



wwPDB X-ray Structure Validation Summary Report ⓘ

Jun 2, 2025 – 12:23 PM JST

PDB ID : 8ZUW / pdb_00008zuw
Title : Crystal structure of Bacteroides thetaiotaomicron glutamate decarboxylase
Authors : Liu, S.; Yulu, W.; Guoming, D.; Boting, W.; Xin, F.
Deposited on : 2024-06-10
Resolution : 2.46 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 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 Phenix2.0rc1
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 2.0rc1
EDS : 3.0
buster-report : 1.1.7 (2018)
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4 : 9.0.006 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.43.1

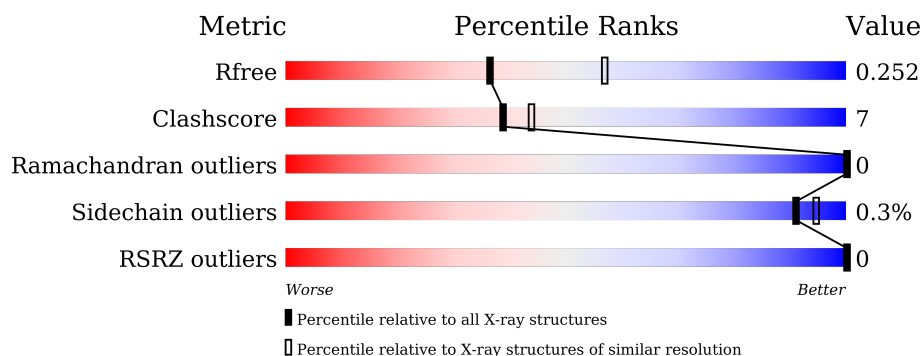
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION





The reported resolution of this entry is 2.46 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}	164625	1096 (2.46-2.46)
Clashscore	180529	1178 (2.46-2.46)
Ramachandran outliers	177936	1170 (2.46-2.46)
Sidechain outliers	177891	1170 (2.46-2.46)
RSRZ outliers	164620	1096 (2.46-2.46)

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	481	 80% 16% 5%
1	B	481	 79% 17% 5%
1	C	481	 80% 15% 5%
1	D	481	 81% 14% 5%
1	E	481	 79% 16% 5%
1	F	481	 78% 17% 5%

2 Entry composition [i](#)

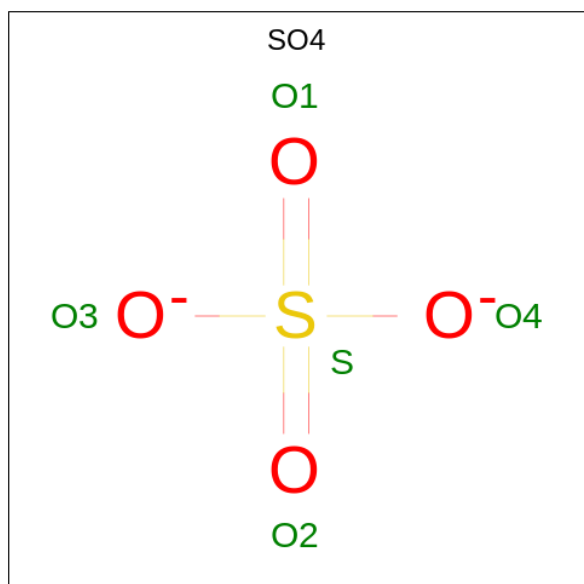
There are 3 unique types of molecules in this entry. The entry contains 22585 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Glutamate decarboxylase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	459	Total	C	N	O	S	4	0	0
			3641	2357	595	669	20			
1	B	459	Total	C	N	O	S	4	0	0
			3629	2349	593	667	20			
1	C	459	Total	C	N	O	S	4	0	0
			3635	2353	594	668	20			
1	D	459	Total	C	N	O	S	4	0	0
			3637	2354	594	669	20			
1	E	459	Total	C	N	O	S	4	0	0
			3641	2358	596	667	20			
1	F	459	Total	C	N	O	S	4	0	0
			3645	2360	596	669	20			

- 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 5 4 1	0	0
2	A	1	Total O S 5 4 1	0	0
2	B	1	Total O S 5 4 1	0	0
2	B	1	Total O S 5 4 1	0	0
2	C	1	Total O S 5 4 1	0	0
2	D	1	Total O S 5 4 1	0	0
2	E	1	Total O S 5 4 1	0	0
2	F	1	Total O S 5 4 1	0	0

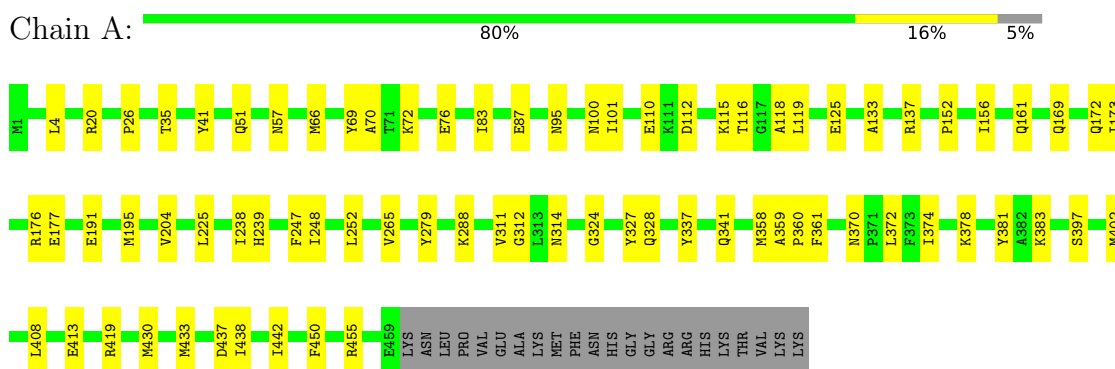
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	121	Total O 121 121	0	0
3	B	144	Total O 144 144	0	0
3	C	124	Total O 124 124	0	0
3	D	103	Total O 103 103	0	0
3	E	118	Total O 118 118	0	0
3	F	107	Total O 107 107	0	0

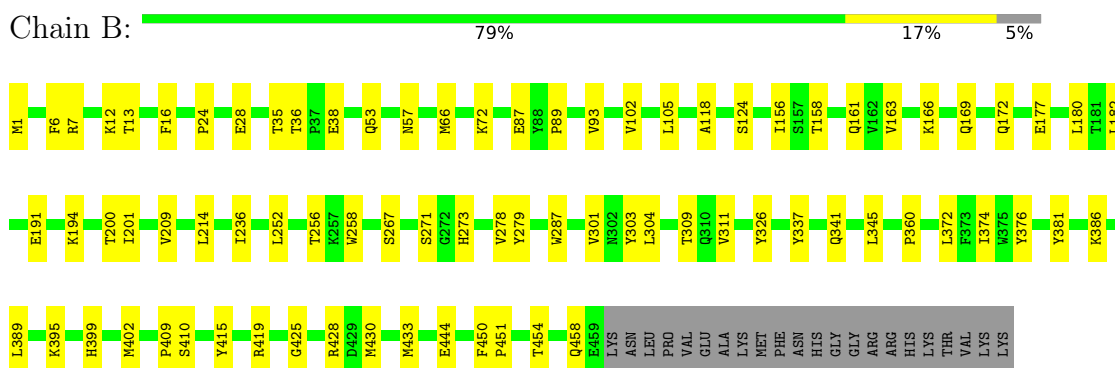
3 Residue-property plots

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.

