



# Full wwPDB X-ray Structure Validation Report ⓘ

Apr 24, 2025 – 12:10 PM JST

PDB ID : 9JTX / pdb\_00009jtx  
Title : Factor inhibiting HIF-1 alpha in complex with Zn(II) and 2-(4-hydroxy-2-oxo-1-(thiazol-4-ylmethoxy)-1,2-dihydroquinoline-3-carboxamido)-2-methylpropionic acid  
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Deposited on : 2024-10-07  
Resolution : 2.08 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](mailto: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>.

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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)	:	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)	:	2.42

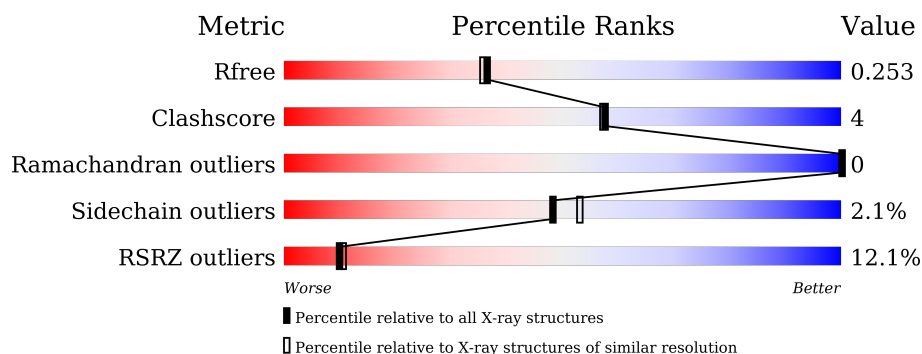
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.08 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	7574 (2.10-2.06)
Clashscore	180529	8325 (2.10-2.06)
Ramachandran outliers	177936	8271 (2.10-2.06)
Sidechain outliers	177891	8272 (2.10-2.06)
RSRZ outliers	164620	7574 (2.10-2.06)

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  $\geq 3$ , 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	349	<div> <div>12%</div> <div>85%</div> <div>12%</div> <div>..</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
3	A1L4V	A	404[A]	-	X	-	-

## 2 Entry composition [i](#)

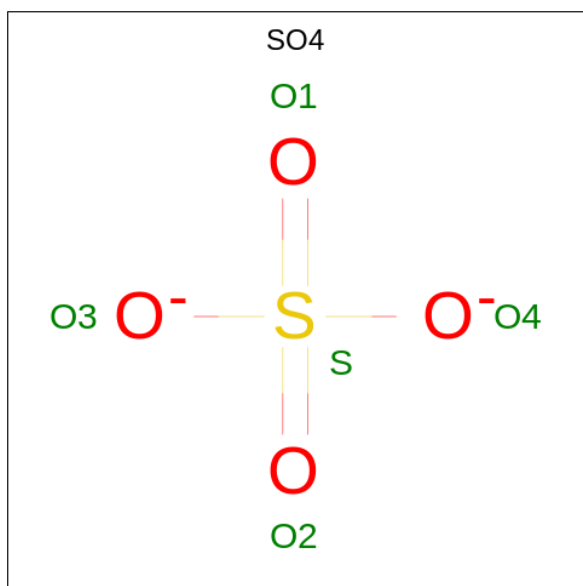
There are 5 unique types of molecules in this entry. The entry contains 2920 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Hypoxia-inducible factor 1-alpha inhibitor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	340	2770	1773	470	516	11	0	2	0

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O<sub>4</sub>S) (labeled as "Ligand of Interest" by depositor).



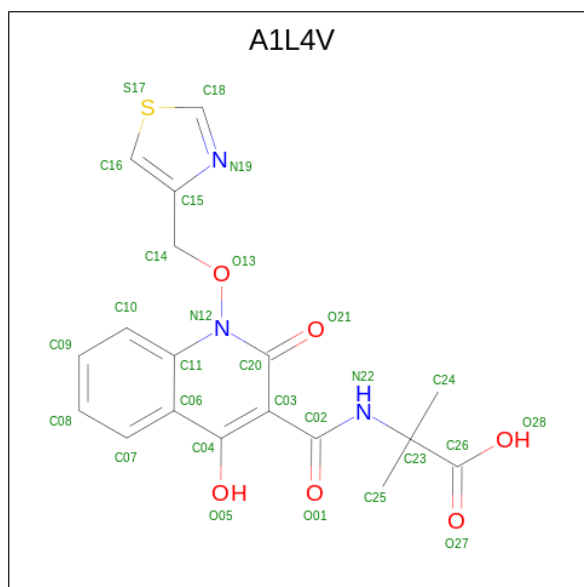
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	O	S		
2	A	1	5	4	1	0	0
2	A	1	5	4	1	0	0
2	A	1	5	4	1	0	0
2	A	1	5	4	1	0	0
2	A	1	5	4	1	0	0

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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 2-methyl-2-[[4-oxidanyl-2-oxidanylidene-1-(1,3-thiazol-4-ylmethoxy)quinolin-3-yl]carbonylamino]propanoic acid (CCD ID: A1L4V) (formula: C<sub>18</sub>H<sub>17</sub>N<sub>3</sub>O<sub>6</sub>S) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	S	
			56	36	6	12	2	1

- Molecule 4 is ZINC ION (CCD ID: ZN) (formula: Zn) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Zn	0	0
			1	1		

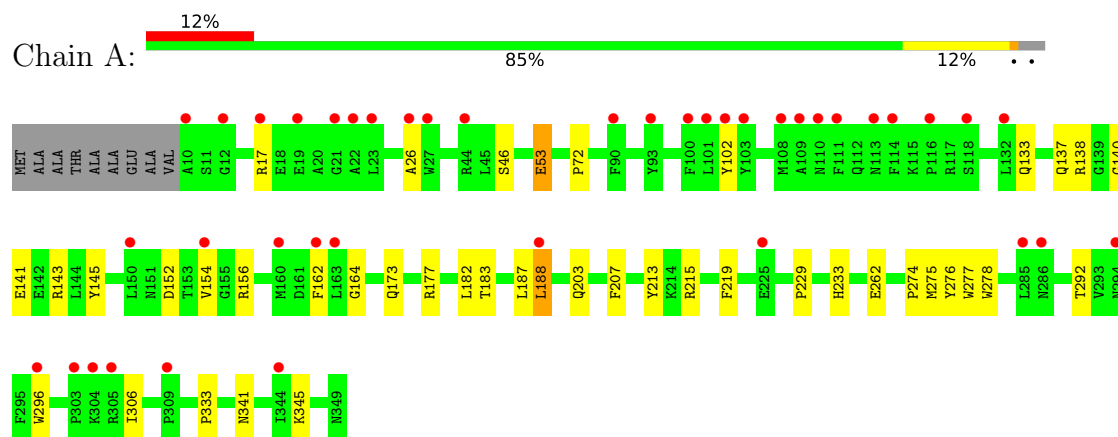
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	58	Total	O	0	0
			58	58		

### 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: Hypoxia-inducible factor 1-alpha inhibitor



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	86.58Å 86.58Å 145.18Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.80 – 2.08 46.80 – 2.08	Depositor EDS
% Data completeness (in resolution range)	93.7 (46.80-2.08) 94.0 (46.80-2.08)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	0.92 (at 2.08Å)	Xtriage
Refinement program	PHENIX 1.17.1 _3660	Depositor
R, $R_{free}$	0.214 , 0.247 0.220 , 0.253	Depositor DCC
$R_{free}$ test set	1798 reflections (5.30%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	64.2	Xtriage
Anisotropy	0.125	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 74.4	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	2920	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	92.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.93% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: ZN, SO4, A1L4V

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	0.25	0/2851	0.44	1/3874 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
1	A	188	LEU	CA-CB-CG	-5.16	103.42	115.30

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	2770	0	2630	24	0
2	A	35	0	0	3	0
3	A	56	0	0	0	0
4	A	1	0	0	0	0
5	A	58	0	0	1	0
All	All	2920	0	2630	24	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 4.



All (24) 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:53:GLU:HG3	1:A:177:ARG:HH22	1.50	0.75
1:A:17:ARG:NH2	5:A:503:HOH:O	2.36	0.59
1:A:341:ASN:HB3	1:A:345:LYS:HE2	1.87	0.57
1:A:187:LEU:C	1:A:188:LEU:HG	2.26	0.56
1:A:138:ARG:HH11	1:A:140:GLY:HA3	1.71	0.55
1:A:133:GLN:OE1	1:A:137:GLN:NE2	2.41	0.54
1:A:274:PRO:HB2	1:A:277:TRP:CD1	2.43	0.54
1:A:219:PHE:HB2	1:A:278:TRP:HB2	1.89	0.53
1:A:183:THR:HG21	1:A:203:GLN:HE22	1.74	0.53
1:A:275:MET:HG2	1:A:276:TYR:CD2	2.45	0.51
1:A:143:ARG:NH1	2:A:401:SO4:O4	2.39	0.50
1:A:143:ARG:NH2	2:A:409:SO4:O1	2.45	0.50
1:A:141[A]:GLU:N	2:A:406:SO4:O1	2.46	0.46
1:A:26:ALA:HB1	1:A:213:TYR:CZ	2.50	0.46
1:A:72:PRO:HG2	1:A:164:GLY:HA3	1.97	0.46
1:A:215:ARG:NE	1:A:262:GLU:OE2	2.46	0.45
1:A:207:PHE:CZ	1:A:292:THR:HG21	2.53	0.43
1:A:173:GLN:O	1:A:177:ARG:HG2	2.19	0.43
1:A:233:HIS:CG	1:A:333:PRO:HB3	2.53	0.42
1:A:187:LEU:O	1:A:188:LEU:HG	2.19	0.42
1:A:182:LEU:HD12	1:A:296:TRP:O	2.20	0.41
1:A:306:ILE:HD13	1:A:306:ILE:HA	1.94	0.40
1:A:229:PRO:HB3	1:A:278:TRP:HH2	1.86	0.40
1:A:152:ASP:C	1:A:154:VAL:H	2.25	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	340/349 (97%)	330 (97%)	10 (3%)	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	294/304 (97%)	288 (98%)	6 (2%)	50	55

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	46	SER
1	A	53	GLU
1	A	102	TYR
1	A	145	TYR
1	A	156	ARG
1	A	162	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	133	GLN
1	A	137	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 [i](#)

There are no oligosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

Of 10 ligands modelled in this entry, 1 is monoatomic - leaving 9 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	SO4	A	403	-	4,4,4	0.14	0	6,6,6	0.05	0
3	A1L4V	A	404[A]	4	26,30,30	4.34	11 (42%)	35,44,44	4.75	20 (57%)
3	A1L4V	A	404[B]	4	26,30,30	4.38	10 (38%)	35,44,44	4.21	16 (45%)
2	SO4	A	409	-	4,4,4	0.14	0	6,6,6	0.04	0
2	SO4	A	406	-	4,4,4	0.14	0	6,6,6	0.05	0
2	SO4	A	407	-	4,4,4	0.14	0	6,6,6	0.05	0
2	SO4	A	408	-	4,4,4	0.14	0	6,6,6	0.05	0
2	SO4	A	401	-	4,4,4	0.14	0	6,6,6	0.05	0
2	SO4	A	402	-	4,4,4	0.14	0	6,6,6	0.04	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	A1L4V	A	404[A]	4	-	14/18/20/20	0/3/3/3
3	A1L4V	A	404[B]	4	-	12/18/20/20	0/3/3/3

