



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 10:35 AM UTC

PDB ID : 6LBG / pdb_00006lbg
Title : Structure of OR51B2 bound FEM1C
Authors : Chen, X.; Liao, S.; Xu, C.
Deposited on : 2019-11-14
Resolution : 2.51 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

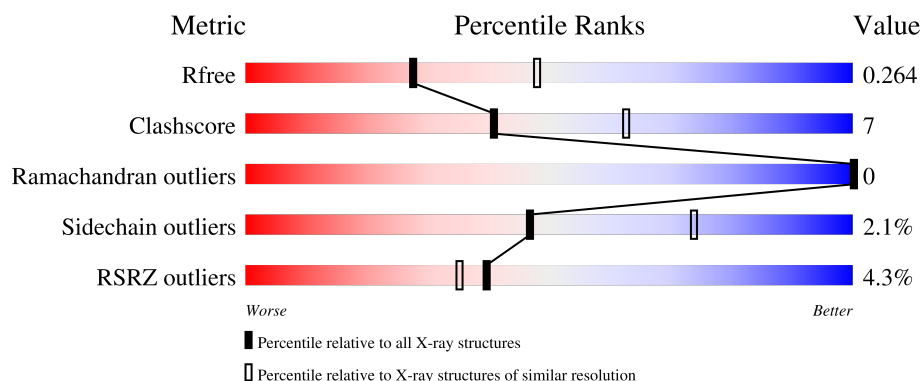
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.51 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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	418	5% 77% 12% • 10%
1	B	418	3% 77% 13% • 9%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5804 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 Protein fem-1 homolog C,Peptide from Olfactory receptor 51B2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	382	Total	C	N	O	S	0	0	0
			2869	1817	494	537	21			
1	A	377	Total	C	N	O	S	0	1	0
			2822	1788	485	528	21			

There are 38 discrepancies between the modelled and reference sequences:

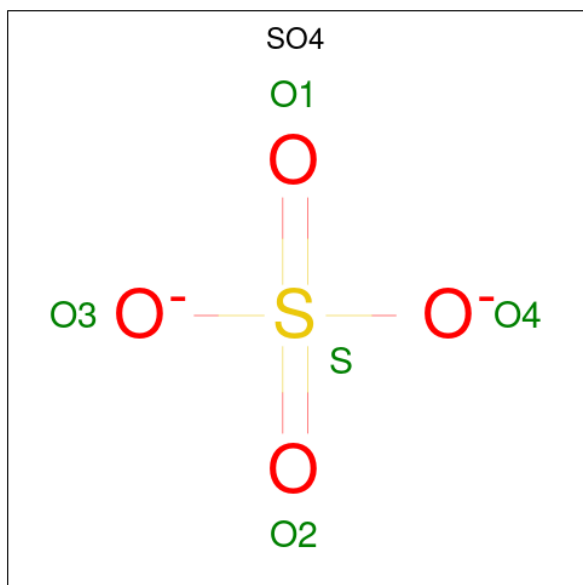
Chain	Residue	Modelled	Actual	Comment	Reference
B	-1	GLY	-	expression tag	UNP Q96JP0
B	0	HIS	-	expression tag	UNP Q96JP0
B	391	GLY	-	linker	UNP Q96JP0
B	392	GLY	-	linker	UNP Q96JP0
B	393	GLY	-	linker	UNP Q96JP0
B	394	SER	-	linker	UNP Q96JP0
B	395	GLY	-	linker	UNP Q96JP0
B	396	GLY	-	linker	UNP Q96JP0
B	397	GLY	-	linker	UNP Q96JP0
B	398	SER	-	linker	UNP Q96JP0
B	399	GLY	-	linker	UNP Q96JP0
B	400	GLY	-	linker	UNP Q96JP0
B	401	GLY	-	linker	UNP Q96JP0
B	402	SER	-	linker	UNP Q96JP0
B	403	GLY	-	linker	UNP Q96JP0
B	404	GLY	-	linker	UNP Q96JP0
B	405	GLY	-	linker	UNP Q96JP0
B	406	SER	-	linker	UNP Q96JP0
B	416	ARG	SER	variant	UNP Q9Y5P1
A	-1	GLY	-	expression tag	UNP Q96JP0
A	0	HIS	-	expression tag	UNP Q96JP0
A	391	GLY	-	linker	UNP Q96JP0
A	392	GLY	-	linker	UNP Q96JP0
A	393	GLY	-	linker	UNP Q96JP0

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Chain	Residue	Modelled	Actual	Comment	Reference
A	394	SER	-	linker	UNP Q96JP0
A	395	GLY	-	linker	UNP Q96JP0
A	396	GLY	-	linker	UNP Q96JP0
A	397	GLY	-	linker	UNP Q96JP0
A	398	SER	-	linker	UNP Q96JP0
A	399	GLY	-	linker	UNP Q96JP0
A	400	GLY	-	linker	UNP Q96JP0
A	401	GLY	-	linker	UNP Q96JP0
A	402	SER	-	linker	UNP Q96JP0
A	403	GLY	-	linker	UNP Q96JP0
A	404	GLY	-	linker	UNP Q96JP0
A	405	GLY	-	linker	UNP Q96JP0
A	406	SER	-	linker	UNP Q96JP0
A	416	ARG	SER	variant	UNP Q9Y5P1

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



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

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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		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	47	Total	O	0	0
			47	47		
3	A	36	Total	O	0	0
			36	36		

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	87.92Å 97.69Å 146.20Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.55 – 2.51 19.55 – 2.51	Depositor EDS
% Data completeness (in resolution range)	99.9 (19.55-2.51) 99.6 (19.55-2.51)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.09 (at 2.50Å)	Xtriage
Refinement program	PHENIX 1.14_3260	Depositor
R, R_{free}	0.196 , 0.231 (Not available) , 0.264	Depositor DCC
R_{free} test set	2204 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	47.3	Xtriage
Anisotropy	0.938	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 46.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	5804	wwPDB-VP
Average B, all atoms (Å ²)	64.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.88% 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: 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	0.38	0/2875	0.51	0/3892
1	B	0.41	0/2920	0.53	0/3953
All	All	0.39	0/5795	0.52	0/7845

There are no bond length outliers.

There are no bond angle outliers.

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	2822	0	2708	34	0
1	B	2869	0	2774	40	0
2	A	15	0	0	1	0
2	B	15	0	0	0	0
3	A	36	0	0	0	0
3	B	47	0	0	2	0
All	All	5804	0	5482	73	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 7.

