



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 03:38 PM UTC

PDB ID : 6K8V / pdb_00006k8v
Title : Crystal structure of N-domain of bacterial malonyl-CoA reductase
Authors : Kim, S.; Kim, K.-J.
Deposited on : 2019-06-13
Resolution : 2.20 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

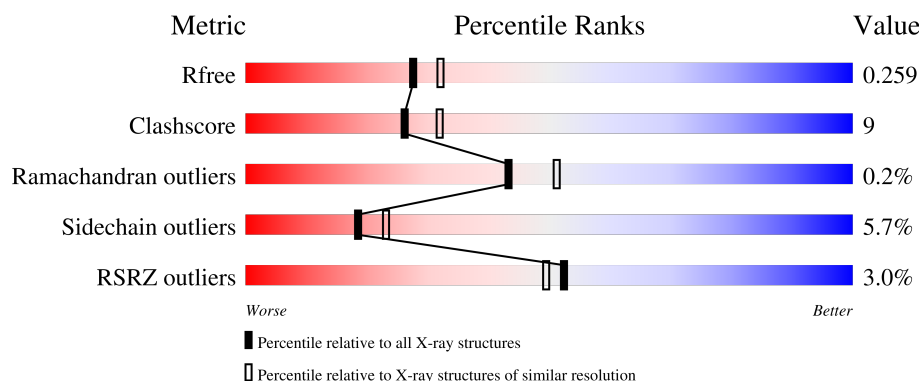
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

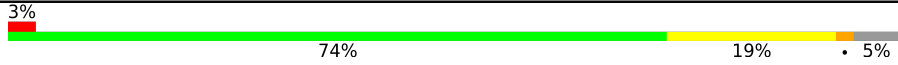
The reported resolution of this entry is 2.20 Å.

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	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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	575	
1	B	575	

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
2	SO4	B	1304	-	-	X	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 8556 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.

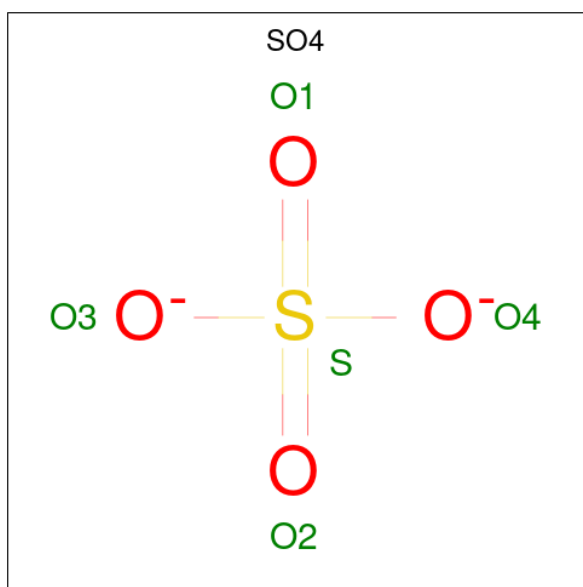
- Molecule 1 is a protein called NAD-dependent epimerase/dehydratase:Short-chain dehydrogenase/reductase SDR.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	545	Total	C	N	O	S	0	0	0
			4074	2541	733	784	16			
1	B	545	Total	C	N	O	S	0	1	0
			4082	2545	734	787	16			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	568	LEU	-	expression tag	UNP A0A1A7BFR5
A	569	GLU	-	expression tag	UNP A0A1A7BFR5
A	570	HIS	-	expression tag	UNP A0A1A7BFR5
A	571	HIS	-	expression tag	UNP A0A1A7BFR5
A	572	HIS	-	expression tag	UNP A0A1A7BFR5
A	573	HIS	-	expression tag	UNP A0A1A7BFR5
A	574	HIS	-	expression tag	UNP A0A1A7BFR5
A	575	HIS	-	expression tag	UNP A0A1A7BFR5
B	568	LEU	-	expression tag	UNP A0A1A7BFR5
B	569	GLU	-	expression tag	UNP A0A1A7BFR5
B	570	HIS	-	expression tag	UNP A0A1A7BFR5
B	571	HIS	-	expression tag	UNP A0A1A7BFR5
B	572	HIS	-	expression tag	UNP A0A1A7BFR5
B	573	HIS	-	expression tag	UNP A0A1A7BFR5
B	574	HIS	-	expression tag	UNP A0A1A7BFR5
B	575	HIS	-	expression tag	UNP A0A1A7BFR5

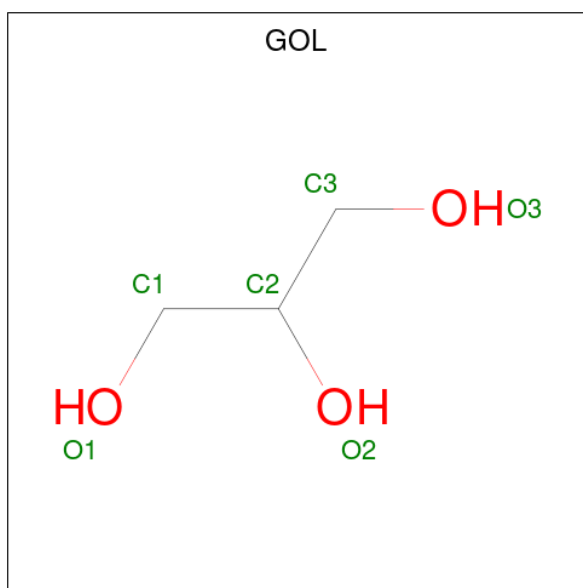
- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S) (labeled as "Ligand of Interest" by depositor).



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

- Molecule 3 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃) (labeled as "Ligand of In-

terest" by depositor).



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

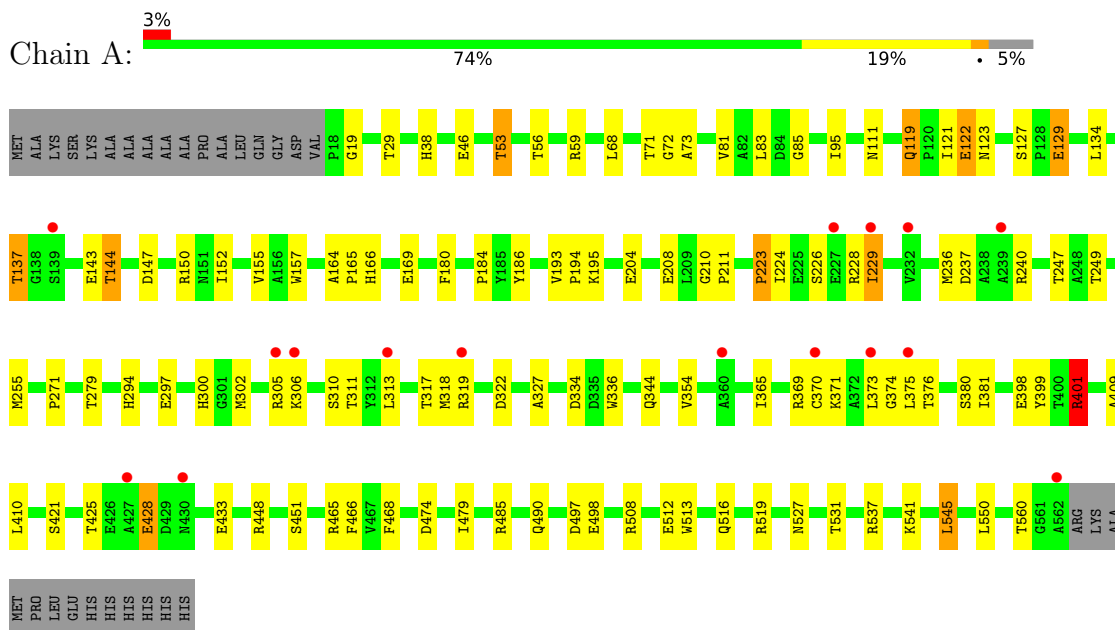
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	160	Total	O	0	0
			160	160		
4	B	169	Total	O	0	0
			169	169		

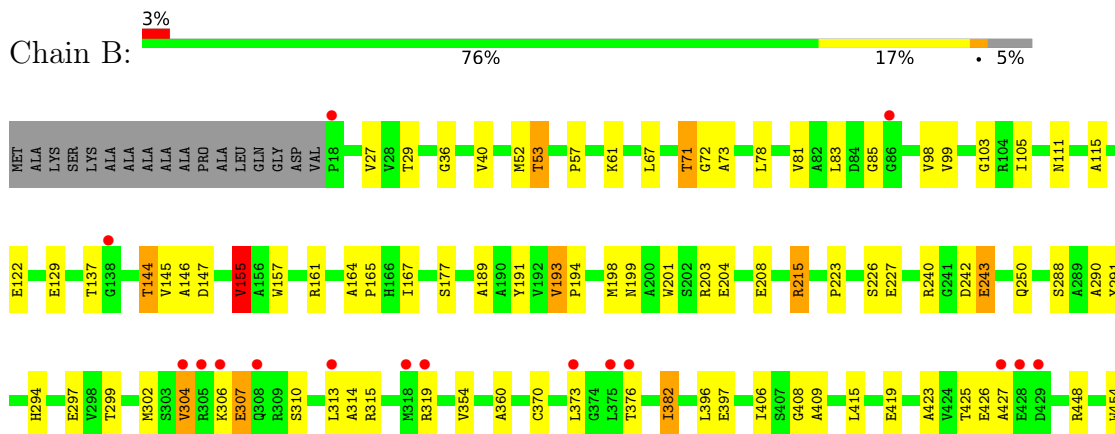
3 Residue-property plots

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

- Molecule 1: NAD-dependent epimerase/dehydratase:Short-chain dehydrogenase/reductase SDR



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4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	85.64Å 109.07Å 129.42Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	32.42 – 2.20 32.42 – 2.20	Depositor EDS
% Data completeness (in resolution range)	96.7 (32.42-2.20) 96.7 (32.42-2.20)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.25 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.8.0238	Depositor
R, R_{free}	0.198 , 0.258 0.208 , 0.259	Depositor DCC
R_{free} test set	3029 reflections (4.88%)	wwPDB-VP
Wilson B-factor (Å ²)	28.3	Xtriage
Anisotropy	0.818	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 30.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	8556	wwPDB-VP
Average B, all atoms (Å ²)	42.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.33% 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: GOL, SO4

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	1.13	1/4145 (0.0%)	1.55	7/5623 (0.1%)
1	B	1.11	0/4153	1.51	13/5634 (0.2%)
All	All	1.12	1/8298 (0.0%)	1.53	20/11257 (0.2%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	327	ALA	C-O	6.74	1.31	1.23

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	177	SER	O-C-N	6.43	126.29	121.47
1	B	36	GLY	CA-C-N	5.95	126.58	119.98
1	B	36	GLY	C-N-CA	5.95	126.58	119.98
1	A	474	ASP	N-CA-C	-5.83	106.22	113.38
1	B	215	ARG	CB-CA-C	-5.75	100.48	109.84
1	A	428	GLU	CA-C-N	5.64	128.11	120.38
1	A	428	GLU	C-N-CA	5.64	128.11	120.38
1	B	72	GLY	CA-C-N	5.61	129.06	120.82
1	B	72	GLY	C-N-CA	5.61	129.06	120.82
1	B	520	PHE	CA-CB-CG	5.49	119.29	113.80
1	B	360	ALA	CA-C-N	5.42	127.85	120.54
1	B	360	ALA	C-N-CA	5.42	127.85	120.54
1	B	155	VAL	N-CA-CB	5.31	116.20	110.62
1	A	401	ARG	N-CA-C	-5.28	105.53	111.28
1	A	237	ASP	CA-CB-CG	5.24	117.84	112.60
1	B	426	GLU	CA-C-N	5.20	128.48	121.05
1	B	426	GLU	C-N-CA	5.20	128.48	121.05
1	B	203	ARG	CB-CG-CD	5.20	123.25	111.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	223	PRO	CA-C-O	-5.10	115.92	121.27
1	A	433	GLU	CB-CA-C	5.10	118.95	110.90

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4074	0	4059	83	0
1	B	4082	0	4062	79	0
2	A	35	0	0	1	0
2	B	30	0	0	2	0
3	B	6	0	8	1	0
4	A	160	0	0	8	0
4	B	169	0	0	4	0
All	All	8556	0	8129	154	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 9.

