



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 09:46 AM UTC

PDB ID : 6WB6 / pdb_00006wb6
Title : 2.05 Å resolution structure of transferrin 1 from Manduca sexta
Authors : Lovell, S.; Kashipathy, M.M.; Battaile, K.P.; Weber, J.J.; Gorman, M.J.
Deposited on : 2020-03-26
Resolution : 2.05 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
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
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

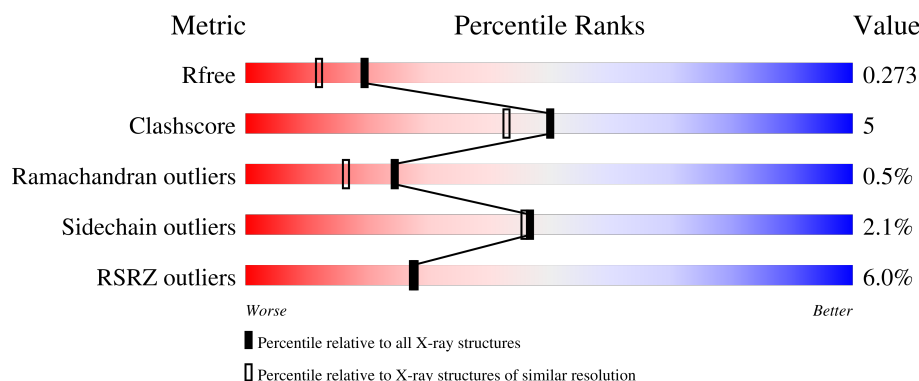
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.05 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	2260 (2.04-2.04)
Clashscore	190562	2333 (2.04-2.04)
Ramachandran outliers	187476	2318 (2.04-2.04)
Sidechain outliers	187428	2318 (2.04-2.04)
RSRZ outliers	180081	2260 (2.04-2.04)

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	663	3% 85% 13% ..
1	B	663	9% 85% 13% ..
2	C	2	100%
2	D	2	100%

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 10434 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 Transferrin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	653	Total	C	N	O	S	0	1	0
			5017	3176	865	941	35			
1	B	653	Total	C	N	O	S	0	0	0
			4908	3099	840	935	34			

- 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
2	C	2	Total	C	N	O	0	0	0
			28	16	2	10			
2	D	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 3 is FE (III) ION (CCD ID: FE) (formula: Fe) (labeled as "Ligand of Interest" by depositor).

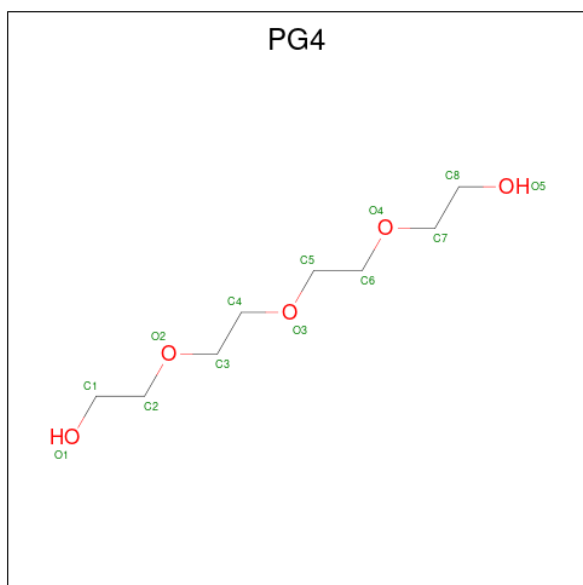
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Fe	0	0
			1	1		
3	B	1	Total	Fe	0	0
			1	1		

- Molecule 4 is CARBONATE ION (CCD ID: CO3) (formula: CO₃) (labeled as "Ligand of Interest" by depositor).



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

- Molecule 5 is TETRAETHYLENE GLYCOL (CCD ID: PG4) (formula: $C_8H_{18}O_5$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total	C	O	0	0
			13	8	5		

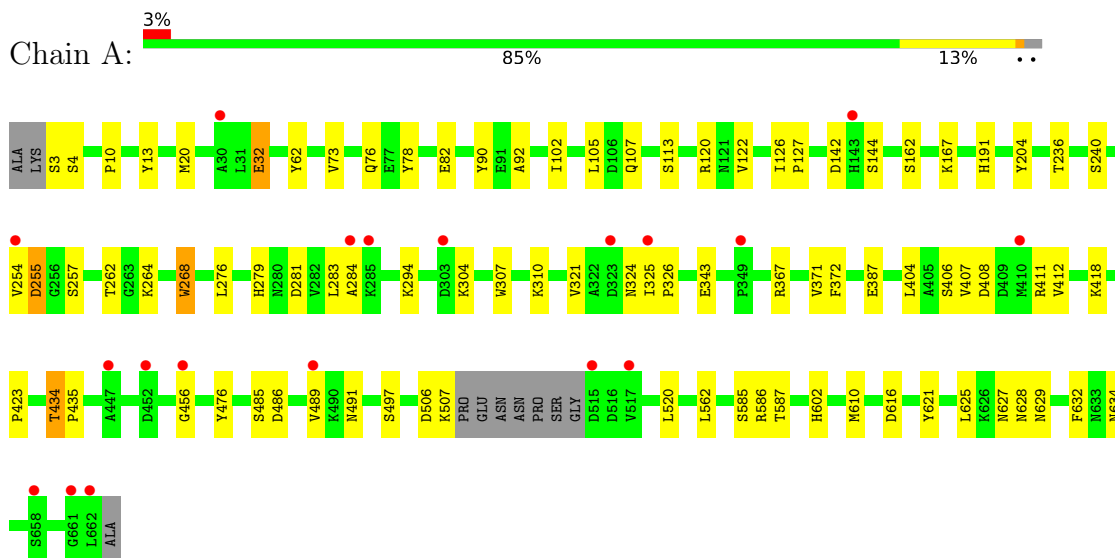
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	295	Total	O	0	0
			295	295		
6	B	127	Total	O	0	0
			127	127		

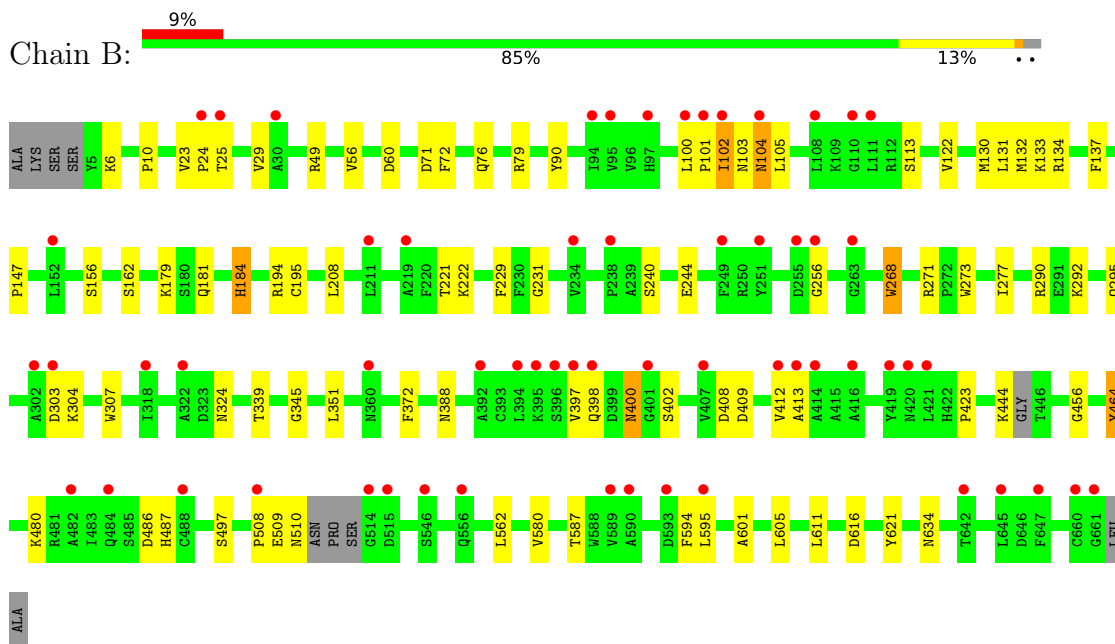
3 Residue-property plots

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: Transferrin



• Molecule 1: Transferrin



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

Chain C:  100%

MAG1
MAG2

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

Chain D:  100%

MAG1
MAG2

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	64.49Å 139.06Å 146.70Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.00 – 2.05 40.00 – 2.05	Depositor EDS
% Data completeness (in resolution range)	99.9 (40.00-2.05) 99.9 (40.00-2.05)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.10 (at 2.05Å)	Xtriage
Refinement program	PHENIX dev_3755	Depositor
R, R_{free}	0.209 , 0.272 0.217 , 0.273	Depositor DCC
R_{free} test set	4065 reflections (4.86%)	wwPDB-VP
Wilson B-factor (Å ²)	26.9	Xtriage
Anisotropy	0.388	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 42.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.42$, $\langle L^2 \rangle = 0.25$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	10434	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 3.90% 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: FE, PG4, NAG, CO3

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	0.46	2/5136 (0.0%)	0.60	0/6972
1	B	0.33	0/5023	0.51	0/6844
All	All	0.40	2/10159 (0.0%)	0.56	0/13816

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	204	TYR	C-O	-5.94	1.17	1.24
1	A	90	TYR	C-O	-5.11	1.16	1.23

There are no bond angle outliers.