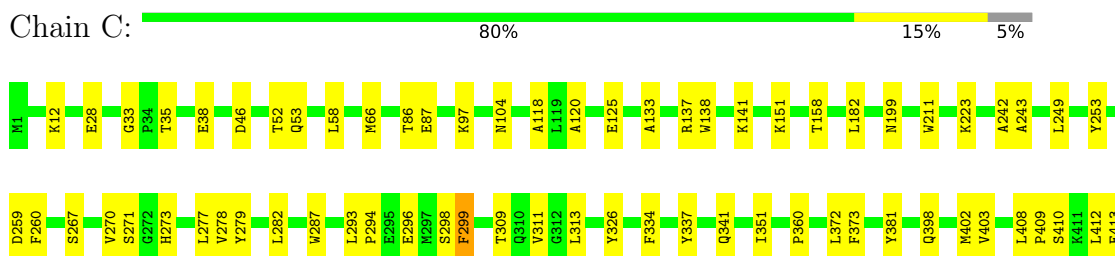
• Molecule 1: Glutamate decarboxylase

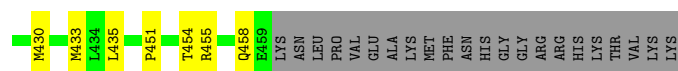


• Molecule 1: Glutamate decarboxylase



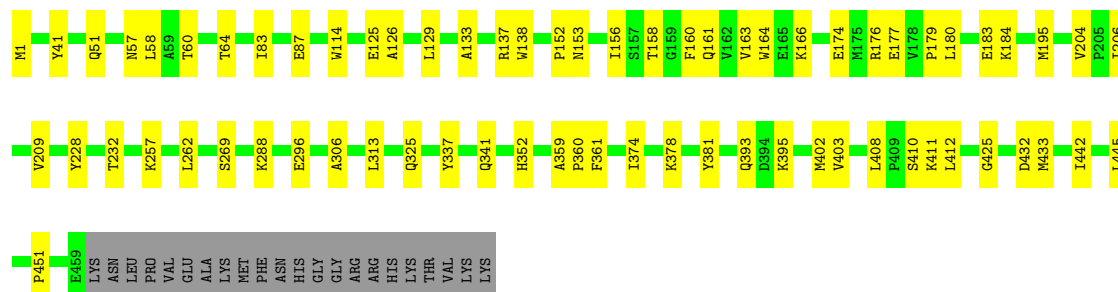
• Molecule 1: Glutamate decarboxylase





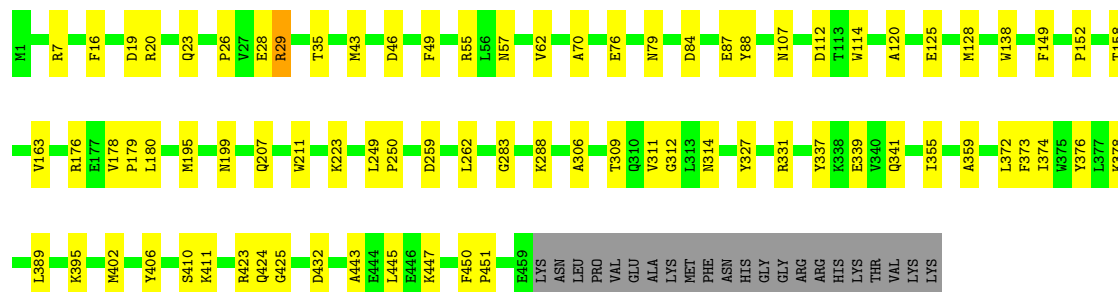
• Molecule 1: Glutamate decarboxylase

Chain D: 81% 14% 5%



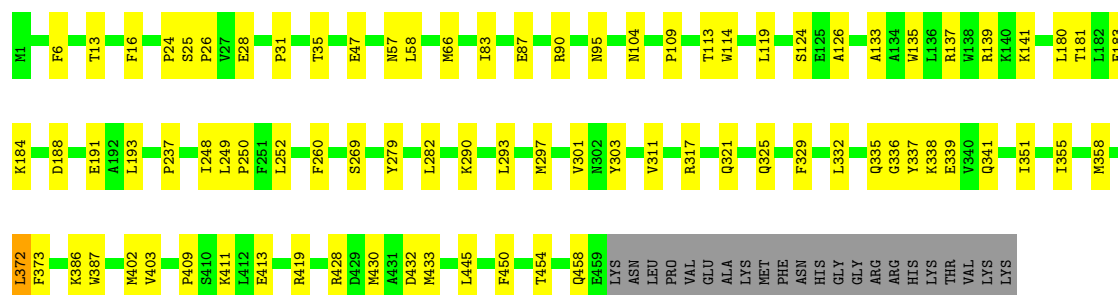
• Molecule 1: Glutamate decarboxylase

Chain E: 79% 16% 5%



• Molecule 1: Glutamate decarboxylase

Chain F: 78% 17% 5%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	109.23Å 124.90Å 109.62Å 90.00° 100.52° 90.00°	Depositor
Resolution (Å)	50.15 – 2.46 50.15 – 2.46	Depositor EDS
% Data completeness (in resolution range)	95.9 (50.15-2.46) 95.0 (50.15-2.46)	Depositor EDS
R_{merge}	0.20	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	673.53 (at 2.45Å)	Xtriage
Refinement program	PHENIX (1.19.1_4122: ???)	Depositor
R, R_{free}	0.181 , 0.234 0.199 , 0.252	Depositor DCC
R_{free} test set	5089 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	31.6	Xtriage
Anisotropy	0.609	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 27.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.40$, $\langle L^2 \rangle = 0.23$	Xtriage
Estimated twinning fraction	0.086 for l,-k,h	Xtriage
Reported twinning fraction	0.120 for l,-k,h	Depositor
Outliers	1 of 100781 reflections (0.001%)	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	22585	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.15% 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 [i](#)

5.1 Standard geometry [i](#)

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	0.42	0/3740	0.62	0/5094
1	B	0.42	0/3728	0.60	0/5081
1	C	0.41	0/3734	0.59	0/5086
1	D	0.42	0/3736	0.61	0/5090
1	E	0.44	1/3740 (0.0%)	0.62	0/5093
1	F	0.38	0/3744	0.57	0/5098
All	All	0.41	1/22422 (0.0%)	0.60	0/30542

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	26	PRO	C-O	-5.28	1.17	1.23

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3641	0	3554	54	0
1	B	3629	0	3528	58	0
1	C	3635	0	3541	55	0
1	D	3637	0	3543	51	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	E	3641	0	3561	58	0
1	F	3645	0	3565	55	0
2	A	10	0	0	0	0
2	B	10	0	0	0	0
2	C	5	0	0	0	0
2	D	5	0	0	0	0
2	E	5	0	0	0	0
2	F	5	0	0	0	0
3	A	121	0	0	5	0
3	B	144	0	0	8	0
3	C	124	0	0	6	0
3	D	103	0	0	3	0
3	E	118	0	0	8	0
3	F	107	0	0	4	0
All	All	22585	0	21292	287	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.

The worst 5 of 287 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:D:180:LEU:O	1:D:411:LYS:HB2	1.80	0.82
1:A:87:GLU:HG3	1:B:402:MET:HG2	1.64	0.80
1:C:138:TRP:HA	1:C:141:LYS:HE3	1.68	0.74
1:A:402:MET:HG2	1:B:87:GLU:HG3	1.72	0.70
1:F:430:MET:HA	1:F:433:MET:HE2	1.75	0.69

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	457/481 (95%)	440 (96%)	17 (4%)	0	100	100
1	B	457/481 (95%)	440 (96%)	17 (4%)	0	100	100
1	C	457/481 (95%)	441 (96%)	16 (4%)	0	100	100
1	D	457/481 (95%)	440 (96%)	17 (4%)	0	100	100
1	E	457/481 (95%)	440 (96%)	17 (4%)	0	100	100
1	F	457/481 (95%)	438 (96%)	19 (4%)	0	100	100
All	All	2742/2886 (95%)	2639 (96%)	103 (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	383/414 (92%)	383 (100%)	0	100	100
1	B	380/414 (92%)	379 (100%)	1 (0%)	91	95
1	C	381/414 (92%)	378 (99%)	3 (1%)	79	88
1	D	382/414 (92%)	382 (100%)	0	100	100
1	E	383/414 (92%)	381 (100%)	2 (0%)	86	92
1	F	384/414 (93%)	382 (100%)	2 (0%)	86	92
All	All	2293/2484 (92%)	2285 (100%)	8 (0%)	91	95

5 of 8 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	F	372	LEU
1	F	35	THR
1	E	29	ARG
1	C	372	LEU
1	E	372	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 24 such sidechains are listed below:

Mol	Chain	Res	Type
1	E	310	GLN
1	E	370	ASN
1	E	363	ASN
1	E	398	GLN
1	B	346	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

8 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	501	-	4,4,4	0.13	0	6,6,6	0.26	0
2	SO4	B	502	-	4,4,4	0.27	0	6,6,6	0.27	0
2	SO4	E	501	-	4,4,4	0.18	0	6,6,6	0.53	0
2	SO4	D	501	-	4,4,4	0.19	0	6,6,6	0.32	0
2	SO4	A	502	-	4,4,4	0.23	0	6,6,6	0.42	0
2	SO4	A	501	-	4,4,4	0.26	0	6,6,6	0.49	0
2	SO4	F	501	-	4,4,4	0.22	0	6,6,6	0.35	0
2	SO4	C	501	-	4,4,4	0.21	0	6,6,6	0.58	0

There are no bond length outliers.

There are no bond angle outliers.

