



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 19, 2026 – 06:45 AM UTC

PDB ID : 7N1R / pdb_00007n1r
Title : A novel and unique ATP hydrolysis to AMP by a human Hsp70 BiP
Authors : Yang, J.; Musayev, F.; Liu, Q.
Deposited on : 2021-05-28
Resolution : 2.03 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

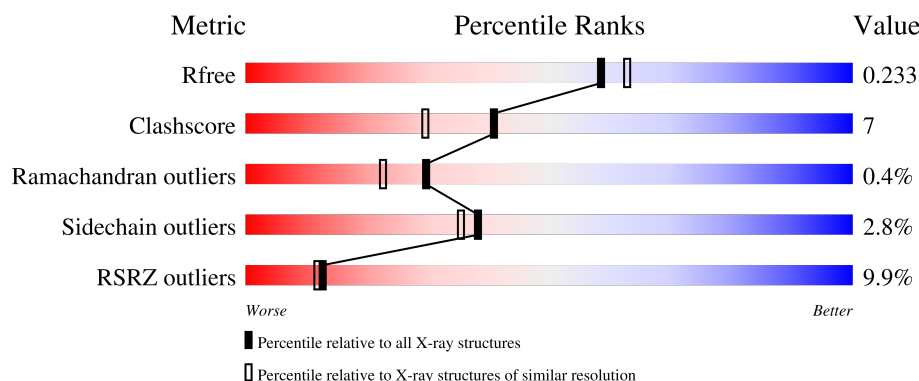
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.03 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	13299 (2.04-2.00)
Clashscore	190562	1022 (2.02-2.02)
Ramachandran outliers	187476	1014 (2.02-2.02)
Sidechain outliers	187428	1014 (2.02-2.02)
RSRZ outliers	180081	13314 (2.04-2.00)

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	606	10% 84% 14% .
1	B	606	10% 86% 12% .

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

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	PO4	A	709	-	-	X	-

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 10811 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 Endoplasmic reticulum chaperone BiP.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	606	Total	C	N	O	S	0	9	0
			4769	2994	812	951	12			
1	B	606	Total	C	N	O	S	0	9	0
			4771	2996	811	952	12			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	24	SER	-	expression tag	UNP P11021
A	453	VAL	THR	conflict	UNP P11021
A	454	GLY	ALA	conflict	UNP P11021
A	455	GLY	SER	conflict	UNP P11021
A	?	-	ASP	deletion	UNP P11021
A	?	-	ASN	deletion	UNP P11021
A	?	-	GLN	deletion	UNP P11021
A	?	-	PRO	deletion	UNP P11021
B	24	SER	-	expression tag	UNP P11021
B	453	VAL	THR	conflict	UNP P11021
B	454	GLY	ALA	conflict	UNP P11021
B	455	GLY	SER	conflict	UNP P11021
B	?	-	ASP	deletion	UNP P11021
B	?	-	ASN	deletion	UNP P11021
B	?	-	GLN	deletion	UNP P11021
B	?	-	PRO	deletion	UNP P11021

- Molecule 2 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



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

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

- Molecule 4 is PHOSPHATE ION (CCD ID: PO4) (formula: O_4P) (labeled as "Ligand of Interest" by depositor).

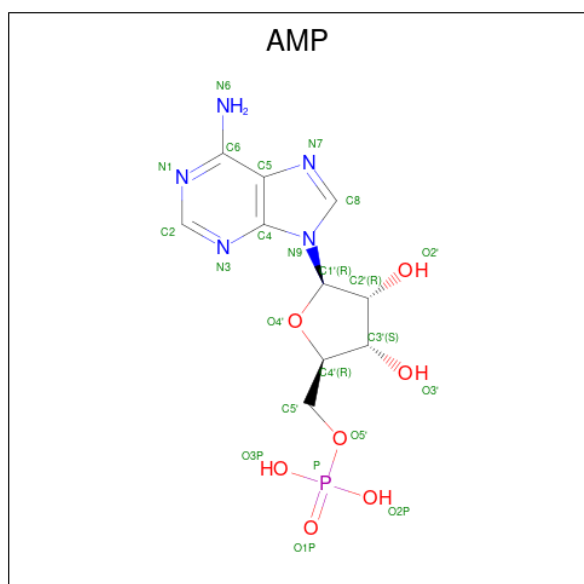


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	O	P	0	0
			5	4	1		
4	A	1	Total	O	P	0	0
			5	4	1		
4	A	1	Total	O	P	0	0
			5	4	1		
4	A	1	Total	O	P	0	0
			5	4	1		
4	A	1	Total	O	P	0	0
			5	4	1		
4	A	1	Total	O	P	0	0
			5	4	1		
4	A	1	Total	O	P	0	0
			5	4	1		
4	B	1	Total	O	P	0	0
			5	4	1		
4	B	1	Total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

- Molecule 5 is MAGNESIUM ION (CCD ID: MG) (formula: Mg) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total	Mg	0	0
			1	1		
5	B	1	Total	Mg	0	0
			1	1		

- Molecule 6 is ADENOSINE MONOPHOSPHATE (CCD ID: AMP) (formula: C₁₀H₁₄N₅O₇P) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
6	A	1	Total	C	N	O	P	0	0
			23	10	5	7	1		
6	B	1	Total	C	N	O	P	0	0
			23	10	5	7	1		

- Molecule 7 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: C₄H₁₀O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total	C	O	0	0
			7	4	3		

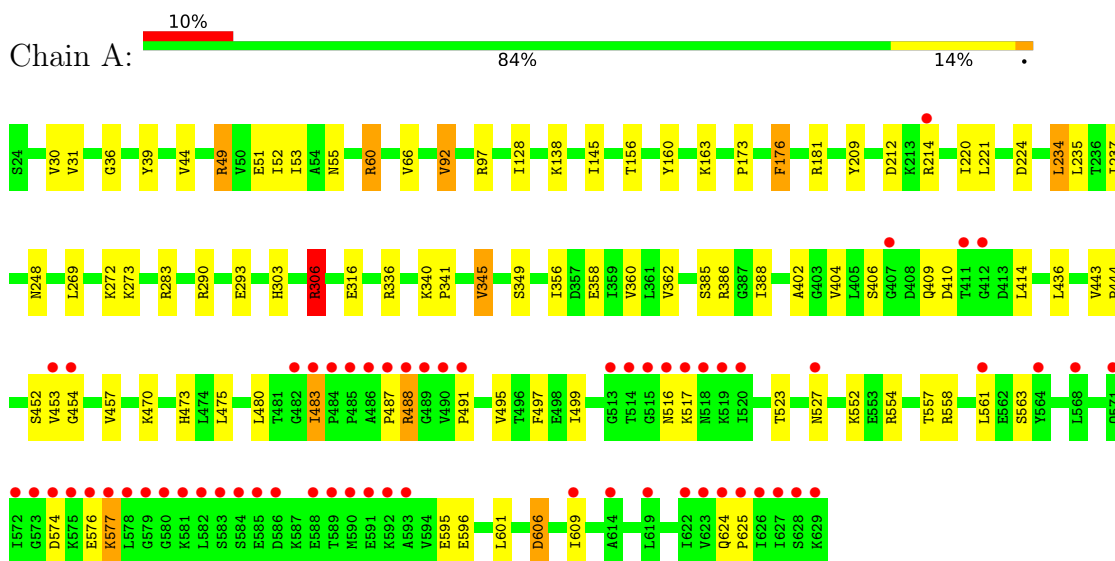
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	537	Total	O	0	0
			537	537		
8	B	531	Total	O	0	0
			531	531		

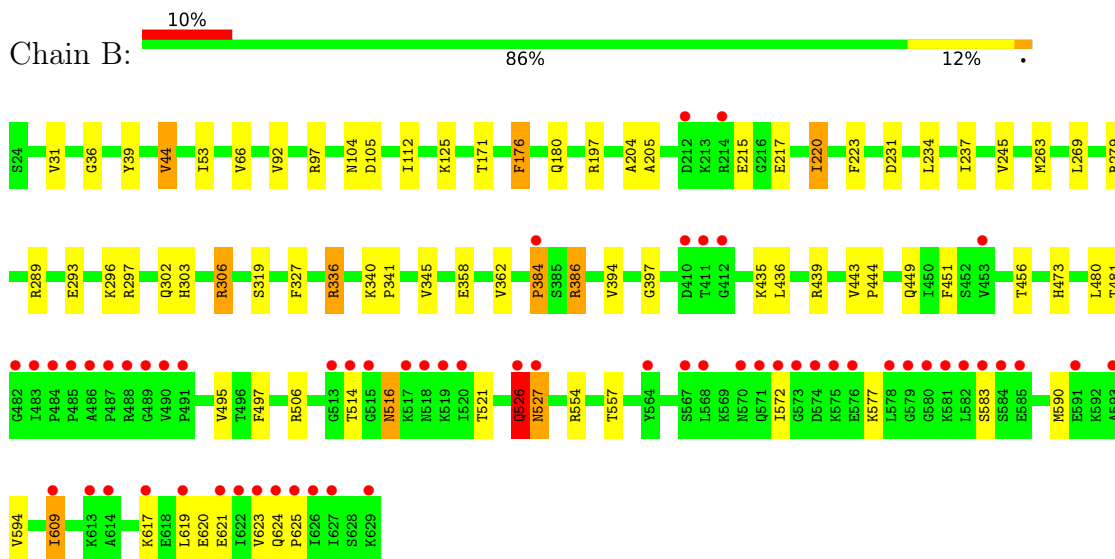
3 Residue-property plots [i](#)

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

• Molecule 1: Endoplasmic reticulum chaperone BiP



• Molecule 1: Endoplasmic reticulum chaperone BiP



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	75.99Å 76.39Å 79.92Å 84.88° 62.49° 63.14°	Depositor
Resolution (Å)	45.31 – 2.03 45.31 – 2.03	Depositor EDS
% Data completeness (in resolution range)	96.8 (45.31-2.03) 96.9 (45.31-2.03)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.51 (at 2.03Å)	Xtriage
Refinement program	REFMAC 5.8.0230	Depositor
R, R_{free}	0.174 , 0.223 0.185 , 0.233	Depositor DCC
R_{free} test set	4408 reflections (4.86%)	wwPDB-VP
Wilson B-factor (Å ²)	18.8	Xtriage
Anisotropy	0.006	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 46.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.011 for -h,-l,-k	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	10811	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.62% 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: PEG, SO4, AMP, GOL, MG, PO4

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	0.87	1/4839 (0.0%)	1.04	2/6533 (0.0%)
1	B	0.87	0/4844	1.04	4/6539 (0.1%)
All	All	0.87	1/9683 (0.0%)	1.04	6/13072 (0.0%)

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	A	0	6
1	B	0	4
All	All	0	10

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	92	VAL	C-O	-5.27	1.18	1.24

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	105	ASP	CA-CB-CG	6.16	118.76	112.60
1	A	306	ARG	CB-CA-C	5.47	118.67	109.48
1	A	181	ARG	CG-CD-NE	-5.31	100.31	112.00
1	B	223	PHE	CA-CB-CG	5.16	118.96	113.80
1	B	231	ASP	CA-CB-CG	5.15	117.75	112.60
1	B	384	PRO	CB-CA-C	-5.15	104.20	110.95

There are no chirality outliers.

All (10) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	212	ASP	Peptide
1	A	290	ARG	Sidechain
1	A	306	ARG	Sidechain
1	A	336	ARG	Sidechain
1	A	60	ARG	Sidechain
1	A	97	ARG	Sidechain
1	B	336	ARG	Sidechain
1	B	386	ARG	Sidechain
1	B	526	GLN	Peptide
1	B	527	ASN	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	4769	0	4796	61	0
1	B	4771	0	4797	62	0
2	A	24	0	32	2	0
2	B	24	0	32	5	0
3	A	10	0	0	0	0
3	B	20	0	0	2	0
4	A	40	0	0	4	0
4	B	30	0	0	0	0
5	A	1	0	0	0	0
5	B	1	0	0	0	0
6	A	23	0	12	0	0
6	B	23	0	12	0	0
7	A	7	0	10	1	0
8	A	537	0	0	17	1
8	B	531	0	0	15	1
All	All	10811	0	9691	127	1

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 7.

All (127) 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:410:ASP:HA	8:A:1070:HOH:O	1.56	1.05
4:A:709:PO4:O3	8:A:801:HOH:O	1.81	0.96
1:A:163:LYS:NZ	8:A:802:HOH:O	2.09	0.86
4:A:709:PO4:P	8:A:801:HOH:O	2.42	0.77
1:B:31:VAL:HG12	1:B:44[B]:VAL:HG22	1.69	0.74
1:A:306:ARG:HG2	1:A:306:ARG:HH11	1.54	0.72
1:B:220:ILE:HD12	1:B:358:GLU:HB2	1.72	0.72
1:A:248:ASN:C	8:A:816:HOH:O	2.32	0.71
1:B:439:ARG:HH22	2:B:704:GOL:H31	1.56	0.70
1:A:409:GLN:O	8:A:804:HOH:O	2.10	0.69
1:B:336:ARG:NH1	8:B:802:HOH:O	2.05	0.68
1:B:31:VAL:HG12	1:B:44[A]:VAL:HG13	1.76	0.66
1:B:514:THR:HG21	1:B:516:ASN:HD22	1.61	0.66
1:A:473:HIS:HE1	8:A:1243:HOH:O	1.78	0.65
1:A:557:THR:HG21	1:A:609:ILE:HD11	1.79	0.64
1:A:220:ILE:HD12	1:A:358:GLU:HB2	1.80	0.64
1:A:410:ASP:N	8:A:814:HOH:O	2.29	0.63
4:A:709:PO4:O2	8:A:805:HOH:O	2.16	0.61
1:B:384:PRO:HD3	8:B:885:HOH:O	2.00	0.60
1:B:557:THR:HG21	1:B:609:ILE:HD11	1.83	0.60
1:A:360:VAL:HG13	1:A:388:ILE:HD13	1.83	0.59
1:A:574:ASP:OD2	1:A:577:LYS:CB	2.50	0.59
1:A:234:LEU:HB2	1:A:345[A]:VAL:CG1	2.33	0.58
1:B:104[A]:ASN:ND2	8:B:815:HOH:O	2.36	0.58
1:A:283:ARG:NH1	8:A:823:HOH:O	2.36	0.57
1:A:454:GLY:HA2	1:A:483:ILE:HD13	1.84	0.57
1:B:269:LEU:HD23	1:B:269:LEU:O	2.05	0.57
1:A:293:GLU:OE2	8:A:806:HOH:O	2.17	0.57
1:B:234[A]:LEU:C	1:B:234[A]:LEU:HD13	2.30	0.57
1:B:234[A]:LEU:HD12	1:B:245:VAL:HB	1.86	0.57
1:B:340:LYS:HB2	1:B:341:PRO:HD3	1.86	0.57
1:A:49:ARG:NH2	7:A:717:PEG:O2	2.39	0.56
1:B:171:THR:HG21	1:B:394:VAL:HG12	1.85	0.56
1:B:44[A]:VAL:HG22	1:B:53:ILE:HD11	1.88	0.56
1:A:340:LYS:HB2	1:A:341:PRO:HD3	1.88	0.56
1:A:31:VAL:HG12	1:A:44[A]:VAL:CG1	2.36	0.56
1:B:340:LYS:NZ	8:B:821:HOH:O	2.39	0.55
1:A:487:PRO:O	1:A:488:ARG:CB	2.54	0.55
1:B:473:HIS:HE1	8:B:1223:HOH:O	1.88	0.55
1:B:44[B]:VAL:HG23	1:B:53:ILE:HD11	1.88	0.55
1:B:97:ARG:NH2	8:B:812:HOH:O	2.39	0.55
1:B:303:HIS:HD2	3:B:708:SO4:O3	1.90	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:234[A]:LEU:HB2	1:B:345[A]:VAL:CG1	2.38	0.54
1:B:205:ALA:HB1	1:B:362:VAL:HG11	1.90	0.54
4:A:707:PO4:O1	8:A:807:HOH:O	2.19	0.53
2:B:705:GOL:H12	8:B:1232:HOH:O	2.08	0.53
1:A:234:LEU:HD21	1:A:356:ILE:CD1	2.38	0.53
1:A:44[B]:VAL:HG23	1:A:53:ILE:HD11	1.91	0.53
1:B:220:ILE:HG12	1:B:237:ILE:HD12	1.90	0.52
1:B:197:ARG:NH1	8:B:804:HOH:O	2.22	0.52
1:A:44[B]:VAL:CG2	1:A:53:ILE:HD11	2.40	0.52
1:B:554:ARG:HA	1:B:609:ILE:HD13	1.93	0.51
1:A:606:ASP:OD1	1:A:606:ASP:C	2.54	0.50
1:B:269:LEU:HD23	1:B:269:LEU:C	2.37	0.49
1:A:303:HIS:HD2	3:B:702:SO4:O4	1.96	0.49
1:B:279:ARG:NH1	8:B:823:HOH:O	2.41	0.49
1:A:214:ARG:HD2	1:B:215:GLU:HB3	1.93	0.49
1:B:204:ALA:O	1:B:397:GLY:HA3	2.12	0.49
1:B:31:VAL:HG12	1:B:44[A]:VAL:CG1	2.41	0.48
1:A:66:VAL:HG22	1:A:92:VAL:HB	1.95	0.48
1:A:552:LYS:NZ	8:A:834:HOH:O	2.44	0.48
1:A:36:GLY:HA3	1:A:39:TYR:O	2.14	0.48
2:B:705:GOL:H11	8:B:1104:HOH:O	2.13	0.47
1:A:173:PRO:HD2	1:A:176:PHE:CD2	2.49	0.47
1:A:552:LYS:NZ	8:A:844:HOH:O	2.47	0.47
1:B:293:GLU:OE2	1:B:297:ARG:NH2	2.48	0.47
1:B:217[A]:GLU:OE1	1:B:217[A]:GLU:C	2.58	0.47
1:A:316:GLU:HG3	8:A:1311:HOH:O	2.15	0.47
1:A:52:ILE:O	1:A:60:ARG:NH1	2.45	0.47
1:A:558:ARG:HH12	2:A:703:GOL:H32	1.80	0.47
1:A:30:VAL:HG12	1:A:402:ALA:HB1	1.97	0.46
1:A:480:LEU:HD11	1:A:495:VAL:HG21	1.98	0.46
1:A:234:LEU:C	1:A:234:LEU:HD13	2.41	0.46
1:A:269:LEU:HD23	1:A:269:LEU:C	2.40	0.46
1:A:209:TYR:O	1:B:386:ARG:NH1	2.48	0.46
1:B:220:ILE:CG1	1:B:237:ILE:HD12	2.45	0.46
1:A:452:SER:HA	1:A:491:PRO:O	2.15	0.46
1:A:443:VAL:HA	1:A:444:PRO:C	2.40	0.45
1:B:436:LEU:HD21	1:B:497:PHE:CD1	2.51	0.45
1:B:624:GLN:HB3	1:B:625:PRO:HD3	1.98	0.45
1:A:386:ARG:HD3	2:A:702:GOL:O3	2.16	0.45
1:B:583:SER:OG	8:B:803:HOH:O	2.20	0.45
1:B:435:LYS:HZ3	2:B:705:GOL:H11	1.82	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:306:ARG:NH1	1:B:319:SER:OG	2.50	0.45
1:A:475:LEU:C	1:A:475:LEU:HD23	2.41	0.44
1:B:279:ARG:NH2	8:B:823:HOH:O	2.42	0.44
1:B:443:VAL:HA	1:B:444:PRO:C	2.43	0.44
1:B:557:THR:OG1	1:B:609:ILE:CD1	2.65	0.44
1:A:453:VAL:CG1	1:A:457:VAL:HG22	2.47	0.44
1:B:449:GLN:HG2	1:B:451:PHE:CZ	2.53	0.44
1:A:345[B]:VAL:O	1:A:349[B]:SER:HB3	2.18	0.44
1:B:44[B]:VAL:CG2	1:B:53:ILE:HD11	2.47	0.43
1:B:125:LYS:CE	1:B:527:ASN:HB3	2.48	0.43
1:B:384:PRO:CD	8:B:885:HOH:O	2.63	0.43
1:B:480:LEU:HD11	1:B:495:VAL:HG21	1.99	0.43
1:A:269:LEU:HD21	1:A:273:LYS:HE3	1.99	0.43
1:B:36:GLY:HA3	1:B:39:TYR:O	2.17	0.43
1:B:296:LYS:HG3	1:B:327:PHE:CZ	2.53	0.43
1:B:572:ILE:HD11	1:B:590:MET:HB3	1.99	0.43
1:A:220:ILE:HG13	1:A:237:ILE:HD12	1.99	0.43
1:A:561:LEU:HG	1:A:601:LEU:HD11	2.01	0.43
1:A:409:GLN:HA	8:A:814:HOH:O	2.19	0.43
1:B:176:PHE:CD2	1:B:180:GLN:HB3	2.53	0.43
1:A:234:LEU:HD13	1:A:235:LEU:N	2.34	0.43
1:A:269:LEU:HD23	1:A:269:LEU:O	2.19	0.42
1:B:439:ARG:NH2	2:B:704:GOL:H31	2.30	0.42
1:B:66:VAL:HG22	1:B:92:VAL:HB	2.02	0.42
1:B:171:THR:HG21	1:B:394:VAL:CG1	2.49	0.42
1:B:514:THR:CG2	1:B:516:ASN:HD22	2.29	0.42
1:B:97:ARG:CZ	8:B:812:HOH:O	2.68	0.42
1:A:128:ILE:HD13	1:A:145:ILE:HG13	2.02	0.42
1:A:55:ASN:HA	1:A:156:THR:HG23	2.02	0.42
1:B:617:LYS:HA	1:B:620:GLU:HB2	2.02	0.42
1:A:444:PRO:HA	1:A:499:ILE:O	2.20	0.42
1:A:624:GLN:HB3	1:A:625:PRO:HD3	2.02	0.42
1:A:51:GLU:HG2	1:A:160:TYR:CZ	2.55	0.41
1:B:263[A]:MET:SD	1:B:289:ARG:HA	2.60	0.41
1:A:138:LYS:NZ	8:A:871:HOH:O	2.53	0.41
1:A:224:ASP:HA	1:A:362:VAL:O	2.21	0.41
1:B:619:LEU:O	1:B:623:VAL:HG23	2.21	0.41
1:A:66:VAL:CG2	1:A:92:VAL:HB	2.50	0.41
1:A:436:LEU:HD21	1:A:497:PHE:CD1	2.56	0.41
1:B:302:GLN:NE2	8:B:858:HOH:O	2.54	0.41
1:B:506:ARG:NH1	1:B:521:THR:OG1	2.53	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:554:ARG:HH12	1:A:606:ASP:HA	1.86	0.40
1:A:404:VAL:HG21	1:A:414:LEU:HD23	2.04	0.40
1:B:112:ILE:HD13	1:B:112:ILE:HG21	1.88	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:A:830:HOH:O	8:B:820:HOH:O[1_565]	1.95	0.25