All (21) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	404[B]	A1L4V	C11-N12	14.90	1.55	1.39
3	A	404[A]	A1L4V	C11-N12	14.12	1.54	1.39
3	A	404[A]	A1L4V	C02-N22	11.09	1.49	1.34

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	404[B]	A1L4V	C02-N22	11.05	1.49	1.34
3	A	404[A]	A1L4V	C20-N12	6.94	1.47	1.38
3	A	404[B]	A1L4V	C20-N12	6.26	1.46	1.38
3	A	404[A]	A1L4V	O13-N12	5.66	1.44	1.38
3	A	404[B]	A1L4V	O13-N12	4.95	1.43	1.38
3	A	404[A]	A1L4V	C03-C02	4.84	1.57	1.48
3	A	404[B]	A1L4V	C03-C02	4.73	1.57	1.48
3	A	404[B]	A1L4V	C06-C04	4.22	1.52	1.45
3	A	404[A]	A1L4V	C03-C04	4.16	1.47	1.39
3	A	404[B]	A1L4V	C03-C04	3.96	1.46	1.39
3	A	404[A]	A1L4V	C06-C04	3.18	1.50	1.45
3	A	404[B]	A1L4V	O13-C14	-2.71	1.40	1.46
3	A	404[A]	A1L4V	O21-C20	-2.52	1.17	1.23
3	A	404[B]	A1L4V	O01-C02	-2.45	1.18	1.23
3	A	404[A]	A1L4V	C14-C15	2.40	1.56	1.50
3	A	404[B]	A1L4V	O21-C20	-2.24	1.18	1.23
3	A	404[A]	A1L4V	C23-C26	-2.10	1.51	1.53
3	A	404[A]	A1L4V	O13-C14	-2.02	1.42	1.46

All (36) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	404[A]	A1L4V	C04-C03-C02	13.31	128.57	118.29
3	A	404[B]	A1L4V	C04-C03-C02	11.84	127.43	118.29
3	A	404[A]	A1L4V	C14-O13-N12	11.72	121.70	109.77
3	A	404[B]	A1L4V	C23-N22-C02	11.33	137.82	122.55
3	A	404[A]	A1L4V	C23-N22-C02	11.01	137.39	122.55
3	A	404[B]	A1L4V	O01-C02-N22	-7.74	110.72	124.29
3	A	404[B]	A1L4V	C14-C15-C16	-7.12	119.03	129.99
3	A	404[A]	A1L4V	O01-C02-N22	-7.11	111.83	124.29
3	A	404[A]	A1L4V	C14-C15-C16	-6.06	120.67	129.99
3	A	404[A]	A1L4V	C25-C23-N22	5.96	121.37	109.11
3	A	404[B]	A1L4V	C25-C23-N22	5.69	120.82	109.11
3	A	404[B]	A1L4V	O01-C02-C03	-5.56	113.01	121.53
3	A	404[A]	A1L4V	C25-C23-C26	-5.33	96.33	108.85
3	A	404[B]	A1L4V	C03-C02-N22	5.26	136.20	117.06
3	A	404[A]	A1L4V	C07-C06-C04	-4.84	114.17	122.50
3	A	404[A]	A1L4V	O21-C20-N12	-4.84	114.92	120.64
3	A	404[B]	A1L4V	O21-C20-N12	-4.82	114.95	120.64
3	A	404[A]	A1L4V	C20-C03-C02	-4.46	116.37	121.51
3	A	404[B]	A1L4V	C25-C23-C26	-4.45	98.40	108.85
3	A	404[B]	A1L4V	C20-C03-C02	-4.38	116.46	121.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	404[A]	A1L4V	C04-C03-C20	-4.00	115.04	120.39
3	A	404[A]	A1L4V	C03-C02-N22	3.86	131.10	117.06
3	A	404[B]	A1L4V	O13-C14-C15	3.81	117.84	107.95
3	A	404[A]	A1L4V	C11-C06-C04	3.61	123.02	119.19
3	A	404[A]	A1L4V	C07-C06-C11	3.46	122.81	118.82
3	A	404[B]	A1L4V	C07-C06-C04	-3.40	116.66	122.50
3	A	404[A]	A1L4V	C25-C23-C24	3.34	115.33	110.41
3	A	404[A]	A1L4V	O05-C04-C03	3.32	127.68	122.71
3	A	404[B]	A1L4V	C04-C03-C20	-3.27	116.02	120.39
3	A	404[A]	A1L4V	O05-C04-C06	-3.01	110.28	115.53
3	A	404[A]	A1L4V	O13-C14-C15	2.93	115.55	107.95
3	A	404[A]	A1L4V	O01-C02-C03	-2.92	117.06	121.53
3	A	404[B]	A1L4V	C07-C06-C11	2.83	122.09	118.82
3	A	404[B]	A1L4V	O13-N12-C20	-2.82	113.33	116.23
3	A	404[B]	A1L4V	C16-S17-C18	2.64	97.76	92.37
3	A	404[A]	A1L4V	C16-S17-C18	2.17	96.80	92.37

There are no chirality outliers.

All (26) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	404[A]	A1L4V	C11-N12-O13-C14
3	A	404[A]	A1L4V	N22-C23-C26-O27
3	A	404[A]	A1L4V	N22-C23-C26-O28
3	A	404[A]	A1L4V	N22-C02-C03-C20
3	A	404[A]	A1L4V	N22-C02-C03-C04
3	A	404[A]	A1L4V	O01-C02-C03-C20
3	A	404[A]	A1L4V	O01-C02-C03-C04
3	A	404[A]	A1L4V	C03-C02-N22-C23
3	A	404[A]	A1L4V	O01-C02-N22-C23
3	A	404[A]	A1L4V	C15-C14-O13-N12
3	A	404[B]	A1L4V	C11-N12-O13-C14
3	A	404[B]	A1L4V	N22-C23-C26-O27
3	A	404[B]	A1L4V	N22-C23-C26-O28
3	A	404[B]	A1L4V	N22-C02-C03-C20
3	A	404[B]	A1L4V	N22-C02-C03-C04
3	A	404[B]	A1L4V	O01-C02-C03-C20
3	A	404[B]	A1L4V	O01-C02-C03-C04
3	A	404[B]	A1L4V	C03-C02-N22-C23
3	A	404[B]	A1L4V	O01-C02-N22-C23
3	A	404[A]	A1L4V	C20-N12-O13-C14
3	A	404[A]	A1L4V	C26-C23-N22-C02

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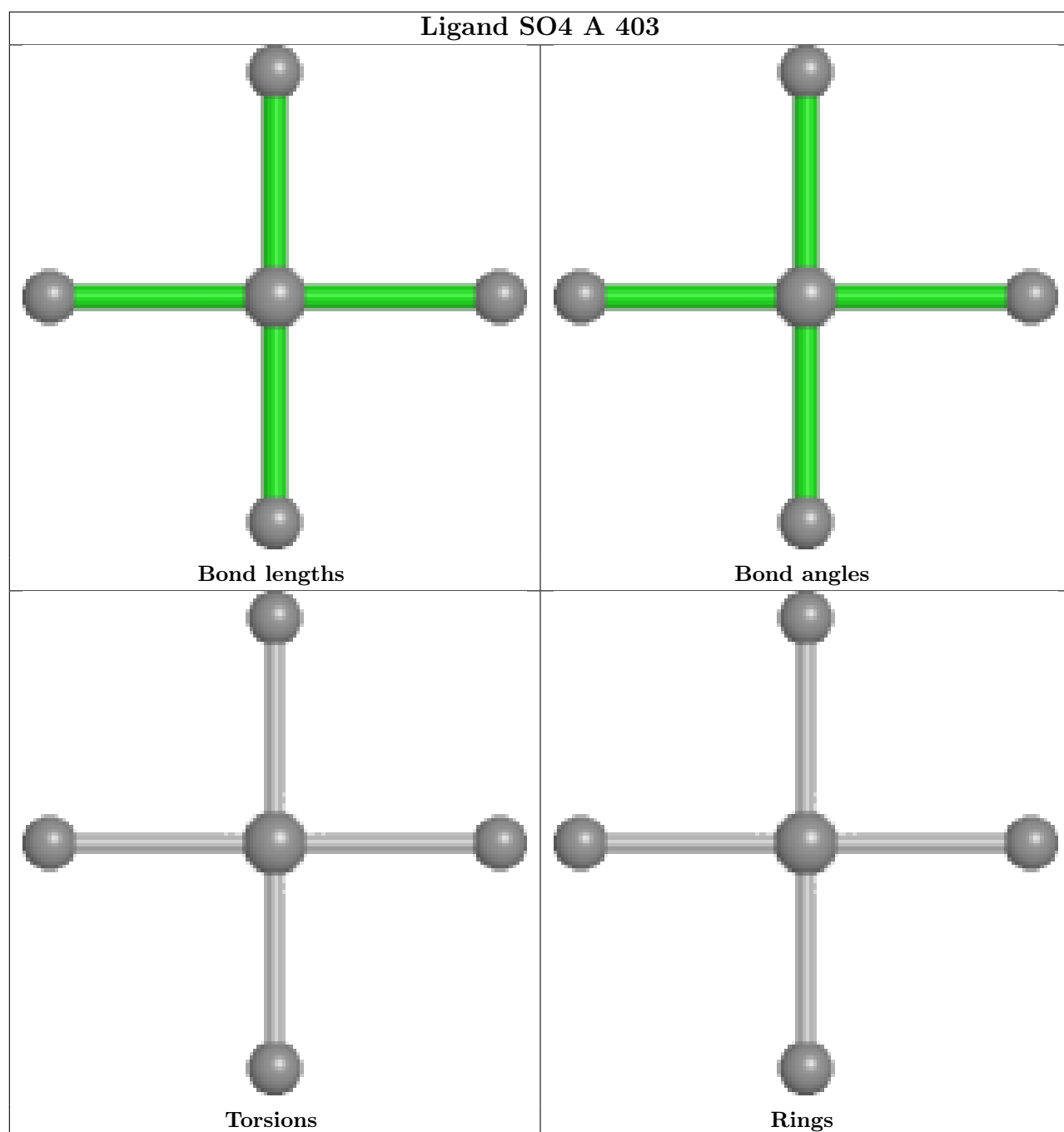
Mol	Chain	Res	Type	Atoms
3	A	404[B]	A1L4V	C26-C23-N22-C02
3	A	404[B]	A1L4V	C20-N12-O13-C14
3	A	404[A]	A1L4V	C24-C23-C26-O27
3	A	404[A]	A1L4V	C24-C23-C26-O28
3	A	404[B]	A1L4V	C24-C23-C26-O27

There are no ring outliers.