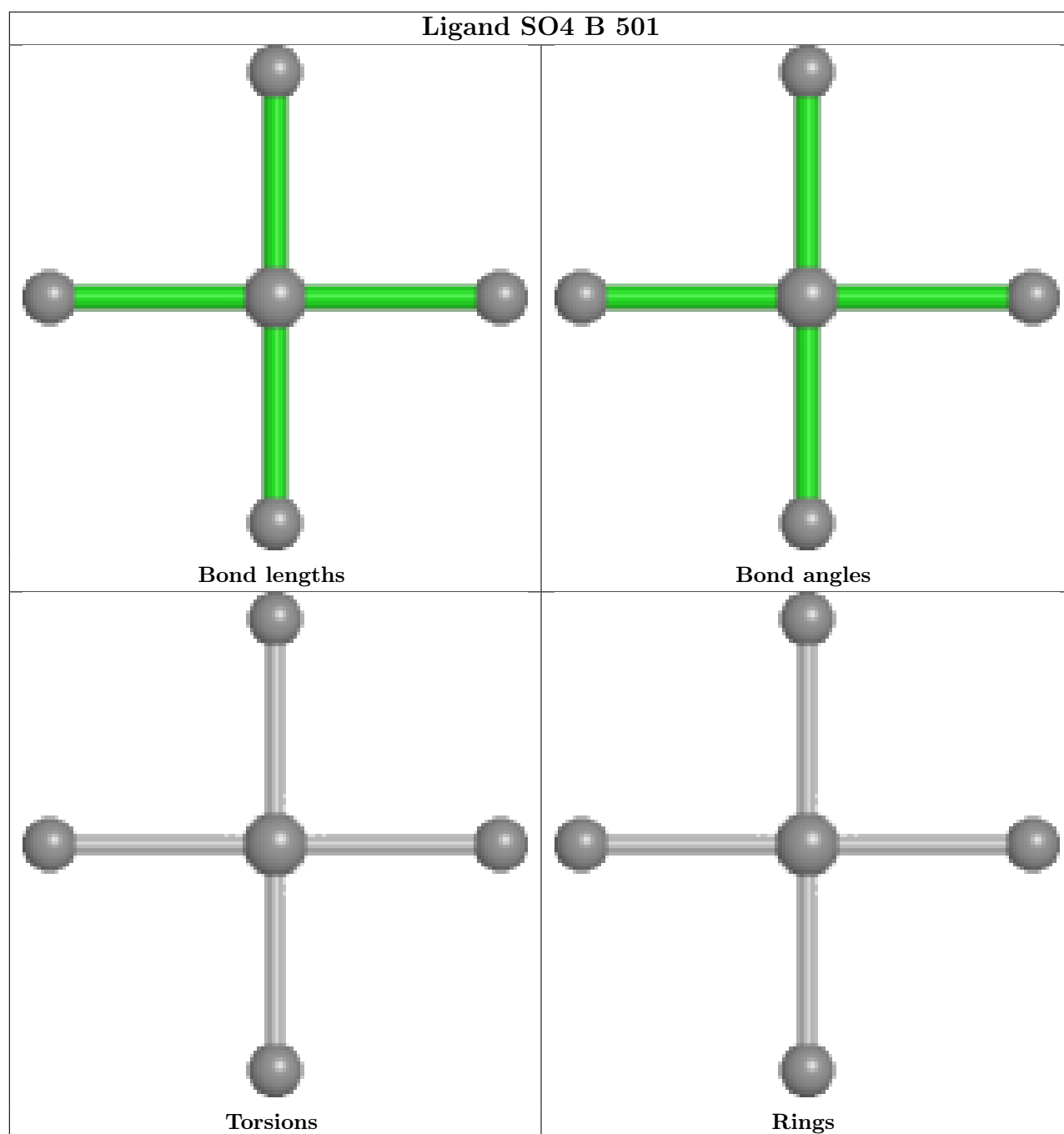
There are no chirality outliers.

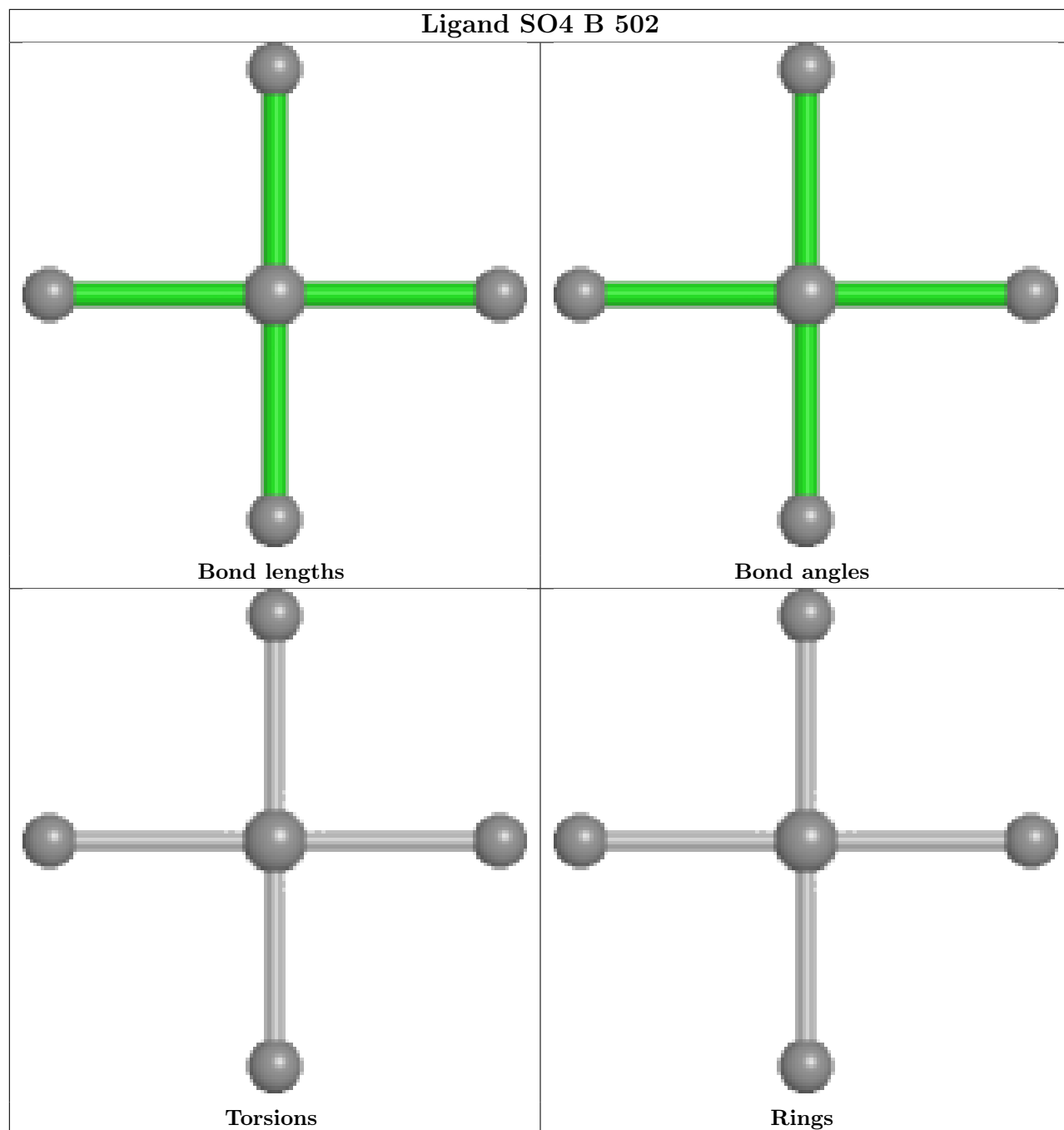
There are no torsion outliers.

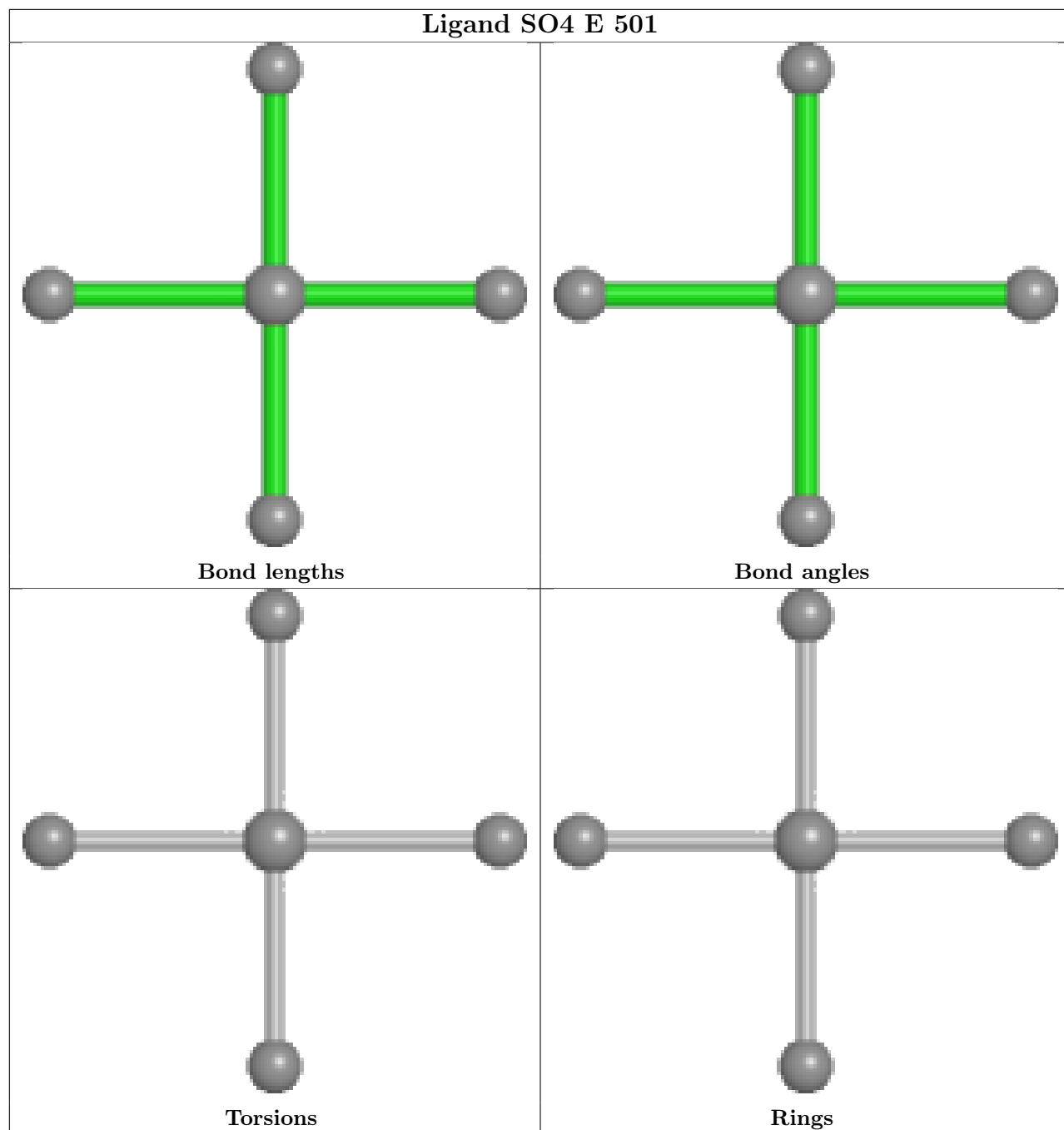
There are no ring outliers.