All (154) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:223:PRO:HD3	1:B:302:MET:HE1	1.24	1.15
1:A:85:GLY:HA3	1:A:155:VAL:HG12	1.24	1.12
1:B:223:PRO:CD	1:B:302:MET:HE1	1.90	0.99
1:A:226:SER:HB2	1:A:229:ILE:HG23	1.52	0.92
1:A:85:GLY:CA	1:A:155:VAL:HG12	2.04	0.87
1:A:344:GLN:HG2	1:A:375:LEU:HD11	1.57	0.86
1:A:344:GLN:CG	1:A:375:LEU:HD11	2.08	0.83
1:A:294:HIS:HE1	1:B:297:GLU:H	1.25	0.83
1:B:71:THR:HG23	1:B:73:ALA:H	1.45	0.80
1:A:208:GLU:OE2	1:A:425:THR:HB	1.82	0.78
1:A:85:GLY:HA3	1:A:155:VAL:CG1	2.10	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:243:GLU:HG2	4:B:1404:HOH:O	1.83	0.77
1:B:315:ARG:CZ	1:B:562:ALA:CB	2.64	0.76
1:B:315:ARG:NE	1:B:562:ALA:HB3	2.01	0.75
1:B:397:GLU:OE1	3:B:1301:GOL:H12	1.87	0.74
1:A:29:THR:HA	1:A:53:THR:HB	1.70	0.72
1:B:215:ARG:HG2	1:B:291:TYR:O	1.88	0.72
1:B:144:THR:HG22	1:B:147:ASP:H	1.52	0.72
1:B:315:ARG:CZ	1:B:562:ALA:HB3	2.20	0.72
1:A:144:THR:HG22	1:A:147:ASP:H	1.55	0.71
1:B:29:THR:HA	1:B:53:THR:HB	1.71	0.71
1:A:344:GLN:HG2	1:A:375:LEU:CD1	2.21	0.70
1:B:315:ARG:NH2	1:B:562:ALA:CB	2.55	0.69
1:B:204:GLU:OE1	1:B:479:ILE:HD12	1.92	0.68
1:A:236:MET:O	1:A:240:ARG:HG2	1.97	0.65
1:B:315:ARG:NH2	1:B:562:ALA:HB3	2.11	0.65
1:A:344:GLN:CG	1:A:375:LEU:CD1	2.73	0.65
1:B:29:THR:O	1:B:111:ASN:HB3	1.98	0.63
1:A:294:HIS:CE1	1:B:297:GLU:H	2.11	0.63
1:A:425:THR:HG22	4:A:1513:HOH:O	1.98	0.63
1:A:166:HIS:HD2	4:A:1544:HOH:O	1.81	0.61
1:A:485:ARG:HA	1:A:516:GLN:HE22	1.65	0.61
1:B:223:PRO:CD	1:B:302:MET:CE	2.75	0.61
1:A:490:GLN:HA	1:A:490:GLN:OE1	2.00	0.61
1:A:519:ARG:HG2	1:A:531:THR:HG21	1.82	0.61
1:B:306:LYS:C	1:B:307:GLU:HG3	2.26	0.60
1:B:98:VAL:HG11	1:B:105:ILE:CD1	2.31	0.60
1:B:193:VAL:HG22	1:B:194:PRO:HD3	1.83	0.60
1:A:369:ARG:HG3	1:A:369:ARG:HH21	1.64	0.60
1:A:38:HIS:HE1	4:A:1434:HOH:O	1.84	0.59
1:B:466:PHE:CE1	1:B:514:GLY:HA3	2.37	0.59
1:A:180:PHE:HZ	1:A:223:PRO:HB3	1.68	0.58
1:B:315:ARG:NH2	1:B:562:ALA:HB1	2.19	0.58
1:A:134:LEU:O	1:A:137:THR:HB	2.04	0.58
1:B:376:THR:HA	4:B:1516:HOH:O	2.03	0.57
1:B:315:ARG:HE	1:B:562:ALA:HB3	1.67	0.57
1:B:315:ARG:NE	1:B:562:ALA:CB	2.66	0.57
1:A:71:THR:C	1:A:73:ALA:H	2.13	0.56
1:A:193:VAL:HB	1:A:194:PRO:HD3	1.88	0.55
1:B:27:VAL:HG21	1:B:105:ILE:HD12	1.89	0.55
1:A:297:GLU:H	1:B:294:HIS:HE1	1.54	0.55
1:B:223:PRO:HD2	1:B:302:MET:HE1	1.84	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:40:VAL:HG11	1:B:67:LEU:HD13	1.89	0.54
1:A:143:GLU:HG3	4:A:1457:HOH:O	2.08	0.53
1:B:415:LEU:HB3	1:B:419:GLU:HG2	1.90	0.53
1:A:223:PRO:O	1:A:271:PRO:HD2	2.09	0.53
1:A:370:CYS:O	1:A:374:GLY:N	2.42	0.53
1:A:550:LEU:HD22	1:B:313:LEU:CD1	2.39	0.53
1:A:297:GLU:H	1:B:294:HIS:CE1	2.26	0.52
1:A:344:GLN:NE2	1:A:373:LEU:HD21	2.25	0.52
1:B:485:ARG:HA	1:B:516:GLN:HE22	1.74	0.52
1:A:204:GLU:OE1	1:A:479:ILE:HD13	2.10	0.51
1:B:208:GLU:OE2	1:B:425:THR:OG1	2.25	0.51
1:A:224:ILE:O	1:A:229:ILE:HG12	2.09	0.51
1:A:310:SER:HB2	1:B:310:SER:OG	2.11	0.51
1:A:537:ARG:HA	1:A:541:LYS:HG3	1.93	0.51
1:A:545:LEU:N	1:A:545:LEU:HD23	2.26	0.51
1:A:164:ALA:HB3	1:A:165:PRO:HD3	1.93	0.50
1:B:354:VAL:CG1	1:B:382:ILE:HD13	2.42	0.49
1:A:421:SER:HB3	4:A:1490:HOH:O	2.12	0.49
1:A:409:ALA:O	1:A:466:PHE:HA	2.13	0.49
1:B:522:ASN:ND2	1:B:524:GLU:OE1	2.43	0.49
1:B:304:VAL:HG12	4:B:1526:HOH:O	2.13	0.48
1:A:111:ASN:ND2	4:A:1407:HOH:O	2.46	0.48
1:A:527:ASN:O	1:A:531:THR:HG23	2.12	0.48
1:A:344:GLN:HG3	1:A:375:LEU:CD1	2.42	0.47
1:B:396:LEU:HD22	1:B:406:ILE:HD13	1.96	0.47
1:B:98:VAL:HG11	1:B:105:ILE:HD11	1.97	0.47
1:B:299:THR:O	1:B:302:MET:HG3	2.14	0.47
1:A:19:GLY:HA3	1:A:46:GLU:O	2.15	0.47
1:B:299:THR:HB	1:B:302:MET:HG3	1.96	0.47
1:A:166:HIS:CD2	4:A:1544:HOH:O	2.63	0.47
1:B:29:THR:HG1	1:B:111:ASN:H	1.63	0.46
1:B:85:GLY:HA3	1:B:155:VAL:CG1	2.45	0.46
1:A:121:ILE:O	1:A:448:ARG:HG3	2.15	0.46
1:A:398:GLU:O	1:A:401:ARG:HB2	2.14	0.46
1:B:223:PRO:HD2	1:B:302:MET:CE	2.42	0.46
1:A:425:THR:HG22	1:A:425:THR:O	2.16	0.46
1:B:53:THR:HG23	1:B:83:LEU:HB3	1.96	0.46
1:A:184:PRO:HB3	1:A:497:ASP:CG	2.40	0.46
1:A:369:ARG:HH21	1:A:369:ARG:CG	2.28	0.45
1:A:53:THR:HG23	1:A:83:LEU:HB3	1.98	0.45
1:A:119:GLN:HG3	1:A:123:ASN:O	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:71:THR:HG23	1:B:73:ALA:CB	2.46	0.45
1:B:315:ARG:HH21	1:B:562:ALA:HB3	1.81	0.45
1:A:150:ARG:O	1:A:155:VAL:HG13	2.17	0.45
1:B:314:ALA:CA	1:B:560:THR:HG22	2.46	0.45
1:A:465:ARG:HA	1:A:513:TRP:O	2.17	0.45
1:A:508:ARG:NH1	1:A:512:GLU:OE2	2.50	0.45
1:A:336:TRP:CZ2	1:A:365:ILE:HG22	2.52	0.44
1:A:344:GLN:HE22	1:A:373:LEU:HD21	1.81	0.44
1:B:189:ALA:O	1:B:193:VAL:HG13	2.16	0.44
1:A:228:ARG:CG	1:A:229:ILE:N	2.81	0.44
1:B:105:ILE:O	1:B:167:ILE:HA	2.17	0.44
1:B:198:MET:O	1:B:199:ASN:C	2.60	0.44
1:B:396:LEU:HD22	1:B:406:ILE:CD1	2.48	0.44
1:B:85:GLY:HA3	1:B:155:VAL:HG13	2.00	0.44
1:B:240:ARG:NH2	1:B:242:ASP:OD2	2.51	0.43
1:B:454:TRP:CD1	1:B:460:LEU:HD11	2.53	0.43
1:A:226:SER:HB2	1:A:229:ILE:CG2	2.34	0.43
1:B:161:ARG:HB2	1:B:425:THR:HA	2.01	0.43
1:B:164:ALA:N	1:B:165:PRO:CD	2.81	0.43
1:A:421:SER:HB2	2:A:1307:SO4:O1	2.19	0.43
1:B:408:GLY:HA2	1:B:465:ARG:O	2.18	0.43
1:A:468:PHE:O	1:A:516:GLN:HA	2.19	0.43
1:B:157:TRP:CE3	1:B:201:TRP:CD1	3.07	0.43
1:A:226:SER:CB	1:A:229:ILE:HG23	2.34	0.42
1:A:157:TRP:CD1	1:A:157:TRP:C	2.98	0.42
1:B:144:THR:HG23	1:B:146:ALA:H	1.83	0.42
1:A:410:LEU:C	1:A:410:LEU:HD23	2.44	0.42
1:B:57:PRO:O	1:B:61:LYS:HB2	2.19	0.42
1:A:71:THR:C	1:A:73:ALA:N	2.78	0.42
1:A:300:HIS:HD2	1:B:290:ALA:HA	1.85	0.42
1:B:215:ARG:NH1	1:B:288:SER:OG	2.53	0.42
1:B:115:ALA:O	1:B:191:TYR:HB2	2.20	0.42
1:B:467:VAL:HG22	1:B:538:ILE:HG21	2.01	0.42
1:A:56:THR:O	1:A:59:ARG:HB3	2.20	0.42
1:A:255:MET:HG2	1:A:302:MET:HG3	2.02	0.42
1:A:83:LEU:C	1:A:83:LEU:HD23	2.45	0.42
1:A:344:GLN:NE2	1:A:373:LEU:CD2	2.82	0.41
1:A:152:ILE:HG23	1:A:195:LYS:HE2	2.02	0.41
1:B:409:ALA:O	1:B:466:PHE:HA	2.20	0.41
1:A:354:VAL:HG22	1:A:399:TYR:CE1	2.55	0.41
1:B:226:SER:HB2	2:B:1304:SO4:O3	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:71:THR:CG2	1:B:73:ALA:HB3	2.51	0.41
1:A:186:TYR:CD2	1:A:498:GLU:HG2	2.55	0.41
1:B:227:GLU:N	2:B:1304:SO4:O3	2.48	0.41
1:B:458:ASP:HA	4:B:1495:HOH:O	2.19	0.41
1:A:425:THR:CG2	4:A:1513:HOH:O	2.61	0.41
1:A:313:LEU:HD13	1:B:545:LEU:HD12	2.03	0.41
1:A:319:ARG:HB2	1:A:322:ASP:OD2	2.21	0.41
1:A:369:ARG:CG	1:A:369:ARG:NH2	2.81	0.41
1:B:466:PHE:CD1	1:B:514:GLY:HA3	2.56	0.41
1:A:373:LEU:HD12	1:A:373:LEU:HA	1.83	0.41
1:B:474:ASP:O	1:B:476:LYS:HG2	2.21	0.41
1:A:68:LEU:O	1:A:71:THR:O	2.39	0.40
1:A:369:ARG:CZ	1:A:369:ARG:HB2	2.50	0.40
1:A:127:SER:HB2	1:A:129:GLU:OE1	2.22	0.40
1:B:189:ALA:HB2	1:B:494:ILE:HG21	2.03	0.40
1:A:210:GLY:N	1:A:211:PRO:CD	2.83	0.40
1:B:52:MET:SD	1:B:78:LEU:HD11	2.62	0.40
1:B:423:ALA:O	1:B:427:ALA:CB	2.69	0.40
1:A:122:GLU:HB2	1:A:451:SER:HB3	2.03	0.40
1:B:99:VAL:O	1:B:103:GLY:N	2.53	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	543/575 (94%)	523 (96%)	18 (3%)	2 (0%)	30	34
1	B	544/575 (95%)	532 (98%)	12 (2%)	0	100	100
All	All	1087/1150 (94%)	1055 (97%)	30 (3%)	2 (0%)	43	51

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	72	GLY
1	A	428	GLU

5.3.2 Protein sidechains ⓘ

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	410/431 (95%)	384 (94%)	26 (6%)	16	19
1	B	411/431 (95%)	390 (95%)	21 (5%)	21	27
All	All	821/862 (95%)	774 (94%)	47 (6%)	18	23

All (47) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	53	THR
1	A	81	VAL
1	A	95	ILE
1	A	119	GLN
1	A	122	GLU
1	A	129	GLU
1	A	137	THR
1	A	144	THR
1	A	169	GLU
1	A	229	ILE
1	A	247	THR
1	A	249	THR
1	A	279	THR
1	A	305	ARG
1	A	306	LYS
1	A	311	THR
1	A	317	THR
1	A	318	MET
1	A	334	ASP
1	A	371	LYS
1	A	376	THR
1	A	380	SER

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Mol	Chain	Res	Type
1	A	381	ILE
1	A	401	ARG
1	A	545	LEU
1	A	560	THR
1	B	53	THR
1	B	71	THR
1	B	81	VAL
1	B	122	GLU
1	B	129	GLU
1	B	137	THR
1	B	144	THR
1	B	145	VAL
1	B	155	VAL
1	B	193	VAL
1	B	243	GLU
1	B	250	GLN
1	B	304	VAL
1	B	307	GLU
1	B	319	ARG
1	B	370	CYS
1	B	373	LEU
1	B	382	ILE
1	B	448	ARG
1	B	485	ARG
1	B	526	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	111	ASN
1	A	166	HIS
1	A	264	ASN
1	A	294	HIS
1	A	344	GLN
1	A	482	HIS
1	A	515	ASN
1	A	516	GLN
1	B	111	ASN
1	B	250	GLN
1	B	264	ASN
1	B	294	HIS
1	B	352	GLN

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Mol	Chain	Res	Type
1	B	471	HIS
1	B	482	HIS
1	B	515	ASN
1	B	516	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](#)

14 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	GOL	B	1301	-	5,5,5	0.08	0	5,5,5	0.37	0
2	SO4	A	1302	-	4,4,4	0.33	0	6,6,6	0.16	0
2	SO4	A	1306	-	4,4,4	0.34	0	6,6,6	0.13	0
2	SO4	A	1301	-	4,4,4	0.40	0	6,6,6	0.18	0
2	SO4	A	1305	-	4,4,4	0.33	0	6,6,6	0.08	0
2	SO4	B	1302	-	4,4,4	0.38	0	6,6,6	0.26	0
2	SO4	B	1306	-	4,4,4	0.29	0	6,6,6	0.19	0
2	SO4	B	1305	-	4,4,4	0.33	0	6,6,6	0.11	0
2	SO4	A	1307	-	4,4,4	0.32	0	6,6,6	0.08	0
2	SO4	B	1304	-	4,4,4	0.30	0	6,6,6	0.16	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	SO4	A	1304	-	4,4,4	0.28	0	6,6,6	0.08	0
2	SO4	B	1303	-	4,4,4	0.34	0	6,6,6	0.19	0
2	SO4	B	1307	-	4,4,4	0.31	0	6,6,6	0.09	0
2	SO4	A	1303	-	4,4,4	0.35	0	6,6,6	0.15	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	GOL	B	1301	-	-	0/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

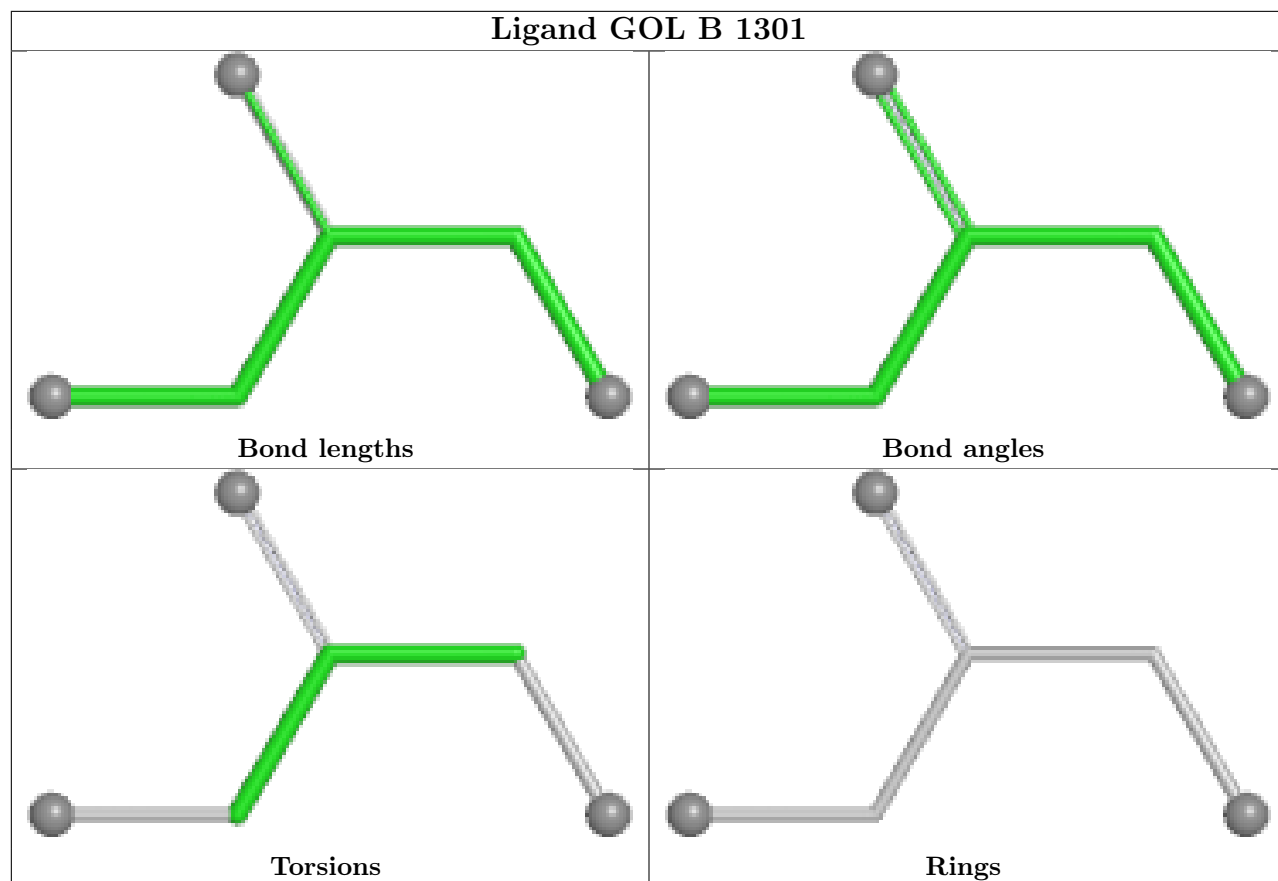
There are no torsion outliers.

