



# Full wwPDB X-ray Structure Validation Report ⓘ

Jun 1, 2026 – 10:18 AM EDT

PDB ID : 9OWZ / pdb\_00009owz  
Title : Structure of *A. thaliana* TBP bound to AG-P nicked DNA site  
Authors : Schumacher, M.A.  
Deposited on : 2025-06-02  
Resolution : 2.91 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>.

---

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0  
Mogul : 2022.3.0, CSD as543be (2022)  
Xtriage (Phenix) : 2.0  
EDS : 3.0  
Buster-report : wwPDB partial adaption of 1.1.7 (2018)  
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)  
CCP4 : 9.0.010 (Gargrove)  
Density-Fitness : 1.0.12  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.49

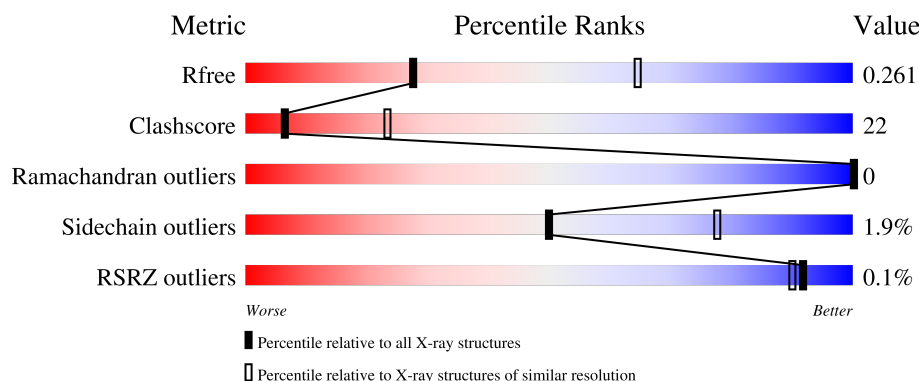
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.91 Å.

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



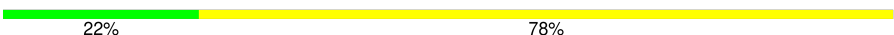
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	180053	2995 (2.94-2.90)
Clashscore	190562	3213 (2.94-2.90)
Ramachandran outliers	187476	3128 (2.94-2.90)
Sidechain outliers	187428	3130 (2.94-2.90)
RSRZ outliers	180081	2995 (2.94-2.90)

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	203	
1	B	203	
1	G	203	
1	J	203	
1	M	203	

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Length	Quality of chain
1	P	203	
2	C	13	
2	E	13	
2	H	13	
2	K	13	
2	N	13	
2	Q	13	
3	D	9	
3	F	9	
3	I	9	
3	L	9	
3	O	9	
3	R	9	

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 11349 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 TATA-box-binding protein 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	187	Total	C	N	O	S	0	0	0
			1425	924	240	253	8			
1	A	187	Total	C	N	O	S	0	0	0
			1427	932	236	251	8			
1	G	185	Total	C	N	O	S	0	0	0
			1405	914	234	249	8			
1	J	187	Total	C	N	O	S	0	0	0
			1458	950	245	255	8			
1	M	188	Total	C	N	O	S	0	0	0
			1453	948	243	254	8			
1	P	187	Total	C	N	O	S	0	0	0
			1427	929	242	248	8			

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-2	GLY	-	expression tag	UNP P28147
B	-1	SER	-	expression tag	UNP P28147
B	0	HIS	-	expression tag	UNP P28147
A	-2	GLY	-	expression tag	UNP P28147
A	-1	SER	-	expression tag	UNP P28147
A	0	HIS	-	expression tag	UNP P28147
G	-2	GLY	-	expression tag	UNP P28147
G	-1	SER	-	expression tag	UNP P28147
G	0	HIS	-	expression tag	UNP P28147
J	-2	GLY	-	expression tag	UNP P28147
J	-1	SER	-	expression tag	UNP P28147
J	0	HIS	-	expression tag	UNP P28147
M	-2	GLY	-	expression tag	UNP P28147
M	-1	SER	-	expression tag	UNP P28147
M	0	HIS	-	expression tag	UNP P28147
P	-2	GLY	-	expression tag	UNP P28147
P	-1	SER	-	expression tag	UNP P28147