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	613/606 (101%)	587 (96%)	23 (4%)	3 (0%)	24	17
1	B	613/606 (101%)	594 (97%)	17 (3%)	2 (0%)	36	31
All	All	1226/1212 (101%)	1181 (96%)	40 (3%)	5 (0%)	30	22

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	488	ARG
1	A	527	ASN
1	A	577	LYS
1	B	526	GLN
1	B	577	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	525/519 (101%)	505 (96%)	20 (4%)	29	23
1	B	525/519 (101%)	513 (98%)	12 (2%)	44	43
All	All	1050/1038 (101%)	1018 (97%)	32 (3%)	38	32

All (32) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	49	ARG
1	A	176	PHE
1	A	221	LEU
1	A	234	LEU
1	A	272	LYS
1	A	345[A]	VAL
1	A	345[B]	VAL
1	A	385	SER
1	A	406	SER
1	A	470	LYS
1	A	483	ILE
1	A	516	ASN
1	A	517	LYS
1	A	523[A]	THR
1	A	523[B]	THR
1	A	563	SER
1	A	576	GLU
1	A	595	GLU
1	A	596	GLU
1	A	606	ASP
1	B	44[A]	VAL
1	B	44[B]	VAL
1	B	176	PHE
1	B	220	ILE
1	B	306	ARG
1	B	456	THR
1	B	481	THR
1	B	516	ASN
1	B	526	GLN
1	B	594	VAL
1	B	609	ILE

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Mol	Chain	Res	Type
1	B	621	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	110	GLN
1	A	302	GLN
1	A	303	HIS
1	A	380	ASN
1	A	473	HIS
1	A	527	ASN
1	A	571	GLN
1	B	182	GLN
1	B	239	ASN
1	B	302	GLN
1	B	303	HIS
1	B	304	GLN
1	B	343	GLN
1	B	516	ASN
1	B	624	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 33 ligands modelled in this entry, 2 are monoatomic - leaving 31 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	SO4	B	708	-	4,4,4	0.46	0	6,6,6	0.62	0
3	SO4	A	706	-	4,4,4	0.53	0	6,6,6	0.56	0
2	GOL	B	704	-	5,5,5	0.43	0	5,5,5	0.73	0
4	PO4	B	710	-	4,4,4	0.78	0	6,6,6	0.77	0
4	PO4	A	712	-	4,4,4	1.26	1 (25%)	6,6,6	0.68	0
4	PO4	A	710	-	4,4,4	0.89	0	6,6,6	0.33	0
4	PO4	A	711	-	4,4,4	1.02	0	6,6,6	0.68	0
2	GOL	A	701	-	5,5,5	0.40	0	5,5,5	0.66	0
4	PO4	B	713	-	4,4,4	0.93	0	6,6,6	0.64	0
3	SO4	B	702	-	4,4,4	0.57	0	6,6,6	0.87	0
6	AMP	A	715	5	25,25,25	1.53	4 (16%)	37,38,38	1.85	10 (27%)
3	SO4	B	709	-	4,4,4	0.45	0	6,6,6	0.56	0
2	GOL	B	701	-	5,5,5	0.44	0	5,5,5	0.94	0
3	SO4	A	705	-	4,4,4	0.56	0	6,6,6	0.34	0
2	GOL	A	704	-	5,5,5	0.59	0	5,5,5	0.34	0
4	PO4	B	712	-	4,4,4	0.71	0	6,6,6	0.62	0
4	PO4	B	716	5	4,4,4	2.16	2 (50%)	6,6,6	0.74	0
4	PO4	A	716	5	4,4,4	2.41	2 (50%)	6,6,6	1.16	0
3	SO4	B	707	-	4,4,4	0.55	0	6,6,6	0.14	0
4	PO4	B	711	-	4,4,4	0.50	0	6,6,6	0.62	0
4	PO4	A	707	-	4,4,4	0.70	0	6,6,6	0.61	0
4	PO4	A	709	-	4,4,4	0.88	0	6,6,6	0.77	0
4	PO4	B	714	-	4,4,4	0.84	0	6,6,6	0.62	0
7	PEG	A	717	-	6,6,6	0.54	0	5,5,5	0.57	0
2	GOL	A	702	-	5,5,5	0.46	0	5,5,5	0.89	0
2	GOL	A	703	-	5,5,5	0.52	0	5,5,5	0.35	0
4	PO4	A	708	-	4,4,4	0.83	0	6,6,6	1.02	0
6	AMP	B	715	5	25,25,25	1.46	4 (16%)	37,38,38	2.57	15 (40%)
2	GOL	B	706	-	5,5,5	0.55	0	5,5,5	0.40	0
4	PO4	A	713	-	4,4,4	0.75	0	6,6,6	0.88	0
2	GOL	B	705	-	5,5,5	0.68	0	5,5,5	0.84	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	GOL	A	701	-	-	4/4/4/4	-
2	GOL	B	701	-	-	2/4/4/4	-
2	GOL	B	704	-	-	4/4/4/4	-
2	GOL	A	704	-	-	4/4/4/4	-
7	PEG	A	717	-	-	2/4/4/4	-
2	GOL	A	702	-	-	2/4/4/4	-
2	GOL	A	703	-	-	2/4/4/4	-
6	AMP	A	715	5	-	0/10/26/26	0/3/3/3
6	AMP	B	715	5	-	0/10/26/26	0/3/3/3
2	GOL	B	706	-	-	4/4/4/4	-
2	GOL	B	705	-	-	2/4/4/4	-

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	A	715	AMP	C5-C4	4.43	1.47	1.39
6	B	715	AMP	C5-C4	3.93	1.46	1.39
4	B	716	PO4	P-O3	-3.22	1.45	1.54
6	A	715	AMP	C2-N1	3.02	1.39	1.33
4	A	716	PO4	P-O3	-3.01	1.45	1.54
4	A	716	PO4	P-O2	-2.97	1.46	1.54
6	B	715	AMP	C8-N7	2.97	1.37	1.31
6	B	715	AMP	C2-N1	2.72	1.38	1.33
6	B	715	AMP	C5-N7	-2.52	1.34	1.39
6	A	715	AMP	C2-N3	2.48	1.38	1.33
4	B	716	PO4	P-O2	-2.29	1.48	1.54
6	A	715	AMP	C4-N9	-2.28	1.33	1.37
4	A	712	PO4	P-O1	2.02	1.55	1.50

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	B	715	AMP	C4-N9-C8	5.62	111.64	105.74
6	B	715	AMP	N3-C4-N9	5.38	136.32	127.17
6	B	715	AMP	C5-C4-N3	-4.99	119.84	126.72
6	A	715	AMP	C4-N9-C8	4.99	110.98	105.74
6	B	715	AMP	O3P-P-O2P	4.90	126.17	107.80
6	B	715	AMP	N9-C8-N7	-4.62	107.38	113.94
6	A	715	AMP	N3-C4-N9	4.42	134.68	127.17
6	B	715	AMP	N3-C2-N1	-3.97	122.58	128.58
6	A	715	AMP	C5-C4-N3	-3.83	121.44	126.72

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	B	715	AMP	C5-C6-N6	-3.82	113.82	123.29
6	B	715	AMP	C2-N3-C4	3.61	120.65	111.83
6	B	715	AMP	C5-N7-C8	3.49	108.93	103.45
6	B	715	AMP	N6-C6-N1	3.30	125.73	118.38
6	A	715	AMP	N3-C2-N1	-3.15	123.81	128.58
6	A	715	AMP	N9-C8-N7	-2.61	110.24	113.94
6	B	715	AMP	C4-N9-C1'	-2.59	120.57	126.63
6	B	715	AMP	O2'-C2'-C3'	2.52	119.91	111.82
6	A	715	AMP	C2-N3-C4	2.44	117.80	111.83
6	A	715	AMP	C5-C6-N6	-2.42	117.29	123.29
6	B	715	AMP	O2P-P-O5'	-2.19	100.95	106.67
6	A	715	AMP	O3P-P-O1P	2.16	119.24	110.83
6	B	715	AMP	O3P-P-O5'	-2.13	101.10	106.67
6	A	715	AMP	C2'-C1'-N9	-2.10	108.10	113.30
6	B	715	AMP	C4-C5-N7	-2.07	108.21	110.58
6	A	715	AMP	N6-C6-N1	2.03	122.90	118.38

There are no chirality outliers.

All (26) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	701	GOL	C1-C2-C3-O3
2	A	704	GOL	O1-C1-C2-C3
2	A	704	GOL	C1-C2-C3-O3
2	B	704	GOL	O1-C1-C2-C3
2	B	705	GOL	C1-C2-C3-O3
2	B	706	GOL	O1-C1-C2-C3
2	B	706	GOL	C1-C2-C3-O3
2	A	704	GOL	O2-C2-C3-O3
2	B	704	GOL	O1-C1-C2-O2
2	B	705	GOL	O2-C2-C3-O3
2	B	706	GOL	O2-C2-C3-O3
2	A	701	GOL	O1-C1-C2-C3
2	A	702	GOL	C1-C2-C3-O3
2	B	701	GOL	O1-C1-C2-C3
2	B	704	GOL	C1-C2-C3-O3
2	A	701	GOL	O2-C2-C3-O3
2	A	702	GOL	O2-C2-C3-O3
2	A	704	GOL	O1-C1-C2-O2
2	B	706	GOL	O1-C1-C2-O2
7	A	717	PEG	O1-C1-C2-O2
2	B	701	GOL	O1-C1-C2-O2

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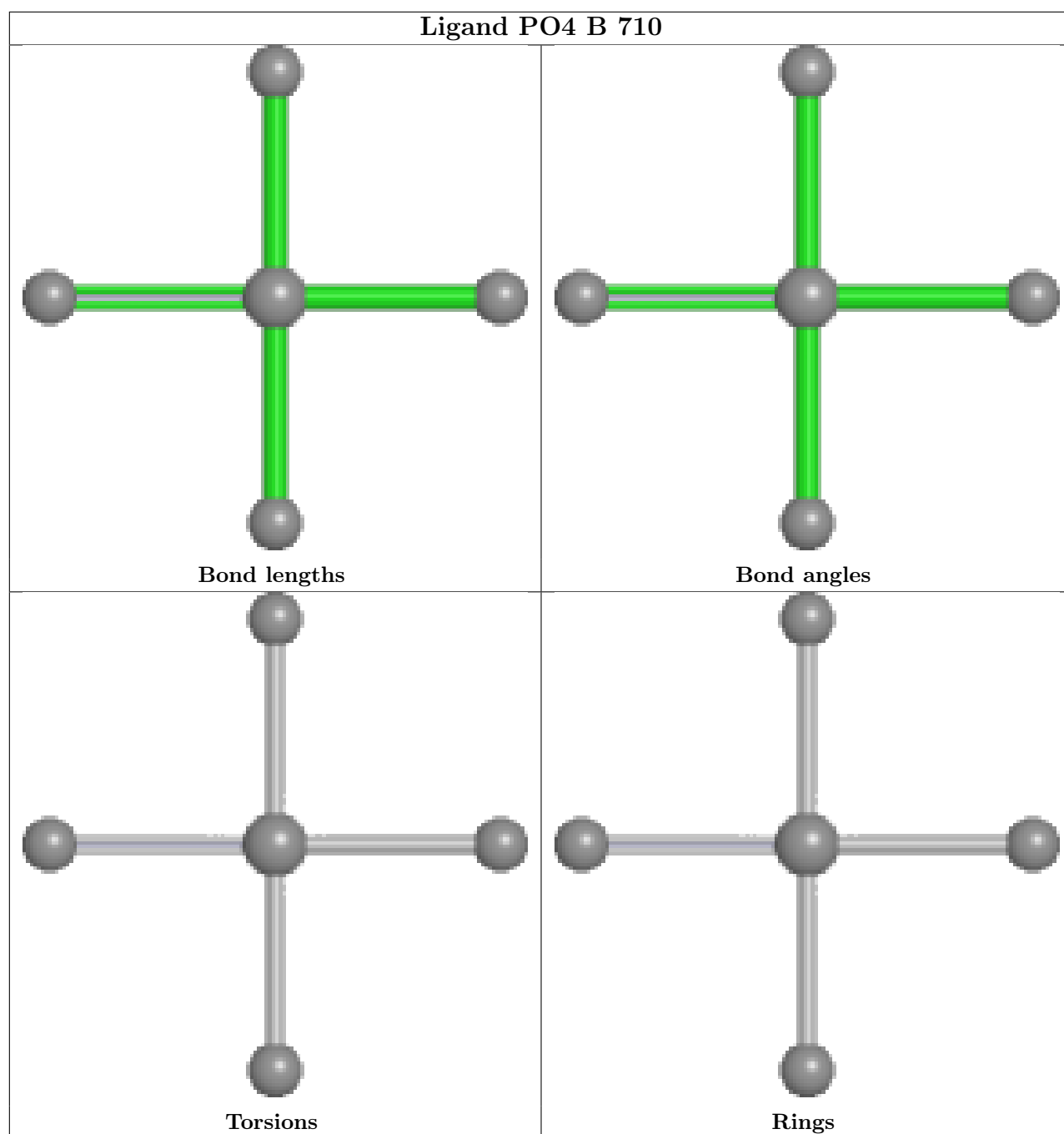
Mol	Chain	Res	Type	Atoms
2	A	701	GOL	O1-C1-C2-O2
7	A	717	PEG	O2-C3-C4-O4
2	B	704	GOL	O2-C2-C3-O3
2	A	703	GOL	C1-C2-C3-O3
2	A	703	GOL	O1-C1-C2-C3

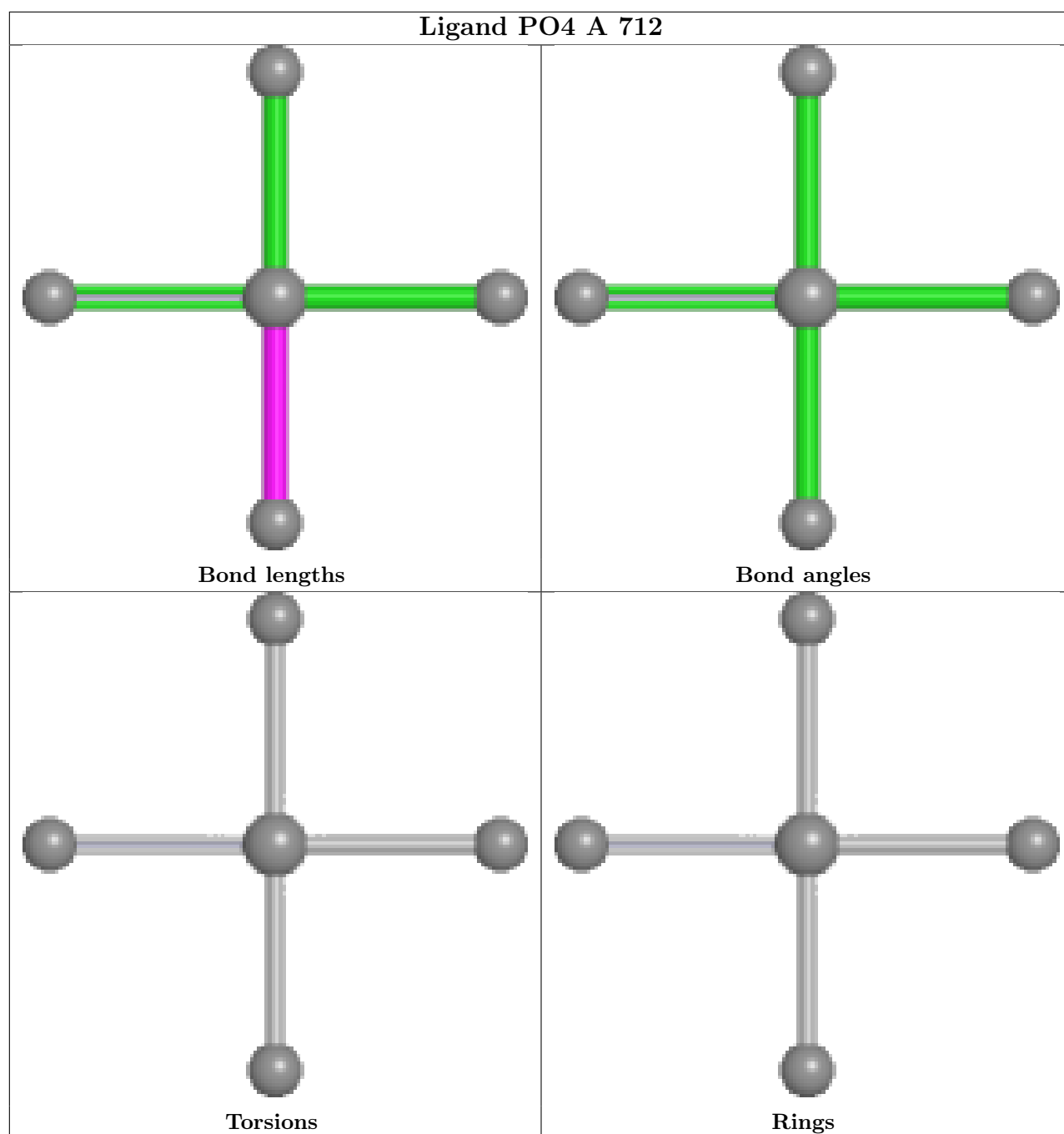
There are no ring outliers.