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	5017	0	4845	54	0
1	B	4908	0	4575	52	0
2	C	28	0	25	0	0
2	D	28	0	25	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	8	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	8	0	0	1	0
5	B	13	0	18	0	0
6	A	295	0	0	6	0
6	B	127	0	0	4	0
All	All	10434	0	9488	104	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 (104) 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:90:TYR:CG	1:B:222:LYS:HD2	2.25	0.72
1:A:434:THR:HG22	1:A:435:PRO:O	1.97	0.65
1:B:132:MET:HE1	1:B:345:GLY:O	2.03	0.59
1:B:290:ARG:NH2	1:B:324:ASN:O	2.36	0.58
1:A:372:PHE:HE2	1:A:610:MET:HE3	1.71	0.56
1:A:78:TYR:OH	1:A:276:LEU:HD21	2.07	0.54
1:A:602:HIS:HE1	6:A:1011:HOH:O	1.90	0.53
1:B:90:TYR:CD2	1:B:222:LYS:HD2	2.43	0.53
1:A:489:VAL:O	1:A:520:LEU:HD11	2.08	0.52
1:B:303:ASP:CB	6:B:911:HOH:O	2.58	0.52
1:B:76:GLN:NE2	1:B:290:ARG:HG2	2.24	0.51
1:B:221:THR:OG1	1:B:222:LYS:N	2.40	0.51
1:B:29:VAL:O	1:B:292:LYS:NZ	2.39	0.51
1:A:407:VAL:HG13	1:A:412:VAL:HG22	1.91	0.51
1:B:100:LEU:O	1:B:102:ILE:N	2.44	0.50
1:B:486:ASP:O	1:B:487:HIS:C	2.55	0.50
1:A:283:LEU:O	1:A:284:ALA:HB3	2.11	0.50
1:A:294:LYS:HG2	1:A:321:VAL:HB	1.93	0.50
1:A:628:ASN:OD1	1:A:634:ASN:ND2	2.45	0.50
1:B:621:TYR:O	1:B:634:ASN:HA	2.12	0.50
1:B:268:TRP:CD1	1:B:268:TRP:H	2.29	0.50
1:B:130:MET:O	1:B:133:LYS:HG2	2.13	0.49
1:B:76:GLN:HE22	1:B:290:ARG:HG2	1.77	0.49
1:A:255:ASP:HB3	1:A:257:SER:H	1.77	0.49
1:A:408:ASP:HB2	1:A:411:ARG:NH2	2.27	0.49
1:A:423:PRO:HB3	1:A:587:THR:HG21	1.93	0.49
1:B:104:ASN:OD1	1:B:105:LEU:N	2.46	0.49
1:B:601:ALA:O	1:B:605:LEU:HG	2.12	0.48
1:A:585:SER:O	1:A:586:ARG:C	2.55	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:90:TYR:CZ	1:B:271:ARG:HD3	2.48	0.48
1:B:194:ARG:NH1	6:B:811:HOH:O	2.40	0.48
1:B:412:VAL:HG23	1:B:413:ALA:H	1.78	0.48
1:A:10:PRO:HG2	1:A:616:ASP:HB2	1.95	0.48
1:A:3:SER:O	1:A:3:SER:OG	2.26	0.48
1:A:62:TYR:OH	1:A:343:GLU:OE2	2.26	0.47
1:A:372:PHE:CE2	1:A:610:MET:HE3	2.48	0.47
1:B:79:ARG:HA	1:B:273:TRP:CE3	2.49	0.47
1:A:113:SER:O	1:A:162:SER:HA	2.15	0.47
1:A:268:TRP:H	1:A:268:TRP:CD1	2.32	0.47
1:B:113:SER:O	1:B:162:SER:HA	2.15	0.47
1:A:387:GLU:OE2	6:A:801:HOH:O	2.21	0.46
1:A:585:SER:HB3	6:A:925:HOH:O	2.15	0.46
1:A:191:HIS:CE1	1:B:24:PRO:HB2	2.50	0.46
1:A:126:ILE:HB	1:A:127:PRO:HD3	1.98	0.46
1:B:398:GLN:C	1:B:400:ASN:H	2.24	0.46
1:A:456:GLY:HA2	1:A:497:SER:OG	2.16	0.45
1:A:279:HIS:CE1	1:A:281:ASP:HB2	2.50	0.45
1:A:13:TYR:HE2	1:A:616:ASP:HB3	1.81	0.45
1:A:325:ILE:O	1:A:326:PRO:C	2.57	0.45
1:B:464:TYR:C	1:B:464:TYR:CD1	2.94	0.45
1:A:324:ASN:OD1	1:A:324:ASN:N	2.50	0.45
1:B:412:VAL:HG23	1:B:413:ALA:N	2.32	0.45
1:A:629:ASN:HA	1:A:632:PHE:O	2.17	0.44
1:B:208:LEU:HD13	1:B:229:PHE:HB2	2.00	0.44
1:B:90:TYR:CB	1:B:222:LYS:HD2	2.47	0.44
1:B:339:THR:HG21	6:B:863:HOH:O	2.17	0.44
1:A:73:VAL:HG11	1:A:283:LEU:CD1	2.48	0.44
1:A:367:ARG:O	1:A:371:VAL:HG23	2.18	0.44
1:A:625:LEU:O	1:A:627:ASN:ND2	2.48	0.44
1:B:304:LYS:CB	6:B:925:HOH:O	2.66	0.44
1:A:142:ASP:OD1	1:A:144:SER:OG	2.34	0.44
1:A:13:TYR:CE1	1:A:310:LYS:HG2	2.53	0.43
1:B:104:ASN:OD1	1:B:104:ASN:C	2.61	0.43
1:A:120:ARG:HB3	4:A:703:CO3:C	2.48	0.43
1:B:456:GLY:HA2	1:B:497:SER:OG	2.19	0.43
1:A:486:ASP:O	1:A:491:ASN:ND2	2.51	0.43
1:A:489:VAL:HG22	1:A:520:LEU:CD1	2.49	0.43
1:B:147:PRO:HD3	1:B:351:LEU:HD23	2.00	0.43
1:B:10:PRO:HG2	1:B:616:ASP:HB2	2.01	0.43
1:A:191:HIS:CD2	1:B:24:PRO:CB	3.02	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:404:LEU:N	1:A:404:LEU:HD23	2.34	0.43
1:B:56:VAL:HB	1:B:60:ASP:HB2	2.01	0.43
1:B:131:LEU:HB3	1:B:137:PHE:CD1	2.53	0.43
1:B:562:LEU:HD12	1:B:580:VAL:HG22	2.01	0.43
1:B:122:VAL:HG22	4:B:702:CO3:C	2.48	0.42
1:A:4:SER:HB2	1:A:32:GLU:CG	2.48	0.42
1:B:231:GLY:HA3	1:B:240:SER:O	2.19	0.42
1:A:406:SER:HA	1:A:587:THR:O	2.18	0.42
1:B:90:TYR:CE1	1:B:122:VAL:HG13	2.54	0.42
1:B:25:THR:HB	1:B:295:GLN:OE1	2.20	0.42
1:B:72:PHE:HB3	1:B:277:ILE:HD11	2.01	0.42
1:A:102:ILE:HG13	1:A:107:GLN:HB2	2.02	0.41
1:A:20:MET:HB2	1:A:307:TRP:CG	2.55	0.41
1:A:616:ASP:OD1	1:A:616:ASP:N	2.53	0.41
1:A:92:ALA:HB2	1:A:122:VAL:HG12	2.02	0.41
1:A:621:TYR:O	1:A:634:ASN:HA	2.20	0.41
1:B:156:SER:HB2	1:B:184:HIS:NE2	2.36	0.41
1:B:103:ASN:O	1:B:256:GLY:HA3	2.21	0.41
1:A:82:GLU:HA	1:A:236:THR:OG1	2.20	0.41
1:B:23:VAL:HG21	1:B:307:TRP:CZ2	2.56	0.41
1:B:105:LEU:HB2	1:B:134:ARG:CZ	2.51	0.41
1:B:372:PHE:CE2	1:B:611:LEU:HD21	2.56	0.41
1:B:49:ARG:NH2	1:B:71:ASP:OD2	2.47	0.41
1:A:610:MET:HE2	1:A:610:MET:HB3	1.95	0.40
1:B:179:LYS:NZ	1:B:195:CYS:O	2.48	0.40
1:B:409:ASP:O	1:B:412:VAL:HG22	2.21	0.40
1:A:476:TYR:CZ	1:A:562:LEU:HB3	2.56	0.40
1:B:423:PRO:HB3	1:B:587:THR:HG21	2.04	0.40
1:A:262:THR:OG1	6:A:802:HOH:O	2.22	0.40
1:A:105:LEU:HD11	1:A:254:VAL:HA	2.04	0.40
1:A:167:LYS:NZ	6:A:850:HOH:O	2.54	0.40
1:A:506:ASP:O	1:A:507:LYS:C	2.64	0.40
1:A:586:ARG:N	6:A:846:HOH:O	2.54	0.40
1:B:397:VAL:HA	1:B:402:SER:O	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	650/663 (98%)	621 (96%)	27 (4%)	2 (0%)	36	30
1	B	647/663 (98%)	609 (94%)	33 (5%)	5 (1%)	16	9
All	All	1297/1326 (98%)	1230 (95%)	60 (5%)	7 (0%)	24	16