*Continued on next page...*

*Continued from previous page...*

Chain	Residue	Modelled	Actual	Comment	Reference
P	0	HIS	-	expression tag	UNP P28147

- Molecule 2 is a DNA chain called DNA (5'-D(\*GP\*CP\*TP\*AP\*TP\*AP\*AP\*AP\*GP\*CP\*GP\*C)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	13	Total 265	C 127	N 53	O 73	P 12	0	0	0
2	E	13	Total 265	C 127	N 53	O 73	P 12	0	0	0
2	H	13	Total 265	C 127	N 53	O 73	P 12	0	0	0
2	K	13	Total 265	C 127	N 53	O 73	P 12	0	0	0
2	N	13	Total 265	C 127	N 53	O 73	P 12	0	0	0
2	Q	13	Total 265	C 127	N 53	O 73	P 12	0	0	0

- Molecule 3 is a DNA chain called DNA (5'-D(P\*TP\*TP\*TP\*TP\*AP\*TP\*AP\*GP\*C)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	D	9	Total 184	C 89	N 28	O 58	P 9	0	0	0
3	F	9	Total 184	C 89	N 28	O 58	P 9	0	0	0
3	I	9	Total 184	C 89	N 28	O 58	P 9	0	0	0
3	L	9	Total 184	C 89	N 28	O 58	P 9	0	0	0
3	O	9	Total 184	C 89	N 28	O 58	P 9	0	0	0
3	R	9	Total 184	C 89	N 28	O 58	P 9	0	0	0

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

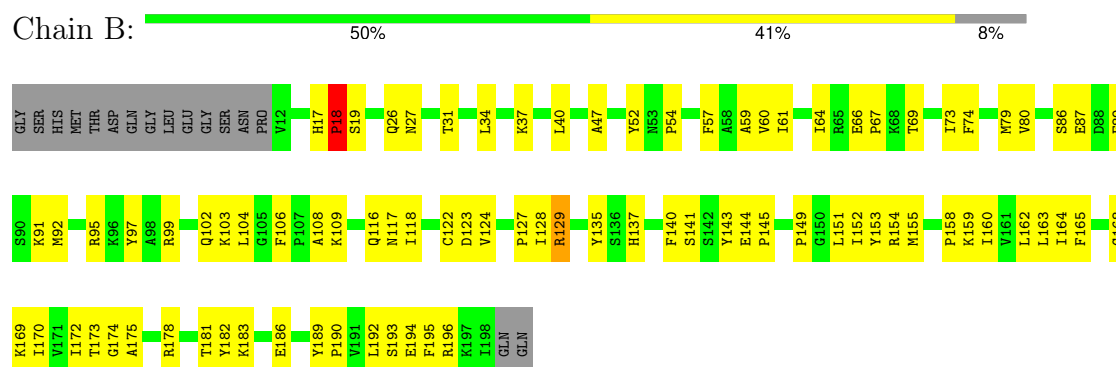


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	O	S	0	0
			5	4	1		
4	B	1	Total	O	S	0	0
			5	4	1		
4	A	1	Total	O	S	0	0
			5	4	1		
4	G	1	Total	O	S	0	0
			5	4	1		
4	G	1	Total	O	S	0	0
			5	4	1		
4	J	1	Total	O	S	0	0
			5	4	1		
4	J	1	Total	O	S	0	0
			5	4	1		
4	M	1	Total	O	S	0	0
			5	4	1		
4	M	1	Total	O	S	0	0
			5	4	1		
4	P	1	Total	O	S	0	0
			5	4	1		
4	P	1	Total	O	S	0	0
			5	4	1		
4	P	1	Total	O	S	0	0
			5	4	1		

