427	GLU	2.3
1	B	250	ARG	2.3
1	A	157	ASP	2.2
1	A	2	VAL	2.2
1	A	418	GLN	2.2
1	B	428	LEU	2.2
1	A	292	ALA	2.1
1	A	341	GLY	2.1
1	A	20	LYS	2.1
1	B	424	ILE	2.1
1	B	420	ARG	2.1
1	B	413	LYS	2.1
1	A	42	ARG	2.1
1	B	362	GLU	2.1
1	B	139	LEU	2.0
1	A	185	GLN	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(\AA^2)	Q<0.9
2	GLC	C	1	11/12	0.81	0.20	71,76,80,82	11

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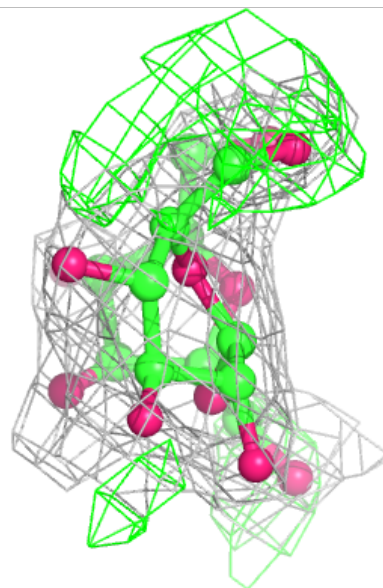
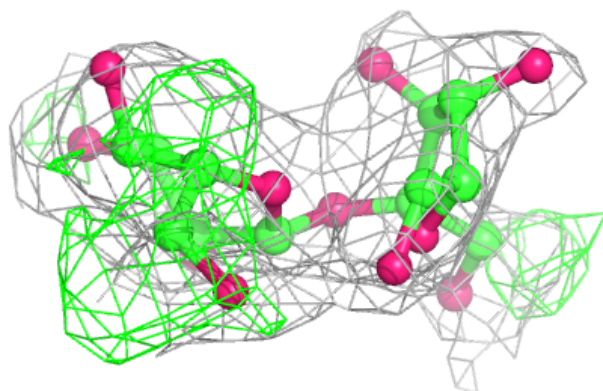
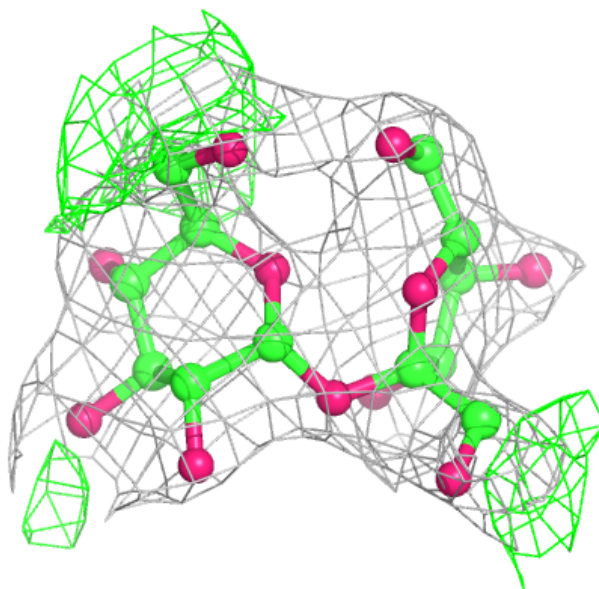
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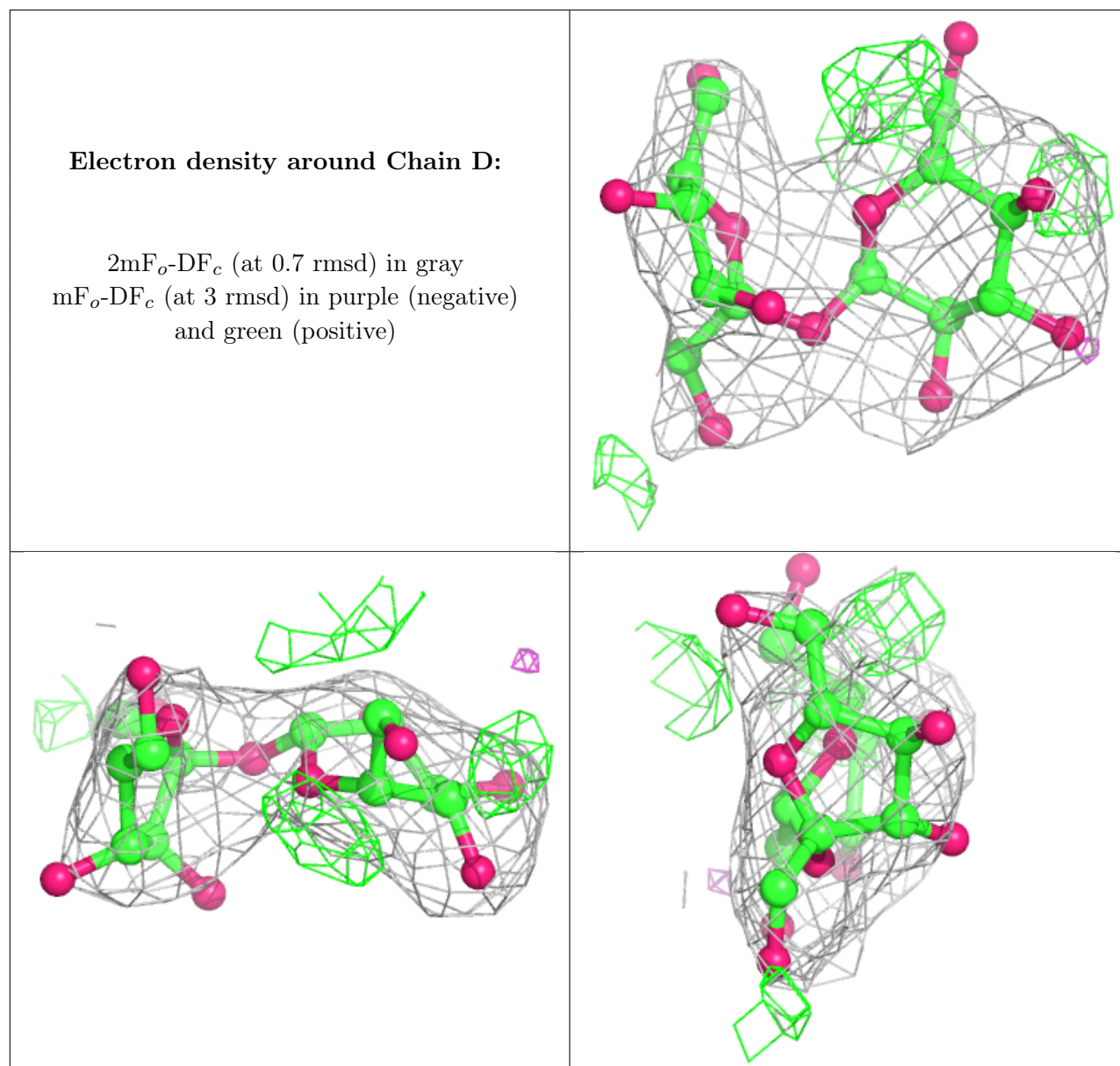
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	FRU	C	2	12/12	0.85	0.26	60,74,80,82	12
2	GLC	D	1	11/12	0.90	0.20	70,78,88,88	11
2	FRU	D	2	12/12	0.93	0.33	65,78,81,87	12

