Full wwPDB X-ray Structure Validation Report ⓘ

Nov 19, 2023 – 05:08 PM JST

PDB ID	:	6LR3
Title	:	Structural and functional insights into macrophage migration inhibitory factor from Oncomelania hupensis, the intermediate host of Schistosoma japonicum
Authors	:	Su, Z.M.; Tian, X.Y.; Li, H.J.; Wei, Z.M.; Chen, L.F.; Ren, H.X.; Peng, W.F.; Tang, C.T.
Deposited on	:	2020-01-15
Resolution	:	1.77 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:

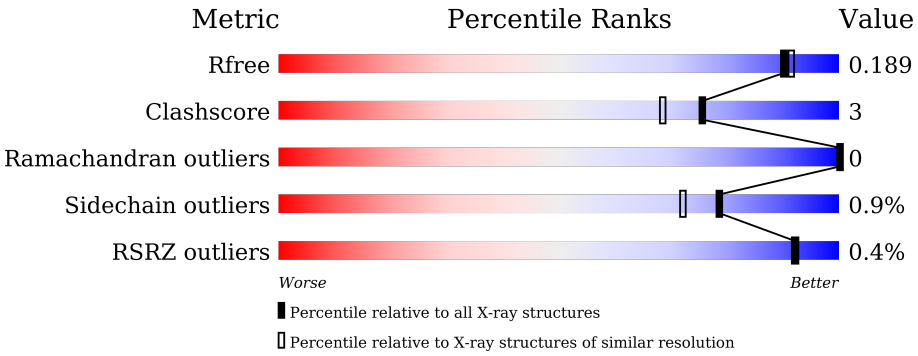
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 1.77 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	9185 (1.80-1.76)
Clashscore	141614	10184 (1.80-1.76)
Ramachandran outliers	138981	10051 (1.80-1.76)
Sidechain outliers	138945	10050 (1.80-1.76)
RSRZ outliers	127900	9032 (1.80-1.76)

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 <=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	131	83% 5% • 11%
1	B	131	% 84% 5% • 11%
1	C	131	% 81% 9% 10%
1	D	131	% 84% • • 12%
1	E	131	83% 5% 11%

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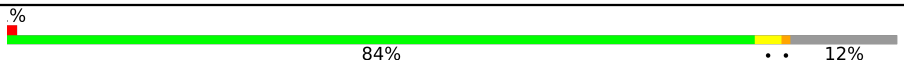
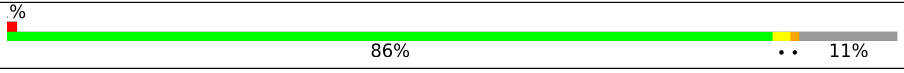
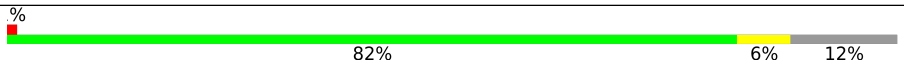
Mol	Chain	Length	Quality of chain
1	F	131	
1	G	131	
1	H	131	
1	I	131	
1	J	131	
1	K	131	
1	L	131	

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 12944 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 Macrophage migration inhibitory factor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	116	Total	C	N	O	S	0	1	0
			906	574	160	161	11			
1	C	118	Total	C	N	O	S	0	0	0
			918	582	161	164	11			
1	D	115	Total	C	N	O	S	0	0	0
			895	568	158	158	11			
1	B	117	Total	C	N	O	S	0	1	0
			914	580	161	162	11			
1	E	116	Total	C	N	O	S	0	0	0
			901	571	159	160	11			
1	F	115	Total	C	N	O	S	0	1	0
			900	571	159	159	11			
1	G	117	Total	C	N	O	S	0	1	0
			913	581	160	161	11			
1	H	114	Total	C	N	O	S	0	0	0
			888	563	157	157	11			
1	I	113	Total	C	N	O	S	0	0	0
			880	559	156	154	11			
1	J	117	Total	C	N	O	S	0	0	0
			909	577	160	161	11			
1	K	114	Total	C	N	O	S	0	0	0
			888	563	157	157	11			
1	L	115	Total	C	N	O	S	0	2	0
			901	573	158	159	11			

There are 132 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	89	ASN	LYS	conflict	UNP A0A1U9W5E8
A	90	VAL	ILE	conflict	UNP A0A1U9W5E8
A	123	TYR	-	expression tag	UNP A0A1U9W5E8
A	124	LEU	-	expression tag	UNP A0A1U9W5E8
A	125	GLU	-	expression tag	UNP A0A1U9W5E8

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Chain	Residue	Modelled	Actual	Comment	Reference
A	126	HIS	-	expression tag	UNP A0A1U9W5E8
A	127	HIS	-	expression tag	UNP A0A1U9W5E8
A	128	HIS	-	expression tag	UNP A0A1U9W5E8
A	129	HIS	-	expression tag	UNP A0A1U9W5E8
A	130	HIS	-	expression tag	UNP A0A1U9W5E8
A	131	HIS	-	expression tag	UNP A0A1U9W5E8
C	89	ASN	LYS	conflict	UNP A0A1U9W5E8
C	90	VAL	ILE	conflict	UNP A0A1U9W5E8
C	123	TYR	-	expression tag	UNP A0A1U9W5E8
C	124	LEU	-	expression tag	UNP A0A1U9W5E8
C	125	GLU	-	expression tag	UNP A0A1U9W5E8
C	126	HIS	-	expression tag	UNP A0A1U9W5E8
C	127	HIS	-	expression tag	UNP A0A1U9W5E8
C	128	HIS	-	expression tag	UNP A0A1U9W5E8
C	129	HIS	-	expression tag	UNP A0A1U9W5E8
C	130	HIS	-	expression tag	UNP A0A1U9W5E8
C	131	HIS	-	expression tag	UNP A0A1U9W5E8
D	89	ASN	LYS	conflict	UNP A0A1U9W5E8
D	90	VAL	ILE	conflict	UNP A0A1U9W5E8
D	123	TYR	-	expression tag	UNP A0A1U9W5E8
D	124	LEU	-	expression tag	UNP A0A1U9W5E8
D	125	GLU	-	expression tag	UNP A0A1U9W5E8
D	126	HIS	-	expression tag	UNP A0A1U9W5E8
D	127	HIS	-	expression tag	UNP A0A1U9W5E8
D	128	HIS	-	expression tag	UNP A0A1U9W5E8
D	129	HIS	-	expression tag	UNP A0A1U9W5E8
D	130	HIS	-	expression tag	UNP A0A1U9W5E8
D	131	HIS	-	expression tag	UNP A0A1U9W5E8
B	89	ASN	LYS	conflict	UNP A0A1U9W5E8
B	90	VAL	ILE	conflict	UNP A0A1U9W5E8
B	123	TYR	-	expression tag	UNP A0A1U9W5E8
B	124	LEU	-	expression tag	UNP A0A1U9W5E8
B	125	GLU	-	expression tag	UNP A0A1U9W5E8
B	126	HIS	-	expression tag	UNP A0A1U9W5E8
B	127	HIS	-	expression tag	UNP A0A1U9W5E8
B	128	HIS	-	expression tag	UNP A0A1U9W5E8
B	129	HIS	-	expression tag	UNP A0A1U9W5E8
B	130	HIS	-	expression tag	UNP A0A1U9W5E8
B	131	HIS	-	expression tag	UNP A0A1U9W5E8
E	89	ASN	LYS	conflict	UNP A0A1U9W5E8
E	90	VAL	ILE	conflict	UNP A0A1U9W5E8
E	123	TYR	-	expression tag	UNP A0A1U9W5E8

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Chain	Residue	Modelled	Actual	Comment	Reference
E	124	LEU	-	expression tag	UNP A0A1U9W5E8
E	125	GLU	-	expression tag	UNP A0A1U9W5E8
E	126	HIS	-	expression tag	UNP A0A1U9W5E8
E	127	HIS	-	expression tag	UNP A0A1U9W5E8
E	128	HIS	-	expression tag	UNP A0A1U9W5E8
E	129	HIS	-	expression tag	UNP A0A1U9W5E8
E	130	HIS	-	expression tag	UNP A0A1U9W5E8
E	131	HIS	-	expression tag	UNP A0A1U9W5E8
F	89	ASN	LYS	conflict	UNP A0A1U9W5E8
F	90	VAL	ILE	conflict	UNP A0A1U9W5E8
F	123	TYR	-	expression tag	UNP A0A1U9W5E8
F	124	LEU	-	expression tag	UNP A0A1U9W5E8
F	125	GLU	-	expression tag	UNP A0A1U9W5E8
F	126	HIS	-	expression tag	UNP A0A1U9W5E8
F	127	HIS	-	expression tag	UNP A0A1U9W5E8
F	128	HIS	-	expression tag	UNP A0A1U9W5E8
F	129	HIS	-	expression tag	UNP A0A1U9W5E8
F	130	HIS	-	expression tag	UNP A0A1U9W5E8
F	131	HIS	-	expression tag	UNP A0A1U9W5E8
G	89	ASN	LYS	conflict	UNP A0A1U9W5E8
G	90	VAL	ILE	conflict	UNP A0A1U9W5E8
G	123	TYR	-	expression tag	UNP A0A1U9W5E8
G	124	LEU	-	expression tag	UNP A0A1U9W5E8
G	125	GLU	-	expression tag	UNP A0A1U9W5E8
G	126	HIS	-	expression tag	UNP A0A1U9W5E8
G	127	HIS	-	expression tag	UNP A0A1U9W5E8
G	128	HIS	-	expression tag	UNP A0A1U9W5E8
G	129	HIS	-	expression tag	UNP A0A1U9W5E8
G	130	HIS	-	expression tag	UNP A0A1U9W5E8
G	131	HIS	-	expression tag	UNP A0A1U9W5E8
H	89	ASN	LYS	conflict	UNP A0A1U9W5E8
H	90	VAL	ILE	conflict	UNP A0A1U9W5E8
H	123	TYR	-	expression tag	UNP A0A1U9W5E8
H	124	LEU	-	expression tag	UNP A0A1U9W5E8
H	125	GLU	-	expression tag	UNP A0A1U9W5E8
H	126	HIS	-	expression tag	UNP A0A1U9W5E8
H	127	HIS	-	expression tag	UNP A0A1U9W5E8
H	128	HIS	-	expression tag	UNP A0A1U9W5E8
H	129	HIS	-	expression tag	UNP A0A1U9W5E8
H	130	HIS	-	expression tag	UNP A0A1U9W5E8
H	131	HIS	-	expression tag	UNP A0A1U9W5E8
I	89	ASN	LYS	conflict	UNP A0A1U9W5E8

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Chain	Residue	Modelled	Actual	Comment	Reference
I	90	VAL	ILE	conflict	UNP A0A1U9W5E8
I	123	TYR	-	expression tag	UNP A0A1U9W5E8
I	124	LEU	-	expression tag	UNP A0A1U9W5E8
I	125	GLU	-	expression tag	UNP A0A1U9W5E8
I	126	HIS	-	expression tag	UNP A0A1U9W5E8
I	127	HIS	-	expression tag	UNP A0A1U9W5E8
I	128	HIS	-	expression tag	UNP A0A1U9W5E8
I	129	HIS	-	expression tag	UNP A0A1U9W5E8
I	130	HIS	-	expression tag	UNP A0A1U9W5E8
I	131	HIS	-	expression tag	UNP A0A1U9W5E8
J	89	ASN	LYS	conflict	UNP A0A1U9W5E8
J	90	VAL	ILE	conflict	UNP A0A1U9W5E8
J	123	TYR	-	expression tag	UNP A0A1U9W5E8
J	124	LEU	-	expression tag	UNP A0A1U9W5E8
J	125	GLU	-	expression tag	UNP A0A1U9W5E8
J	126	HIS	-	expression tag	UNP A0A1U9W5E8
J	127	HIS	-	expression tag	UNP A0A1U9W5E8
J	128	HIS	-	expression tag	UNP A0A1U9W5E8
J	129	HIS	-	expression tag	UNP A0A1U9W5E8
J	130	HIS	-	expression tag	UNP A0A1U9W5E8
J	131	HIS	-	expression tag	UNP A0A1U9W5E8
K	89	ASN	LYS	conflict	UNP A0A1U9W5E8
K	90	VAL	ILE	conflict	UNP A0A1U9W5E8
K	123	TYR	-	expression tag	UNP A0A1U9W5E8
K	124	LEU	-	expression tag	UNP A0A1U9W5E8
K	125	GLU	-	expression tag	UNP A0A1U9W5E8
K	126	HIS	-	expression tag	UNP A0A1U9W5E8
K	127	HIS	-	expression tag	UNP A0A1U9W5E8
K	128	HIS	-	expression tag	UNP A0A1U9W5E8
K	129	HIS	-	expression tag	UNP A0A1U9W5E8
K	130	HIS	-	expression tag	UNP A0A1U9W5E8
K	131	HIS	-	expression tag	UNP A0A1U9W5E8
L	89	ASN	LYS	conflict	UNP A0A1U9W5E8
L	90	VAL	ILE	conflict	UNP A0A1U9W5E8
L	123	TYR	-	expression tag	UNP A0A1U9W5E8
L	124	LEU	-	expression tag	UNP A0A1U9W5E8
L	125	GLU	-	expression tag	UNP A0A1U9W5E8
L	126	HIS	-	expression tag	UNP A0A1U9W5E8
L	127	HIS	-	expression tag	UNP A0A1U9W5E8
L	128	HIS	-	expression tag	UNP A0A1U9W5E8
L	129	HIS	-	expression tag	UNP A0A1U9W5E8
L	130	HIS	-	expression tag	UNP A0A1U9W5E8