All (73) 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:186:LEU:HD13	1:A:208:MET:HE3	1.70	0.72
1:B:169:LEU:CD2	1:B:175:VAL:HG12	2.23	0.68
1:B:163:GLU:H	1:B:163:GLU:CD	2.03	0.67
1:B:15:LYS:NZ	3:B:602:HOH:O	2.29	0.65
1:B:245:ILE:HG13	1:B:275:MET:HE2	1.81	0.62
1:B:283:ILE:H	1:B:283:ILE:HD12	1.65	0.61
1:A:290:GLN:H	1:A:290:GLN:CD	2.06	0.61
1:B:126:ASP:OD2	1:B:416:ARG:NH1	2.34	0.60
1:A:7:VAL:CG1	1:A:8:PHE:N	2.64	0.60
1:B:169:LEU:HD21	1:B:175:VAL:HG12	1.83	0.59
1:A:236:HIS:O	1:A:244:ARG:NH1	2.35	0.59
1:B:83:GLY:HA3	1:B:114:LEU:HG	1.88	0.54
1:B:220:SER:O	1:B:224:THR:HG23	2.07	0.54
1:B:319:ARG:HD2	1:B:344:ARG:CZ	2.37	0.54
1:A:52:GLY:HA2	1:A:98:VAL:HG21	1.89	0.53
1:A:7:VAL:HG13	1:A:8:PHE:N	2.22	0.53
1:A:186:LEU:HD13	1:A:208:MET:CE	2.38	0.53
1:B:16:LEU:HA	1:B:56:MET:HE1	1.91	0.53
1:A:112:THR:HG22	1:A:118:THR:HG22	1.91	0.53
1:B:68:ILE:HG23	1:B:69:GLU:HG2	1.90	0.52
1:B:416:ARG:OXT	1:B:416:ARG:HG3	2.10	0.52
1:B:52:GLY:HA2	1:B:98:VAL:HG21	1.92	0.52
1:B:315:PRO:O	1:B:319:ARG:HG3	2.11	0.51
1:A:291:THR:HG22	1:A:292:LEU:O	2.09	0.51
1:B:346:ALA:HB1	1:A:386:PHE:HB3	1.93	0.50
1:B:7:VAL:HG13	1:B:19:LEU:HD11	1.93	0.50
1:A:68:ILE:HG23	1:A:69:GLU:HG2	1.94	0.50
1:A:342:ARG:NH1	1:A:368:GLN:OE1	2.44	0.50
1:A:223:VAL:HG23	1:A:250:LEU:HD21	1.94	0.50
1:A:100:GLN:HG2	1:A:134:TYR:CZ	2.47	0.49
1:A:56:MET:O	1:A:60:LEU:HG	2.14	0.48
1:B:100:GLN:HG2	1:B:134:TYR:CZ	2.49	0.48
1:B:343:TYR:OH	3:B:601:HOH:O	2.20	0.48
1:A:35:ILE:O	1:A:44:PRO:HD2	2.14	0.48
1:A:200:MET:HA	1:A:203:MET:HE2	1.96	0.48
1:B:244:ARG:HB2	1:B:275:MET:HE1	1.95	0.47
1:B:169:LEU:HD11	1:B:201:LEU:HD23	1.97	0.47
1:A:367:MET:HG3	1:A:368:GLN:N	2.29	0.47
1:A:73:SER:OG	1:A:82:GLU:OE1	2.31	0.47
1:A:208:MET:HE2	1:A:217:PRO:HG2	1.97	0.47
1:B:311:LEU:HD23	1:B:311:LEU:HA	1.76	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:342:ARG:HD2	1:A:384:LEU:HD23	1.97	0.46
1:A:317:GLU:O	1:A:321:GLN:HG2	2.16	0.46
1:B:18:LEU:O	1:B:22:LEU:HG	2.15	0.46
1:A:59:PHE:O	1:A:63:GLN:HG2	2.17	0.45
1:B:82:GLU:HB2	1:B:114:LEU:HD12	1.98	0.45
1:A:16:LEU:N	2:A:501:SO4:O3	2.50	0.44
1:B:266:LEU:HD12	1:B:266:LEU:HA	1.90	0.44
1:B:215:MET:HE1	1:B:223:VAL:HG21	1.98	0.44
1:B:186:LEU:HD13	1:B:208:MET:HG3	1.99	0.44
1:A:84:ALA:HA	1:A:113:THR:HG22	1.99	0.44
1:B:195:LEU:HG	1:B:199:LYS:HE3	2.00	0.44
1:B:349:ALA:HB2	1:B:357:CYS:SG	2.59	0.43
1:A:38:LYS:CB	1:A:44:PRO:HD3	2.49	0.43
1:A:159:LYS:HA	1:A:159:LYS:HD2	1.82	0.43
1:B:301:GLU:HA	1:B:321:GLN:OE1	2.19	0.43
1:A:46:LEU:HD21	1:A:74:VAL:HG13	2.00	0.43
1:B:83:GLY:O	1:B:113:THR:HA	2.19	0.42
1:B:179:SER:OG	1:B:183:ASN:HB3	2.19	0.42
1:A:272:ALA:O	1:A:276:ARG:HG3	2.18	0.42
1:B:151:THR:OG1	1:B:154:MET:HG3	2.19	0.42
1:A:291:THR:CG2	1:A:292:LEU:N	2.83	0.41
1:A:263:LEU:H	1:A:263:LEU:HD12	1.85	0.41
1:B:32:SER:HB3	1:B:64:CYS:SG	2.61	0.41
1:A:125:PHE:O	1:A:159:LYS:HG3	2.21	0.41
1:B:125:PHE:CD2	1:B:416:ARG:HA	2.56	0.41
1:B:215:MET:HE3	1:B:215:MET:HB3	1.92	0.41
1:A:167:TYR:O	1:A:171:LYS:HG2	2.21	0.41
1:B:127:GLY:HA2	1:B:164:ILE:CD1	2.50	0.41
1:A:276:ARG:HB3	1:A:284:ILE:O	2.21	0.40
1:B:56:MET:HE2	1:B:56:MET:HB2	1.63	0.40
1:B:94:GLY:HA2	1:B:131:ILE:HD12	2.03	0.40
1:B:317:GLU:O	1:B:321:GLN:HG2	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

of similar resolution.

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	370/418 (88%)	358 (97%)	12 (3%)	0	100	100
1	B	376/418 (90%)	362 (96%)	14 (4%)	0	100	100
All	All	746/836 (89%)	720 (96%)	26 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	279/343 (81%)	272 (98%)	7 (2%)	42	69
1	B	287/343 (84%)	282 (98%)	5 (2%)	53	78
All	All	566/686 (82%)	554 (98%)	12 (2%)	47	74

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	3	LEU
1	B	22	LEU
1	B	32	SER
1	B	163	GLU
1	B	183	ASN
1	A	3	LEU
1	A	208	MET
1	A	284	ILE
1	A	290	GLN
1	A	311	LEU
1	A	316	ASP
1	A	336	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such

sidechains are listed below:

Mol	Chain	Res	Type
1	B	303	ASN
1	B	371	ASN
1	A	104	ASN
1	A	183	ASN
1	A	353	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

6 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	$\# Z > 2$
2	SO4	B	502	-	4,4,4	0.27	0	6,6,6	0.16	0
2	SO4	A	502	-	4,4,4	0.24	0	6,6,6	0.18	0
2	SO4	A	501	-	4,4,4	0.23	0	6,6,6	0.12	0
2	SO4	B	503	-	4,4,4	0.25	0	6,6,6	0.19	0
2	SO4	A	503	-	4,4,4	0.22	0	6,6,6	0.14	0
2	SO4	B	501	-	4,4,4	0.26	0	6,6,6	0.25	0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