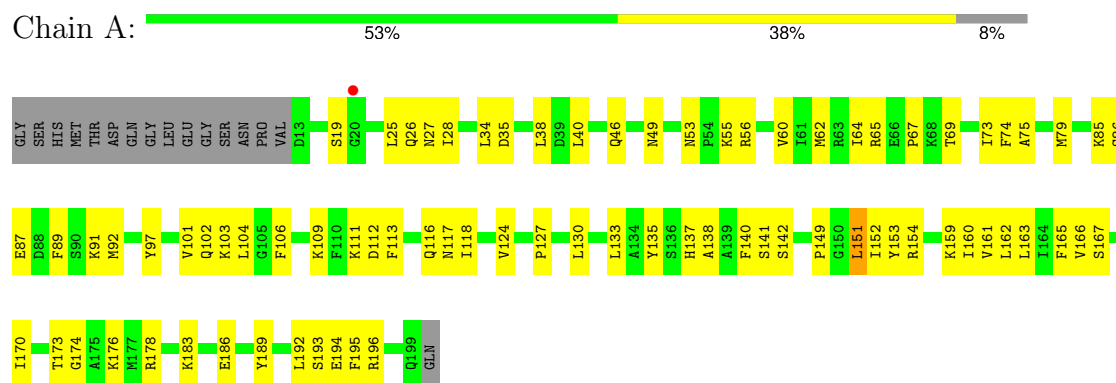
### 3 Residue-property plots [i](#)

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.

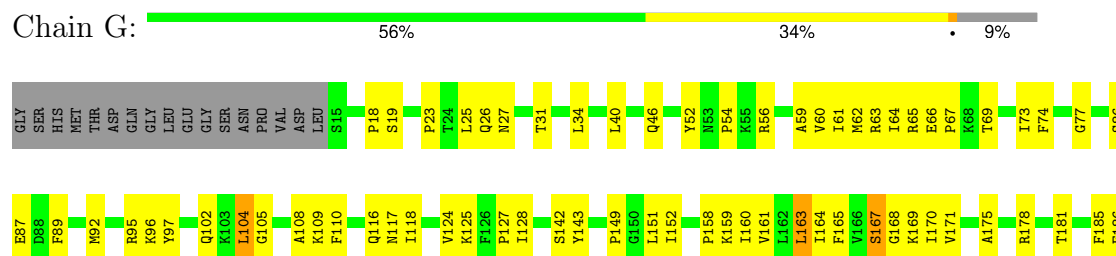
#### • Molecule 1: TATA-box-binding protein 1



#### • Molecule 1: TATA-box-binding protein 1



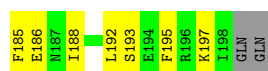
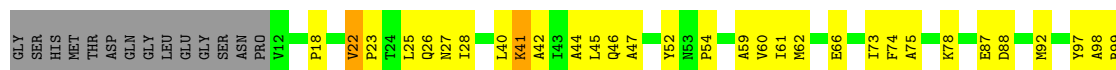
#### • Molecule 1: TATA-box-binding protein 1





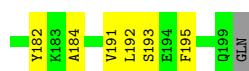
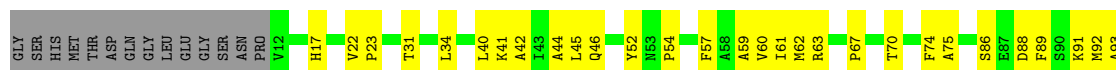
- Molecule 1: TATA-box-binding protein 1

Chain J: 53% 37% 8%



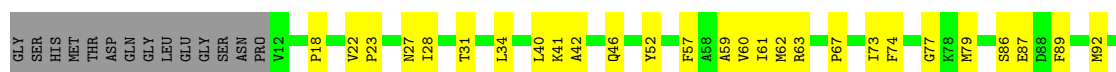
- Molecule 1: TATA-box-binding protein 1