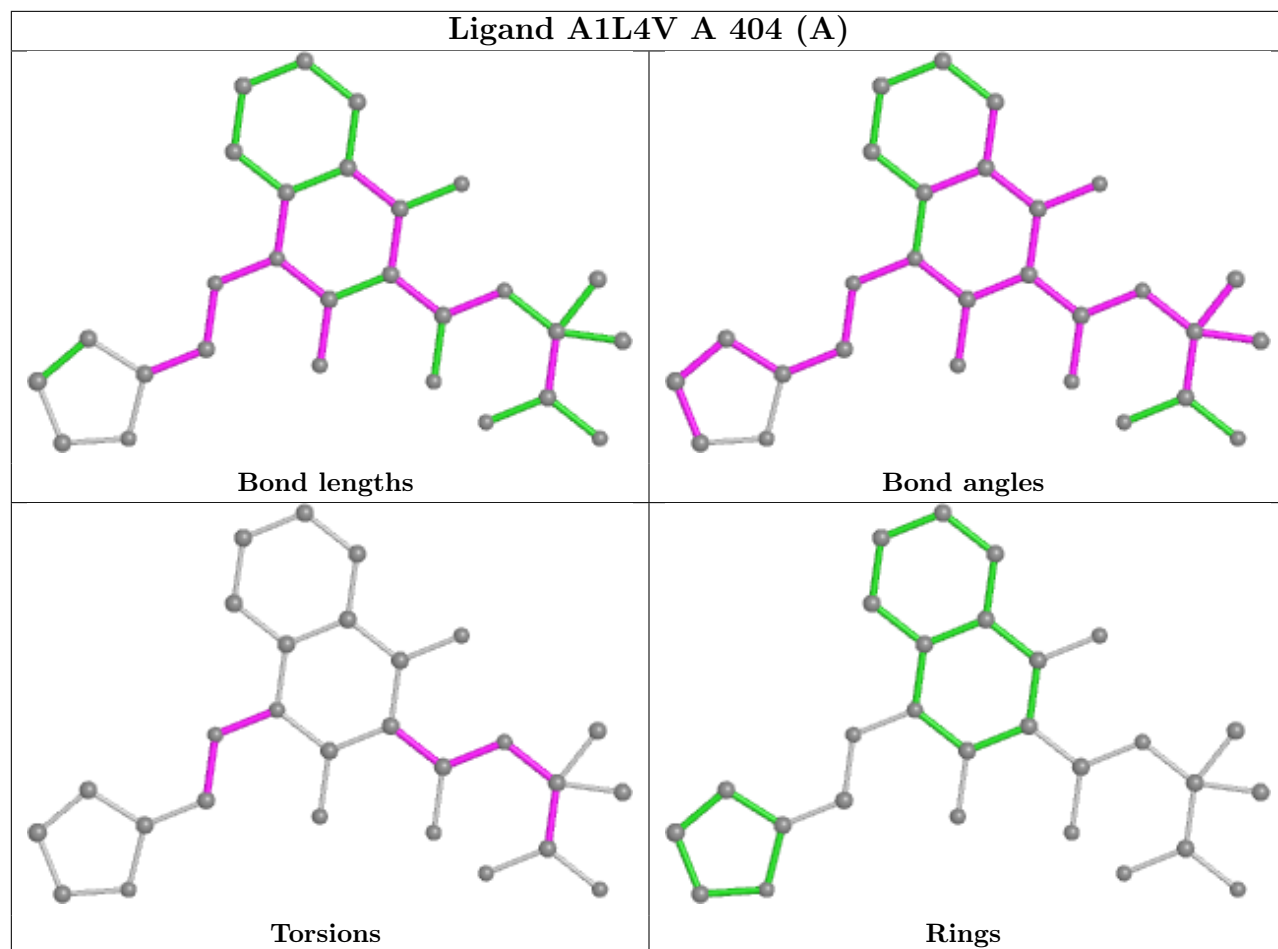
3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	409	SO4	1	0
2	A	406	SO4	1	0
2	A	401	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.

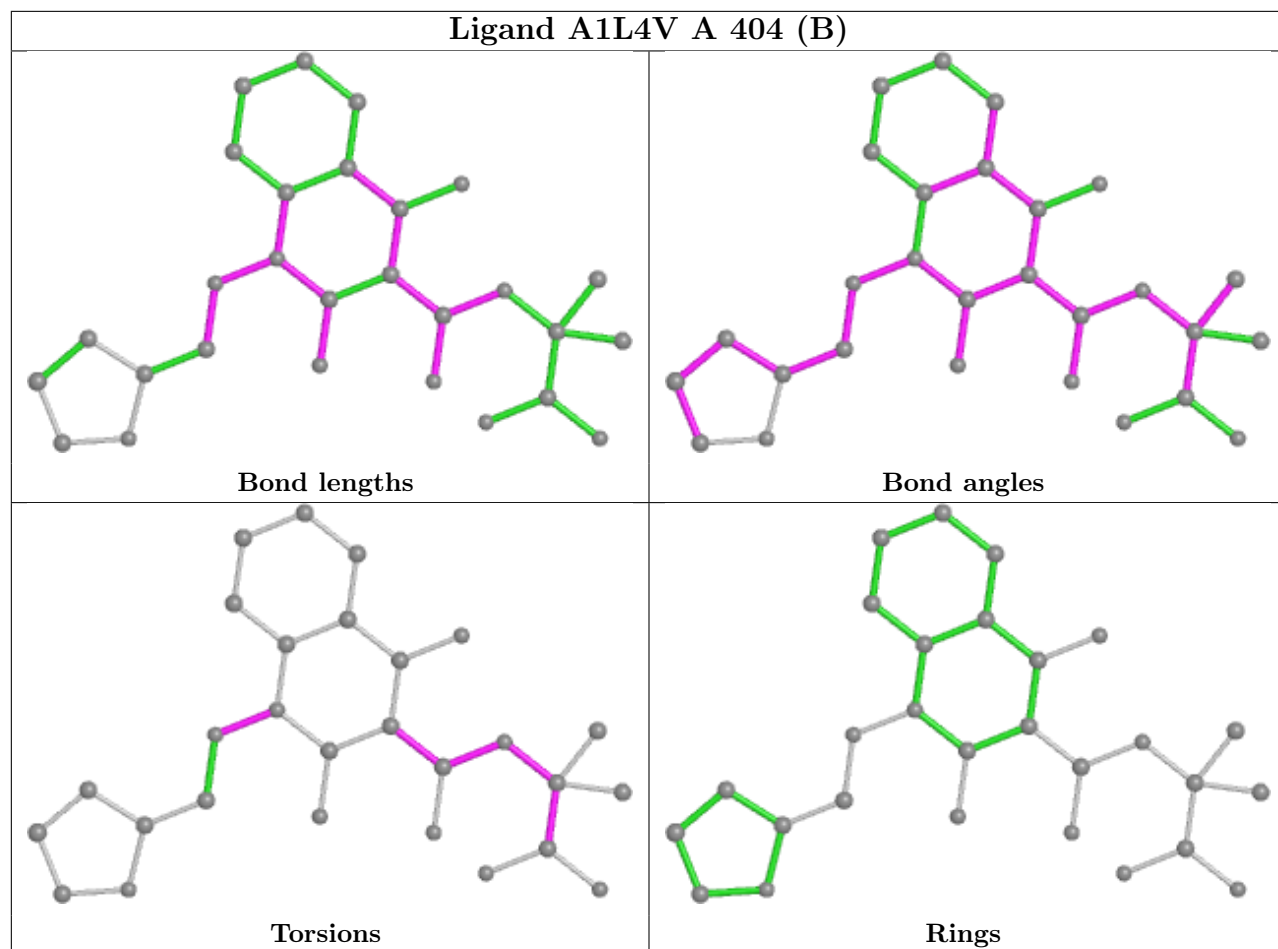


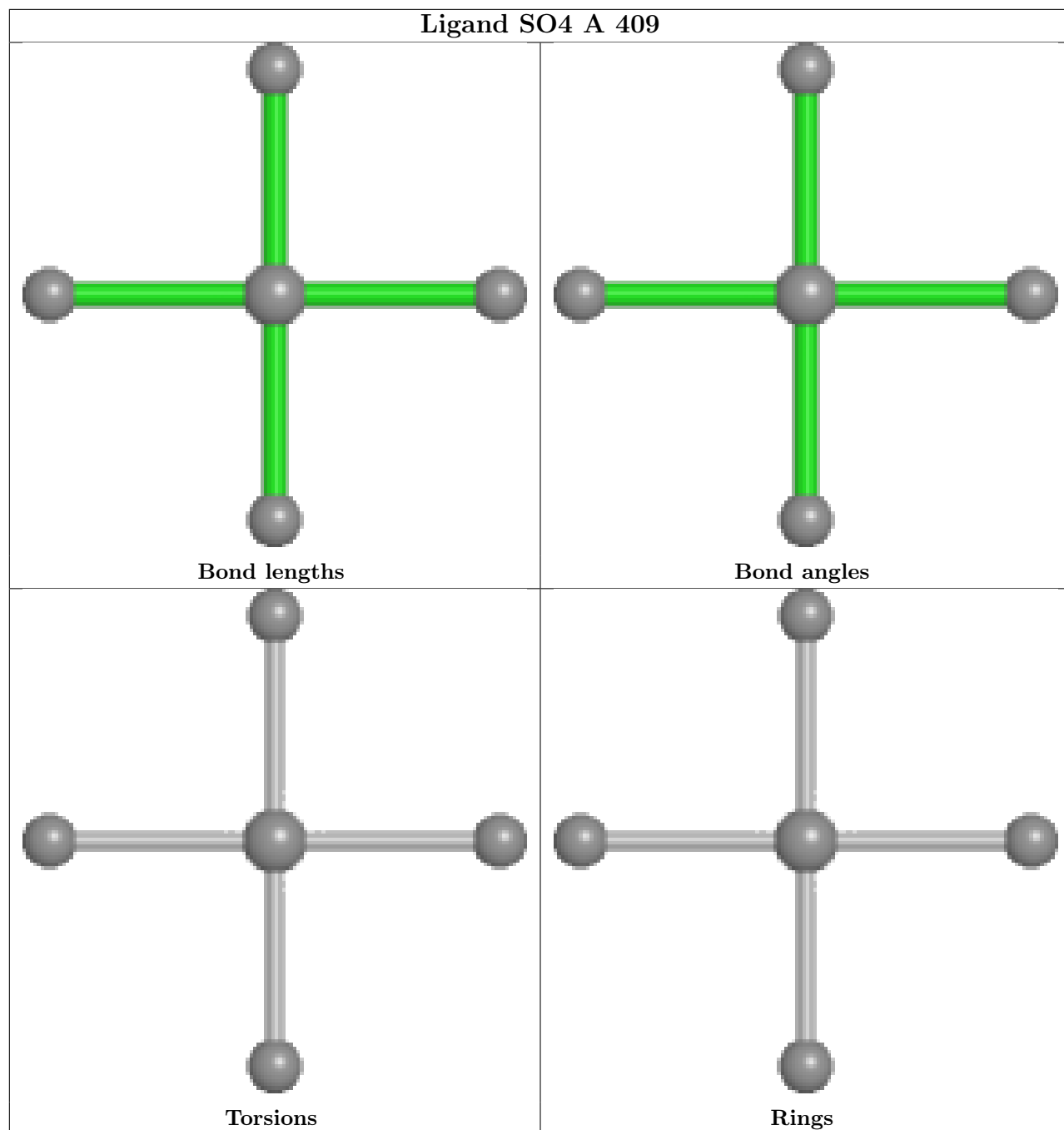
## Ligand A1L4V A 404 (A)