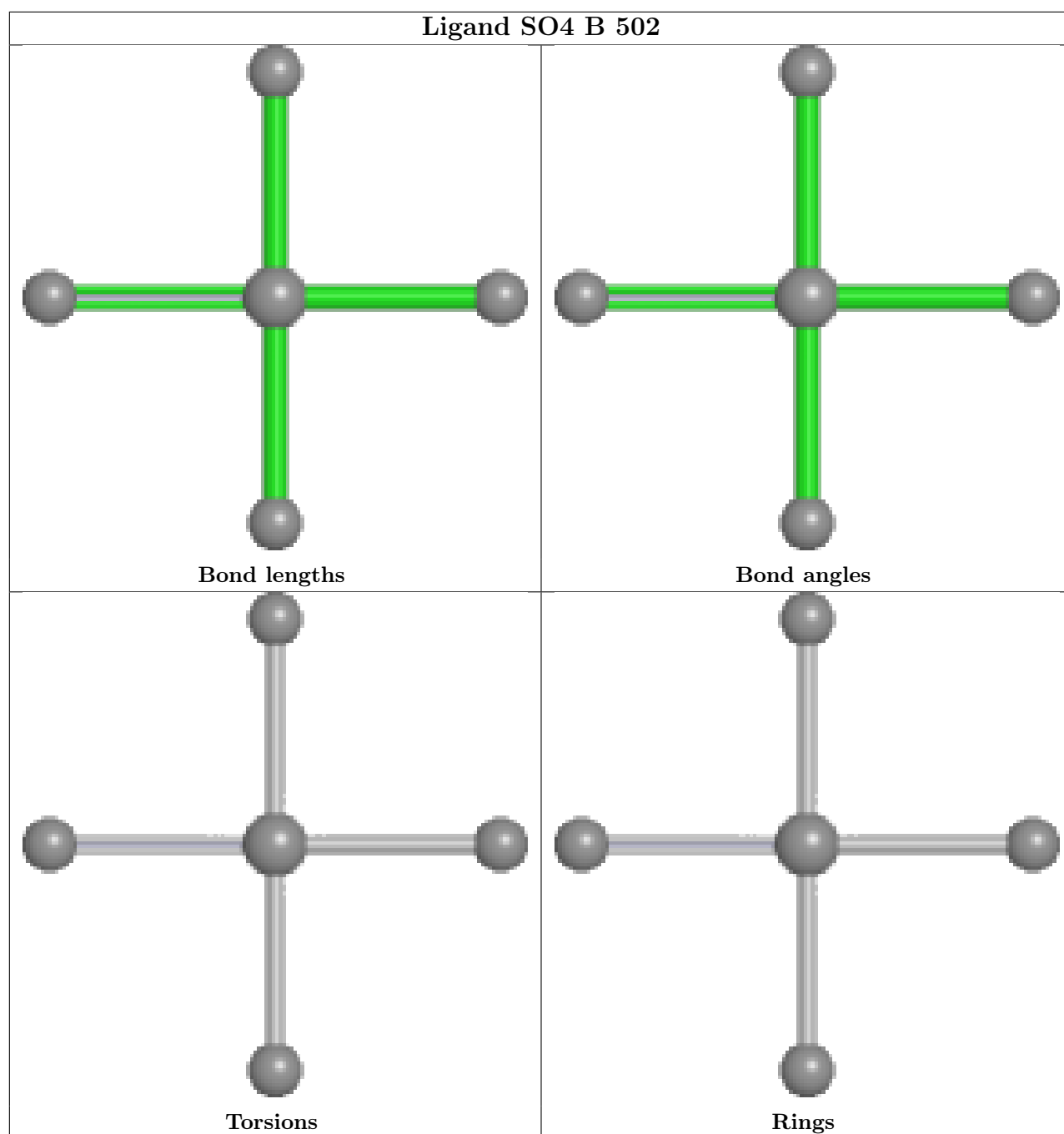
There are no torsion outliers.

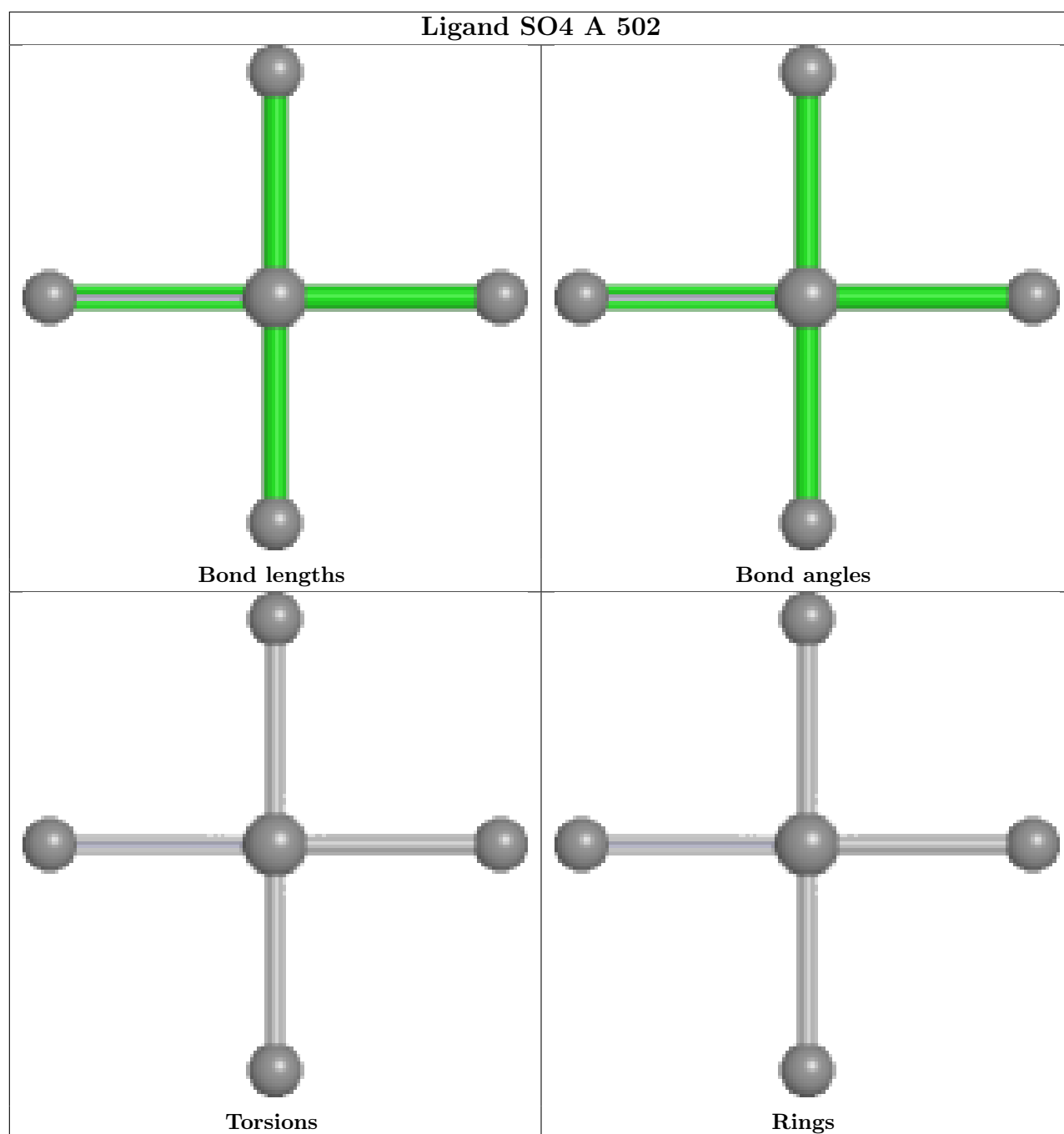
There are no ring outliers.