Chain M: 57% 35% 7%



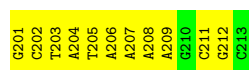
- Molecule 1: TATA-box-binding protein 1

Chain P: 55% 37% 8%



- Molecule 2: DNA (5'-D(\*GP\*CP\*TP\*AP\*TP\*AP\*AP\*AP\*AP\*GP\*CP\*GP\*C)-3')

Chain C: 15% 85%





- Molecule 2: DNA (5'-D(\*GP\*CP\*TP\*AP\*TP\*AP\*AP\*AP\*AP\*GP\*CP\*GP\*C)-3')

Chain E:  31% 69%



- Molecule 2: DNA (5'-D(\*GP\*CP\*TP\*AP\*TP\*AP\*AP\*AP\*AP\*GP\*CP\*GP\*C)-3')

Chain H:  38% 62%



- Molecule 2: DNA (5'-D(\*GP\*CP\*TP\*AP\*TP\*AP\*AP\*AP\*AP\*GP\*CP\*GP\*C)-3')

Chain K:  38% 62%



- Molecule 2: DNA (5'-D(\*GP\*CP\*TP\*AP\*TP\*AP\*AP\*AP\*AP\*GP\*CP\*GP\*C)-3')

Chain N:  38% 62%



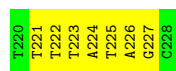
- Molecule 2: DNA (5'-D(\*GP\*CP\*TP\*AP\*TP\*AP\*AP\*AP\*AP\*GP\*CP\*GP\*C)-3')

Chain Q:  38% 62%



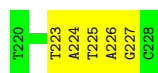
- Molecule 3: DNA (5'-D(P\*TP\*TP\*TP\*TP\*AP\*TP\*AP\*GP\*C)-3')

Chain D:  22% 78%




- Molecule 3: DNA (5'-D(P\*TP\*TP\*TP\*TP\*AP\*TP\*AP\*GP\*C)-3')

Chain F:  44% 56%



- Molecule 3: DNA (5'-D(P\*TP\*TP\*TP\*TP\*AP\*TP\*AP\*GP\*C)-3')

Chain I:  100%

T220  
T221  
T222  
T223  
A224  
T225  
A226  
G227  
C228

- Molecule 3: DNA (5'-D(P\*TP\*TP\*TP\*TP\*AP\*TP\*AP\*GP\*C)-3')

Chain L:  33% 67%

T220  
T221  
T222  
T223  
A224  
T225  
A226  
G227  
C228

- Molecule 3: DNA (5'-D(P\*TP\*TP\*TP\*TP\*AP\*TP\*AP\*GP\*C)-3')

Chain O:  22% 78%

T220  
T221  
T222  
T223  
A224  
T225  
A226  
G227  
C228

- Molecule 3: DNA (5'-D(P\*TP\*TP\*TP\*TP\*AP\*TP\*AP\*GP\*C)-3')

Chain R:  56% 44%

T220  
T221  
T222  
T223  
A224  
T225  
A226  
G227  
C228

## 4 Data and refinement statistics

Property	Value	Source
Space group	I 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	122.73Å 106.21Å 161.92Å 90.00° 101.10° 90.00°	Depositor
Resolution (Å)	79.66 – 2.91 79.66 – 2.91	Depositor EDS
% Data completeness (in resolution range)	97.0 (79.66-2.91) 96.2 (79.66-2.91)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.24 (at 2.91Å)	Xtriage
Refinement program	PHENIX 1.20.1_4487	Depositor
R, $R_{free}$	0.234 , 0.253 0.243 , 0.261	Depositor DCC
$R_{free}$ test set	2017 reflections (4.44%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	79.9	Xtriage
Anisotropy	0.396	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.38 , 112.7	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.428 for $-1/2^*h+1/2^*k+1/2^*l, 1/2^*h-1/2^*k+1/2^*l, h+k$ 0.419 for $-1/2^*h-1/2^*k+1/2^*l, -1/2^*h-1/2^*k-1/2^*l, h-k$	Xtriage
$F_o, F_c$ correlation	0.97	EDS
Total number of atoms	11349	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	83.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.64% 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

