



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2026 – 01:06 PM UTC

PDB ID : 7LOX / pdb_00007lox
Title : The structure of Agmatinase from E. Coli at 3.2 Å displaying guanidine in the active site
Authors : Maturana, P.; Figueroa, M.; Gonzalez-Ordenes, F.; Villalobos, P.; Martinez-Oyanedel, J.; Uribe, E.A.; Castro-Fernandez, V.
Deposited on : 2021-02-11
Resolution : 3.20 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

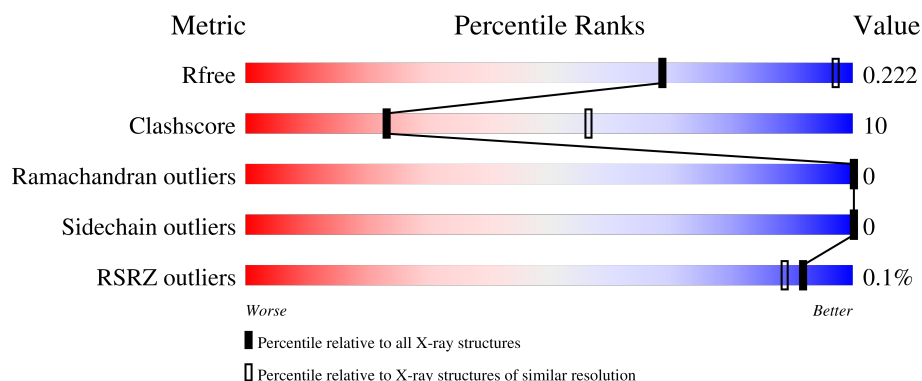
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.20 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	1466 (3.20-3.20)
Clashscore	190562	1573 (3.20-3.20)
Ramachandran outliers	187476	1548 (3.20-3.20)
Sidechain outliers	187428	1547 (3.20-3.20)
RSRZ outliers	180081	1466 (3.20-3.20)

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	306	 75% 17% 7%
1	B	306	 66% 26% 8%
1	C	306	 75% 17% 7%

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
3	GAI	A	403	-	X	-	-
3	GAI	A	404	-	X	-	-
3	GAI	C	403	-	X	-	-

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6539 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.

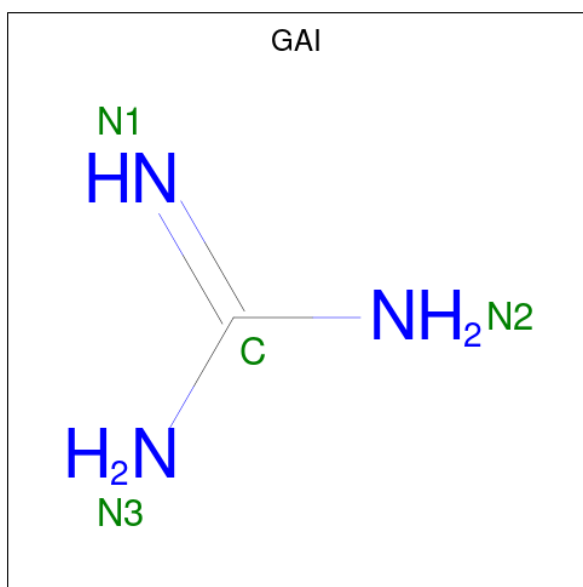
- Molecule 1 is a protein called Agmatinase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	284	Total	C	N	O	S	0	0	0
			2176	1385	374	404	13			
1	B	283	Total	C	N	O	S	0	0	0
			2156	1371	371	401	13			
1	C	285	Total	C	N	O	S	0	0	0
			2185	1390	374	408	13			

- Molecule 2 is MANGANESE (II) ION (CCD ID: MN) (formula: Mn) (labeled as "Ligand of Interest" by depositor).

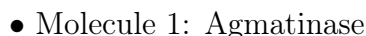
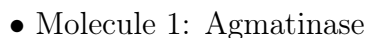
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total	Mn	0	0
			2	2		
2	B	2	Total	Mn	0	0
			2	2		
2	C	2	Total	Mn	0	0
			2	2		

- Molecule 3 is GUANIDINE (CCD ID: GAI) (formula: CH₅N₃) (labeled as "Ligand of Interest" by depositor).



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

- Molecule 1: Agmatinase



GLU

4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	129.03Å 129.03Å 88.29Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	64.51 – 3.20 64.51 – 3.20	Depositor EDS
% Data completeness (in resolution range)	99.9 (64.51-3.20) 95.8 (64.51-3.20)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.02 (at 2.86Å)	Xtriage
Refinement program	PHENIX 1.14 _3260	Depositor
R, R_{free}	0.172 , 0.218 0.176 , 0.222	Depositor DCC
R_{free} test set	920 reflections (4.67%)	wwPDB-VP
Wilson B-factor (Å ²)	69.8	Xtriage
Anisotropy	0.293	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 50.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.033 for -h,-k,l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6539	wwPDB-VP
Average B, all atoms (Å ²)	81.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.83% 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: MN, GAI

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	0.70	4/2226 (0.2%)	0.94	10/3023 (0.3%)
1	B	0.61	1/2206 (0.0%)	0.87	5/2997 (0.2%)
1	C	0.65	1/2236 (0.0%)	0.91	9/3038 (0.3%)
All	All	0.66	6/6668 (0.1%)	0.91	24/9058 (0.3%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	C	0	1

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	244	THR	C-N	13.51	1.50	1.33
1	C	244	THR	C-N	12.73	1.50	1.33
1	A	244	THR	C-N	12.50	1.49	1.33
1	A	41	PRO	C-O	7.63	1.29	1.24
1	A	41	PRO	CA-C	-6.30	1.46	1.52
1	A	39	GLY	CA-C	5.18	1.55	1.52

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	244	THR	CA-C-N	10.54	130.49	119.85
1	C	244	THR	C-N-CA	10.54	130.49	119.85
1	A	244	THR	CA-C-N	9.86	129.81	119.85
1	A	244	THR	C-N-CA	9.86	129.81	119.85

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	244	THR	CA-C-N	9.63	130.34	120.14
1	B	244	THR	C-N-CA	9.63	130.34	120.14
1	C	80	ARG	CG-CD-NE	7.50	128.49	112.00
1	C	148	PHE	CA-C-O	-7.35	112.50	120.36
1	A	149	ASP	CA-CB-CG	7.15	119.75	112.60
1	C	149	ASP	CA-CB-CG	6.48	119.08	112.60
1	A	148	PHE	CA-C-O	-6.40	113.27	120.43
1	A	243	GLY	CA-C-N	-6.33	115.37	123.16
1	A	243	GLY	C-N-CA	-6.33	115.37	123.16
1	B	243	GLY	CA-C-N	-6.26	115.57	123.15
1	B	243	GLY	C-N-CA	-6.26	115.57	123.15
1	B	149	ASP	CA-CB-CG	6.11	118.71	112.60
1	A	244	THR	CA-CB-OG1	-6.01	100.59	109.60
1	A	202	CYS	CA-C-N	5.88	128.44	120.44
1	A	202	CYS	C-N-CA	5.88	128.44	120.44
1	C	243	GLY	CA-C-N	-5.66	115.82	123.46
1	C	243	GLY	C-N-CA	-5.66	115.82	123.46
1	C	244	THR	CA-CB-OG1	-5.60	101.20	109.60
1	A	41	PRO	O-C-N	-5.60	118.75	122.73
1	C	244	THR	O-C-N	-5.33	116.51	121.36

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	167	PHE	Peptide

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	2176	0	2103	32	0
1	B	2156	0	2069	69	0
1	C	2185	0	2102	34	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
2	C	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	8	0	8	1	0
3	B	4	0	3	0	0
3	C	4	0	4	0	0
All	All	6539	0	6289	132	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 10.