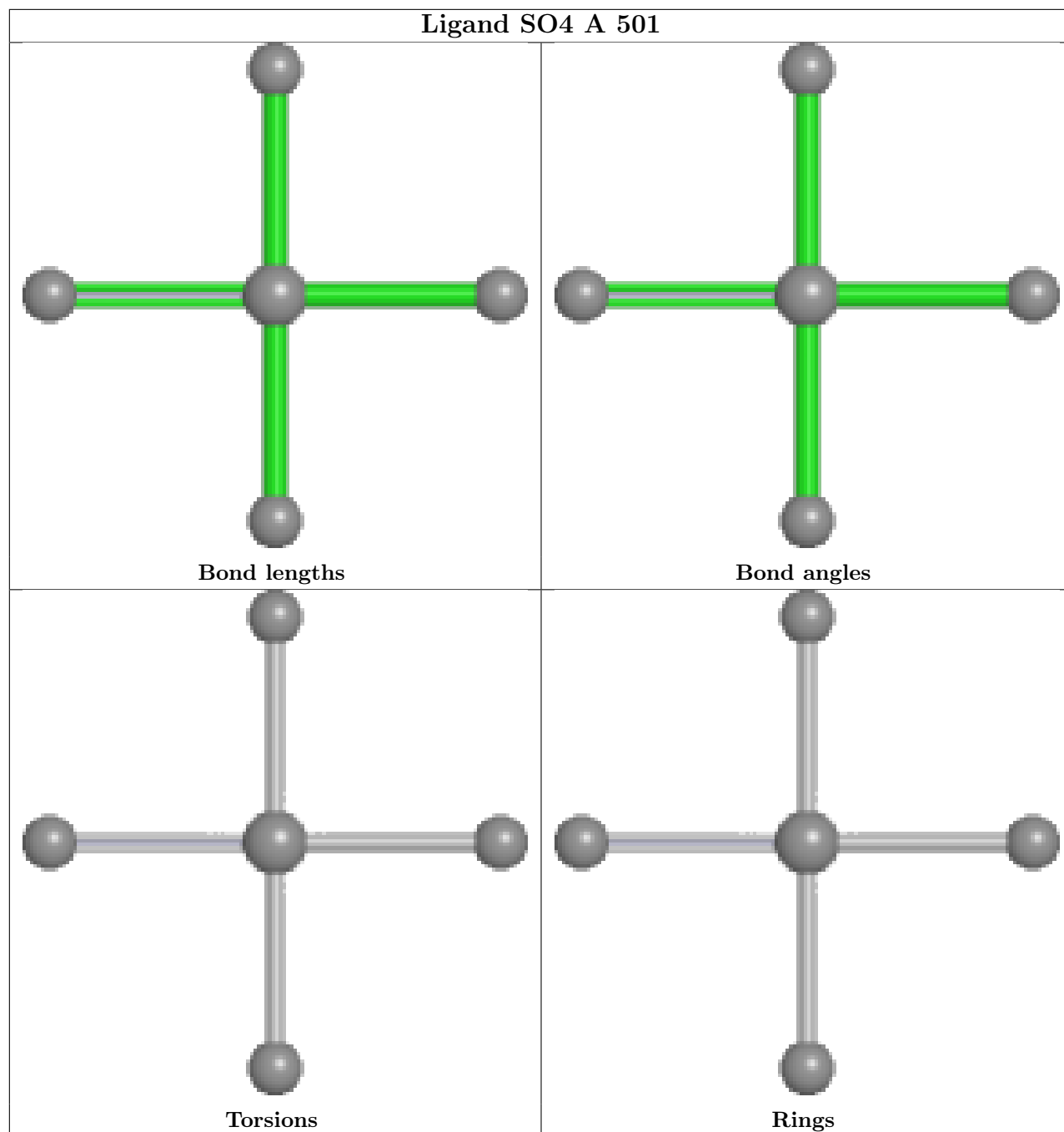
1 monomer is involved in 1 short contact:

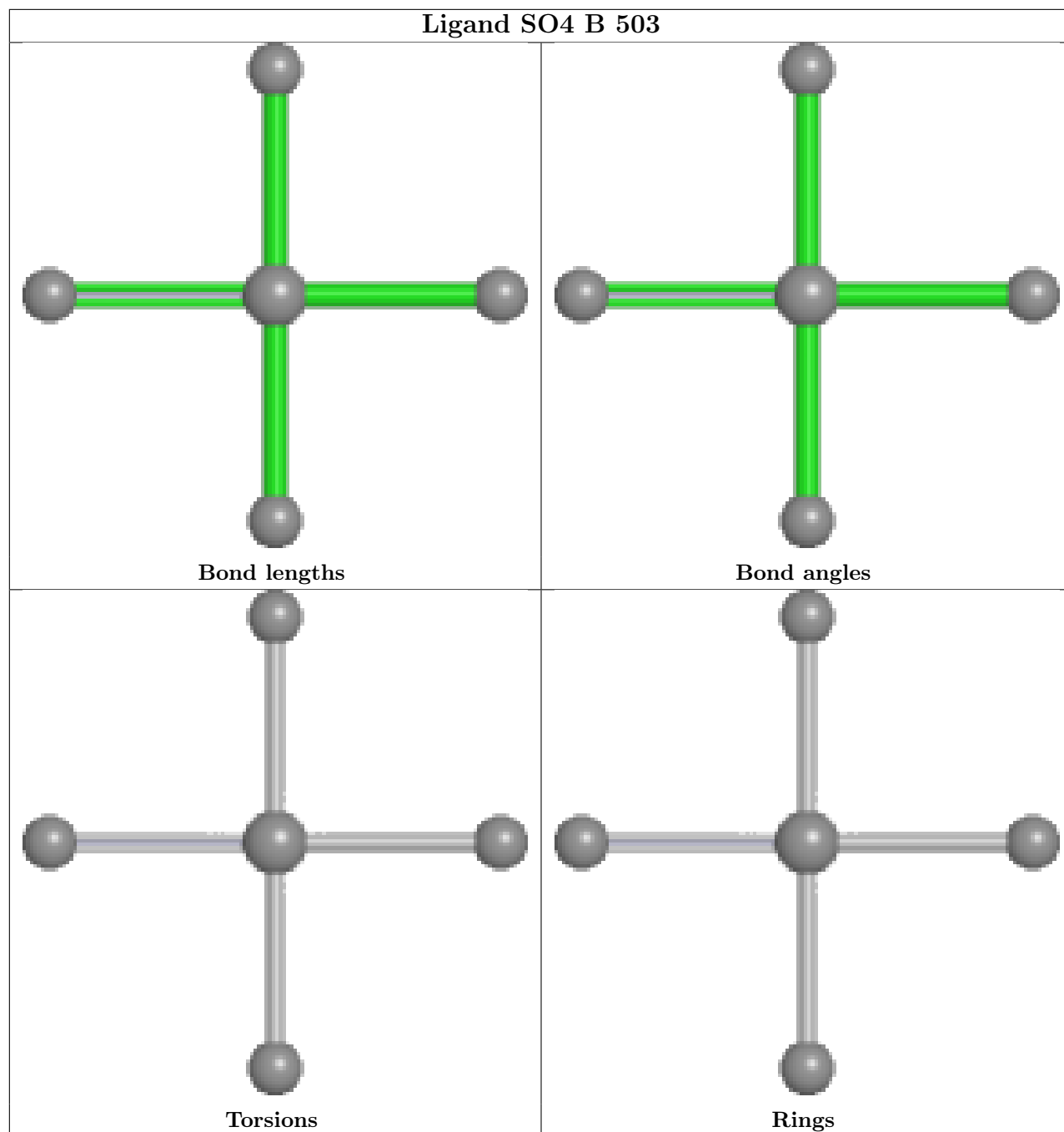
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	501	SO4	1	0