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	400	ASN
1	B	509	GLU
1	A	255	ASP
1	B	594	PHE
1	B	508	PRO
1	B	101	PRO
1	A	304	LYS

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	539/569 (95%)	531 (98%)	8 (2%)	57	58
1	B	509/569 (90%)	495 (97%)	14 (3%)	38	34
All	All	1048/1138 (92%)	1026 (98%)	22 (2%)	47	46

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	32	GLU
1	A	76	GLN
1	A	240	SER
1	A	264	LYS
1	A	268	TRP
1	A	418	LYS
1	A	434	THR
1	A	485	SER
1	B	6	LYS
1	B	102	ILE
1	B	104	ASN
1	B	181	GLN
1	B	184	HIS
1	B	244	GLU
1	B	268	TRP
1	B	388	ASN
1	B	408	ASP
1	B	444	LYS
1	B	464	TYR
1	B	480	LYS
1	B	510	ASN
1	B	595	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	76	GLN
1	A	103	ASN
1	A	191	HIS
1	A	280	ASN
1	A	602	HIS
1	A	609	GLN
1	B	97	HIS
1	B	471	HIS
1	B	609	GLN
1	B	633	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$
2	NAG	C	1	2,1	14,14,15	0.42	0	17,19,21	1.30	1 (5%)
2	NAG	C	2	2	14,14,15	0.67	1 (7%)	17,19,21	1.26	1 (5%)
2	NAG	D	1	2,1	14,14,15	0.30	0	17,19,21	1.09	2 (11%)
2	NAG	D	2	2	14,14,15	0.44	0	17,19,21	1.09	1 (5%)

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	C	1	2,1	-	0/6/23/26	0/1/1/1
2	NAG	C	2	2	-	2/6/23/26	0/1/1/1
2	NAG	D	1	2,1	-	1/6/23/26	0/1/1/1
2	NAG	D	2	2	-	1/6/23/26	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	2	NAG	C1-C2	2.11	1.55	1.52

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	1	NAG	C1-O5-C5	2.95	116.14	112.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	2	NAG	C4-C3-C2	2.59	114.82	111.02
2	C	1	NAG	O4-C4-C5	-2.49	103.19	109.32
2	D	1	NAG	O4-C4-C3	-2.44	104.62	110.38
2	D	2	NAG	O4-C4-C3	-2.14	105.34	110.38

There are no chirality outliers.

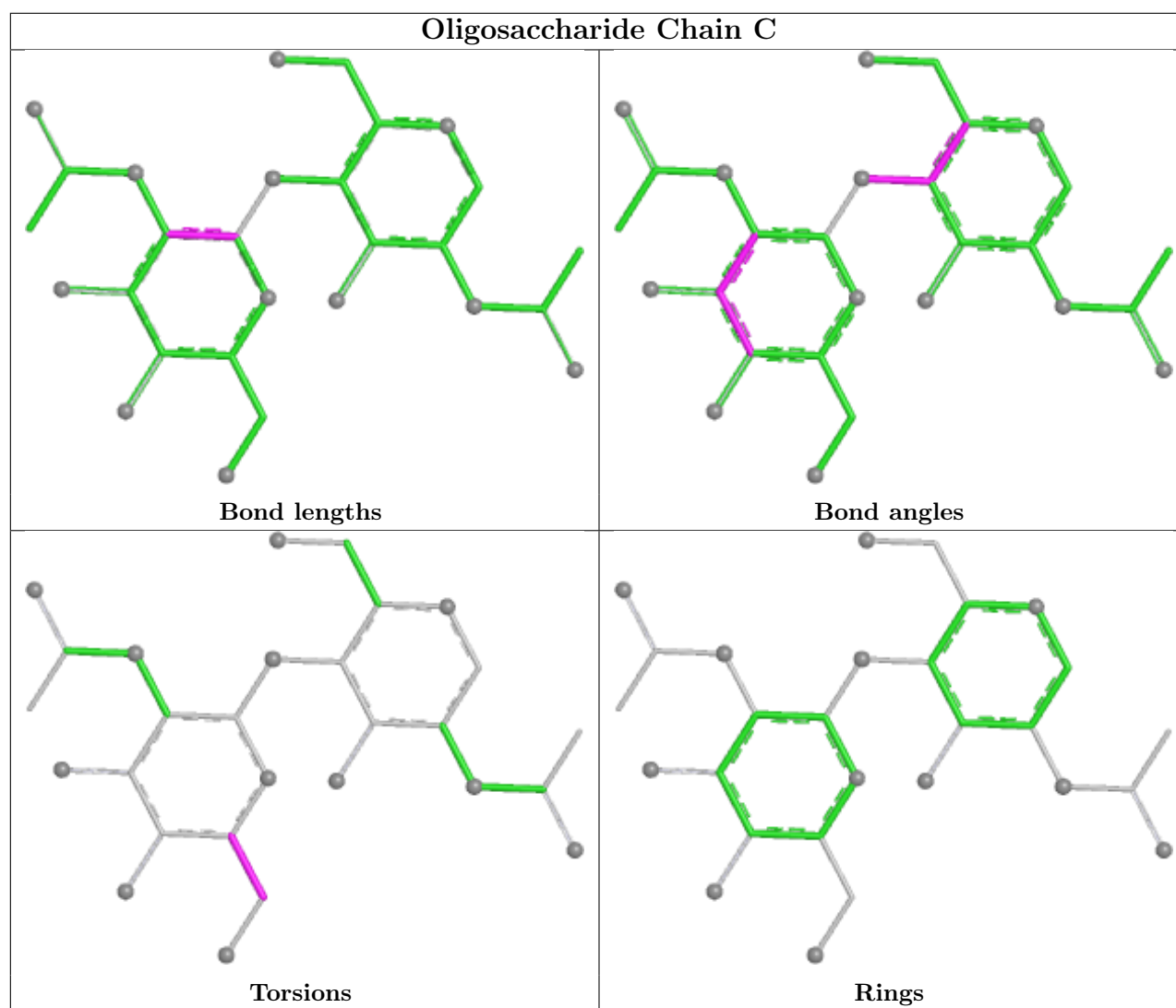
All (4) torsion outliers are listed below:

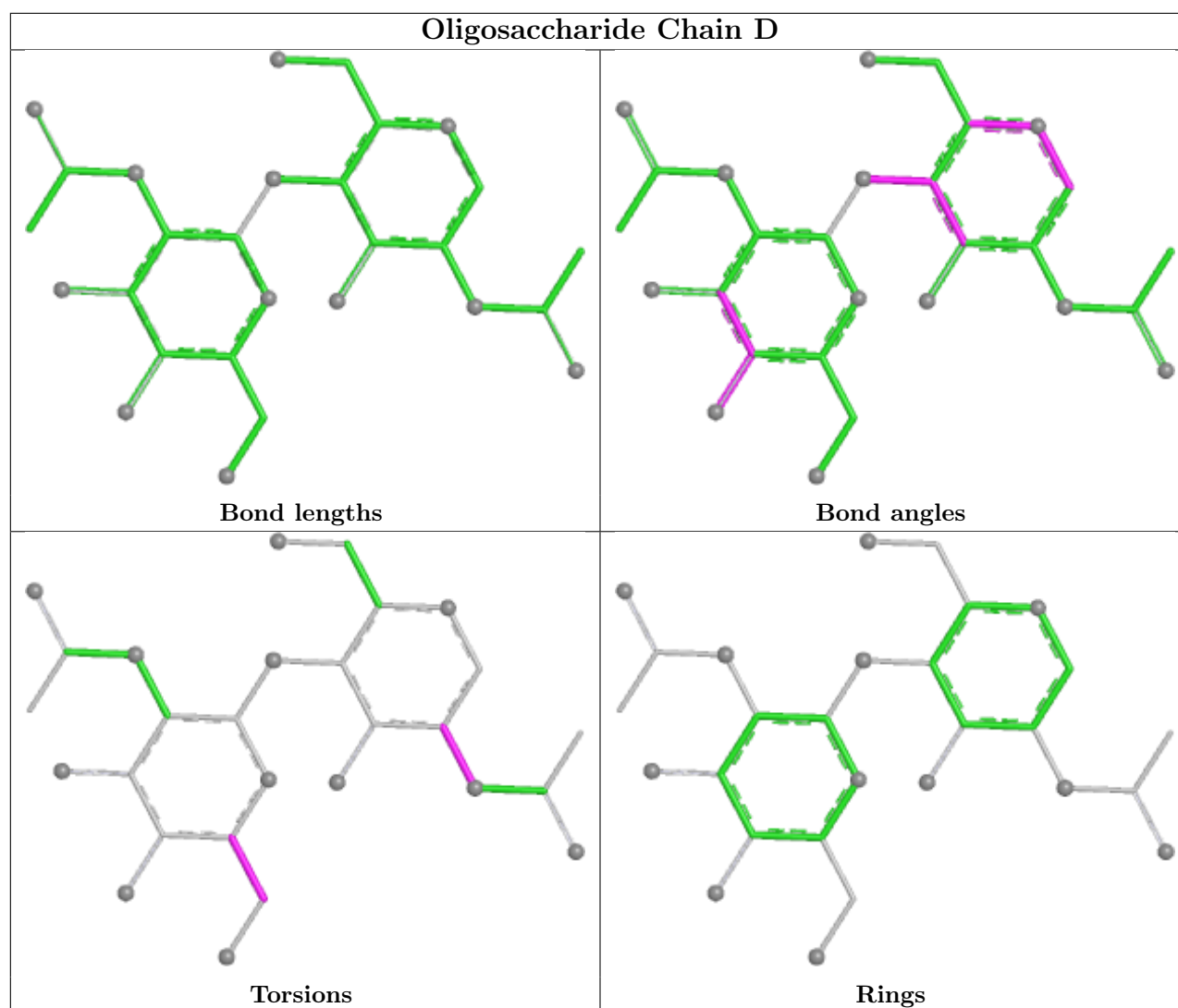
Mol	Chain	Res	Type	Atoms
2	C	2	NAG	C4-C5-C6-O6
2	C	2	NAG	O5-C5-C6-O6
2	D	1	NAG	C1-C2-N2-C7
2	D	2	NAG	C4-C5-C6-O6