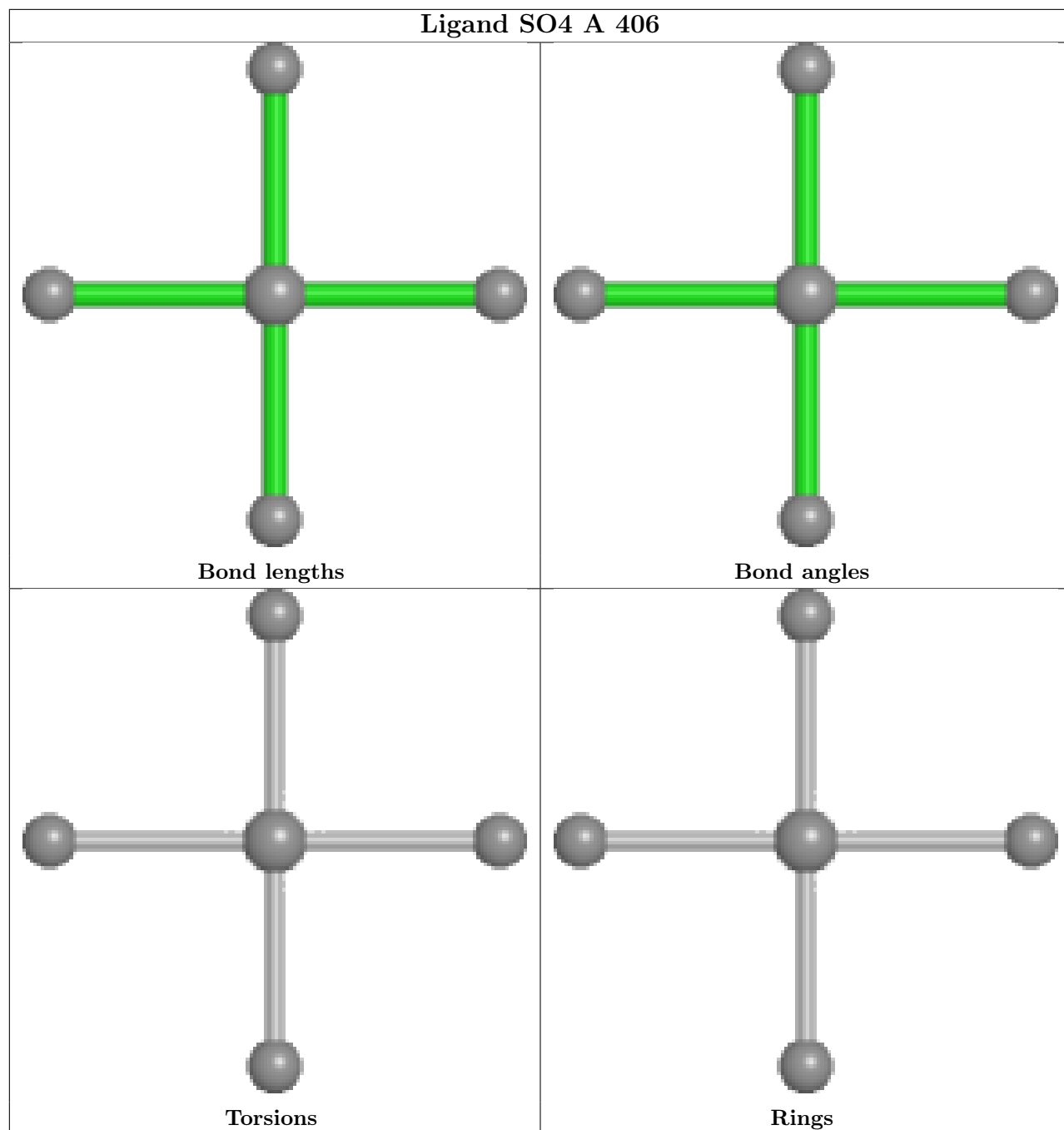


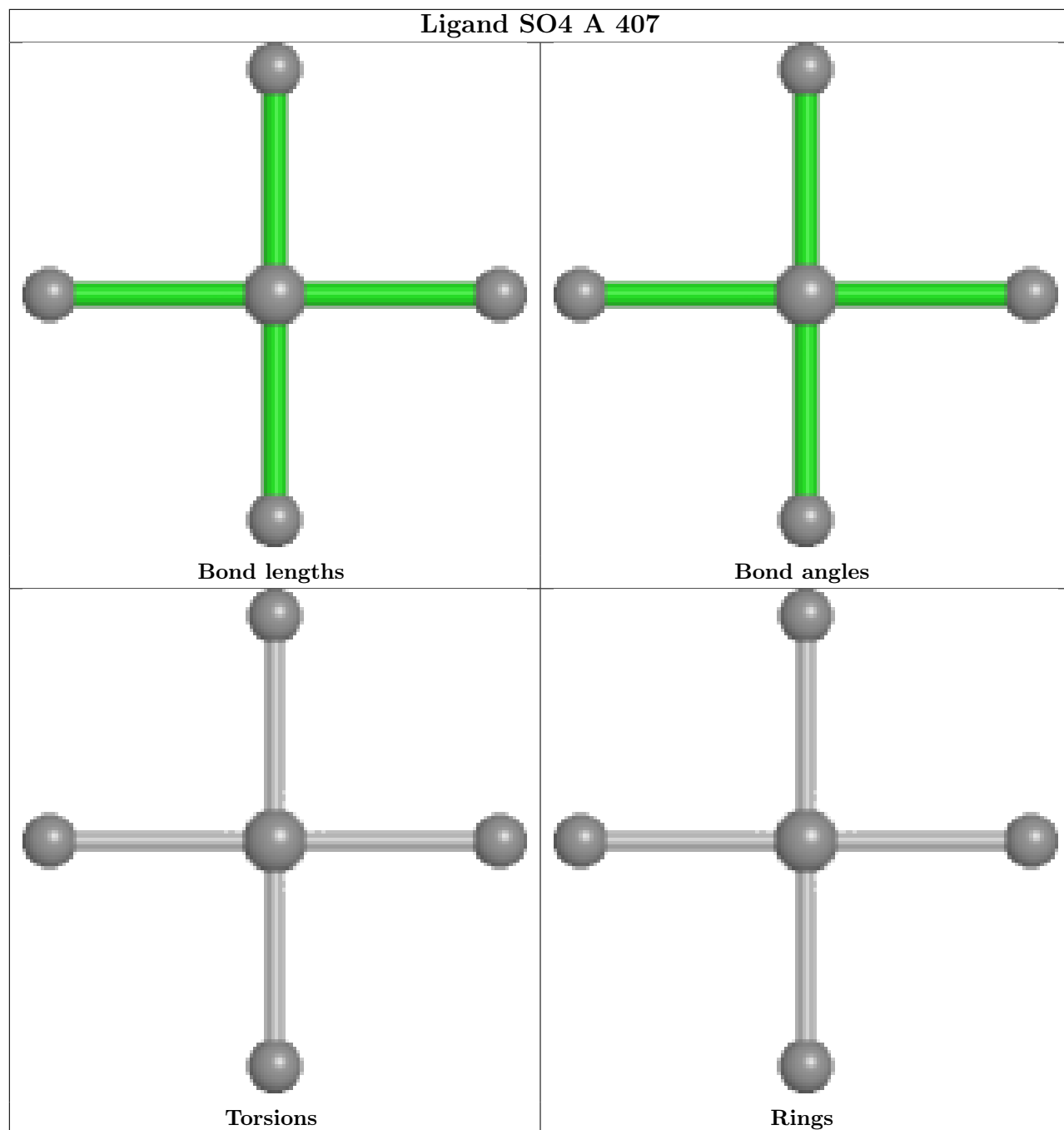


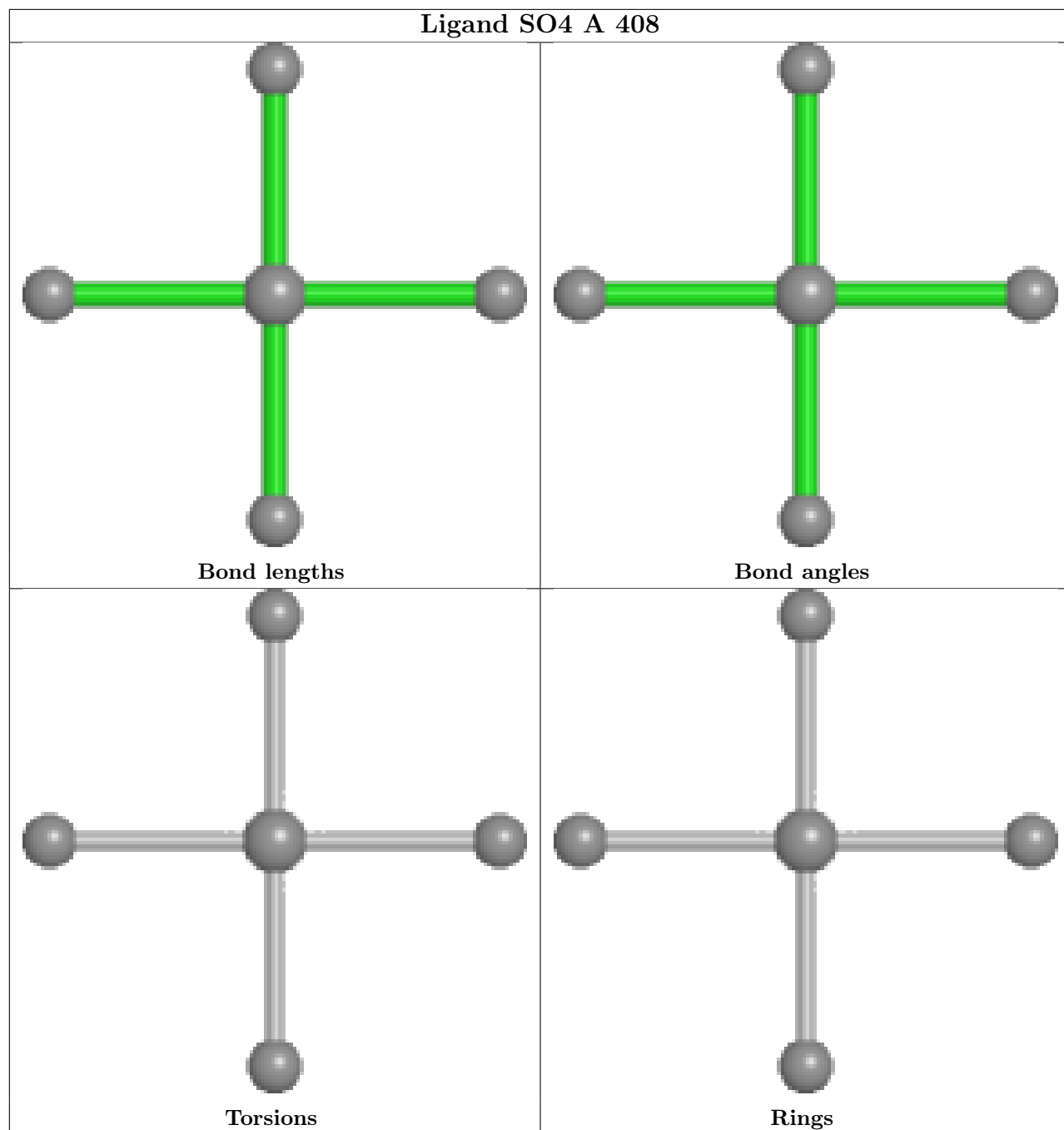
## Ligand A1L4V A 404 (B)