### 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	0.57	0/1456	0.71	0/1969
1	B	0.63	1/1454 (0.1%)	0.88	5/1967 (0.3%)
1	G	0.56	0/1433	0.74	0/1940
1	J	0.38	0/1488	0.62	0/2008
1	M	0.44	1/1483 (0.1%)	0.65	0/2003
1	P	0.43	0/1456	0.63	0/1970
2	C	0.49	0/298	0.66	0/458
2	E	0.45	0/298	0.69	0/458
2	H	0.48	0/298	0.62	0/458
2	K	0.33	0/298	0.56	0/458
2	N	0.38	0/298	0.60	0/458
2	Q	0.43	0/298	0.59	0/458
3	D	0.62	0/204	0.68	0/311
3	F	0.66	0/204	0.71	0/311
3	I	0.65	0/204	0.69	0/311
3	L	0.48	0/204	0.68	0/311
3	O	0.55	0/204	0.68	0/311
3	R	0.52	0/204	0.66	0/311
All	All	0.51	2/11782 (0.0%)	0.69	5/16471 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	M	31	THR	C-N	6.25	1.39	1.33
1	B	129	ARG	CB-CG	5.14	1.67	1.52

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	18	PRO	N-CD-CG	-10.60	87.30	103.20
1	B	18	PRO	CA-CB-CG	-10.35	84.84	104.50
1	B	18	PRO	CB-CG-CD	-6.00	86.91	106.10
1	B	18	PRO	N-CA-CB	-5.99	96.96	103.25
1	B	18	PRO	N-CA-C	5.13	123.05	112.47

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	129	ARG	Sidechain

## 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	1427	0	1461	63	0
1	B	1425	0	1436	91	0
1	G	1405	0	1422	62	0
1	J	1458	0	1514	68	0
1	M	1453	0	1501	63	0
1	P	1427	0	1462	74	0
2	C	265	0	147	17	0
2	E	265	0	147	12	0
2	H	265	0	147	12	0
2	K	265	0	147	10	0
2	N	265	0	147	6	0
2	Q	265	0	147	6	0
3	D	184	0	105	9	0
3	F	184	0	105	7	0
3	I	184	0	105	10	0
3	L	184	0	105	5	0

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	O	184	0	105	6	0
3	R	184	0	105	4	0
4	A	5	0	0	1	0
4	B	10	0	0	1	0
4	G	10	0	0	1	0
4	J	10	0	0	1	0
4	M	10	0	0	0	0
4	P	15	0	0	0	0
All	All	11349	0	10308	455	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 22.