All (132) 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:163:HIS:O	1:A:166:MET:SD	2.04	1.15
1:B:153:ASP:C	1:B:164:GLY:O	1.95	1.09
1:B:164:GLY:H	1:B:166:MET:HE2	1.17	1.07
1:B:127:PHE:CE2	1:B:161:PHE:HA	1.99	0.96
1:A:270:MET:HE1	1:A:300:GLN:HE21	1.43	0.82
1:B:232:ASP:HB3	1:B:242:THR:HG21	1.61	0.81
1:B:200:ASP:OD1	1:B:203:GLN:HG2	1.79	0.81
1:A:223:MET:HG2	1:A:224:PRO:HD2	1.64	0.78
1:A:147:HIS:HD2	1:A:183:GLN:HG2	1.49	0.76
1:B:38:THR:HG21	1:B:108:HIS:NE2	2.00	0.76
1:B:256:ILE:HD11	1:B:295:GLU:HG3	1.68	0.76
1:B:187:ARG:NH2	1:B:232:ASP:O	2.20	0.74
1:B:127:PHE:HE2	1:B:161:PHE:HA	1.49	0.74
1:A:147:HIS:CD2	1:A:183:GLN:HG2	2.22	0.73
1:B:43:ASP:OD2	1:B:53:ARG:HG2	1.89	0.72
1:B:153:ASP:CA	1:B:164:GLY:O	2.36	0.72
1:C:239:ALA:HB1	1:C:285:THR:HG23	1.69	0.72
1:B:38:THR:HG22	1:B:88:CYS:SG	2.30	0.71
1:B:144:ALA:HB3	1:B:225:VAL:HG12	1.74	0.69
1:C:41:PRO:O	1:C:93:TYR:OH	2.12	0.68
1:B:256:ILE:HD12	1:B:292:LEU:HD23	1.75	0.67
1:B:217:LYS:HG2	1:B:265:LEU:HD21	1.77	0.66
1:B:252:SER:HB2	1:C:247:ILE:HD12	1.77	0.65
1:B:270:MET:HE3	1:B:300:GLN:HE21	1.61	0.64
1:A:223:MET:CG	1:A:224:PRO:HD2	2.30	0.61
1:C:106:GLN:OE1	1:C:134:ARG:HB3	2.00	0.60
1:C:216:VAL:O	1:C:220:VAL:HG23	2.01	0.60
1:C:129:THR:HG22	1:C:133:LEU:HD21	1.85	0.59
1:B:153:ASP:CB	1:B:164:GLY:HA2	2.32	0.59
1:C:129:THR:HG22	1:C:133:LEU:CD2	2.33	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:170:ALA:HA	1:B:175:LEU:HD12	1.85	0.58
1:B:41:PRO:O	1:B:93:TYR:OH	2.14	0.58
1:B:147:HIS:ND1	1:B:183:GLN:HG2	2.19	0.57
1:B:33:ALA:HA	1:B:84:ASN:HB3	1.86	0.57
1:B:153:ASP:O	1:B:164:GLY:O	2.22	0.56
1:B:164:GLY:N	1:B:166:MET:HE2	2.03	0.56
1:C:270:MET:SD	1:C:296:MET:HB3	2.47	0.55
1:A:270:MET:SD	1:A:296:MET:HB3	2.47	0.55
1:A:259:VAL:O	1:A:299:ILE:HD11	2.06	0.55
1:B:153:ASP:HA	1:B:164:GLY:O	2.06	0.55
1:A:270:MET:HE1	1:A:300:GLN:NE2	2.17	0.55
1:B:153:ASP:CG	1:B:164:GLY:HA2	2.31	0.55
1:B:77:PHE:N	1:B:77:PHE:CD2	2.74	0.55
1:B:36:VAL:HG21	1:B:112:LEU:CD1	2.37	0.54
1:B:24:MET:HA	1:B:87:ASP:O	2.08	0.54
1:A:256:ILE:HD12	1:A:292:LEU:HD23	1.89	0.54
1:A:187:ARG:HD3	1:A:249:GLY:HA2	1.89	0.54
1:C:187:ARG:HD3	1:C:249:GLY:HA2	1.89	0.54
1:A:17:PHE:HB3	1:A:24:MET:HG2	1.90	0.53
1:B:144:ALA:O	1:B:225:VAL:HA	2.08	0.53
1:B:126:HIS:CE1	1:B:230:ASP:HB2	2.44	0.53
1:B:227:LEU:O	1:B:270:MET:HA	2.07	0.53
1:B:75:TRP:CD1	1:B:298:TYR:CD2	2.97	0.52
1:C:106:GLN:OE1	1:C:134:ARG:CB	2.57	0.52
1:B:256:ILE:CD1	1:B:295:GLU:HG3	2.38	0.51
1:C:125:ASP:OD2	1:C:163:HIS:HB3	2.10	0.51
1:C:130:LEU:HG	1:C:166:MET:HB2	1.93	0.51
1:B:294:LEU:HD11	1:B:298:TYR:HE1	1.76	0.50
1:C:95:PHE:CD1	1:C:95:PHE:C	2.89	0.50
1:C:127:PHE:CE1	1:C:161:PHE:HA	2.46	0.50
1:B:229:PHE:CE2	1:B:292:LEU:HD13	2.47	0.50
1:C:227:LEU:O	1:C:270:MET:HA	2.12	0.50
1:C:244:THR:O	1:C:244:THR:OG1	2.24	0.50
1:A:223:MET:HG2	1:A:224:PRO:CD	2.40	0.49
1:B:73:PHE:C	1:B:73:PHE:CD1	2.90	0.49
1:B:98:ALA:HB3	1:B:161:PHE:CD2	2.48	0.49
1:B:129:THR:HA	1:B:132:LEU:HD12	1.94	0.49
1:B:29:TYR:CG	1:B:111:LYS:HE3	2.47	0.49
1:B:71:ASN:O	1:C:189:GLU:HB2	2.12	0.49
1:C:145:LEU:HD22	1:C:176:ILE:HD13	1.93	0.49
1:B:43:ASP:HB2	1:B:53:ARG:HH21	1.77	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:159:CYS:SG	1:C:160:GLU:N	2.85	0.48
1:B:126:HIS:CD2	1:B:273:VAL:HG21	2.48	0.48
1:B:153:ASP:HA	1:B:164:GLY:HA2	1.95	0.48
1:A:203:GLN:HG2	1:A:207:ARG:HD3	1.95	0.47
1:A:37:ILE:HA	1:A:120:LEU:O	2.15	0.47
1:C:200:ASP:OD1	1:C:203:GLN:HG2	2.15	0.47
1:A:57:ALA:O	1:A:61:GLN:HG3	2.15	0.46
1:A:79:MET:HB2	1:A:298:TYR:CE1	2.51	0.46
1:C:37:ILE:HA	1:C:120:LEU:O	2.15	0.46
1:B:29:TYR:CD1	1:B:111:LYS:HG2	2.50	0.46
1:B:35:TRP:CZ3	1:B:118:ARG:HB3	2.50	0.46
1:A:235:ASP:OD1	1:A:236:PRO:HD2	2.16	0.45
1:B:166:MET:H	1:B:166:MET:HG3	1.49	0.45
1:A:228:THR:HA	1:A:271:ASP:OD1	2.17	0.45
1:B:127:PHE:CE2	1:B:161:PHE:CA	2.86	0.45
1:B:130:LEU:HB3	1:B:131:PRO:HD3	1.99	0.45
1:A:126:HIS:CD2	1:A:273:VAL:HG21	2.51	0.45
1:C:126:HIS:CD2	1:C:273:VAL:HG21	2.51	0.45
1:B:35:TRP:CD2	1:B:118:ARG:HB2	2.52	0.45
1:B:123:GLY:HA3	1:B:273:VAL:HG12	1.98	0.45
1:A:69:GLU:OE2	1:B:188:THR:HG22	2.17	0.45
1:B:105:LEU:HD23	1:B:131:PRO:HB2	1.98	0.45
1:C:95:PHE:C	1:C:95:PHE:HD1	2.26	0.44
1:B:154:THR:N	1:B:164:GLY:O	2.48	0.44
1:C:253:ASP:OD1	1:C:253:ASP:C	2.60	0.44
1:A:238:PHE:O	1:A:284:ILE:HG22	2.18	0.44
1:B:69:GLU:O	1:B:80:ARG:NH2	2.50	0.44
1:B:127:PHE:HE2	1:B:162:ASP:H	1.64	0.44
1:A:153:ASP:OD2	3:A:403:GAI:N3	2.51	0.43
1:C:16:ALA:HA	1:C:21:ARG:O	2.19	0.43
1:A:29:TYR:CD1	1:A:111:LYS:HG2	2.53	0.43
1:B:31:SER:O	1:B:117:LYS:NZ	2.49	0.43
1:B:36:VAL:HG21	1:B:112:LEU:HD12	1.99	0.43
1:A:24:MET:HA	1:A:87:ASP:O	2.19	0.43
1:B:72:ARG:HB3	1:B:298:TYR:CZ	2.53	0.43
1:B:223:MET:CG	1:B:224:PRO:HD2	2.49	0.43
1:B:185:GLY:O	1:B:187:ARG:HG3	2.17	0.42
1:B:127:PHE:CE2	1:B:162:ASP:N	2.87	0.42
1:C:142:LYS:HD3	1:C:177:ASP:HB2	2.02	0.42
1:C:78:ASP:OD1	1:C:80:ARG:HG2	2.20	0.42
1:C:253:ASP:O	1:C:256:ILE:HG22	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:32:ASP:OD1	1:A:32:ASP:N	2.48	0.41
1:B:187:ARG:NH1	1:B:233:CYS:O	2.44	0.41
1:B:108:HIS:O	1:B:112:LEU:HD23	2.19	0.41
1:A:105:LEU:HD11	1:A:132:LEU:HD21	2.01	0.41
1:B:123:GLY:HA2	1:B:124:GLY:HA3	1.80	0.41
1:C:43:ASP:OD1	1:C:43:ASP:N	2.53	0.41
1:A:41:PRO:O	1:A:93:TYR:OH	2.25	0.41
1:A:155:TYR:HB3	1:A:165:THR:HG22	2.03	0.41
1:C:24:MET:HA	1:C:87:ASP:O	2.20	0.41
1:A:72:ARG:HB3	1:A:298:TYR:CZ	2.56	0.41
1:A:229:PHE:C	1:A:229:PHE:CD1	2.99	0.41
1:B:215:GLN:O	1:B:219:ILE:HG12	2.20	0.41
1:C:75:TRP:CD1	1:C:298:TYR:CD2	3.09	0.41
1:A:125:ASP:OD2	1:A:163:HIS:HB3	2.21	0.40
1:C:162:ASP:OD1	1:C:163:HIS:N	2.55	0.40
1:B:37:ILE:HA	1:B:120:LEU:O	2.21	0.40
1:B:242:THR:HG22	1:B:244:THR:H	1.87	0.40
1:C:127:PHE:CD2	1:C:127:PHE:C	2.99	0.40
1:C:126:HIS:CE1	1:C:230:ASP:HB2	2.56	0.40
1:B:94:ALA:O	1:B:95:PHE:C	2.64	0.40

