



wwPDB X-ray Structure Validation Summary Report ⓘ

Jan 3, 2022 – 12:06 PM EST

PDB ID : 7N1R
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 wwPDB X-ray Structure Validation Summary 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.25
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.25

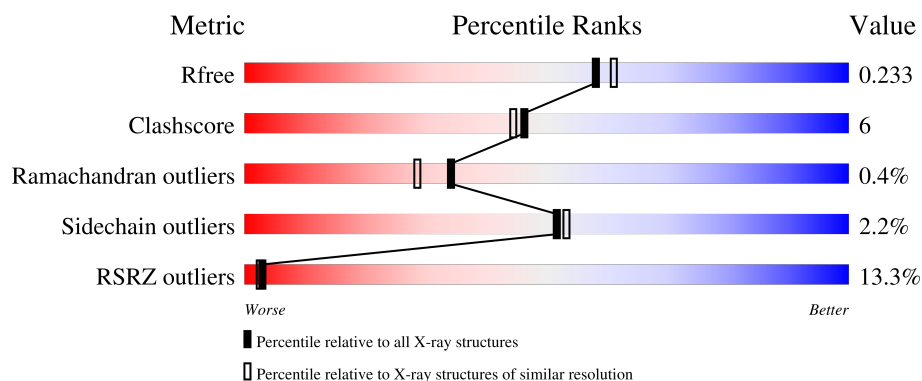
1 Overall quality at a glance

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}	130704	10434 (2.04-2.00)
Clashscore	141614	11643 (2.04-2.00)
Ramachandran outliers	138981	11493 (2.04-2.00)
Sidechain outliers	138945	11492 (2.04-2.00)
RSRZ outliers	127900	10220 (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	<div> <div>13%</div> <div>84%</div> <div>15%</div> </div>
1	B	606	<div> <div>13%</div> <div>87%</div> <div>11%</div> </div>

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:

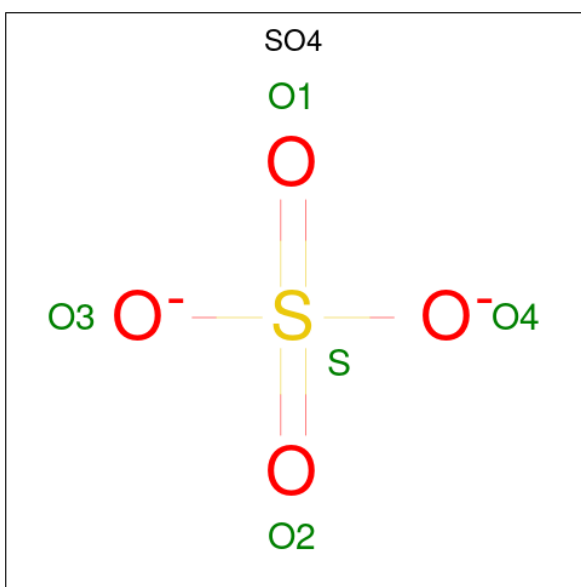
Chain	Residue	Modelled	Actual	Comment	Reference
A	24	SER	-	expression tag	UNP P11021
A	453	VAL	THR	conflict	UNP P11021
A	454	GLY	ALA	conflict	UNP P11021
A	455	GLY	SER	conflict	UNP P11021
A	?	-	ASP	deletion	UNP P11021
A	?	-	ASN	deletion	UNP P11021
A	?	-	GLN	deletion	UNP P11021
A	?	-	PRO	deletion	UNP P11021
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 (three-letter code: GOL) (formula: C₃H₈O₃).



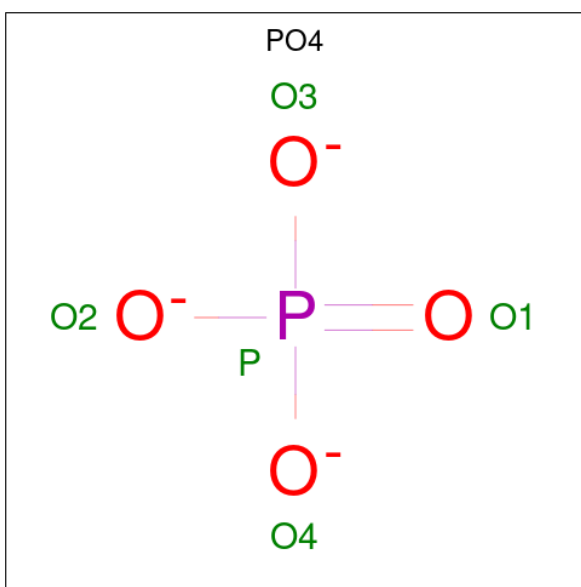
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			6	3	3		
2	A	1	Total	C	O	0	0
			6	3	3		
2	A	1	Total	C	O	0	0
			6	3	3		
2	A	1	Total	C	O	0	0
			6	3	3		
2	B	1	Total	C	O	0	0
			6	3	3		
2	B	1	Total	C	O	0	0
			6	3	3		
2	B	1	Total	C	O	0	0
			6	3	3		

- Molecule 3 is SULFATE ION (three-letter code: 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 (three-letter code: PO4) (formula: O₄P) (labeled as "Ligand of Interest" by depositor).

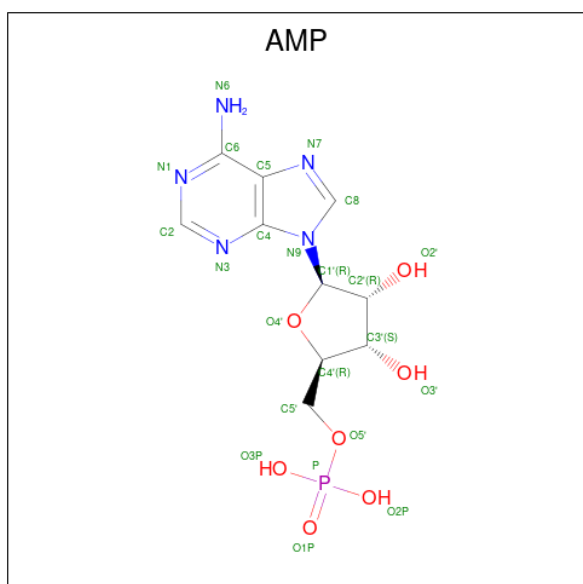


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	O	P	0	0
			5	4	1		
4	A	1	Total	O	P	0	0
			5	4	1		
4	A	1	Total	O	P	0	0
			5	4	1		
4	A	1	Total	O	P	0	0
			5	4	1		
4	A	1	Total	O	P	0	0
			5	4	1		
4	A	1	Total	O	P	0	0
			5	4	1		
4	A	1	Total	O	P	0	0
			5	4	1		
4	B	1	Total	O	P	0	0
			5	4	1		
4	B	1	Total	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 (three-letter code: MG) (formula: Mg) (labeled as "Ligand of Interest" by depositor).

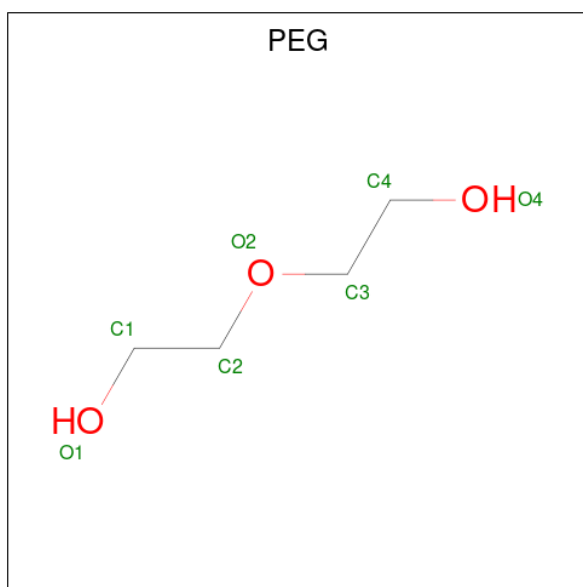
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total	Mg	0	0
			1	1		
5	B	1	Total	Mg	0	0
			1	1		

- Molecule 6 is ADENOSINE MONOPHOSPHATE (three-letter code: 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 (three-letter code: PEG) (formula: C₄H₁₀O₃).



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

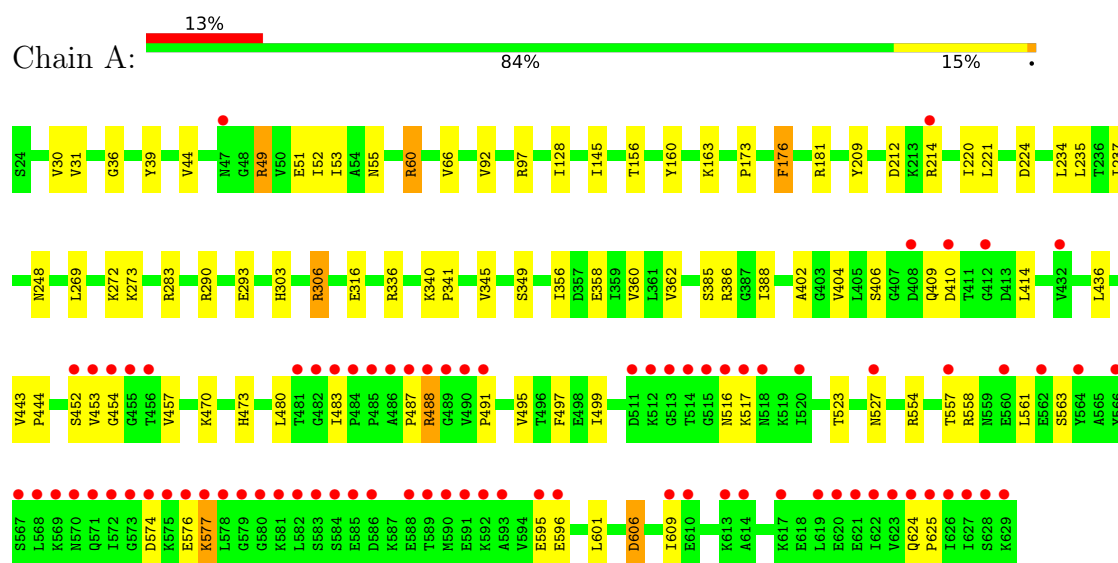
- Molecule 8 is water.

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

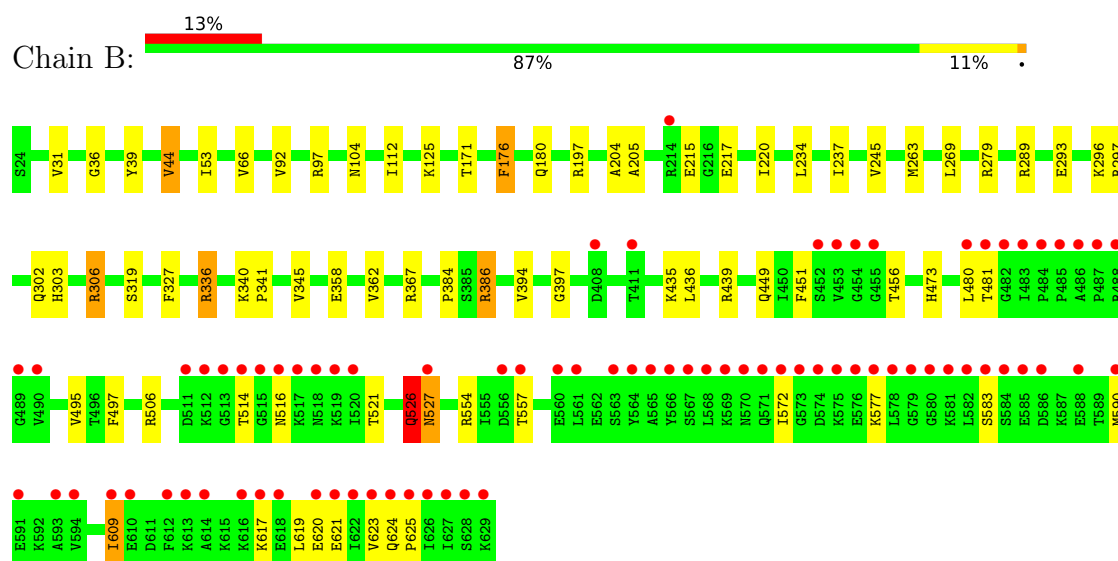
3 Residue-property plots [i](#)

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 39.63 – 2.03	Depositor EDS
% Data completeness (in resolution range)	96.8 (45.31-2.03) 96.9 (39.63-2.03)	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 (5.01%)	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 , 53.4	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: MG, AMP, PO4, PEG, SO4, GOL

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	0.57	0/4839	0.71	1/6533 (0.0%)
1	B	0.58	0/4844	0.70	1/6539 (0.0%)
All	All	0.58	0/9683	0.71	2/13072 (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

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	181	ARG	CG-CD-NE	-5.47	100.31	111.80
1	B	367	ARG	NE-CZ-NH2	-5.34	117.63	120.30

There are no chirality outliers.

5 of 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	60	ARG	Sidechain

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Group
1	A	97	ARG	Sidechain

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4769	0	4796	57	0
1	B	4771	0	4797	61	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	14	1
8	B	531	0	0	15	1
All	All	10811	0	9691	122	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 6.

The worst 5 of 122 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

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the sym-

metry 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%)	29	22
1	B	613/606 (101%)	594 (97%)	17 (3%)	2 (0%)	41	36
All	All	1226/1212 (101%)	1181 (96%)	40 (3%)	5 (0%)	34	28

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%)	509 (97%)	16 (3%)	41	40
1	B	525/519 (101%)	516 (98%)	9 (2%)	60	63

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	1050/1038 (101%)	1025 (98%)	25 (2%)	52 49

5 of 25 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	596	GLU
1	B	44[B]	VAL
1	B	621	GLU
1	B	44[A]	VAL
1	B	176	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 11 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	303	HIS
1	B	516	ASN
1	B	624	GLN
1	B	526	GLN
1	A	527	ASN

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 monosaccharides 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
2	GOL	B	704	-	5,5,5	0.41	0	5,5,5	0.69	0
2	GOL	A	702	-	5,5,5	0.44	0	5,5,5	0.83	0
2	GOL	A	703	-	5,5,5	0.50	0	5,5,5	0.34	0
4	PO4	A	711	-	4,4,4	0.97	0	6,6,6	0.64	0
4	PO4	A	712	-	4,4,4	1.19	0	6,6,6	0.66	0
4	PO4	B	710	-	4,4,4	0.73	0	6,6,6	0.73	0
3	SO4	B	702	-	4,4,4	0.47	0	6,6,6	0.88	0
3	SO4	A	706	-	4,4,4	0.44	0	6,6,6	0.53	0
2	GOL	B	701	-	5,5,5	0.43	0	5,5,5	0.88	0
4	PO4	B	713	-	4,4,4	0.88	0	6,6,6	0.61	0
2	GOL	A	704	-	5,5,5	0.57	0	5,5,5	0.33	0
3	SO4	B	709	-	4,4,4	0.36	0	6,6,6	0.61	0
7	PEG	A	717	-	6,6,6	0.53	0	5,5,5	0.58	0
4	PO4	A	716	5	4,4,4	2.34	2 (50%)	6,6,6	1.13	0
4	PO4	B	711	-	4,4,4	0.47	0	6,6,6	0.59	0
4	PO4	B	716	5	4,4,4	2.09	2 (50%)	6,6,6	0.71	0
4	PO4	A	708	-	4,4,4	0.80	0	6,6,6	0.97	0
4	PO4	A	710	-	4,4,4	0.85	0	6,6,6	0.31	0
3	SO4	B	707	-	4,4,4	0.46	0	6,6,6	0.23	0
4	PO4	A	707	-	4,4,4	0.65	0	6,6,6	0.57	0
4	PO4	A	709	-	4,4,4	0.85	0	6,6,6	0.73	0
2	GOL	B	706	-	5,5,5	0.53	0	5,5,5	0.40	0
4	PO4	B	712	-	4,4,4	0.68	0	6,6,6	0.59	0
3	SO4	B	708	-	4,4,4	0.36	0	6,6,6	0.69	0
3	SO4	A	705	-	4,4,4	0.47	0	6,6,6	0.32	0
2	GOL	B	705	-	5,5,5	0.66	0	5,5,5	0.79	0
4	PO4	A	713	-	4,4,4	0.71	0	6,6,6	0.85	0
2	GOL	A	701	-	5,5,5	0.38	0	5,5,5	0.62	0
4	PO4	B	714	-	4,4,4	0.79	0	6,6,6	0.59	0
6	AMP	B	715	5	22,25,25	1.15	2 (9%)	25,38,38	2.10	9 (36%)
6	AMP	A	715	5	22,25,25	1.45	4 (18%)	25,38,38	1.26	5 (20%)

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	B	701	-	-	2/4/4/4	-
2	GOL	A	701	-	-	4/4/4/4	-
2	GOL	A	702	-	-	2/4/4/4	-
2	GOL	B	704	-	-	4/4/4/4	-
2	GOL	B	706	-	-	4/4/4/4	-
2	GOL	A	704	-	-	4/4/4/4	-
2	GOL	A	703	-	-	2/4/4/4	-
7	PEG	A	717	-	-	2/4/4/4	-
6	AMP	B	715	5	-	0/6/26/26	0/3/3/3
6	AMP	A	715	5	-	0/6/26/26	0/3/3/3
2	GOL	B	705	-	-	2/4/4/4	-

The worst 5 of 10 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	A	715	AMP	C2-N3	3.88	1.38	1.32
4	B	716	PO4	P-O3	-3.11	1.45	1.54
6	A	715	AMP	C2'-C1'	-3.11	1.49	1.53
4	A	716	PO4	P-O3	-2.90	1.45	1.54
6	A	715	AMP	C2-N1	2.90	1.39	1.33

The worst 5 of 14 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	B	715	AMP	O3P-P-O2P	4.85	126.17	107.64
6	B	715	AMP	C5-C6-N6	-4.30	113.82	120.35
6	B	715	AMP	N3-C2-N1	-3.90	122.58	128.68
6	B	715	AMP	C1'-N9-C4	-3.46	120.57	126.64
6	B	715	AMP	N6-C6-N1	3.45	125.73	118.57

There are no chirality outliers.