All (455) 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:P:130:LEU:HB3	1:P:143:TYR:HE2	1.24	1.02
1:J:130:LEU:HB3	1:J:143:TYR:HE1	1.25	0.97
1:A:162:LEU:HD21	1:A:170:ILE:HD11	1.47	0.96
1:G:63:ARG:NH1	3:I:222:DT:OP1	2.02	0.93
1:J:153:TYR:HE1	1:J:155:MET:HB2	1.37	0.88
1:P:155:MET:HE1	1:P:184:ALA:HA	1.54	0.87
1:A:152:ILE:HD12	2:E:205:DT:H5''	1.57	0.85
1:J:127:PRO:HB2	1:J:197:LYS:HG2	1.61	0.83
1:P:41:LYS:H	1:P:41:LYS:HD3	1.43	0.83
1:J:78:LYS:HE2	2:K:209:DA:H5''	1.60	0.82
1:A:92:MET:HE1	1:P:92:MET:HE1	1.60	0.82
1:J:109:LYS:HE3	1:J:110:PHE:H	1.43	0.82
1:M:143:TYR:HE1	1:M:145:PRO:HG3	1.43	0.82
1:M:40:LEU:HD13	1:M:60:VAL:HG23	1.59	0.81
1:G:117:ASN:HD22	1:G:118:ILE:N	1.78	0.81
1:M:143:TYR:CE1	1:M:145:PRO:HG3	2.16	0.80
1:P:67:PRO:HD3	1:P:92:MET:HE2	1.65	0.79
1:P:130:LEU:HB3	1:P:143:TYR:CE2	2.15	0.77
1:P:109:LYS:HA	1:P:109:LYS:HE2	1.66	0.77
1:B:170:ILE:HD13	1:B:172:ILE:HD11	1.67	0.77
2:N:201:DG:N2	3:O:228:DC:O2	2.19	0.75
1:B:124:VAL:HG11	1:B:192:LEU:HD13	1.67	0.74
1:B:137:HIS:HB3	1:B:140:PHE:HB2	1.67	0.74
1:G:65:ARG:N	4:G:301:SO4:O3	2.22	0.73
1:B:135:TYR:OH	1:A:183:LYS:HD2	1.89	0.73

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:P:164:ILE:HD12	1:P:192:LEU:HD21	1.70	0.73
1:J:130:LEU:HB3	1:J:143:TYR:CE1	2.17	0.72
1:G:92:MET:HE1	1:J:92:MET:HE1	1.72	0.72
1:P:62:MET:HE1	1:P:100:ILE:HG21	1.72	0.72
1:A:130:LEU:HB3	1:A:151:LEU:HD13	1.73	0.71
1:A:161:VAL:HG11	2:E:205:DT:H4'	1.73	0.71
1:G:66:GLU:OE1	1:J:99:ARG:NH2	2.24	0.70
1:P:40:LEU:HG	1:P:60:VAL:HG23	1.73	0.70
1:M:131:GLU:OE1	1:M:131:GLU:N	2.25	0.70
1:G:125:LYS:HD2	1:G:125:LYS:O	1.91	0.69
1:G:163:LEU:HB2	1:G:171:VAL:HB	1.75	0.69
2:C:204:DA:H2''	2:C:205:DT:H5'	1.72	0.69
1:J:22:VAL:HG23	1:J:182:TYR:HE2	1.58	0.69
1:P:31:THR:HG23	1:P:116:GLN:HG3	1.75	0.69
1:P:148:PHE:HD1	1:P:149:PRO:HD2	1.57	0.69
1:M:114:LYS:HG2	1:M:116:GLN:HG2	1.74	0.68
1:A:141:SER:HA	1:A:153:TYR:HA	1.76	0.68
1:M:67:PRO:HD3	1:M:92:MET:HE3	1.75	0.68
1:B:19:SER:OG	1:B:186:GLU:OE2	2.09	0.68
1:G:86:SER:HB3	1:G:89:PHE:CD2	2.29	0.67
2:Q:210:DG:H2''	2:Q:211:DC:C5	2.30	0.67
1:G:40:LEU:HD22	1:G:60:VAL:HG23	1.77	0.67
1:A:87:GLU:OE1	1:A:178:ARG:NH1	2.26	0.66
1:B:118:ILE:HB	1:B:175:ALA:HB3	1.77	0.66
1:P:63:ARG:NH2	3:R:222:DT:OP1	2.28	0.66
1:J:117:ASN:HD21	2:K:206:DA:H1'	1.60	0.66
1:A:174:GLY:HA2	2:E:206:DA:H2''	1.78	0.65
1:J:59:ALA:HB2	1:J:74:PHE:CE1	2.32	0.65
1:M:86:SER:HB3	1:M:89:PHE:CD2	2.32	0.65
1:P:169:LYS:HD2	3:R:225:DT:H5''	1.79	0.65
1:P:102:GLN:HG3	1:P:108:ALA:HB3	1.77	0.65
2:E:203:DT:O2	3:F:227:DG:N2	2.30	0.64
1:A:154:ARG:HG2	1:A:161:VAL:HG22	1.79	0.64
2:N:210:DG:H2''	2:N:211:DC:C6	2.32	0.64
1:J:41:LYS:O	1:J:45:LEU:HG	1.98	0.64
1:J:18:PRO:HD2	1:J:186:GLU:HG3	1.78	0.64
1:A:162:LEU:CD2	1:A:170:ILE:HD11	2.25	0.64
2:H:201:DG:H2'	2:H:202:DC:C6	2.32	0.64
1:J:61:ILE:HD13	3:L:221:DT:H5''	1.80	0.64
1:G:158:PRO:HG2	1:G:160:ILE:HG12	1.80	0.63
1:B:143:TYR:HB2	1:B:151:LEU:HD12	1.79	0.63