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](#)

Of 7 ligands modelled in this entry, 2 are monoatomic - leaving 5 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	CO3	B	703	3	3,3,3	1.33	0	2,3,3	0.63	0
4	CO3	A	703	3	3,3,3	0.85	0	2,3,3	1.50	0
5	PG4	B	706	-	12,12,12	0.15	0	11,11,11	0.14	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	CO3	A	702	3	3,3,3	0.76	0	2,3,3	1.34	0
4	CO3	B	702	3	3,3,3	1.36	0	2,3,3	0.17	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
5	PG4	B	706	-	-	2/10/10/10	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) torsion outliers are listed below:

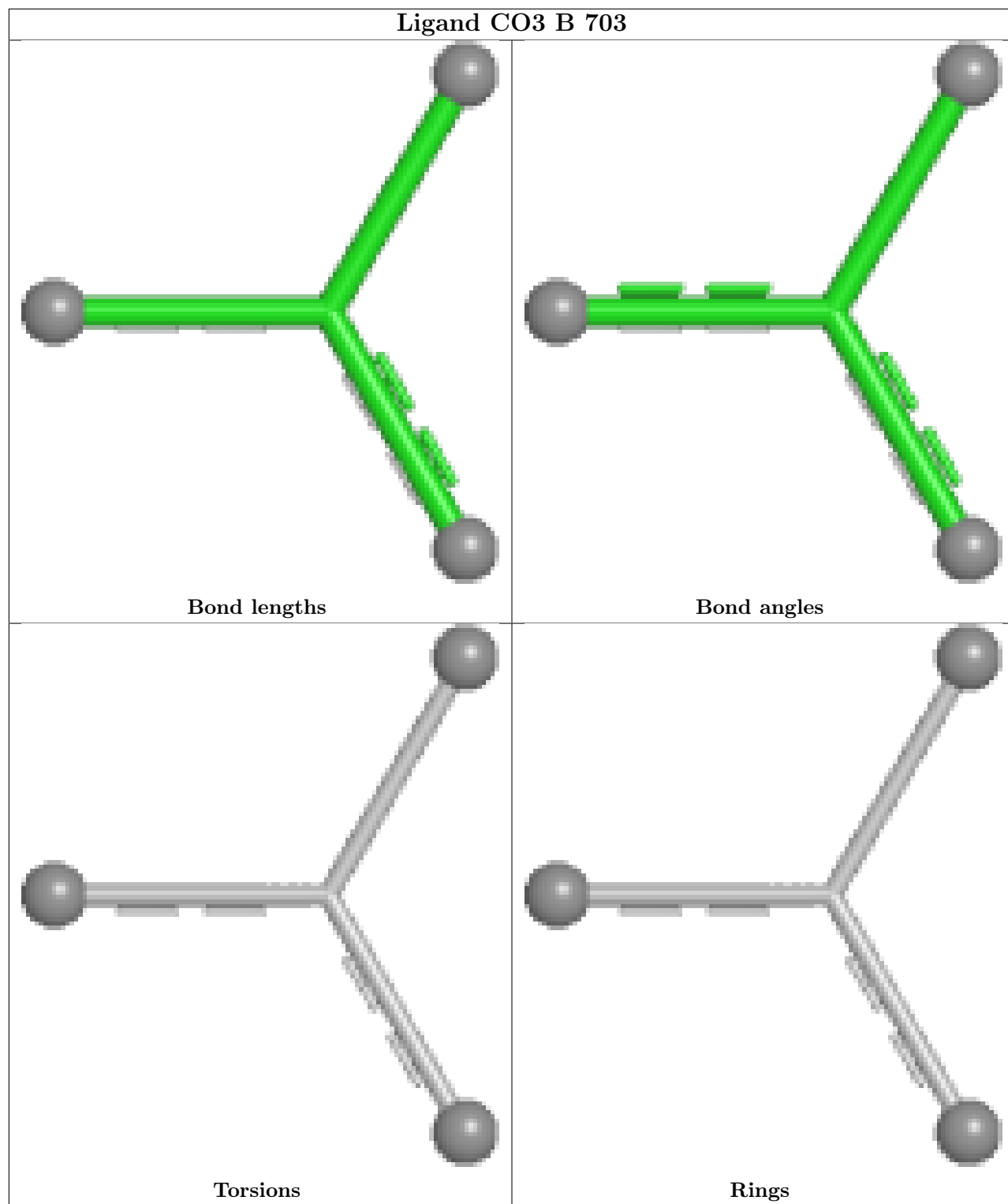
Mol	Chain	Res	Type	Atoms
5	B	706	PG4	C5-C6-O4-C7
5	B	706	PG4	C4-C3-O2-C2

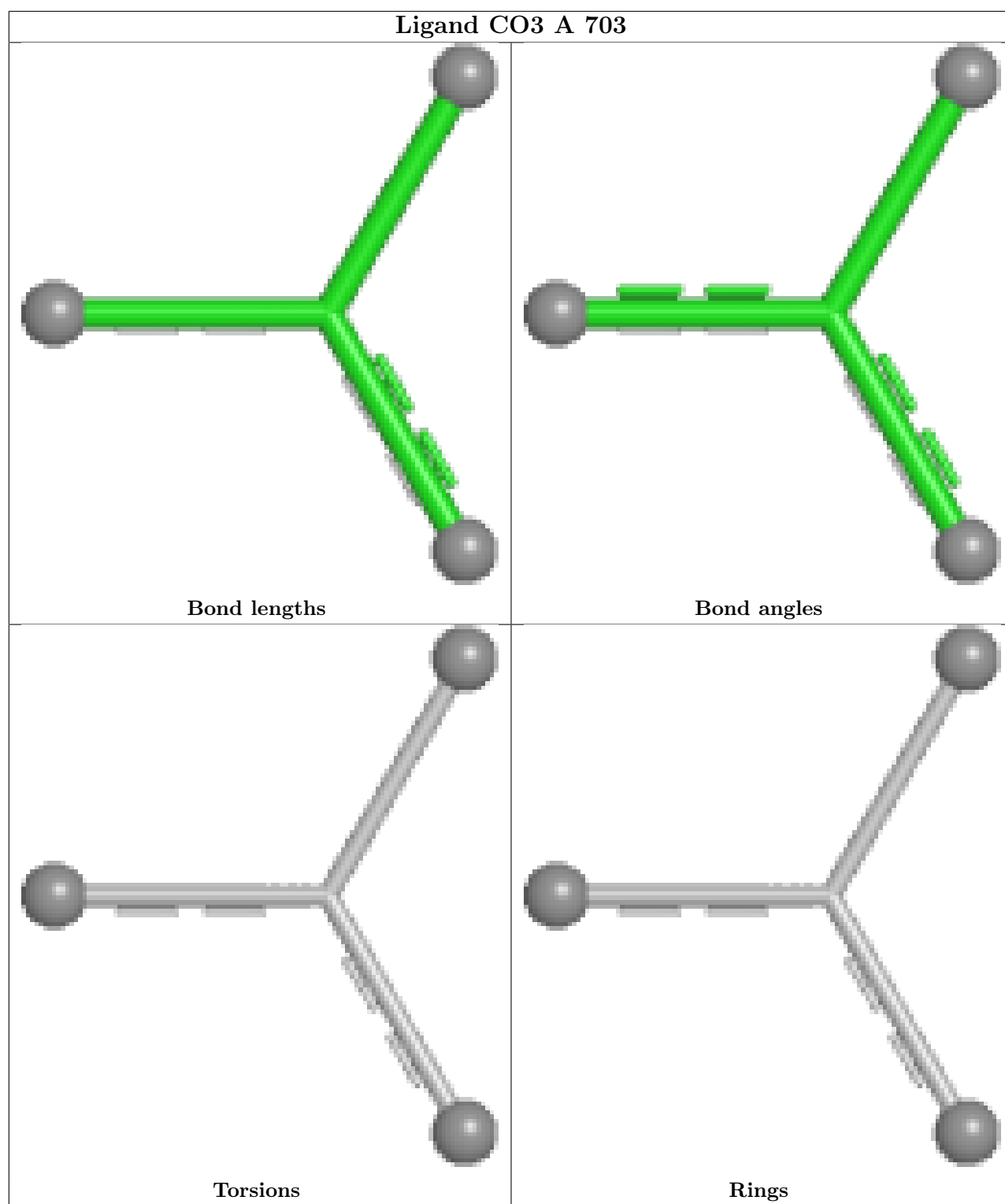
There are no ring outliers.