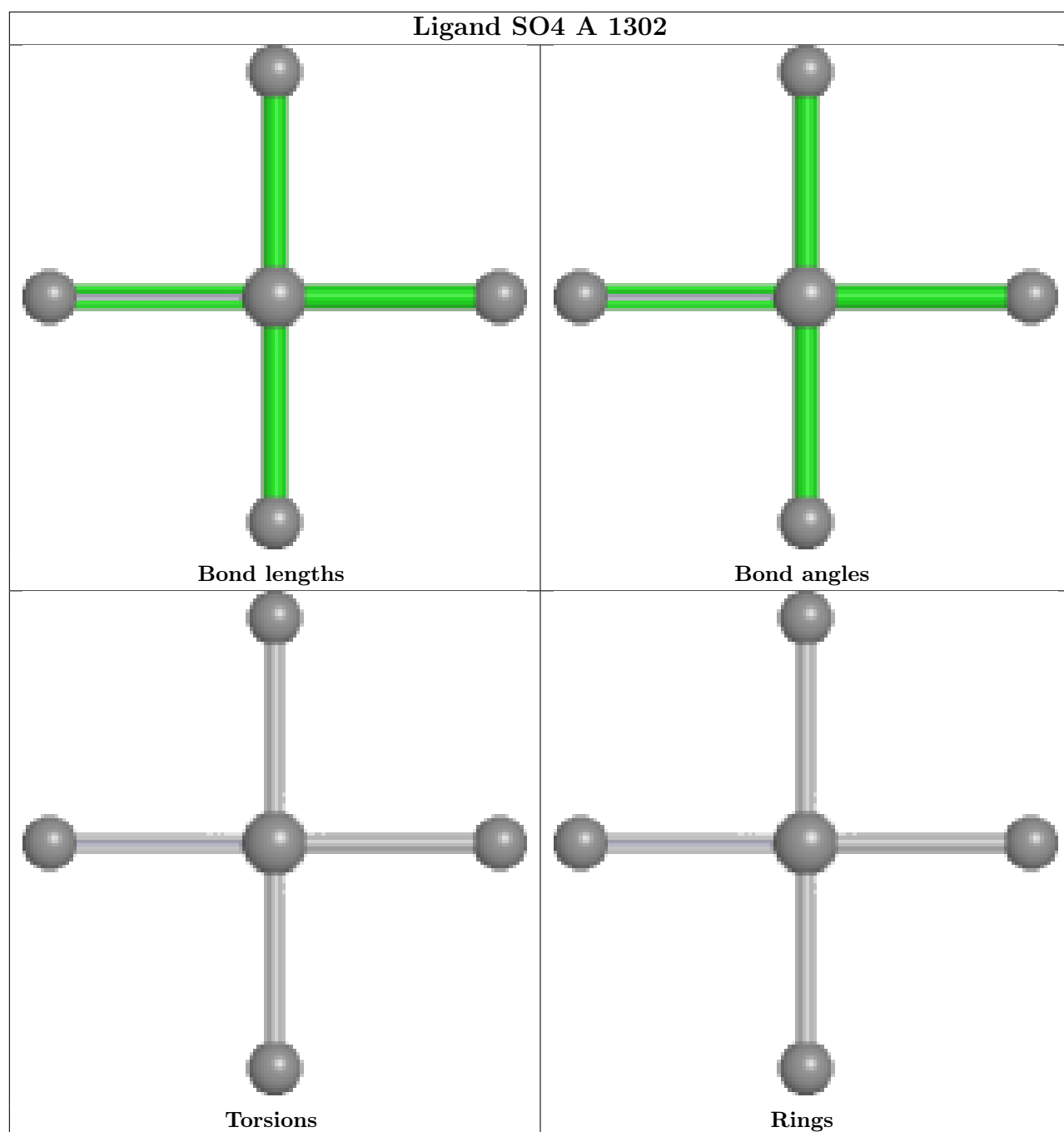
There are no ring outliers.

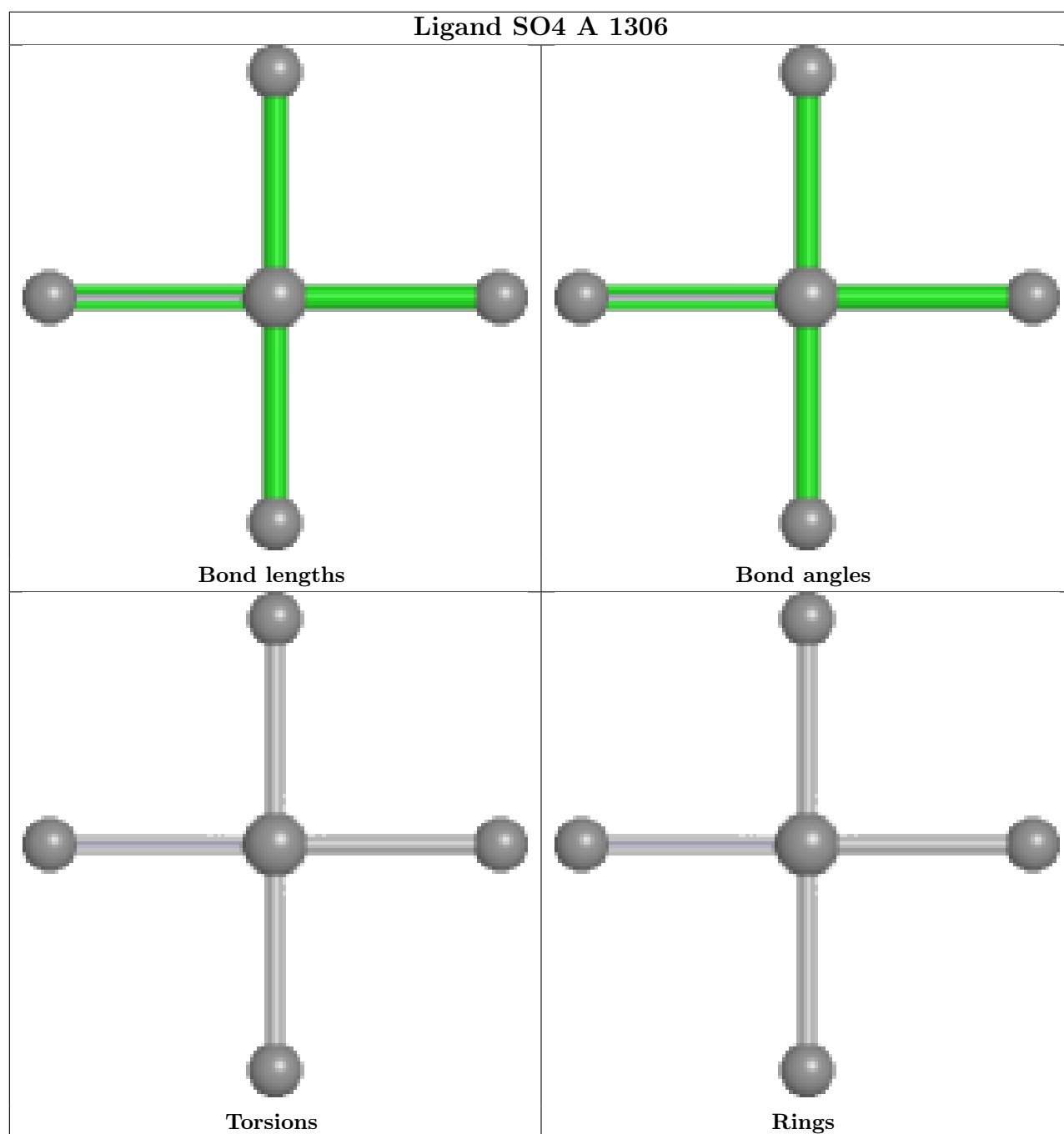
3 monomers are involved in 4 short contacts:

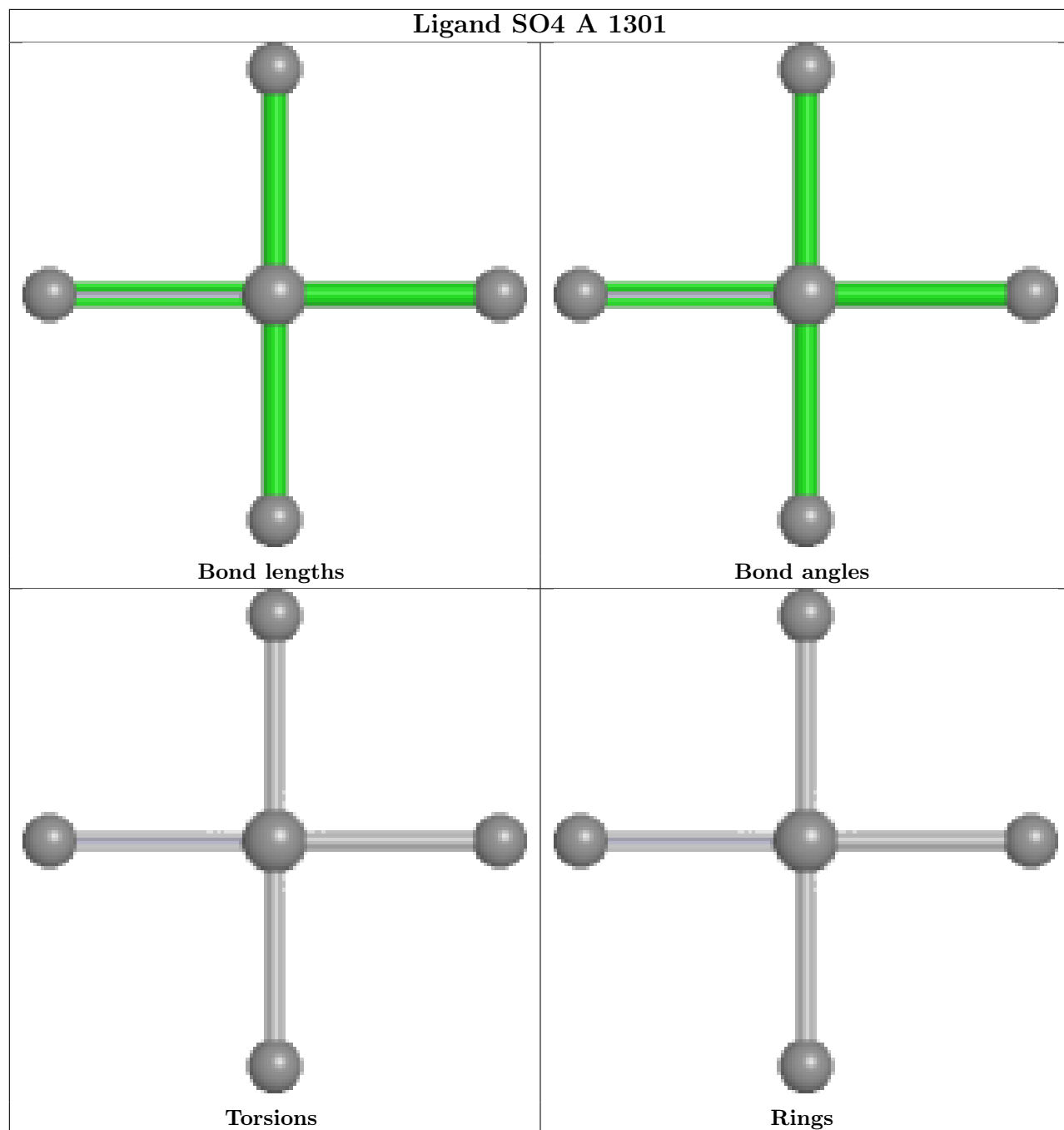
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	1301	GOL	1	0
2	A	1307	SO4	1	0
2	B	1304	SO4	2	0

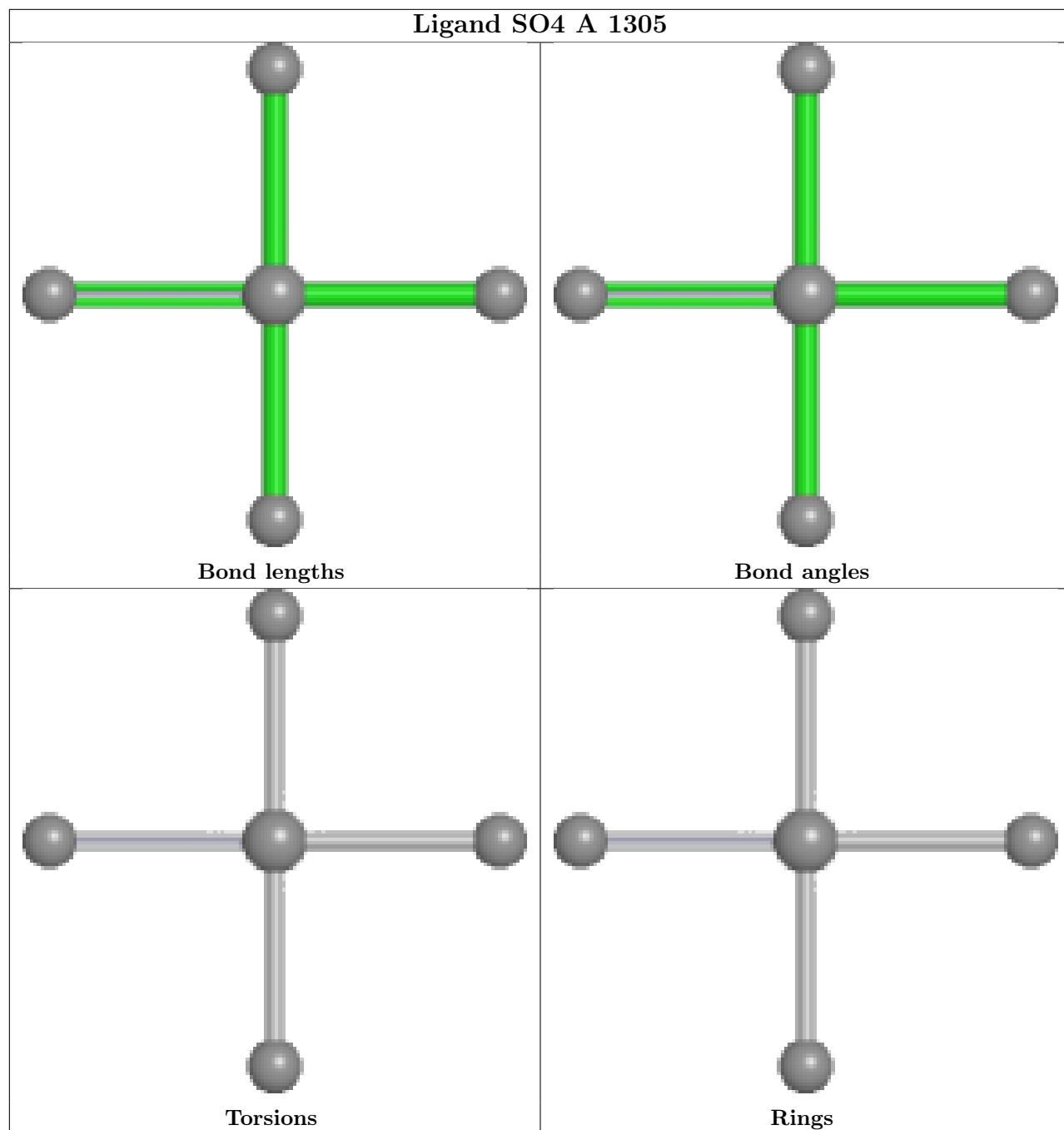
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