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	280/306 (92%)	275 (98%)	5 (2%)	0	100	100
1	B	279/306 (91%)	274 (98%)	5 (2%)	0	100	100
1	C	281/306 (92%)	277 (99%)	4 (1%)	0	100	100
All	All	840/918 (92%)	826 (98%)	14 (2%)	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	222/247 (90%)	222 (100%)	0	100	100
1	B	219/247 (89%)	219 (100%)	0	100	100
1	C	224/247 (91%)	224 (100%)	0	100	100
All	All	665/741 (90%)	665 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

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	70	HIS
1	A	147	HIS
1	A	163	HIS
1	A	281	GLN
1	A	300	GLN
1	B	70	HIS
1	B	300	GLN
1	C	163	HIS
1	C	218	GLN

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](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 10 ligands modelled in this entry, 6 are monoatomic - leaving 4 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
3	GAI	C	403	2	3,3,3	4.55	3 (100%)	3,3,3	1.09	0
3	GAI	A	403	2	3,3,3	4.56	3 (100%)	3,3,3	1.10	0
3	GAI	B	403	2	3,3,3	3.28	2 (66%)	3,3,3	1.09	0
3	GAI	A	404	-	3,3,3	2.93	3 (100%)	3,3,3	1.58	1 (33%)

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	403	GAI	C-N1	5.68	1.45	1.30
3	A	403	GAI	C-N1	5.67	1.45	1.30
3	C	403	GAI	C-N3	-3.95	1.25	1.35
3	A	403	GAI	C-N2	-3.94	1.25	1.35
3	B	403	GAI	C-N3	3.83	1.45	1.35
3	A	403	GAI	C-N3	3.83	1.45	1.35
3	B	403	GAI	C-N2	3.82	1.45	1.35
3	C	403	GAI	C-N2	3.80	1.45	1.35
3	A	404	GAI	C-N2	3.18	1.44	1.35
3	A	404	GAI	C-N3	3.04	1.43	1.35
3	A	404	GAI	C-N1	-2.55	1.23	1.30

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	404	GAI	N3-C-N2	2.35	121.67	116.18

There are no chirality outliers.

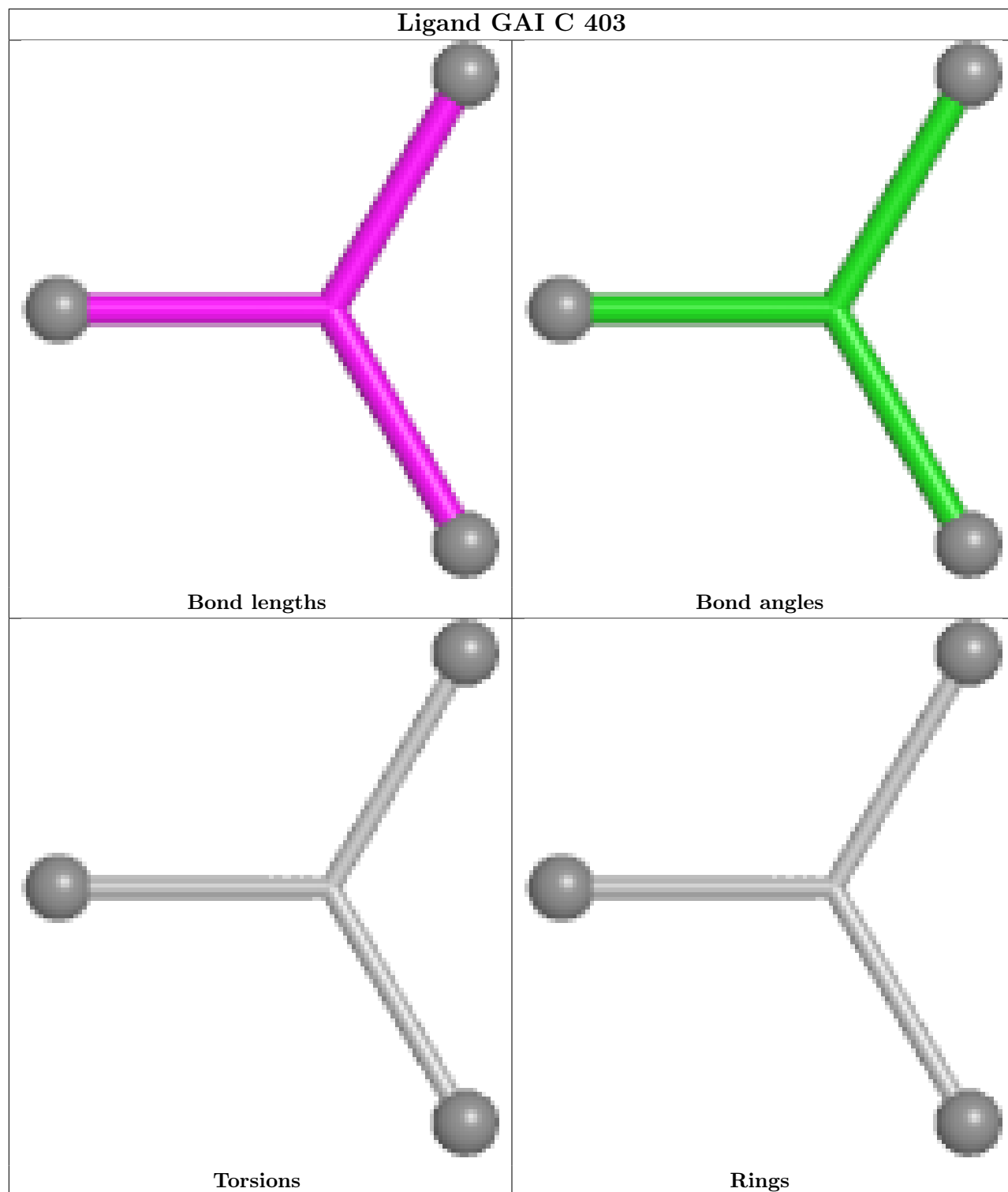
There are no torsion outliers.

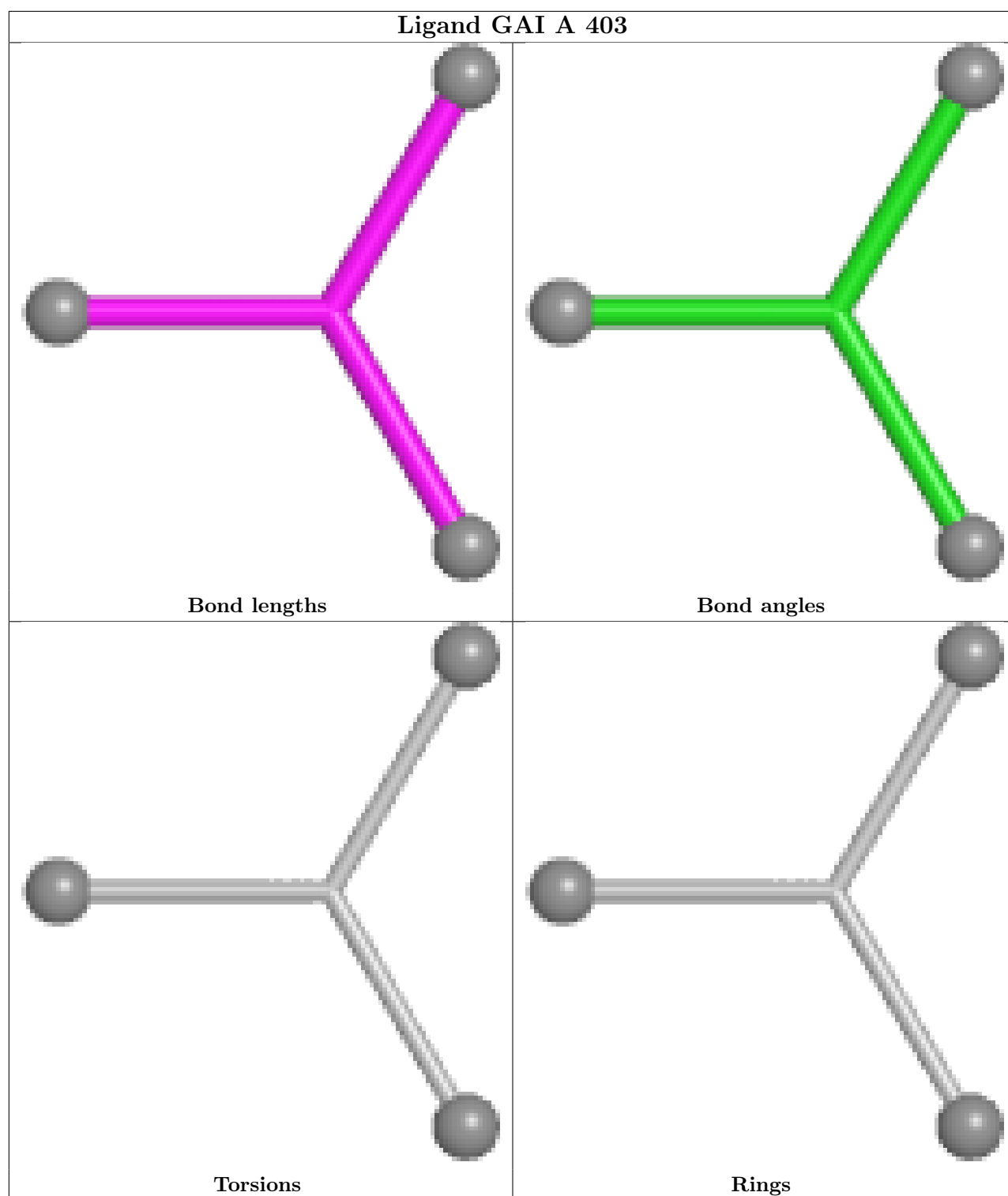
There are no ring outliers.