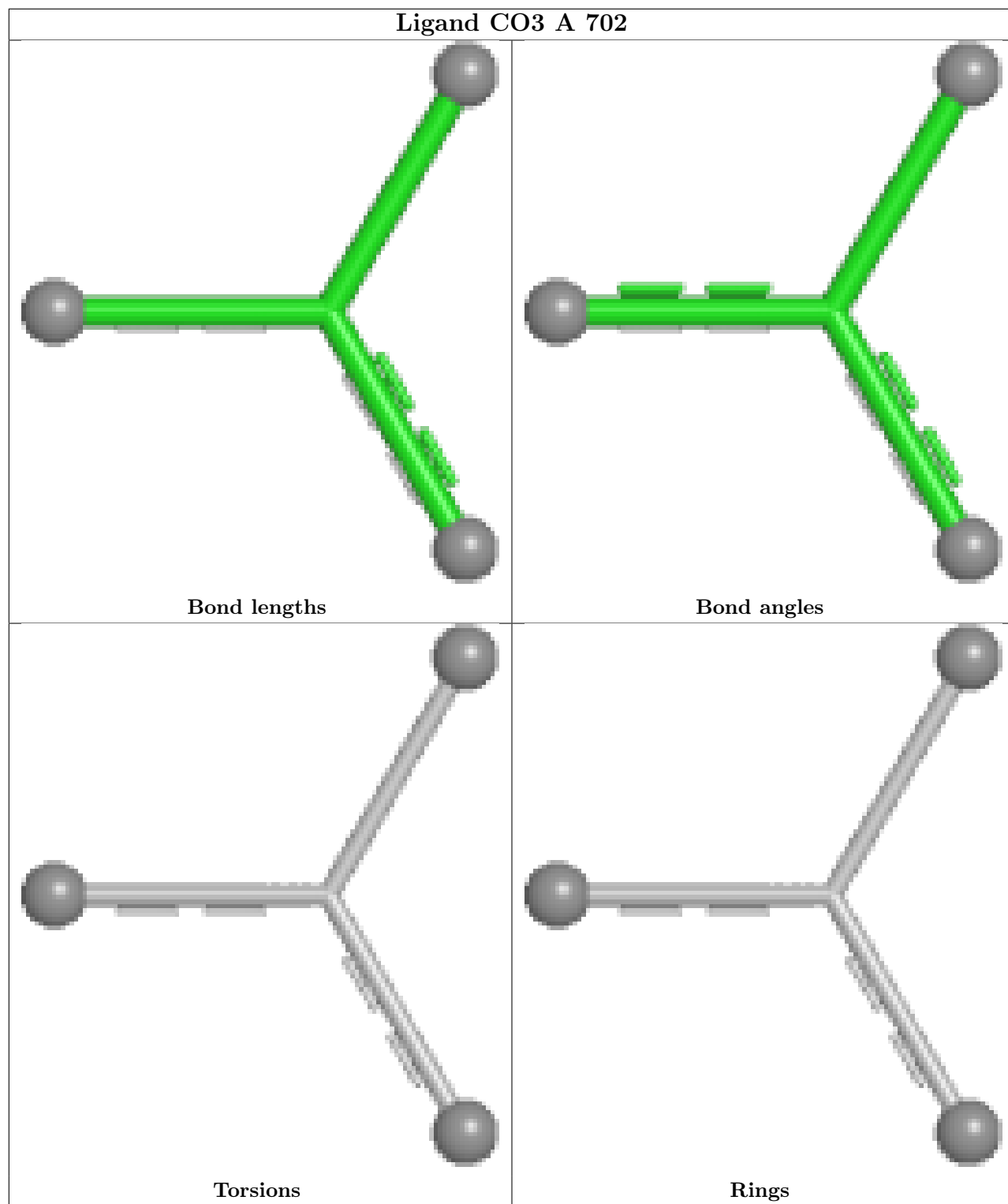
2 monomers are involved in 2 short contacts:

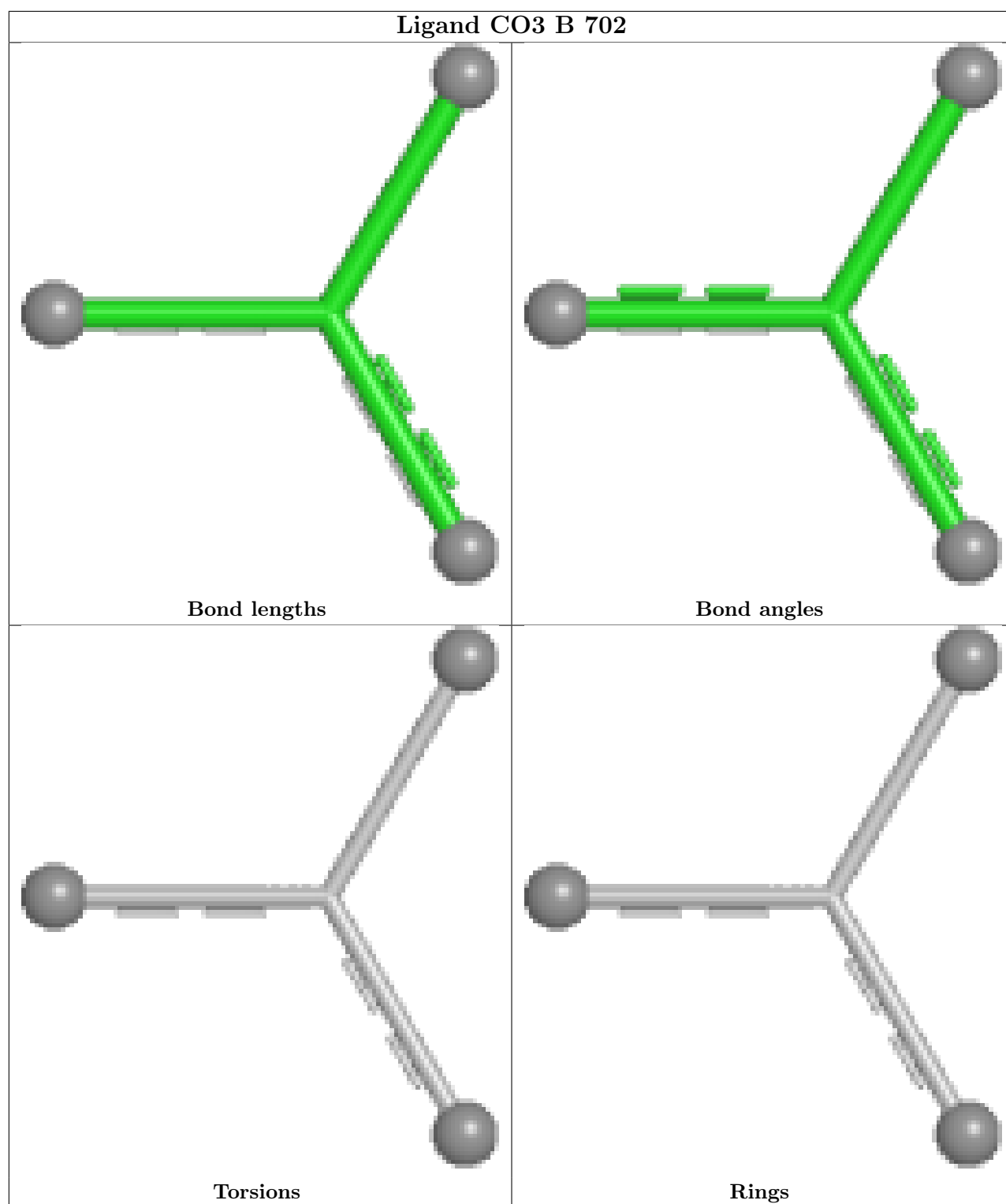
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	703	CO3	1	0
4	B	702	CO3	1	0