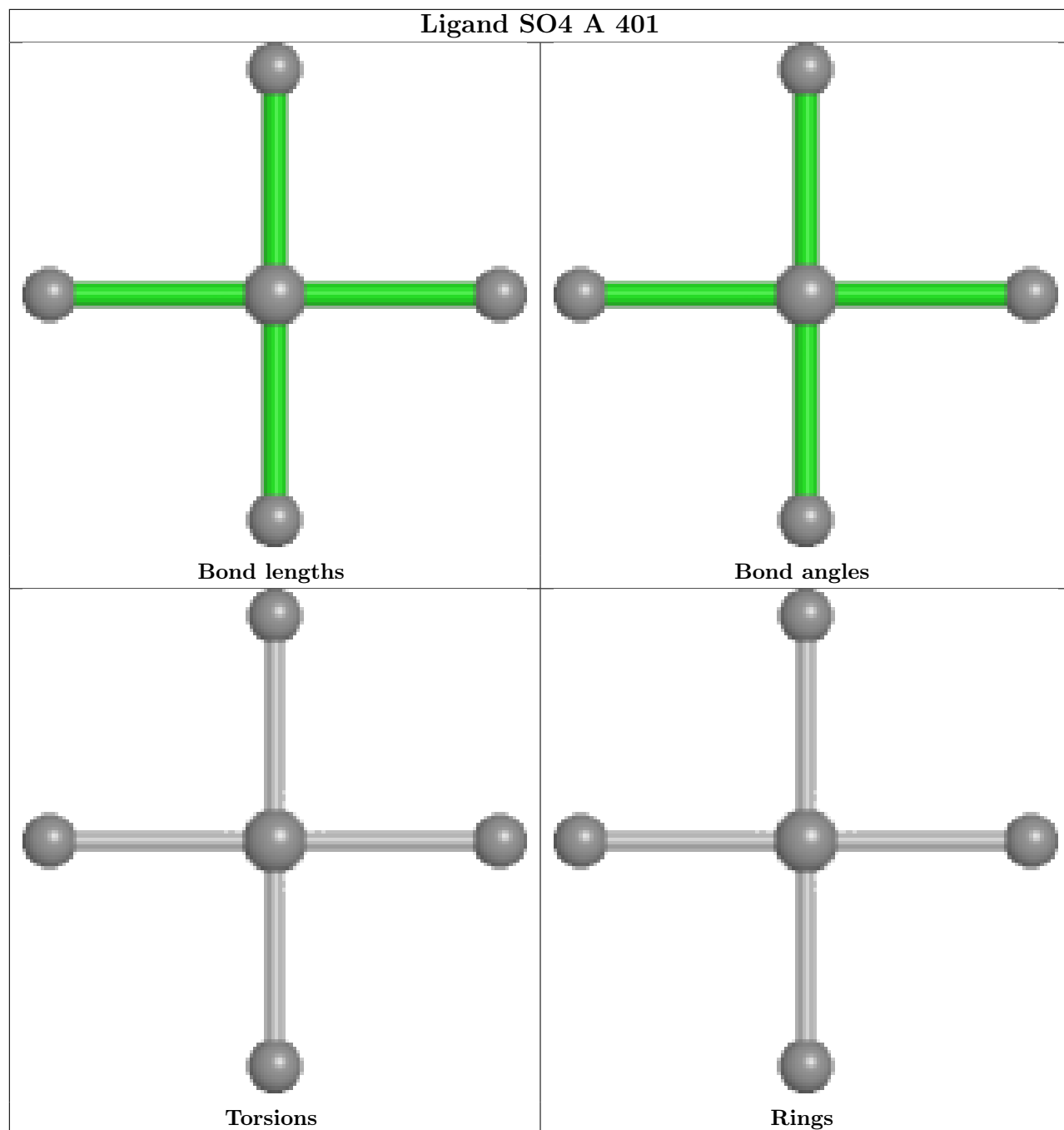


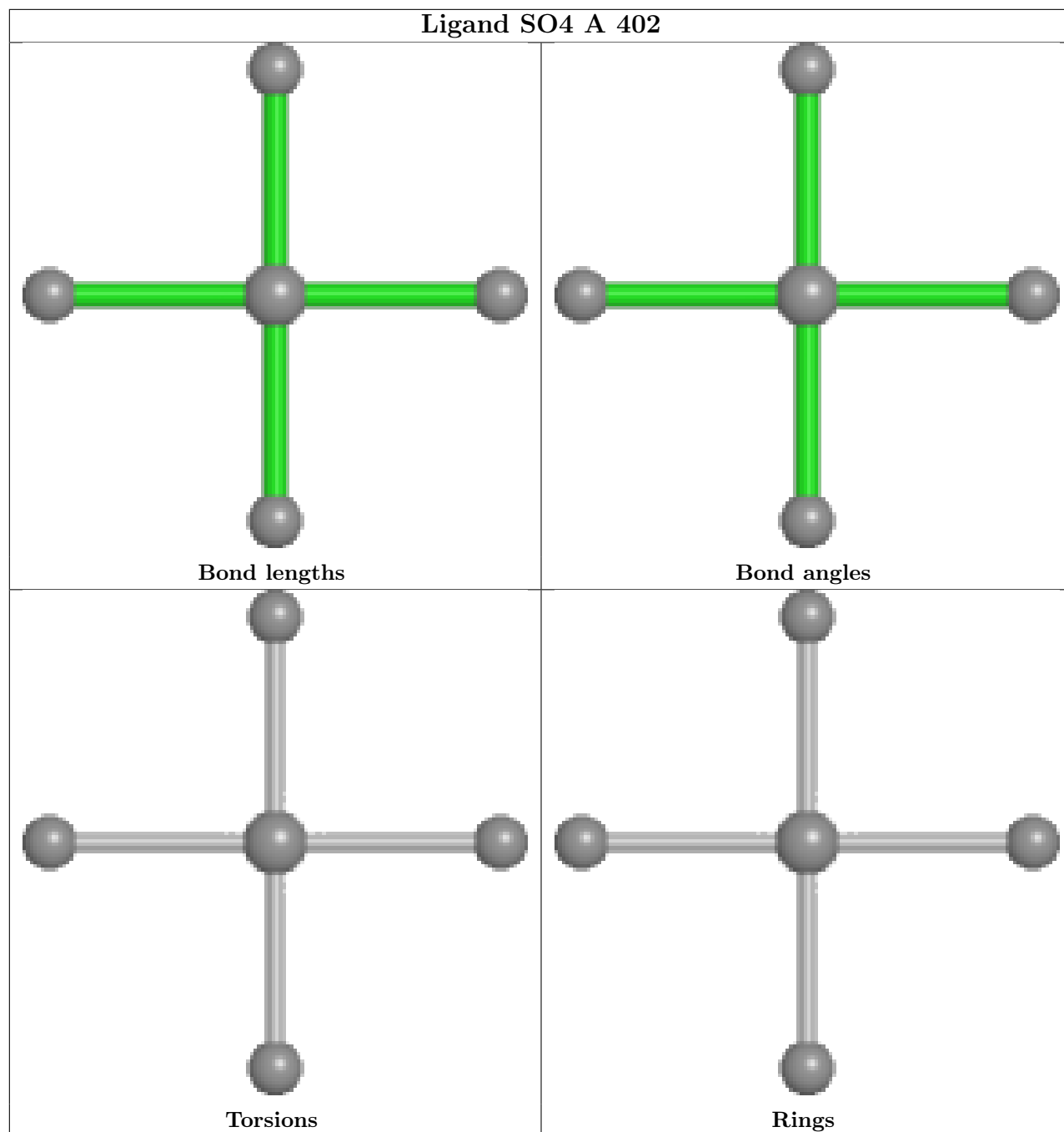












## 5.7 Other polymers [i](#)

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	340/349 (97%)	0.95	41 (12%)	10 10	40, 88, 135, 152	2 (0%)

All (41) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	103	TYR	5.4
1	A	114	PHE	4.8
1	A	162	PHE	4.0
1	A	10	ALA	3.8
1	A	101	LEU	3.7
1	A	21	GLY	3.6
1	A	118	SER	3.6
1	A	296	TRP	3.6
1	A	113	ASN	3.5
1	A	23	LEU	3.5
1	A	102	TYR	3.3
1	A	12	GLY	3.2
1	A	309	PRO	3.2
1	A	132	LEU	2.9
1	A	188	LEU	2.9
1	A	111	PHE	2.9
1	A	22	ALA	2.8
1	A	109	ALA	2.6
1	A	27	TRP	2.6
1	A	116	PRO	2.6
1	A	304	LYS	2.6
1	A	303	PRO	2.5
1	A	150	LEU	2.4
1	A	110	ASN	2.4
1	A	344	ILE	2.4
1	A	19	GLU	2.3
1	A	163	LEU	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	160	MET	2.3
1	A	154	VAL	2.3
1	A	286	ASN	2.3
1	A	93	TYR	2.3
1	A	305	ARG	2.3
1	A	90	PHE	2.2
1	A	100	PHE	2.2
1	A	26	ALA	2.2
1	A	225	GLU	2.2
1	A	108	MET	2.2
1	A	17	ARG	2.1
1	A	285	LEU	2.1
1	A	294	ASN	2.0
1	A	44	ARG	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 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
2	SO4	A	409	5/5	0.54	0.13	190,190,190,190	0
2	SO4	A	408	5/5	0.58	0.09	166,167,168,168	0
3	A1L4V	A	404[A]	28/28	0.62	0.38	129,153,170,170	28
3	A1L4V	A	404[B]	28/28	0.62	0.38	129,153,169,170	28
2	SO4	A	401	5/5	0.65	0.09	166,166,167,168	0
2	SO4	A	403	5/5	0.71	0.11	159,159,160,160	0
2	SO4	A	406	5/5	0.75	0.09	150,151,151,152	0
2	SO4	A	407	5/5	0.75	0.11	140,141,142,143	0
2	SO4	A	402	5/5	0.86	0.21	162,162,163,163	0