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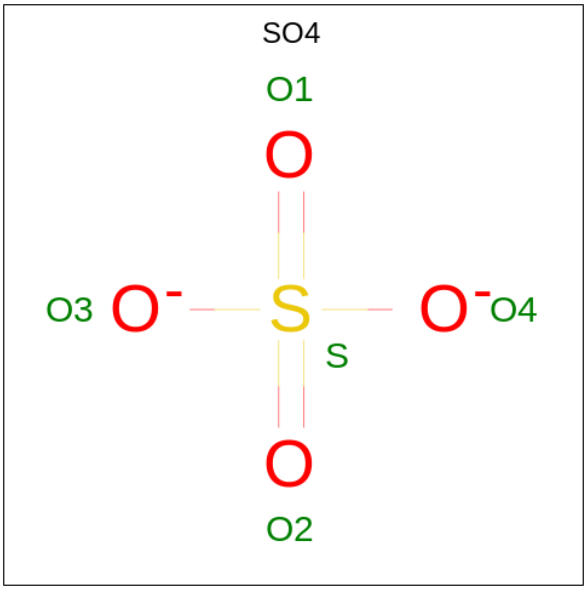
Chain	Residue	Modelled	Actual	Comment	Reference
L	131	HIS	-	expression tag	UNP A0A1U9W5E8

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O_4S) (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	D	1	Total	O	S		0	0
			5	4	1			
2	B	1	Total	O	S		0	0
			5	4	1			
2	F	1	Total	O	S		0	0
			5	4	1			
2	G	1	Total	O	S		0	0
			5	4	1			
2	G	1	Total	O	S		0	0
			5	4	1			
2	J	1	Total	O	S		0	0
			5	4	1			
2	J	1	Total	O	S		0	0
			5	4	1			

- Molecule 3 is water.

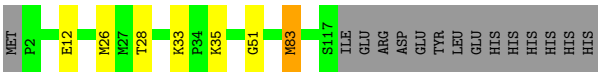
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	198	Total O 198 198	0	0
3	C	167	Total O 167 167	0	0
3	D	167	Total O 167 167	0	0
3	B	196	Total O 196 196	0	0
3	E	182	Total O 182 182	0	0
3	F	165	Total O 165 165	0	0
3	G	192	Total O 192 192	0	0
3	H	157	Total O 157 157	0	0
3	I	170	Total O 170 170	0	0
3	J	185	Total O 185 185	0	0
3	K	145	Total O 145 145	0	0
3	L	167	Total O 167 167	0	0

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.

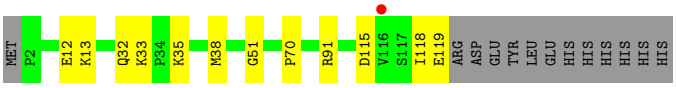
• Molecule 1: Macrophage migration inhibitory factor

Chain A:



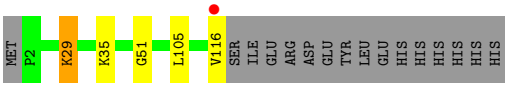
• Molecule 1: Macrophage migration inhibitory factor

Chain C:



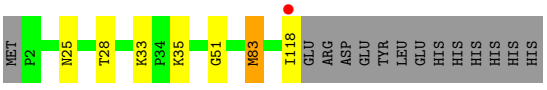
• Molecule 1: Macrophage migration inhibitory factor

Chain D:



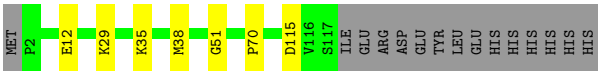
• Molecule 1: Macrophage migration inhibitory factor

Chain B:



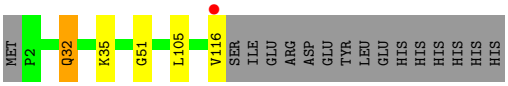
• Molecule 1: Macrophage migration inhibitory factor

Chain E:



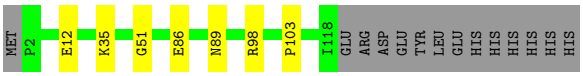
• Molecule 1: Macrophage migration inhibitory factor

Chain F:



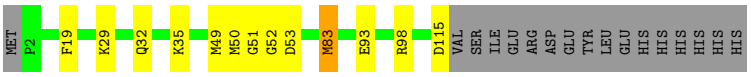
• Molecule 1: Macrophage migration inhibitory factor

Chain G:



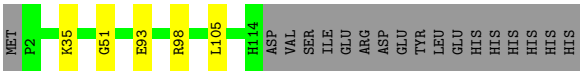
• Molecule 1: Macrophage migration inhibitory factor

Chain H:



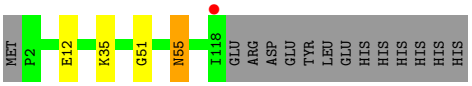
• Molecule 1: Macrophage migration inhibitory factor

Chain I:



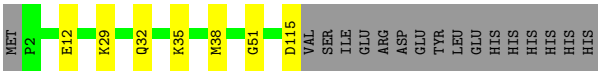
• Molecule 1: Macrophage migration inhibitory factor

Chain J:



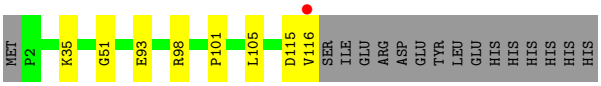
• Molecule 1: Macrophage migration inhibitory factor

Chain K:



• Molecule 1: Macrophage migration inhibitory factor

Chain L:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	49.30Å 139.48Å 105.67Å 90.00° 89.81° 90.00°	Depositor
Resolution (Å)	49.35 – 1.77 49.30 – 1.77	Depositor EDS
% Data completeness (in resolution range)	96.2 (49.35-1.77) 95.3 (49.30-1.77)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	4.41 (at 1.77Å)	Xtriage
Refinement program	REFMAC 5.8.0258	Depositor
R, R_{free}	0.150 , 0.180 0.161 , 0.189	Depositor DCC
R_{free} test set	6675 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å^2)	18.5	Xtriage
Anisotropy	0.033	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.33 , 36.6	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	0.478 for h,-k,-l	Xtriage
F_o,F_c correlation	0.97	EDS
Total number of atoms	12944	wwPDB-VP
Average B, all atoms (Å^2)	21.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.02% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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	0.73	0/925	0.87	0/1250
1	B	0.73	0/933	0.86	0/1261
1	C	0.67	0/934	0.87	0/1262
1	D	0.71	0/911	0.87	0/1231
1	E	0.67	0/917	0.88	0/1239
1	F	0.72	0/919	0.86	0/1242
1	G	0.68	0/932	0.83	1/1260 (0.1%)
1	H	0.72	1/904 (0.1%)	0.88	2/1221 (0.2%)
1	I	0.75	1/896 (0.1%)	0.84	1/1210 (0.1%)
1	J	0.68	0/925	0.82	0/1250
1	K	0.71	0/904	0.88	0/1221
1	L	0.75	1/923 (0.1%)	0.84	1/1247 (0.1%)
All	All	0.71	3/11023 (0.0%)	0.86	5/14894 (0.0%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	L	93	GLU	CD-OE1	5.44	1.31	1.25
1	I	93	GLU	CD-OE1	5.11	1.31	1.25
1	H	93	GLU	CD-OE1	5.07	1.31	1.25

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	G	98	ARG	NE-CZ-NH2	-5.57	117.52	120.30
1	H	98	ARG	NE-CZ-NH2	-5.53	117.53	120.30
1	H	83	MET	CG-SD-CE	-5.35	91.64	100.20
1	L	98	ARG	NE-CZ-NH2	-5.19	117.71	120.30
1	I	98	ARG	NE-CZ-NH2	-5.13	117.74	120.30

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	906	0	926	6	0
1	B	914	0	937	5	0
1	C	918	0	937	16	0
1	D	895	0	915	5	0
1	E	901	0	920	10	0
1	F	900	0	921	7	0
1	G	913	0	940	6	0
1	H	888	0	906	9	0
1	I	880	0	902	8	0
1	J	909	0	931	4	0
1	K	888	0	906	5	0
1	L	901	0	924	10	0
2	A	5	0	0	0	0
2	B	5	0	0	0	0
2	D	5	0	0	0	0
2	F	5	0	0	0	0
2	G	10	0	0	0	0
2	J	10	0	0	0	0
3	A	198	0	0	1	0
3	B	196	0	0	1	0
3	C	167	0	0	5	0
3	D	167	0	0	2	0
3	E	182	0	0	5	0
3	F	165	0	0	4	0
3	G	192	0	0	3	0
3	H	157	0	0	6	0
3	I	170	0	0	0	0
3	J	185	0	0	1	0
3	K	145	0	0	3	0
3	L	167	0	0	1	0
All	All	12944	0	11065	67	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 3.