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)





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
6	CL	A	515	1/1	0.79	0.10	98,98,98,98	0
4	SO4	B	506	5/5	0.82	0.22	79,79,101,101	5
4	SO4	B	508	5/5	0.84	0.19	69,71,86,88	5
4	SO4	A	507	5/5	0.84	0.15	75,87,108,130	0
6	CL	B	512	1/1	0.85	0.59	90,90,90,90	0

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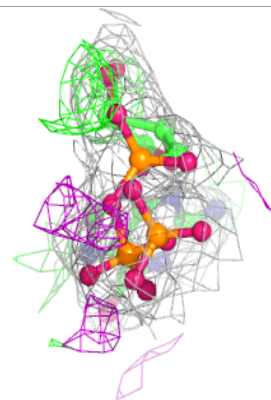
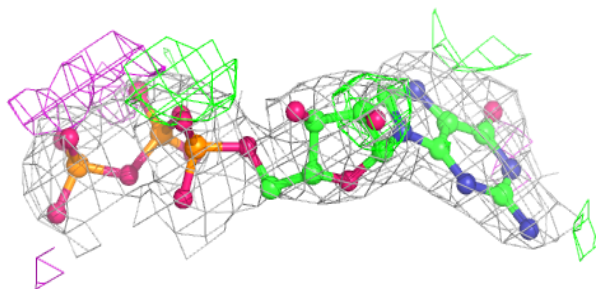
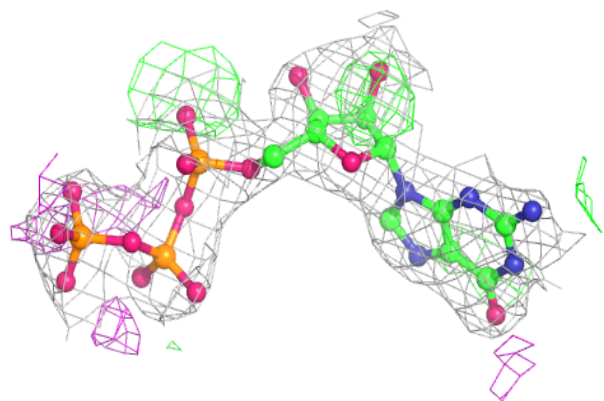
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
4	SO4	A	504	5/5	0.86	0.19	88,96,105,119	5
4	SO4	A	510	5/5	0.86	0.16	87,97,116,119	0
4	SO4	A	508	5/5	0.87	0.14	73,77,92,95	5
4	SO4	B	509	5/5	0.88	0.19	85,91,105,111	5
3	GTP	A	501	32/32	0.91	0.22	52,60,76,80	32
4	SO4	A	512	5/5	0.92	0.27	62,67,75,77	5
4	SO4	B	504	5/5	0.92	0.17	74,74,88,95	5
4	SO4	A	511	5/5	0.92	0.12	92,97,105,116	0
4	SO4	B	507	5/5	0.92	0.20	69,70,91,93	5
4	SO4	A	505	5/5	0.93	0.11	73,75,91,95	5
4	SO4	B	505	5/5	0.93	0.22	81,89,103,116	0
4	SO4	A	506	5/5	0.94	0.17	80,92,104,116	0
4	SO4	A	509	5/5	0.95	0.09	80,91,97,106	5
3	GTP	B	501	32/32	0.95	0.25	54,60,68,70	18
4	SO4	A	503	5/5	0.97	0.23	55,66,76,84	0
5	MN	B	510	1/1	0.98	0.14	61,61,61,61	0
5	MN	B	511	1/1	0.98	0.15	54,54,54,54	1
4	SO4	B	502	5/5	0.98	0.30	53,57,68,70	5
4	SO4	B	503	5/5	0.98	0.22	50,58,74,77	5
6	CL	B	513	1/1	0.98	0.20	78,78,78,78	0
5	MN	A	514	1/1	0.99	0.17	66,66,66,66	0
4	SO4	A	502	5/5	0.99	0.17	44,49,56,60	5
5	MN	A	513	1/1	1.00	0.13	63,63,63,63	0

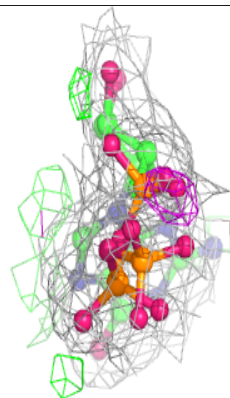
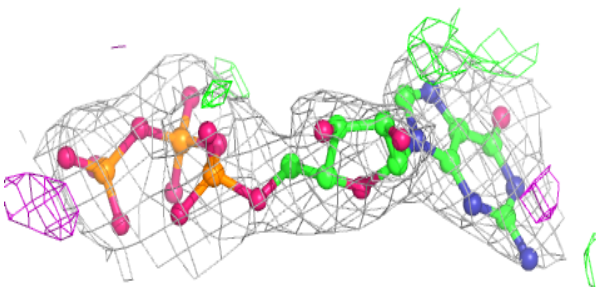
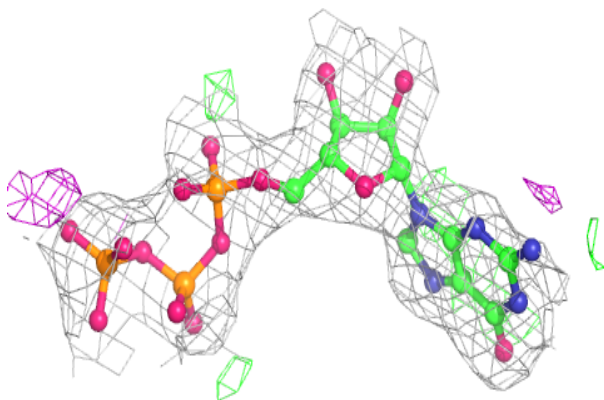
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