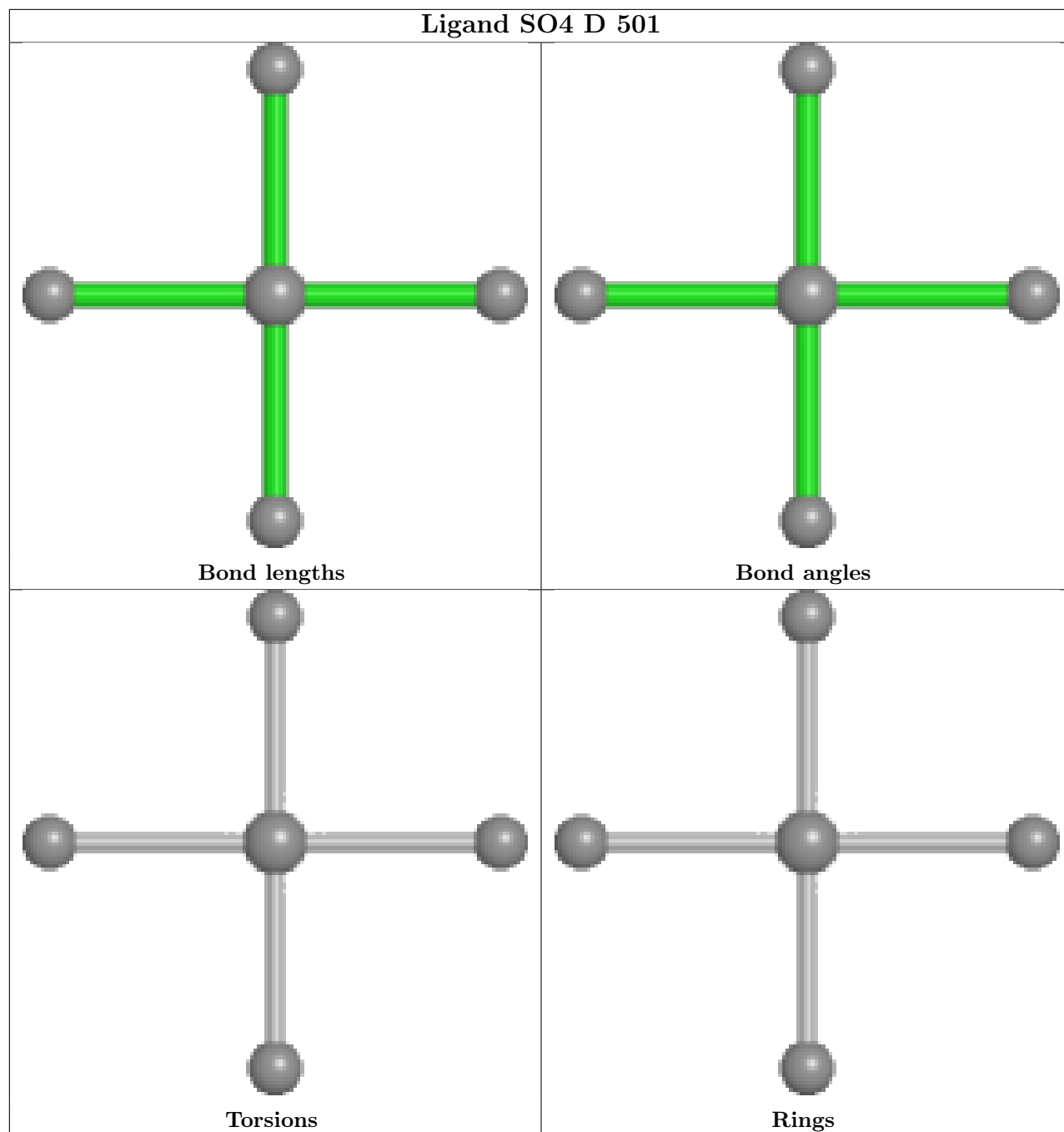
No monomer is involved in short contacts.

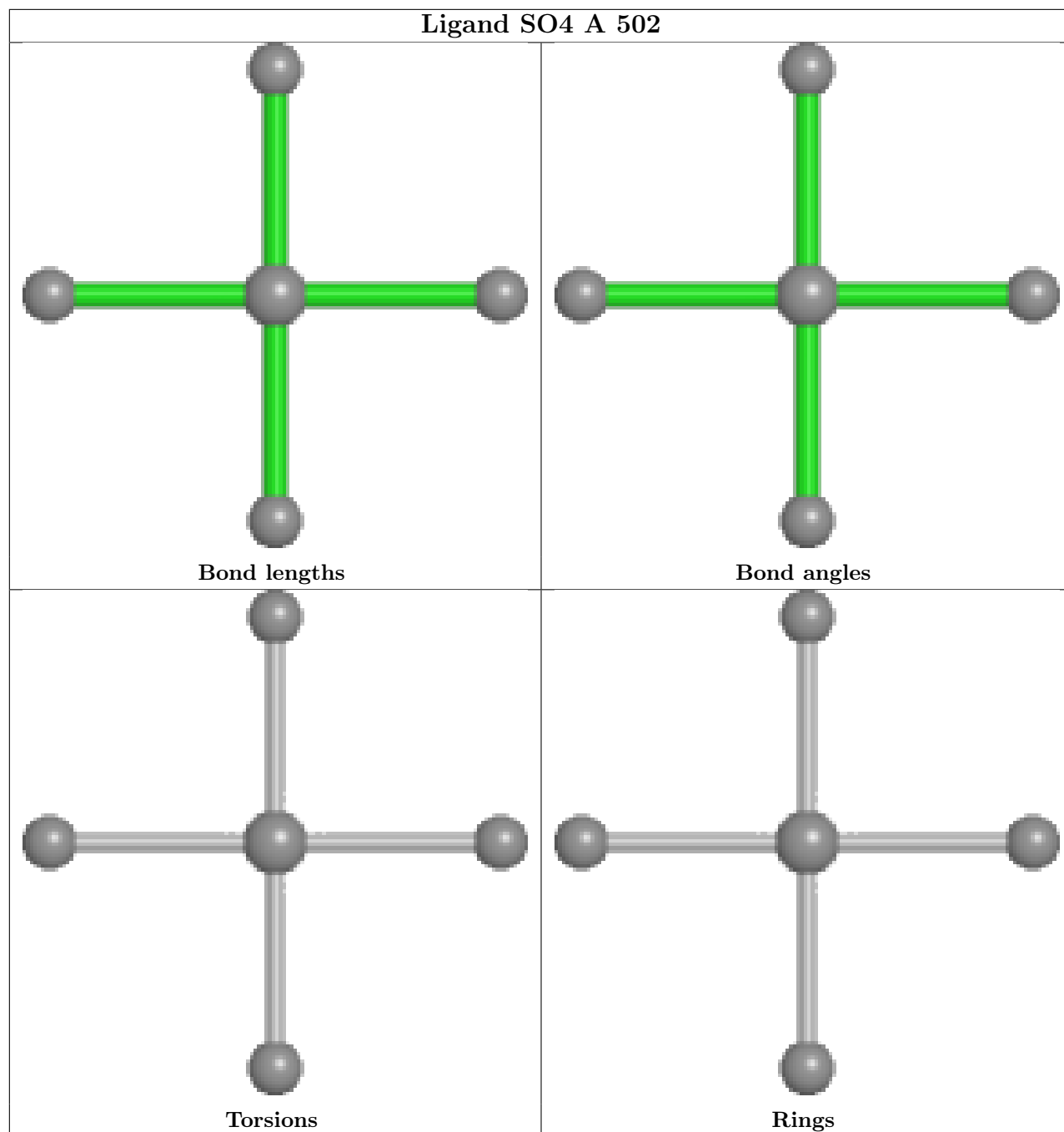
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