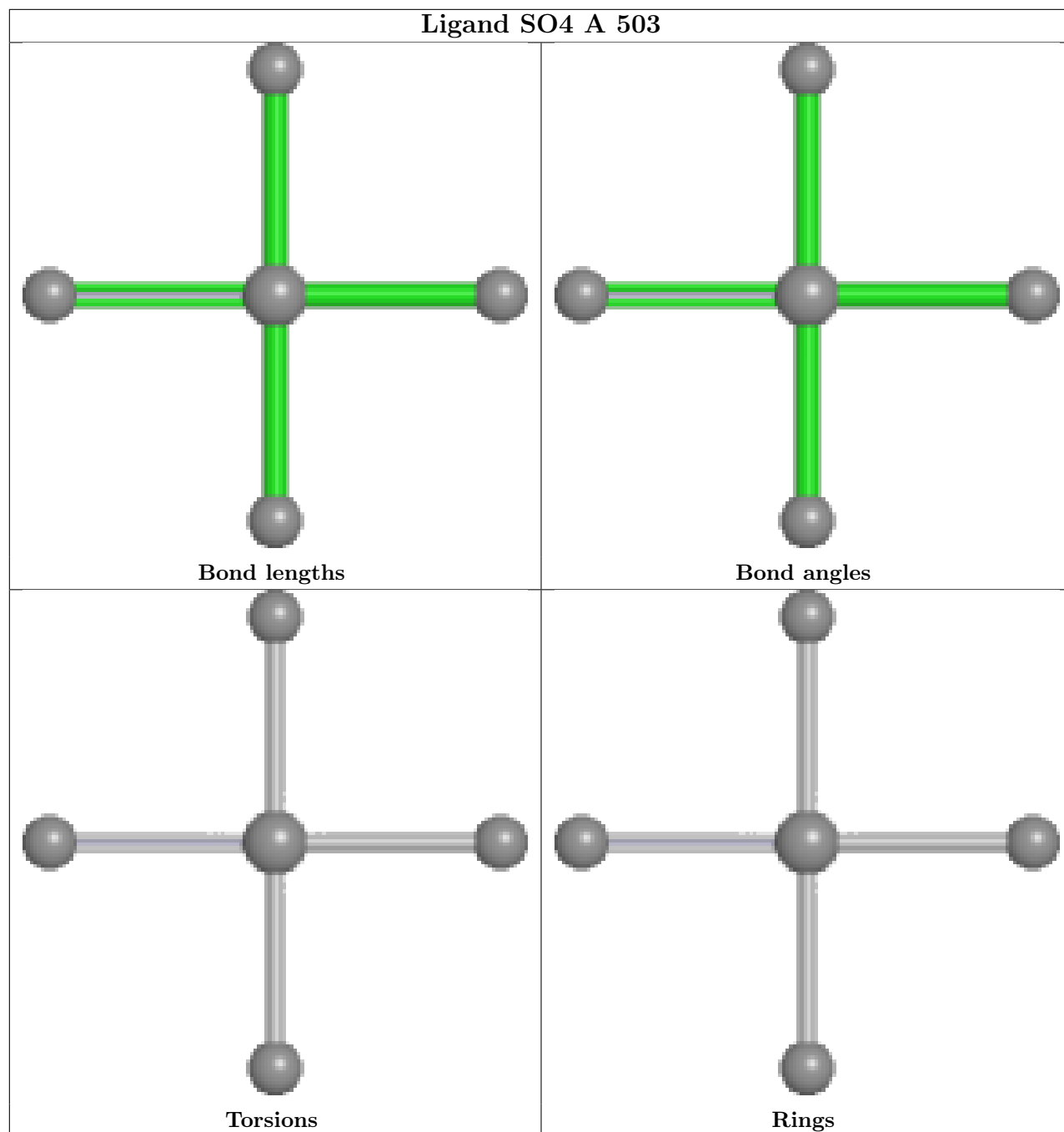
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