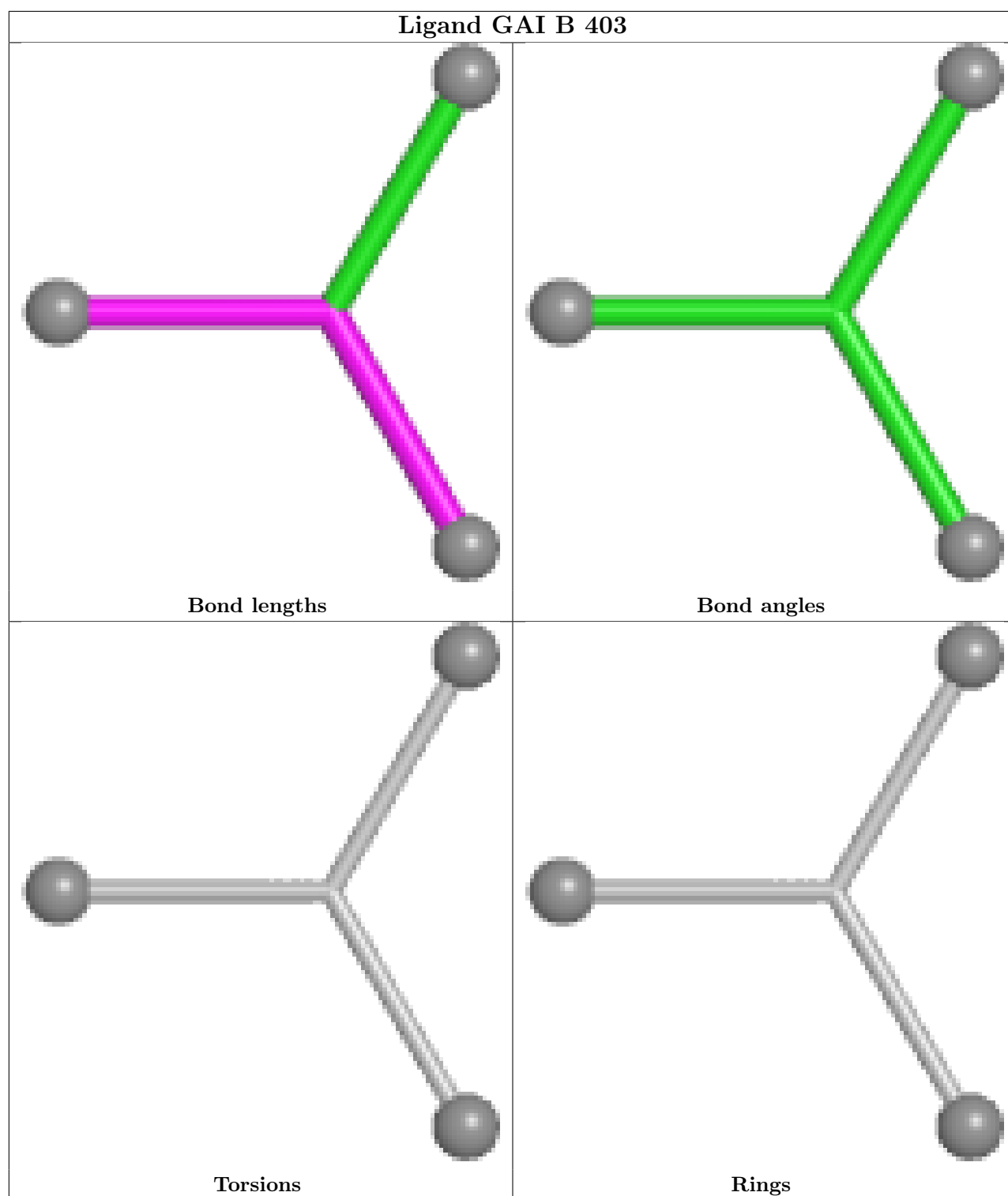
1 monomer is involved in 1 short contact:

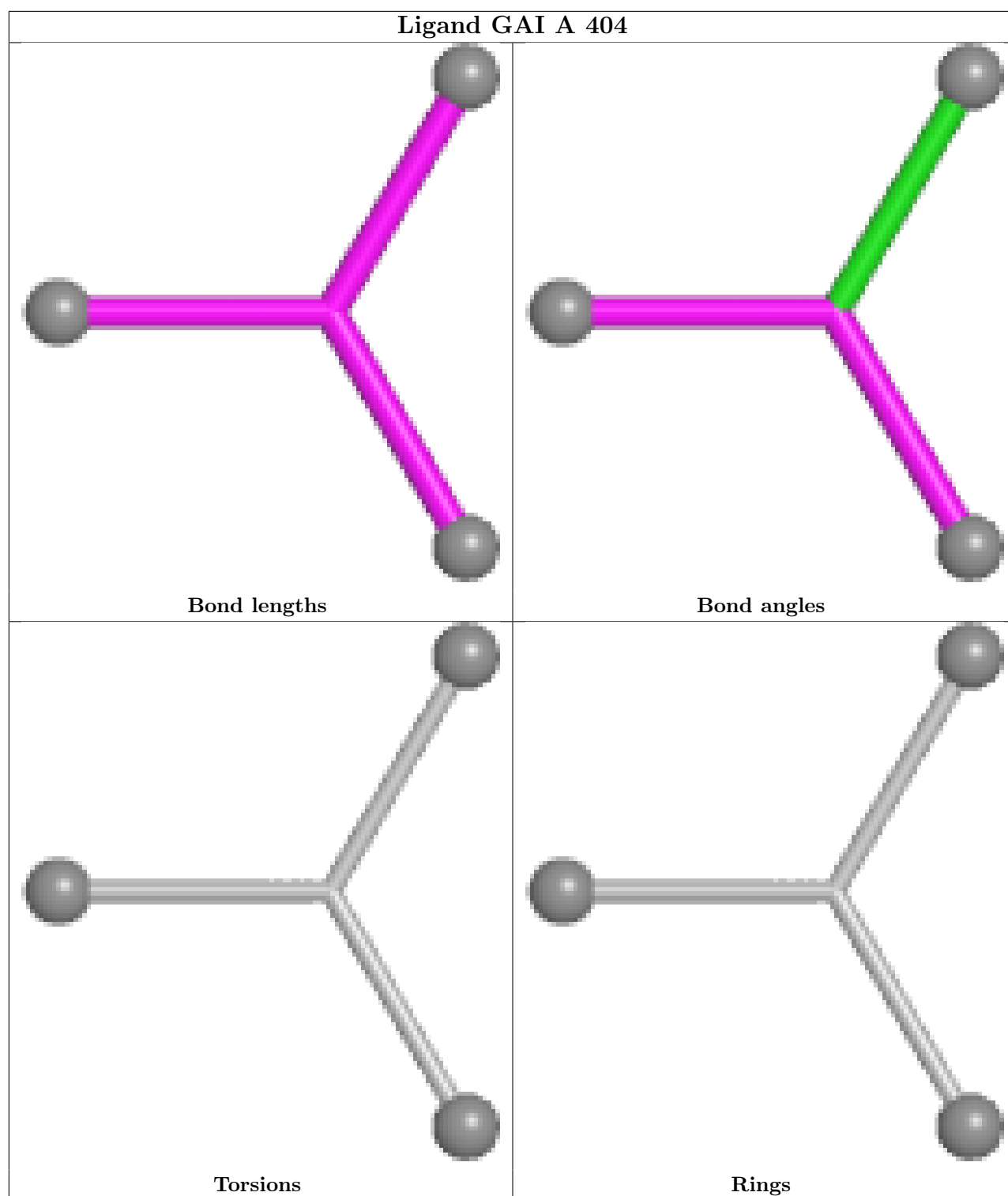
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	403	GAI	1	0

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 [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	284/306 (92%)	-0.64	0	100 100	61, 73, 96, 112	0
1	B	283/306 (92%)	-0.49	0	100 100	64, 89, 110, 122	0
1	C	285/306 (93%)	-0.60	1 (0%)	88 79	56, 76, 104, 119	0
All	All	852/918 (92%)	-0.58	1 (0%)	92 89	56, 79, 104, 122	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	44	MET	2.6

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](#)

There are no oligosaccharides in this entry.