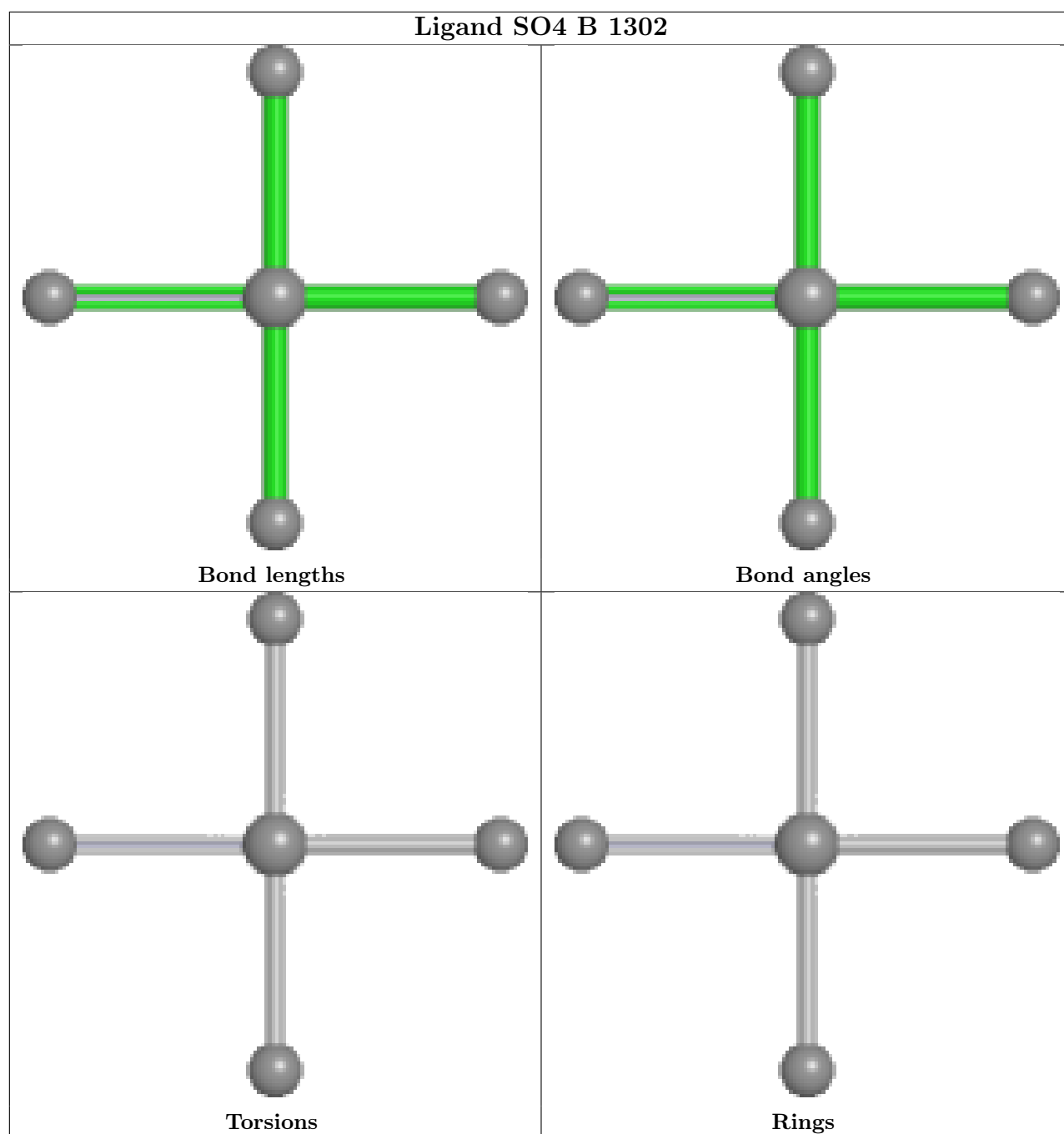


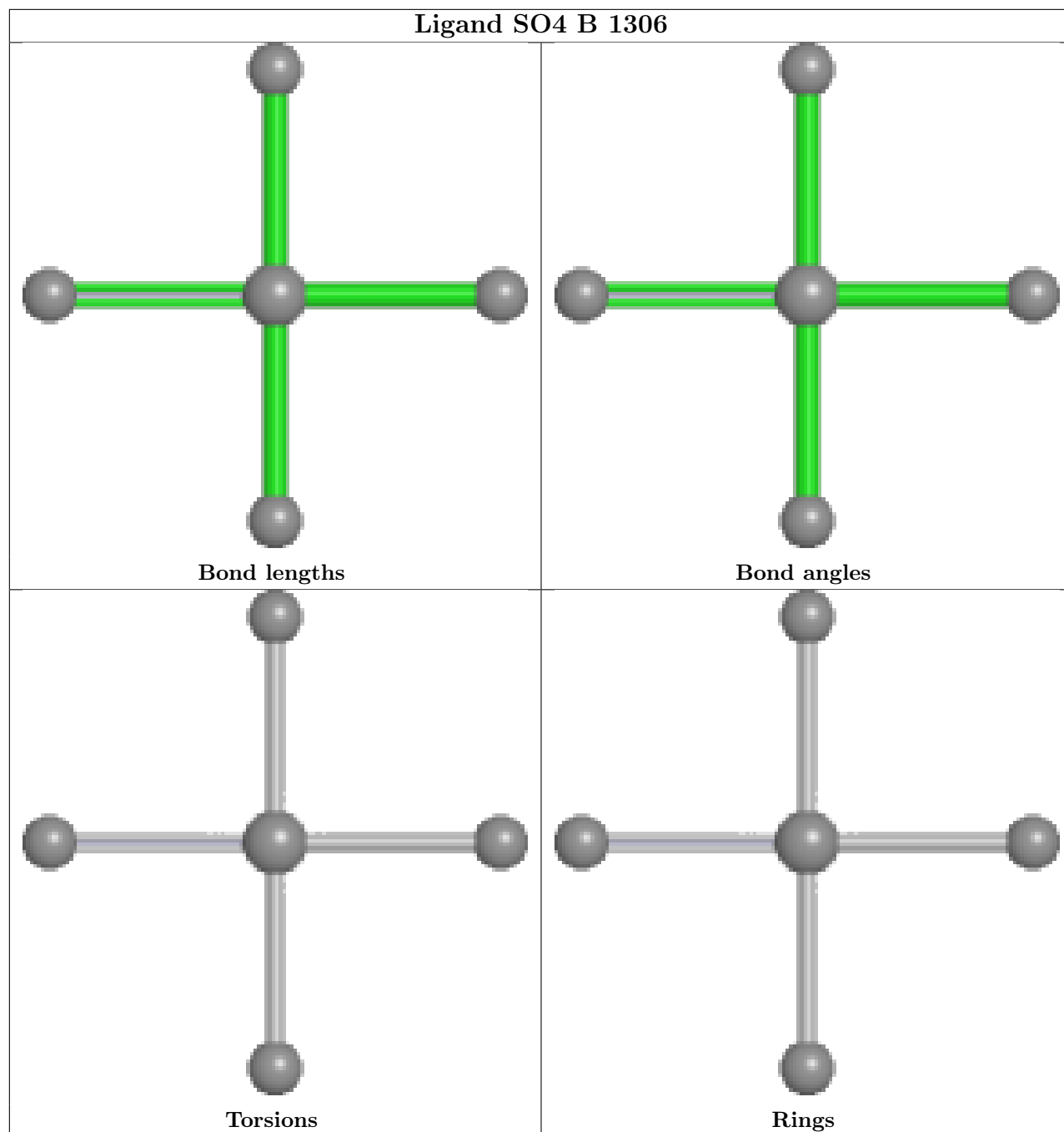


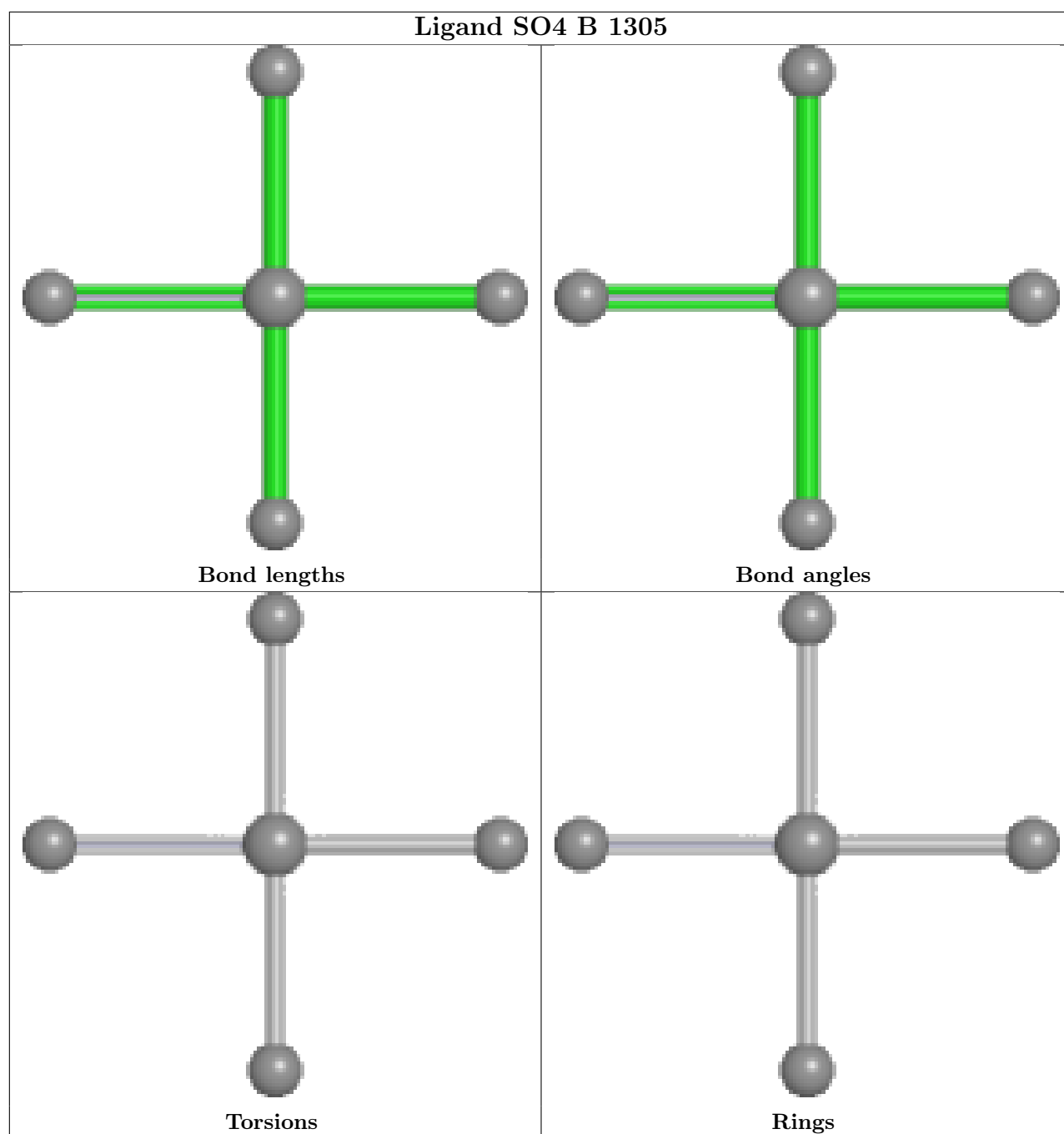


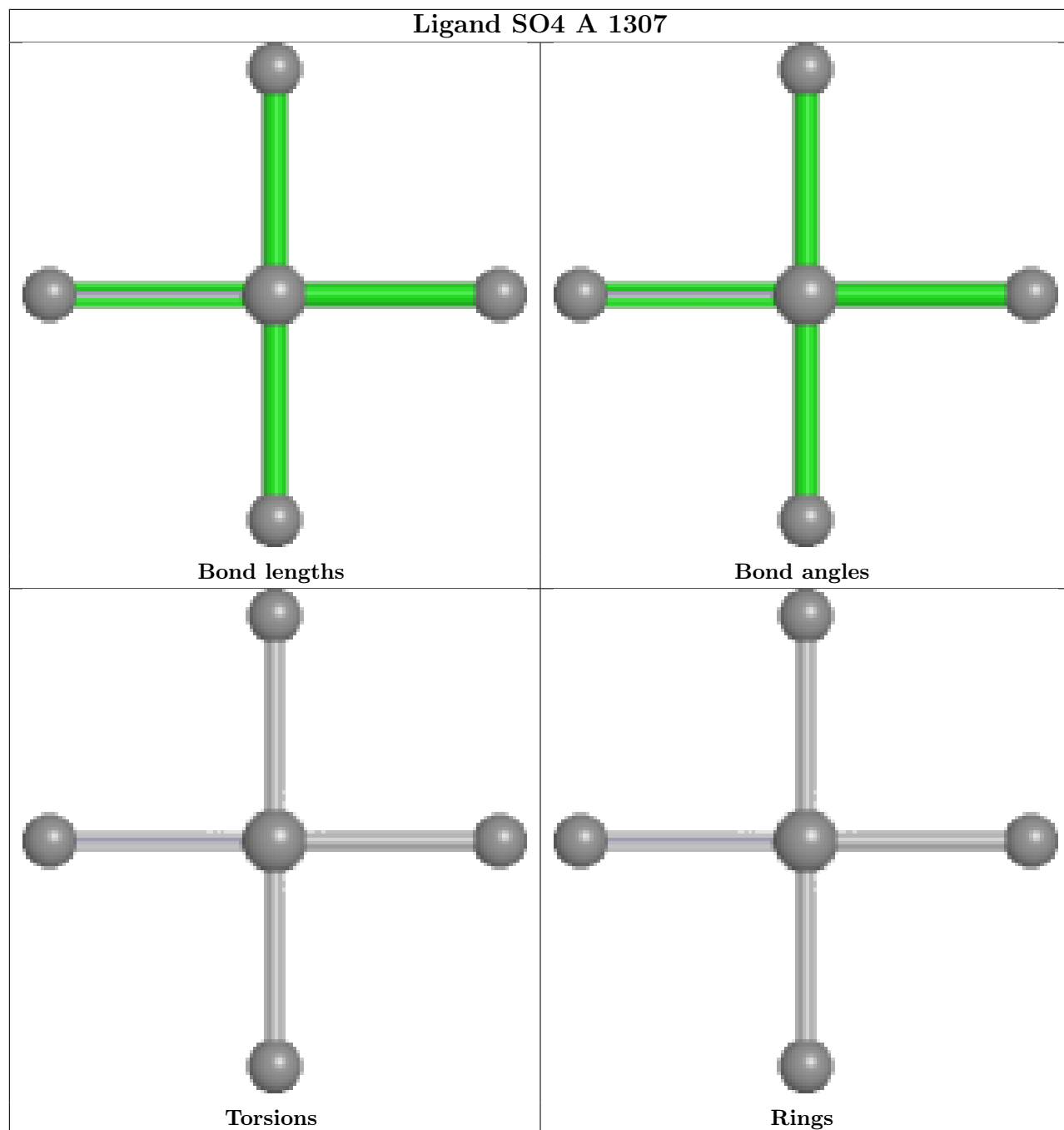


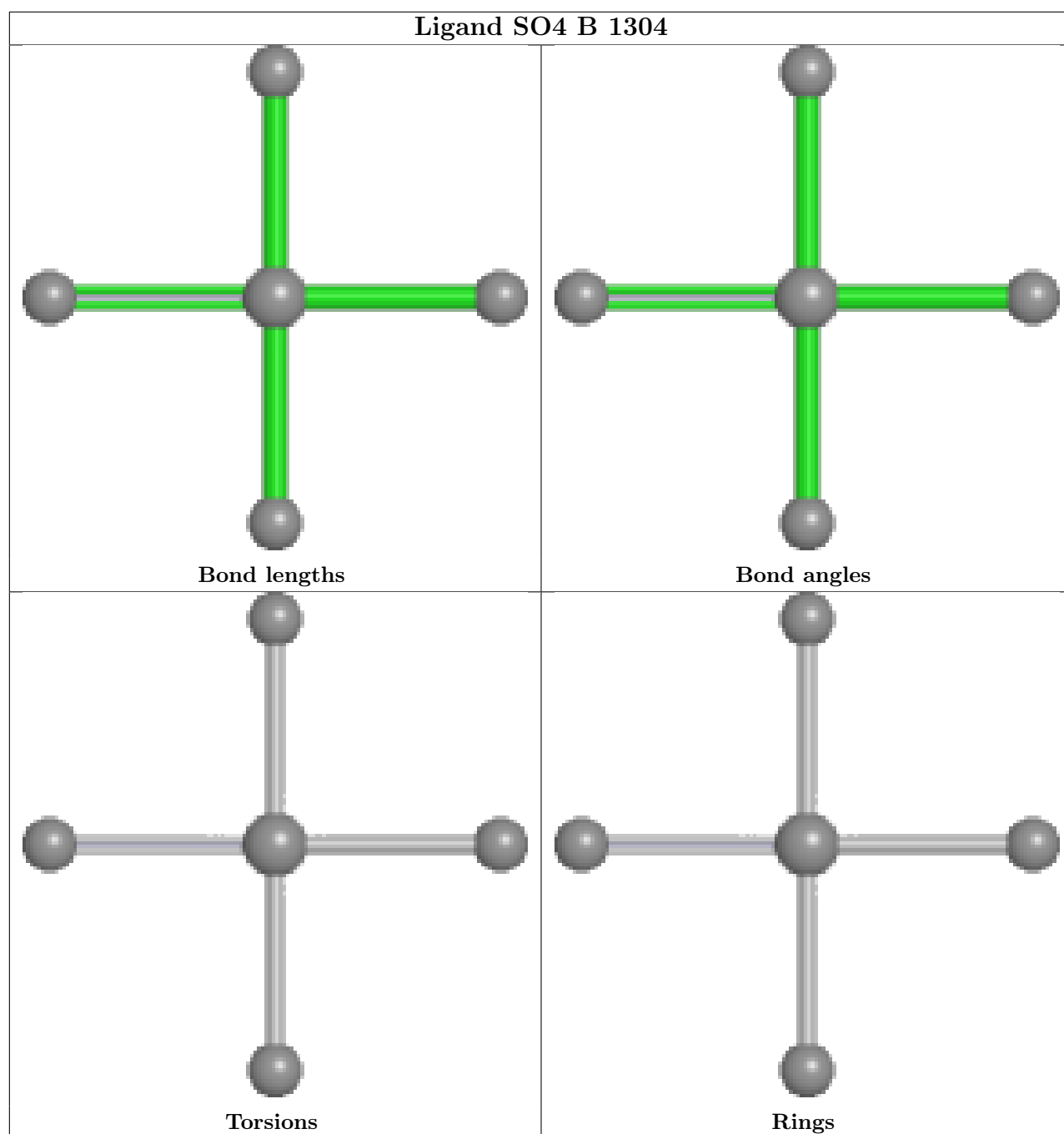


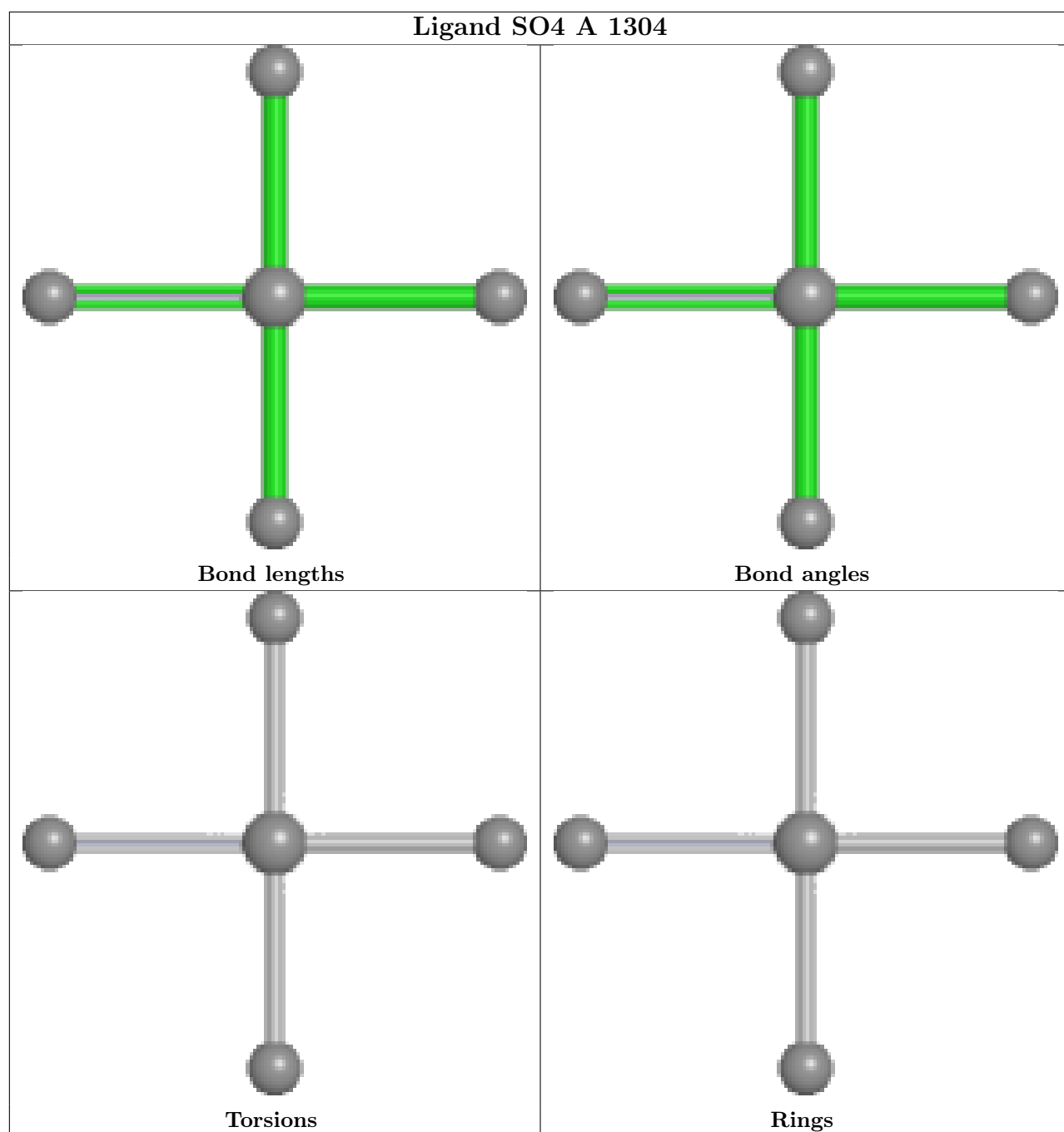


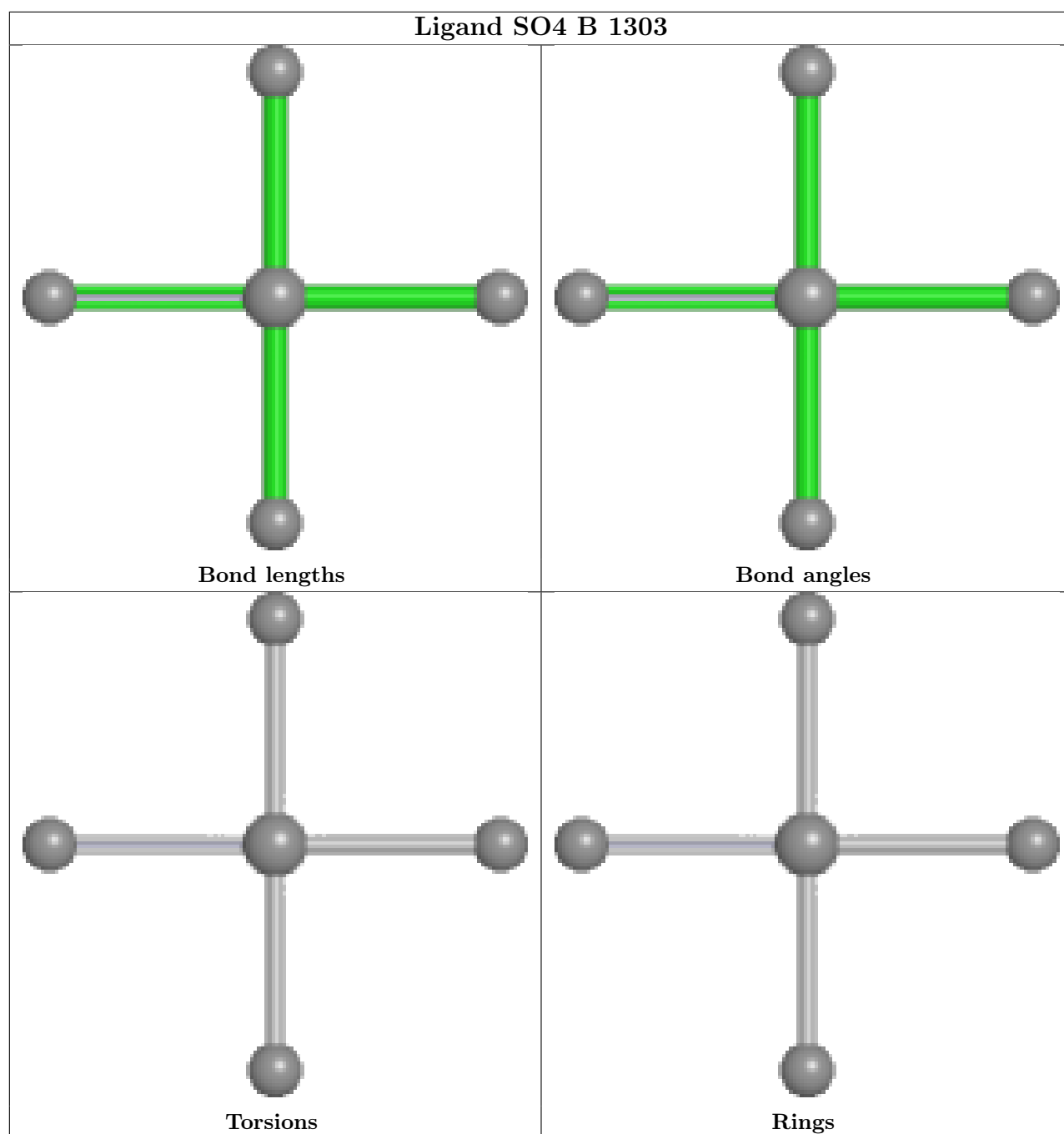


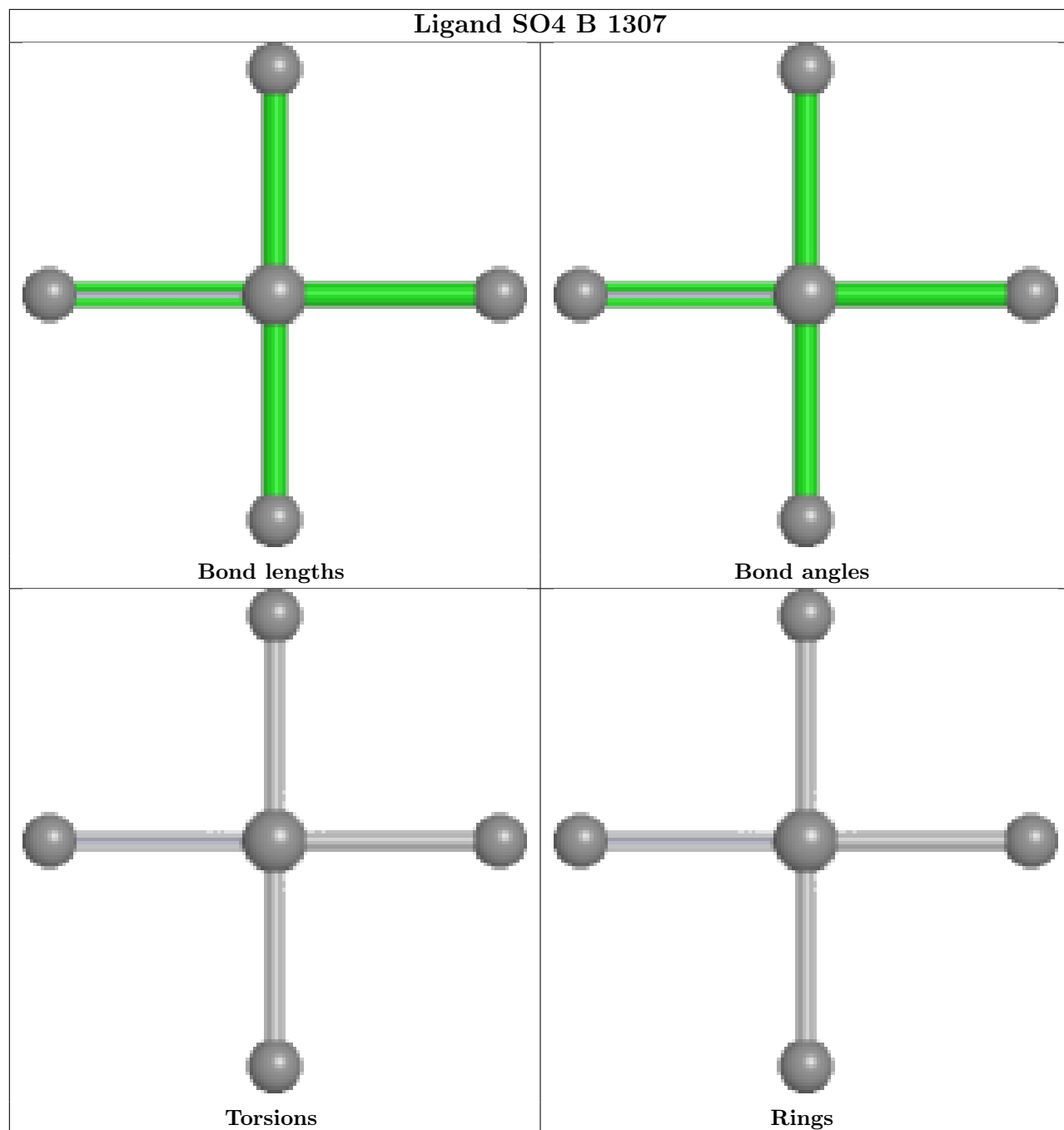


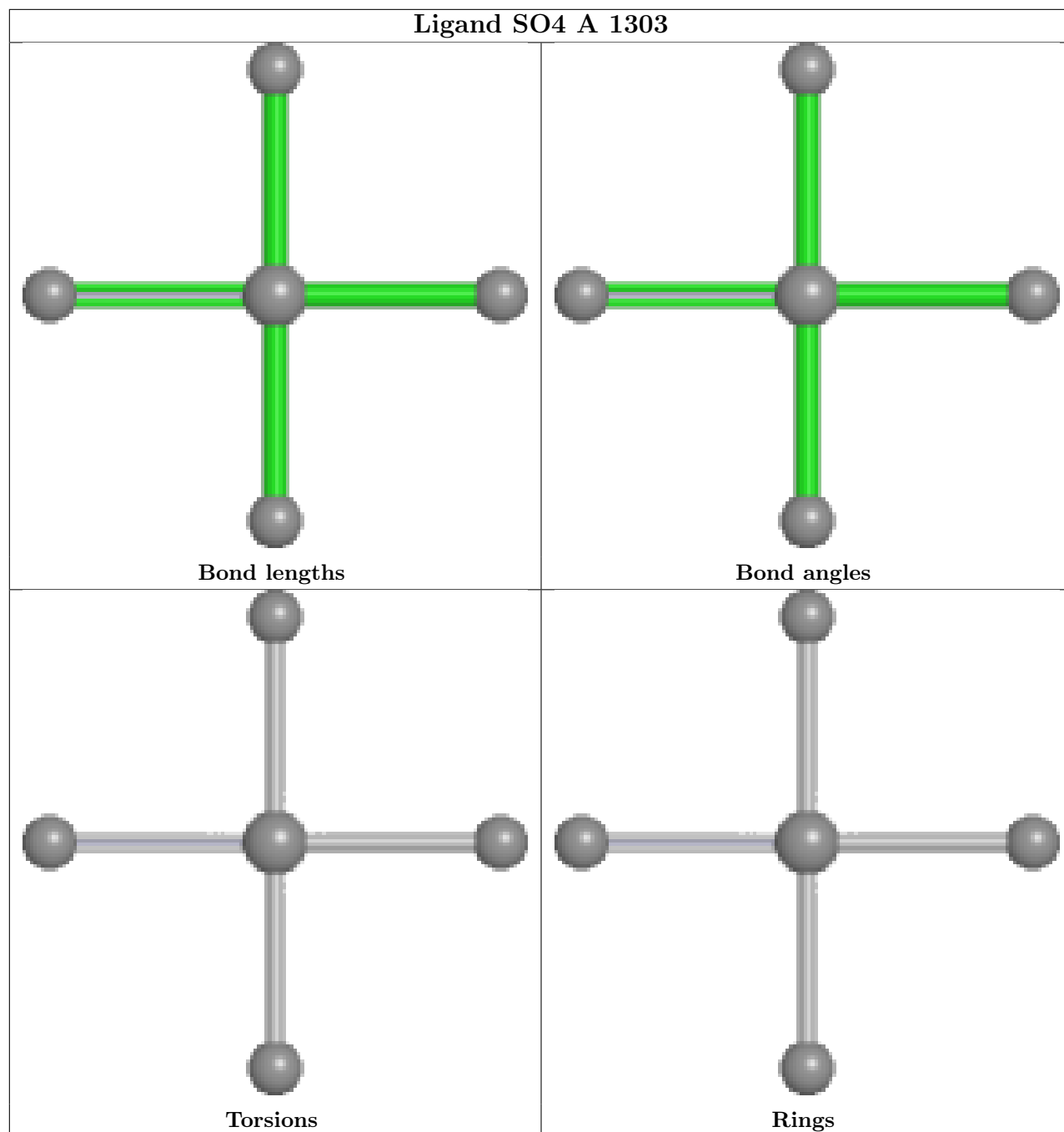












5.7 Other polymers [i](#)

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	545/575 (94%)	0.06	16 (2%) 53 50	20, 37, 83, 128	0
1	B	545/575 (94%)	-0.04	17 (3%) 51 48	21, 36, 80, 146	1 (0%)
All	All	1090/1150 (94%)	0.01	33 (3%) 52 49	20, 36, 82, 146	1 (0%)

All (33) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	427	ALA	5.4
1	A	427	ALA	4.1
1	B	375	LEU	3.9
1	A	306	LYS	3.5
1	A	229	ILE	3.2
1	A	370	CYS	3.2
1	B	318	MET	3.1
1	A	430	ASN	3.1