All (67) 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:12:GLU:HG3	3:A:359:HOH:O	1.43	1.19
1:C:32:GLN:NE2	1:I:105:LEU:HD12	1.56	1.17
1:E:12:GLU:HG3	3:E:295:HOH:O	1.57	1.05
1:C:32:GLN:HE22	1:I:105:LEU:CD1	1.73	1.01
1:C:12:GLU:HG3	3:C:276:HOH:O	1.64	0.98
1:H:32:GLN:HG3	3:H:251:HOH:O	1.62	0.97
1:C:32:GLN:NE2	1:I:105:LEU:CD1	2.29	0.94
1:C:32:GLN:HE22	1:I:105:LEU:HD12	1.26	0.83
1:K:32:GLN:HG3	3:K:262:HOH:O	1.78	0.81
1:G:86:GLU:HG2	3:G:437:HOH:O	1.86	0.76
3:E:373:HOH:O	1:L:116:VAL:HG21	1.86	0.74
1:C:32:GLN:HE22	1:I:105:LEU:CG	2.02	0.71
1:C:38:MET:HE2	3:C:283:HOH:O	1.90	0.71
1:C:91:ARG:NH2	3:C:201:HOH:O	2.26	0.69
1:B:25[B]:ASN:ND2	3:B:301:HOH:O	2.26	0.65
1:K:38:MET:HE2	3:K:222:HOH:O	1.97	0.65
1:H:53:ASP:N	3:H:201:HOH:O	2.30	0.63
1:F:116:VAL:HA	3:F:370:HOH:O	1.97	0.63
1:A:26:MET:SD	1:A:83:MET:CE	2.87	0.63
1:G:89:ASN:HB3	3:G:403:HOH:O	1.99	0.63
1:A:35:LYS:O	1:D:51:GLY:HA2	2.00	0.61
1:C:33:LYS:NZ	1:C:119:GLU:OE1	2.33	0.60
1:E:38:MET:HE2	3:E:288:HOH:O	2.02	0.59
1:B:35:LYS:O	1:F:51:GLY:HA2	2.03	0.58
1:J:12:GLU:HG2	3:J:439:HOH:O	2.03	0.58
1:E:29:LYS:HD3	1:L:116:VAL:CG1	2.35	0.56
1:E:38:MET:CE	3:E:288:HOH:O	2.53	0.56
1:G:12:GLU:HG3	3:G:443:HOH:O	2.05	0.56
1:B:51:GLY:HA2	1:E:35:LYS:O	2.05	0.55
1:A:51:GLY:HA2	1:C:35:LYS:O	2.06	0.54
1:J:55:ASN:N	1:J:55:ASN:HD22	2.06	0.54
1:C:51:GLY:HA2	1:D:35:LYS:O	2.08	0.53
1:L:105:LEU:CD2	1:L:116:VAL:HG22	2.38	0.53
1:E:51:GLY:HA2	1:F:35:LYS:O	2.10	0.52
1:C:115:ASP:HB2	3:C:320:HOH:O	2.09	0.52
1:E:29:LYS:HD3	1:L:116:VAL:HG11	1.92	0.52
1:H:115:ASP:HA	3:H:295:HOH:O	2.10	0.51
1:D:29:LYS:NZ	3:D:302:HOH:O	2.45	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:32:GLN:HE22	1:I:105:LEU:HG	1.76	0.50
1:G:51:GLY:HA2	1:H:35:LYS:O	2.12	0.50
1:H:49:MET:HB2	3:H:201:HOH:O	2.13	0.49
1:L:105:LEU:HD21	1:L:116:VAL:HG22	1.94	0.49
1:E:115:ASP:HB2	3:E:335:HOH:O	2.11	0.48
1:J:51:GLY:HA2	1:K:35:LYS:O	2.14	0.48
1:G:35:LYS:O	1:I:51:GLY:HA2	2.14	0.48
1:F:32:GLN:HG3	3:F:307:HOH:O	2.12	0.47
1:J:35:LYS:O	1:L:51:GLY:HA2	2.15	0.47
1:K:51:GLY:HA2	1:L:35:LYS:O	2.15	0.46
1:F:116:VAL:CA	3:F:370:HOH:O	2.61	0.46
1:C:118:ILE:HD11	1:G:103:PRO:HG2	1.98	0.46
1:H:50:MET:N	3:H:201:HOH:O	2.48	0.46
1:H:51:GLY:HA2	1:I:35:LYS:O	2.16	0.46
1:F:116:VAL:C	3:F:370:HOH:O	2.54	0.45
1:D:116:VAL:HA	3:D:381:HOH:O	2.17	0.45
1:E:29:LYS:HE2	1:L:115:ASP:CB	2.47	0.44
1:C:38:MET:CE	3:C:283:HOH:O	2.57	0.43
1:L:116:VAL:C	3:L:275:HOH:O	2.57	0.43
1:D:105:LEU:HD11	1:E:70:PRO:HG2	2.00	0.43
1:A:28:THR:HG23	1:A:33:LYS:O	2.19	0.42
1:B:83:MET:HE2	1:B:83:MET:HB2	1.88	0.42
1:K:12:GLU:HG3	3:K:314:HOH:O	2.19	0.42
1:C:70:PRO:HG2	1:F:105:LEU:HD11	2.02	0.41
1:A:26:MET:SD	1:A:83:MET:HE3	2.60	0.41
1:B:28:THR:HG23	1:B:33:LYS:O	2.20	0.41
1:L:105:LEU:HD22	1:L:116:VAL:HG22	2.03	0.41
1:H:52:GLY:N	3:H:201:HOH:O	2.53	0.41
1:H:19:PHE:CD1	1:H:83:MET:CE	3.05	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	115/131 (88%)	114 (99%)	1 (1%)	0	100	100
1	B	116/131 (88%)	115 (99%)	1 (1%)	0	100	100
1	C	116/131 (88%)	114 (98%)	2 (2%)	0	100	100
1	D	113/131 (86%)	112 (99%)	1 (1%)	0	100	100
1	E	114/131 (87%)	112 (98%)	2 (2%)	0	100	100
1	F	114/131 (87%)	113 (99%)	1 (1%)	0	100	100
1	G	116/131 (88%)	115 (99%)	1 (1%)	0	100	100
1	H	112/131 (86%)	111 (99%)	1 (1%)	0	100	100
1	I	111/131 (85%)	110 (99%)	1 (1%)	0	100	100
1	J	115/131 (88%)	114 (99%)	1 (1%)	0	100	100
1	K	112/131 (86%)	111 (99%)	1 (1%)	0	100	100
1	L	115/131 (88%)	114 (99%)	1 (1%)	0	100	100
All	All	1369/1572 (87%)	1355 (99%)	14 (1%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains (i)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	102/116 (88%)	101 (99%)	1 (1%)	76	68
1	B	103/116 (89%)	101 (98%)	2 (2%)	57	43
1	C	103/116 (89%)	102 (99%)	1 (1%)	76	68
1	D	100/116 (86%)	99 (99%)	1 (1%)	76	68
1	E	101/116 (87%)	101 (100%)	0	100	100
1	F	101/116 (87%)	100 (99%)	1 (1%)	76	68
1	G	103/116 (89%)	103 (100%)	0	100	100
1	H	99/116 (85%)	98 (99%)	1 (1%)	76	68
1	I	98/116 (84%)	98 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	J	102/116 (88%)	101 (99%)	1 (1%)	76 68
1	K	99/116 (85%)	97 (98%)	2 (2%)	55 40
1	L	101/116 (87%)	100 (99%)	1 (1%)	76 68
All	All	1212/1392 (87%)	1201 (99%)	11 (1%)	78 72

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	83	MET
1	C	13	LYS
1	D	29	LYS
1	B	83	MET
1	B	118	ILE
1	F	32	GLN
1	H	29	LYS
1	J	55	ASN
1	K	29	LYS
1	K	115	ASP
1	L	101	PRO

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	32	GLN
1	I	114	HIS
1	J	55	ASN
1	L	114	HIS

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

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	G	202	-	4,4,4	0.73	0	6,6,6	0.26	0
2	SO4	A	201	-	4,4,4	0.60	0	6,6,6	0.33	0
2	SO4	B	201	-	4,4,4	0.68	0	6,6,6	0.24	0
2	SO4	D	201	-	4,4,4	0.38	0	6,6,6	0.28	0
2	SO4	J	202	-	4,4,4	0.56	0	6,6,6	0.24	0
2	SO4	G	201	-	4,4,4	0.29	0	6,6,6	0.12	0
2	SO4	F	201	-	4,4,4	0.26	0	6,6,6	0.19	0
2	SO4	J	201	-	4,4,4	0.31	0	6,6,6	0.23	0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

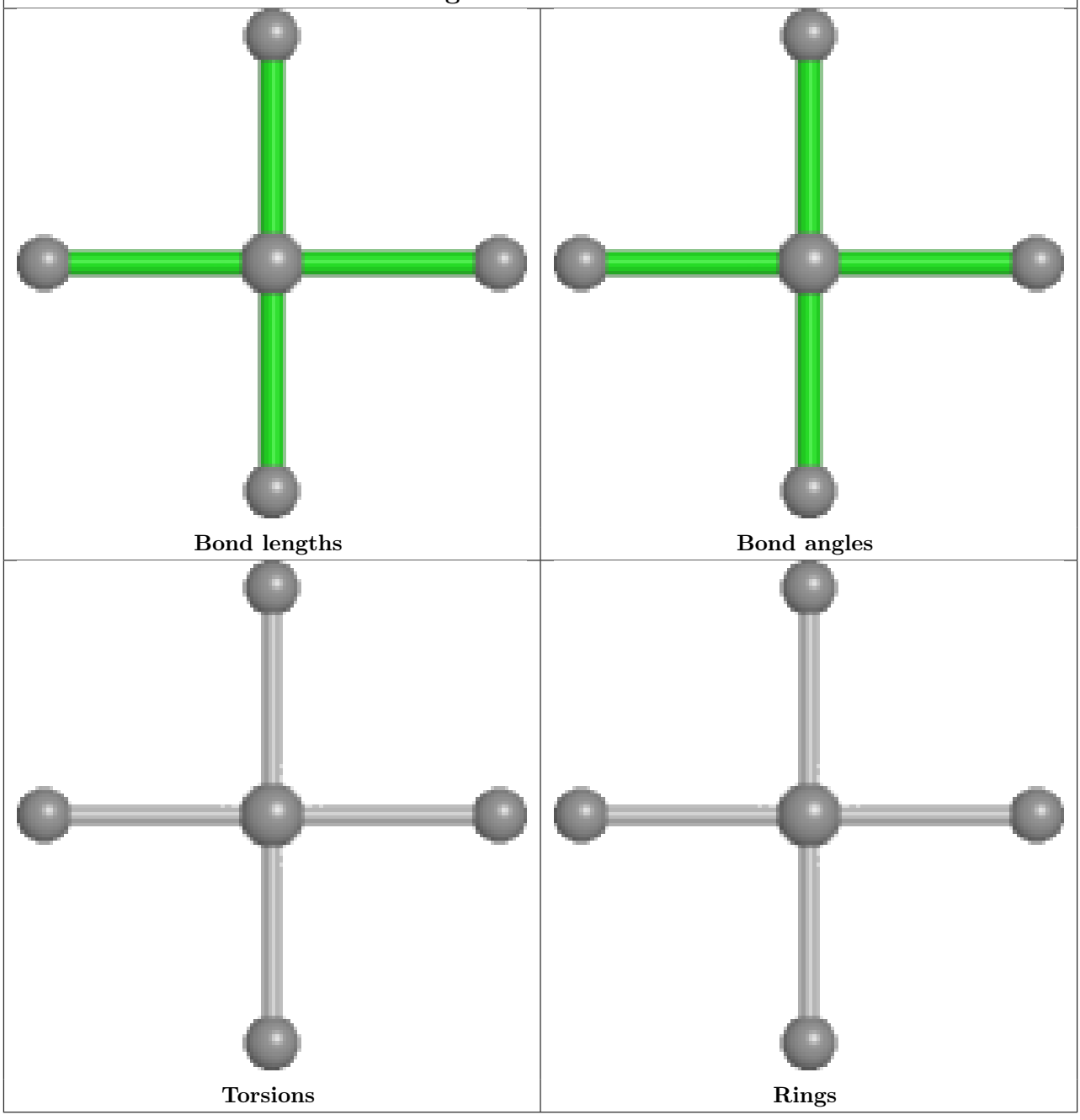
There are no ring outliers.

No monomer is involved in short contacts.

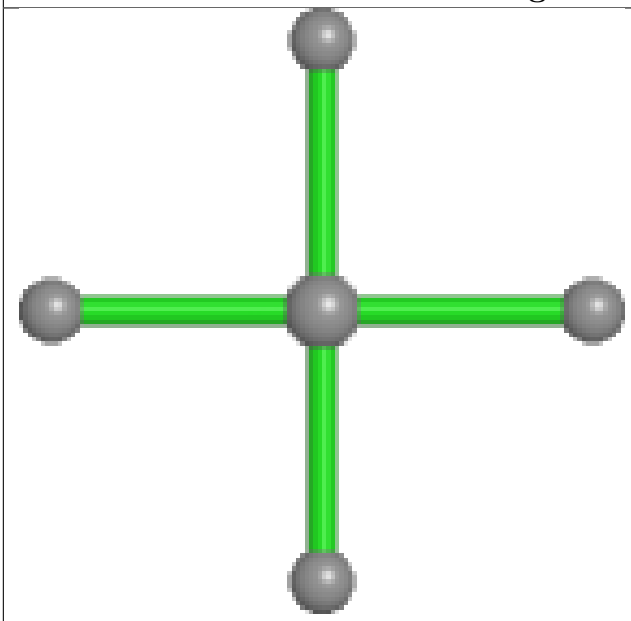
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