5 of 26 torsion outliers are listed below:

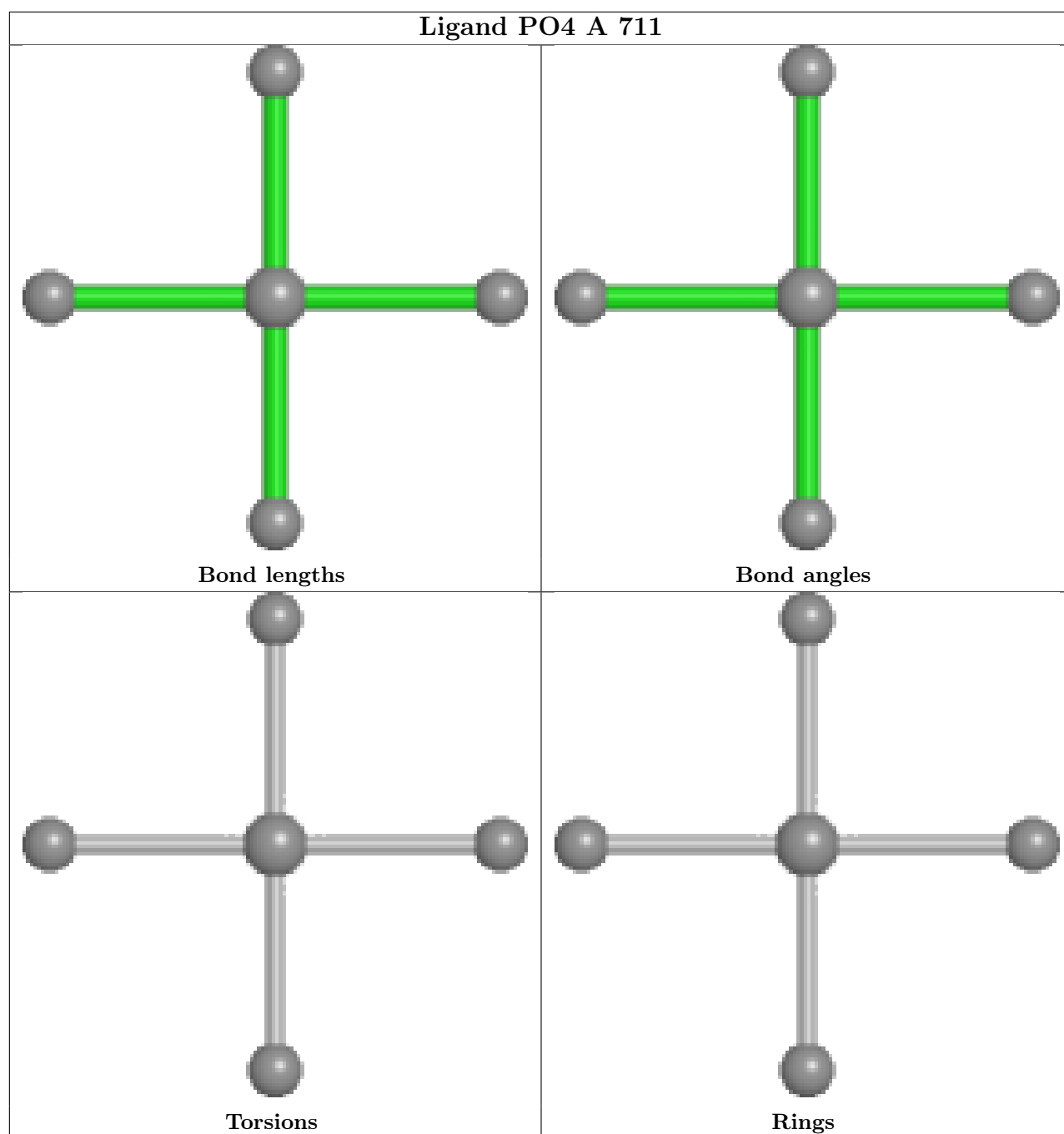
Mol	Chain	Res	Type	Atoms
2	A	701	GOL	O1-C1-C2-C3
2	A	701	GOL	C1-C2-C3-O3
2	A	704	GOL	O1-C1-C2-C3
2	A	704	GOL	C1-C2-C3-O3
2	B	704	GOL	O1-C1-C2-C3

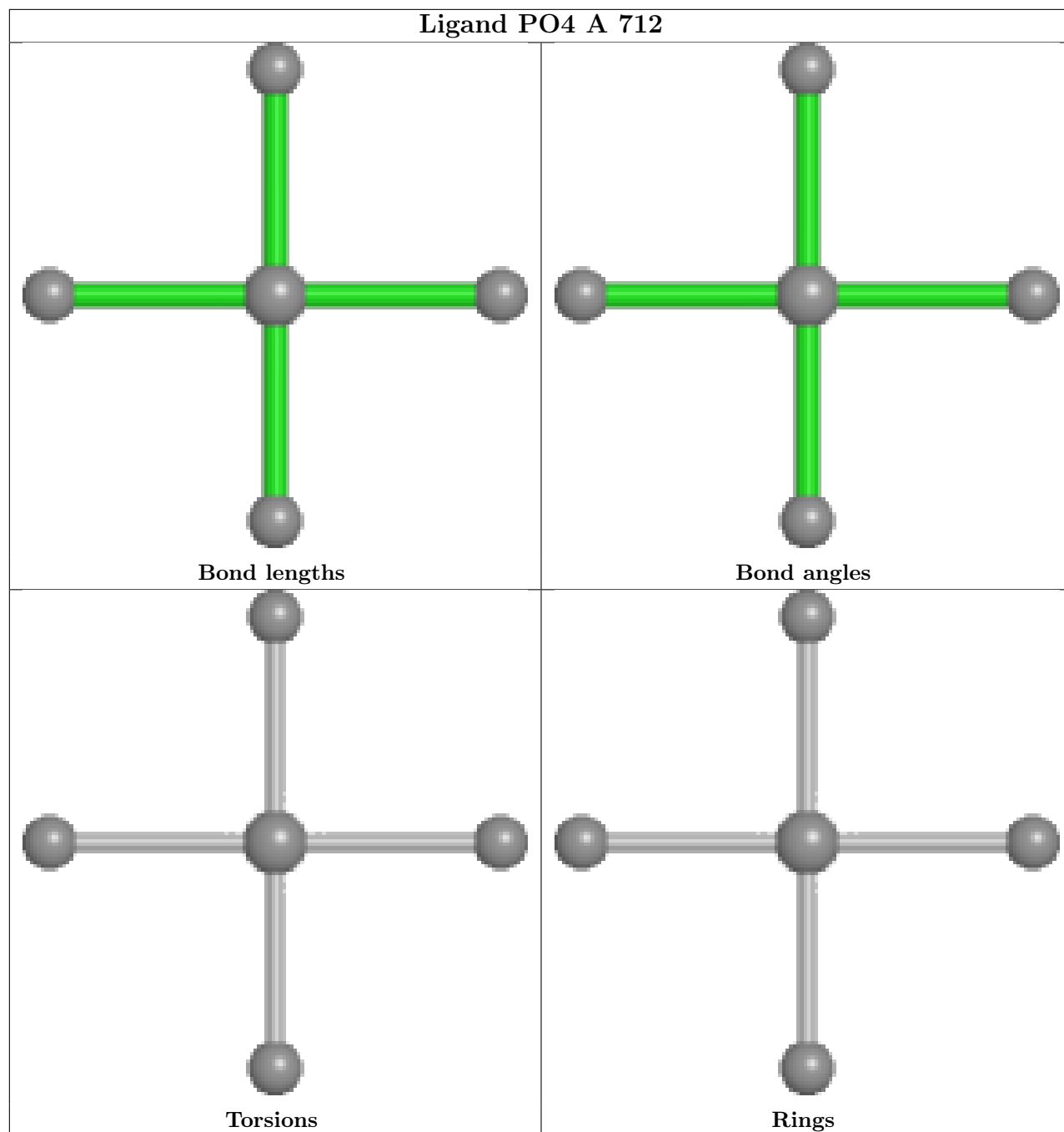
There are no ring outliers.

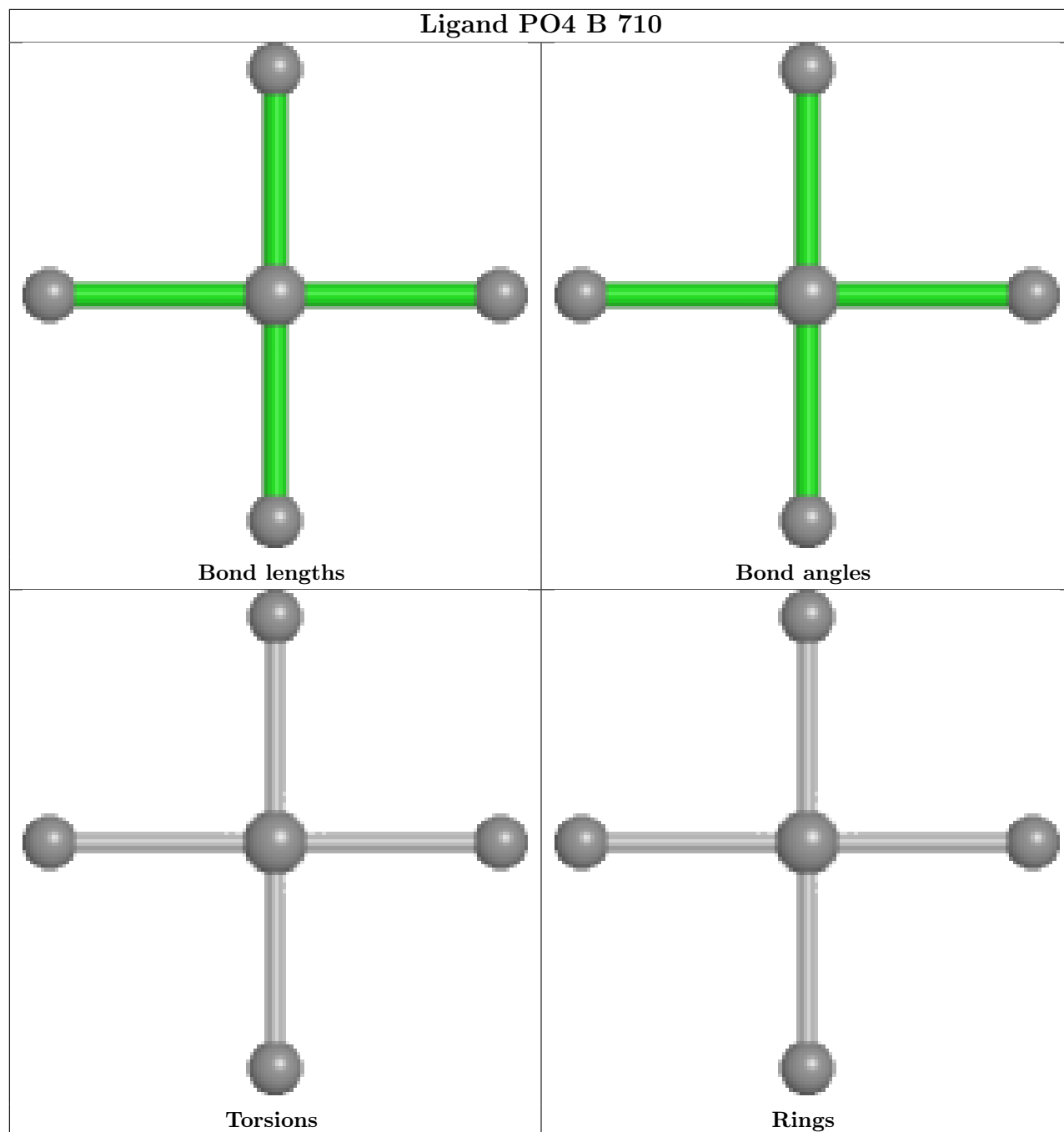
9 monomers are involved in 14 short contacts:

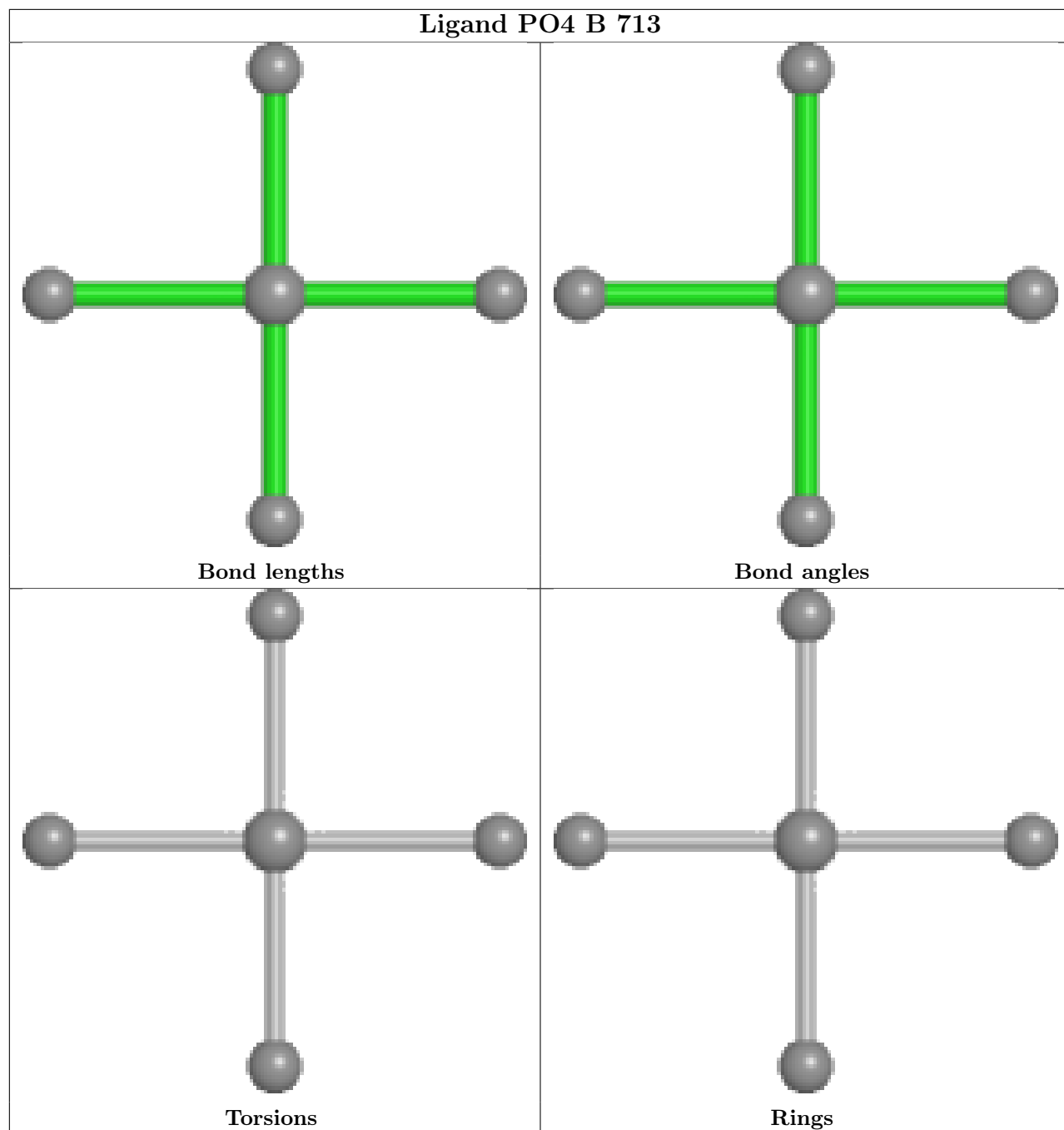
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	704	GOL	2	0
2	A	702	GOL	1	0
2	A	703	GOL	1	0
3	B	702	SO4	1	0
7	A	717	PEG	1	0
4	A	707	PO4	1	0
4	A	709	PO4	3	0
3	B	708	SO4	1	0
2	B	705	GOL	3	0

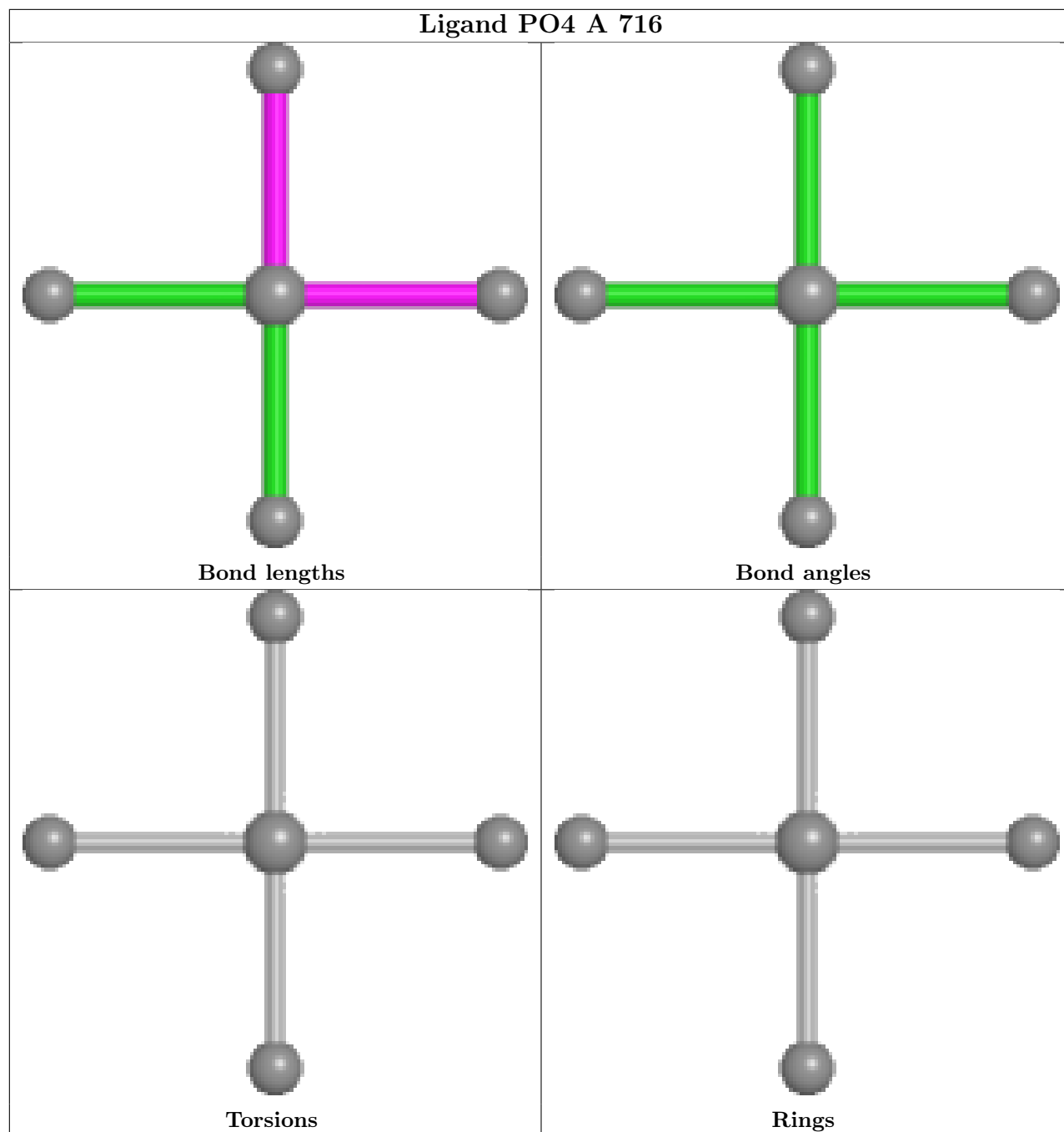
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