1	B	18	PRO	3.0
1	A	239	ALA	2.8
1	A	562	ALA	2.8
1	B	305	ARG	2.8
1	B	138	GLY	2.6
1	B	306	LYS	2.4
1	B	376	THR	2.4
1	A	313	LEU	2.4
1	A	375	LEU	2.4
1	A	305	ARG	2.4
1	A	232	VAL	2.3
1	B	429	ASP	2.3
1	A	319	ARG	2.2
1	B	561	GLY	2.2
1	A	360	ALA	2.2
1	B	86	GLY	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	139	SER	2.2
1	B	373	LEU	2.1
1	A	227	GLU	2.1
1	B	304	VAL	2.1
1	A	373	LEU	2.1
1	B	428	GLU	2.1
1	B	308	GLN	2.0
1	B	319	ARG	2.0
1	B	313	LEU	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](#)

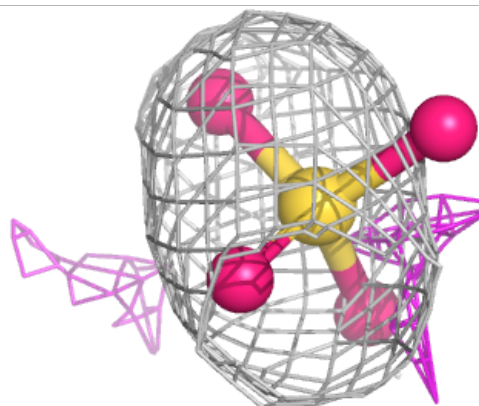
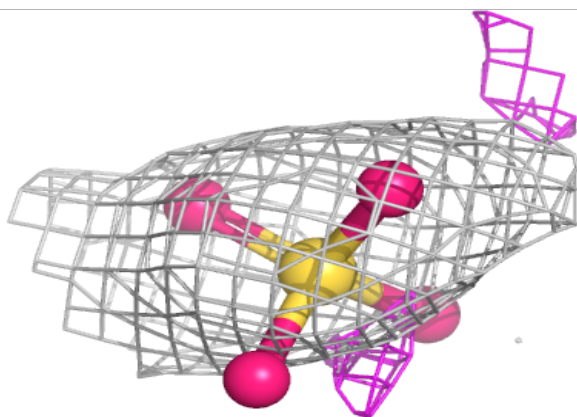
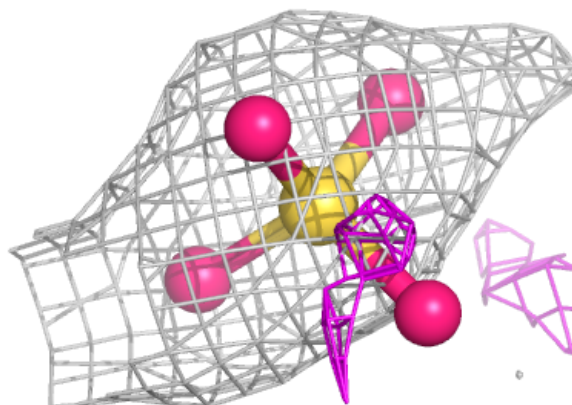
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	SO4	B	1306	5/5	0.62	0.14	96,100,119,132	0
2	SO4	A	1305	5/5	0.69	0.09	97,105,112,113	0
2	SO4	A	1304	5/5	0.69	0.12	83,86,98,103	0
2	SO4	B	1305	5/5	0.76	0.10	100,102,104,109	0