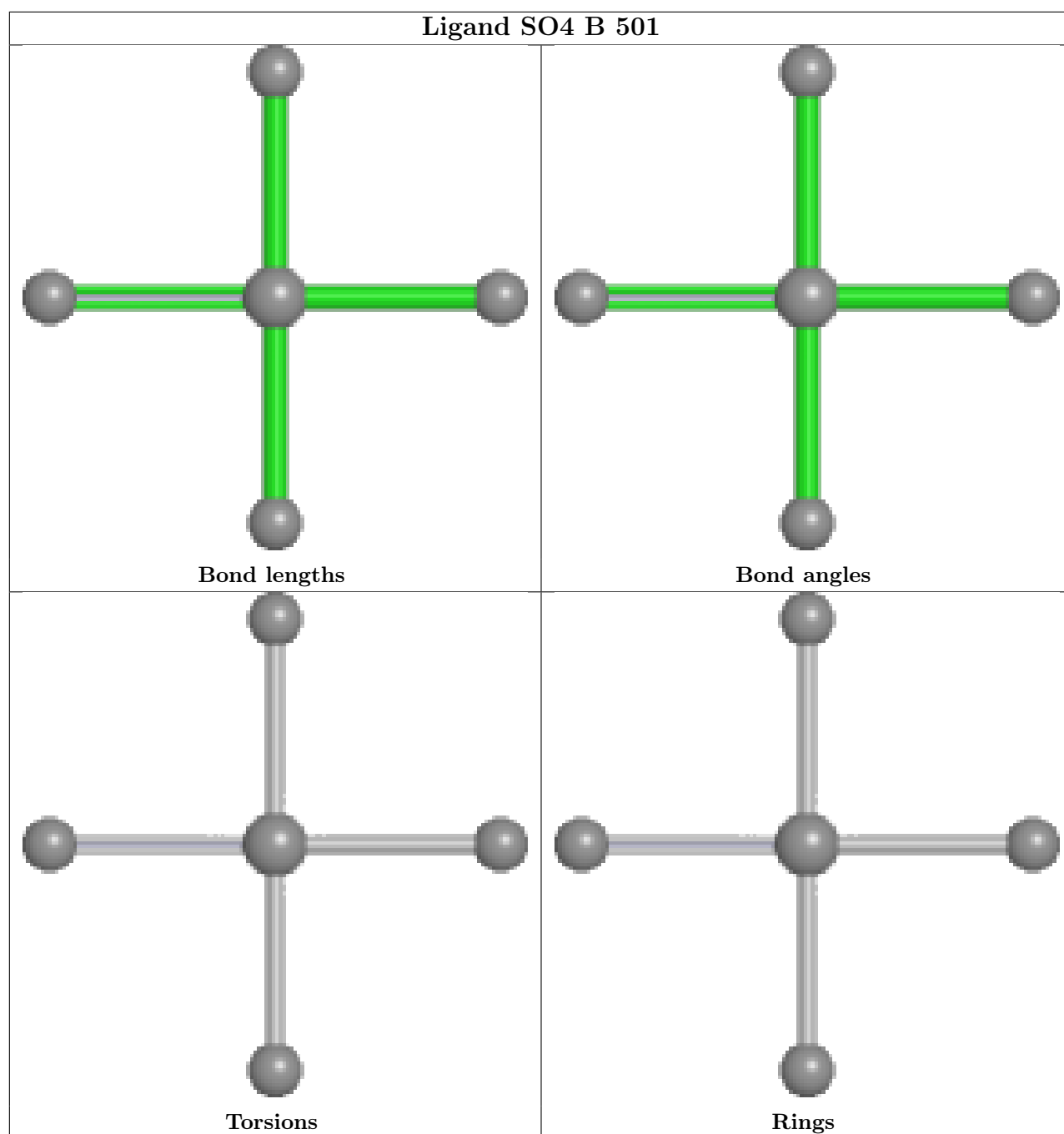












5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [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	377/418 (90%)	0.26	19 (5%) 34 30	33, 61, 105, 146	1 (0%)
1	B	382/418 (91%)	0.20	14 (3%) 45 40	40, 60, 109, 143	0
All	All	759/836 (90%)	0.23	33 (4%) 40 35	33, 60, 105, 146	1 (0%)

All (33) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	372	LEU	6.3
1	B	34	LEU	4.2
1	A	3	LEU	4.0
1	A	8	PHE	3.7
1	A	35	ILE	3.7
1	A	28	LYS	3.7
1	A	27	SER	3.7
1	B	25	SER	3.6
1	A	22	LEU	3.3
1	B	2	ASP	3.2
1	A	30	GLU	3.1
1	B	23	LEU	3.0
1	B	67	SER	2.9
1	B	415	SER	2.8
1	A	32	SER	2.8
1	A	36	SER	2.7
1	A	371	ASN	2.4
1	A	43	THR	2.4
1	B	77	ASP	2.4
1	A	34	LEU	2.3
1	B	414	PHE	2.3
1	B	384	LEU	2.2
1	A	15	LYS	2.2
1	A	31	VAL	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	412	HIS	2.2
1	B	28	LYS	2.2
1	A	80	THR	2.2
1	A	366	ASP	2.2
1	B	183	ASN	2.1
1	A	40	ASN	2.1
1	A	67	SER	2.0
1	B	306	GLU	2.0
1	B	39	THR	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](#)

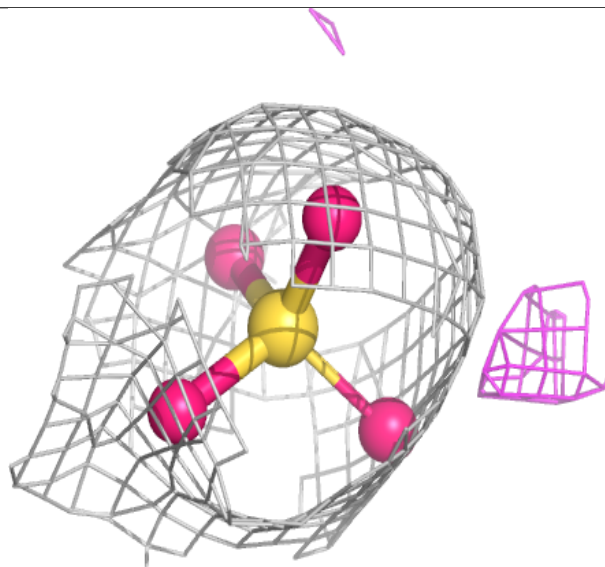
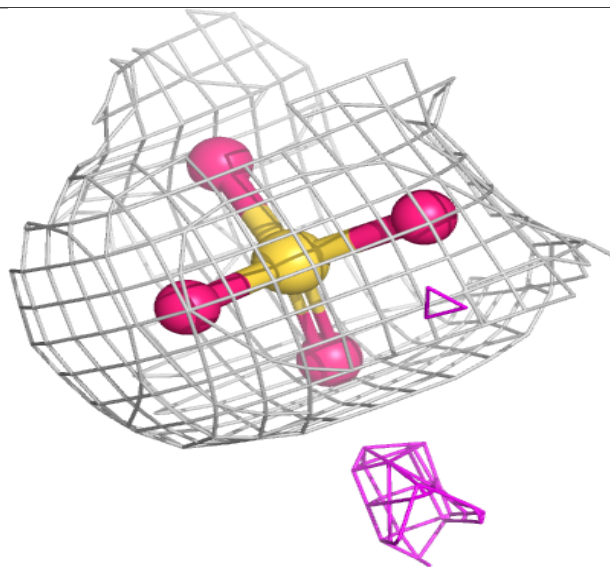
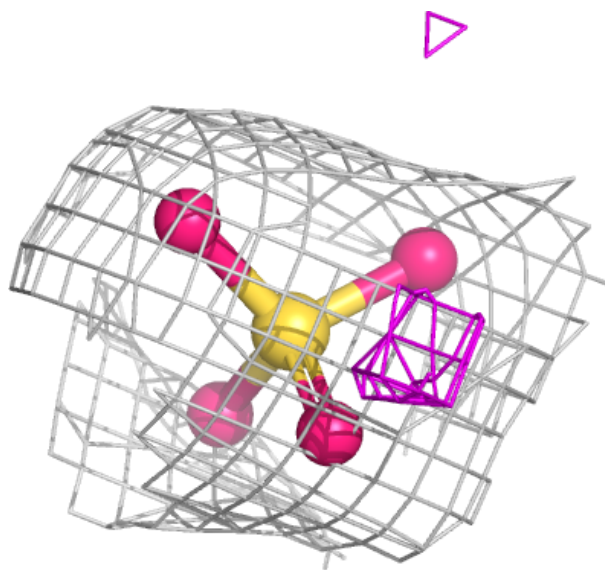
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	A	501	5/5	0.73	0.13	120,121,122,123	0
2	SO4	A	503	5/5	0.83	0.20	129,130,131,131	0
2	SO4	B	503	5/5	0.85	0.13	94,95,97,98	0
2	SO4	B	501	5/5	0.86	0.10	85,87,90,90	0
2	SO4	B	502	5/5	0.93	0.09	74,74,78,79	0
2	SO4	A	502	5/5	0.95	0.07	56,59,62,65	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.