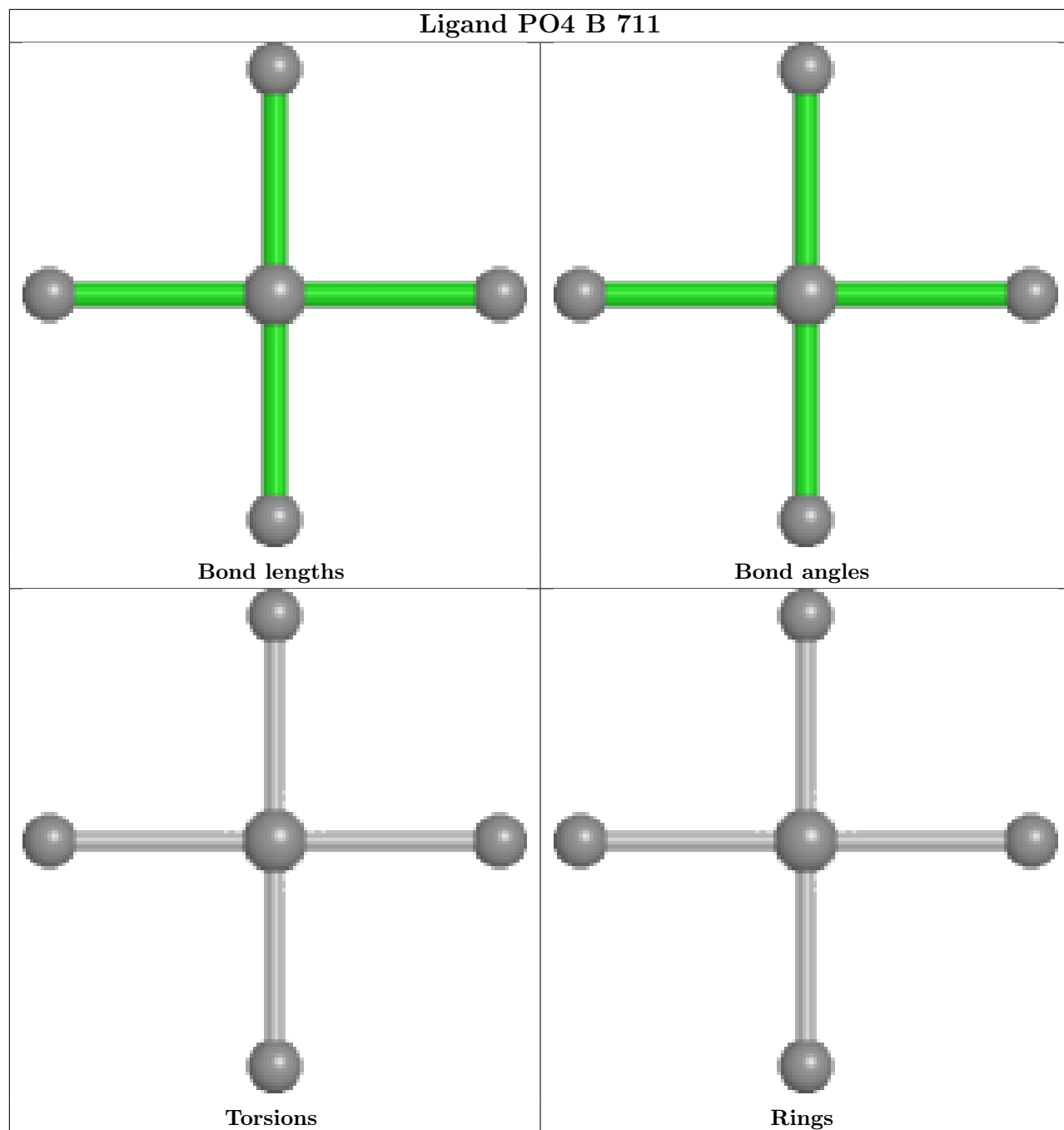


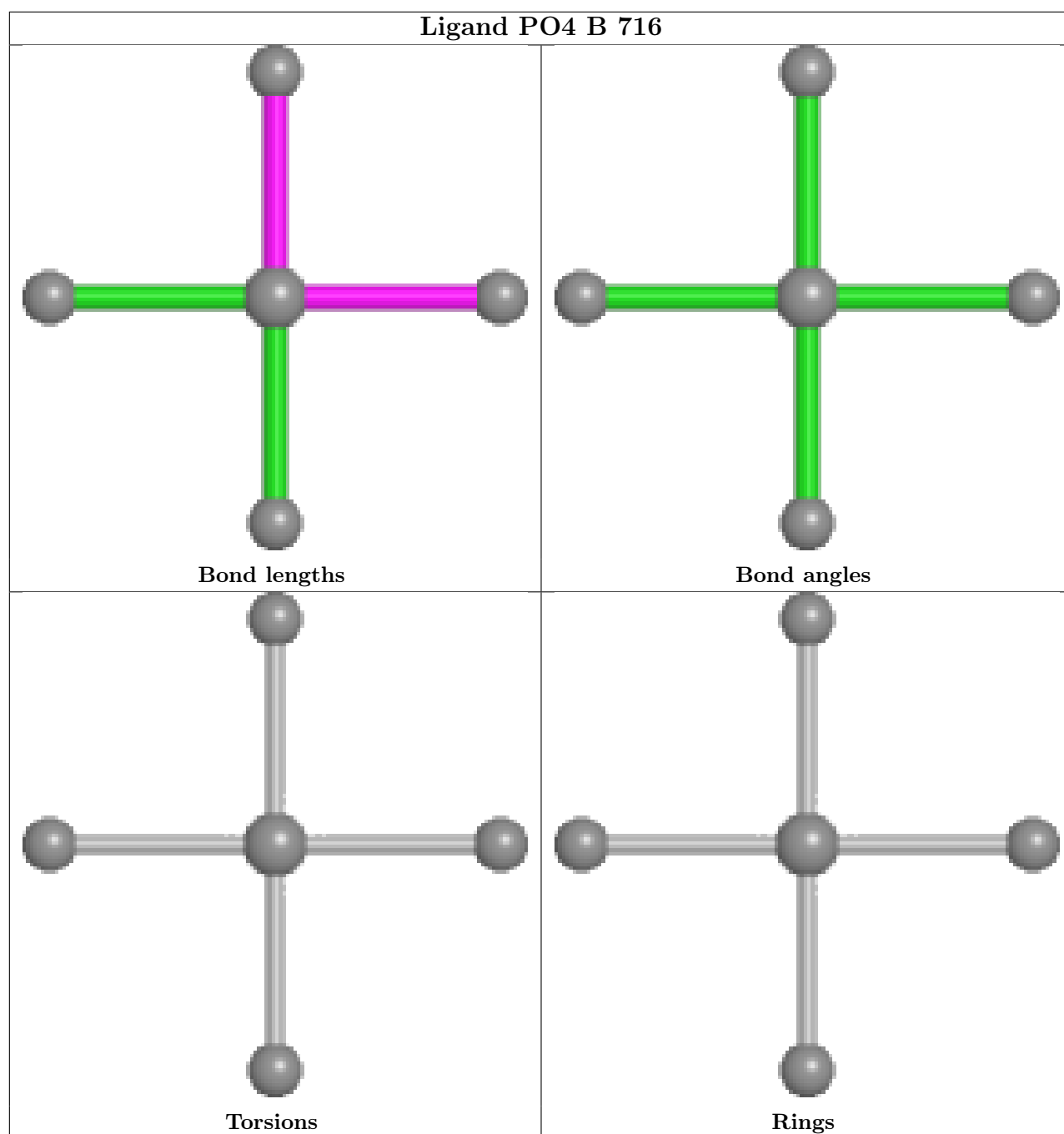


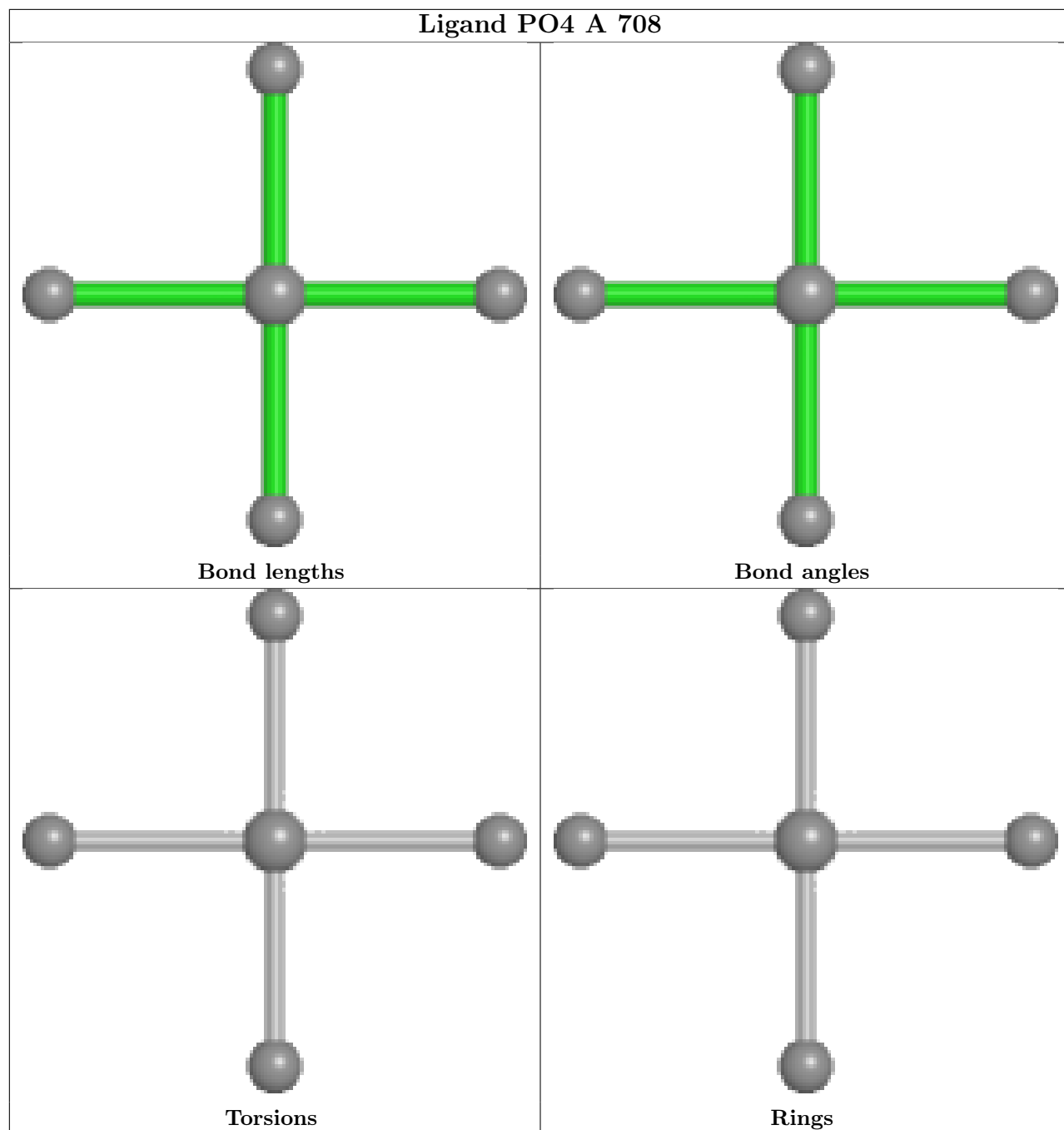


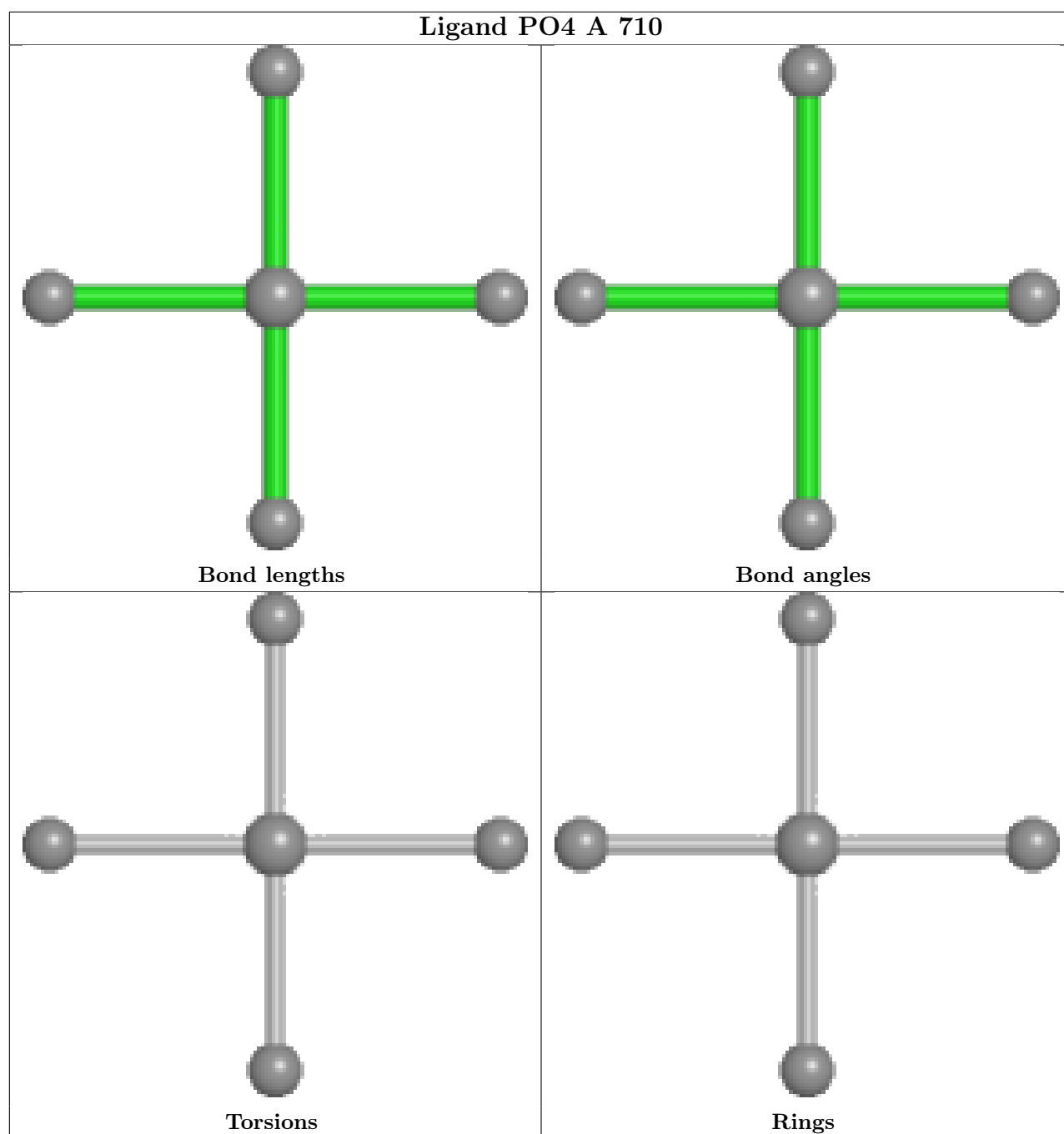


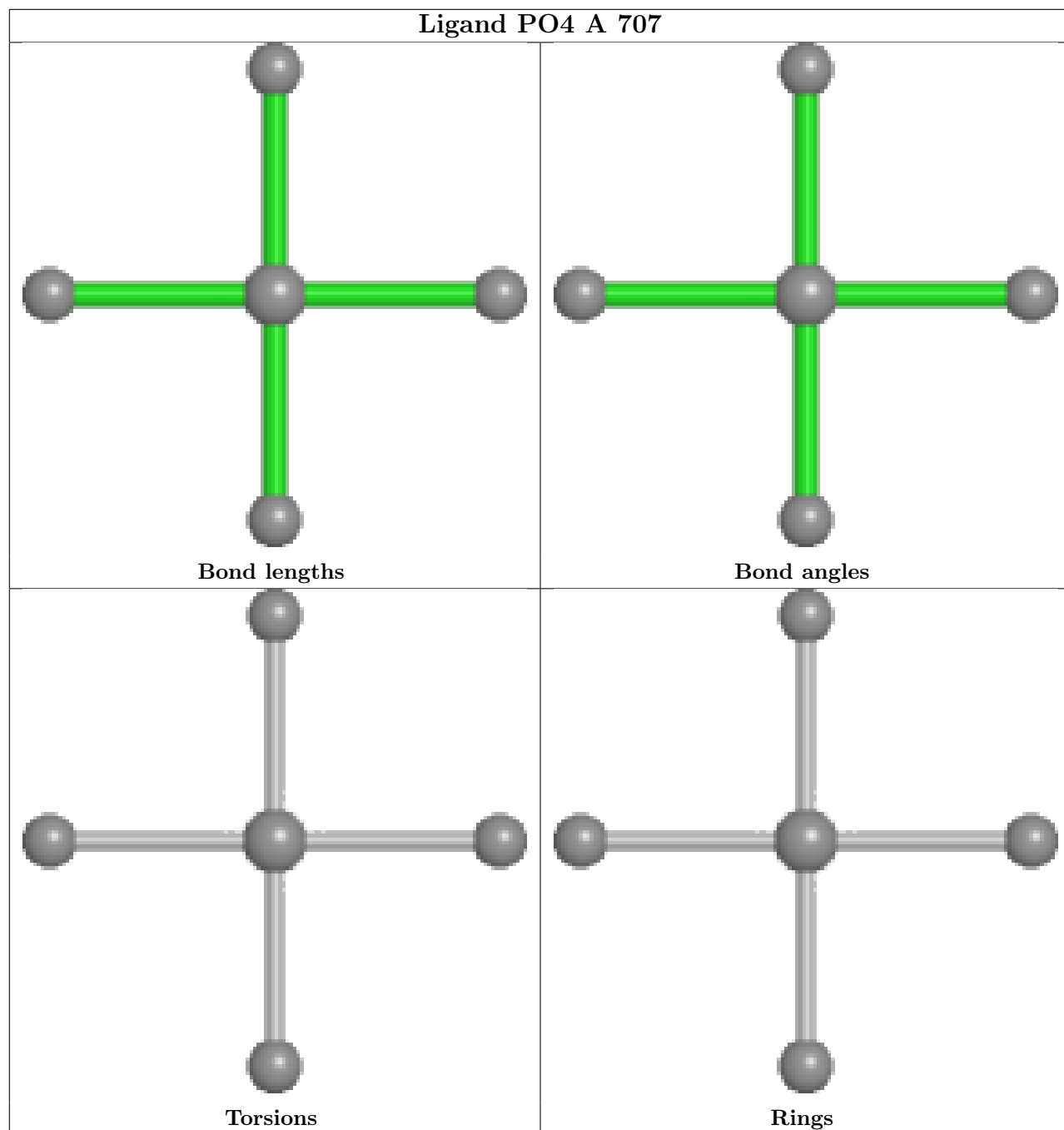


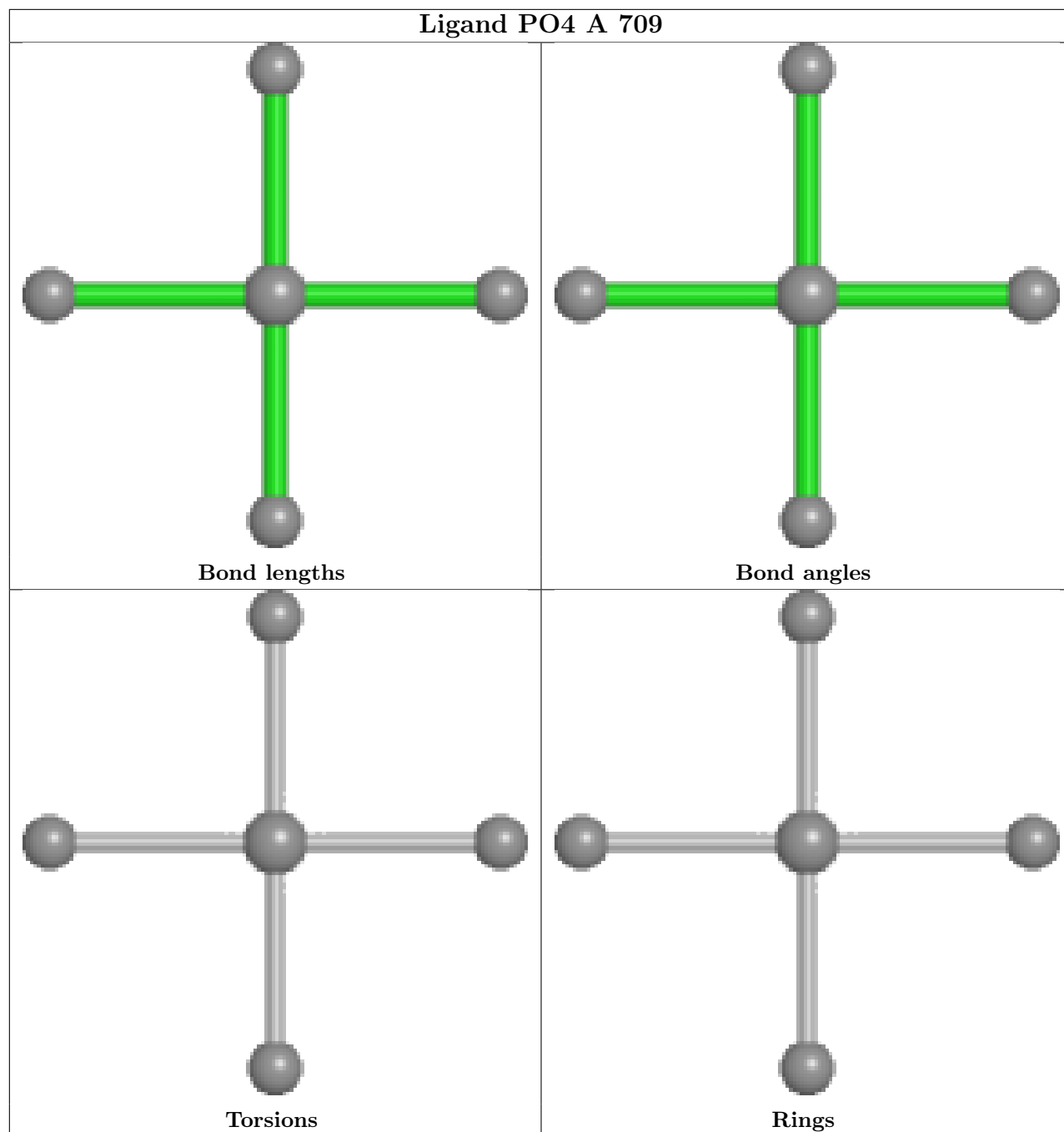


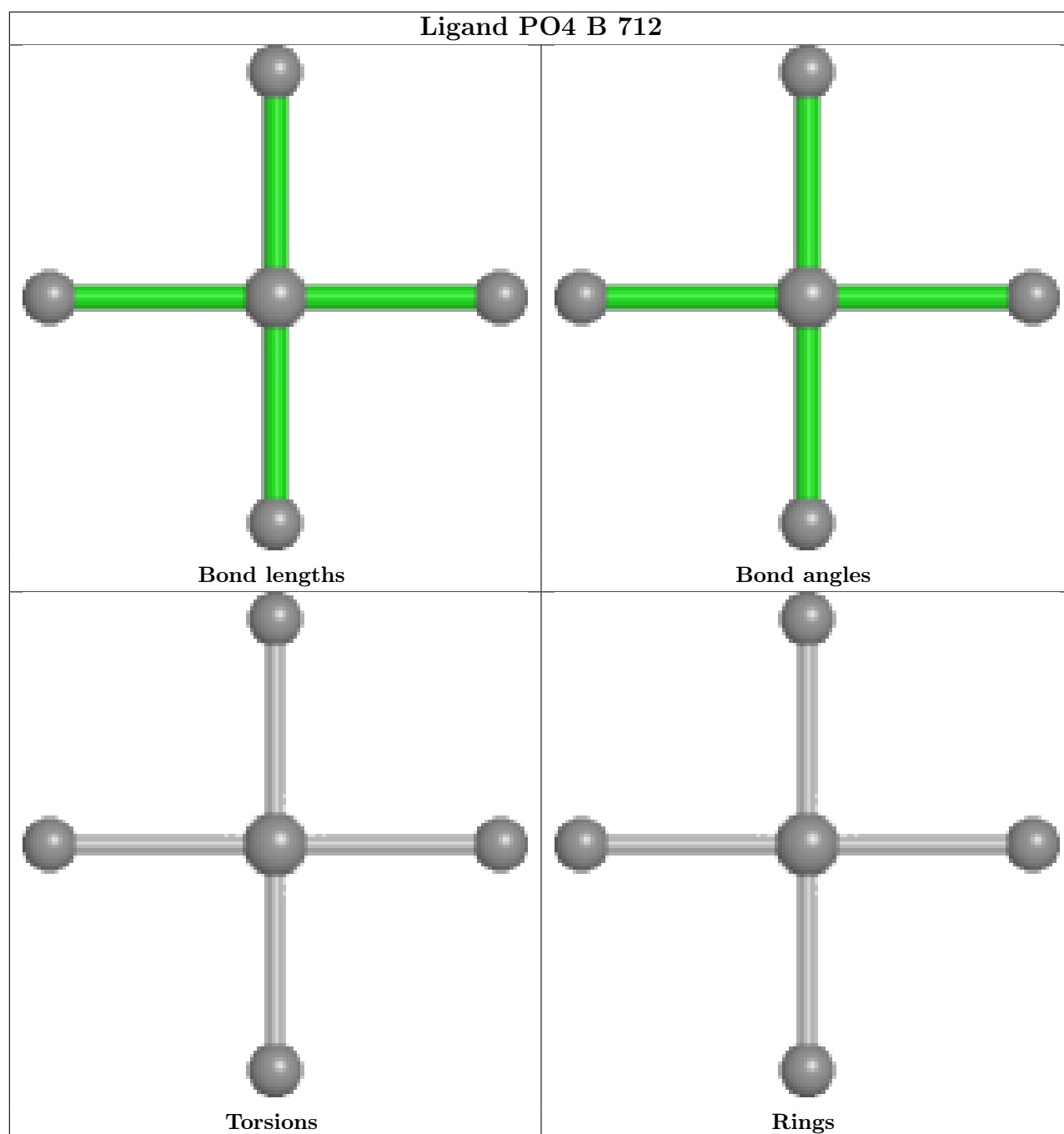


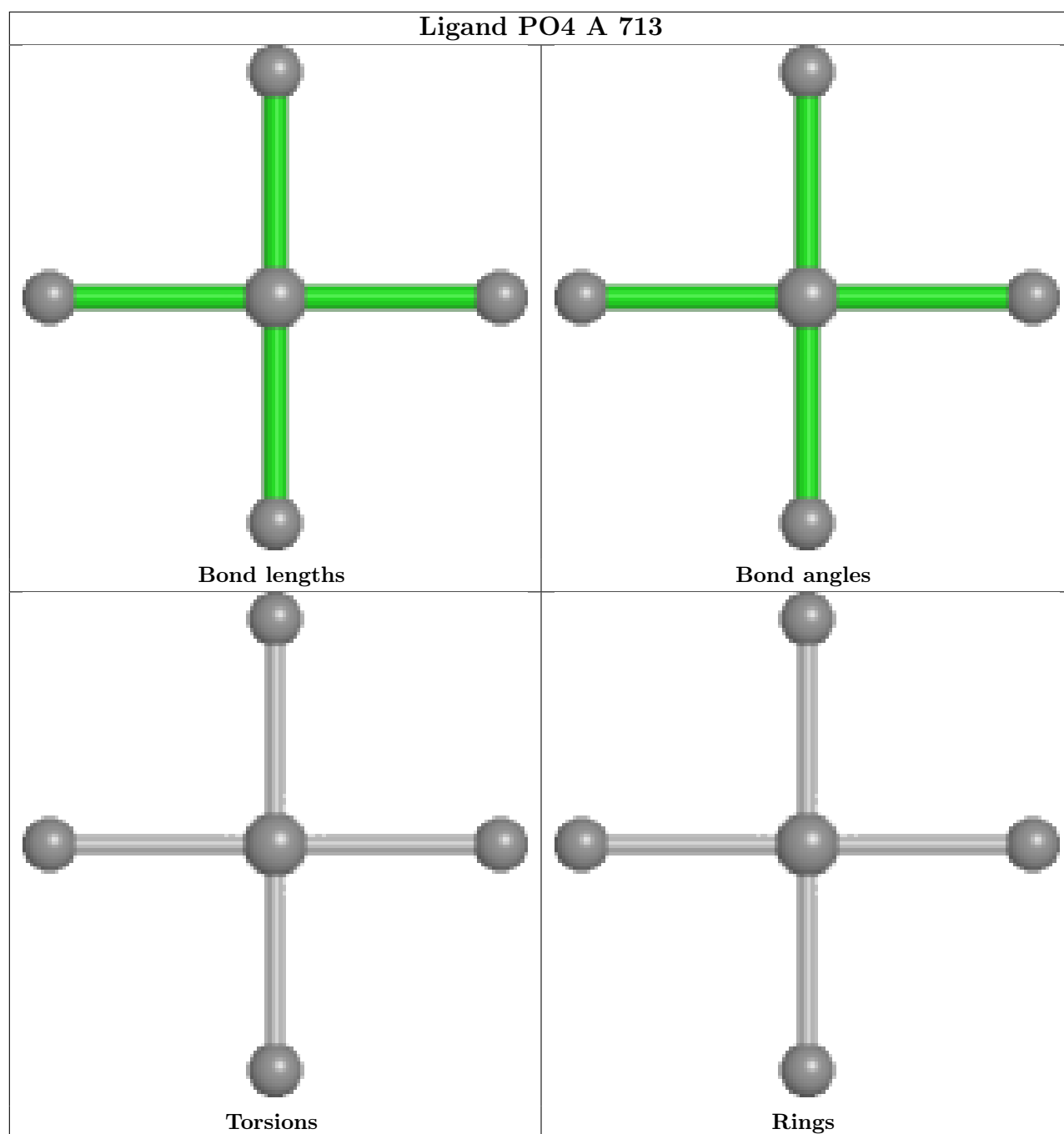


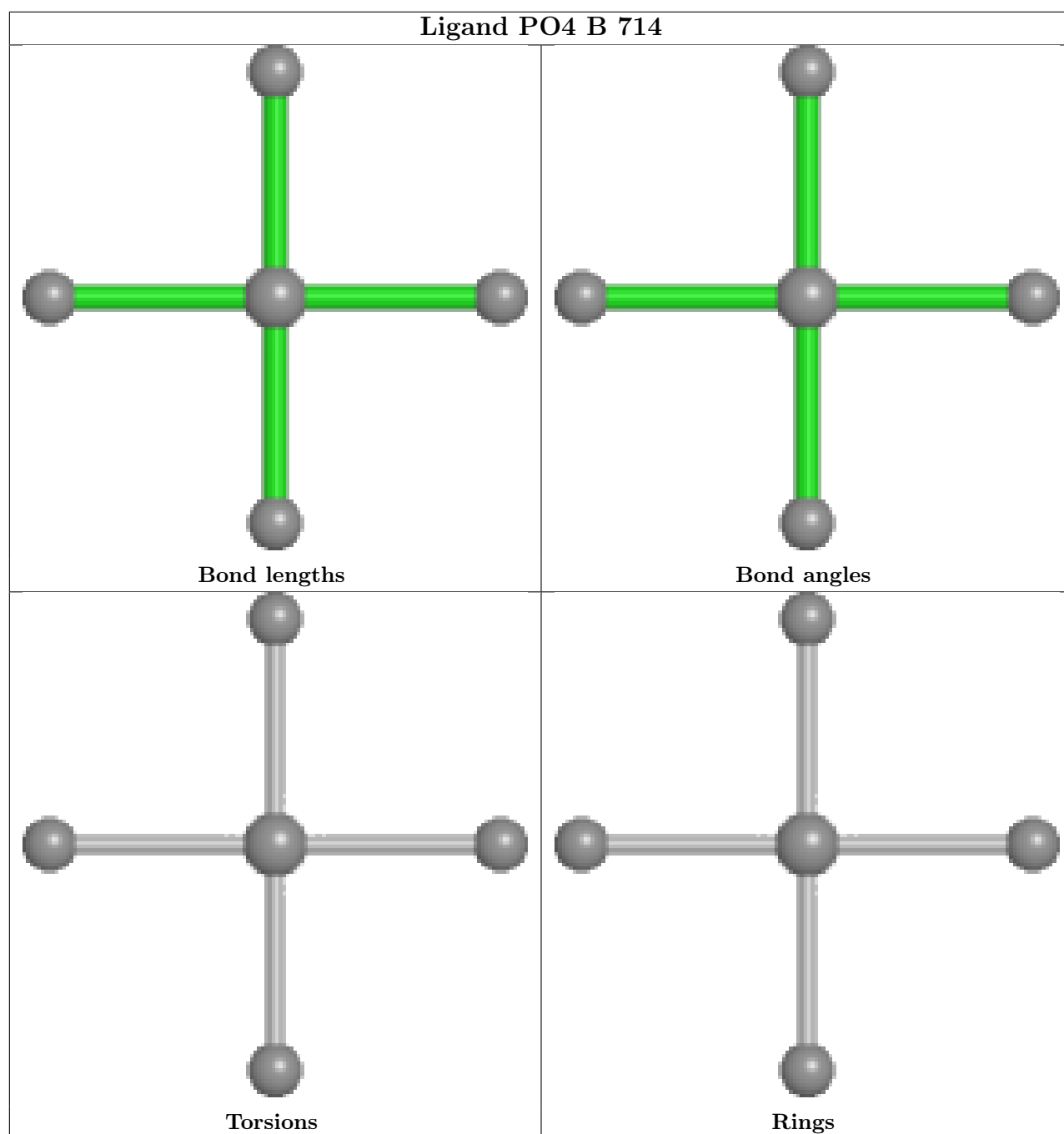


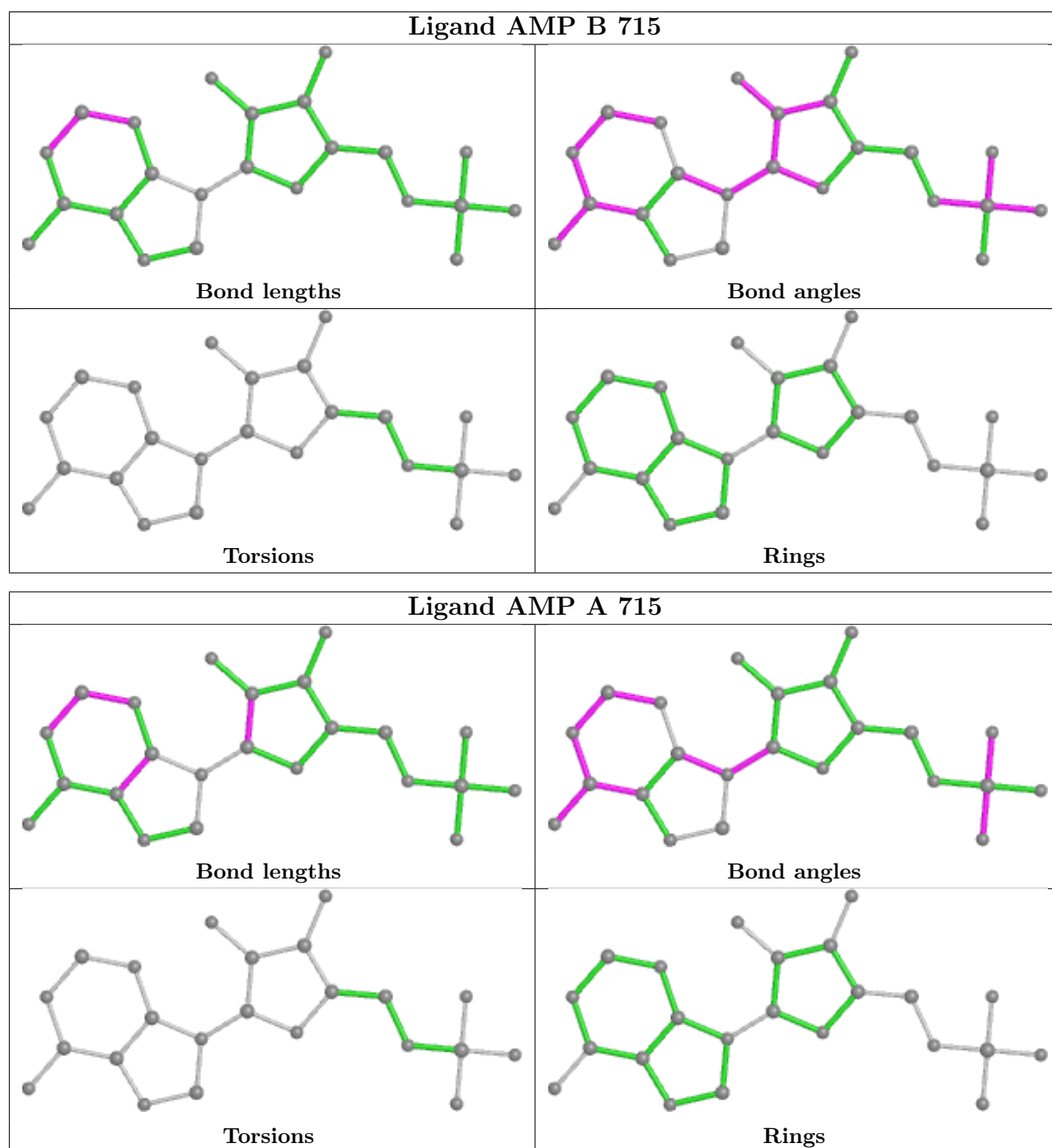












5.7 Other polymers [i](#)

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 [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	606/606 (100%)	0.48	81 (13%) 3 2	8, 20, 114, 188	0
1	B	606/606 (100%)	0.43	80 (13%) 3 2	8, 21, 105, 182	0
All	All	1212/1212 (100%)	0.46	161 (13%) 3 2	8, 21, 112, 188	0

The worst 5 of 161 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	579	GLY	18.9
1	B	487	PRO	18.9
1	A	578	LEU	17.5
1	B	515	GLY	15.4
1	A	485	PRO	14.6

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 monosaccharides in this entry.

6.4 Ligands [i](#)

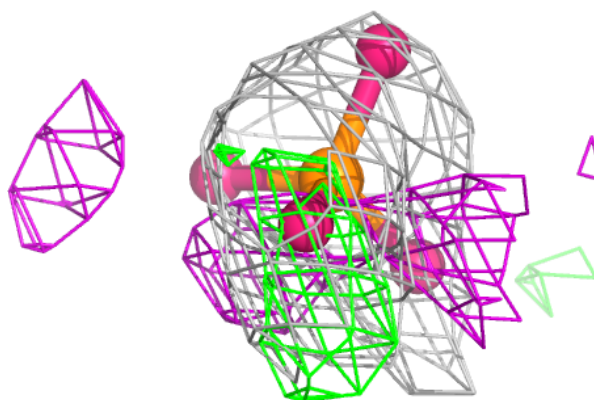
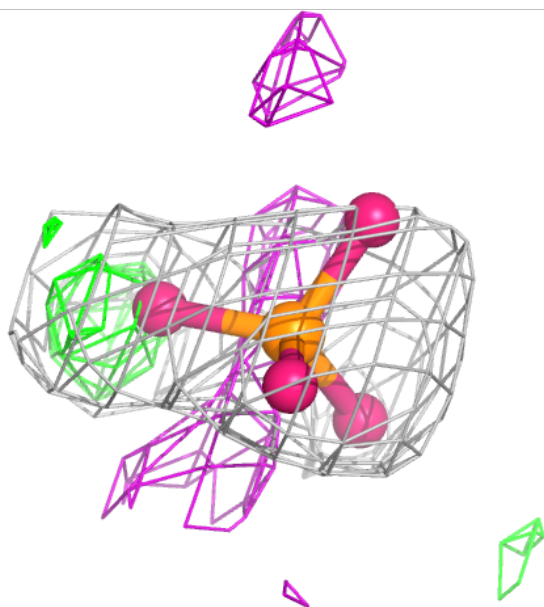
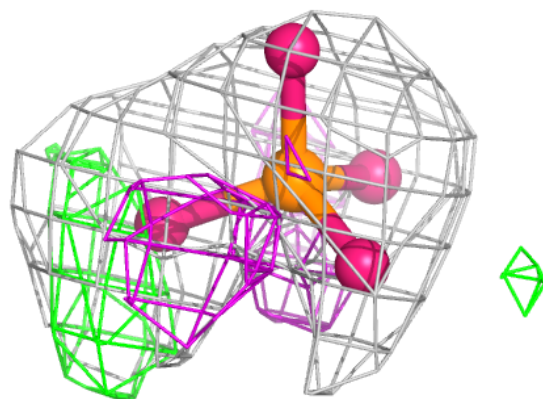
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q < 0.9’ lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
4	PO4	A	712	5/5	0.65	0.27	57,67,76,79	0
2	GOL	A	701	6/6	0.71	0.22	57,60,63,69	0