2	SO4	A	1303	5/5	0.80	0.08	73,75,85,91	0
2	SO4	A	1307	5/5	0.80	0.10	100,103,113,115	0
2	SO4	B	1304	5/5	0.84	0.10	77,77,101,101	0
2	SO4	B	1307	5/5	0.85	0.11	83,86,93,104	0
3	GOL	B	1301	6/6	0.88	0.13	37,42,45,46	0
2	SO4	B	1303	5/5	0.89	0.09	49,55,68,74	0
2	SO4	A	1306	5/5	0.93	0.09	37,45,56,58	0
2	SO4	B	1302	5/5	0.94	0.09	50,53,55,58	0
2	SO4	A	1302	5/5	0.97	0.10	38,38,46,51	0
2	SO4	A	1301	5/5	0.98	0.05	40,45,49,52	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.

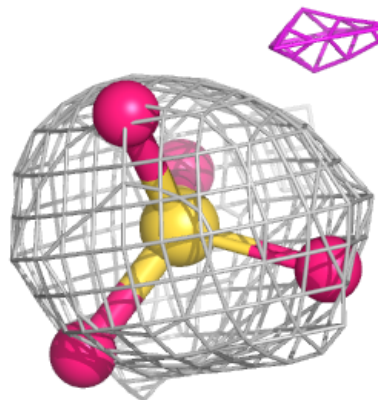
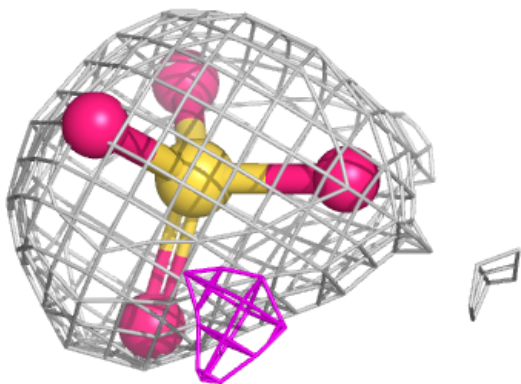
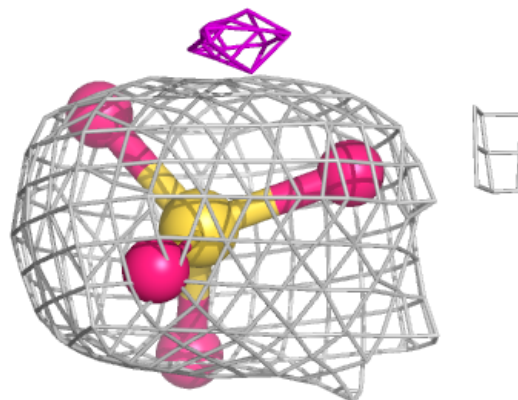
Electron density around SO4 B 1306:

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



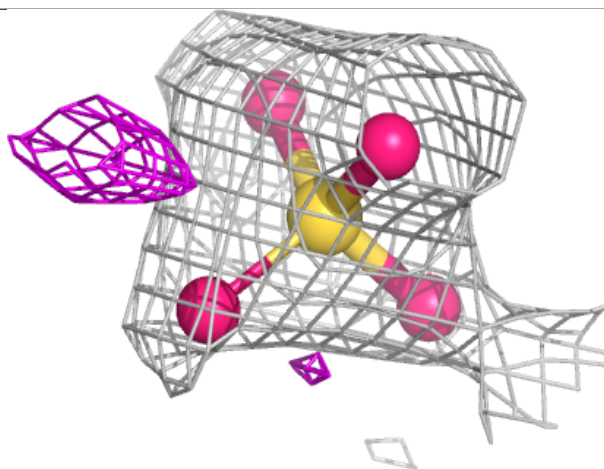
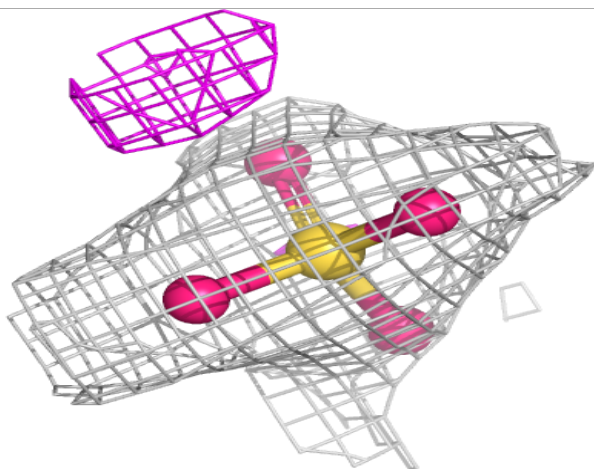
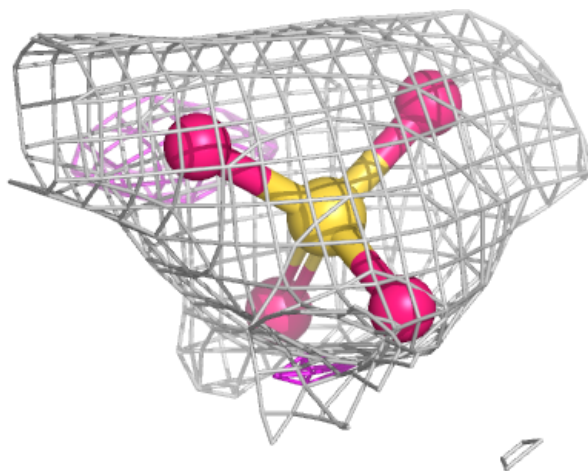
Electron density around SO4 A 1305:

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



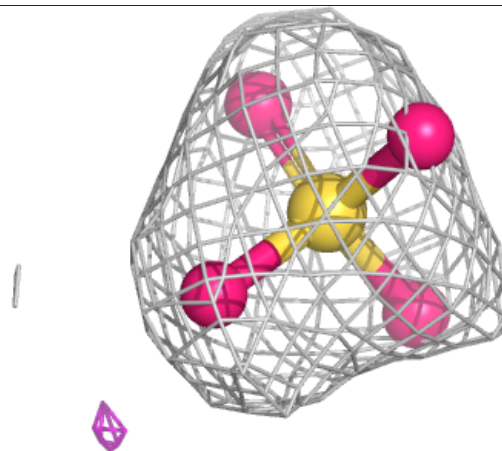
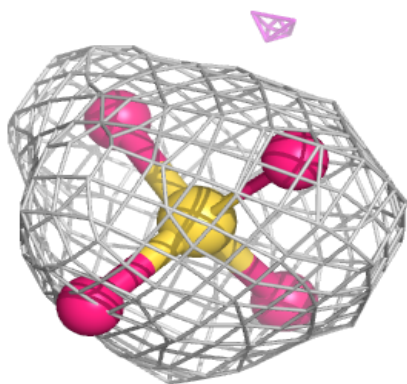
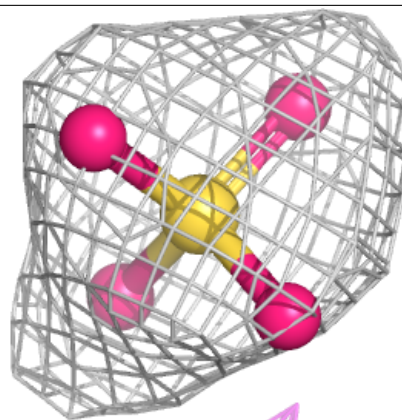
Electron density around SO4 A 1304:

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



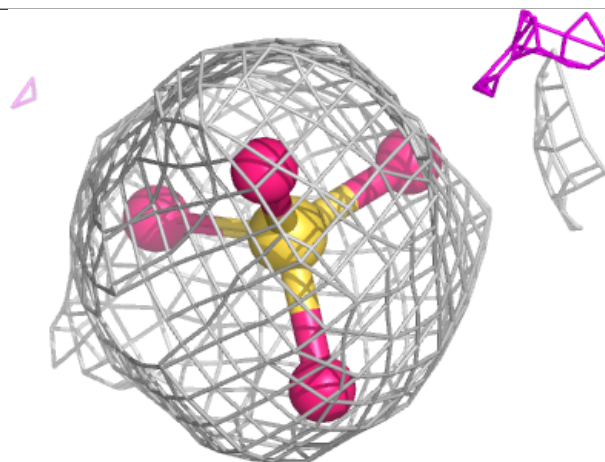
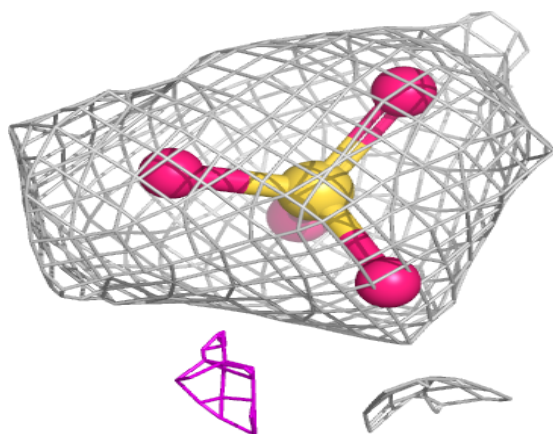
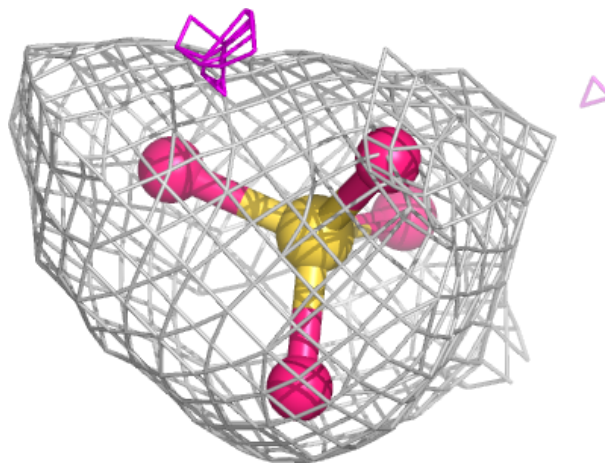
Electron density around SO4 B 1305:

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



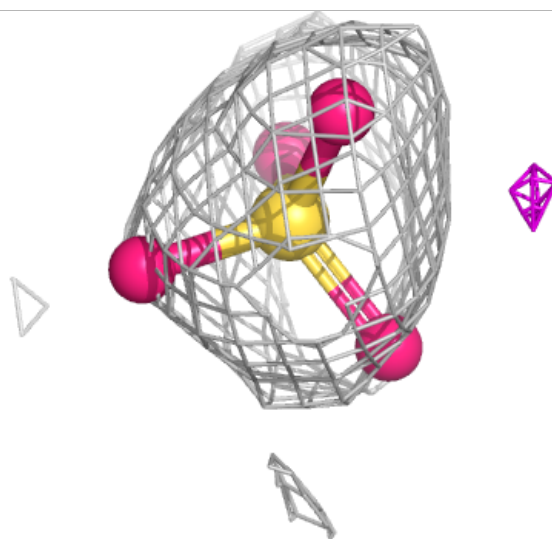
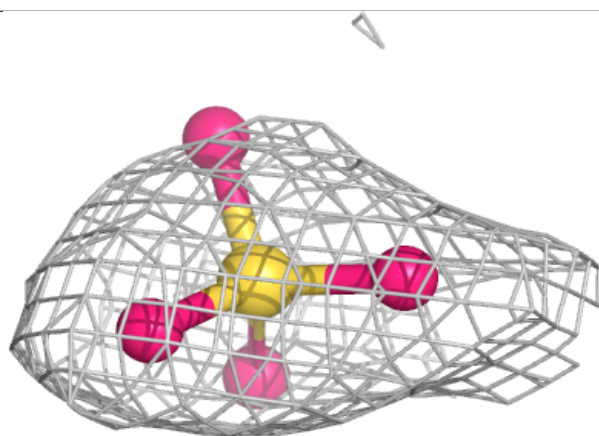
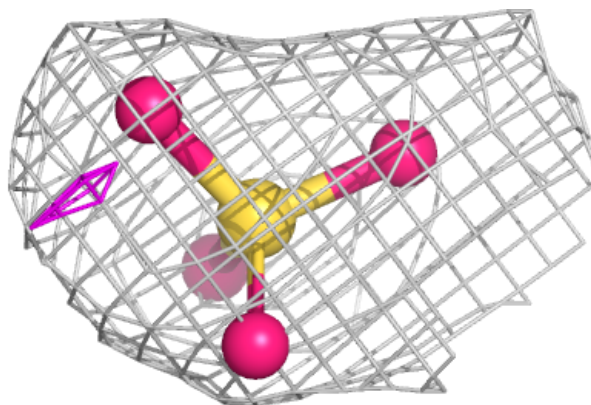
Electron density around SO4 A 1303:

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



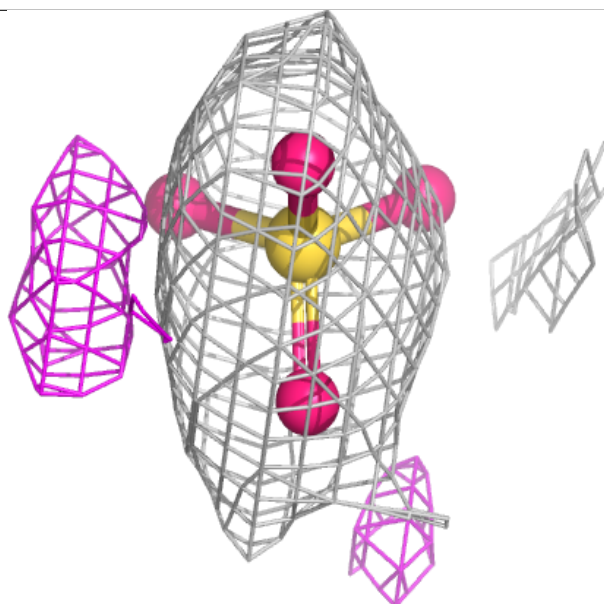
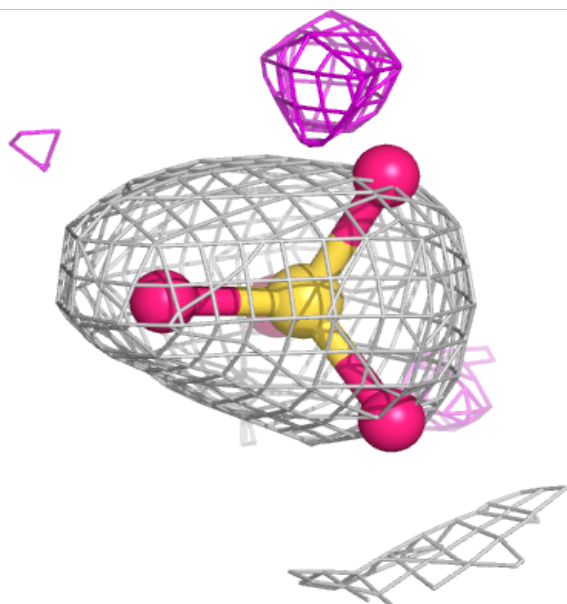
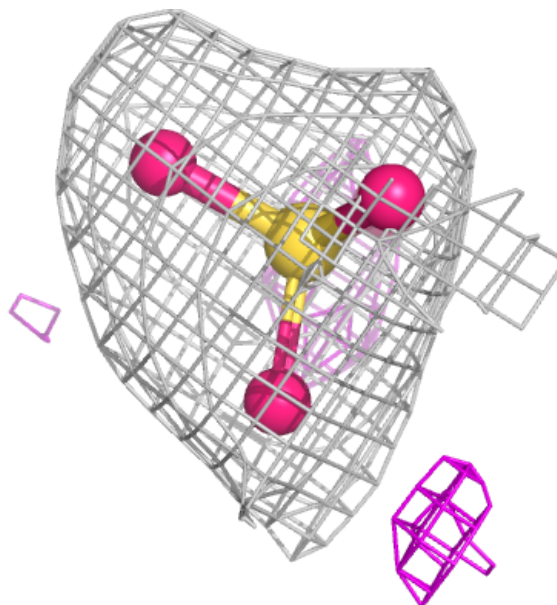
Electron density around SO4 A 1307:

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



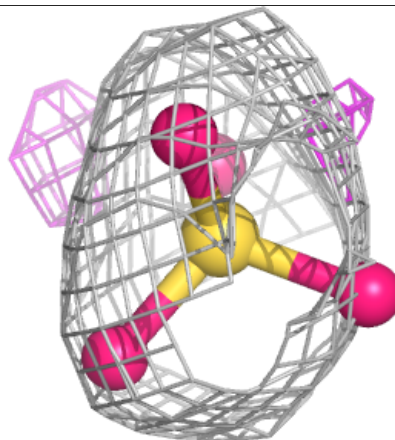
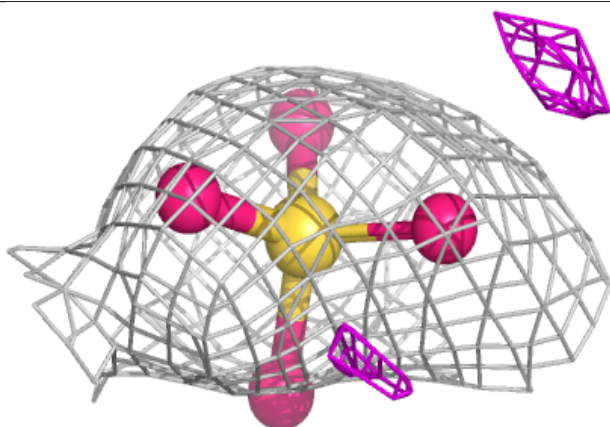
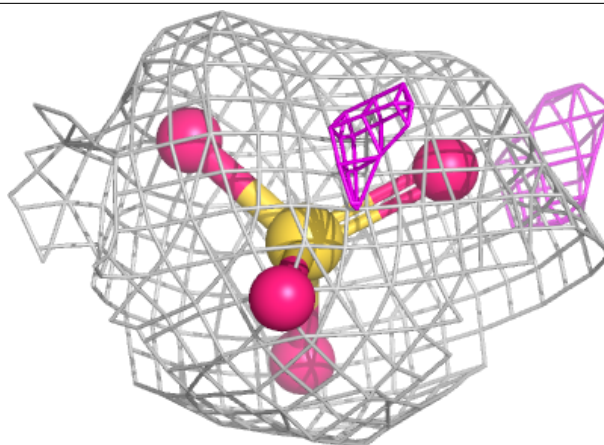
Electron density around SO4 B 1304:

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



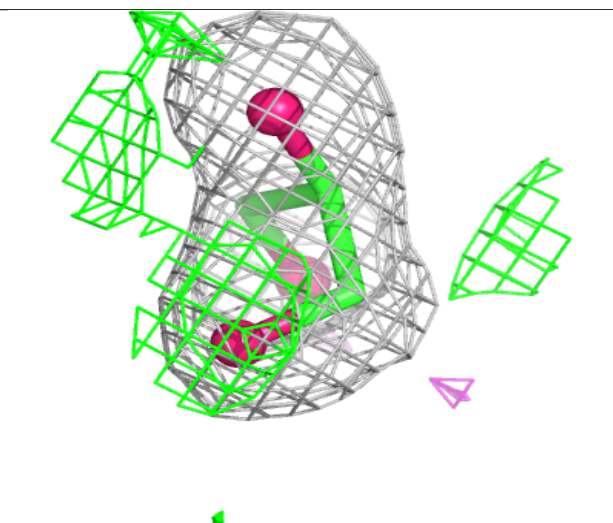
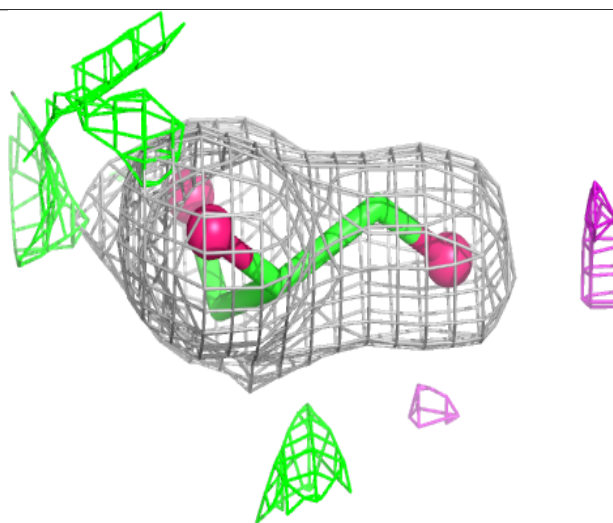
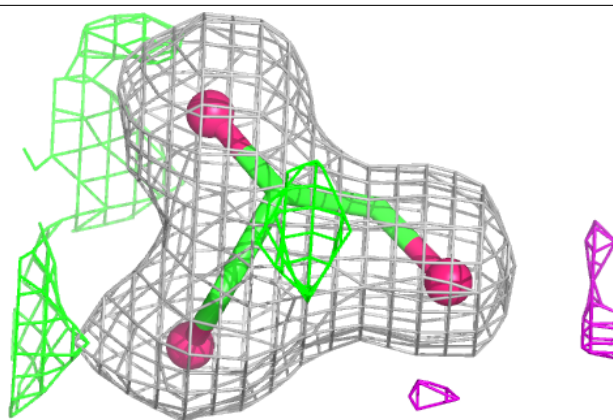
Electron density around SO4 B 1307:

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



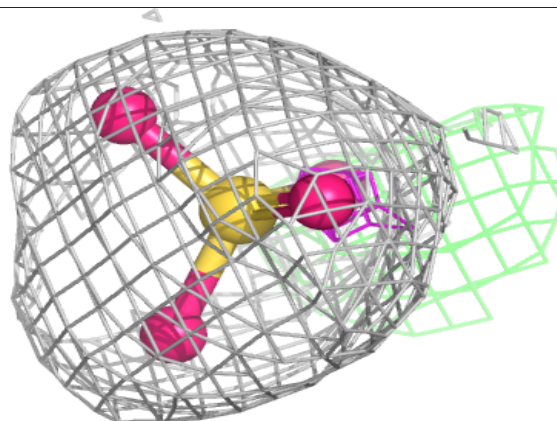
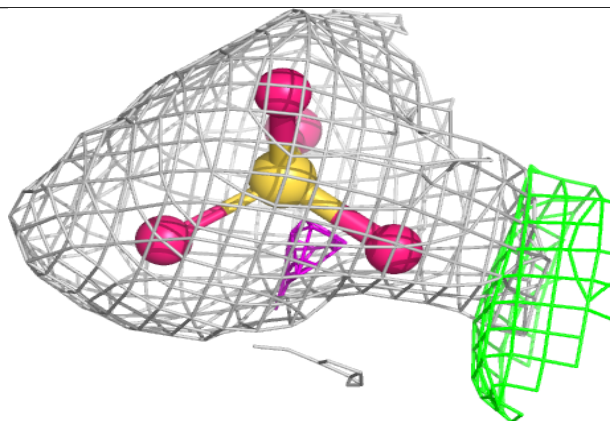
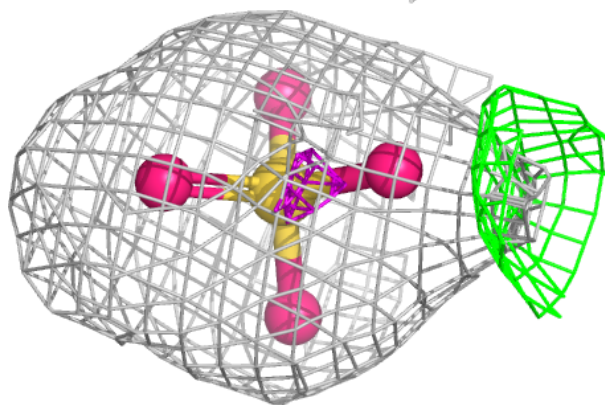
Electron density around GOL B 1301:

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



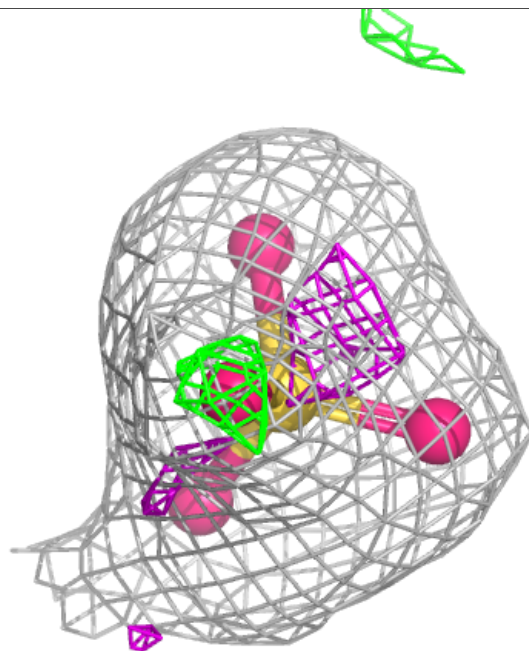
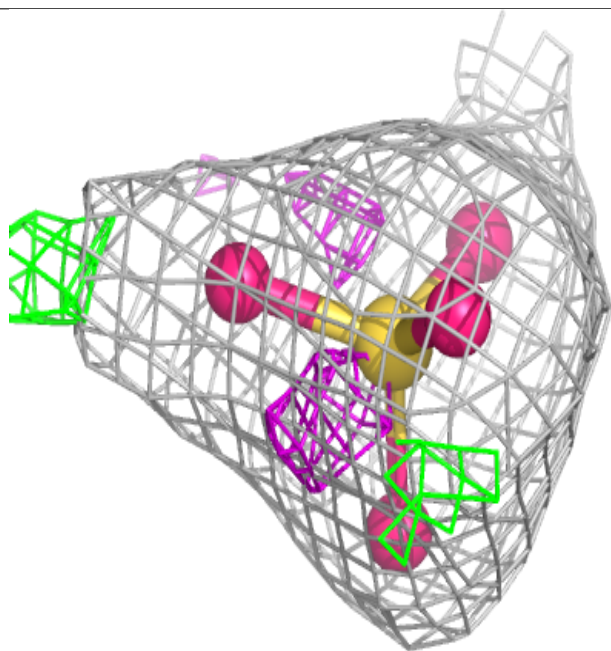
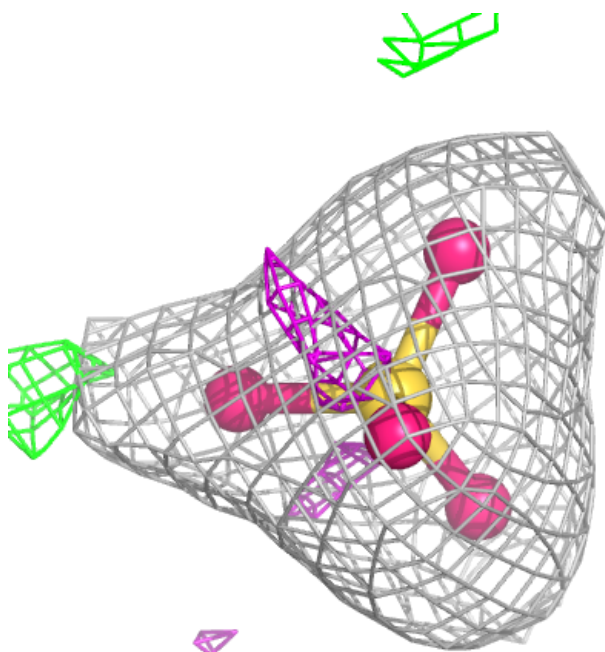
Electron density around SO4 B 1303:

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



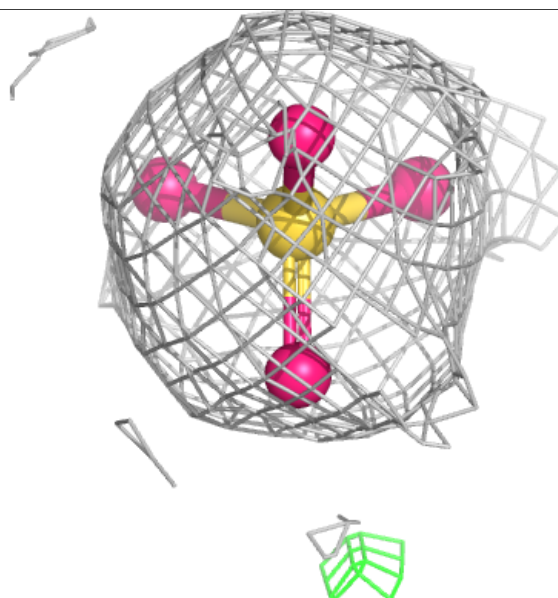
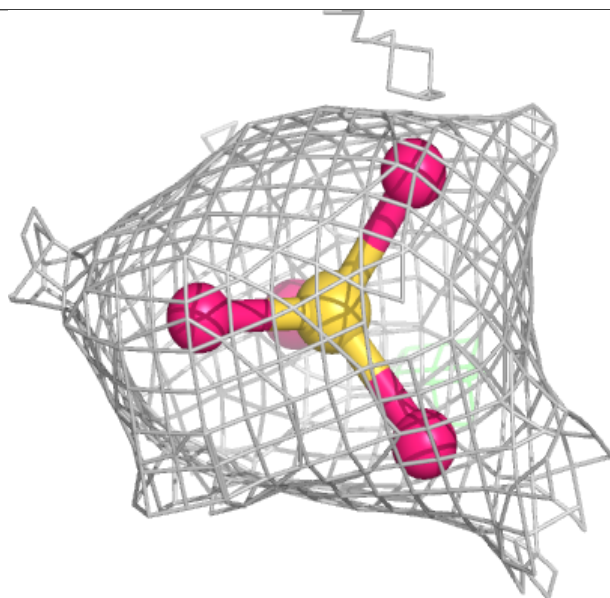
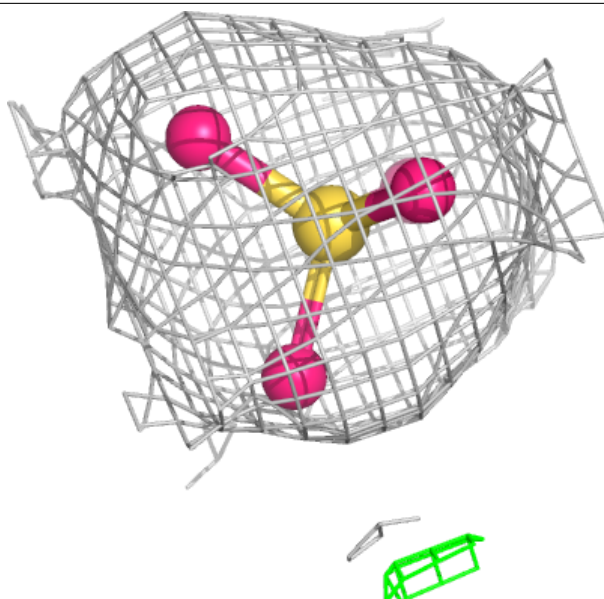
Electron density around SO4 A 1306:

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



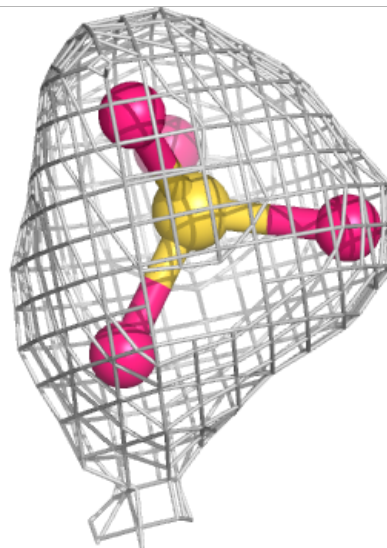
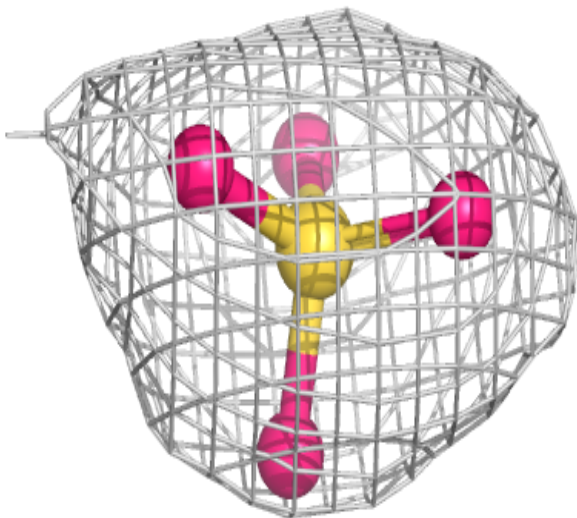
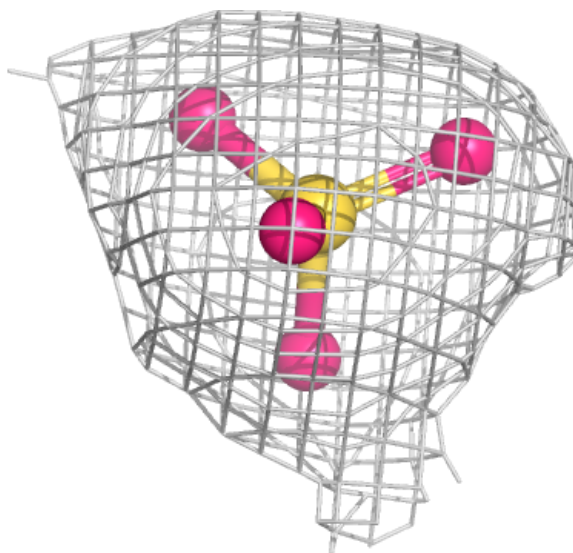
Electron density around SO4 B 1302:

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



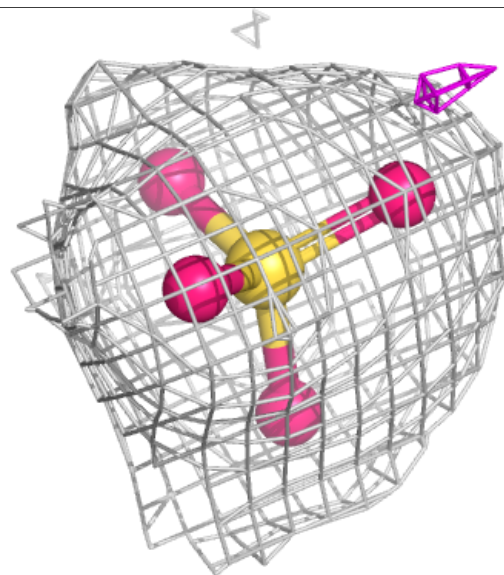
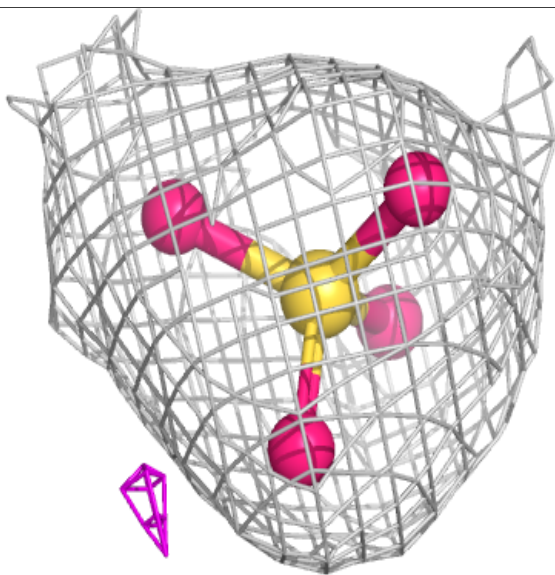
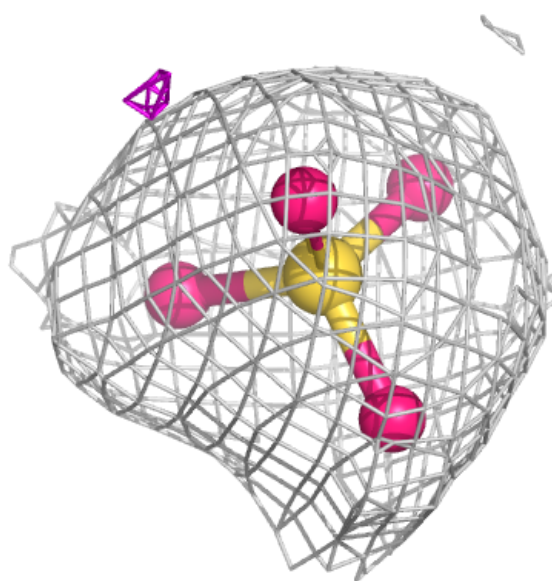
Electron density around SO4 A 1302:

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



Electron density around SO4 A 1301:

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



6.5 Other polymers [i](#)

There are no such residues in this entry.