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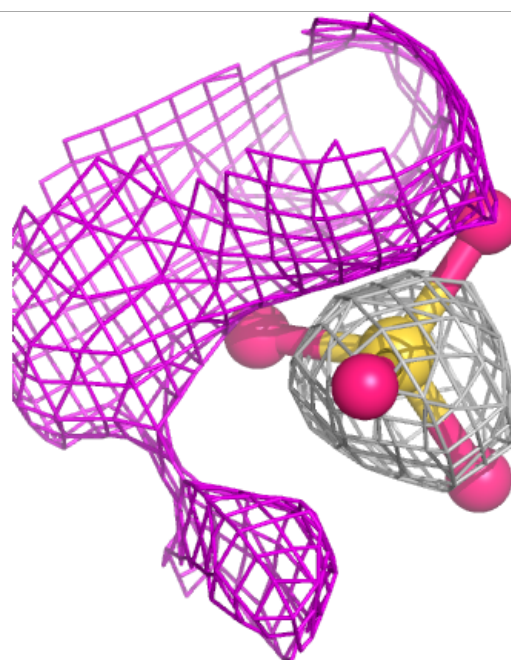
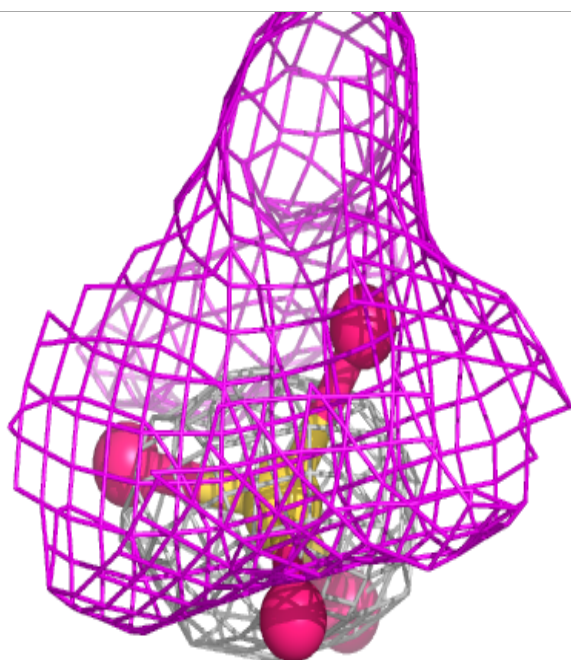
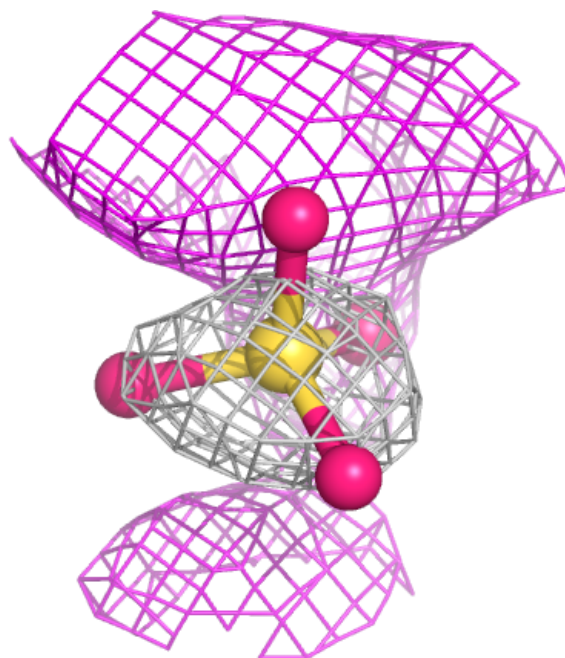
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
4	ZN	A	405	1/1	0.99	0.04	76,76,76,76	0

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.

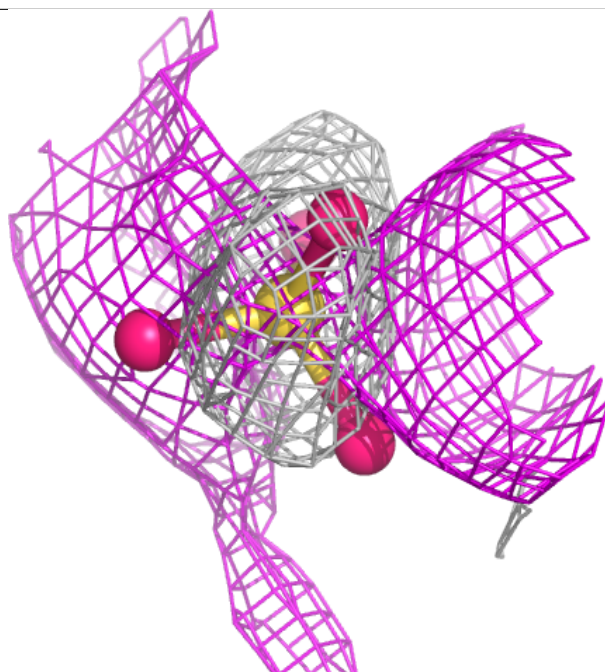
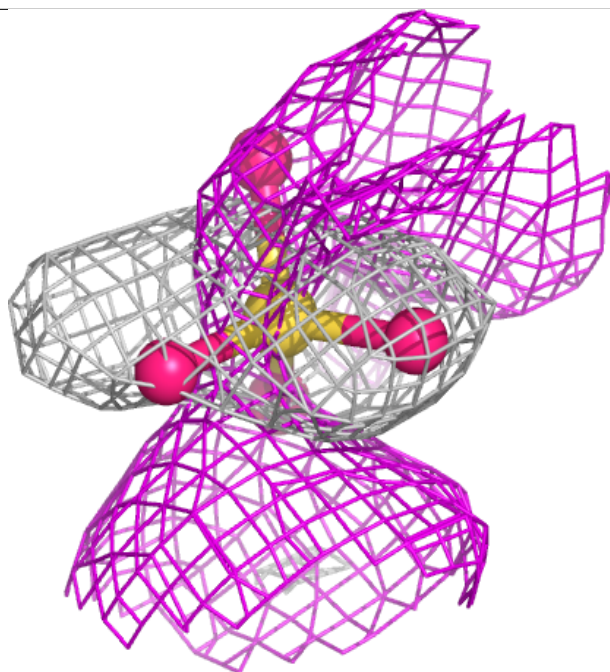
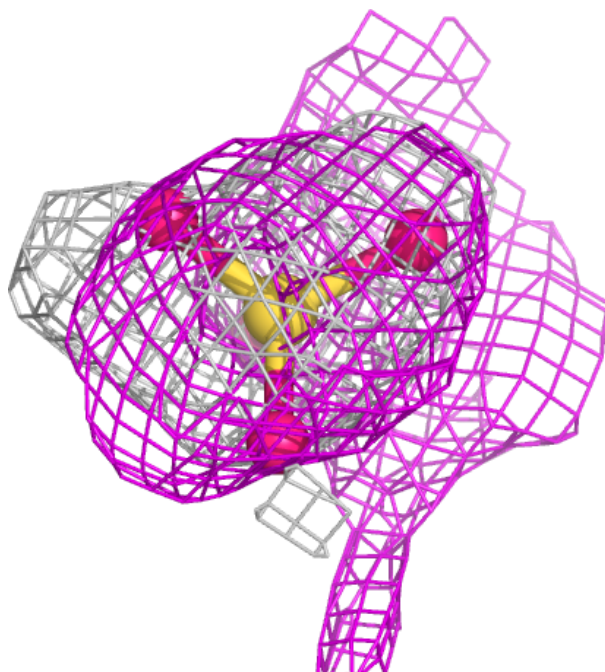
**Electron density around SO4 A 409:**

2mF<sub>o</sub>-DF<sub>c</sub> (at 0.7 rmsd) in gray  
mF<sub>o</sub>-DF<sub>c</sub> (at 3 rmsd) in purple (negative)  
and green (positive)



**Electron density around SO4 A 408:**

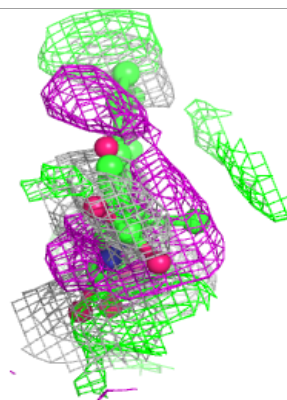
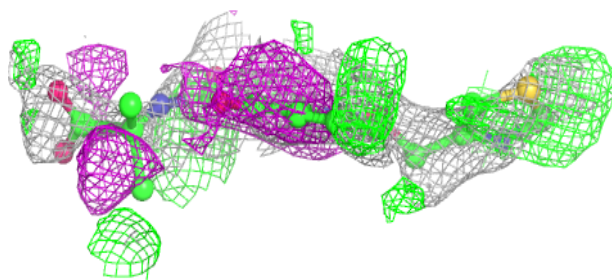
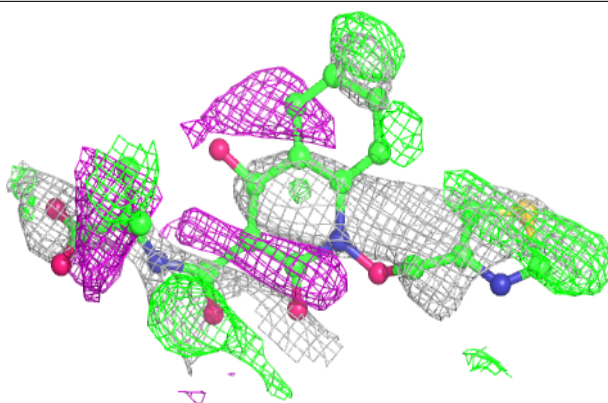
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