2	GOL	A	703	6/6	0.72	0.21	53,54,57,58	0
2	GOL	A	702	6/6	0.73	0.20	38,40,45,47	0
2	GOL	A	704	6/6	0.75	0.21	48,52,53,55	0
2	GOL	B	706	6/6	0.79	0.24	50,56,66,73	0
7	PEG	A	717	7/7	0.82	0.15	43,49,52,53	0
4	PO4	B	713	5/5	0.86	0.26	75,75,84,84	0
2	GOL	B	701	6/6	0.87	0.18	30,40,43,45	0
2	GOL	B	704	6/6	0.89	0.14	34,34,39,44	0
2	GOL	B	705	6/6	0.90	0.16	28,32,33,42	0
3	SO4	B	707	5/5	0.91	0.16	66,70,77,81	0
4	PO4	B	714	5/5	0.91	0.17	75,77,81,81	0
4	PO4	B	710	5/5	0.91	0.10	42,44,51,59	0
4	PO4	B	712	5/5	0.92	0.14	50,54,60,63	0
4	PO4	A	707	5/5	0.92	0.10	41,45,48,59	0
3	SO4	A	705	5/5	0.93	0.23	47,51,57,64	0
3	SO4	A	706	5/5	0.93	0.18	62,62,68,75	0
4	PO4	A	713	5/5	0.94	0.15	42,42,46,53	0
4	PO4	A	708	5/5	0.95	0.09	41,41,46,49	0
4	PO4	A	709	5/5	0.95	0.15	54,55,58,60	0
3	SO4	B	709	5/5	0.95	0.18	35,46,50,56	0
4	PO4	A	710	5/5	0.96	0.15	52,56,59,66	0
4	PO4	A	711	5/5	0.97	0.13	33,37,40,41	0
4	PO4	B	711	5/5	0.97	0.13	41,43,51,54	0
6	AMP	B	715	23/23	0.98	0.10	11,12,14,16	0
3	SO4	B	708	5/5	0.98	0.06	40,43,47,47	0
4	PO4	B	716	5/5	0.99	0.10	13,13,16,17	0
5	MG	A	714	1/1	0.99	0.06	23,23,23,23	0
5	MG	B	703	1/1	0.99	0.05	22,22,22,22	0
6	AMP	A	715	23/23	0.99	0.09	11,13,13,14	0
4	PO4	A	716	5/5	0.99	0.11	13,15,17,19	0
3	SO4	B	702	5/5	0.99	0.08	36,37,38,39	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.