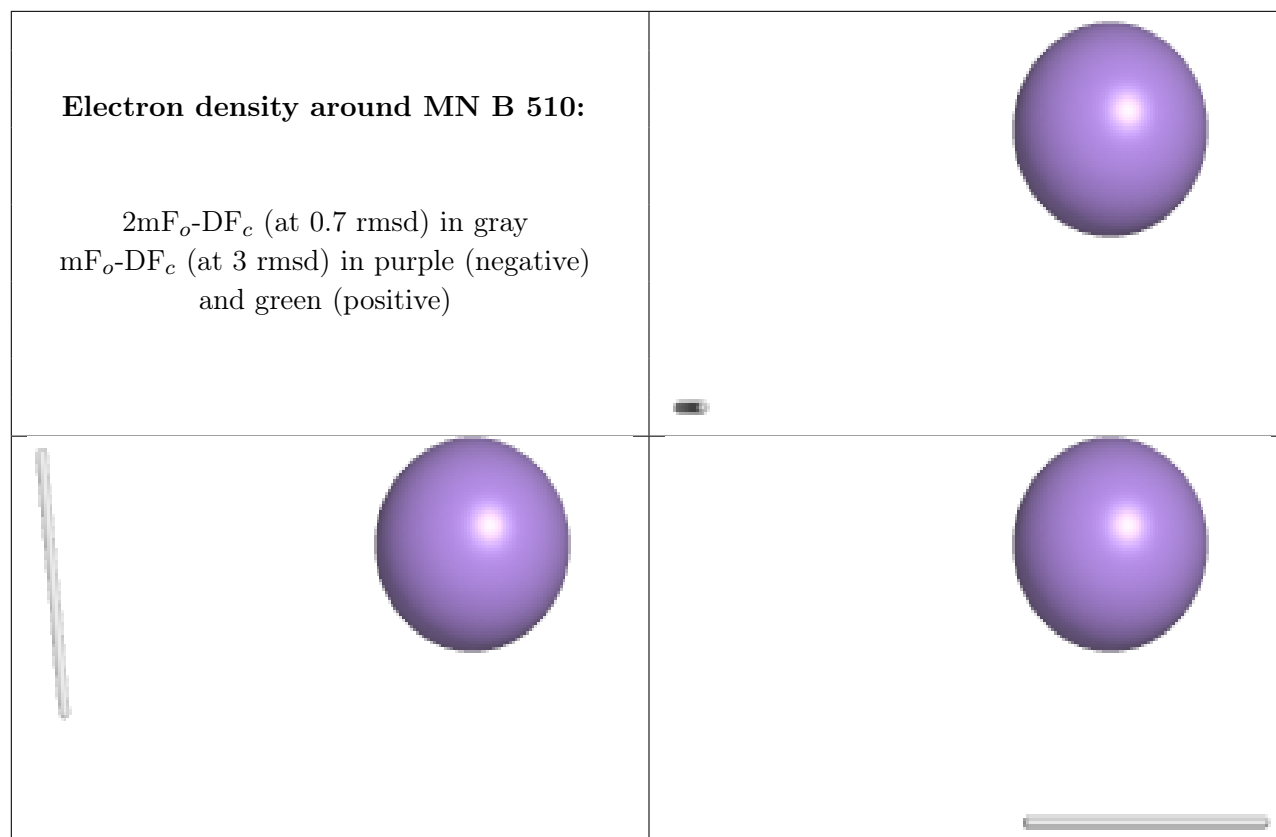
Electron density around GTP A 501:

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

**Electron density around GTP B 501:**

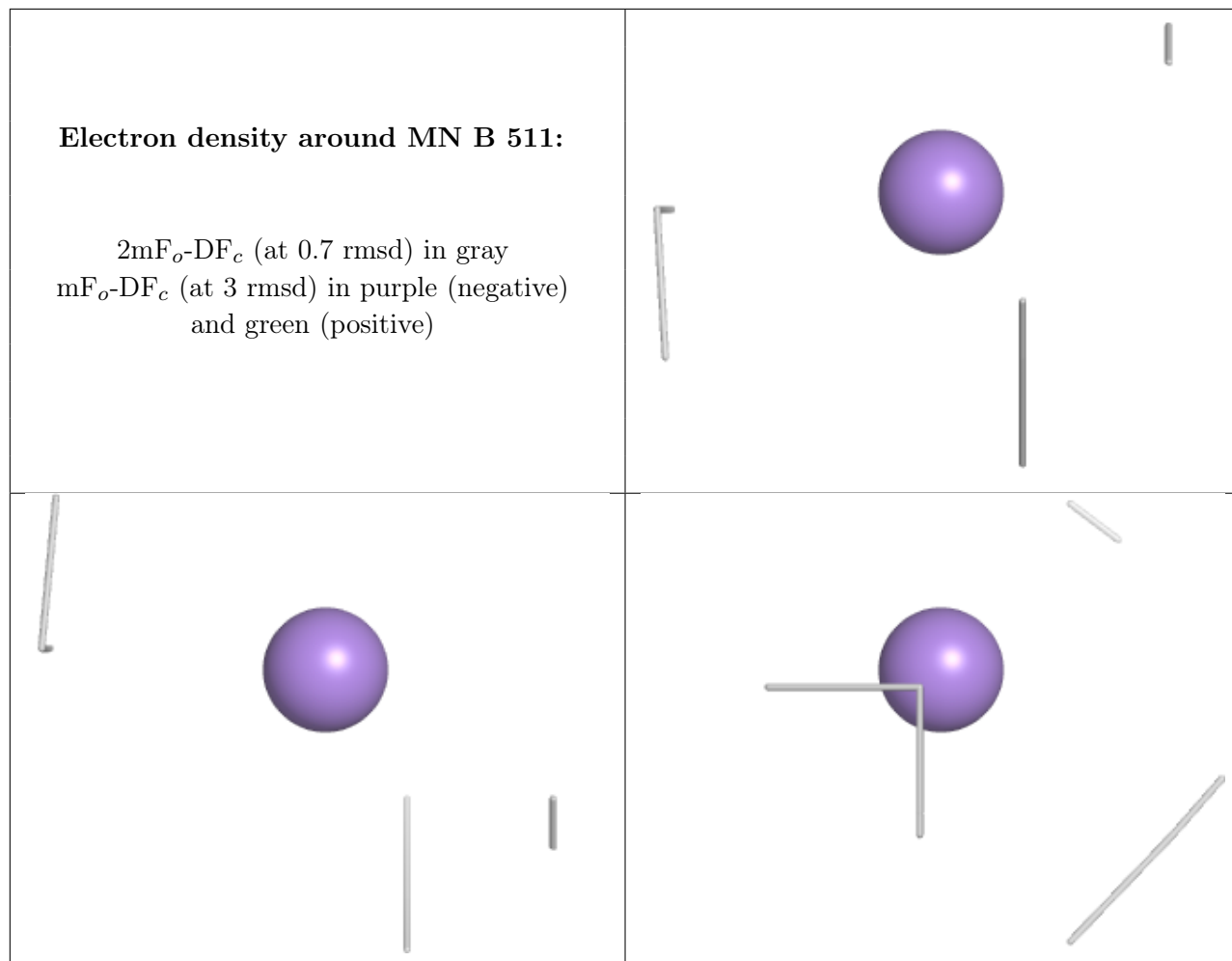
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)