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 ⓘ

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	653/663 (98%)	0.21	19 (2%) 53 55	14, 27, 43, 59	1 (0%)
1	B	653/663 (98%)	0.82	60 (9%) 14 14	23, 38, 56, 70	0
All	All	1306/1326 (98%)	0.51	79 (6%) 27 28	14, 33, 53, 70	1 (0%)

All (79) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	303	ASP	4.7
1	B	484	GLN	3.5
1	A	285	LYS	3.5
1	A	489	VAL	3.3
1	B	416	ALA	3.2
1	B	421	LEU	3.2
1	B	414	ALA	3.2
1	B	30	ALA	3.1
1	B	515	ASP	3.0
1	B	590	ALA	2.9
1	B	645	LEU	2.9
1	B	647	PHE	2.9
1	B	407	VAL	2.9
1	B	211	LEU	2.9
1	B	482	ALA	2.8
1	B	397	VAL	2.8
1	B	392	ALA	2.8
1	A	662	LEU	2.7
1	A	447	ALA	2.7
1	B	396	SER	2.7
1	B	303	ASP	2.7
1	B	514	GLY	2.6
1	A	349	PRO	2.6
1	B	589	VAL	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	256	GLY	2.6
1	B	401	GLY	2.6
1	A	410	MET	2.6
1	B	24	PRO	2.6
1	A	456	GLY	2.5
1	B	488	CYS	2.5
1	A	30	ALA	2.5
1	B	263	GLY	2.5
1	B	94	ILE	2.5
1	B	593	ASP	2.4
1	B	25	THR	2.4
1	B	413	ALA	2.4
1	B	420	ASN	2.4
1	A	517	VAL	2.4
1	B	660	CYS	2.4
1	B	302	ALA	2.4
1	B	100	LEU	2.4
1	B	394	LEU	2.4
1	B	152	LEU	2.3
1	B	95	VAL	2.3
1	A	661	GLY	2.3
1	B	255	ASP	2.3
1	B	419	TYR	2.3
1	B	318	ILE	2.3
1	B	661	GLY	2.3
1	B	101	PRO	2.2
1	B	642	THR	2.2
1	B	219	ALA	2.2
1	B	111	LEU	2.2
1	B	595	LEU	2.2
1	B	251	TYR	2.2
1	B	104	ASN	2.2
1	B	234	VAL	2.2
1	B	102	ILE	2.2
1	B	238	PRO	2.2
1	B	412	VAL	2.2
1	A	284	ALA	2.2
1	B	546	SER	2.2
1	B	322	ALA	2.1
1	B	395	LYS	2.1
1	A	323	ASP	2.1
1	A	452	ASP	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	398	GLN	2.1
1	B	110	GLY	2.1
1	A	658	SER	2.1
1	B	508	PRO	2.1
1	A	325	ILE	2.1
1	B	249	PHE	2.1
1	B	556	GLN	2.0
1	B	97	HIS	2.0
1	A	254	VAL	2.0
1	B	360	ASN	2.0
1	B	108	LEU	2.0
1	A	515	ASP	2.0
1	A	143	HIS	2.0

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 ⓘ

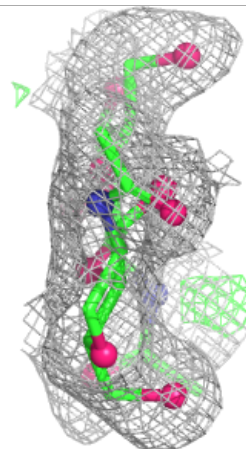
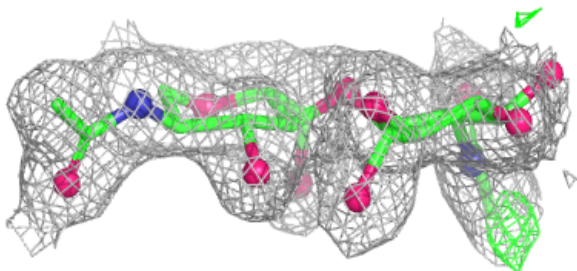
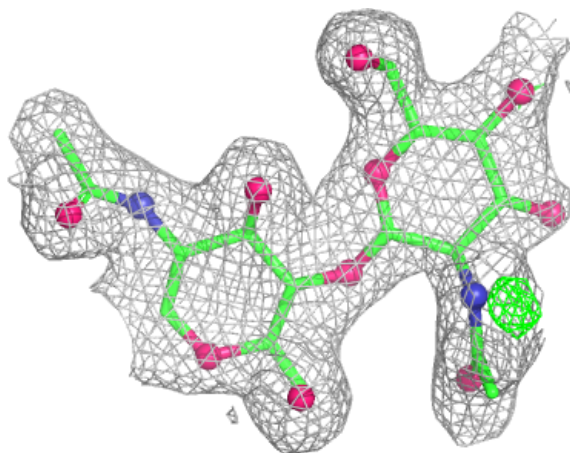
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

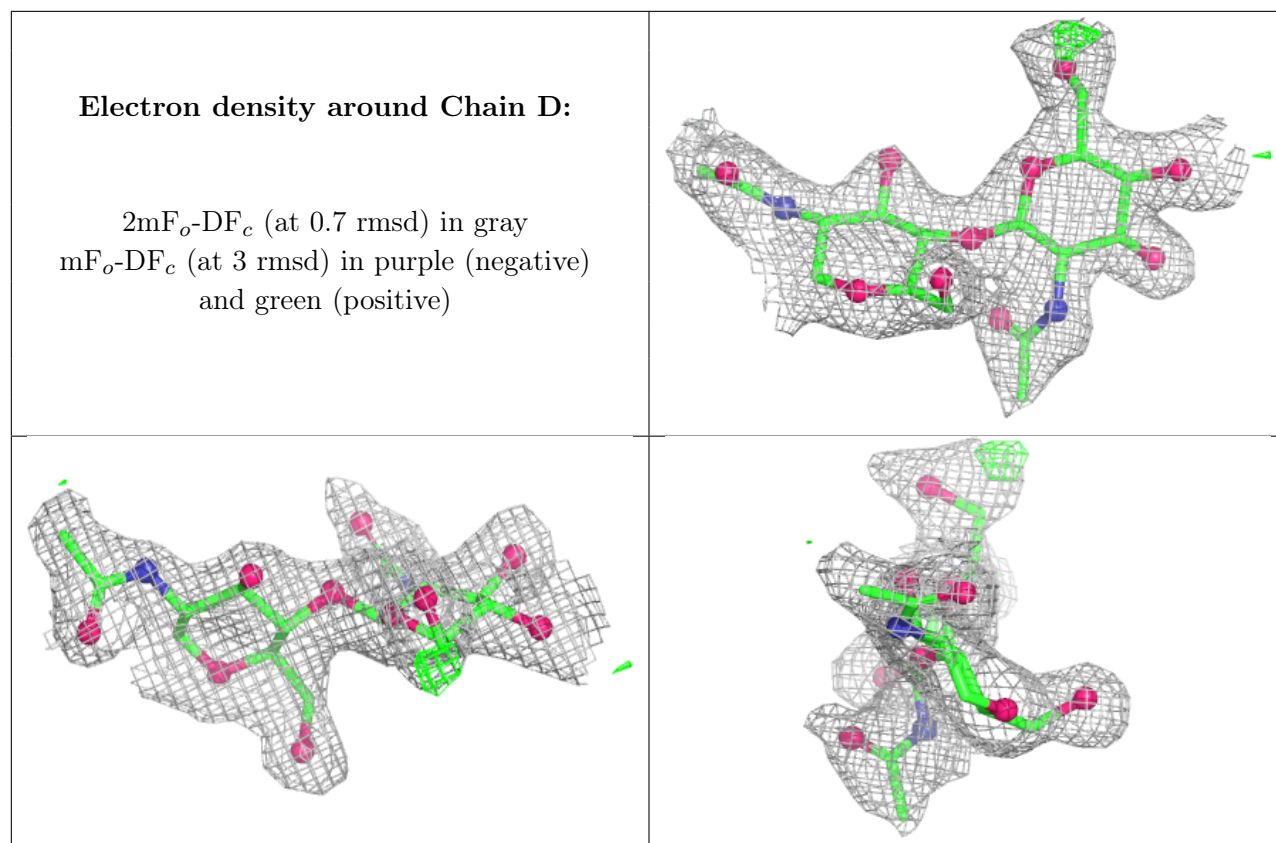
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	NAG	C	2	14/15	0.79	0.12	38,45,52,57	0
2	NAG	D	2	14/15	0.81	0.12	33,46,52,57	0
2	NAG	D	1	14/15	0.88	0.11	34,38,50,50	0
2	NAG	C	1	14/15	0.93	0.08	24,31,36,36	0

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)