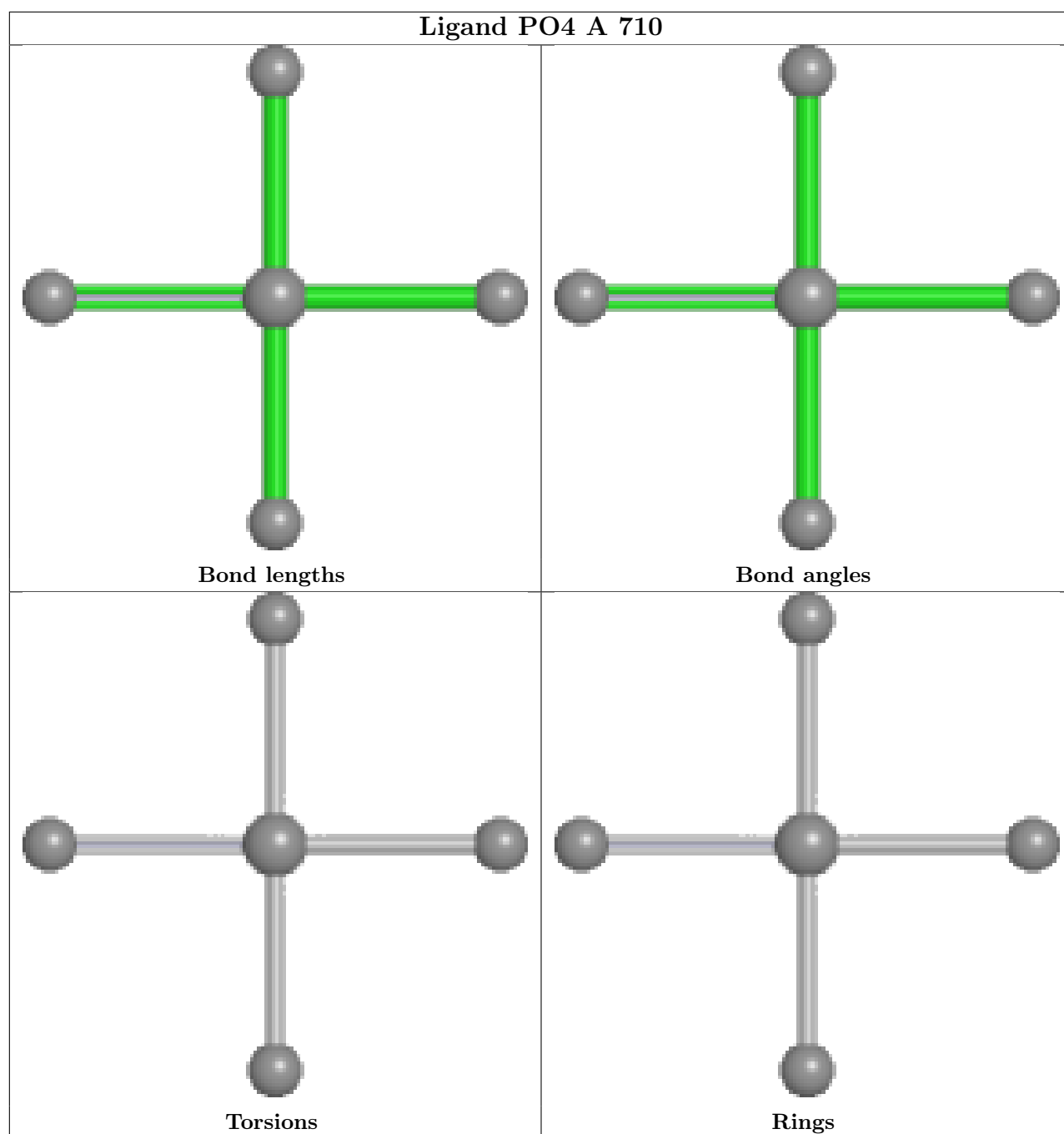
9 monomers are involved in 14 short contacts:

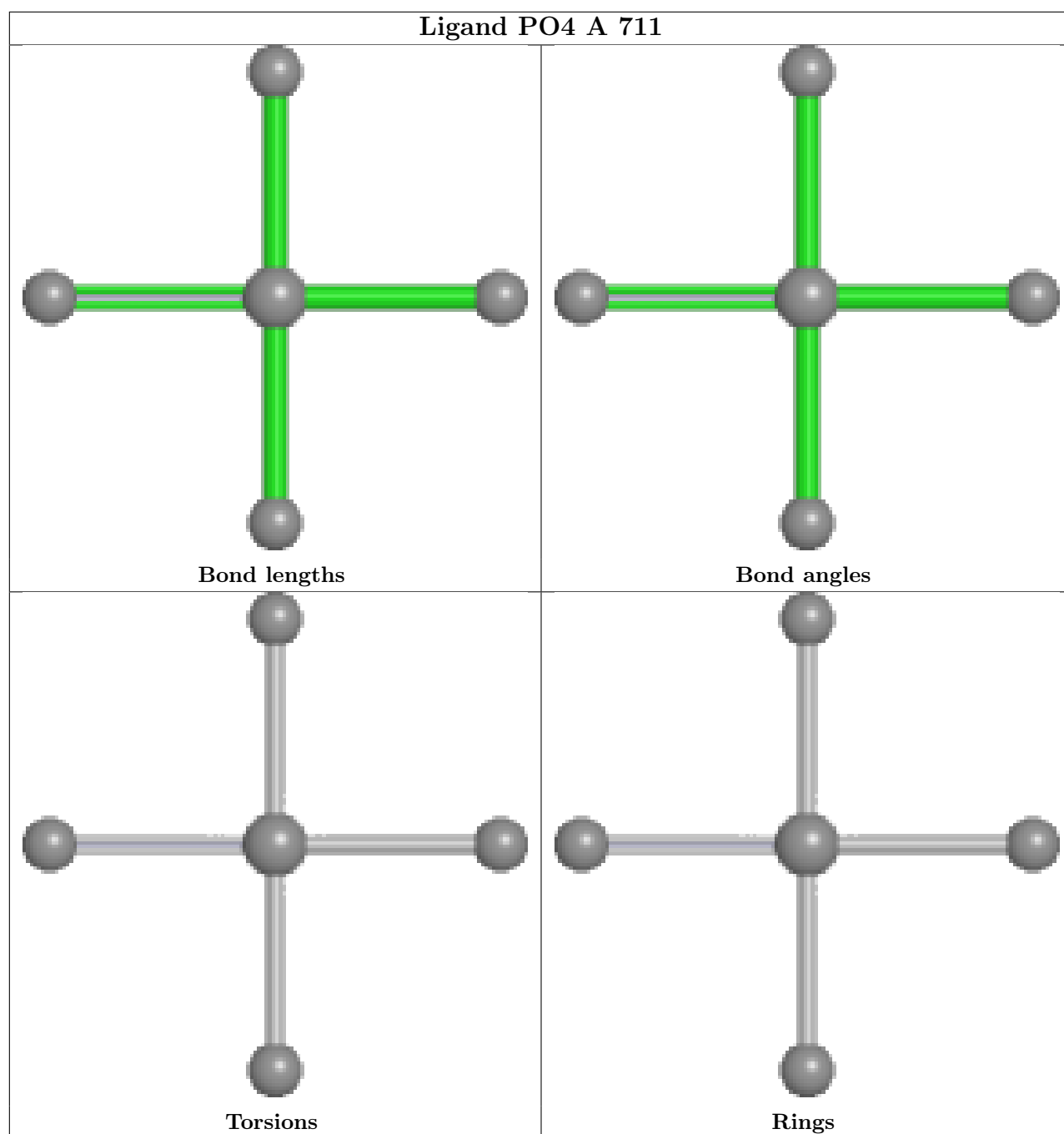
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	708	SO4	1	0
2	B	704	GOL	2	0
3	B	702	SO4	1	0
4	A	707	PO4	1	0
4	A	709	PO4	3	0
7	A	717	PEG	1	0
2	A	702	GOL	1	0
2	A	703	GOL	1	0
2	B	705	GOL	3	0

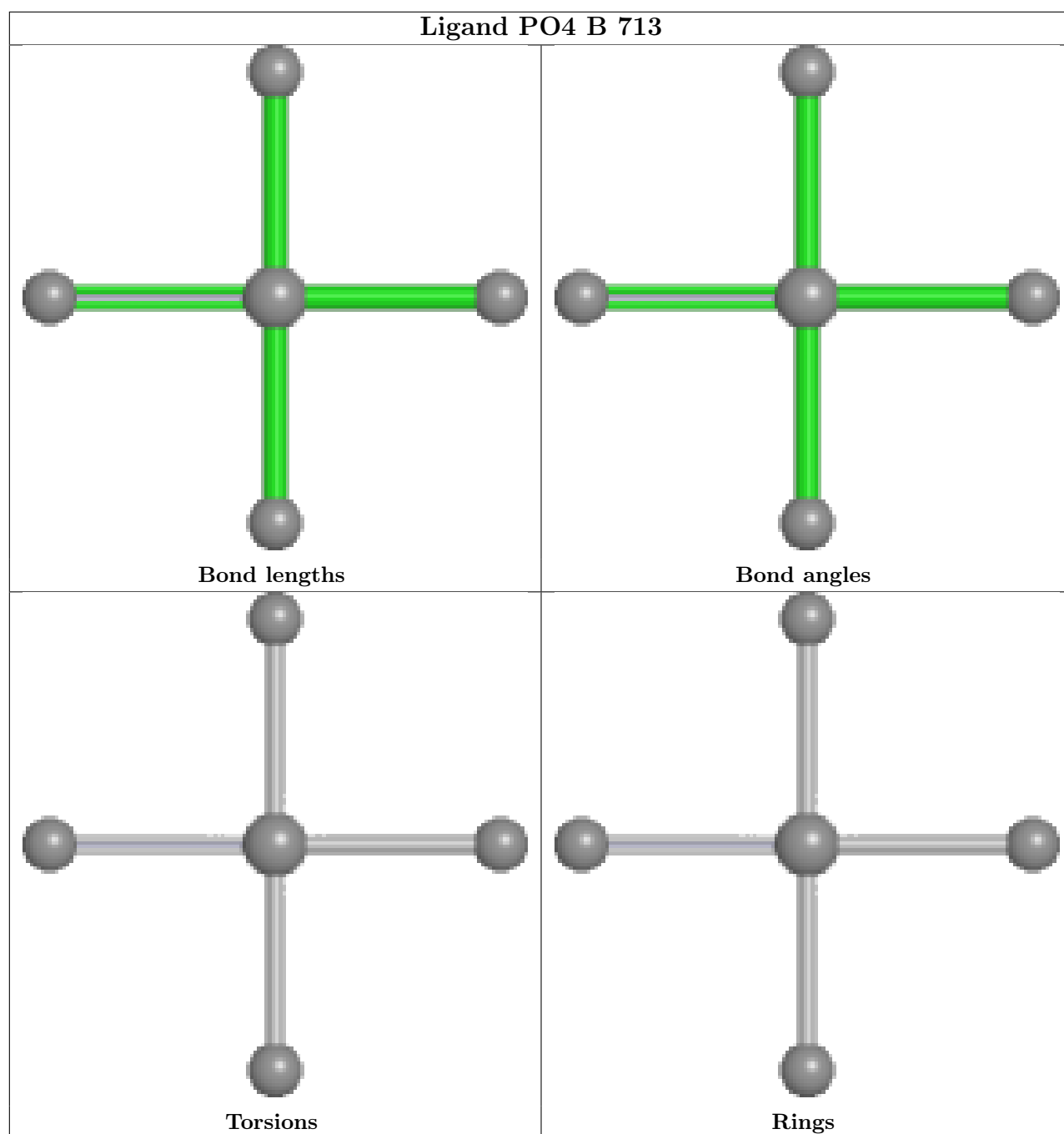
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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.