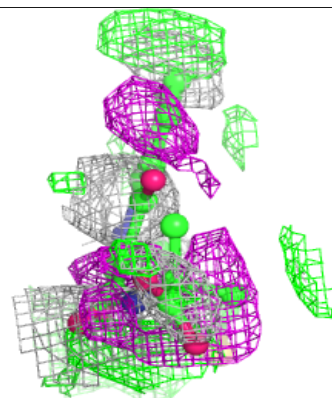
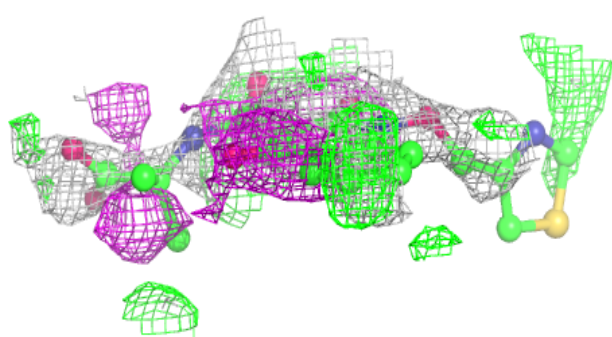
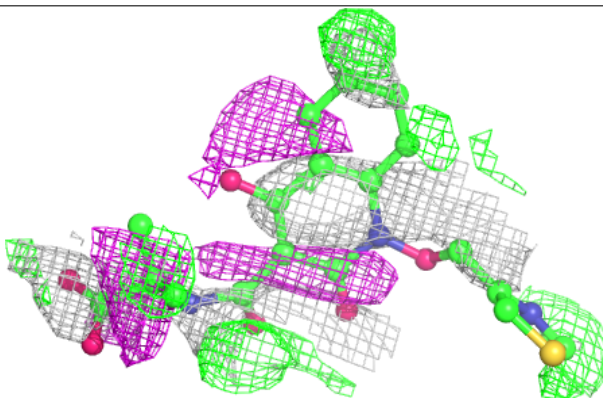


**Electron density around A1L4V A 404 (A):**

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

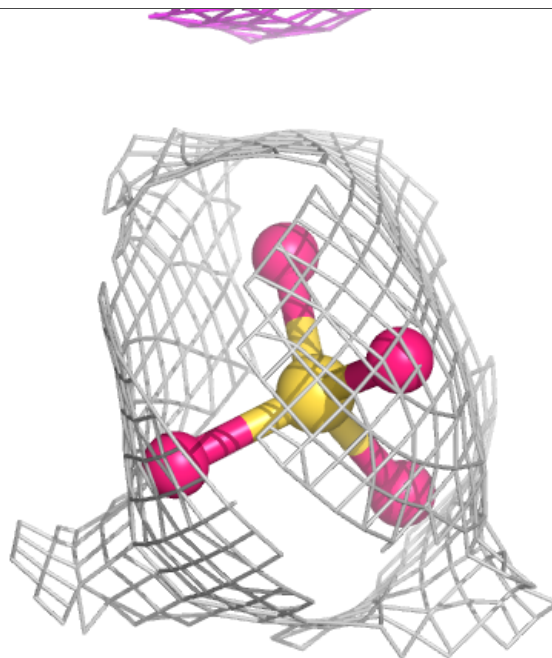
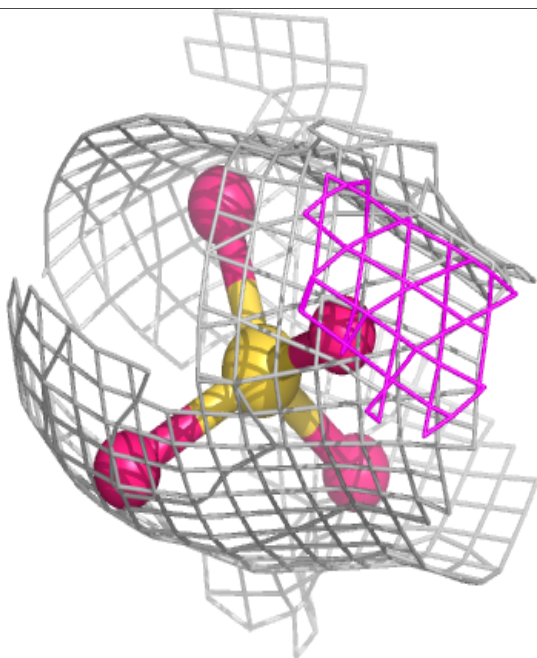
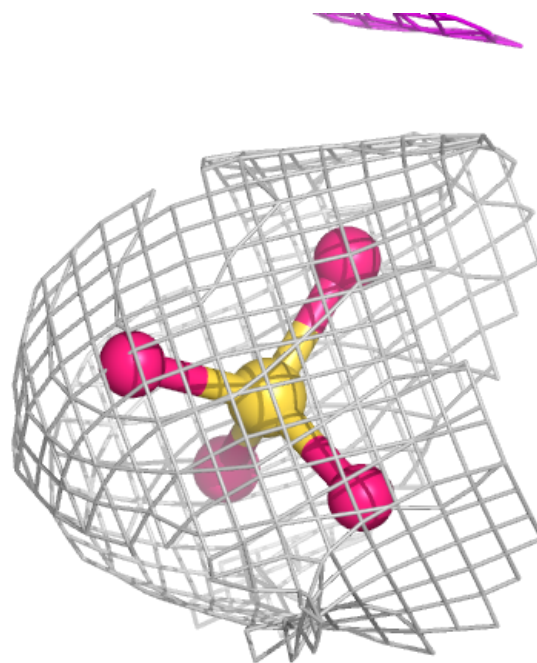
**Electron density around A1L4V A 404 (B):**

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



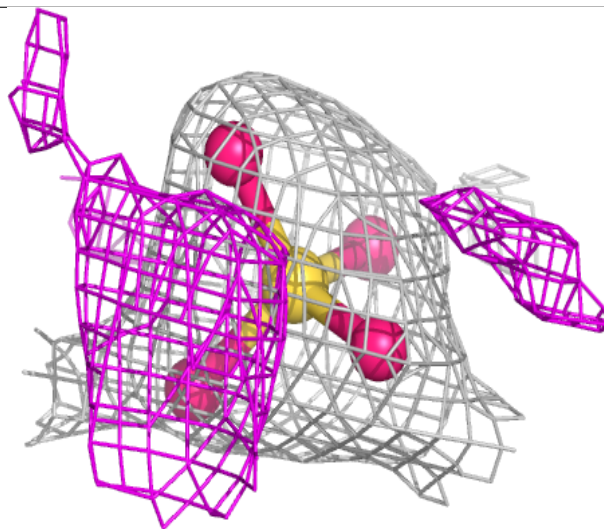
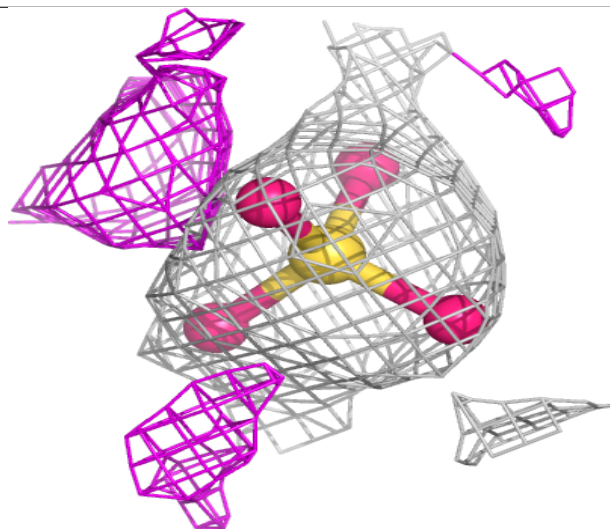
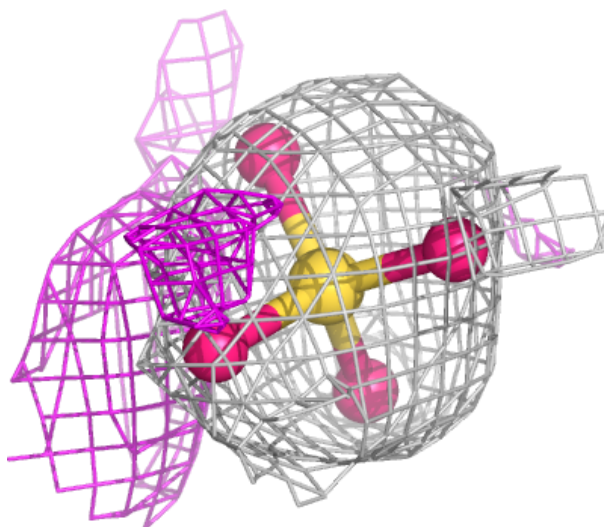
**Electron density around SO4 A 401:**

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



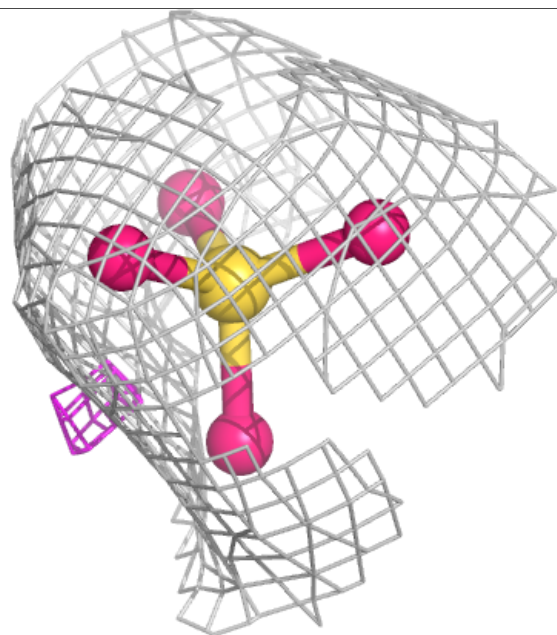
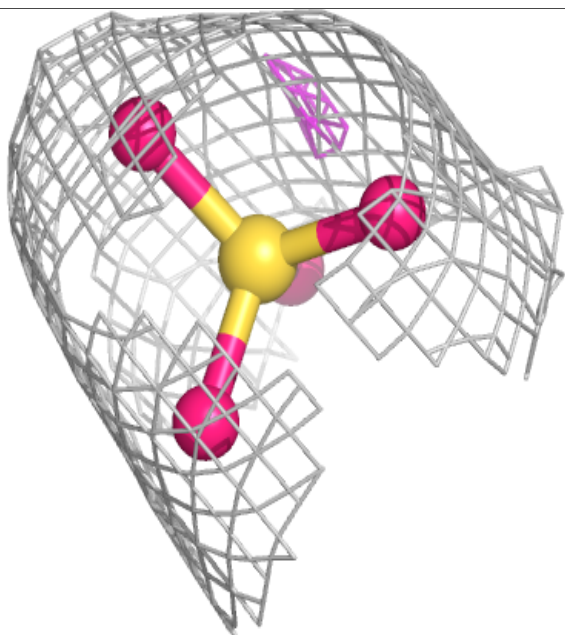
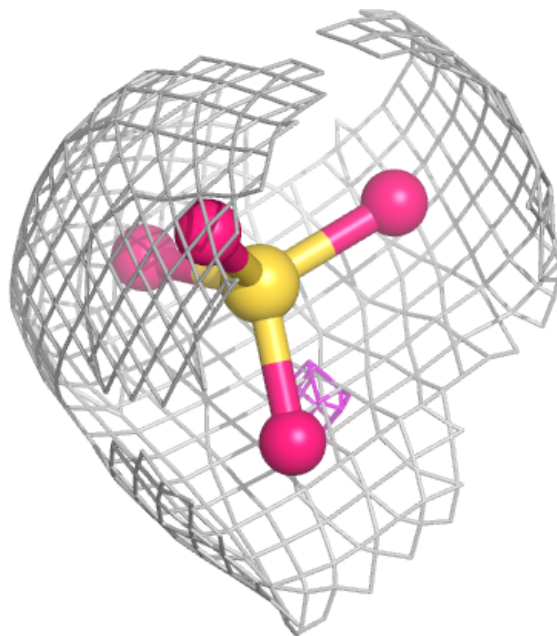
**Electron density around SO4 A 403:**