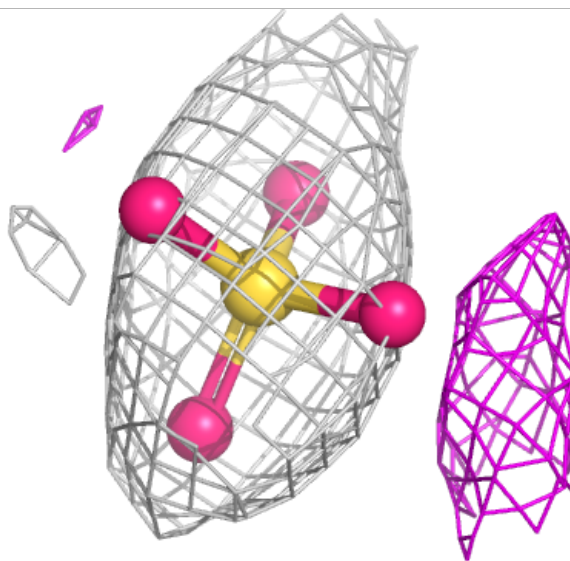
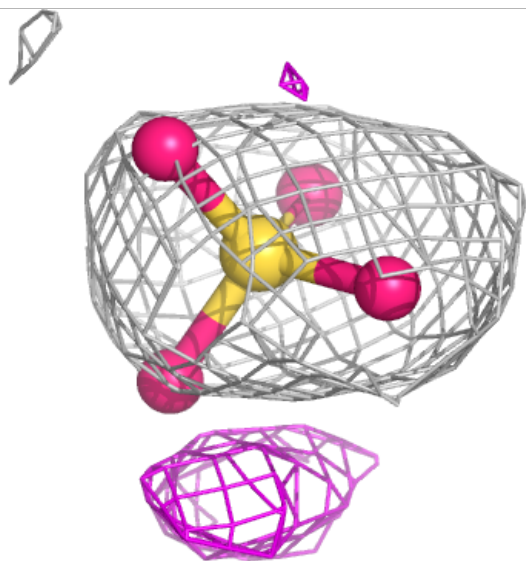
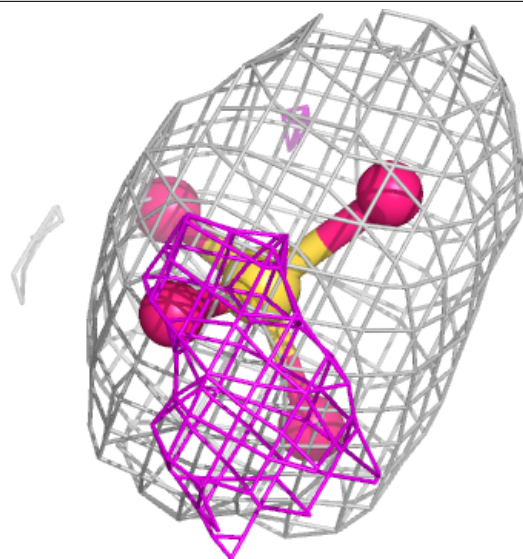
Electron density around SO4 A 501:

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



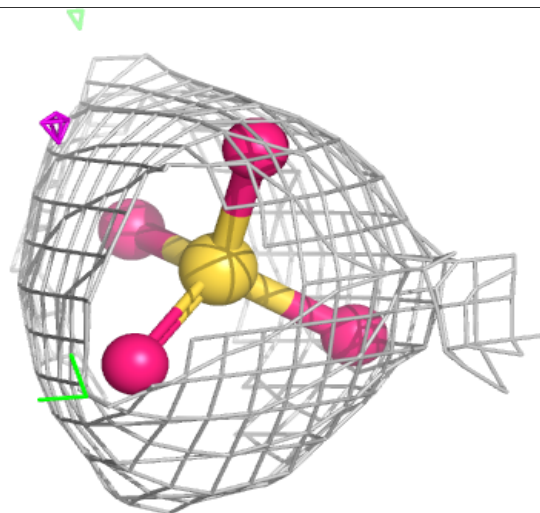
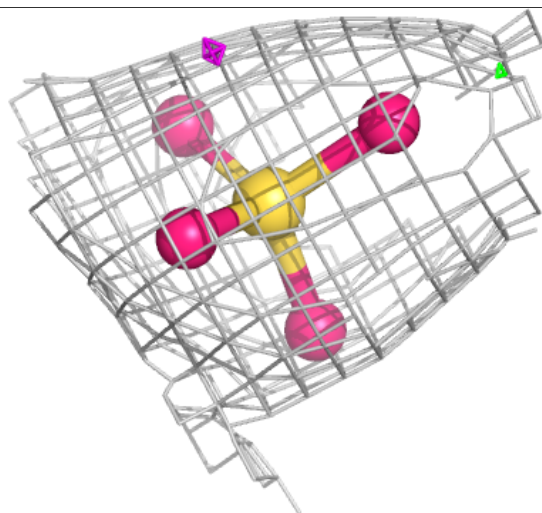
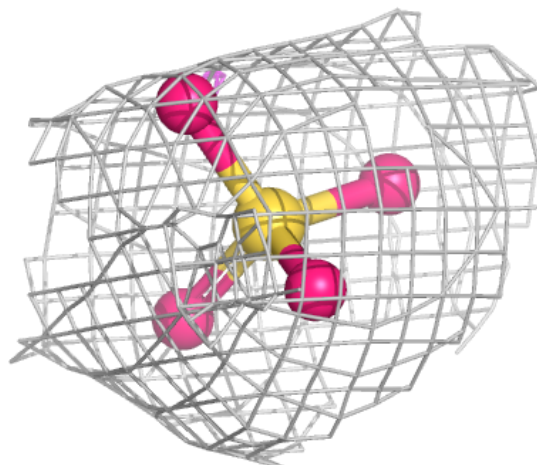
Electron density around SO4 A 503:

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



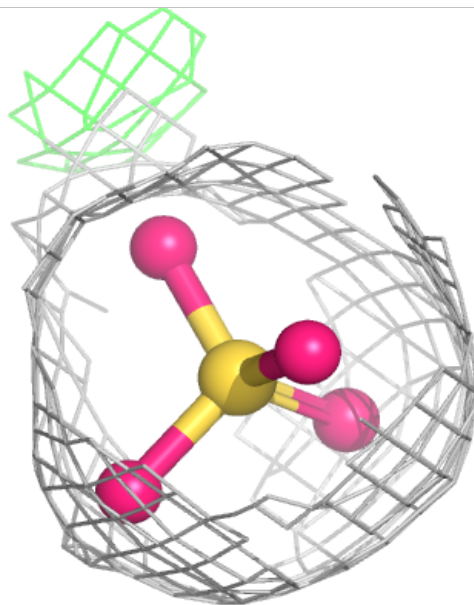
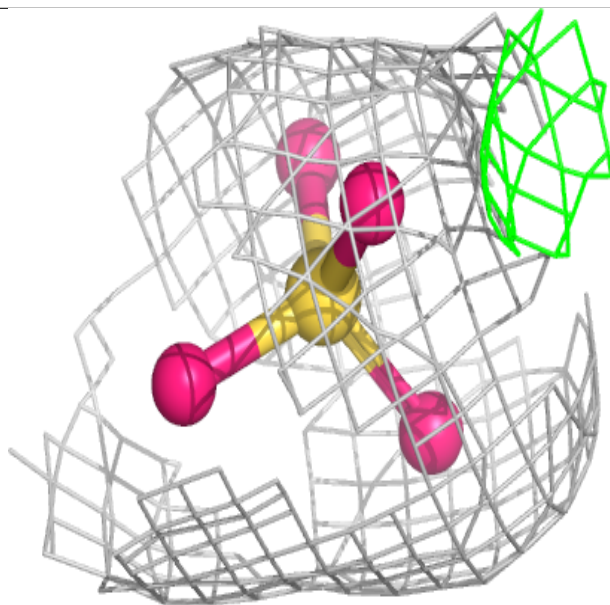
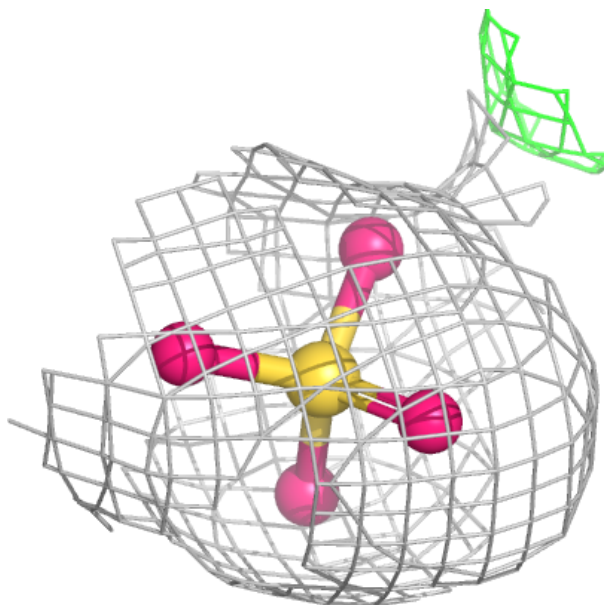
Electron density around SO4 B 503:

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



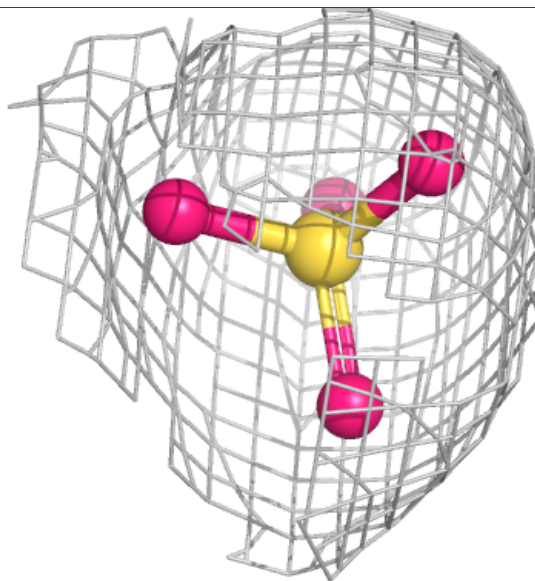
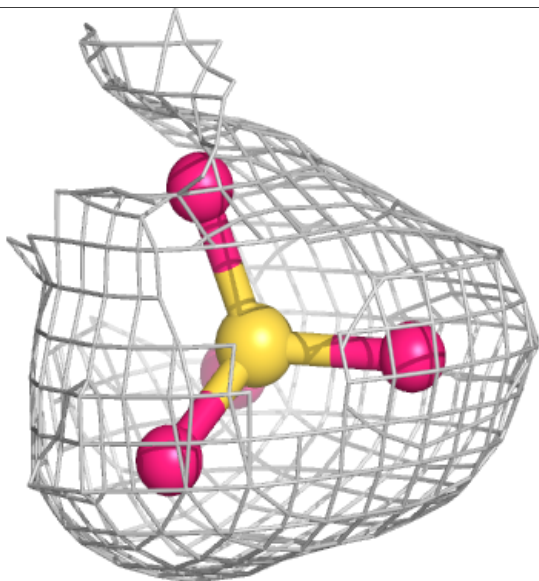
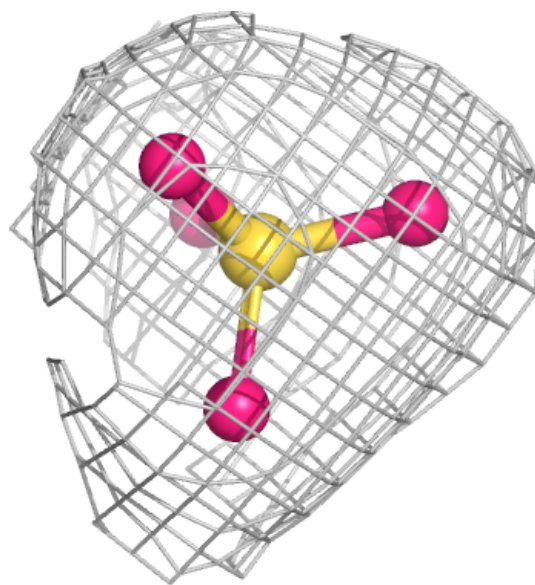
Electron density around SO4 B 501:

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and green (positive)



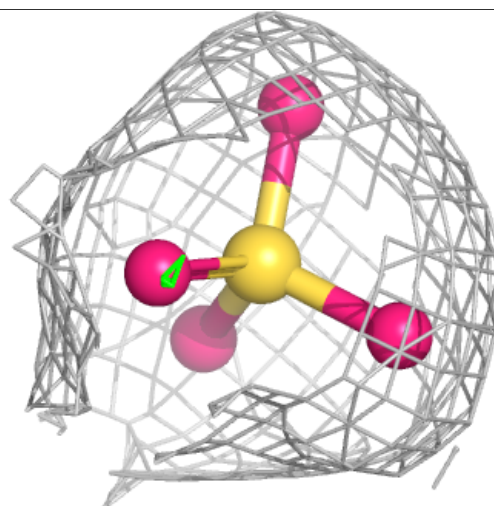
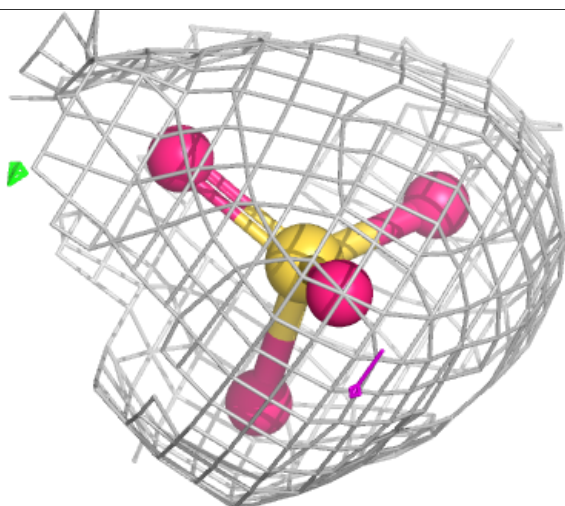
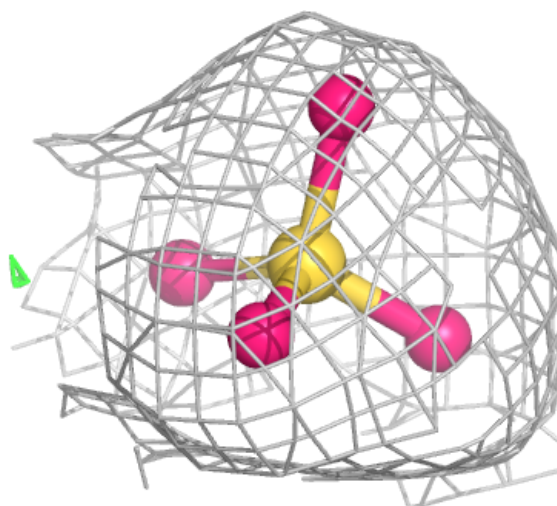
Electron density around SO4 B 502:

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



Electron density around SO4 A 502:

$2mF_o-DF_c$ (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.