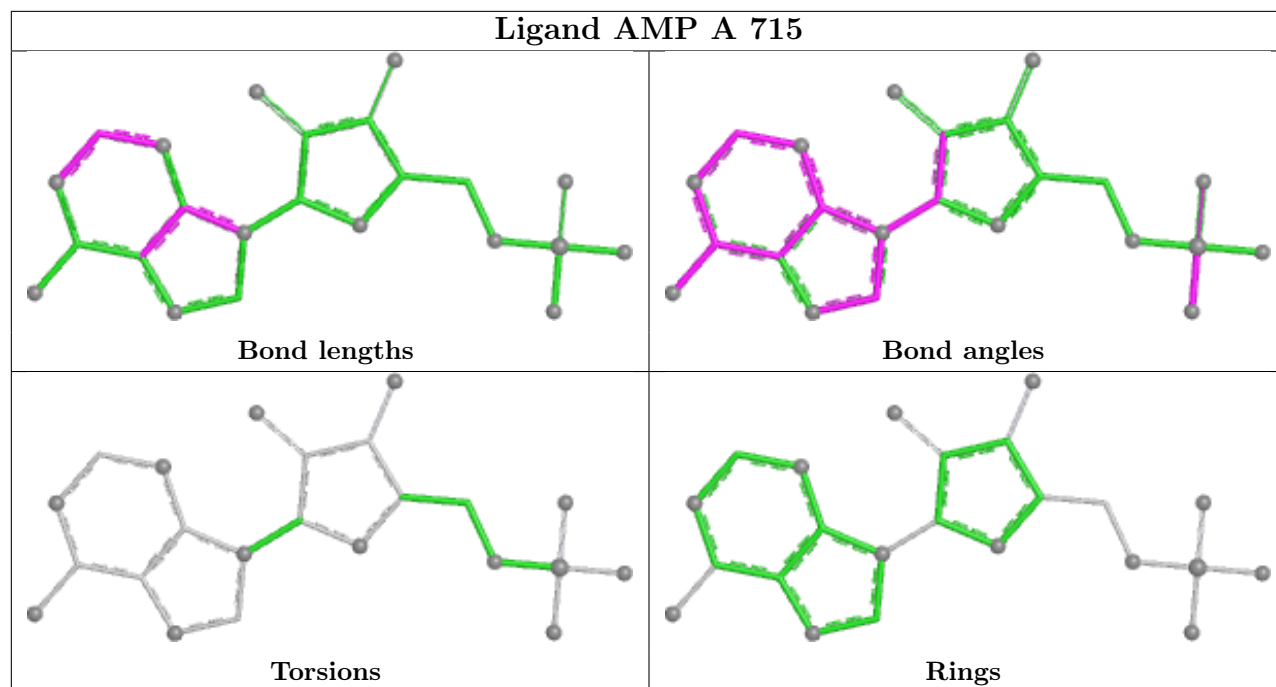


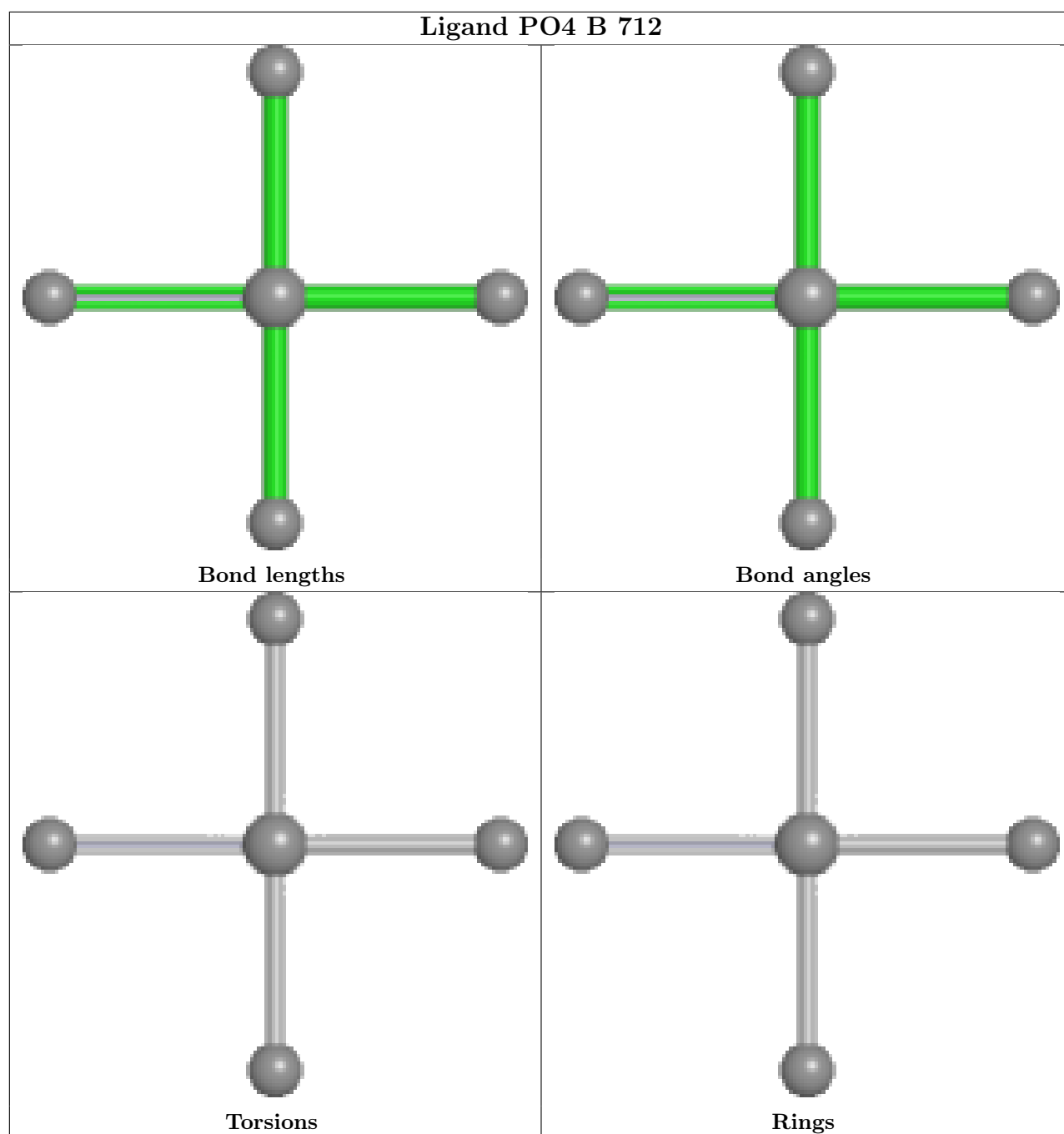


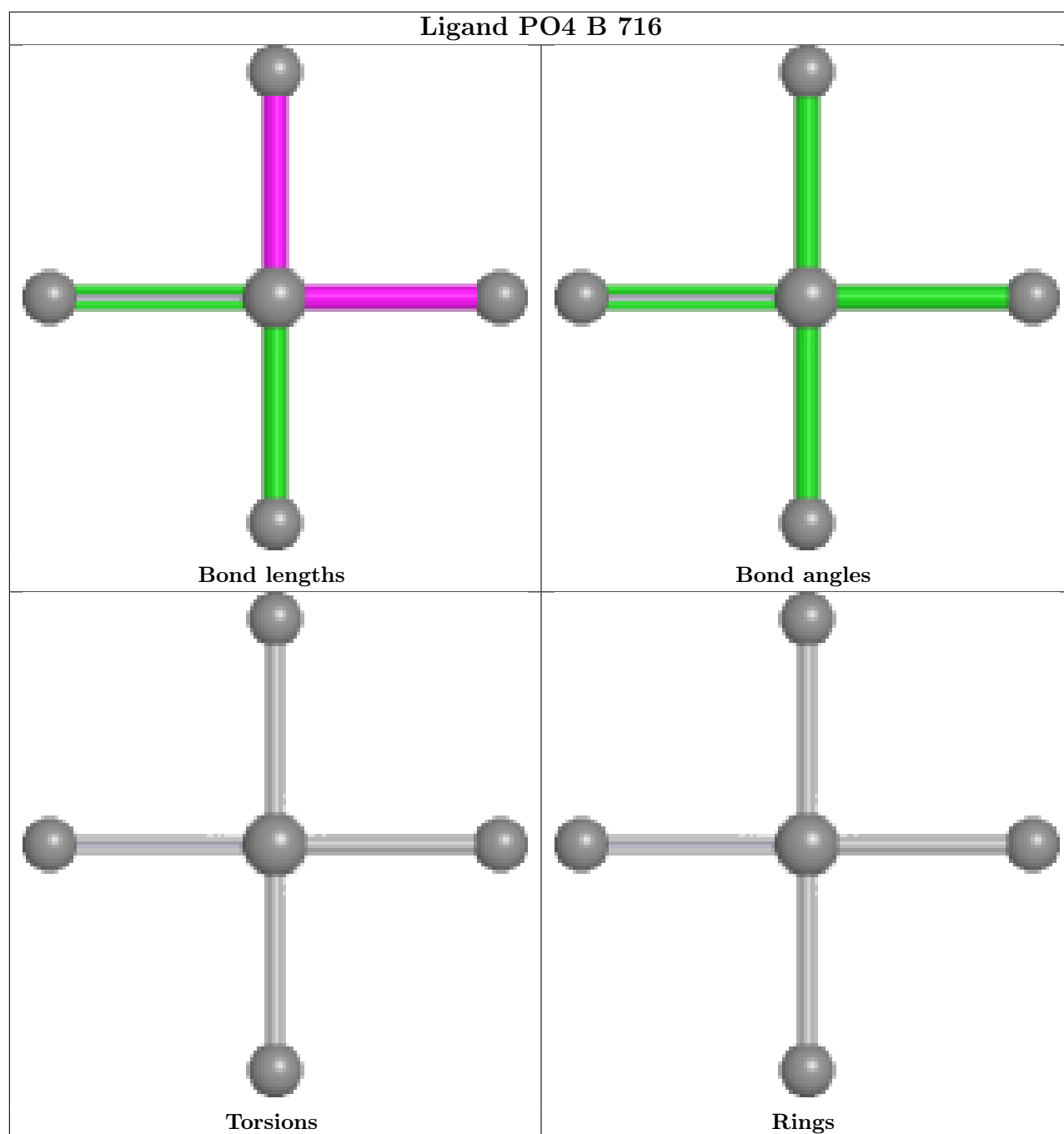


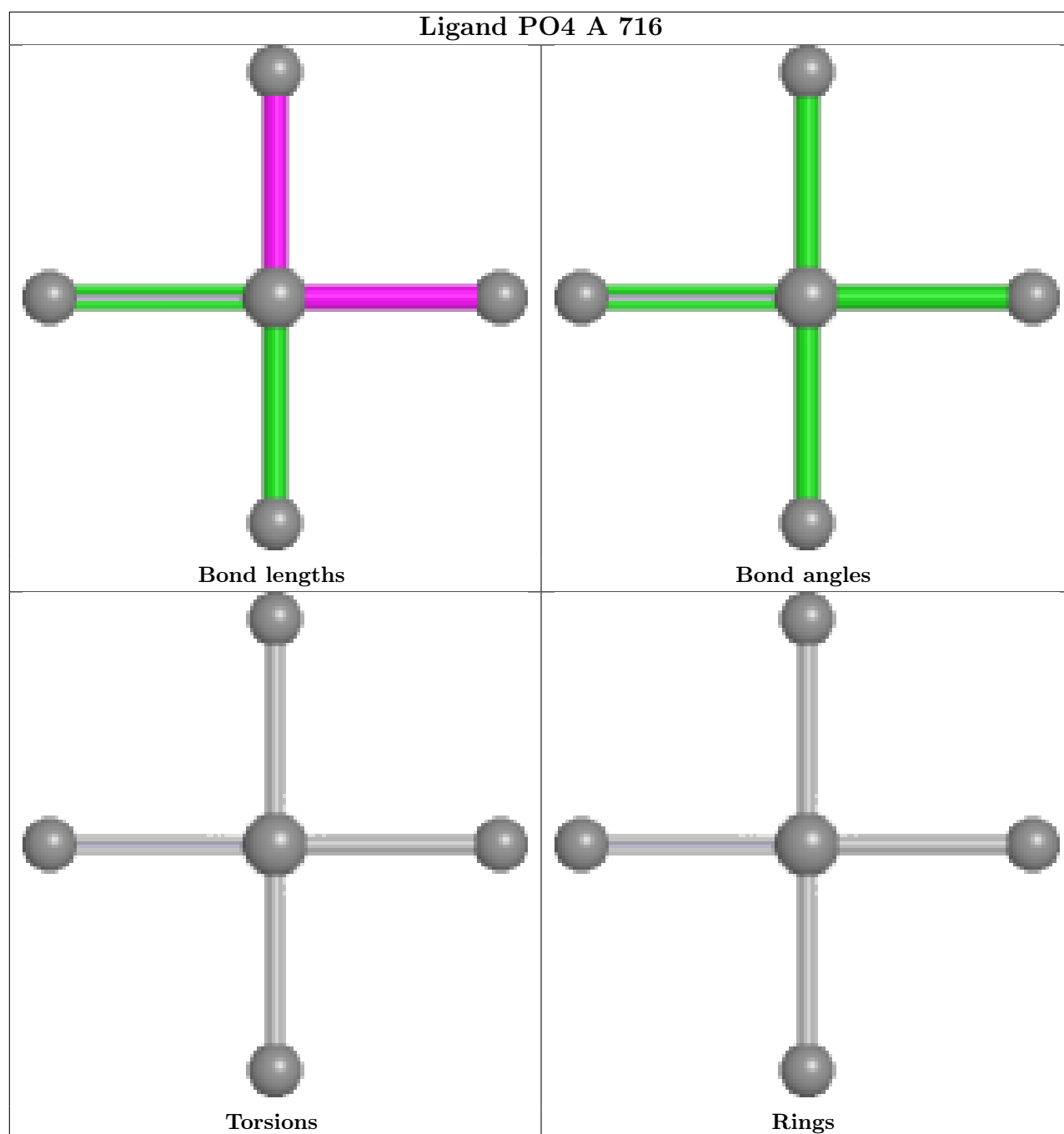


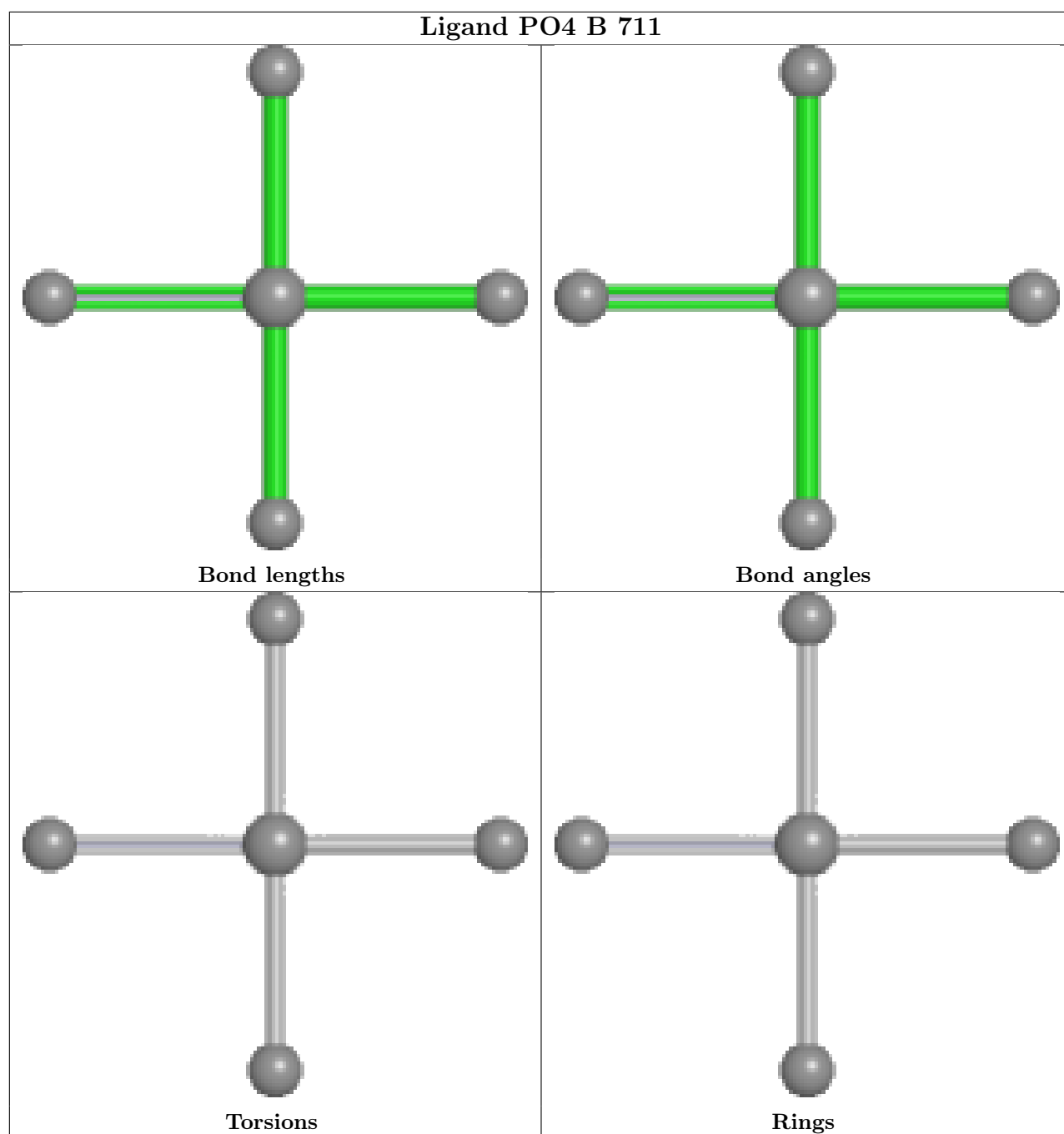


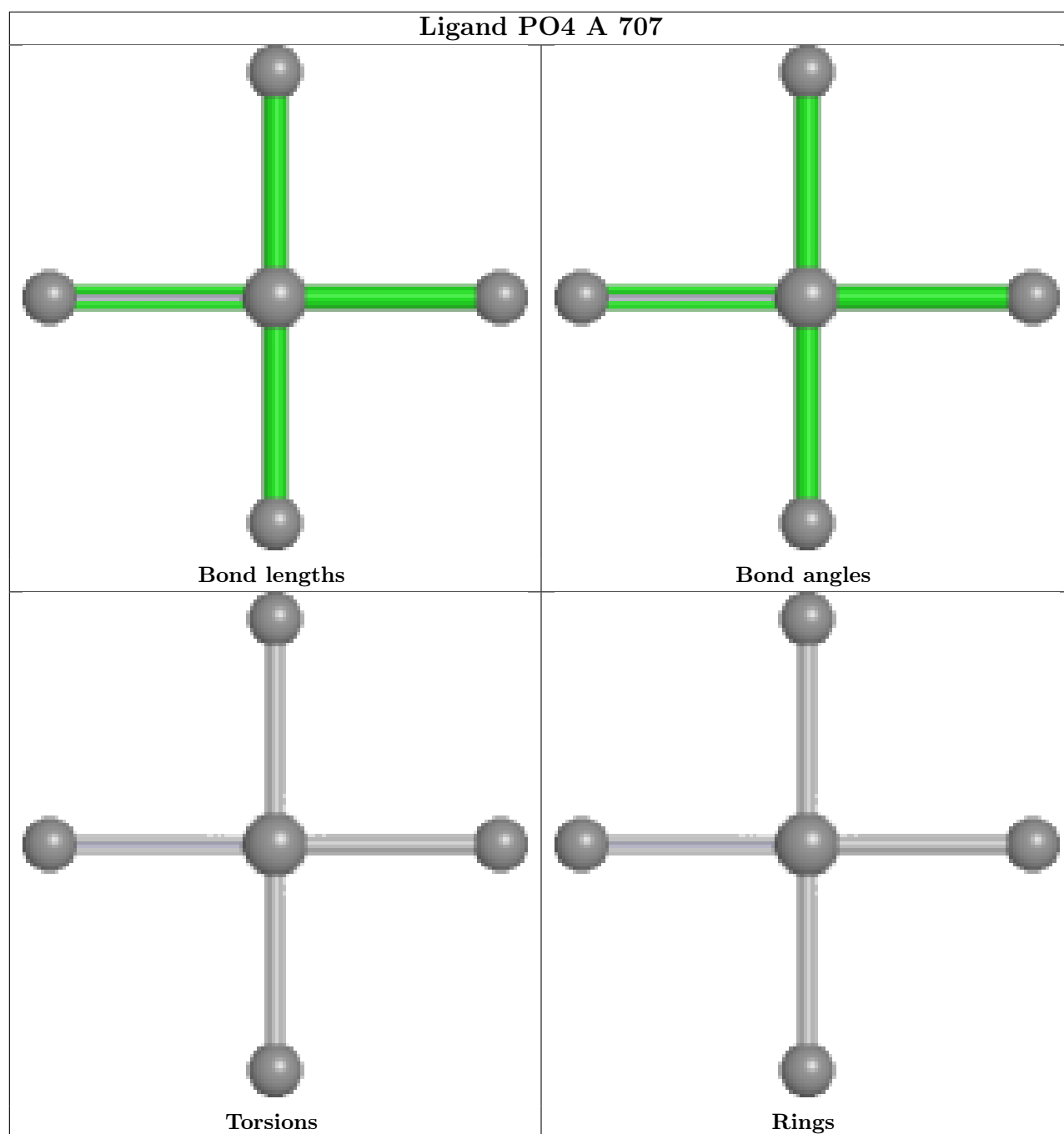


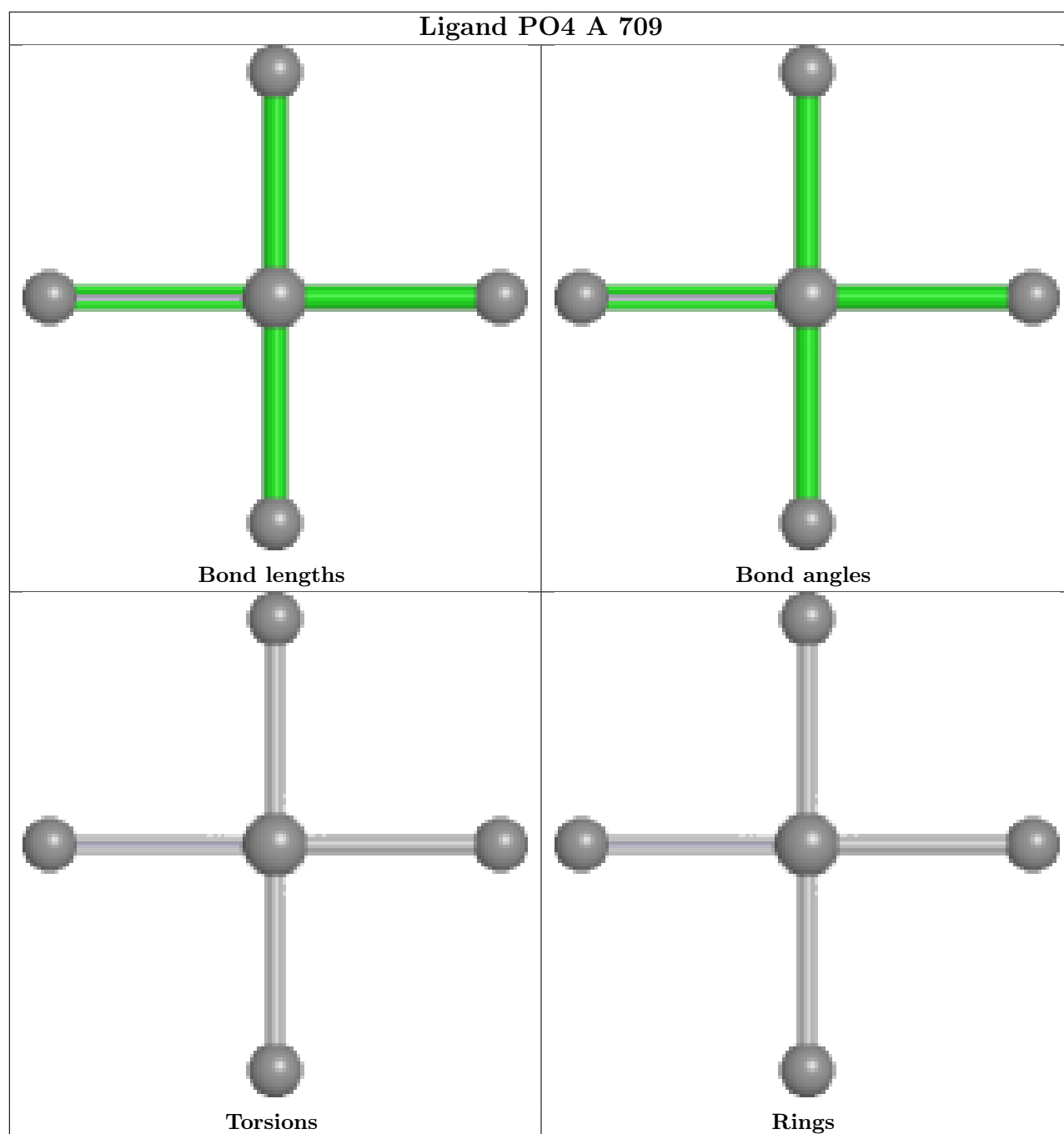


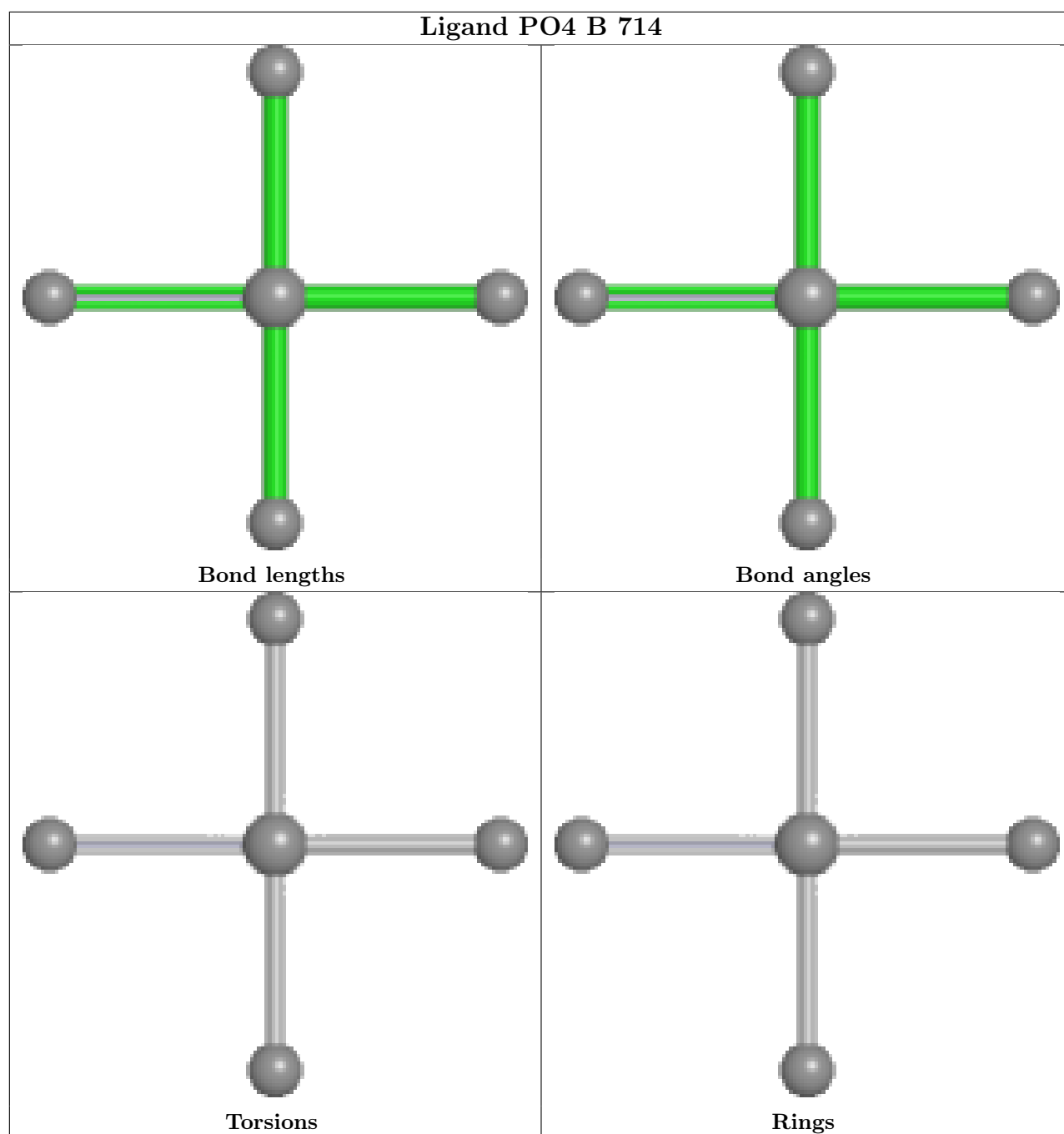


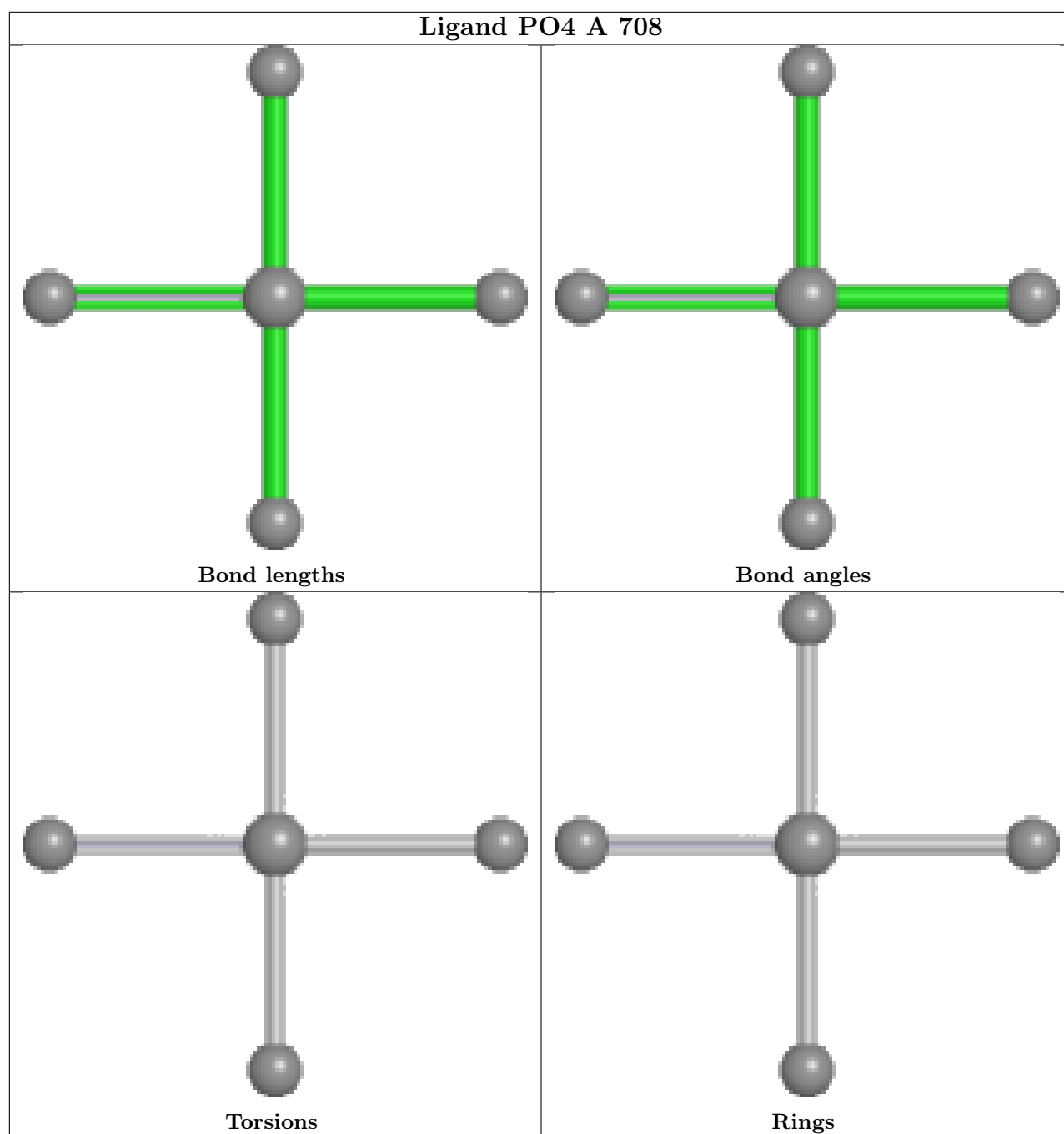


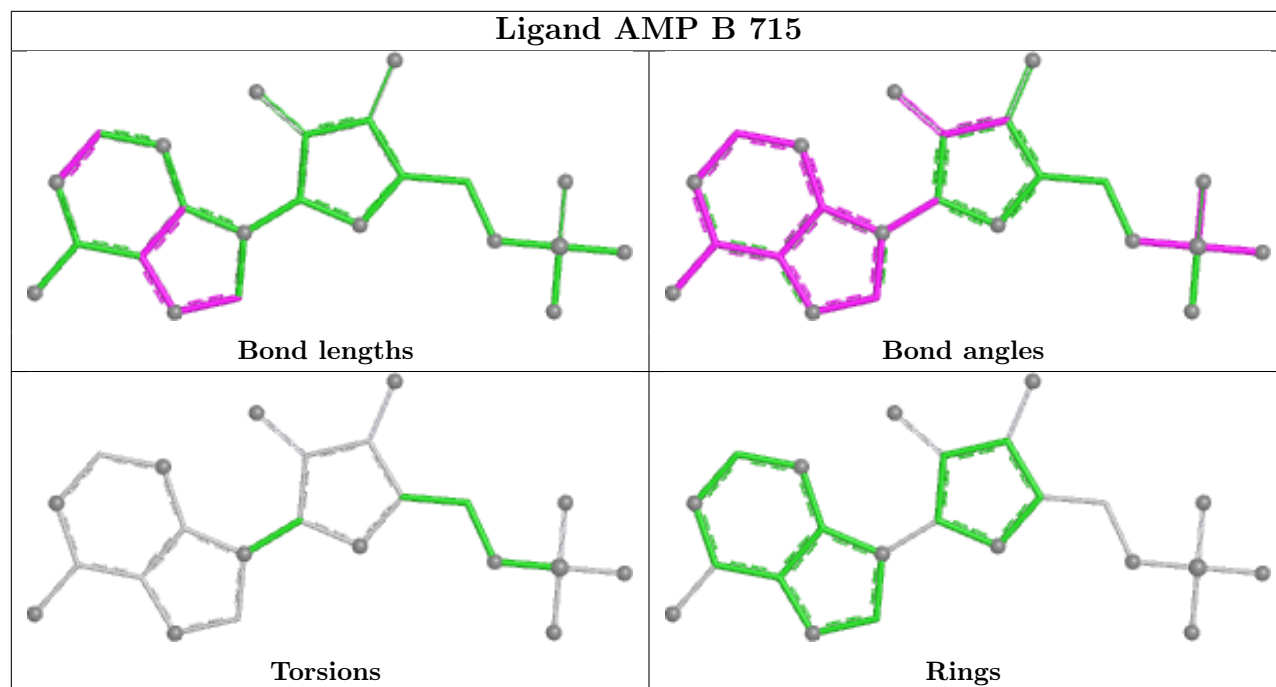


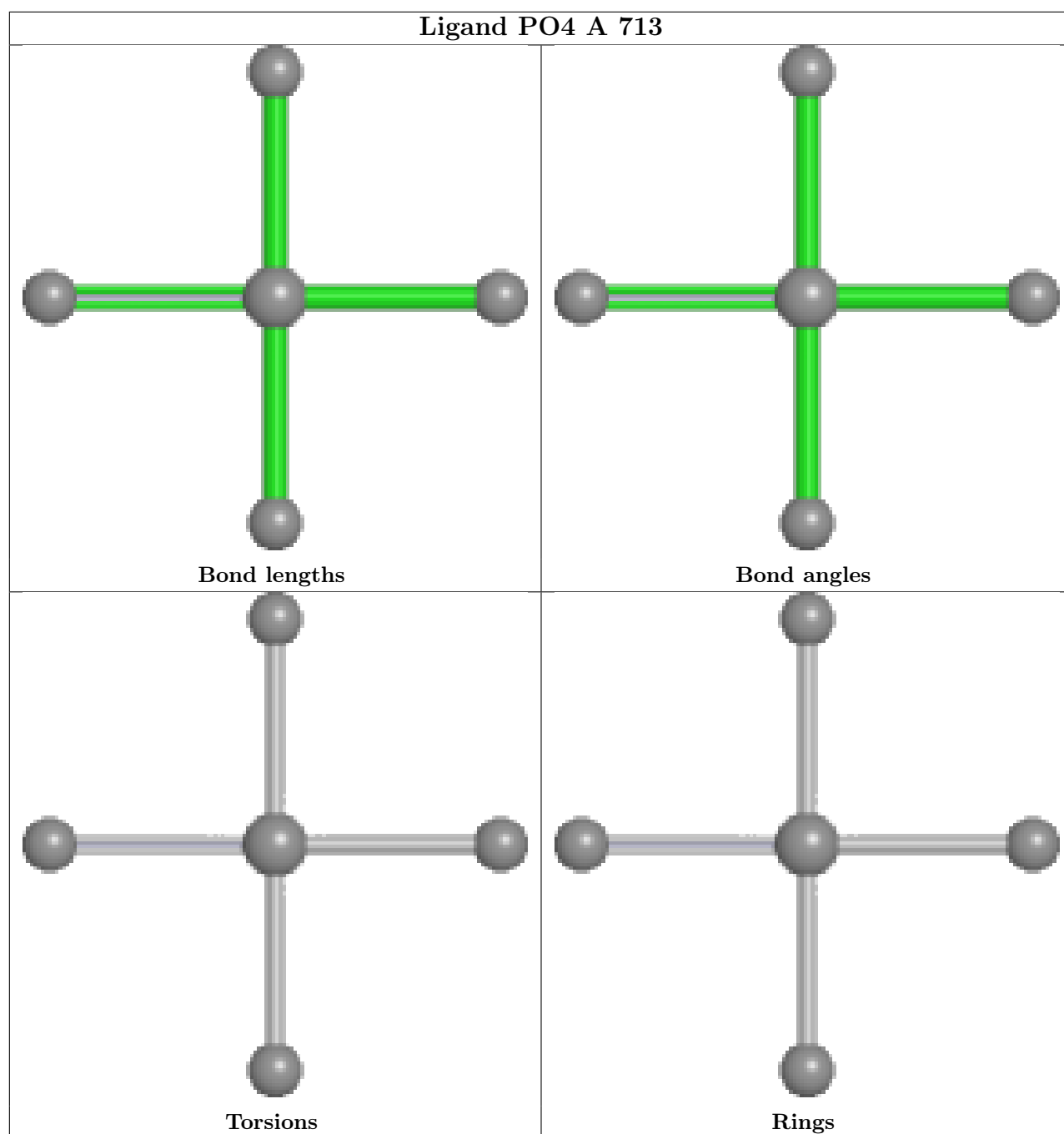












5.7 Other polymers [i](#)

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	606/606 (100%)	0.03	61 (10%)	12 11	5, 20, 114, 188	9 (1%)
1	B	606/606 (100%)	0.03	59 (9%)	13 12	5, 21, 105, 182	9 (1%)
All	All	1212/1212 (100%)	0.03	120 (9%)	13 12	5, 20, 112, 188	18 (1%)

All (120) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	626	ILE	6.2
1	A	578	LEU	6.0
1	A	582	LEU	5.8
1	B	485	PRO	5.7
1	A	589	THR	5.5
1	B	490	VAL	5.1
1	B	582	LEU	5.0
1	A	487	PRO	4.9
1	A	517	LYS	4.9
1	B	487	PRO	4.9
1	B	572	ILE	4.8
1	A	579	GLY	4.7
1	B	579	GLY	4.7
1	B	578	LEU	4.6
1	A	575	LYS	4.6
1	A	490	VAL	4.5
1	A	483	ILE	4.3
1	A	491	PRO	4.3
1	A	514	THR	4.1
1	B	626	ILE	4.1
1	B	486	ALA	4.1
1	B	515	GLY	4.1
1	B	453	VAL	4.0
1	A	484	PRO	4.0