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and green (positive)



**Electron density around SO4 A 406:**

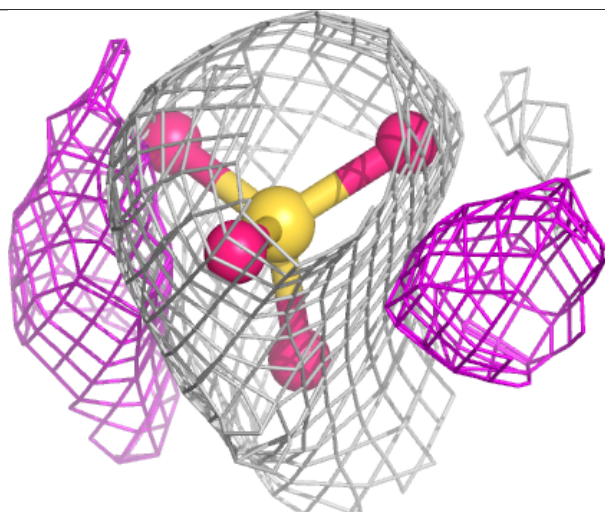
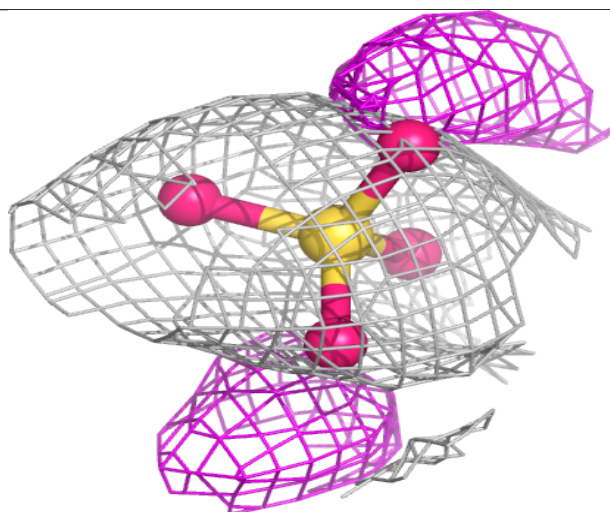
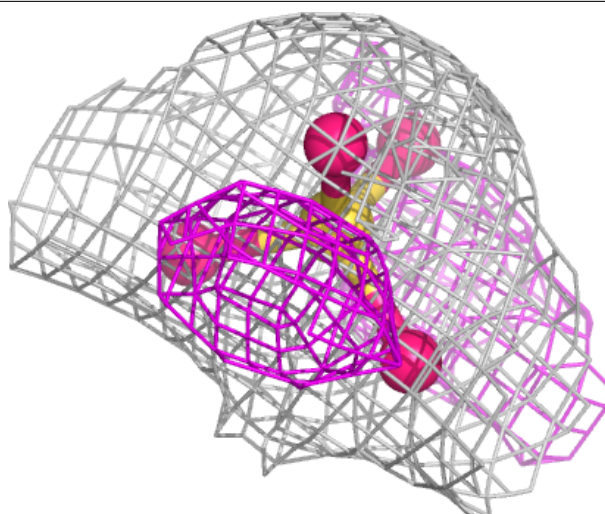
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 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





**Electron density around SO4 A 407:**

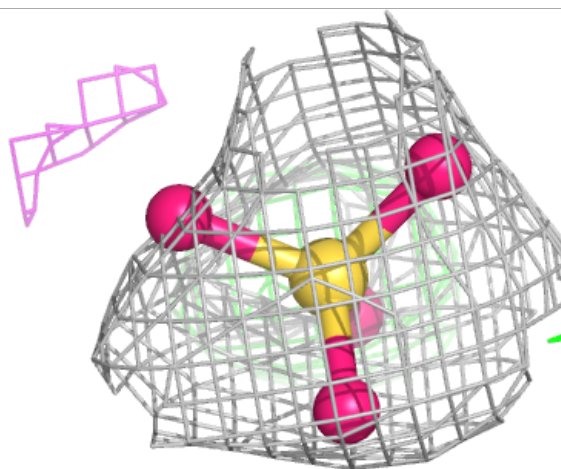
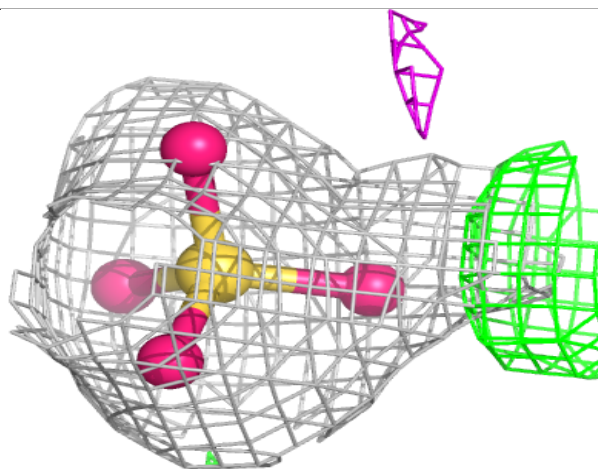
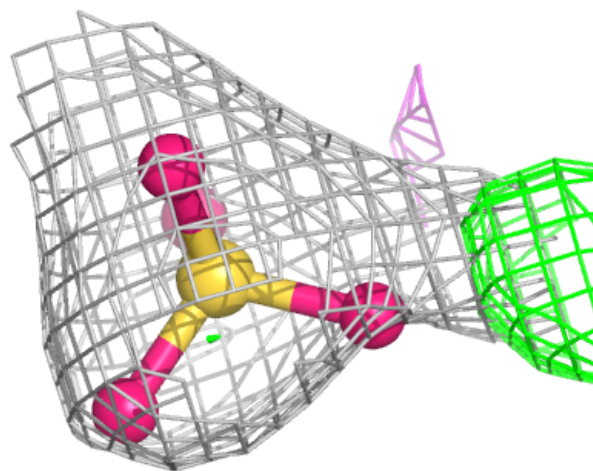
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 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





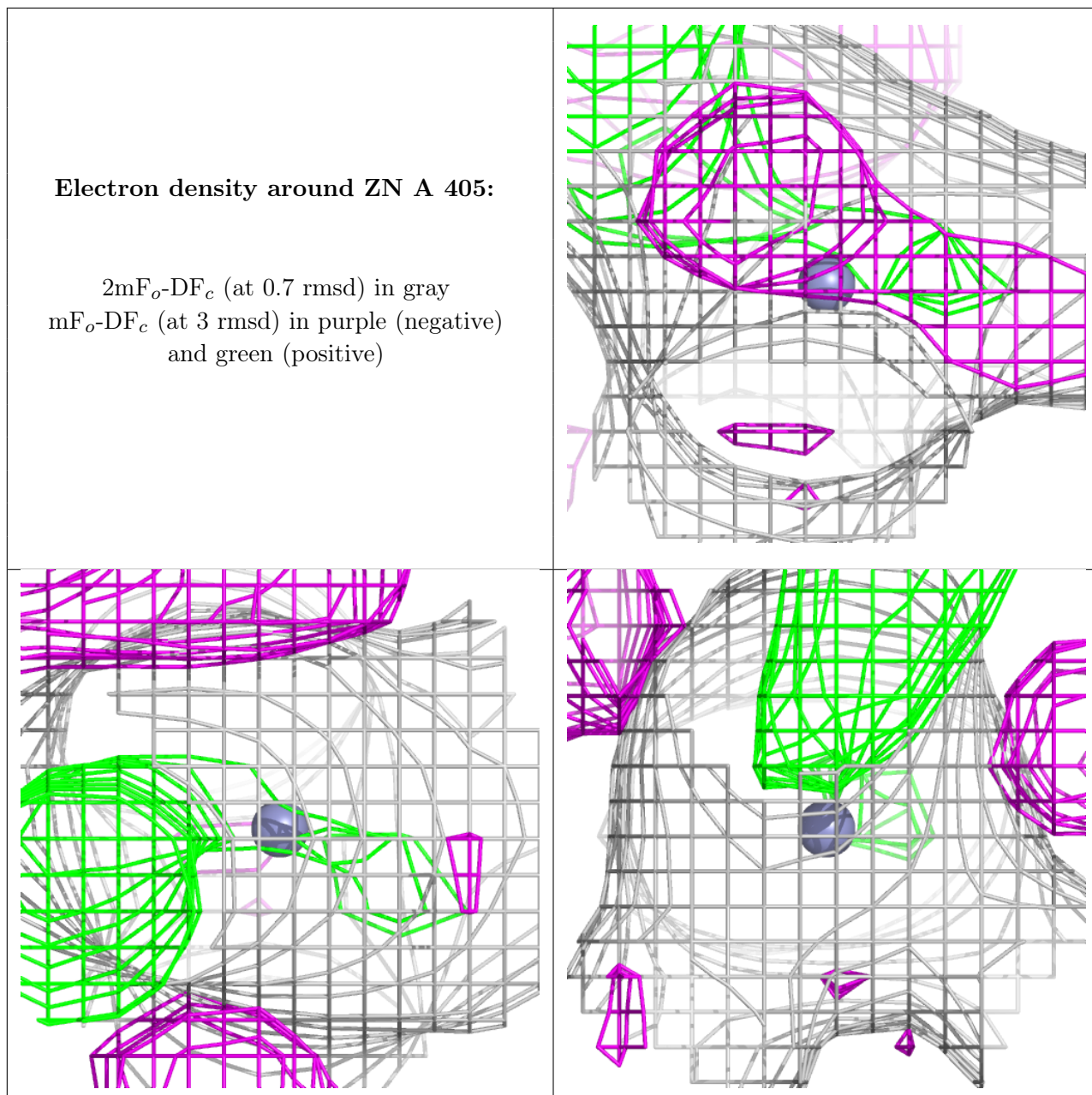
**Electron density around SO4 A 402:**

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



**Electron density around ZN A 405:**

$2mF_o-DF_c$  (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.