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
3	GAI	C	403	4/4	0.78	0.14	74,80,81,83	0
3	GAI	A	404	4/4	0.85	0.12	70,73,81,84	0

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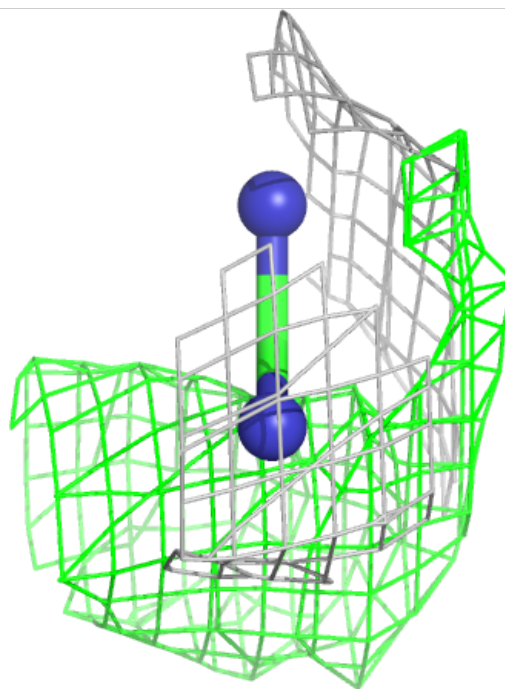
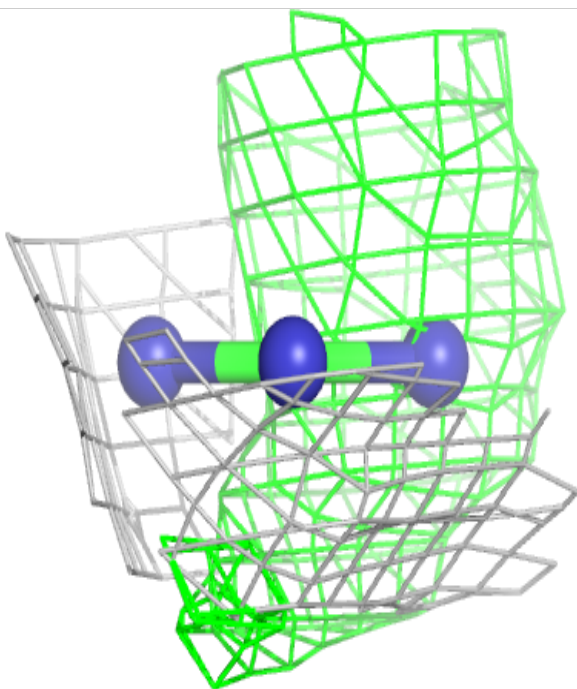
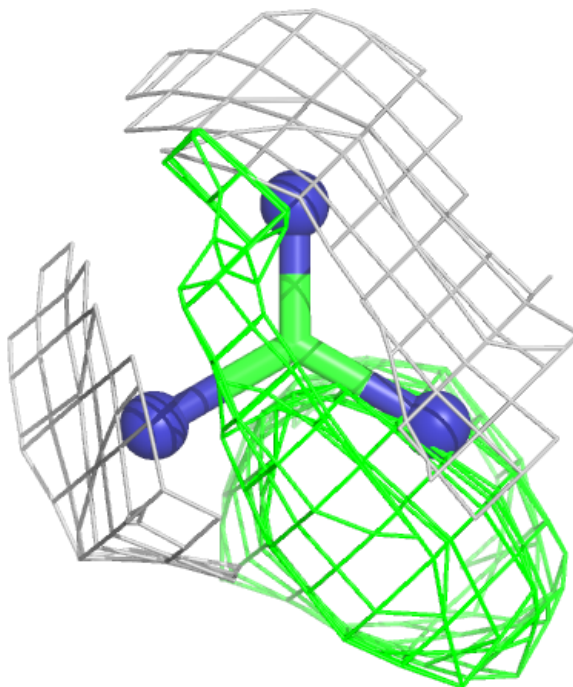
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	GAI	B	403	4/4	0.86	0.15	83,83,84,85	0
3	GAI	A	403	4/4	0.88	0.12	58,63,64,73	0
2	MN	C	401	1/1	0.95	0.06	83,83,83,83	1
2	MN	B	401	1/1	0.96	0.06	79,79,79,79	1
2	MN	A	401	1/1	0.98	0.06	68,68,68,68	1
2	MN	A	402	1/1	0.99	0.04	65,65,65,65	1
2	MN	C	402	1/1	0.99	0.05	82,82,82,82	0
2	MN	B	402	1/1	0.99	0.07	79,79,79,79	1

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.

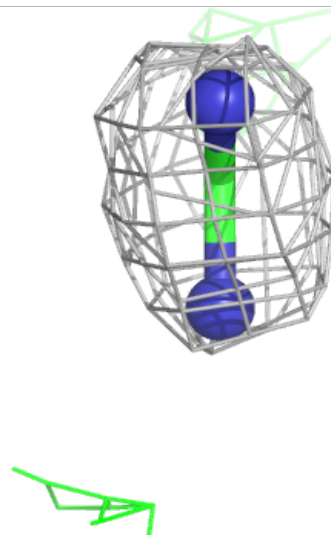
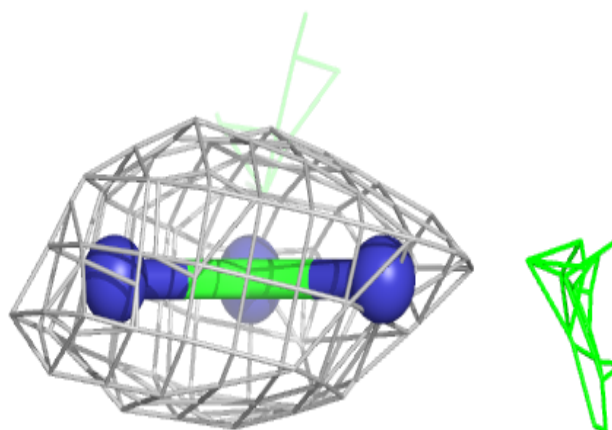
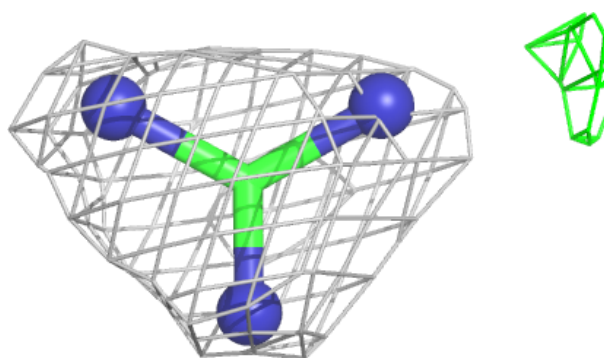
Electron density around GAI C 403:

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



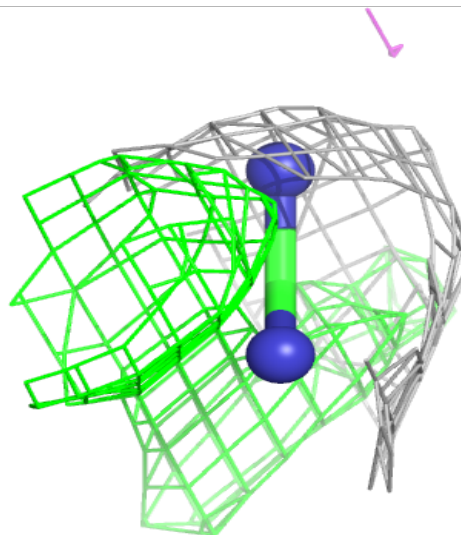
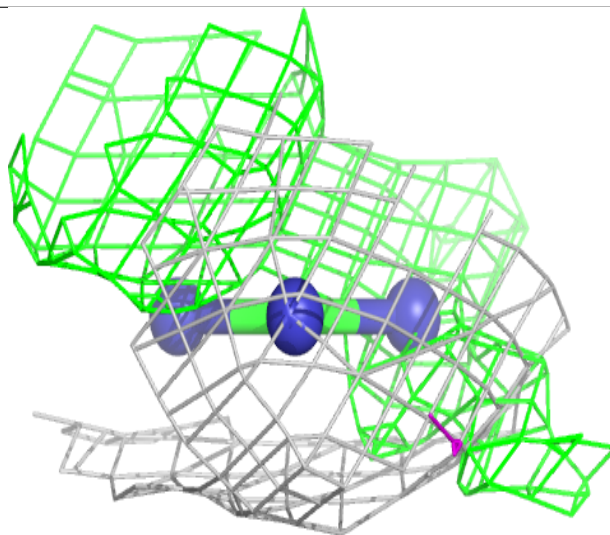
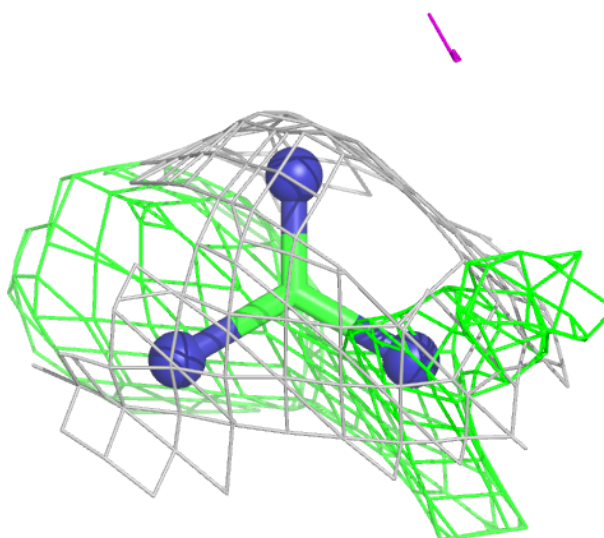
Electron density around GAI A 404:

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



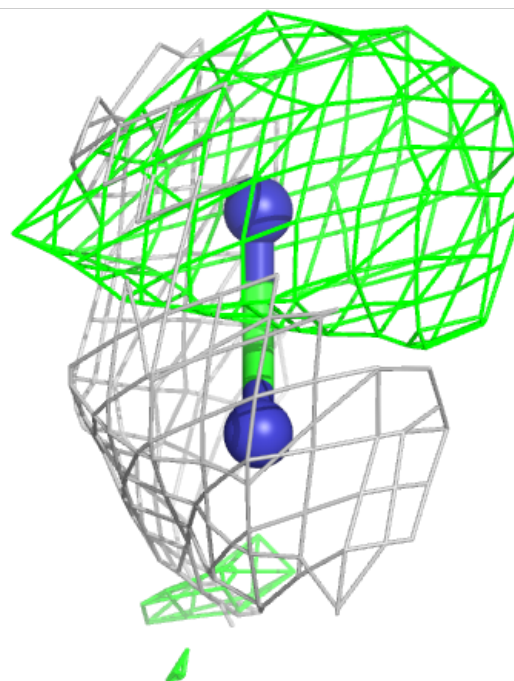
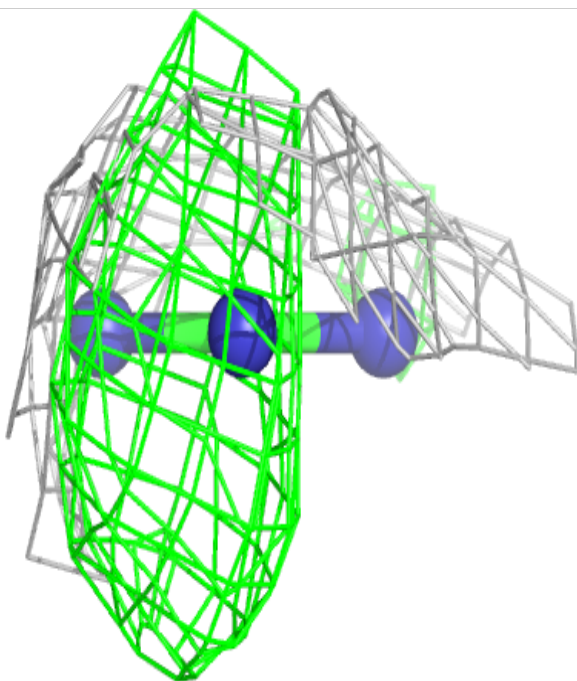
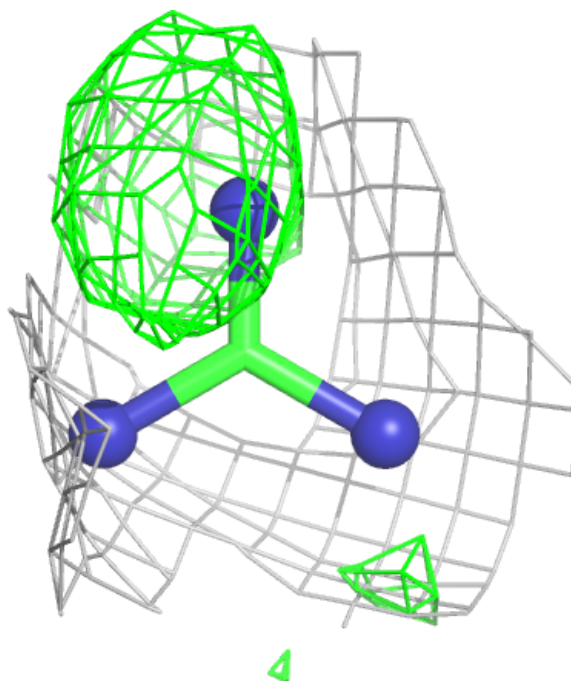
Electron density around GAI B 403:

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and green (positive)



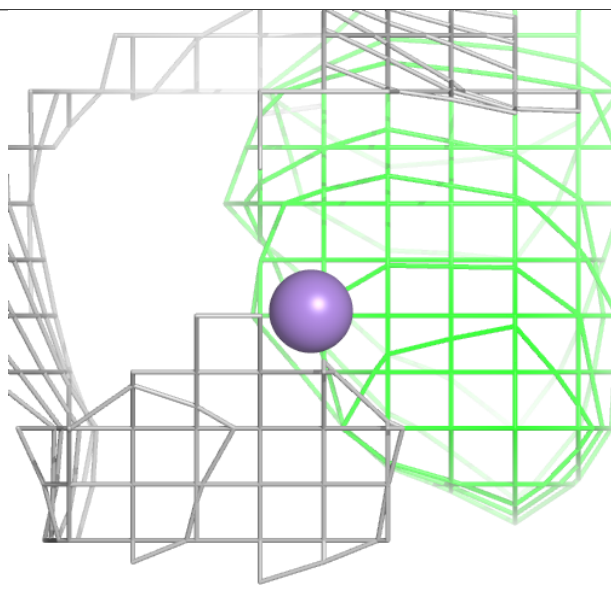
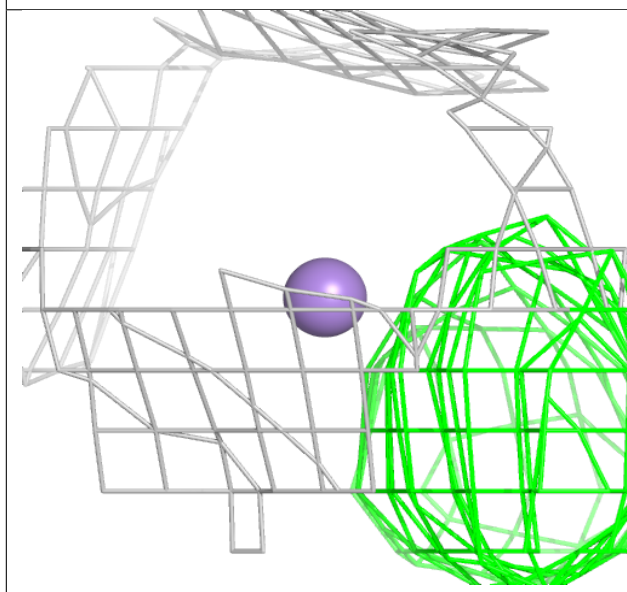
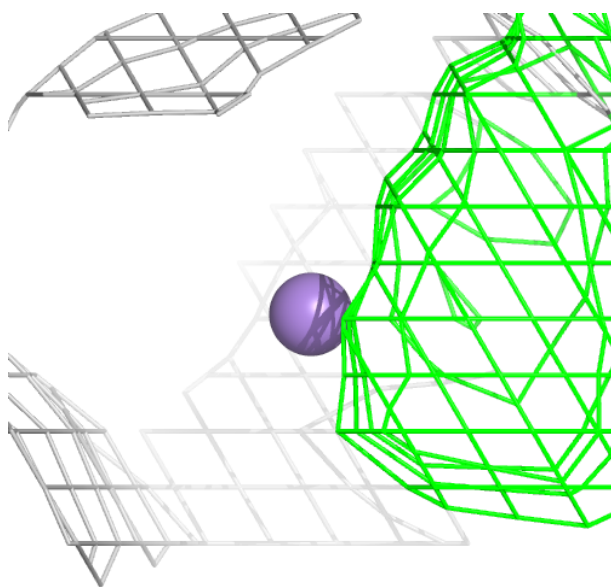
Electron density around GAI A 403:

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



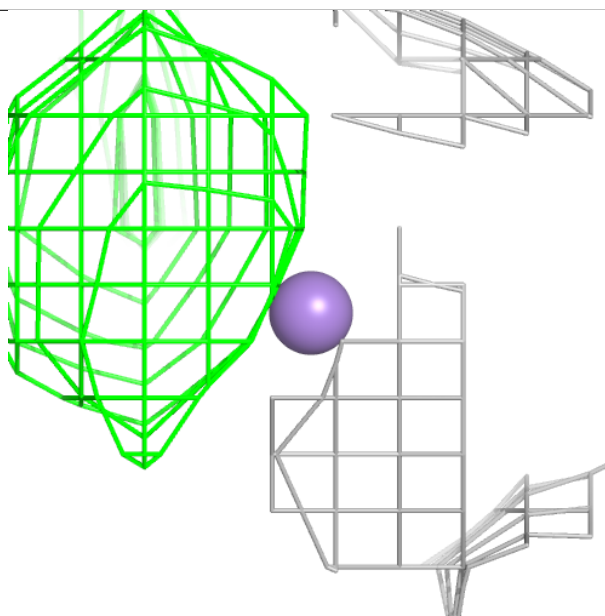
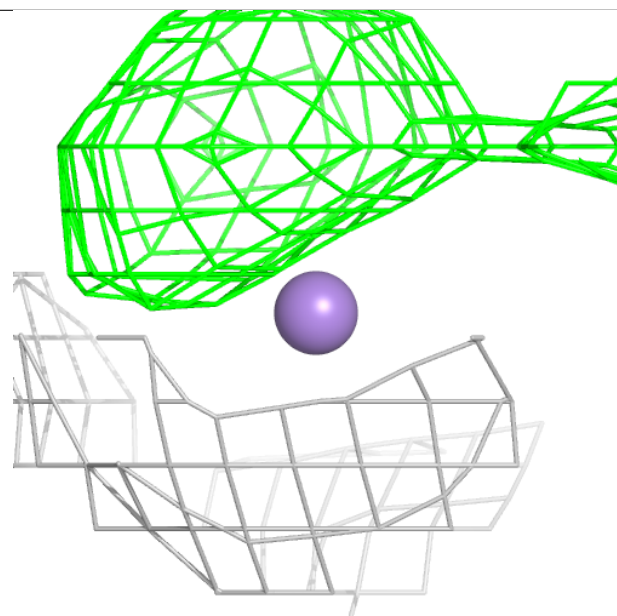
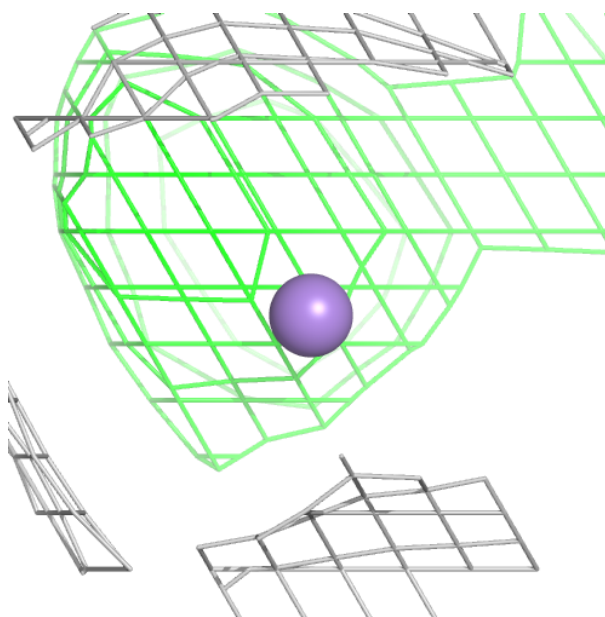
Electron density around MN C 401:

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



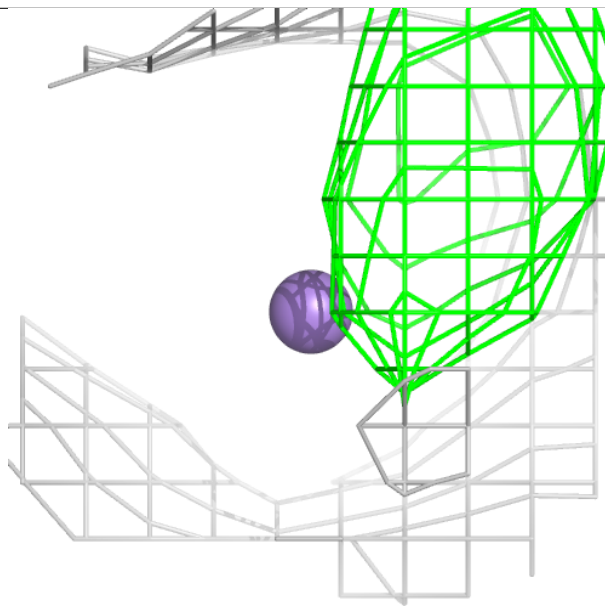
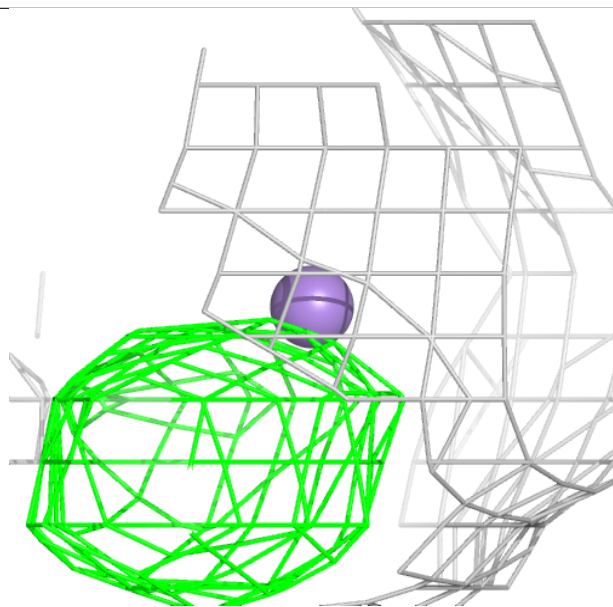
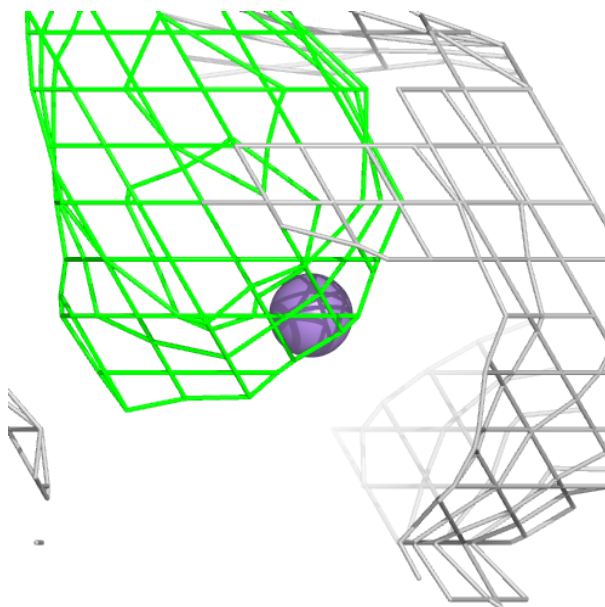
Electron density around MN B 401:

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and green (positive)



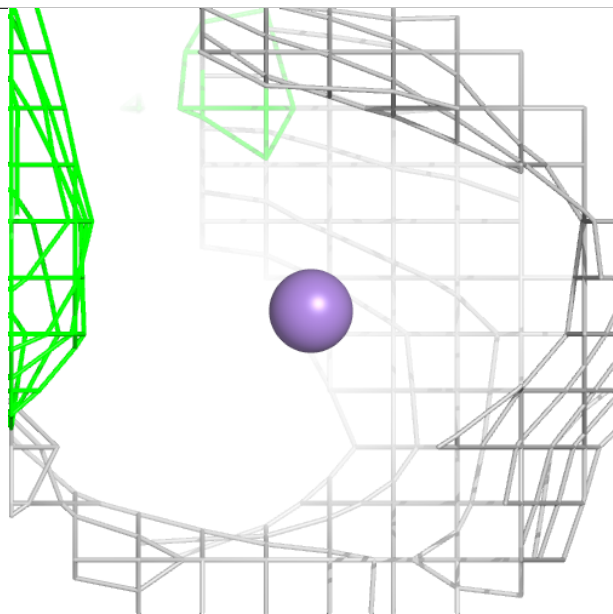
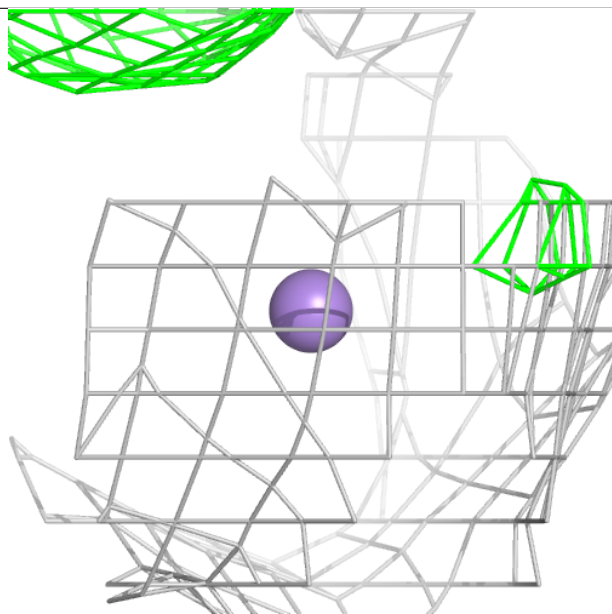
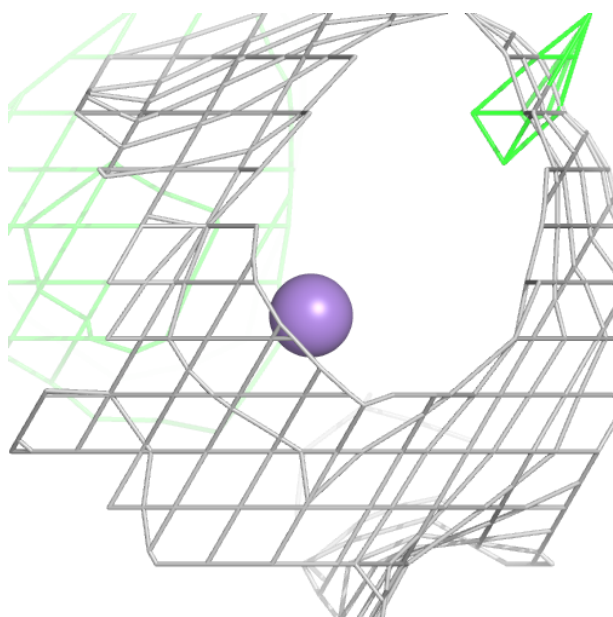
Electron density around MN A 401:

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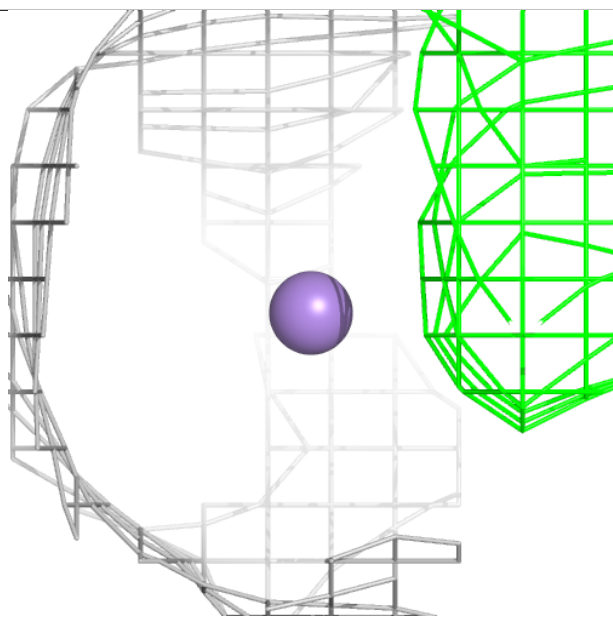
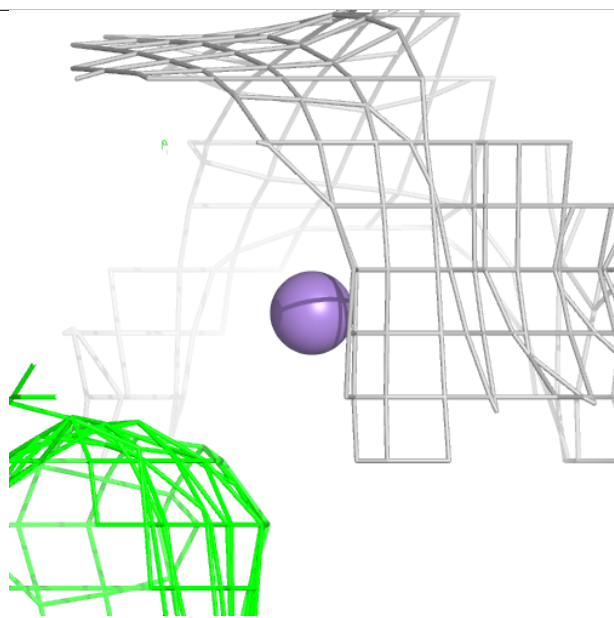
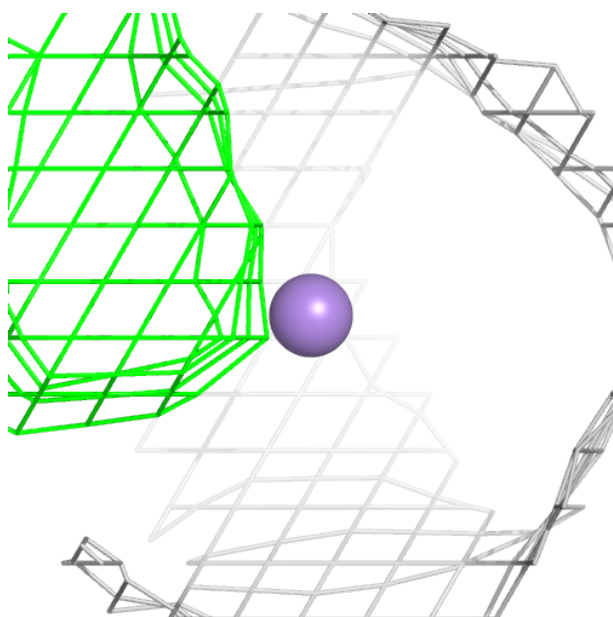
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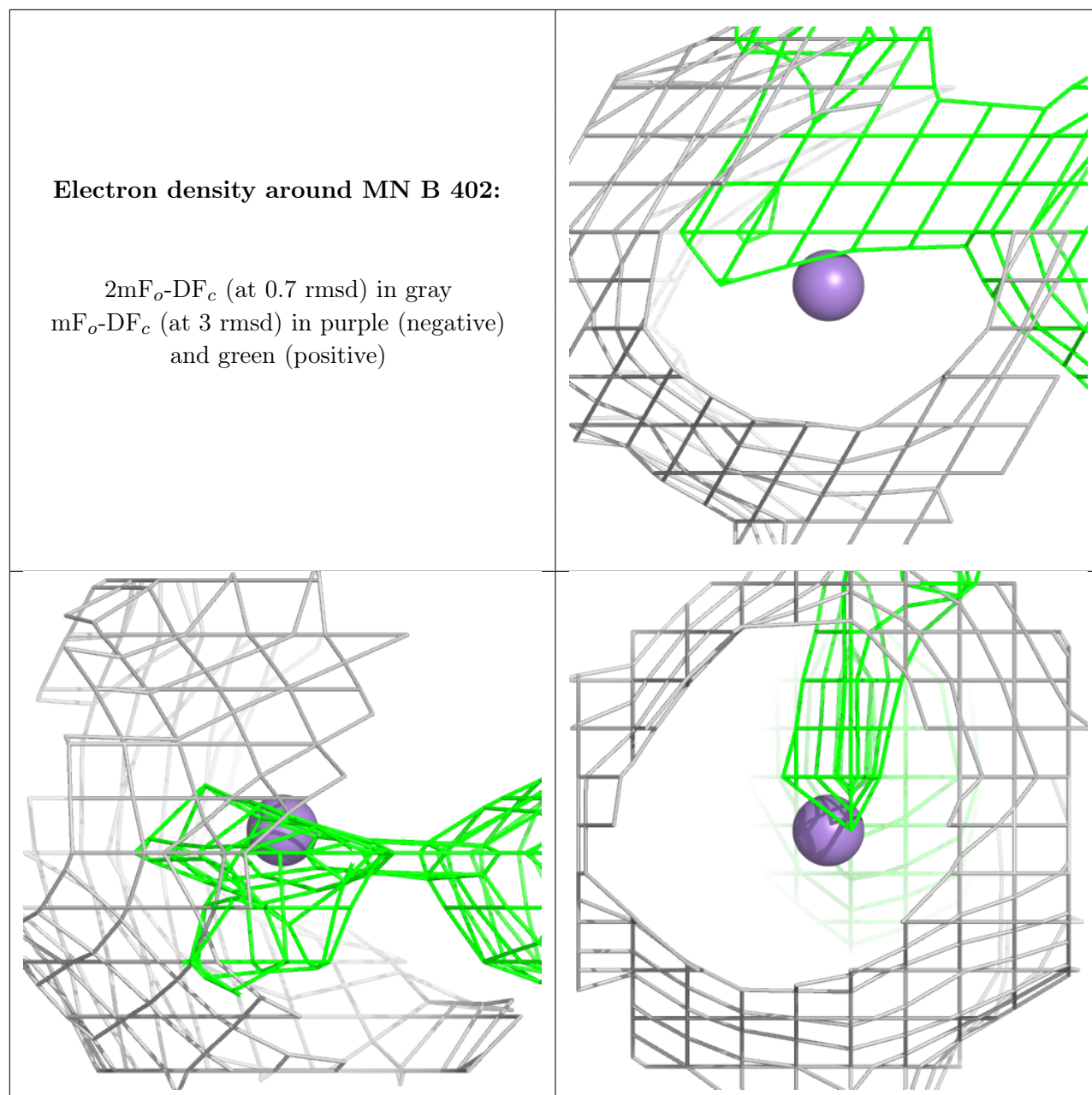
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around MN C 402:

$2mF_o-DF_c$ (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.