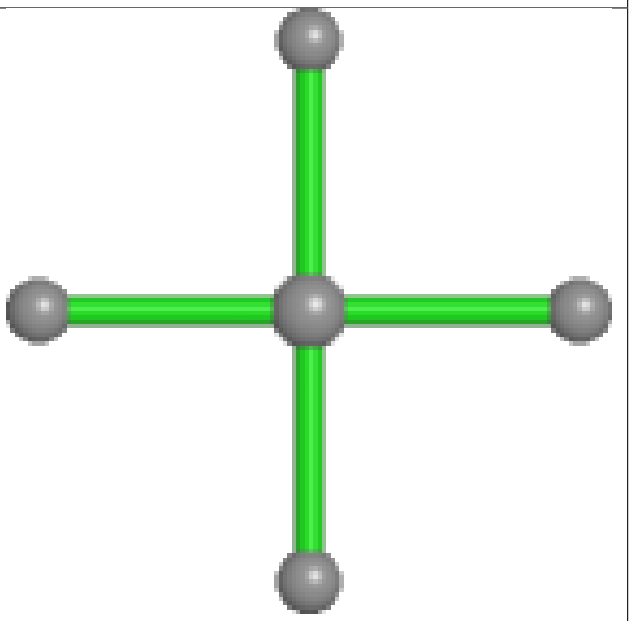
Ligand SO4 G 202



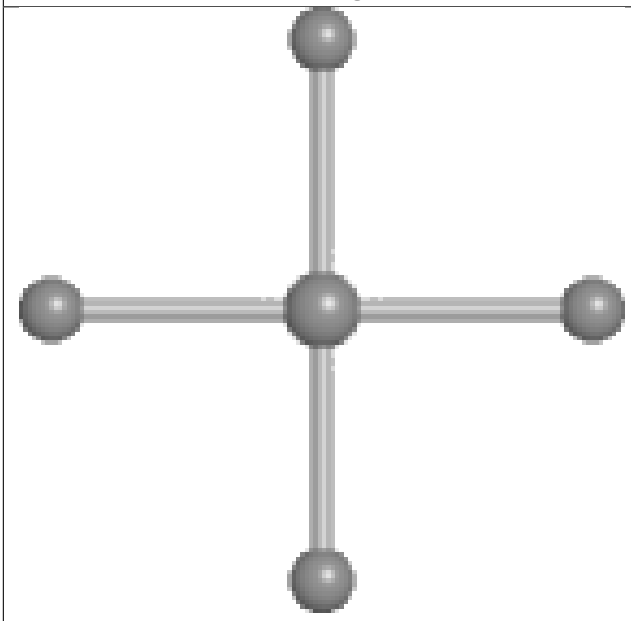
Ligand SO4 A 201



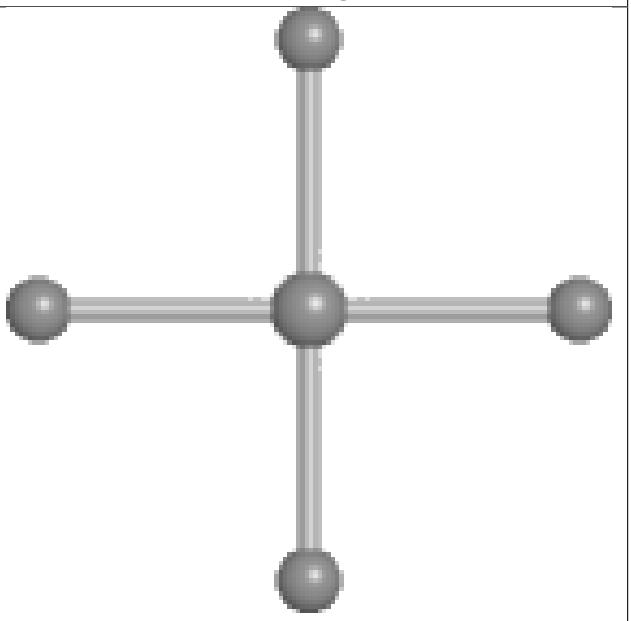
Bond lengths



Bond angles

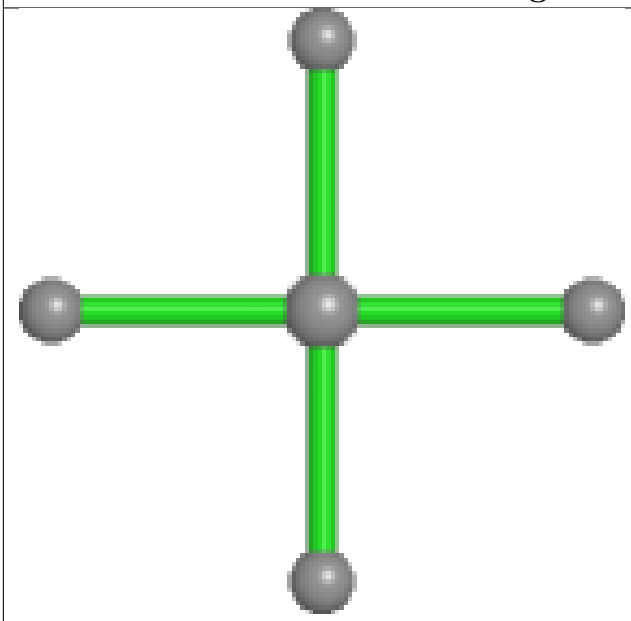


Torsions

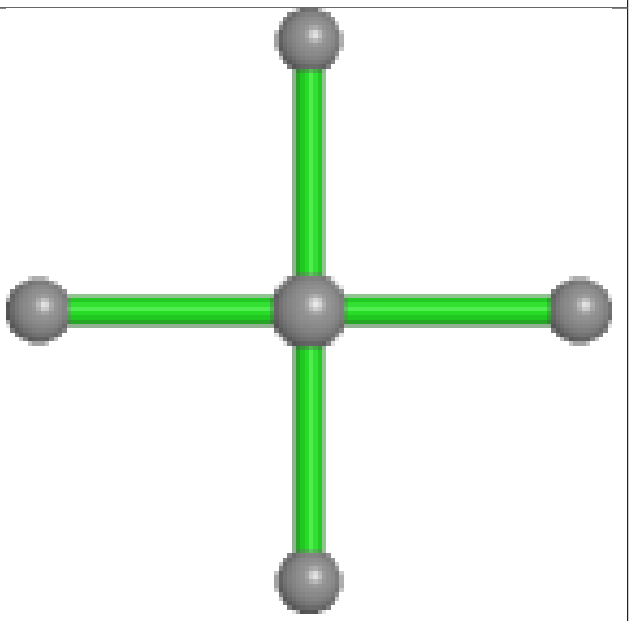


Rings

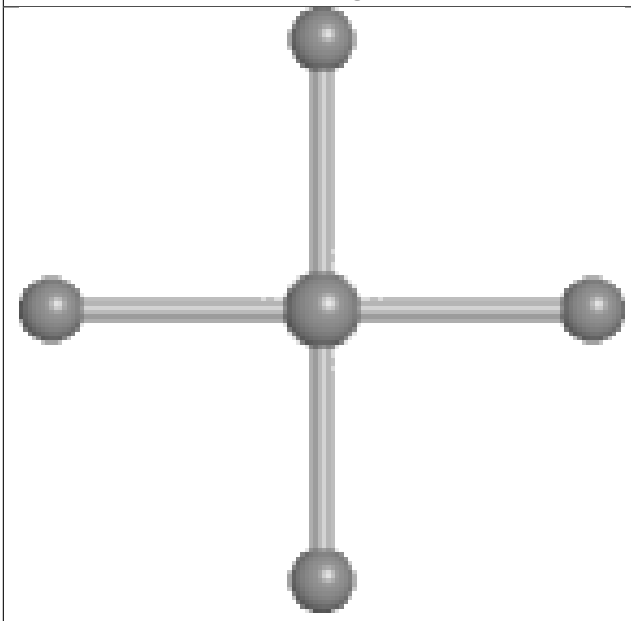
Ligand SO4 B 201



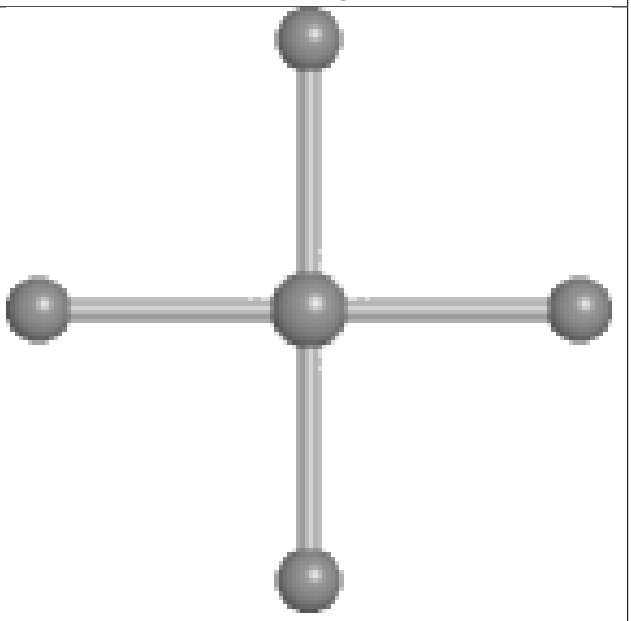
Bond lengths



Bond angles

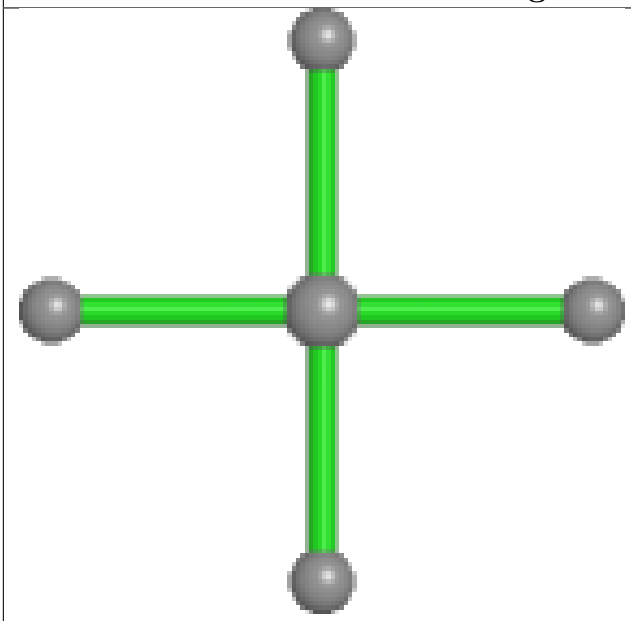


Torsions

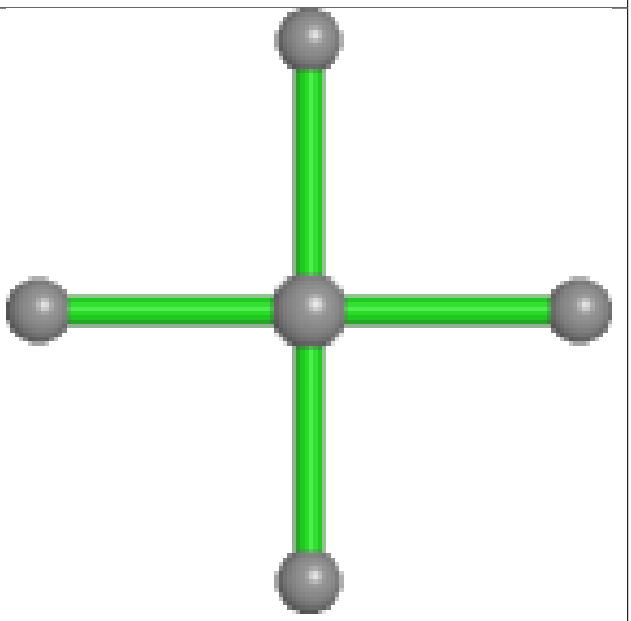


Rings

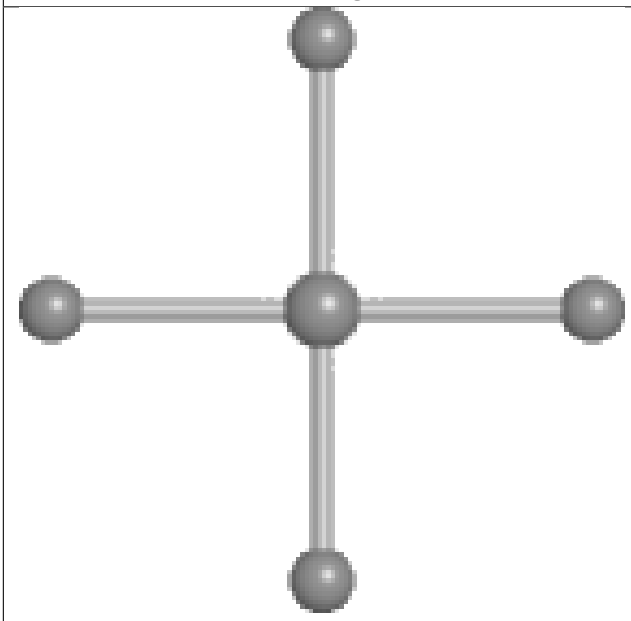
Ligand SO4 D 201



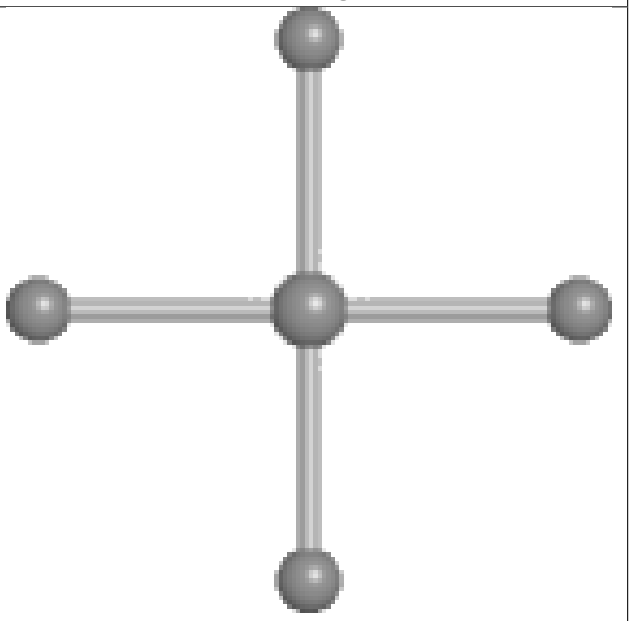
Bond lengths



Bond angles

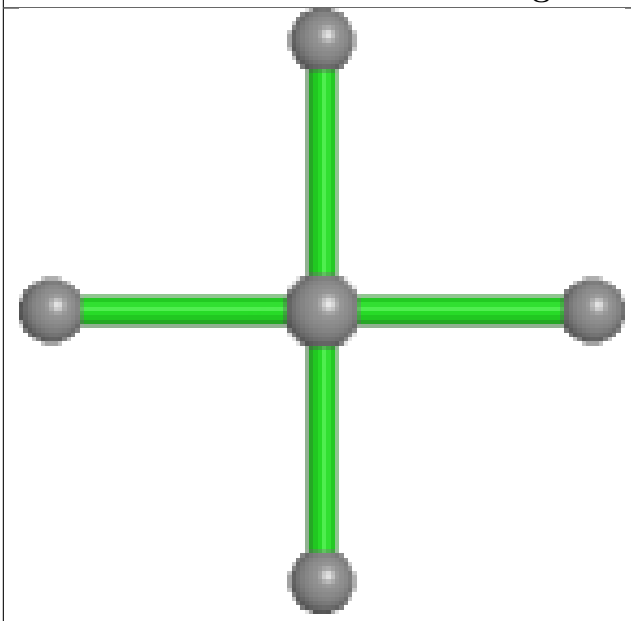


Torsions

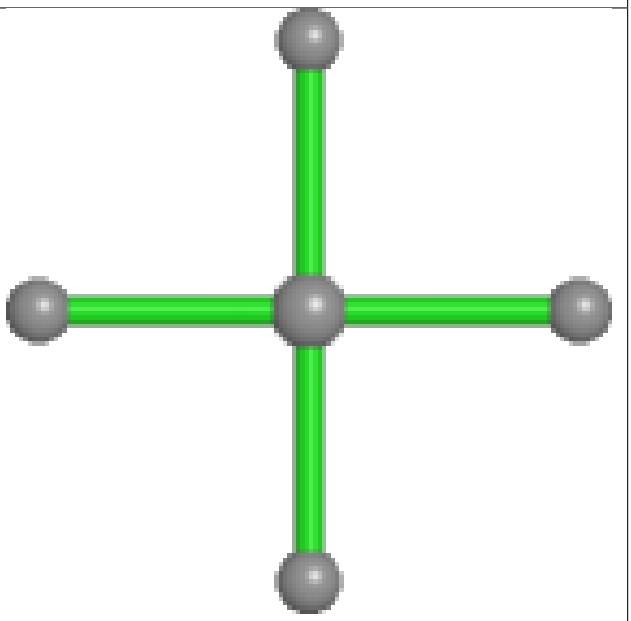


Rings

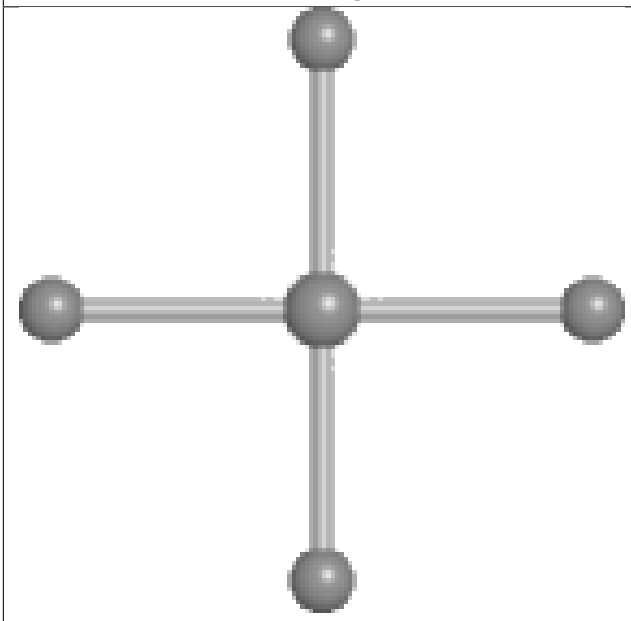
Ligand SO4 J 202



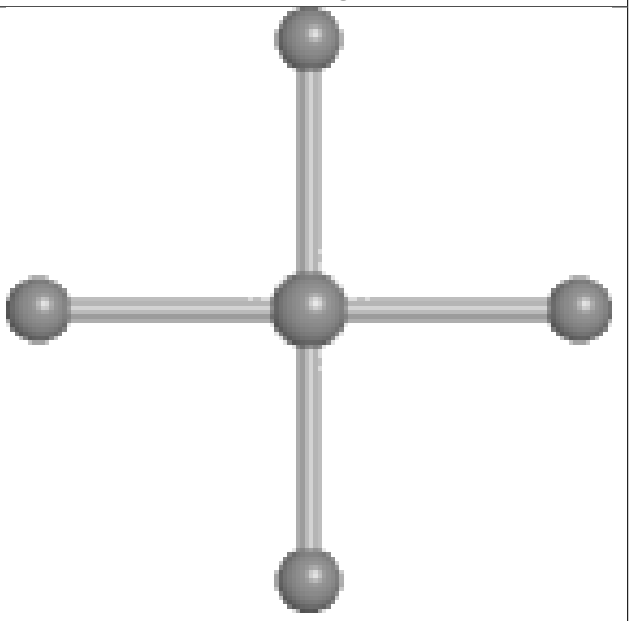
Bond lengths



Bond angles

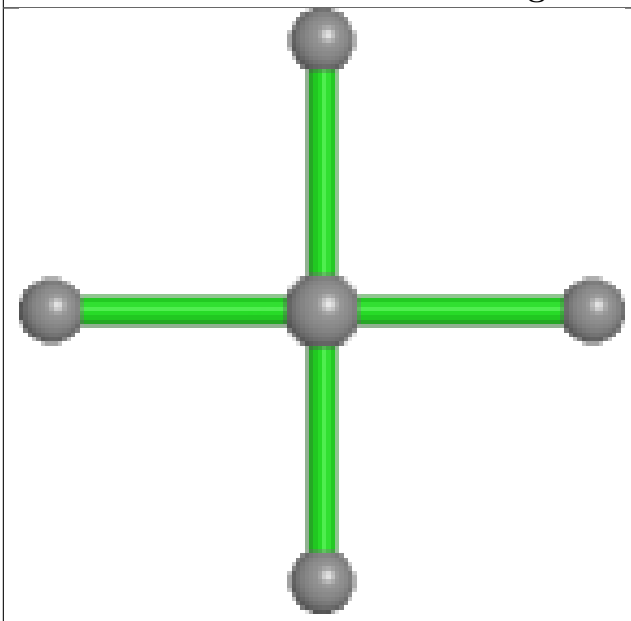


Torsions

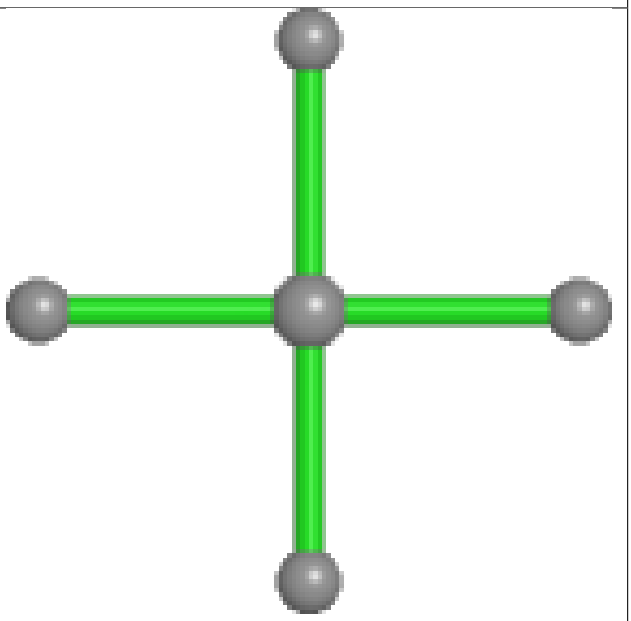


Rings