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Mol	Chain	Res	Type	RSRZ
1	A	520	ILE	4.0
1	B	514	THR	4.0
1	A	583	SER	3.9
1	A	485	PRO	3.9
1	B	517	LYS	3.8
1	A	625	PRO	3.8
1	B	627	ILE	3.8
1	A	588	GLU	3.8
1	B	519	LYS	3.7
1	B	580	GLY	3.7
1	A	622	ILE	3.6
1	B	520	ILE	3.6
1	A	515	GLY	3.6
1	B	484	PRO	3.5
1	B	575	LYS	3.5
1	B	568	LEU	3.5
1	A	623	VAL	3.5
1	B	483	ILE	3.5
1	B	625	PRO	3.5
1	A	627	ILE	3.4
1	B	488	ARG	3.4
1	B	491	PRO	3.4
1	A	584	SER	3.4
1	A	592	LYS	3.3
1	A	572	ILE	3.3
1	A	486	ALA	3.3
1	B	581	LYS	3.2
1	B	623	VAL	3.1
1	A	580	GLY	3.1
1	A	519	LYS	3.1
1	A	412	GLY	3.0
1	B	482	GLY	3.0
1	A	624	GLN	3.0
1	A	214	ARG	2.9
1	A	585	GLU	2.9
1	A	453	VAL	2.9
1	A	574	ASP	2.9
1	A	488	ARG	2.9
1	B	489	GLY	2.9
1	B	571	GLN	2.9
1	B	527	ASN	2.9
1	B	513	GLY	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	614	ALA	2.8
1	B	614	ALA	2.8
1	A	518	ASN	2.8
1	A	586	ASP	2.8
1	B	573	GLY	2.8
1	B	384	PRO	2.8
1	B	518	ASN	2.8
1	A	590	MET	2.7
1	A	629	LYS	2.7
1	A	482	GLY	2.7
1	A	577	LYS	2.7
1	A	591	GLU	2.6
1	B	412	GLY	2.6
1	B	410	ASP	2.6
1	B	576	GLU	2.6
1	B	591	GLU	2.6
1	A	581	LYS	2.6
1	A	407	GLY	2.6
1	B	526	GLN	2.5
1	A	513	GLY	2.5
1	A	516	ASN	2.5
1	A	593	ALA	2.5
1	A	619	LEU	2.5
1	A	489	GLY	2.5
1	B	411	THR	2.5
1	A	628	SER	2.4
1	B	617	LYS	2.4
1	A	527	ASN	2.3
1	A	454	GLY	2.3
1	B	212	ASP	2.3
1	A	568	LEU	2.3
1	B	564	TYR	2.3
1	A	576	GLU	2.3
1	A	561	LEU	2.3
1	B	570	ASN	2.2
1	B	567	SER	2.2
1	B	609	ILE	2.2
1	A	564	TYR	2.2
1	B	583	SER	2.2
1	B	613	LYS	2.2
1	B	621	GLU	2.2
1	B	214	ARG	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	411	THR	2.2
1	B	624	GLN	2.1
1	B	584	SER	2.1
1	B	629	LYS	2.1
1	A	609	ILE	2.1
1	B	622	ILE	2.1
1	B	619	LEU	2.1
1	A	571	GLN	2.1
1	B	585	GLU	2.1
1	B	574	ASP	2.1
1	B	593	ALA	2.1
1	A	573	GLY	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](#)

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
2	GOL	A	704	6/6	0.74	0.15	48,52,53,55	0
4	PO4	A	712	5/5	0.75	0.21	57,67,76,79	0
2	GOL	A	701	6/6	0.80	0.18	57,60,63,69	0
4	PO4	B	713	5/5	0.80	0.20	75,75,84,84	0
2	GOL	A	703	6/6	0.81	0.12	53,54,57,58	0
2	GOL	B	706	6/6	0.83	0.15	50,56,66,73	0
4	PO4	B	714	5/5	0.84	0.12	75,77,81,81	0
2	GOL	A	702	6/6	0.85	0.13	38,40,45,47	0
7	PEG	A	717	7/7	0.86	0.14	43,49,52,53	0
2	GOL	B	704	6/6	0.87	0.10	34,34,39,44	0
2	GOL	B	701	6/6	0.88	0.12	30,40,43,45	0

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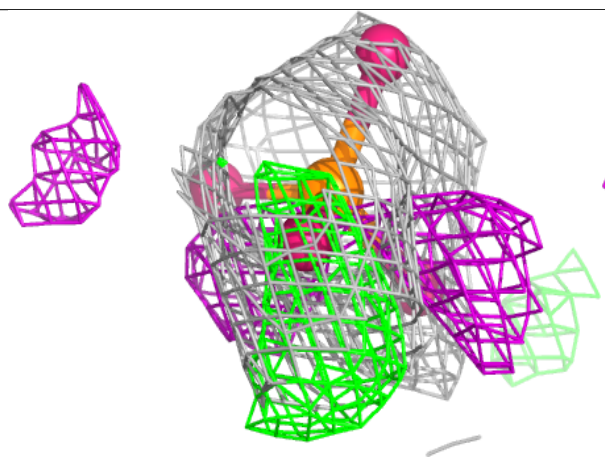
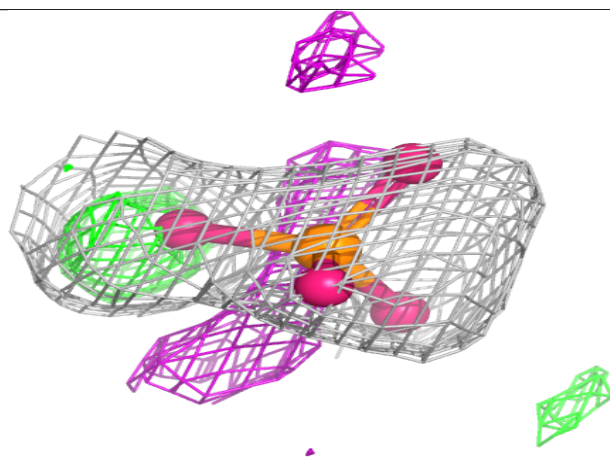
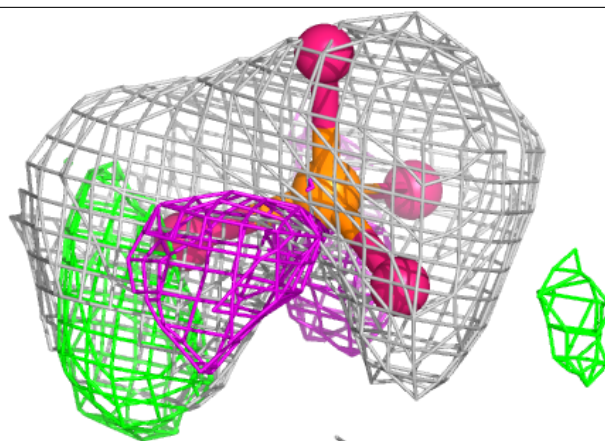
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	SO4	B	707	5/5	0.88	0.11	66,70,77,81	0
2	GOL	B	705	6/6	0.89	0.12	28,32,33,42	0
3	SO4	A	705	5/5	0.91	0.10	47,51,57,64	0
3	SO4	A	706	5/5	0.91	0.11	62,62,68,75	0
4	PO4	A	713	5/5	0.91	0.12	42,42,46,53	0
4	PO4	A	708	5/5	0.92	0.10	41,41,46,49	0
4	PO4	A	709	5/5	0.92	0.12	54,55,58,60	0
4	PO4	B	712	5/5	0.92	0.13	50,54,60,63	0
4	PO4	A	707	5/5	0.93	0.09	41,45,48,59	0
4	PO4	B	710	5/5	0.93	0.10	42,44,51,59	0
4	PO4	B	711	5/5	0.94	0.13	41,43,51,54	0
3	SO4	B	709	5/5	0.95	0.09	35,46,50,56	0
4	PO4	A	710	5/5	0.95	0.15	52,56,59,66	0
4	PO4	A	711	5/5	0.96	0.12	33,37,40,41	0
3	SO4	B	708	5/5	0.97	0.09	40,43,47,47	0
6	AMP	A	715	23/23	0.98	0.04	11,13,13,14	0
6	AMP	B	715	23/23	0.98	0.05	11,12,14,16	0
3	SO4	B	702	5/5	0.98	0.06	36,37,38,39	0
5	MG	B	703	1/1	0.99	0.05	22,22,22,22	0
4	PO4	A	716	5/5	0.99	0.04	13,15,17,19	0
4	PO4	B	716	5/5	0.99	0.04	13,13,16,17	0
5	MG	A	714	1/1	0.99	0.04	23,23,23,23	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.

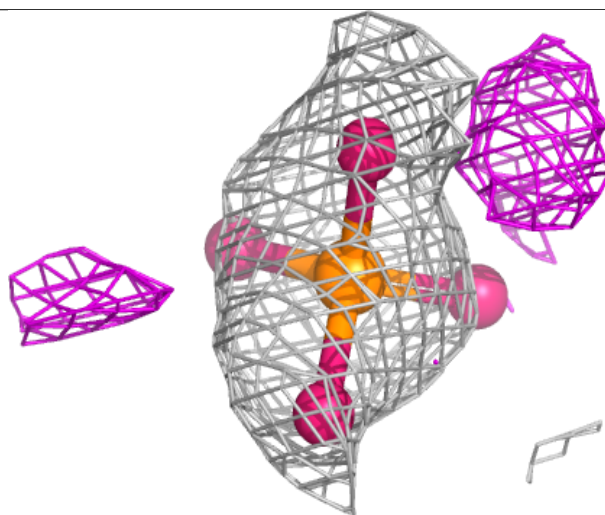
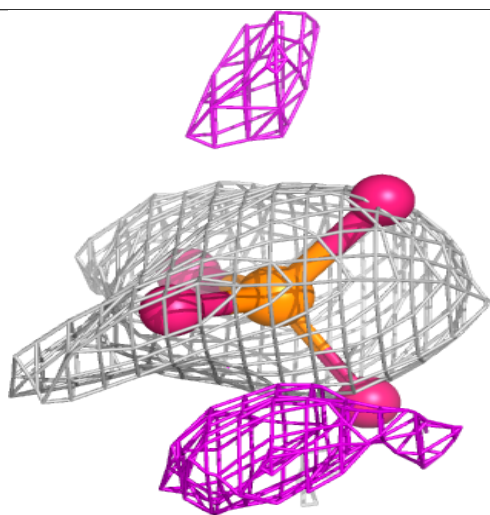
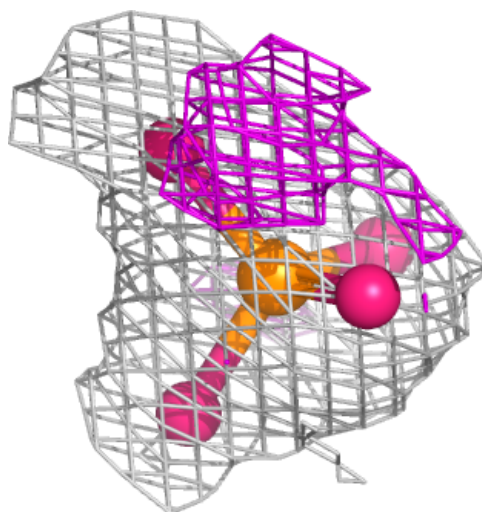
Electron density around PO4 A 712:

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



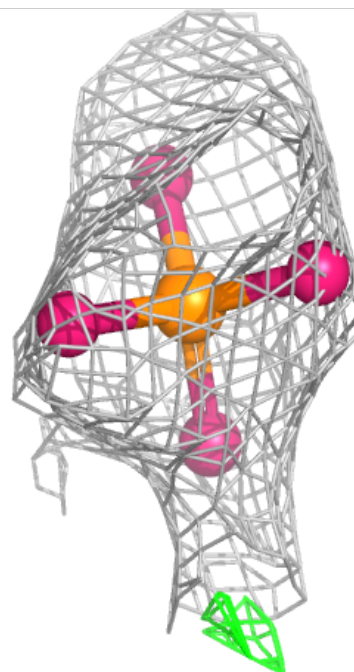
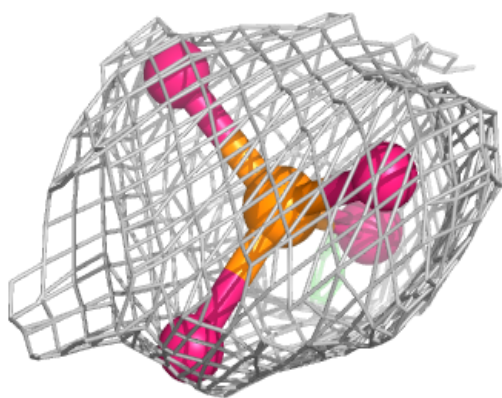
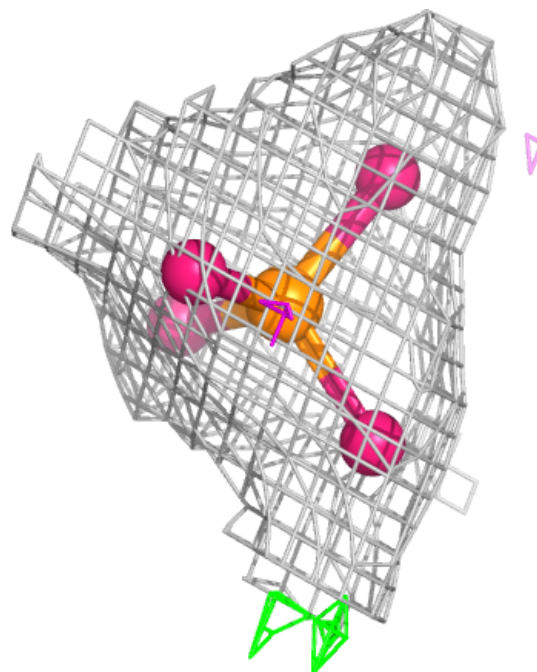
Electron density around PO4 B 713:

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



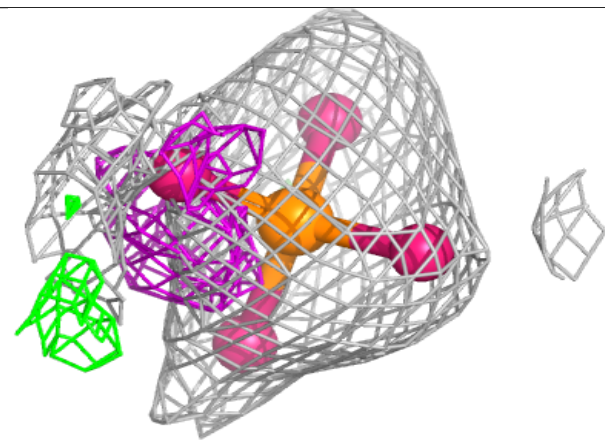
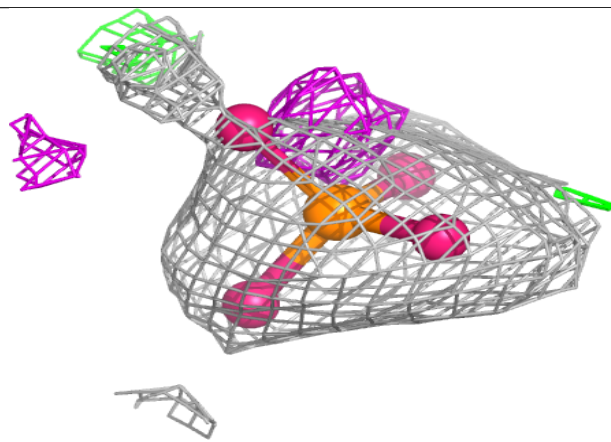
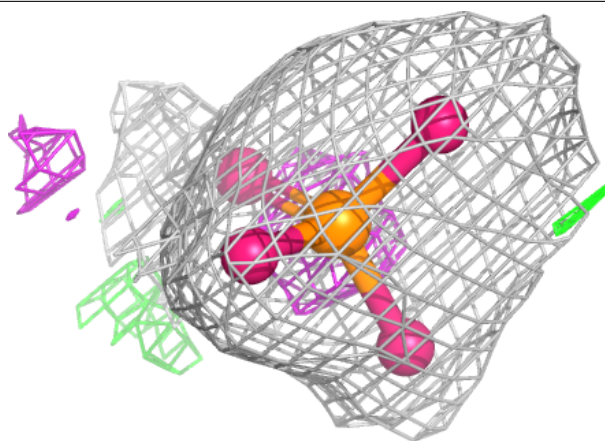
Electron density around PO4 B 714:

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and green (positive)



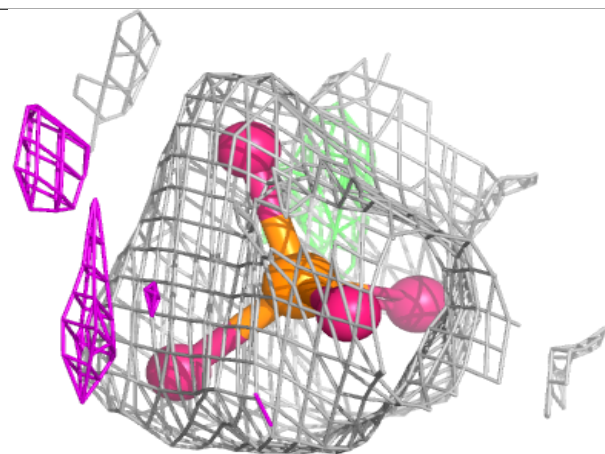
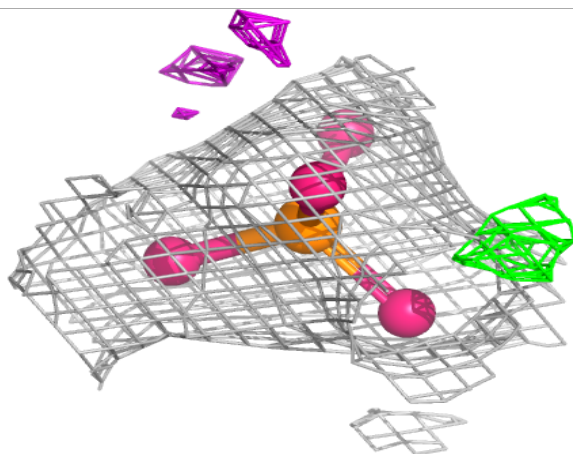
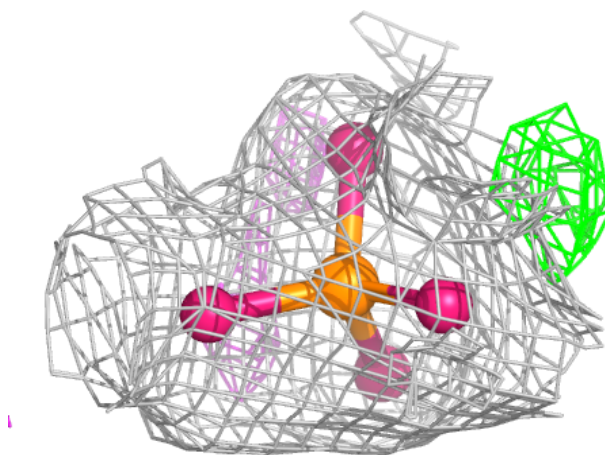
Electron density around PO4 A 713:

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



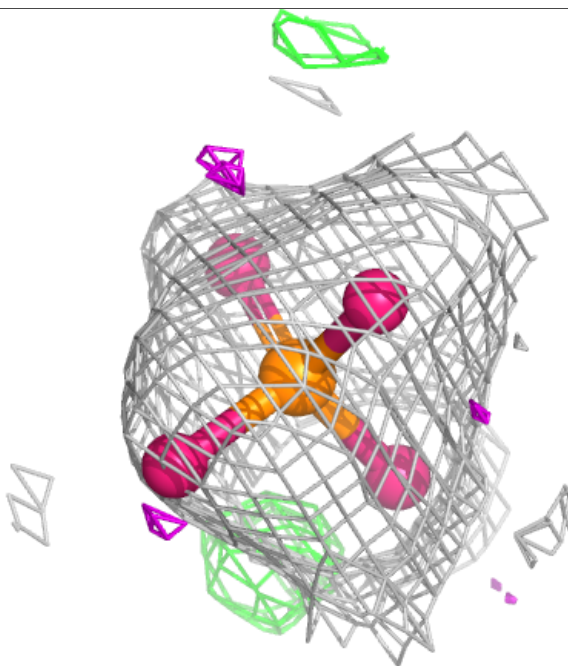
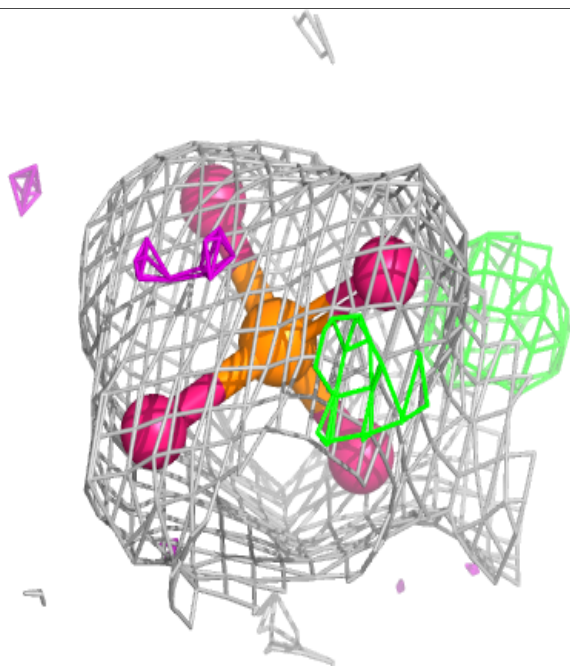
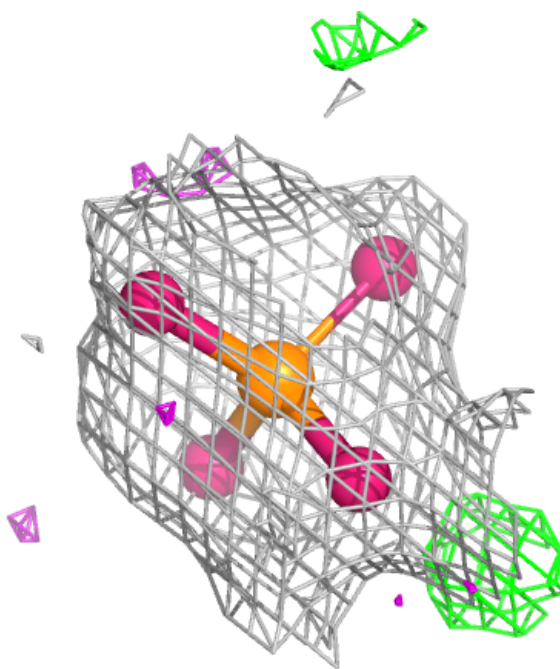
Electron density around PO4 A 708:

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



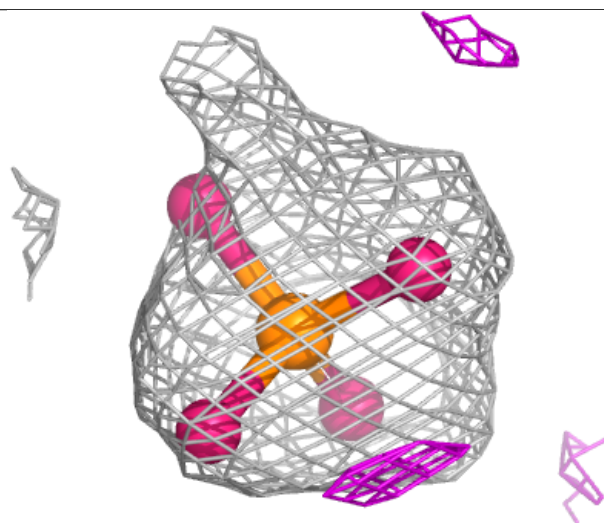
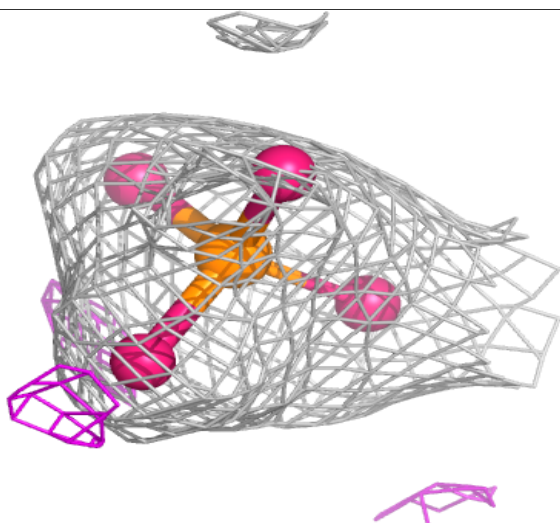
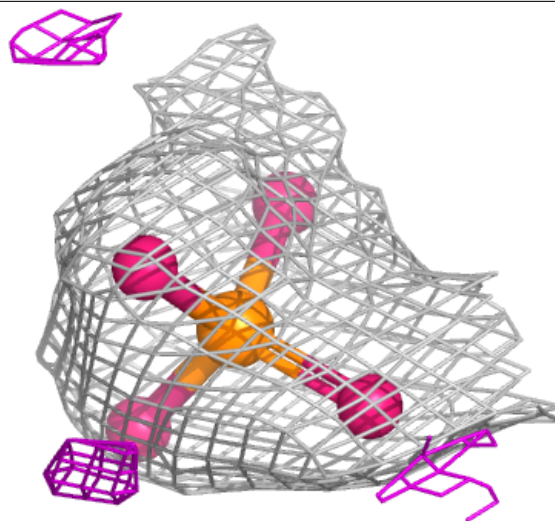
Electron density around PO4 A 709:

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



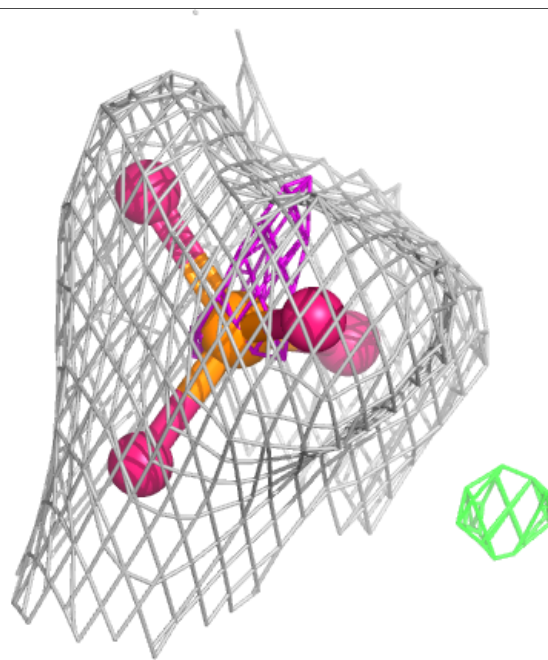
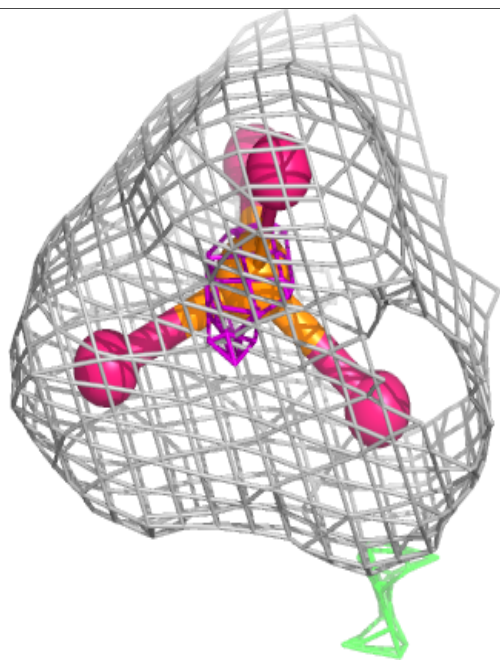
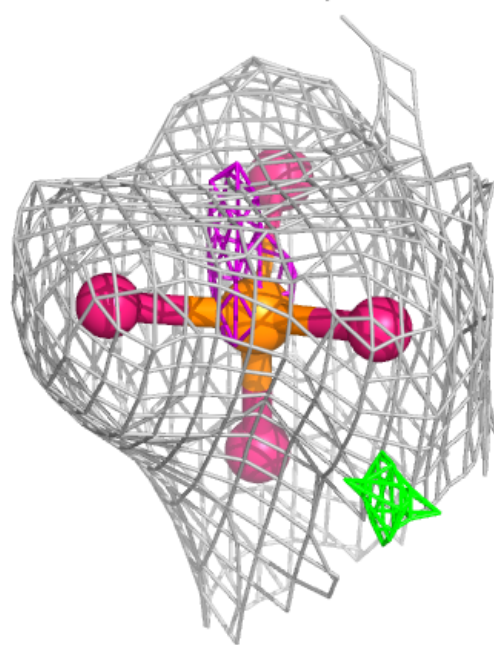
Electron density around PO4 B 712:

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



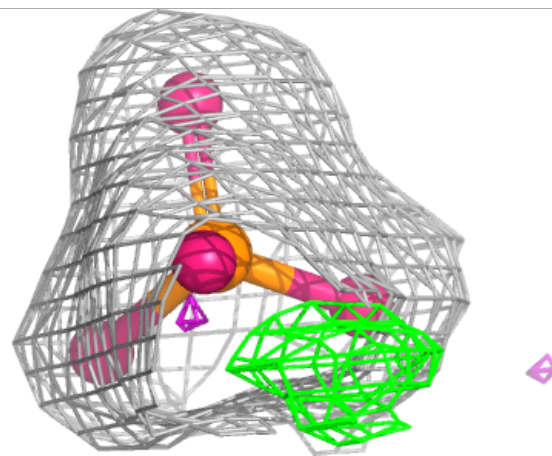
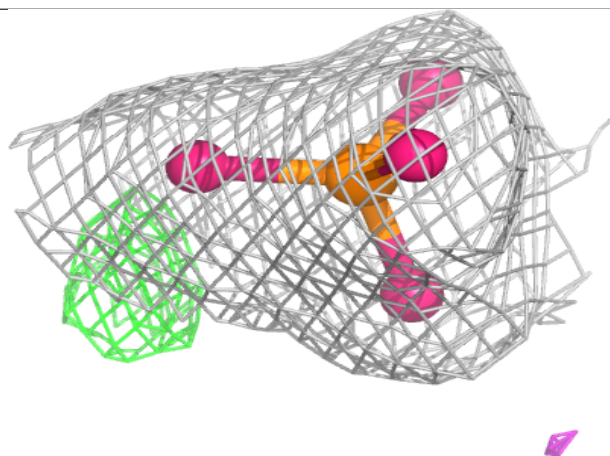
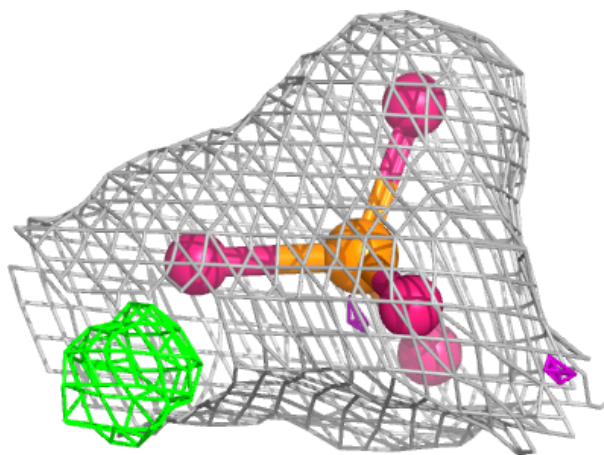
Electron density around PO4 A 707:

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and green (positive)



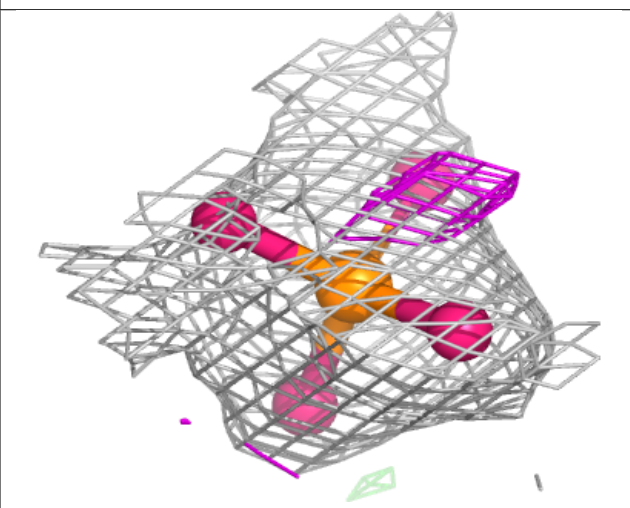
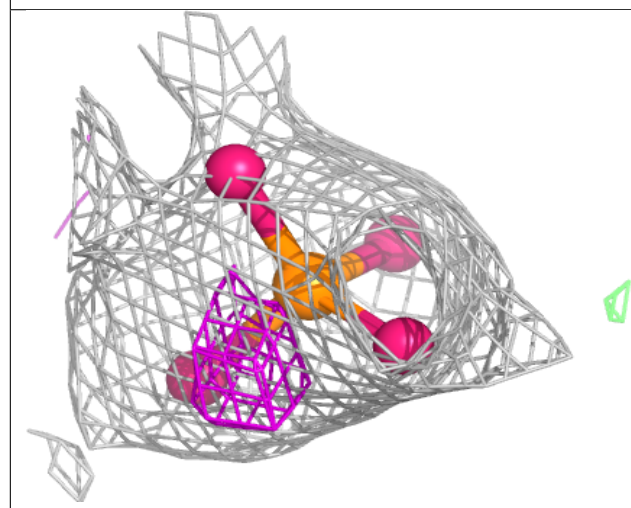
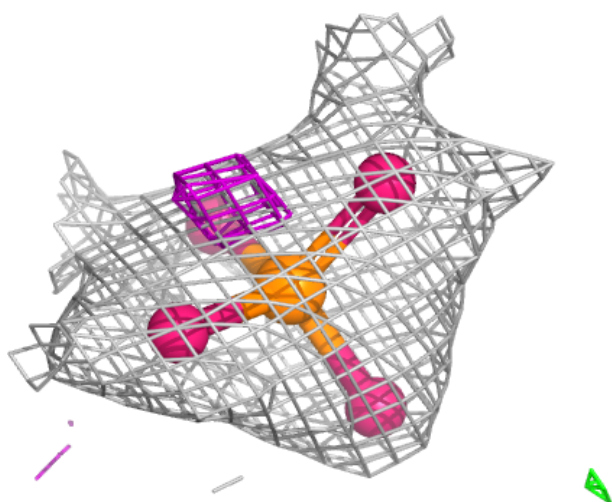
Electron density around PO4 B 710:

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and green (positive)



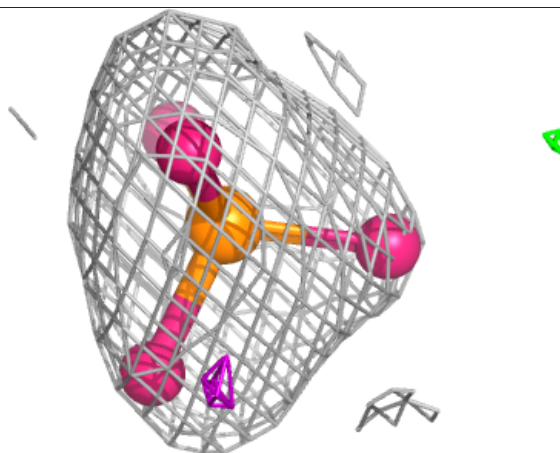
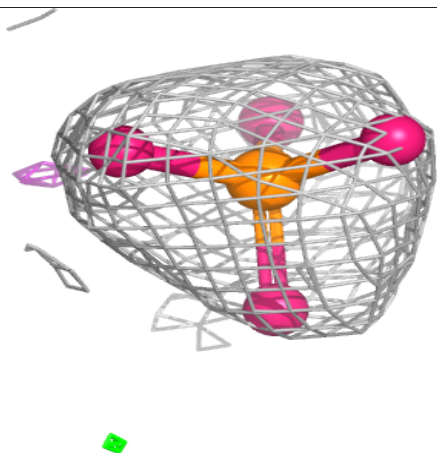
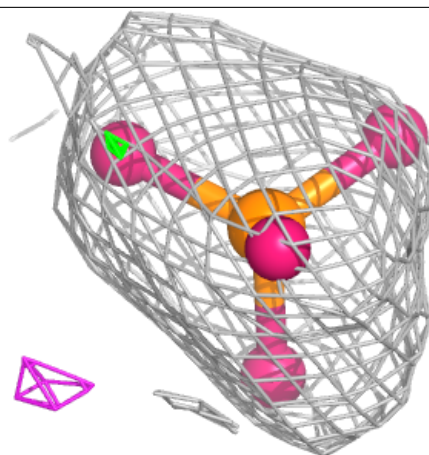
Electron density around PO4 B 711:

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



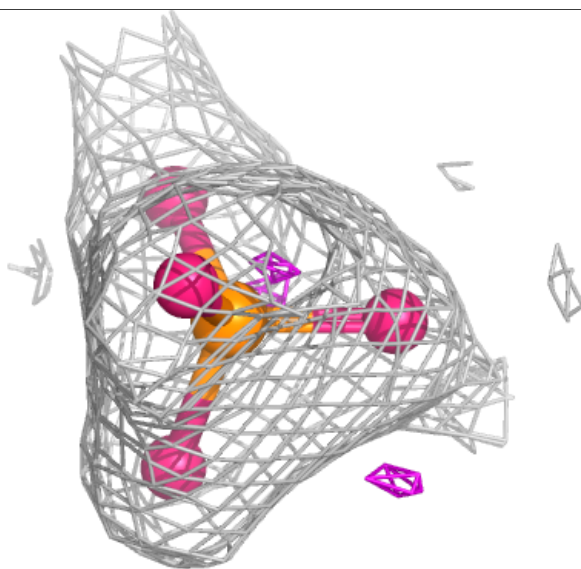
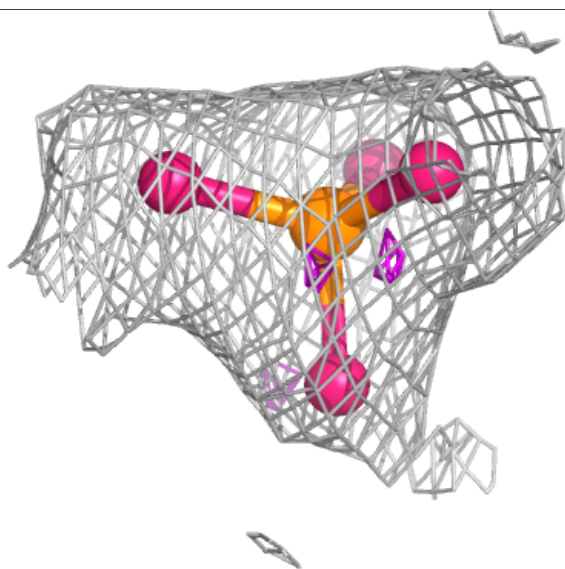
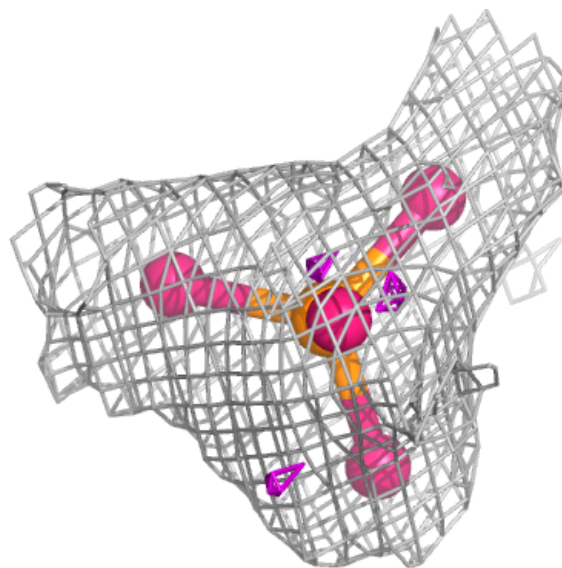
Electron density around PO4 A 710:

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and green (positive)



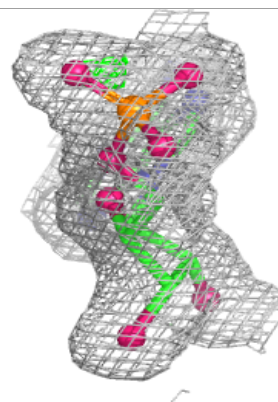
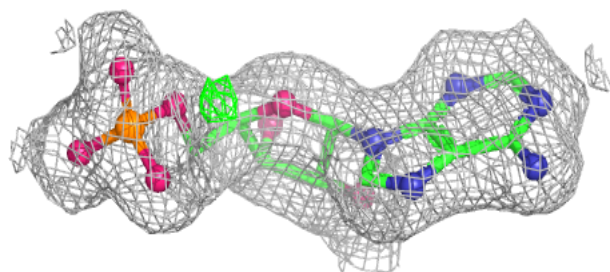
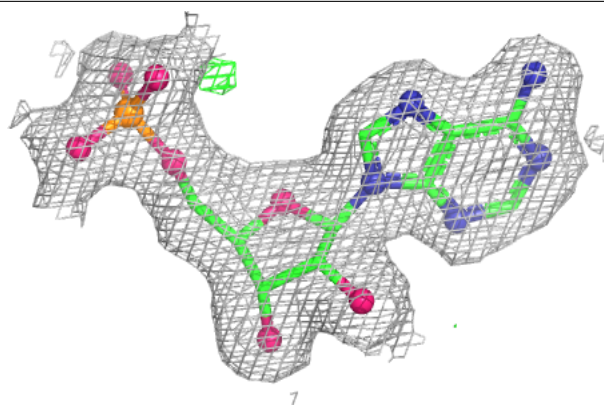
Electron density around PO4 A 711:

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and green (positive)

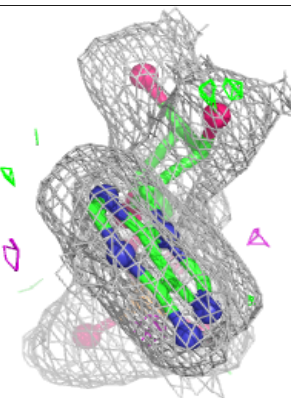
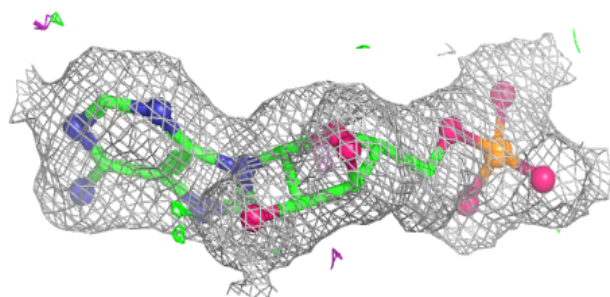
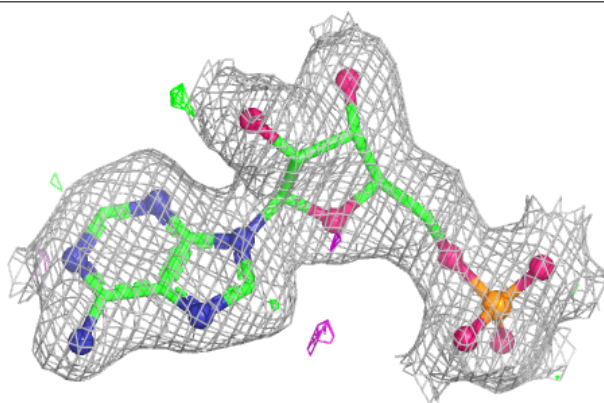


Electron density around AMP A 715:

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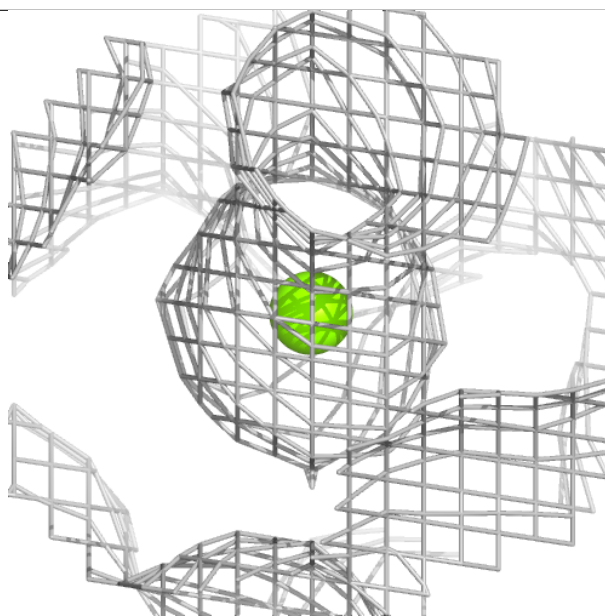
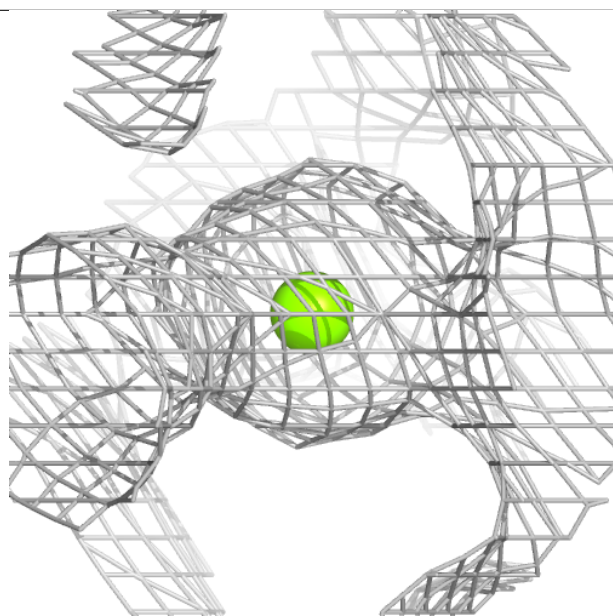
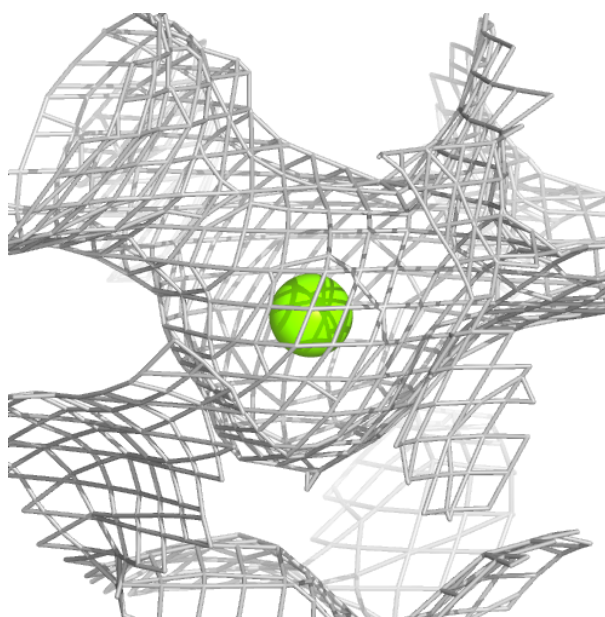
**Electron density around AMP B 715:**

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and green (positive)



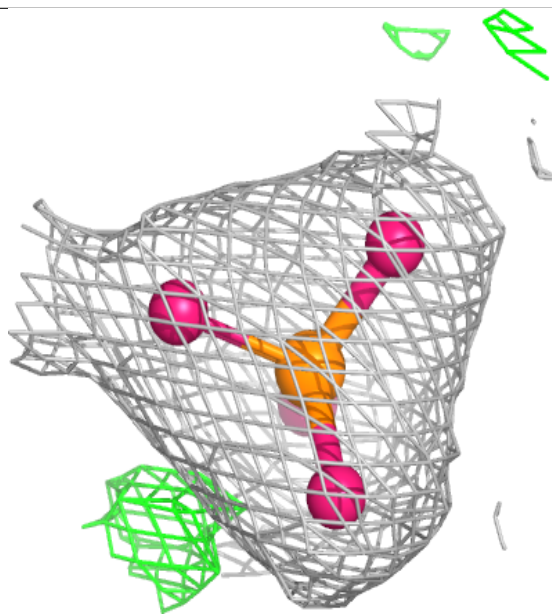
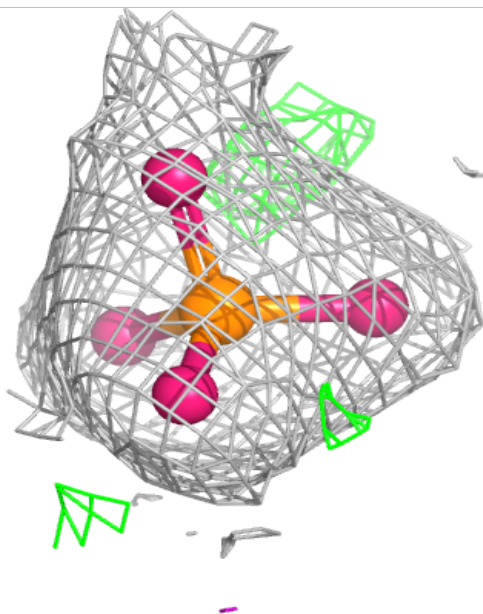
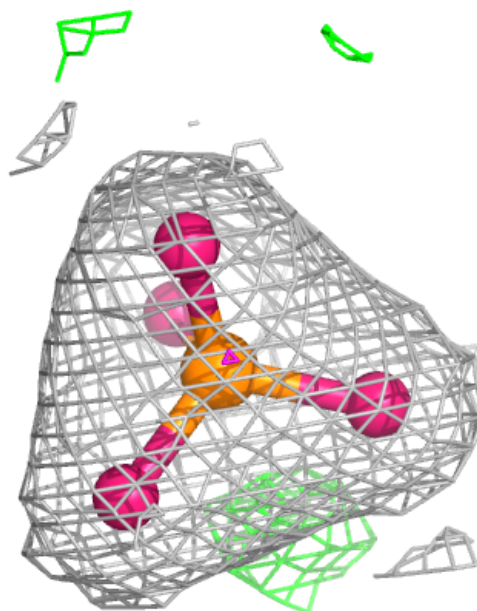
Electron density around MG B 703:

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and green (positive)



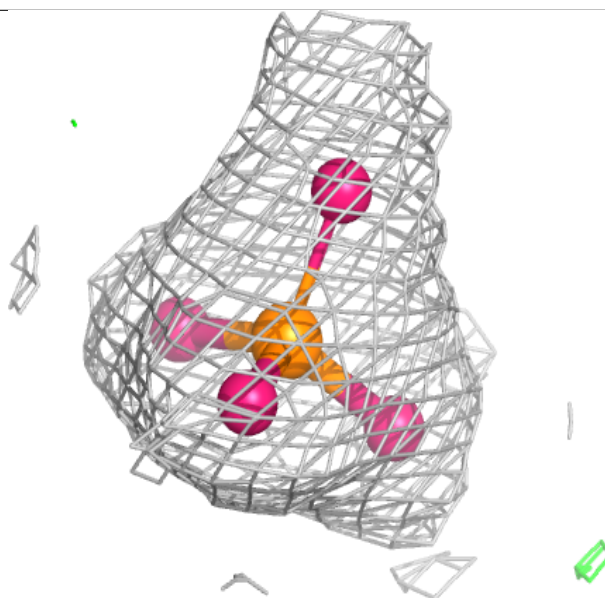
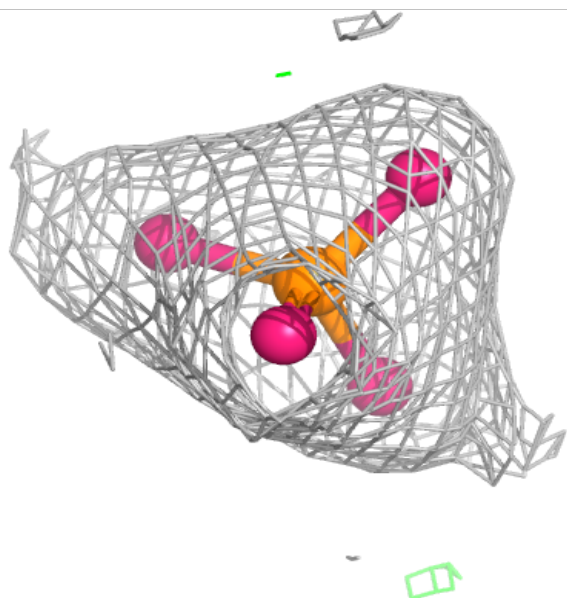
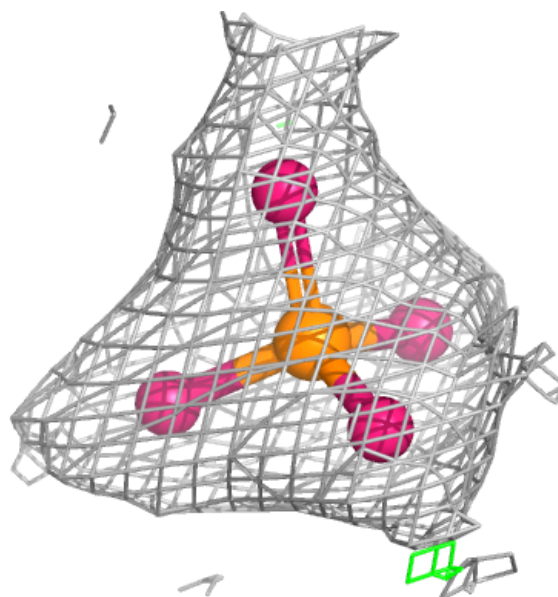
Electron density around PO4 A 716:

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



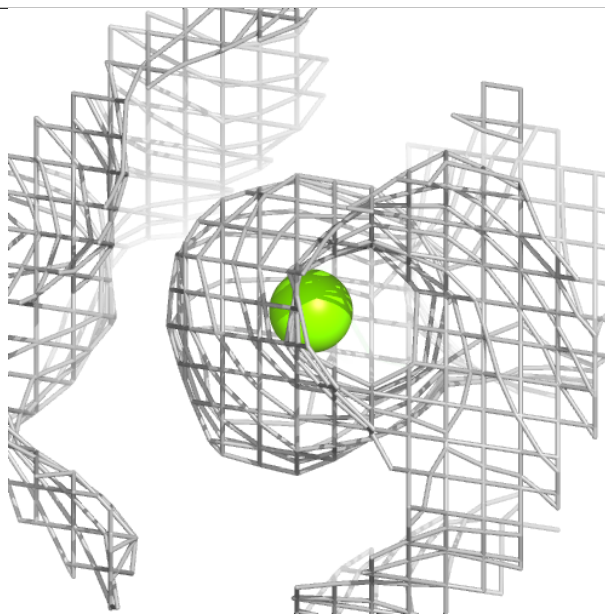
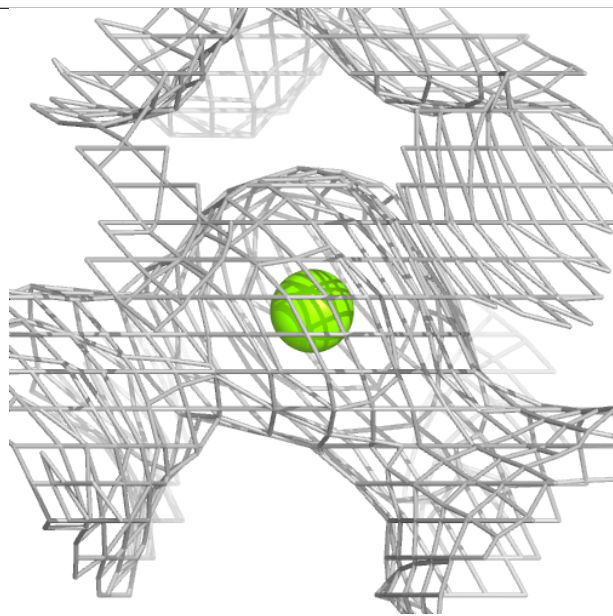
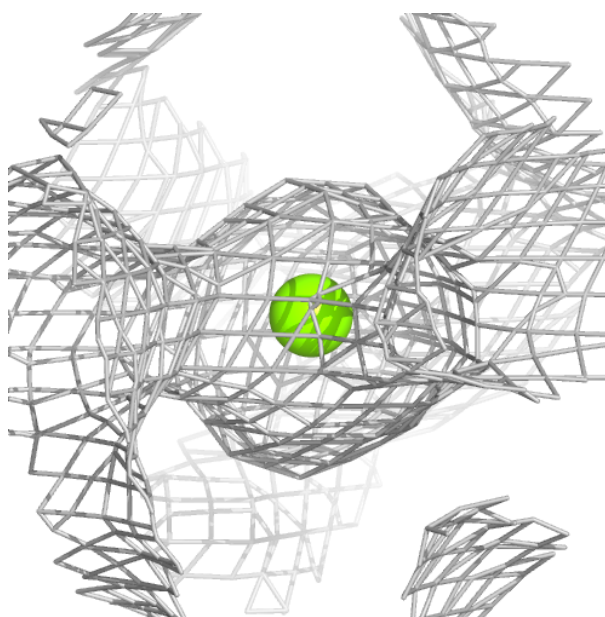
Electron density around PO4 B 716:

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



Electron density around MG A 714:

$2mF_o-DF_c$ (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.