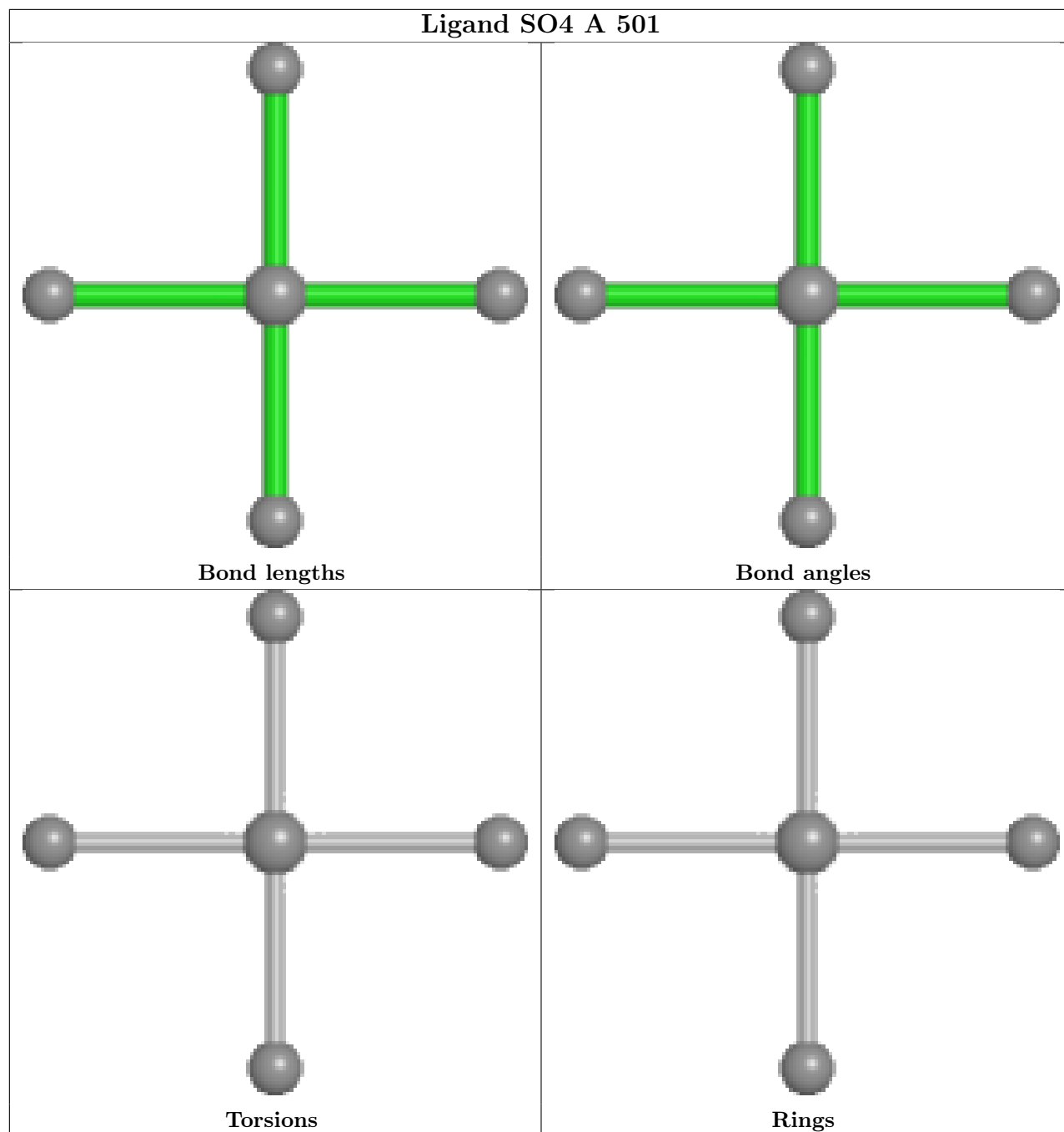


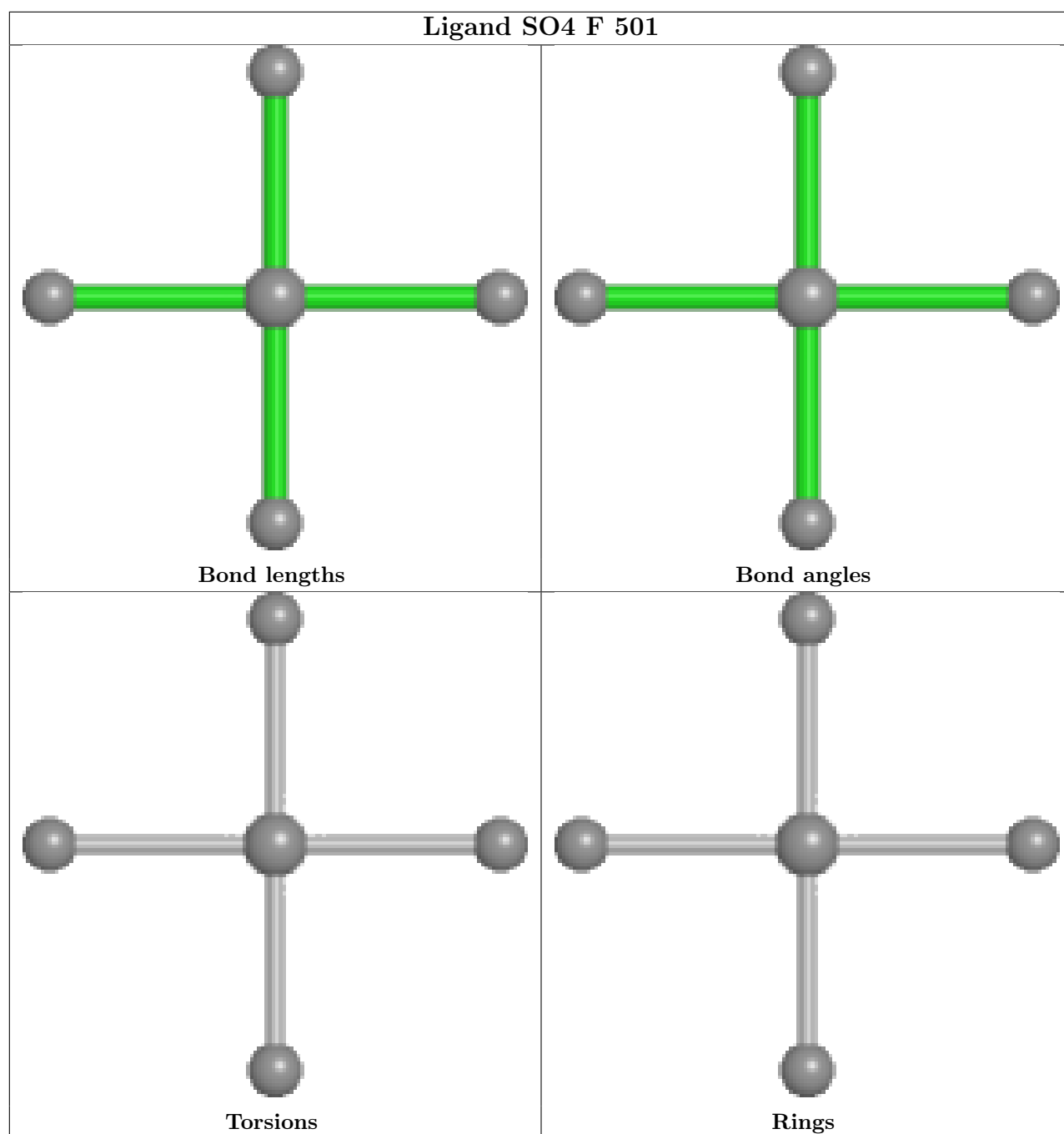


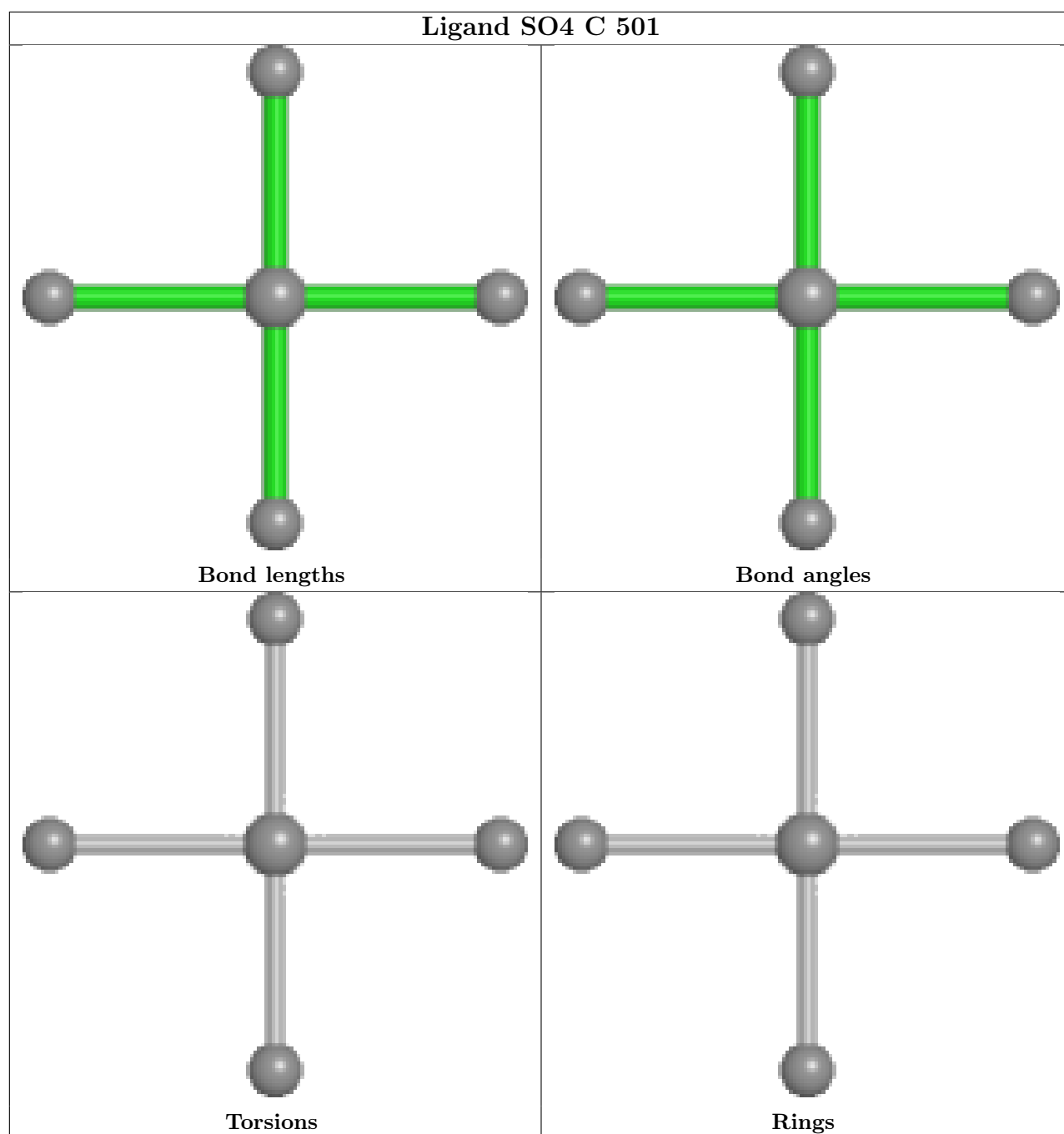












5.7 Other polymers [i](#)

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	459/481 (95%)	-1.55	0 100 100	15, 30, 44, 79	1 (0%)
1	B	459/481 (95%)	-1.57	0 100 100	15, 30, 44, 57	1 (0%)
1	C	459/481 (95%)	-1.55	0 100 100	17, 32, 46, 66	1 (0%)
1	D	459/481 (95%)	-1.56	0 100 100	15, 31, 45, 55	1 (0%)
1	E	459/481 (95%)	-1.52	0 100 100	17, 31, 46, 59	1 (0%)
1	F	459/481 (95%)	-1.50	0 100 100	16, 35, 51, 65	1 (0%)
All	All	2754/2886 (95%)	-1.54	0 100 100	15, 31, 46, 79	6 (0%)

There are no RSRZ outliers to report.

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(Å ²)	Q<0.9
2	SO4	B	502	5/5	0.99	0.04	37,37,39,40	0