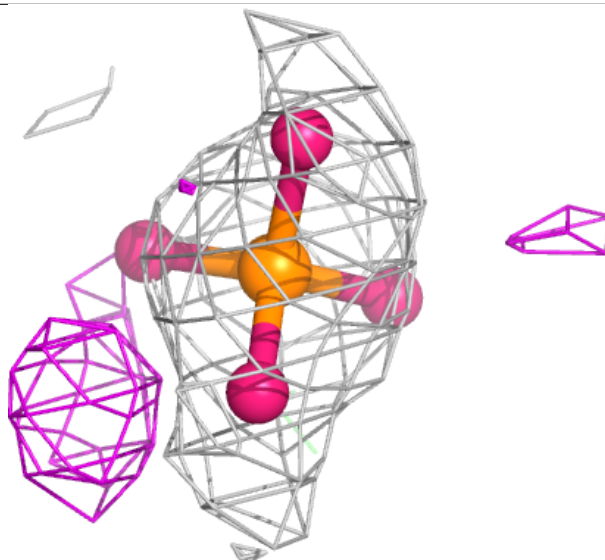
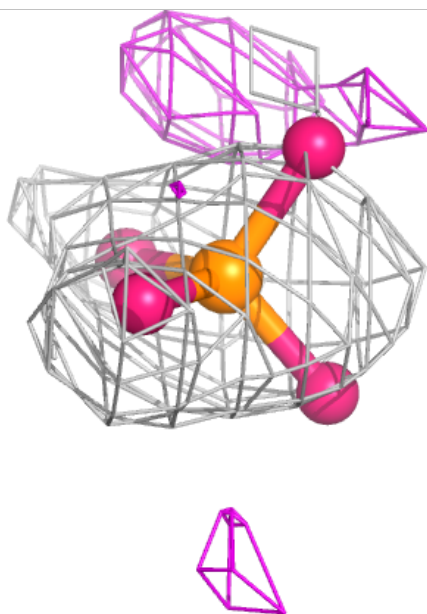
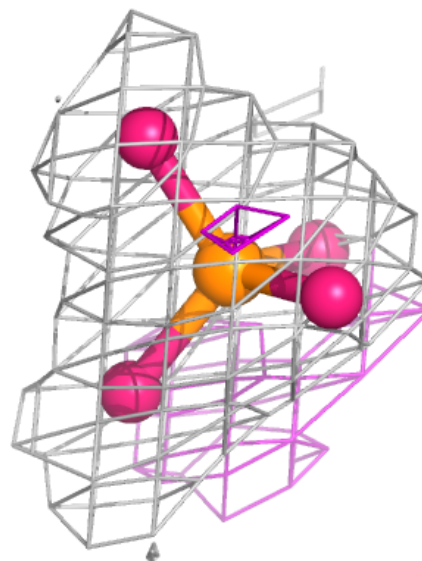
Electron density around PO4 A 712:

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



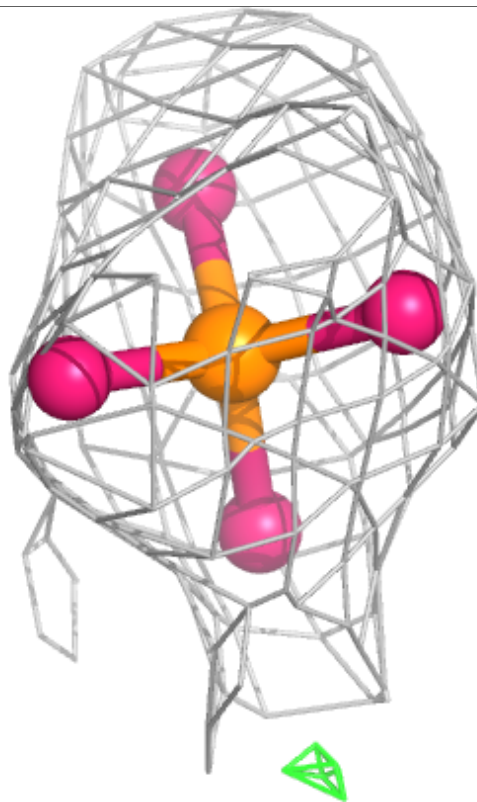
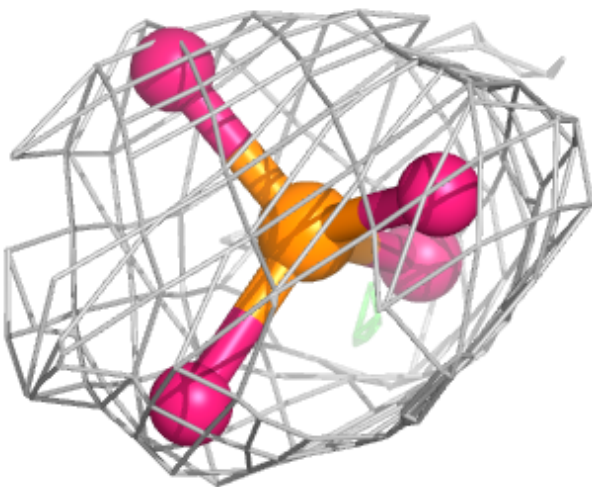
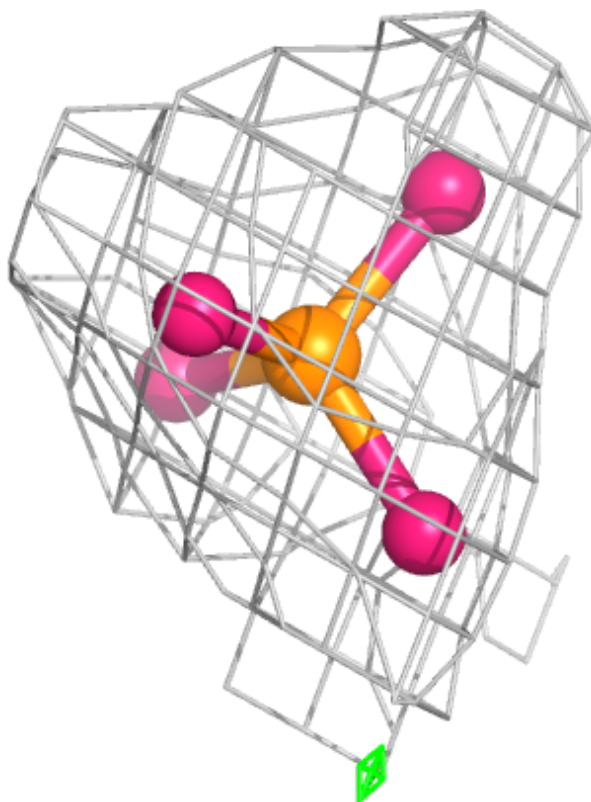
Electron density around PO4 B 713:

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



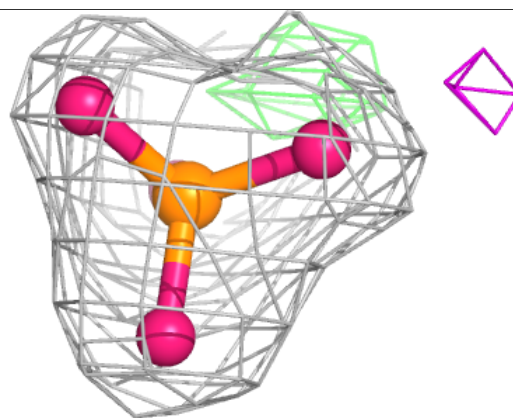
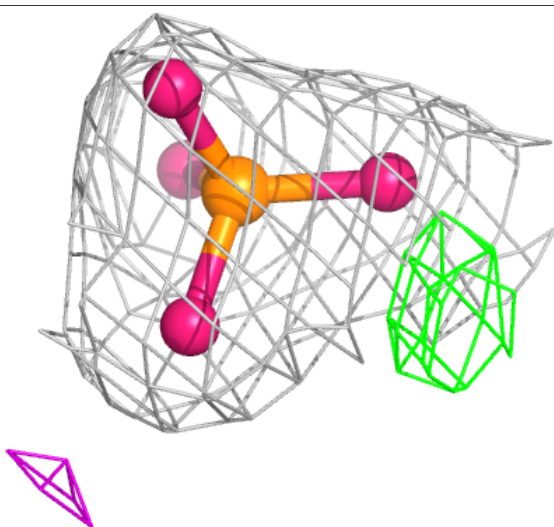
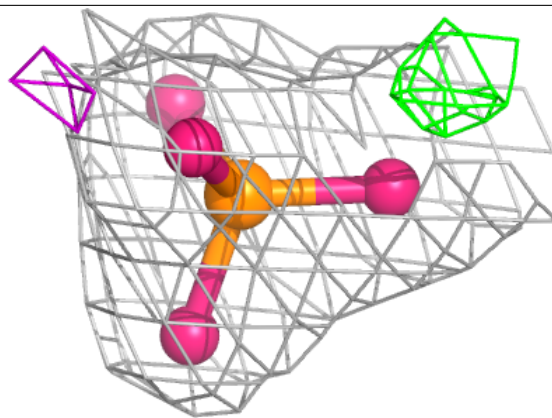
Electron density around PO4 B 714:

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



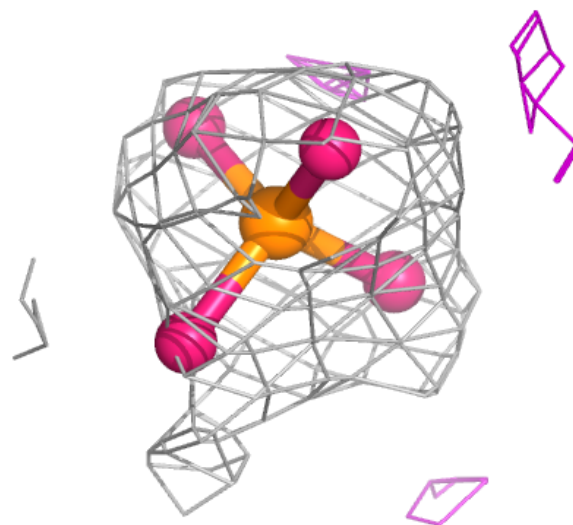
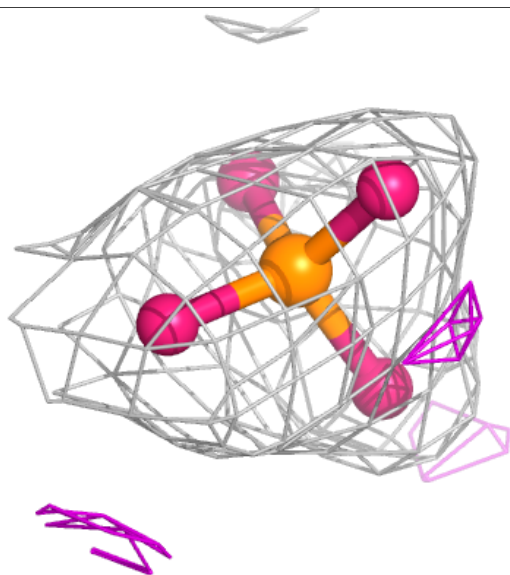
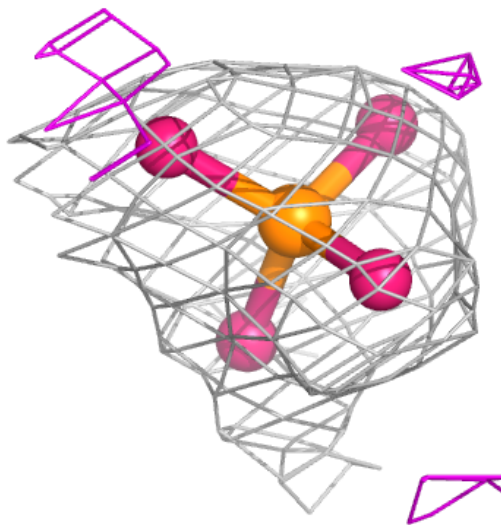
Electron density around PO4 B 710:

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



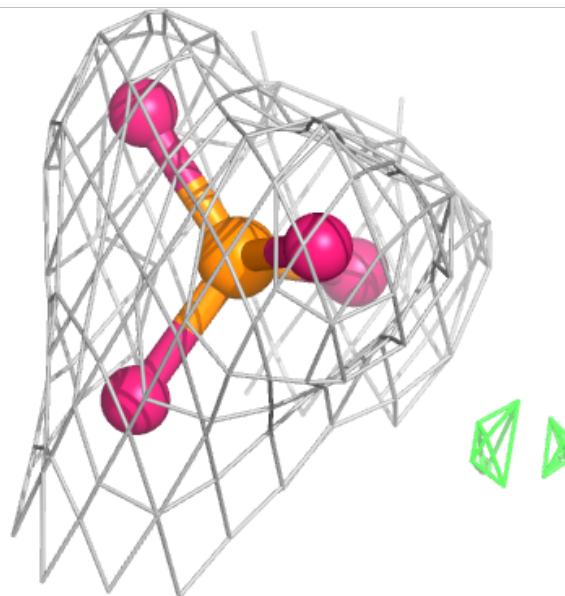
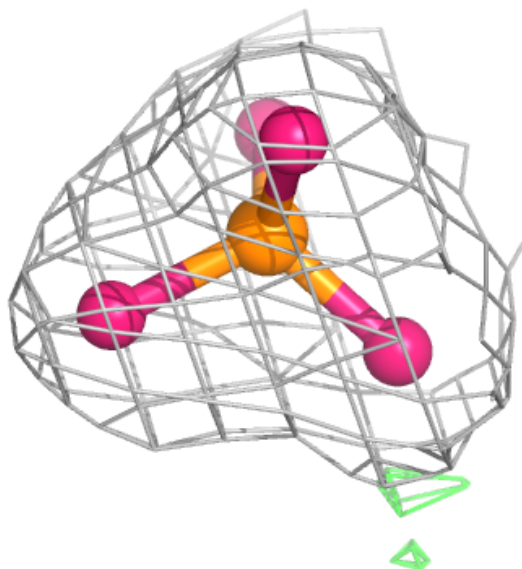
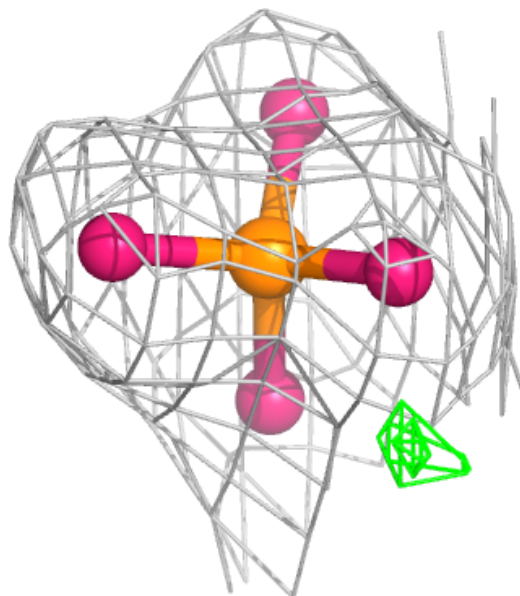
Electron density around PO4 B 712:

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



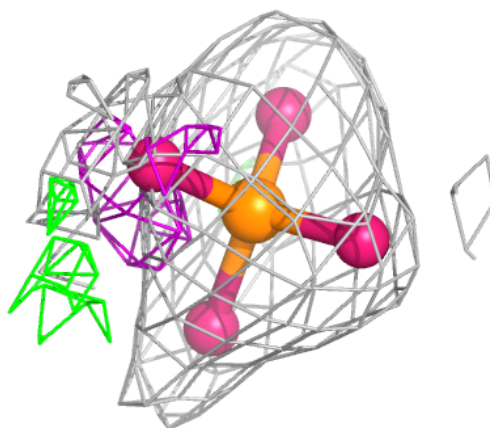
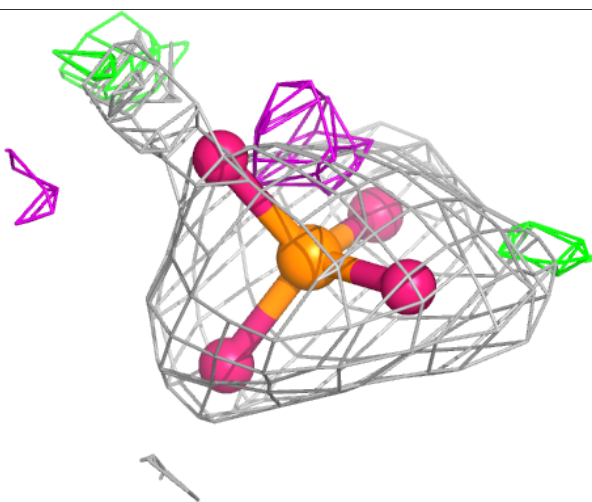
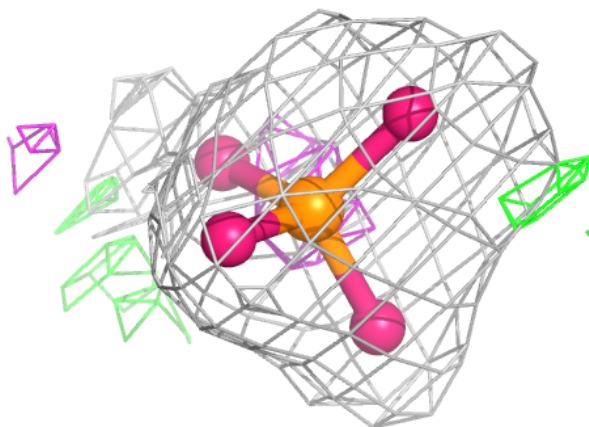
Electron density around PO4 A 707:

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



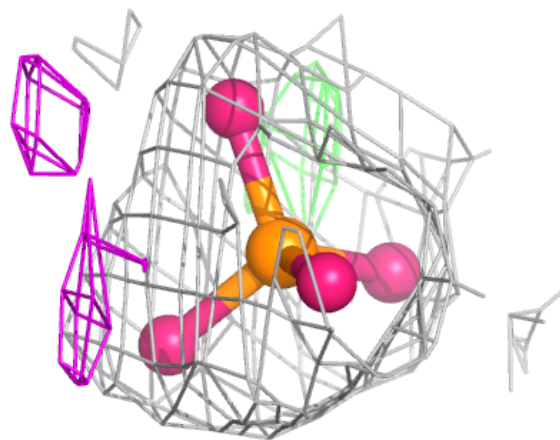
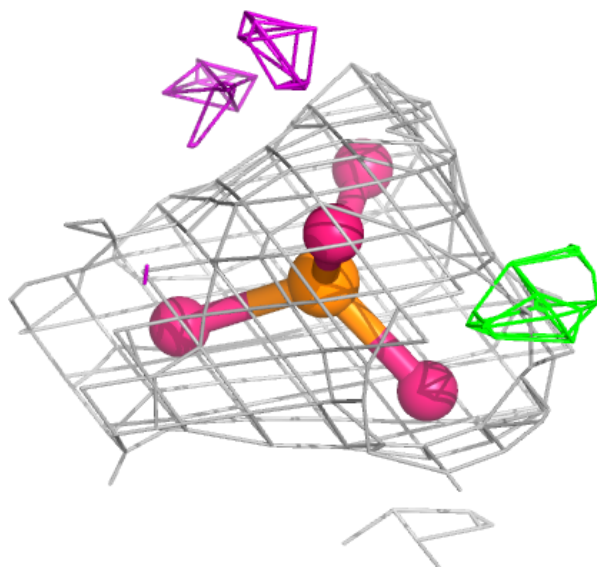
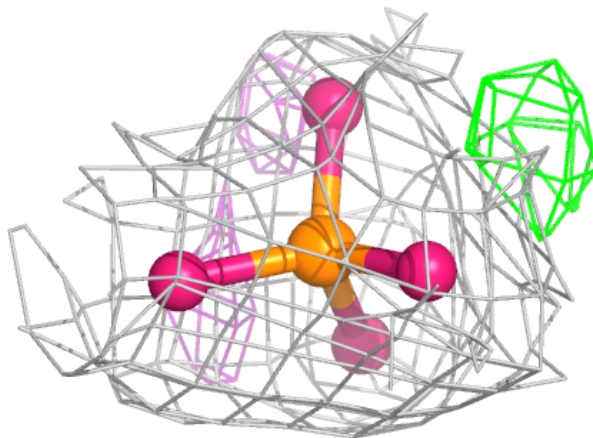
Electron density around PO4 A 713:

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



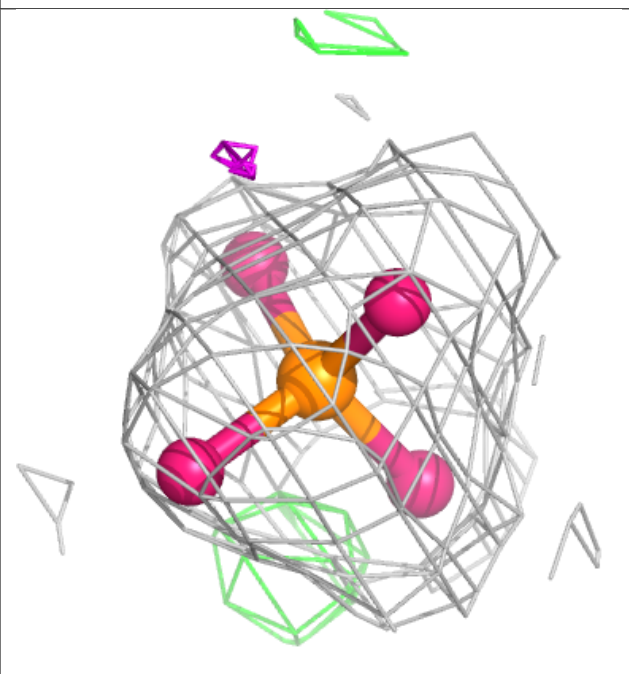
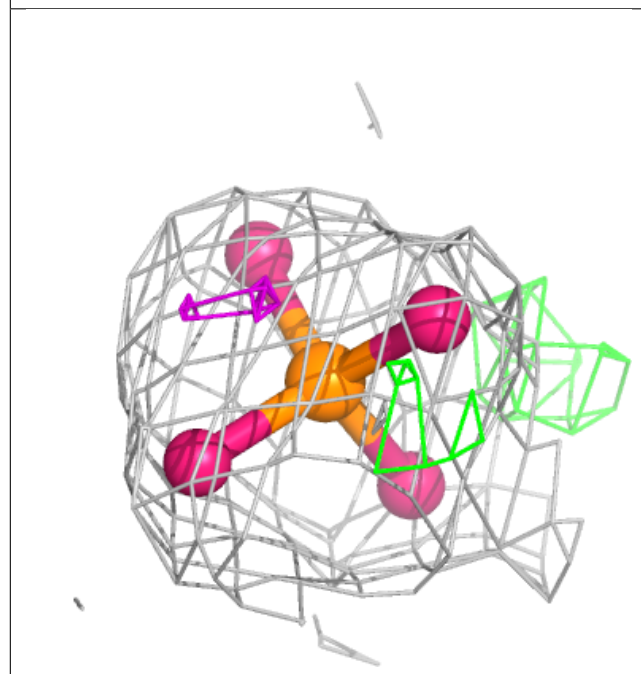
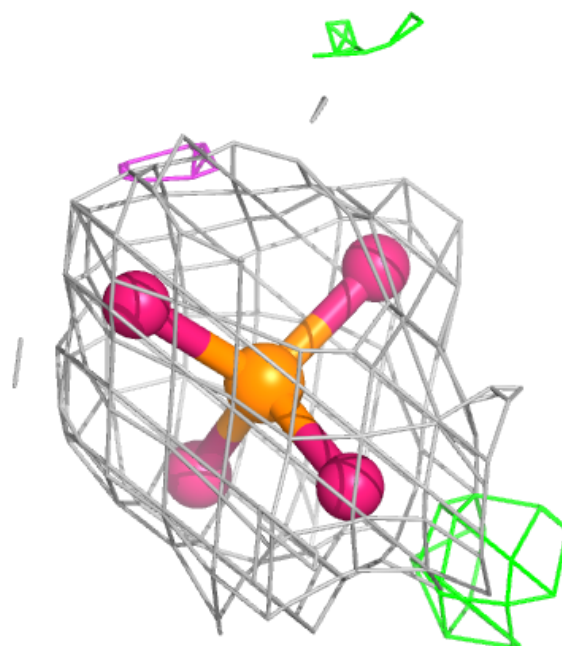
Electron density around PO4 A 708:

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



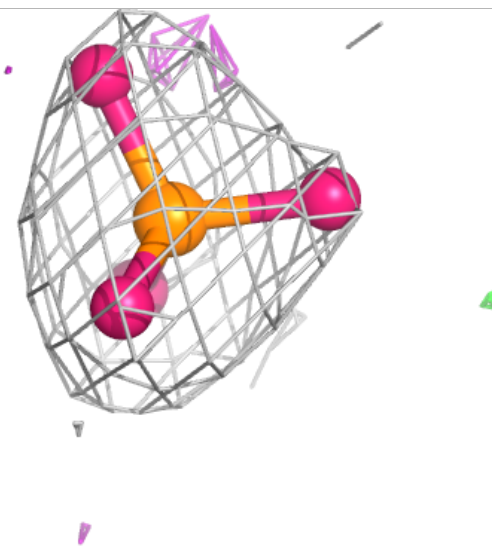
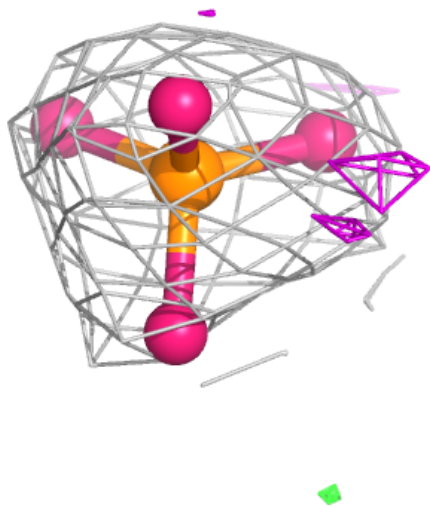
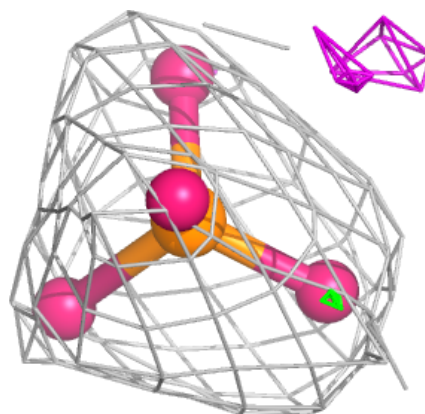
Electron density around PO4 A 709:

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



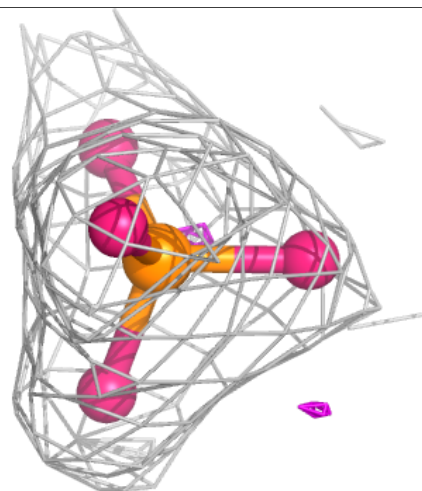
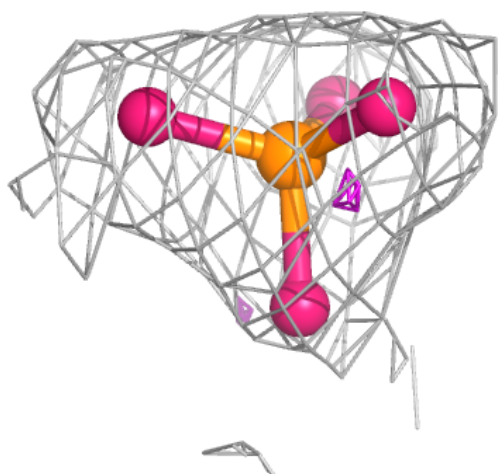
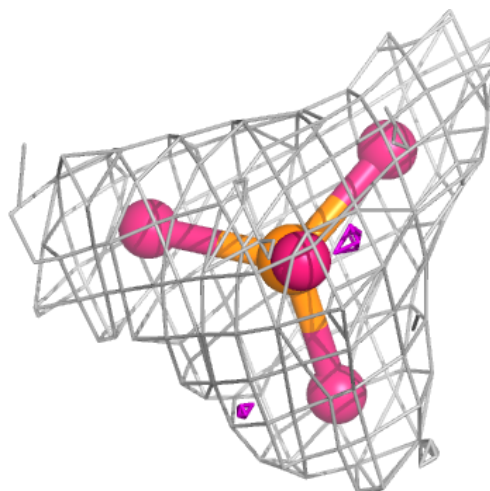
Electron density around PO4 A 710:

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



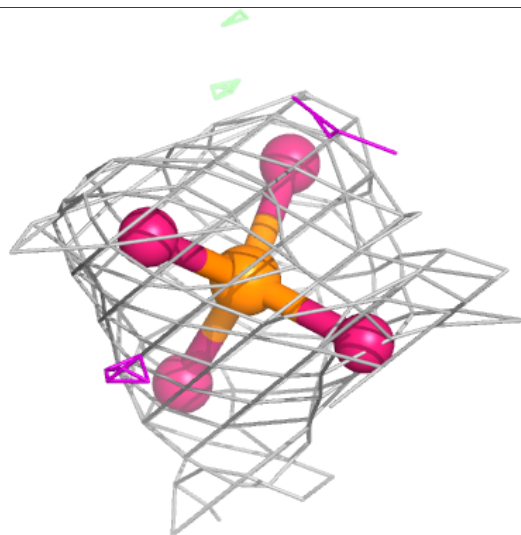
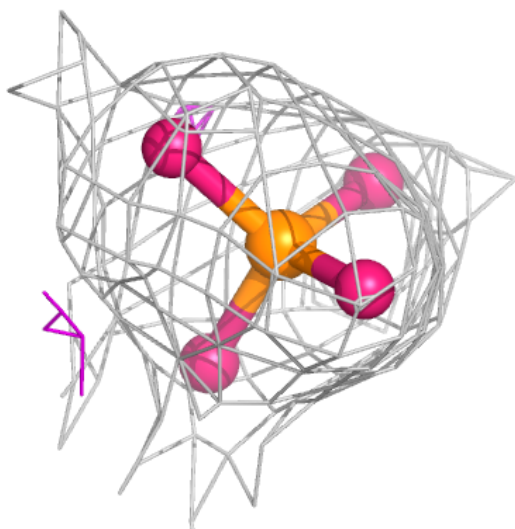
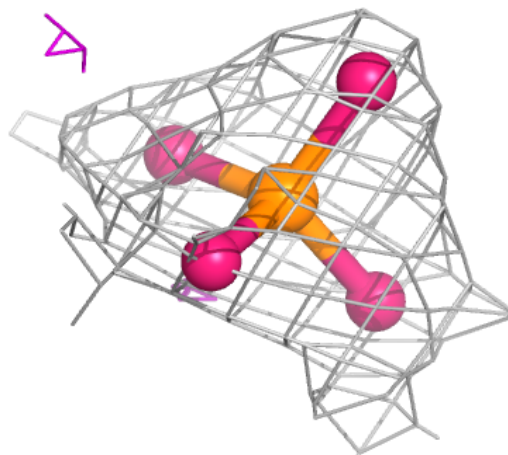
Electron density around PO4 A 711:

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



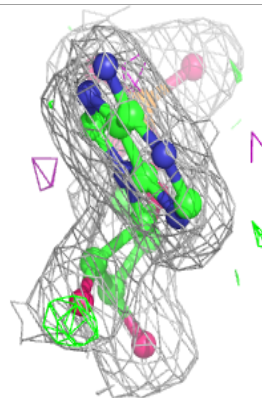
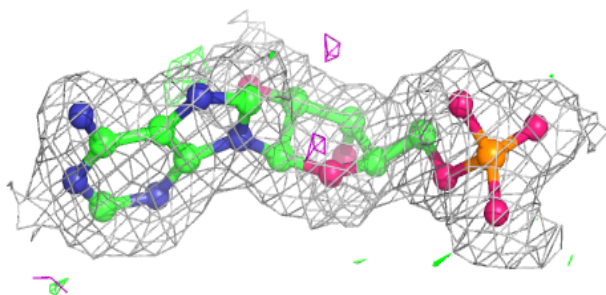
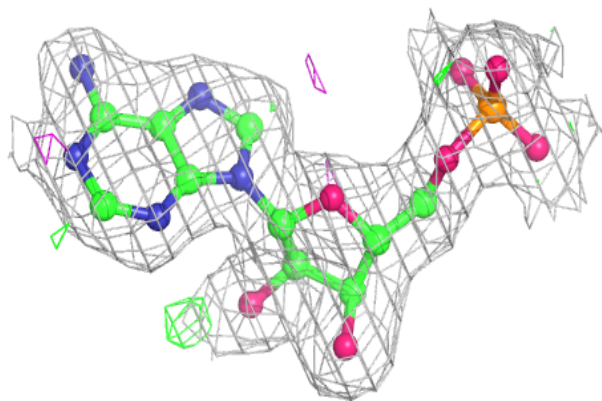
Electron density around PO4 B 711:

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



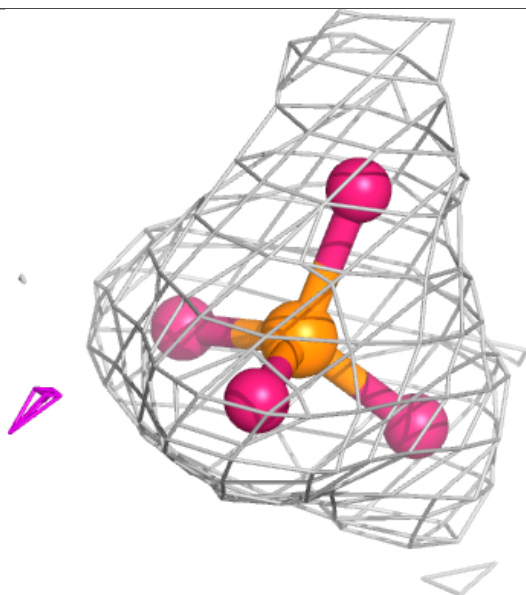
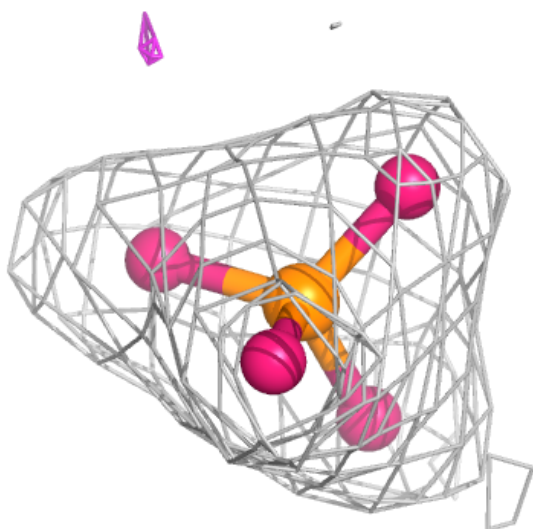
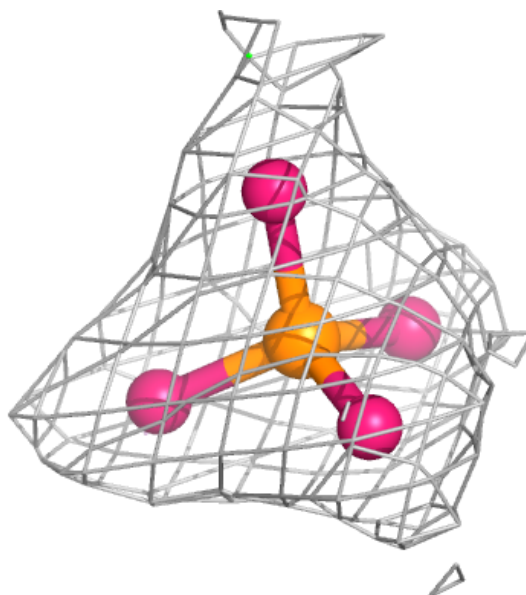
Electron density around AMP B 715:

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



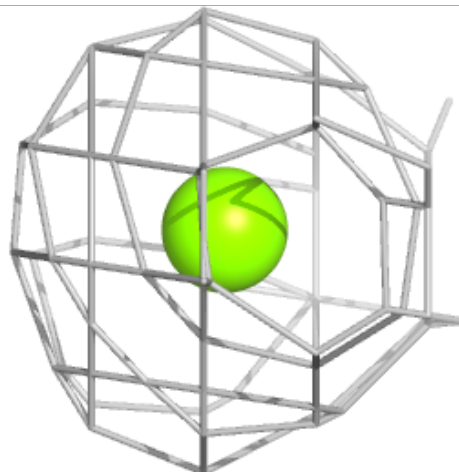
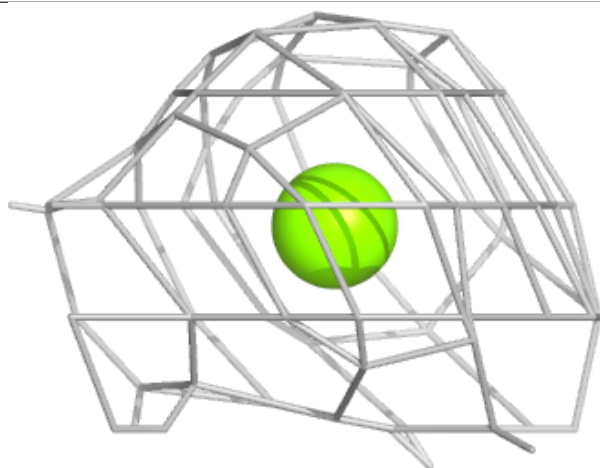
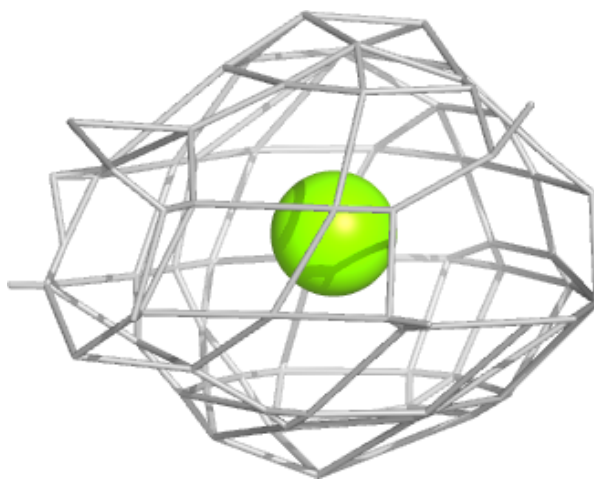
Electron density around PO4 B 716:

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



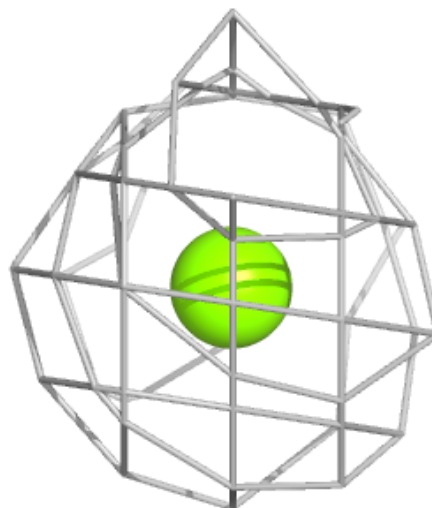
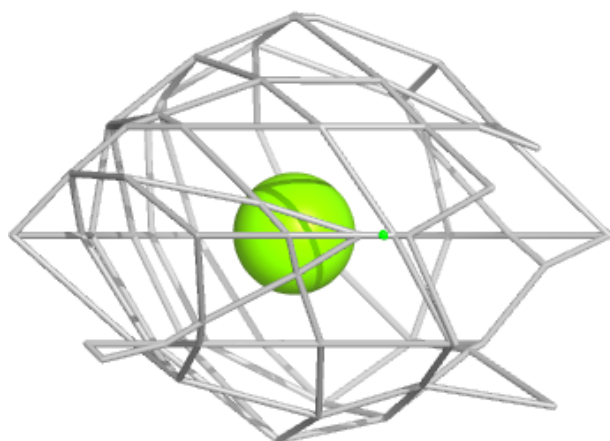
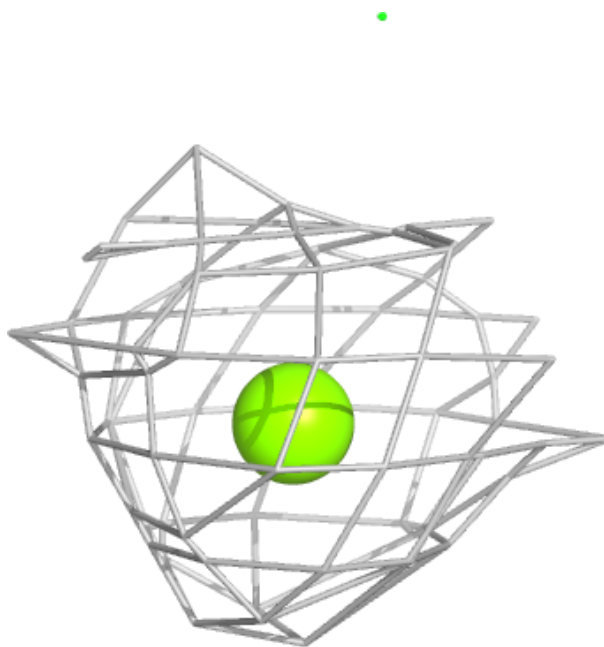
Electron density around MG A 714:

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



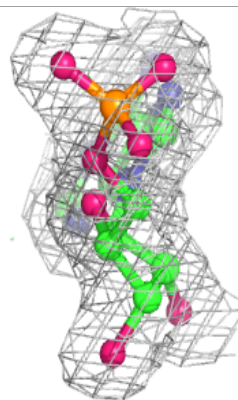
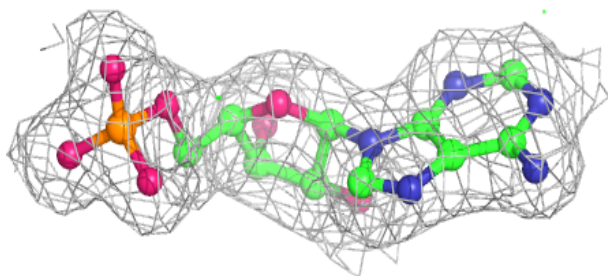
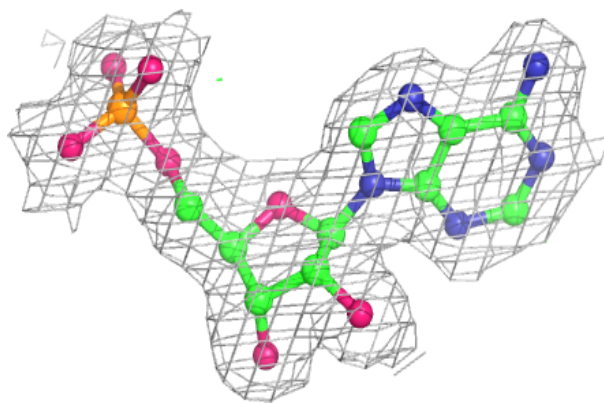
Electron density around MG B 703:

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



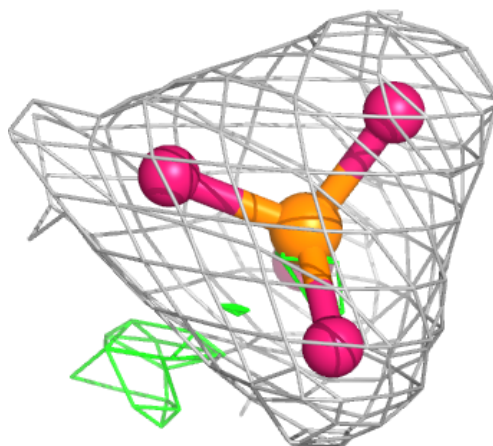
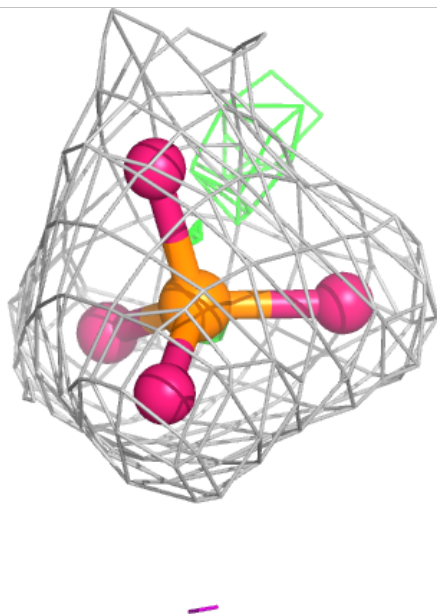
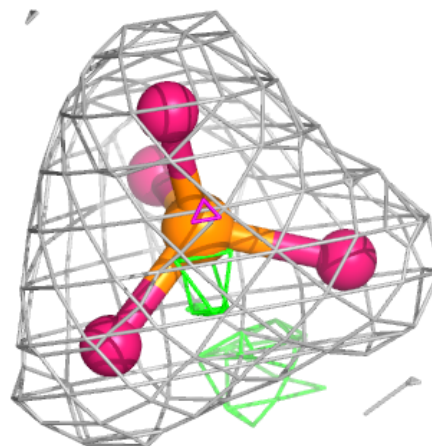
Electron density around AMP A 715:

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



Electron density around PO4 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)



6.5 Other polymers ⓘ

There are no such residues in this entry.