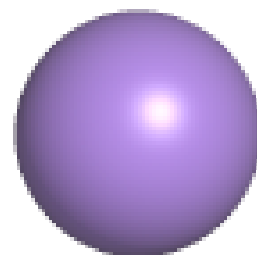
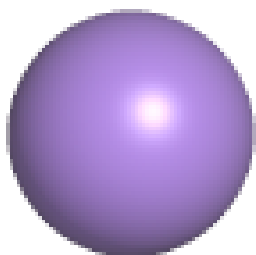
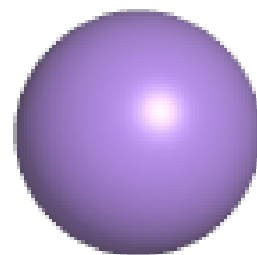
Electron density around MN B 511:

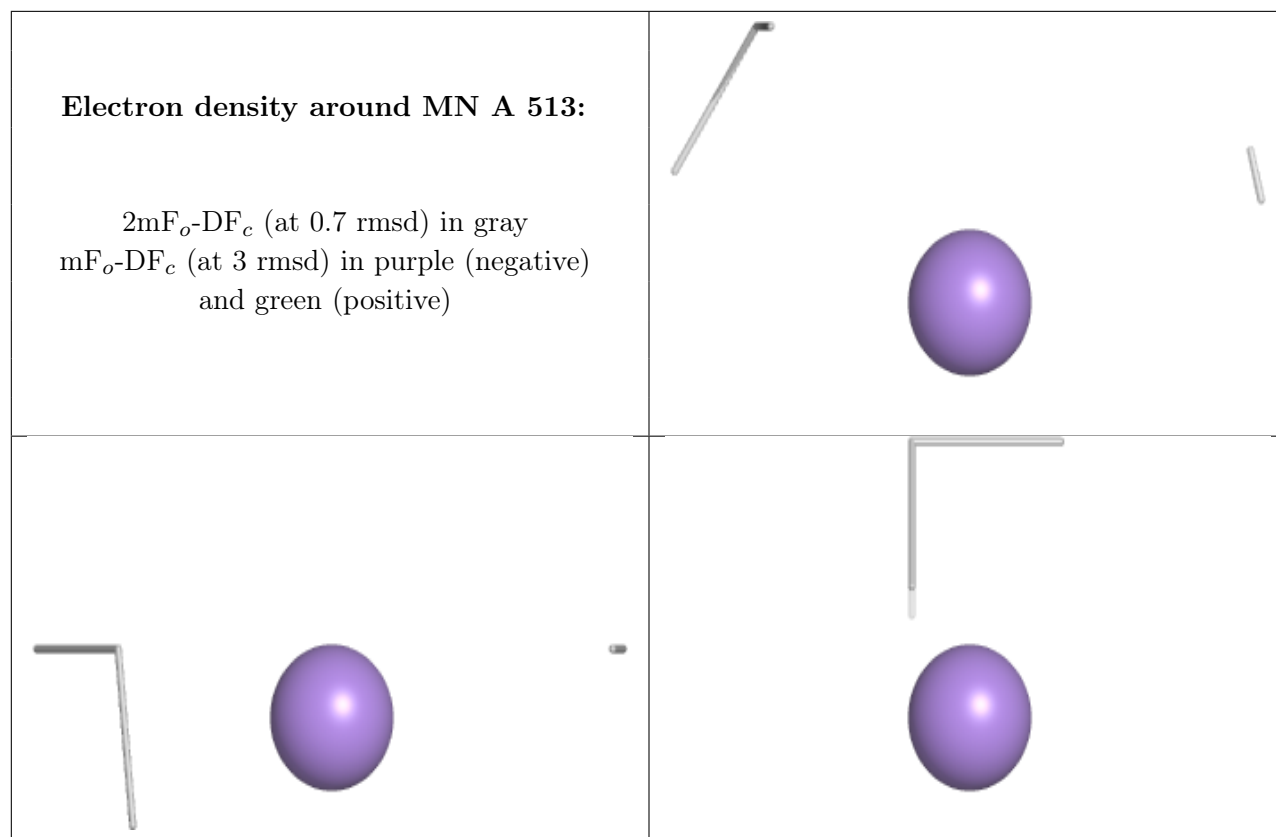
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around MN A 514:

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





6.5 Other polymers [i](#)

There are no such residues in this entry.