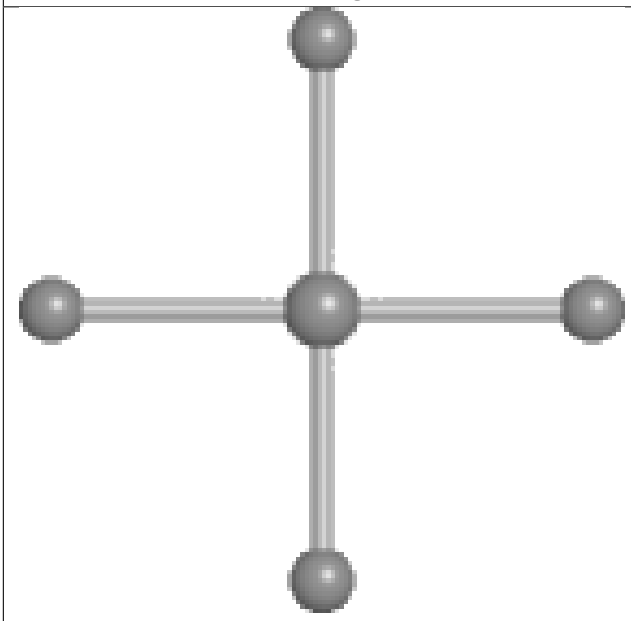
Ligand SO4 G 201



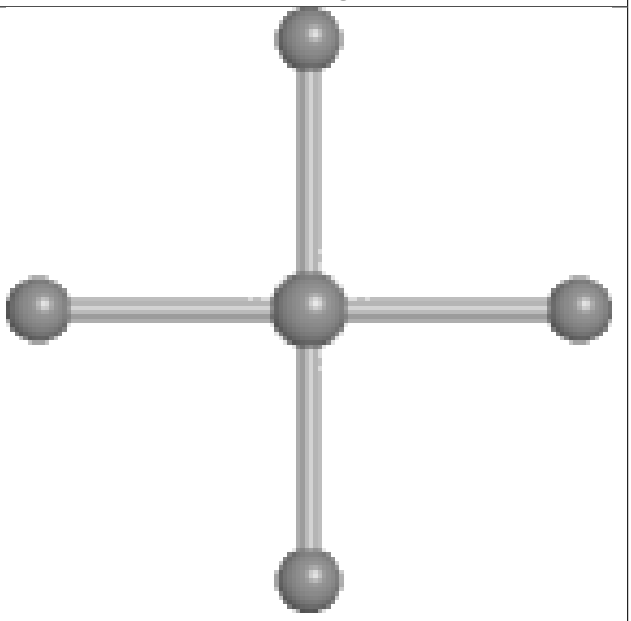
Bond lengths



Bond angles

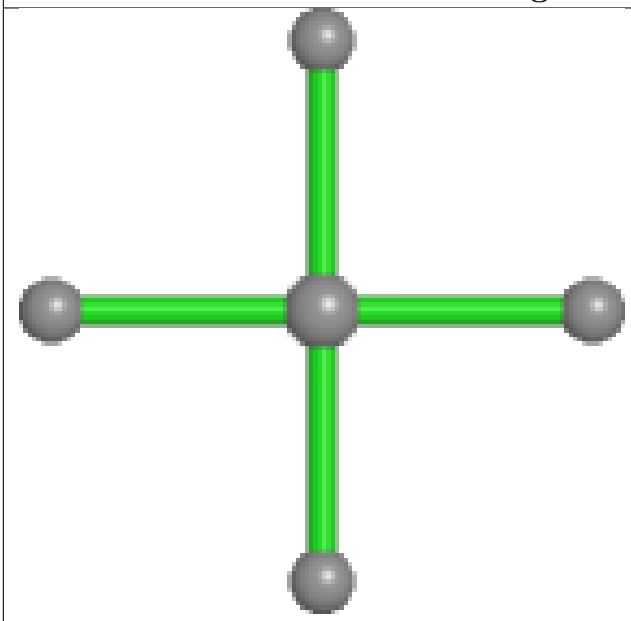


Torsions

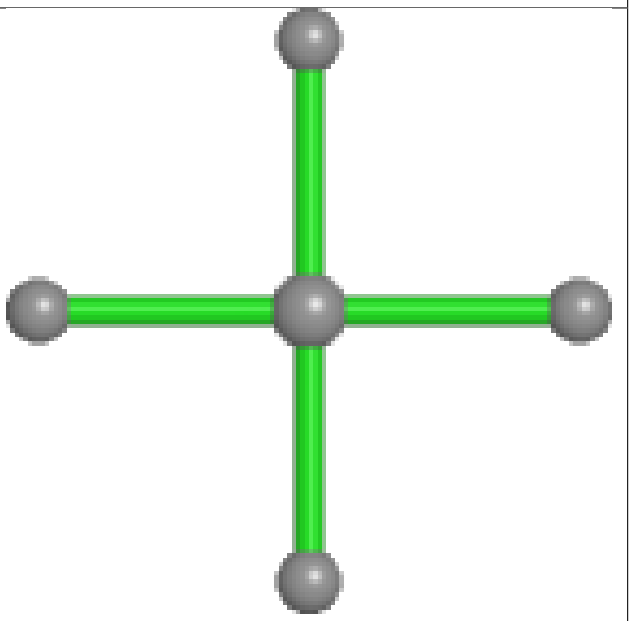


Rings

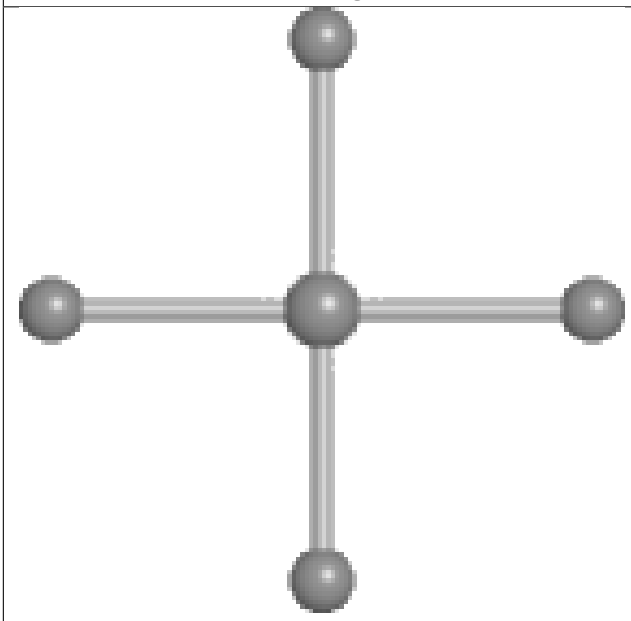
Ligand SO4 F 201



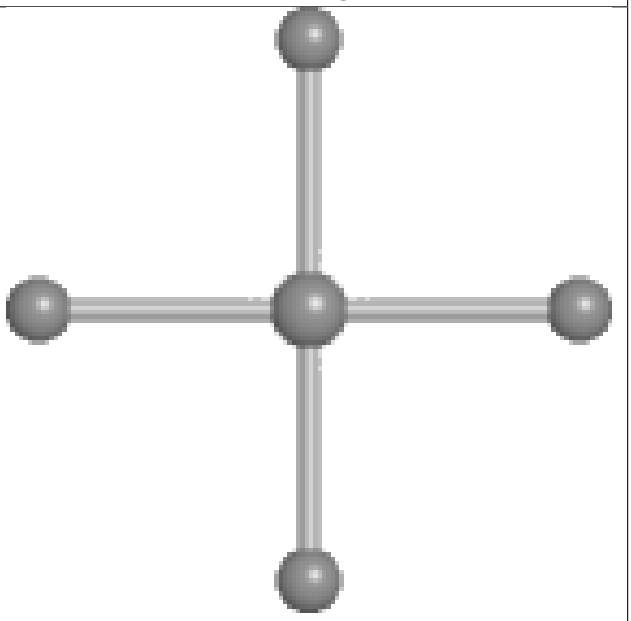
Bond lengths



Bond angles

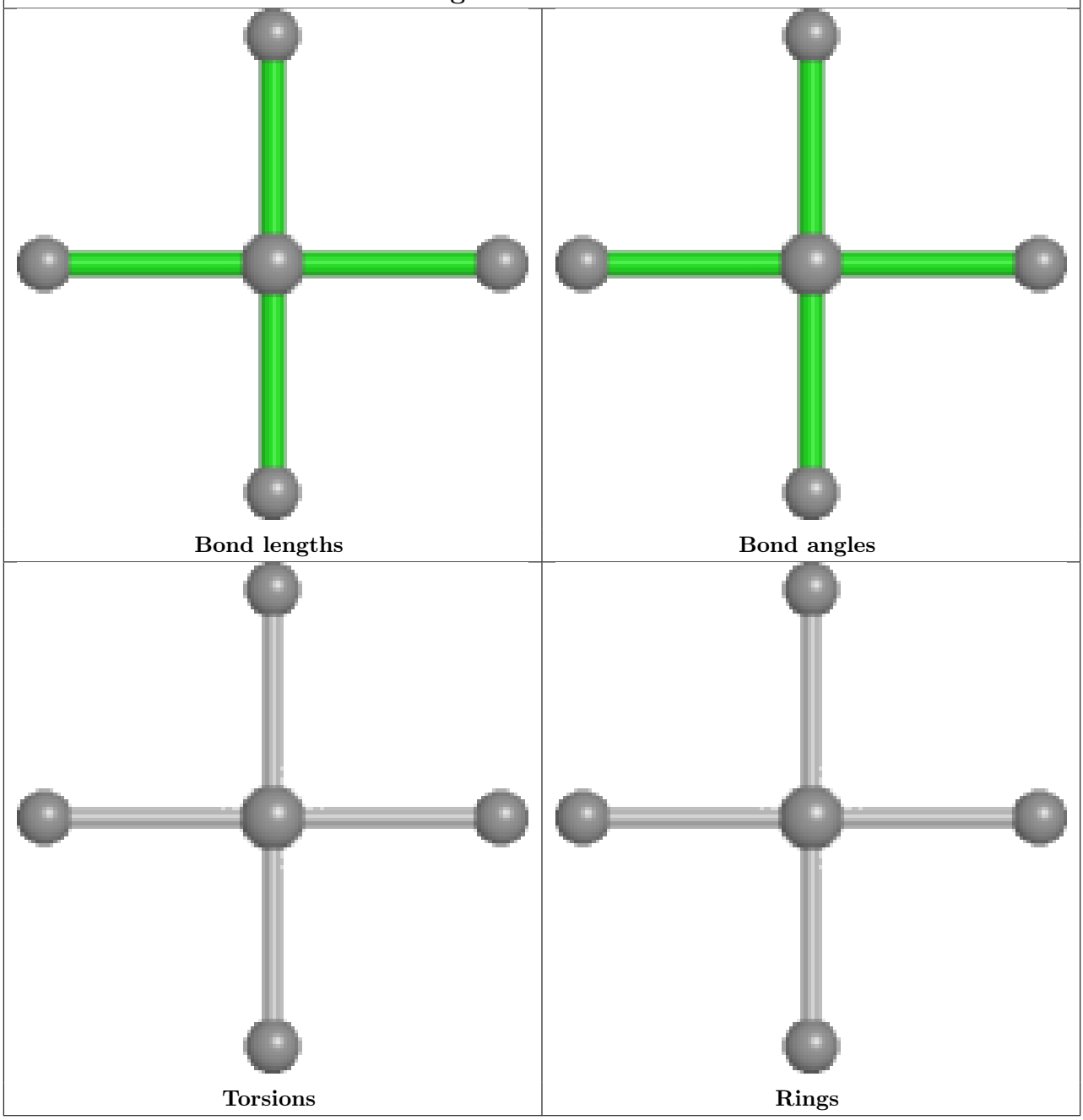


Torsions



Rings

Ligand SO4 J 201



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^{th} 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	116/131 (88%)	-0.60	0 100 100	12, 16, 26, 40	0
1	B	117/131 (89%)	-0.58	1 (0%) 84 84	12, 17, 26, 56	0
1	C	118/131 (90%)	-0.55	1 (0%) 86 86	13, 17, 36, 47	0
1	D	115/131 (87%)	-0.49	1 (0%) 84 84	13, 19, 31, 52	0
1	E	116/131 (88%)	-0.59	0 100 100	13, 18, 29, 47	0
1	F	115/131 (87%)	-0.53	1 (0%) 84 84	12, 19, 30, 52	0
1	G	117/131 (89%)	-0.56	0 100 100	13, 18, 29, 48	0
1	H	114/131 (87%)	-0.52	0 100 100	14, 20, 32, 65	0
1	I	113/131 (86%)	-0.55	0 100 100	13, 18, 27, 42	0