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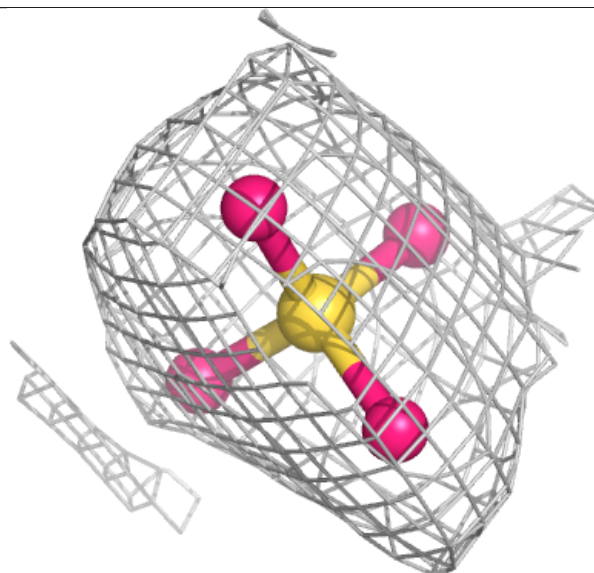
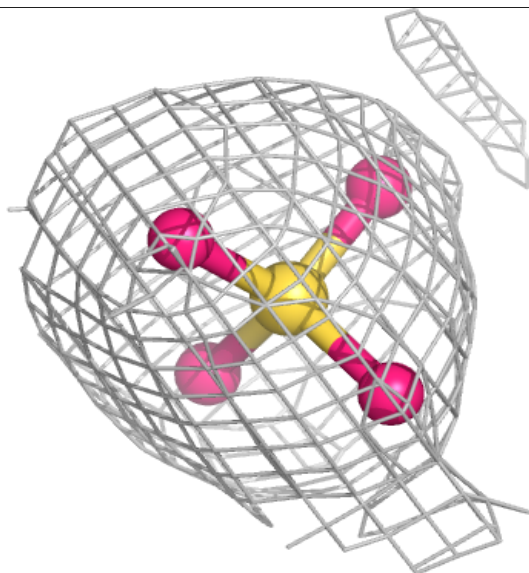
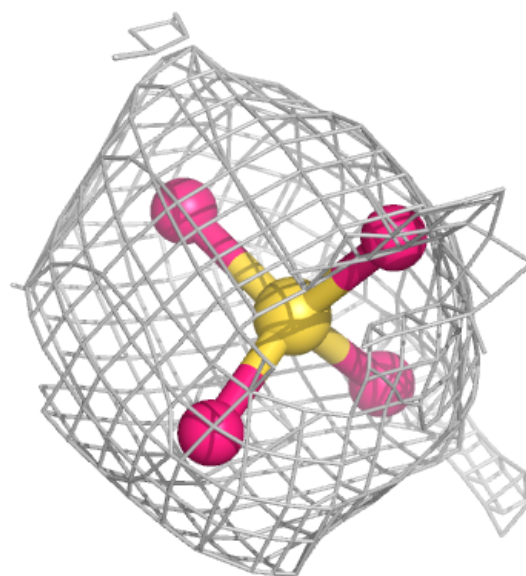
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	SO4	A	502	5/5	1.00	0.03	29,31,32,35	0
2	SO4	B	501	5/5	1.00	0.02	27,28,30,34	0
2	SO4	A	501	5/5	1.00	0.02	25,25,27,30	0
2	SO4	C	501	5/5	1.00	0.03	30,31,35,36	0
2	SO4	D	501	5/5	1.00	0.02	26,28,29,30	0
2	SO4	E	501	5/5	1.00	0.02	25,27,30,30	0
2	SO4	F	501	5/5	1.00	0.02	32,35,36,37	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.

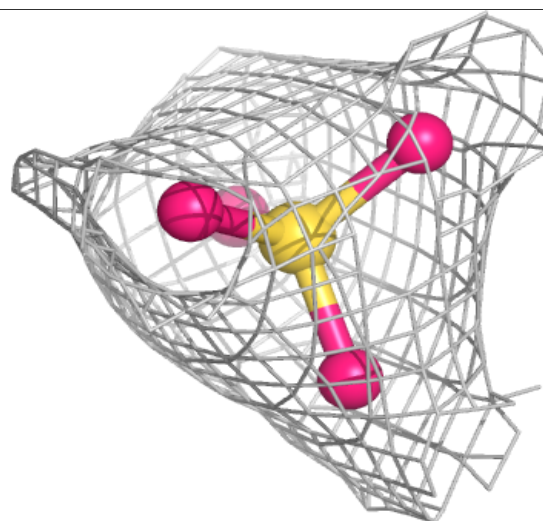
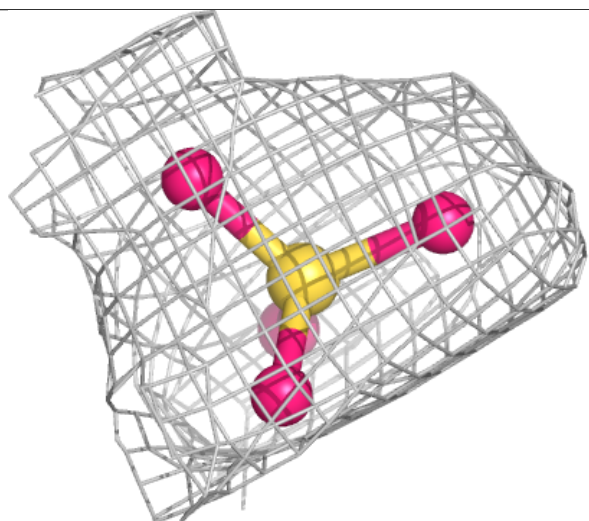
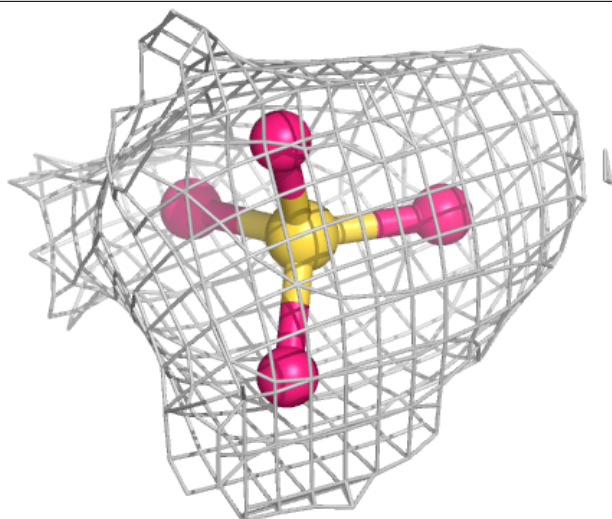
Electron density around SO4 B 502:

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



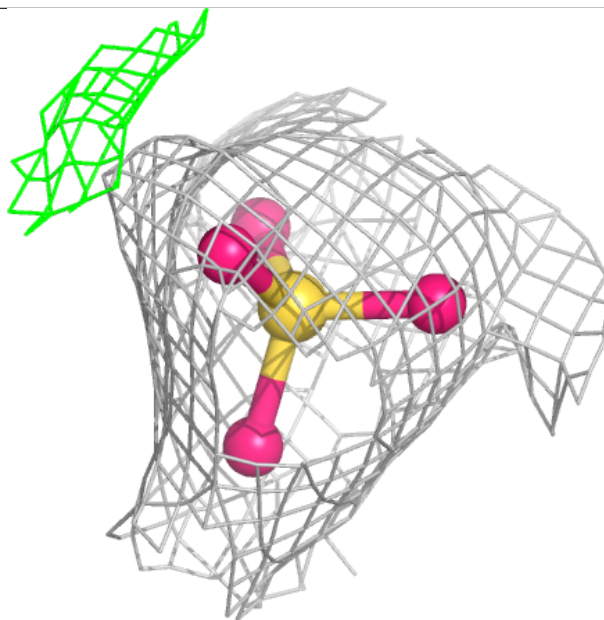
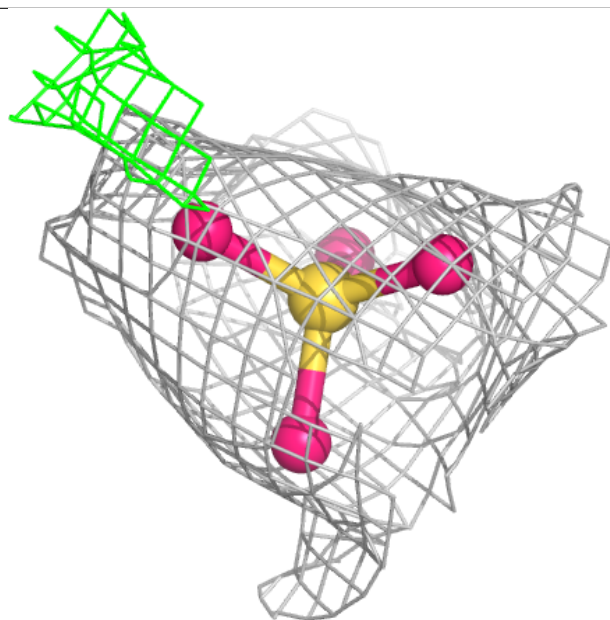
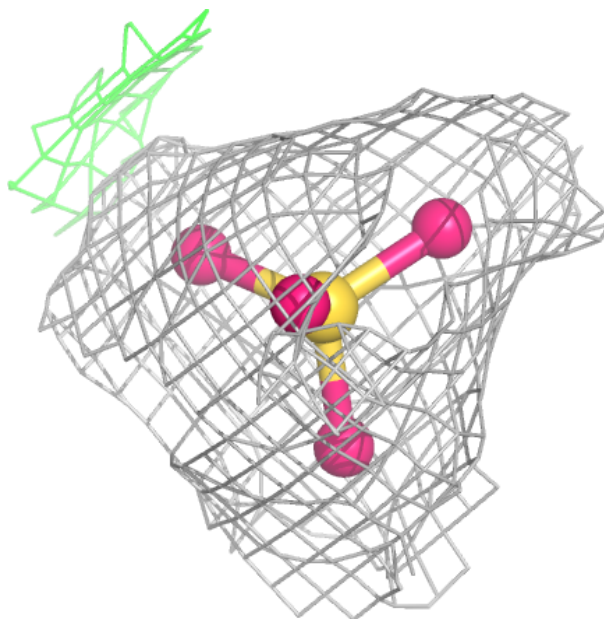
Electron density around SO4 A 502:

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



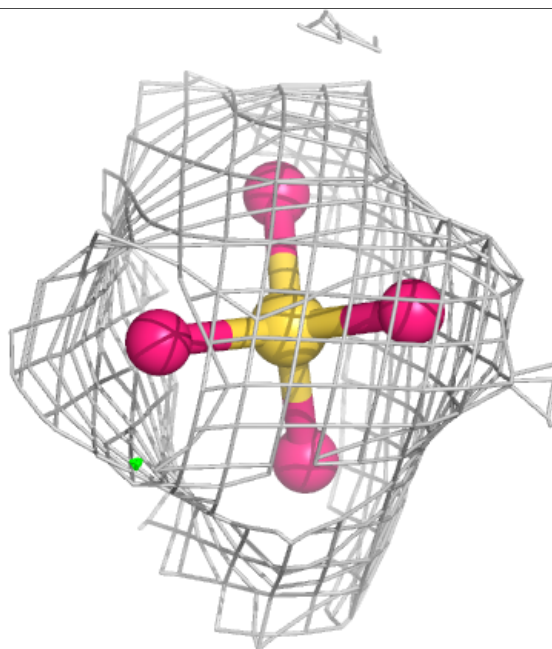
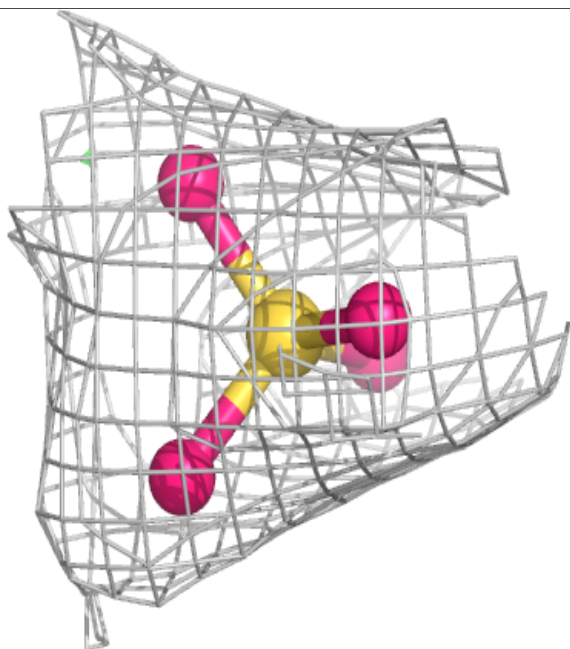
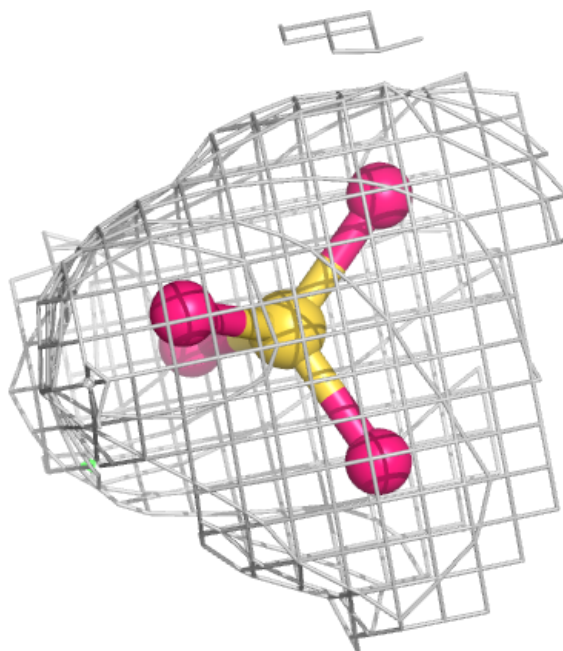
Electron density around SO4 B 501:

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



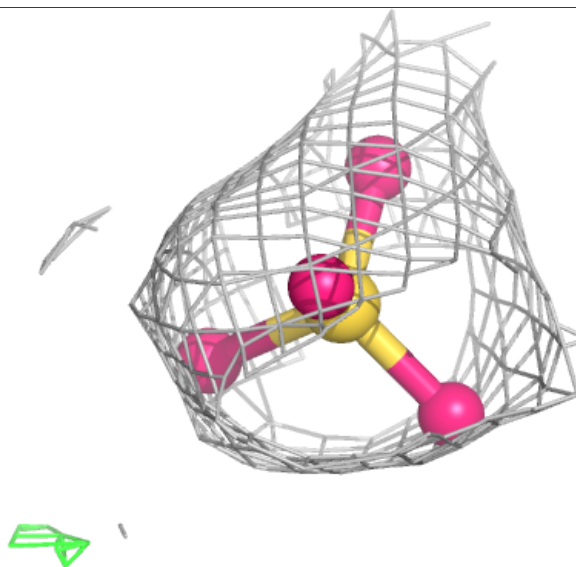
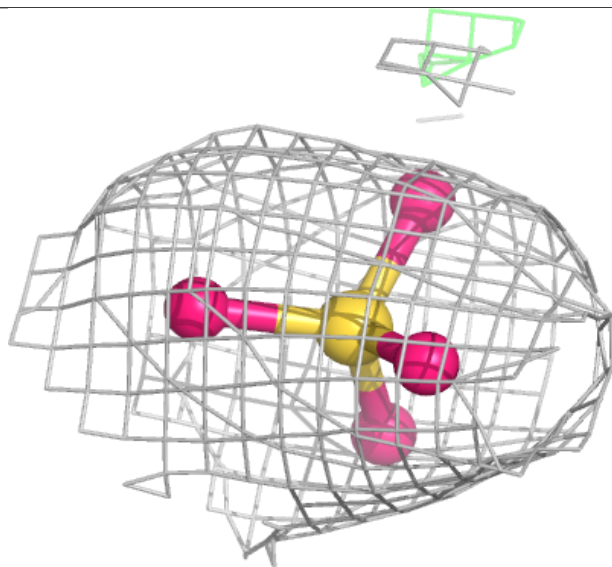
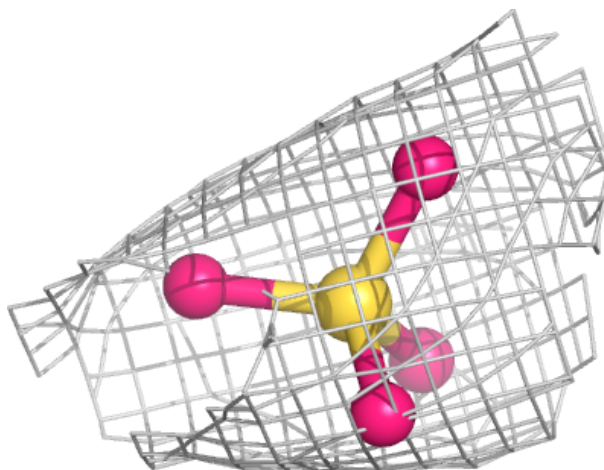
Electron density around SO4 A 501:

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



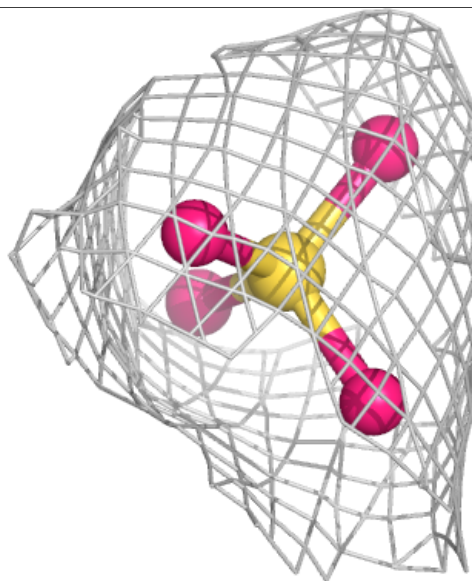
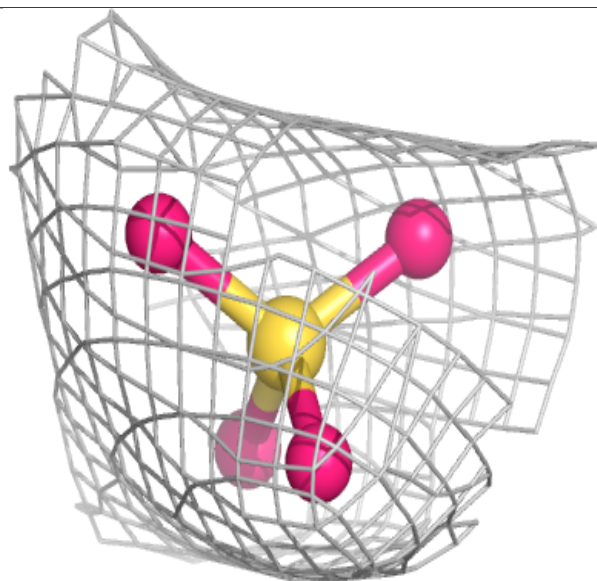
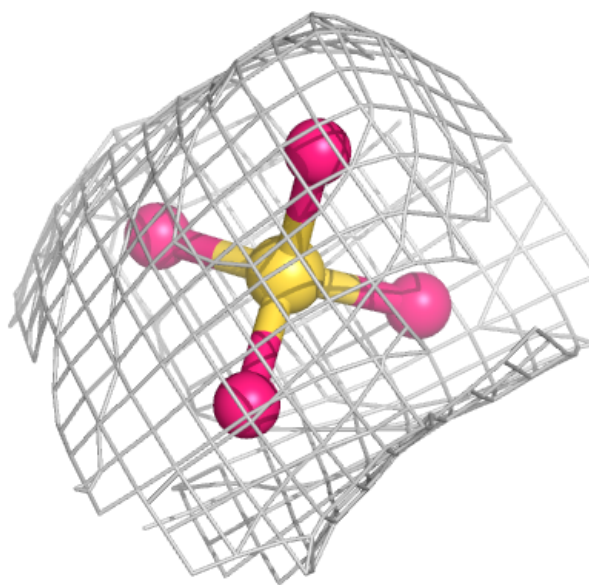
Electron density around SO4 C 501:

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



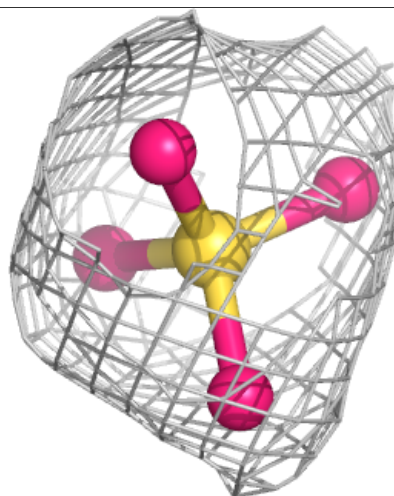
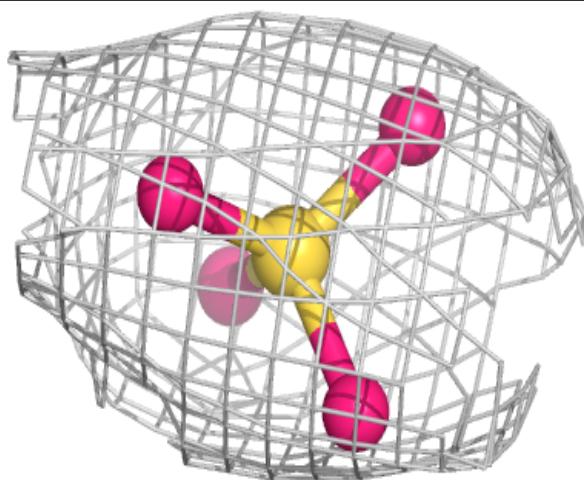
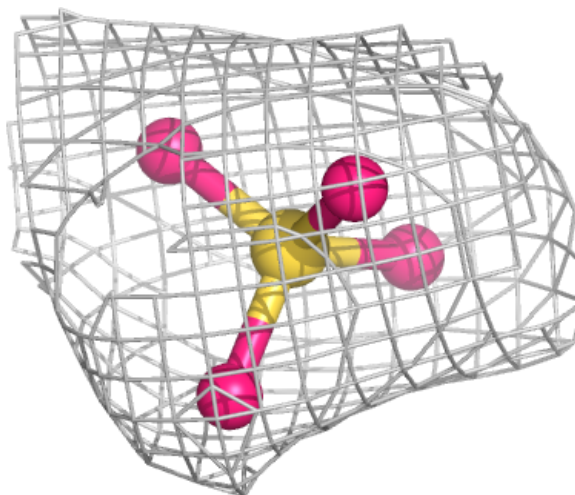
Electron density around SO4 D 501:

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



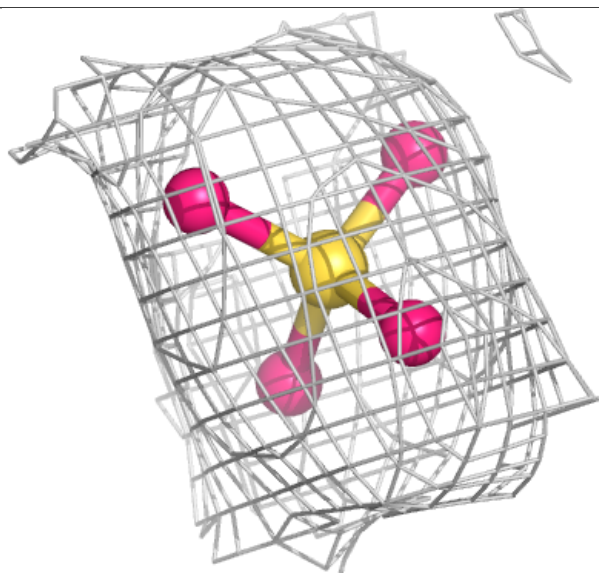
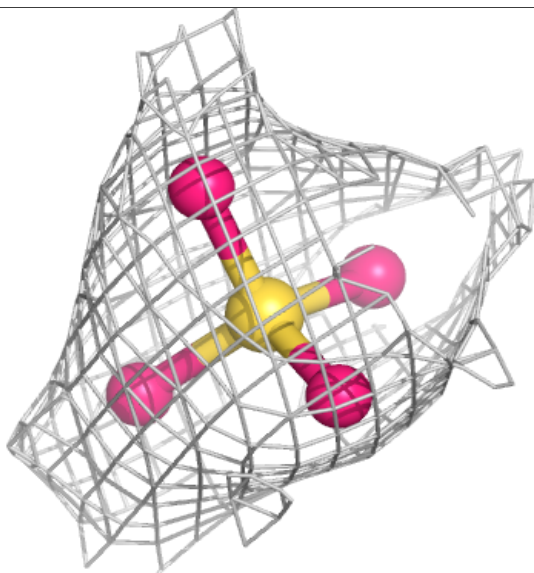
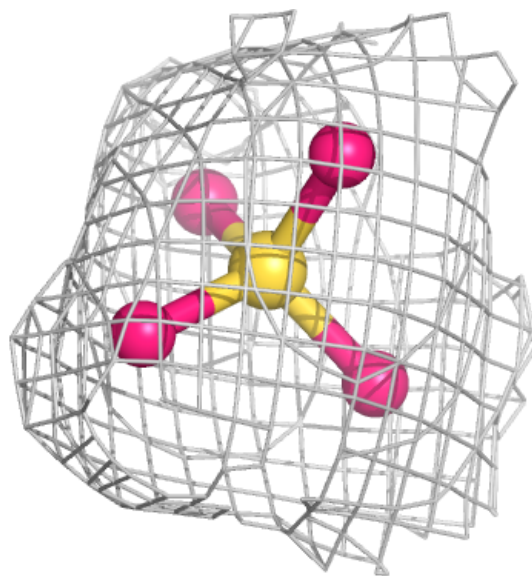
Electron density around SO4 E 501:

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



Electron density around SO4 F 501:

$2mF_o-DF_c$ (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.