6.4 Ligands [i](#)

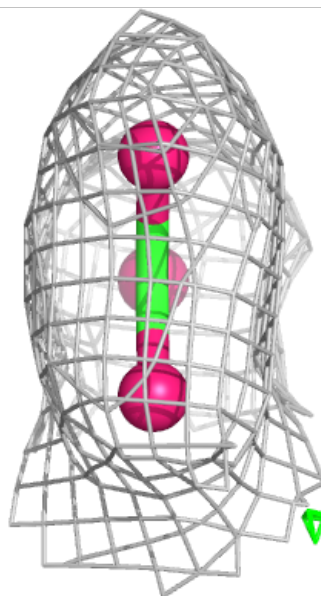
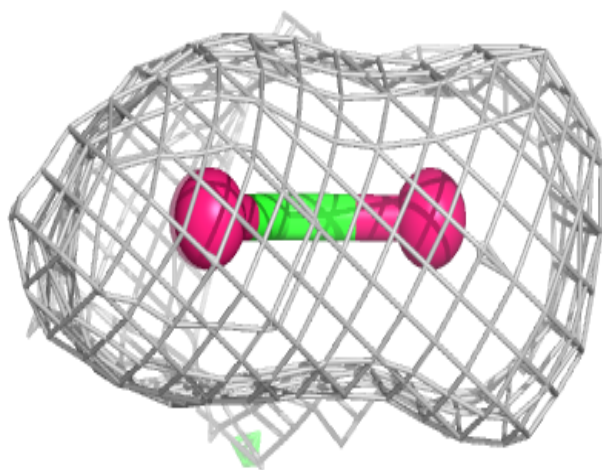
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	PG4	B	706	13/13	0.85	0.14	35,43,46,52	0
4	CO3	B	703	4/4	0.91	0.09	24,26,27,30	0
4	CO3	A	702	4/4	0.96	0.07	13,14,26,30	0
4	CO3	B	702	4/4	0.96	0.07	20,21,25,29	0
4	CO3	A	703	4/4	0.97	0.06	17,17,17,18	0
3	FE	B	701	1/1	0.99	0.02	23,23,23,23	0
3	FE	A	701	1/1	0.99	0.03	13,13,13,13	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.

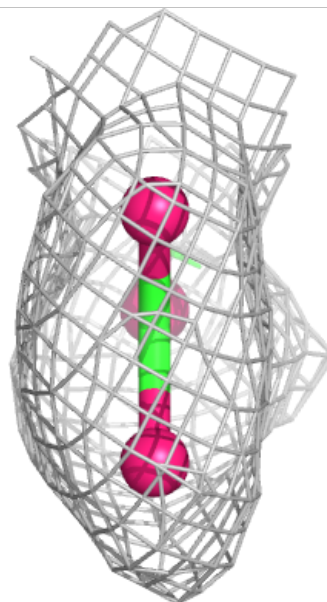
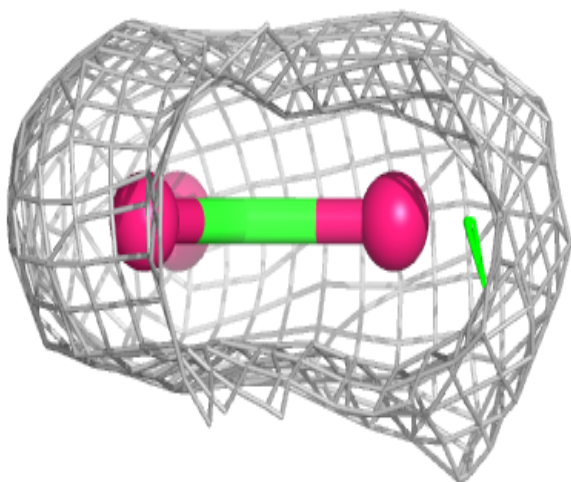
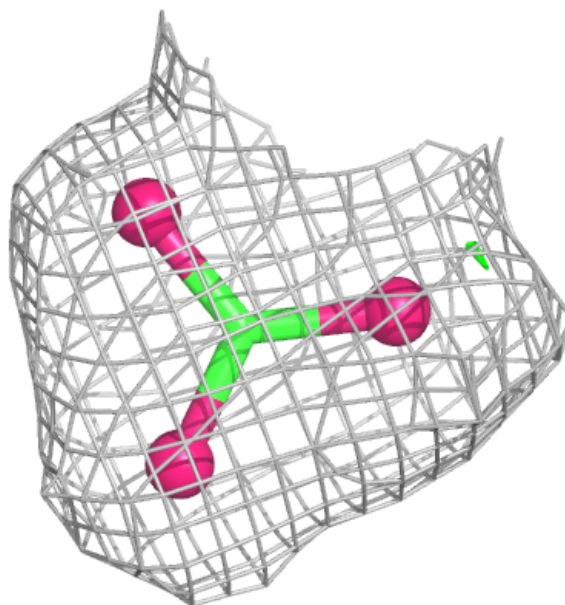
Electron density around CO3 B 703:

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



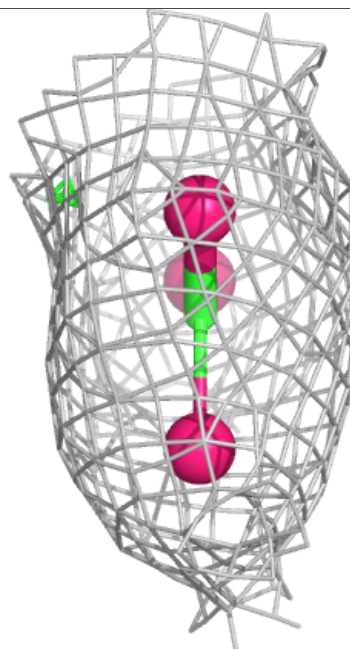
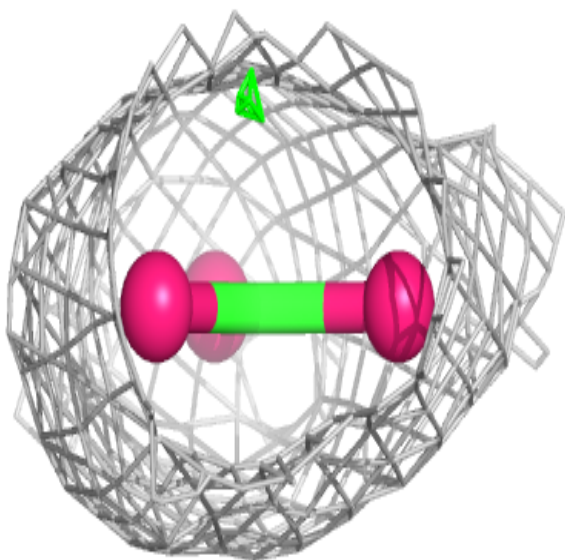
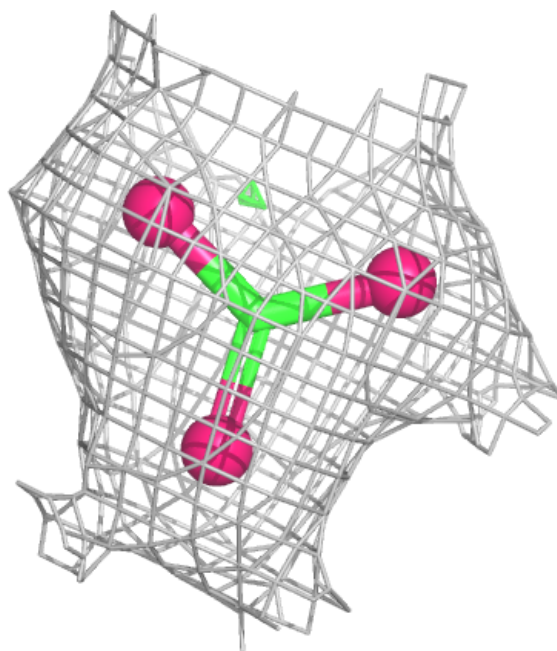
Electron density around CO3 A 702:

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



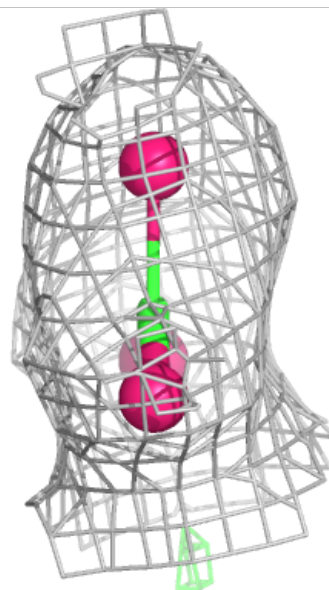
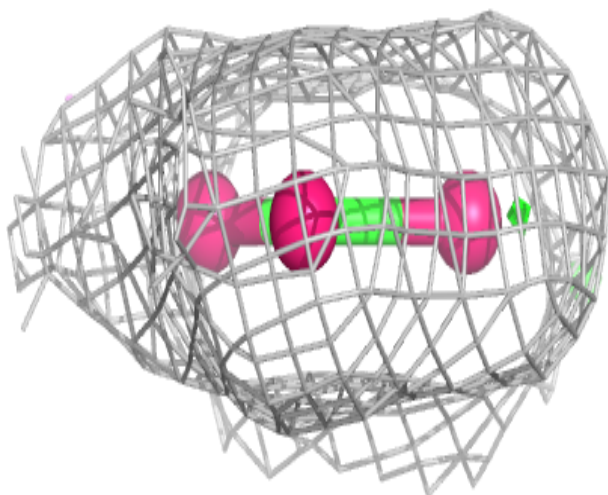
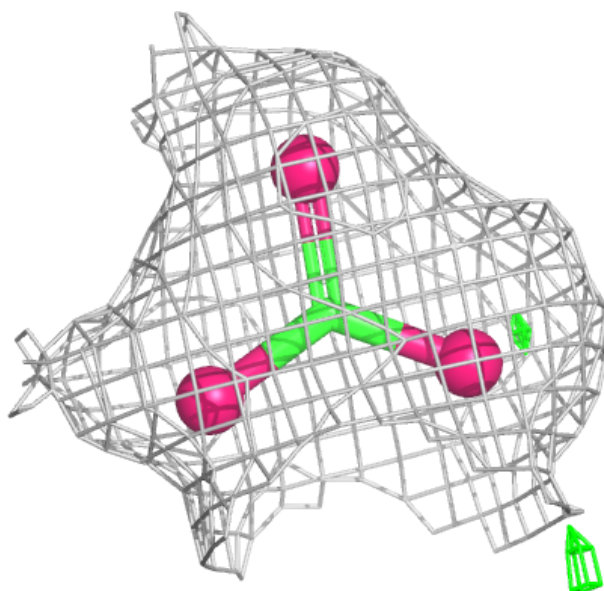
Electron density around CO3 B 702:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



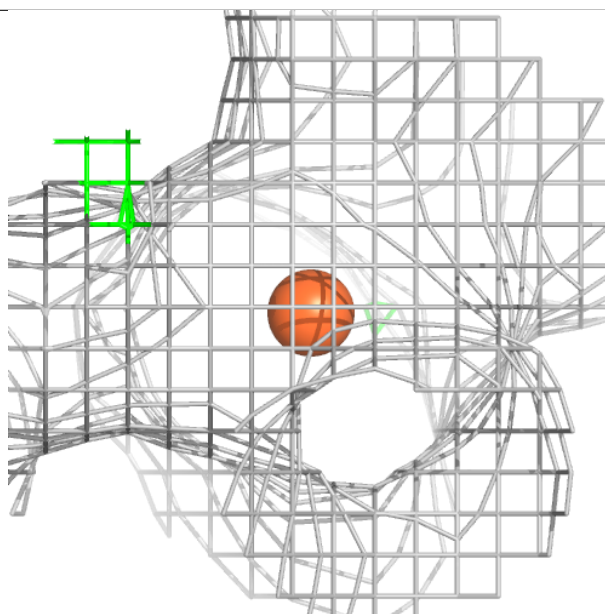
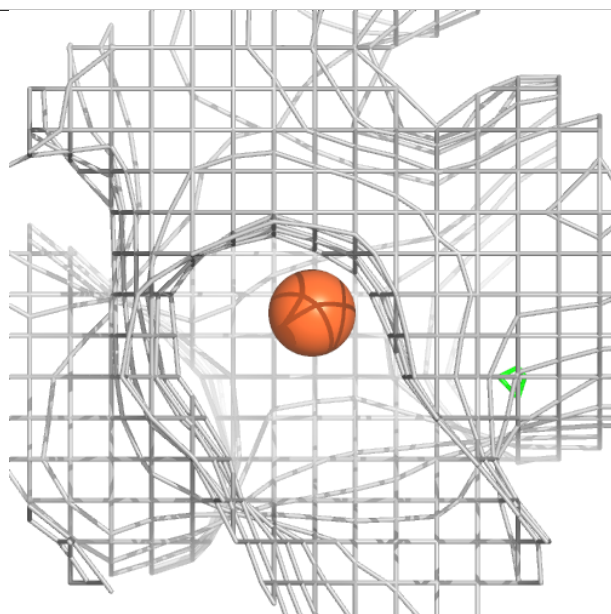
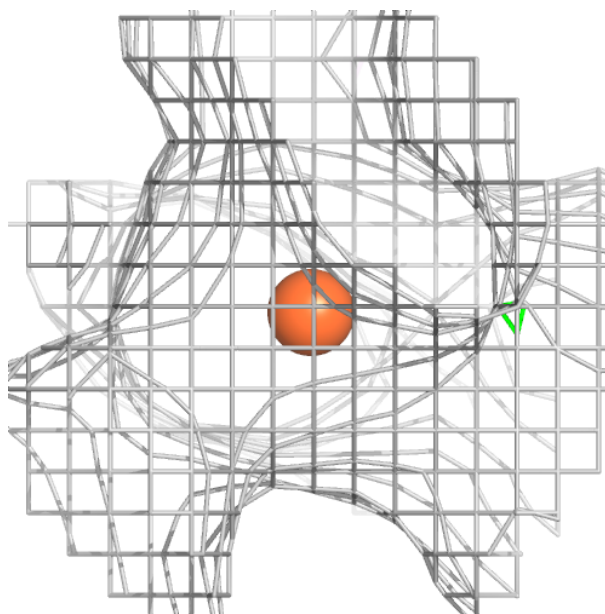
Electron density around CO3 A 703:

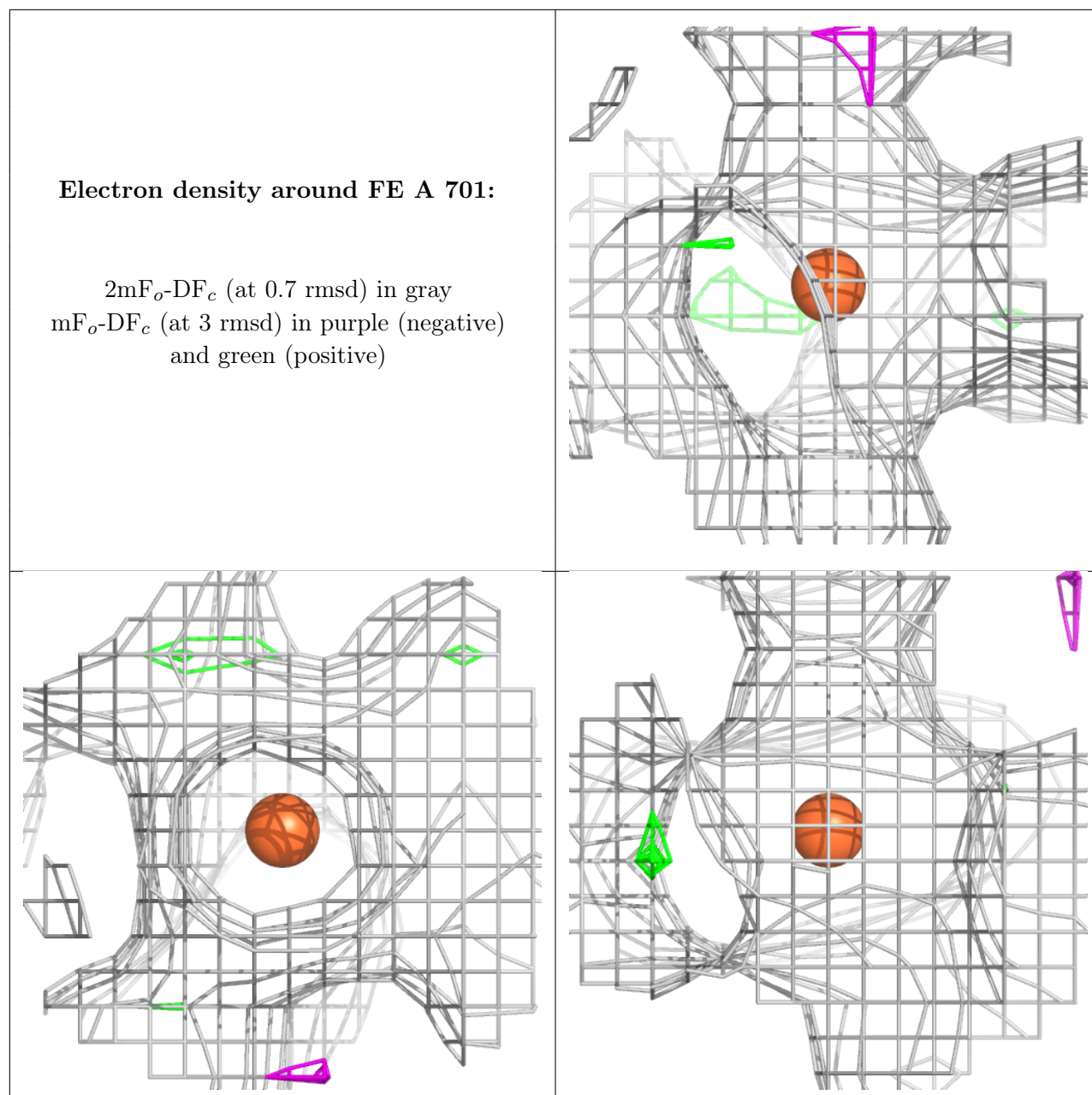
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around FE B 701:

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





6.5 Other polymers ⓘ

There are no such residues in this entry.