1	J	117/131 (89%)	-0.58	1 (0%) 84 84	13, 18, 29, 49	0
1	K	114/131 (87%)	-0.55	0 100 100	14, 19, 31, 58	0
1	L	115/131 (87%)	-0.49	1 (0%) 84 84	13, 17, 28, 52	0
All	All	1387/1572 (88%)	-0.55	6 (0%) 92 92	12, 18, 29, 65	0

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	116	VAL	4.4
1	B	118	ILE	3.5
1	L	116	VAL	3.3
1	J	118	ILE	2.7
1	F	116	VAL	2.4
1	C	116	VAL	2.1

6.2 Non-standard residues in protein, DNA, RNA chains

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)

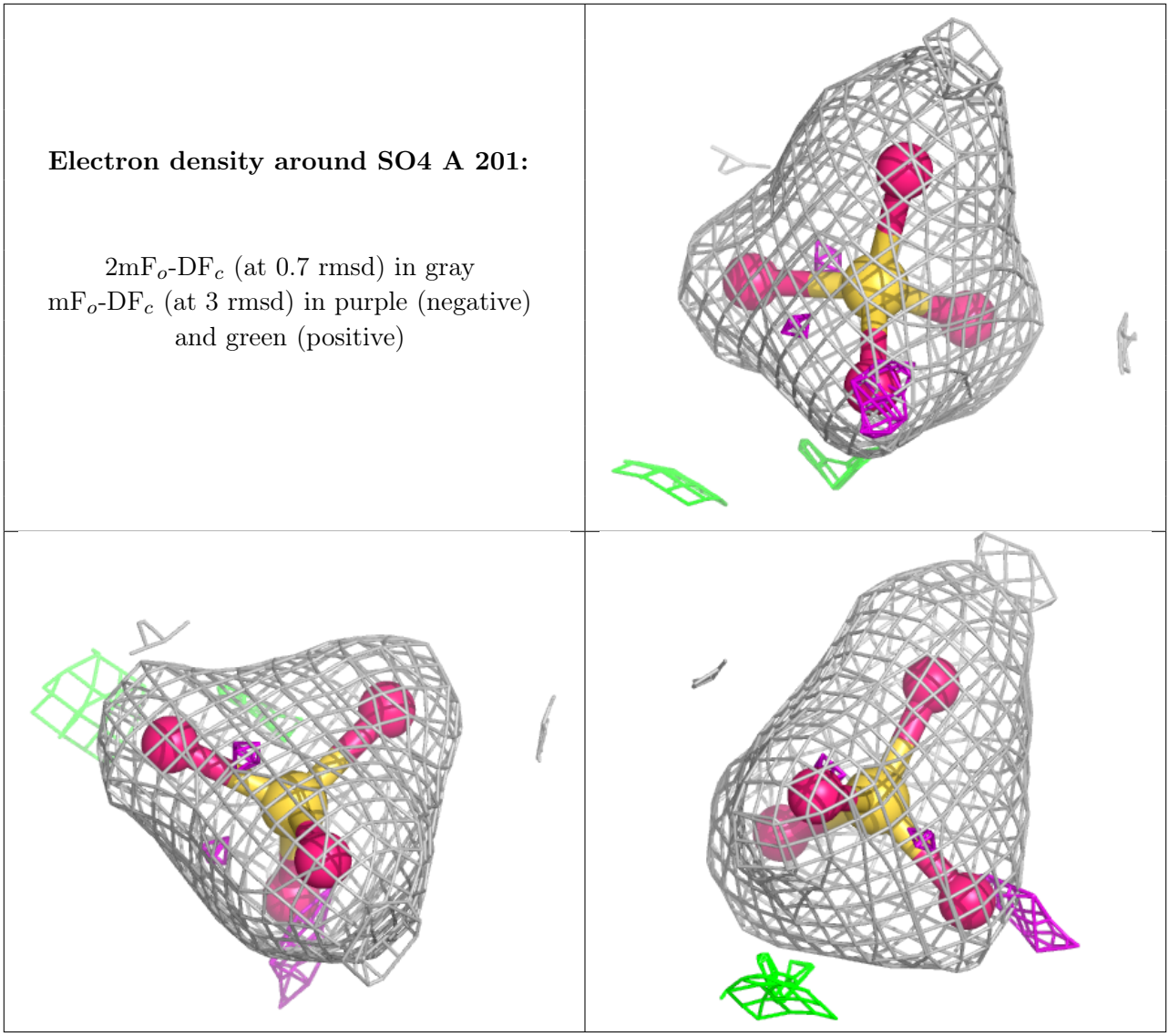
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' 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	201	5/5	0.97	0.10	23,27,30,31	0
2	SO4	B	201	5/5	0.97	0.11	23,26,31,32	0
2	SO4	J	202	5/5	0.98	0.11	27,28,34,35	0
2	SO4	F	201	5/5	0.99	0.06	18,19,20,22	0
2	SO4	G	201	5/5	0.99	0.06	21,21,22,22	0
2	SO4	G	202	5/5	0.99	0.12	27,28,32,36	0
2	SO4	J	201	5/5	0.99	0.07	22,22,22,22	0
2	SO4	D	201	5/5	0.99	0.07	18,18,19,20	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.

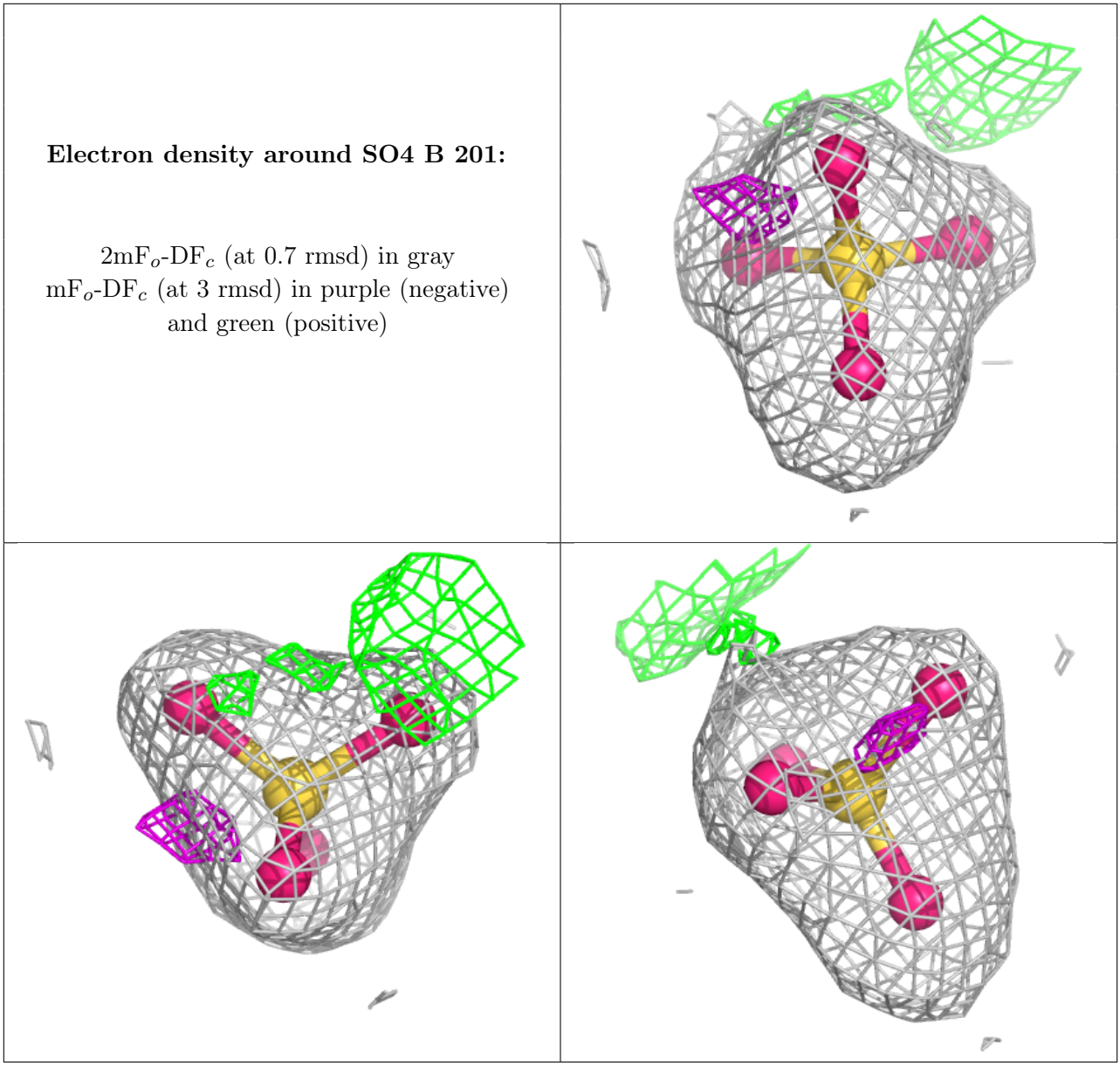
Electron density around SO4 A 201:

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



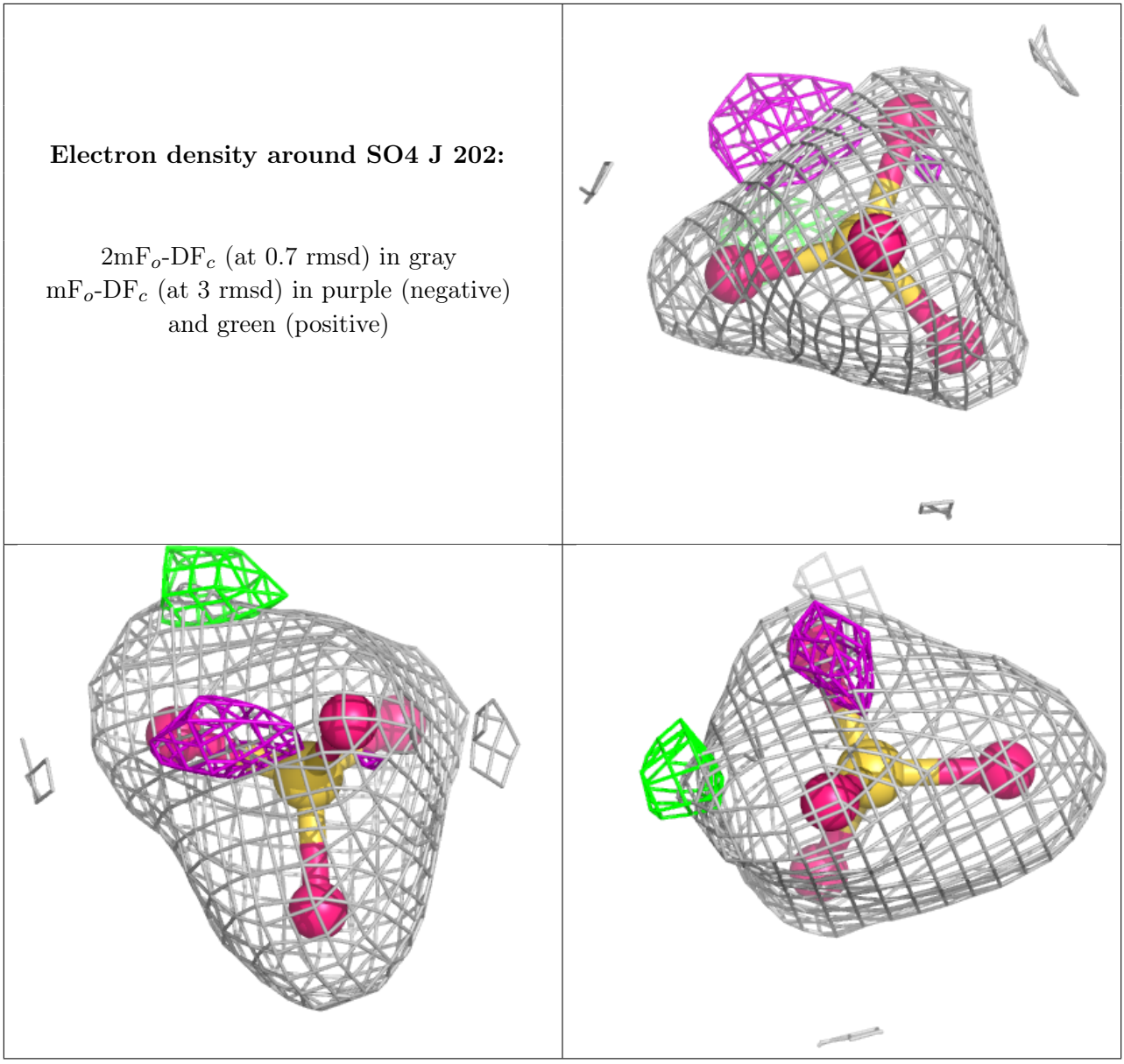
Electron density around SO4 B 201:

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



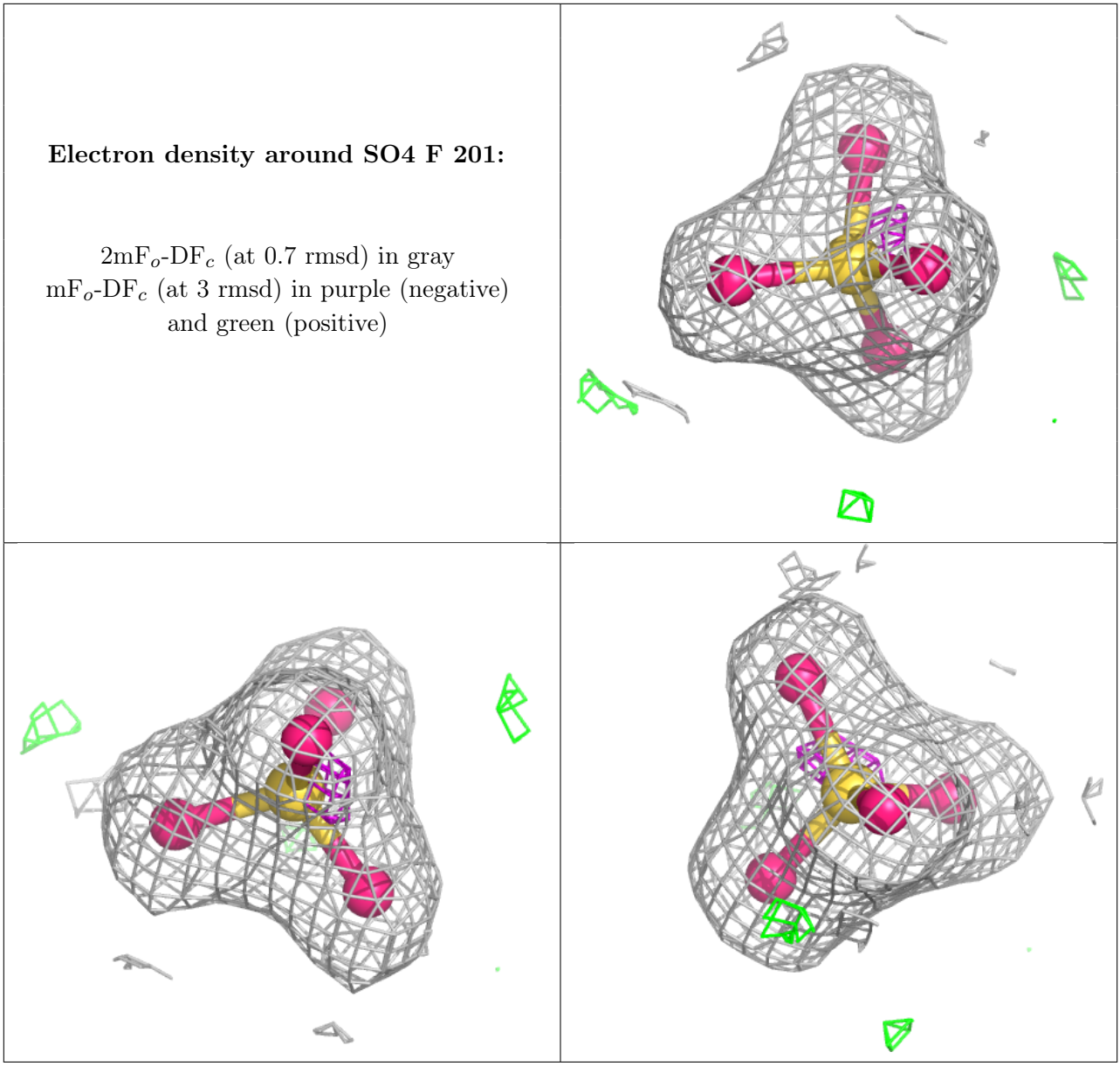
Electron density around SO4 J 202:

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



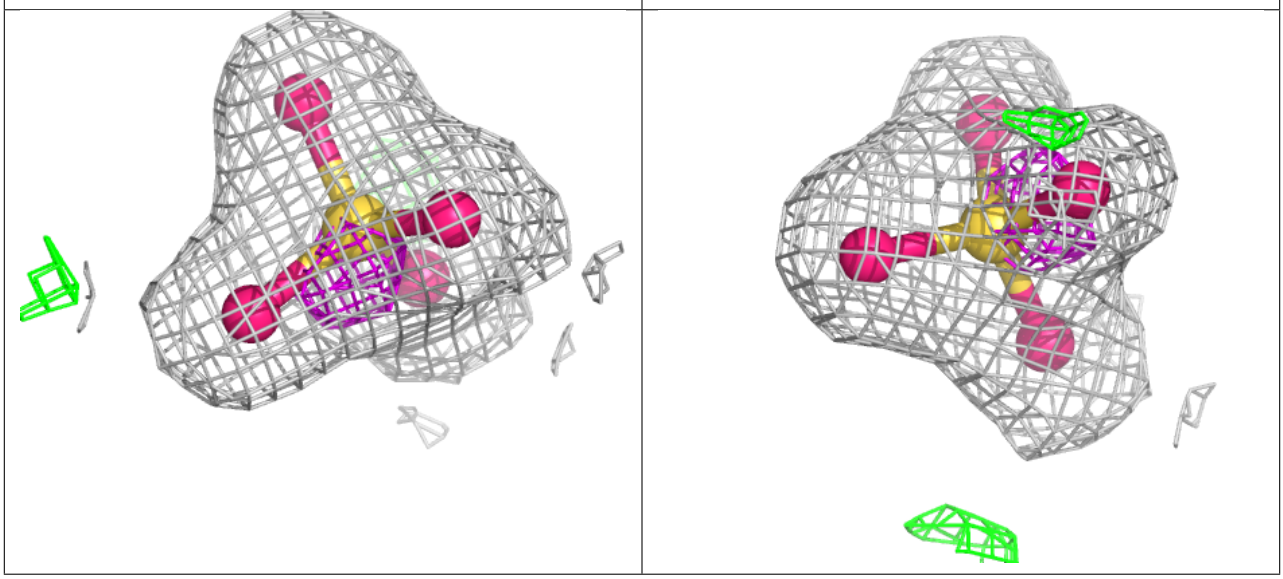
Electron density around SO4 F 201:

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



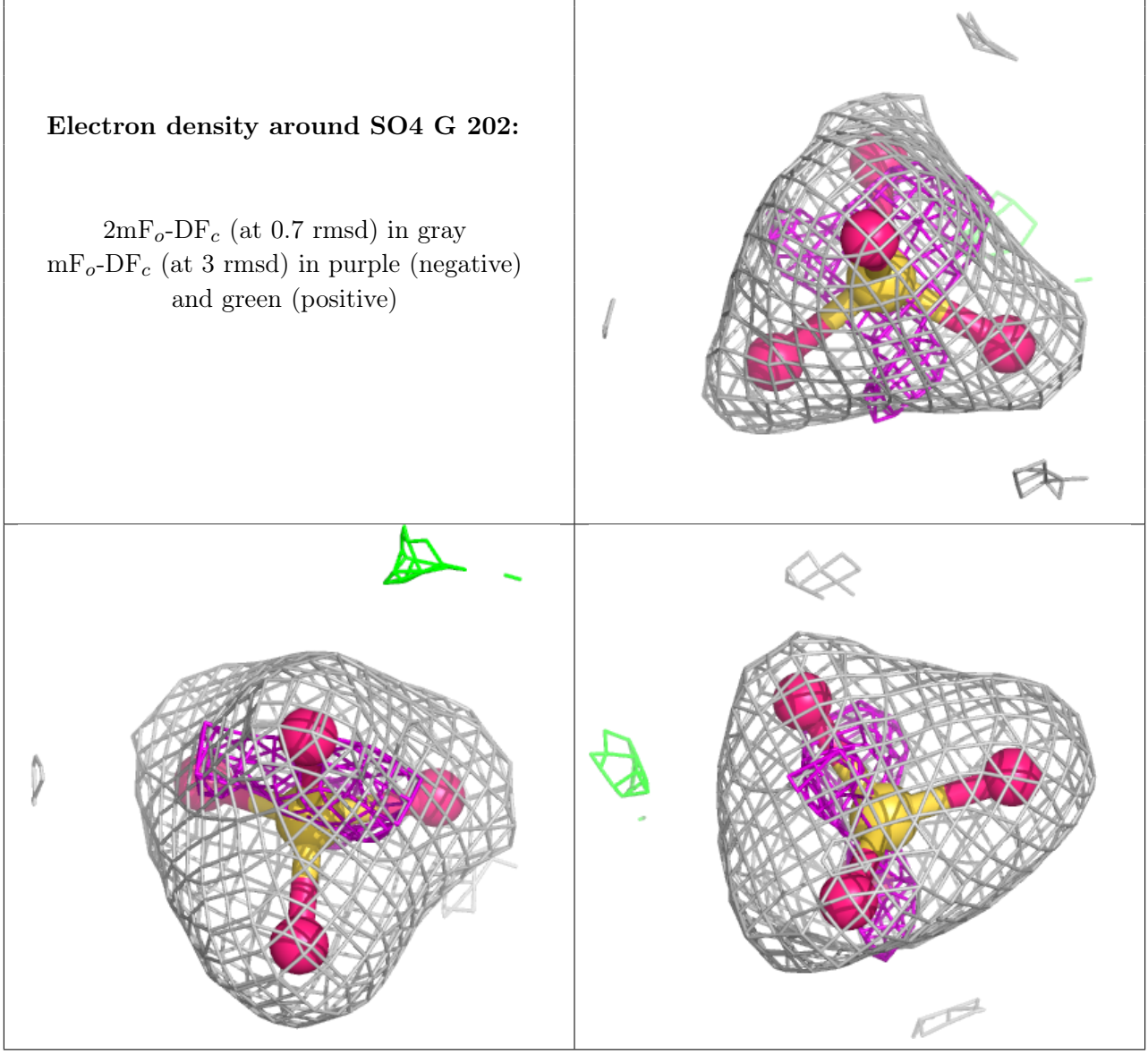
Electron density around SO4 G 201:

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



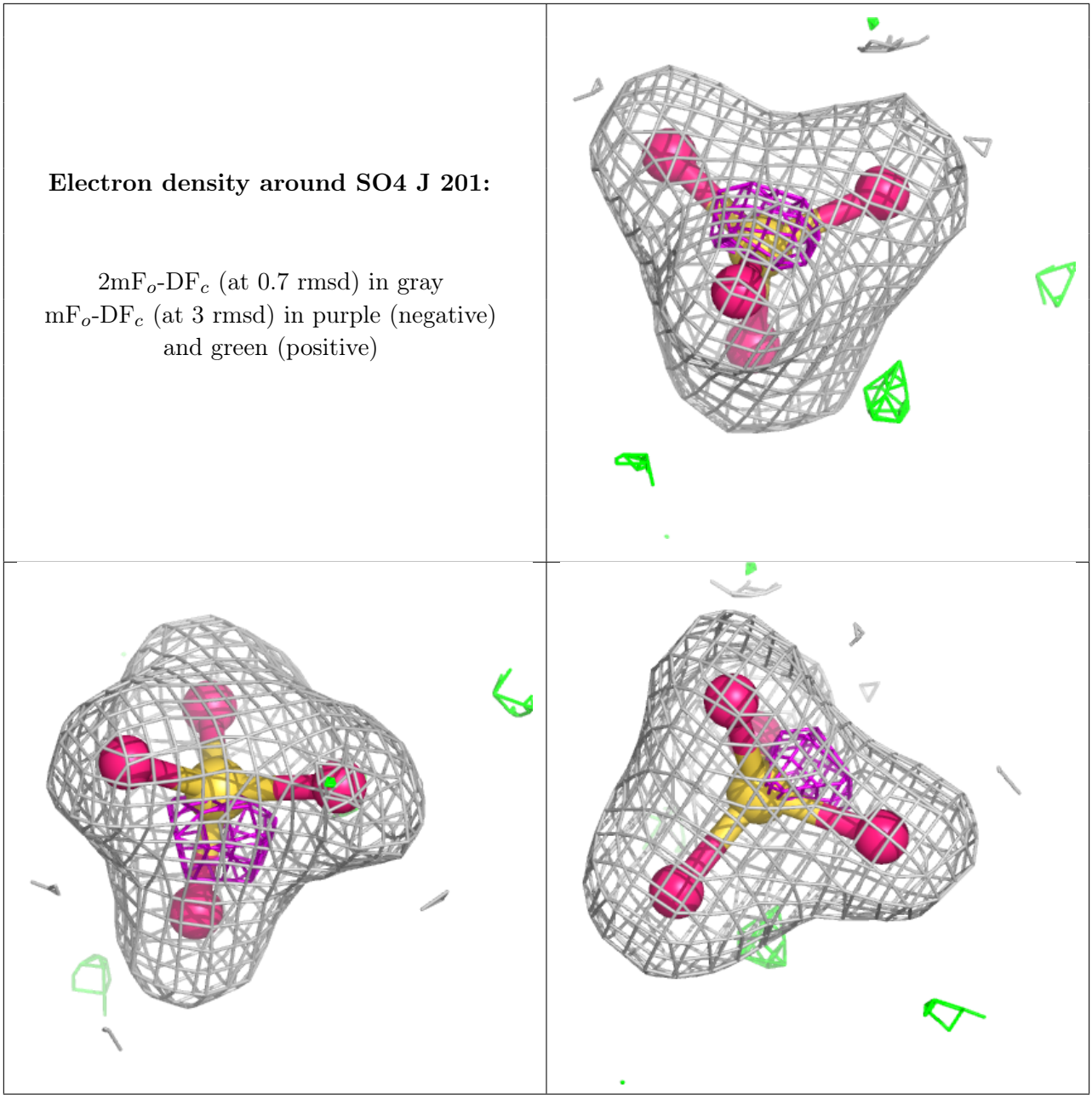
Electron density around SO4 G 202:

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



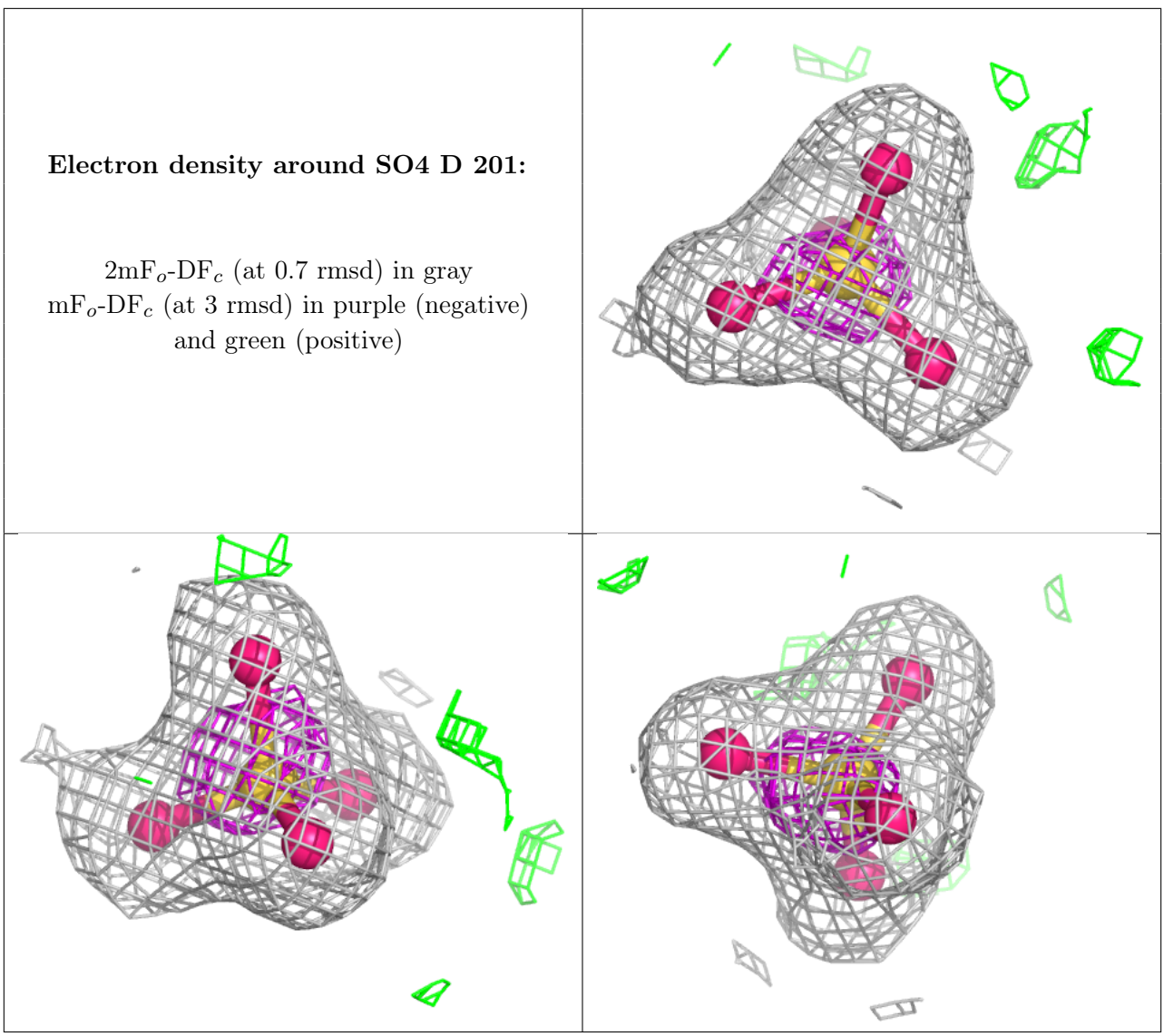
Electron density around SO4 J 201:

$2mF_o$-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 201:

$2mF_o$-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.