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:34:LEU:HG	1:G:108:ALA:HB1	1.80	0.63
1:J:141:SER:HB3	1:J:151:LEU:HD21	1.80	0.63
2:C:202:DC:H2'	2:C:203:DT:C6	2.34	0.63
1:G:46:GLN:HB3	1:G:104:LEU:HD11	1.81	0.63
1:P:86:SER:HB3	1:P:89:PHE:CD2	2.34	0.63
1:M:100:ILE:O	1:M:104:LEU:HD23	1.99	0.62
1:J:44:ALA:HB2	1:J:52:TYR:HB3	1.81	0.62
1:P:143:TYR:HE1	1:P:145:PRO:HB3	1.64	0.62
1:B:92:MET:SD	1:M:92:MET:HE2	2.40	0.61
1:B:143:TYR:CE1	1:B:145:PRO:HD3	2.35	0.61
1:J:62:MET:HE1	1:J:100:ILE:HG21	1.81	0.61
1:G:74:PHE:CE2	2:H:209:DA:H1'	2.35	0.61
1:P:73:ILE:HG12	1:P:79:MET:HG3	1.81	0.61
1:P:158:PRO:HB2	1:P:160:ILE:HD13	1.81	0.61
2:N:212:DG:H2''	2:N:213:DC:O4'	2.00	0.61
2:K:211:DC:H2'	2:K:212:DG:C8	2.36	0.61
1:M:44:ALA:HB2	1:M:52:TYR:HB3	1.83	0.60
1:A:142:SER:HB2	1:A:152:ILE:HG23	1.84	0.59
1:G:128:ILE:HD11	1:G:168:GLY:HA2	1.83	0.59
1:G:86:SER:HB3	1:G:89:PHE:HD2	1.66	0.59
1:M:63:ARG:HG2	1:M:70:THR:CG2	2.32	0.59
1:B:122:CYS:SG	1:B:123:ASP:N	2.76	0.59
1:B:189:TYR:HA	1:B:192:LEU:HD12	1.84	0.58
1:B:17:HIS:HA	1:B:186:GLU:HG2	1.85	0.58
1:P:141:SER:HB3	1:P:151:LEU:HD21	1.85	0.58
1:M:123:ASP:OD1	1:M:169:LYS:HG3	2.03	0.58
1:M:40:LEU:CD1	1:M:60:VAL:HG23	2.33	0.58
2:E:201:DG:H2'	2:E:202:DC:C6	2.39	0.57
1:G:175:ALA:HB2	1:G:181:THR:HG23	1.85	0.57
1:J:103:LYS:NZ	4:J:301:SO4:O1	2.24	0.57
1:A:64:ILE:HG23	1:A:97:TYR:CE1	2.39	0.57
1:A:193:SER:O	1:A:196:ARG:HG3	2.04	0.57
1:G:124:VAL:HG13	1:G:170:ILE:HD11	1.87	0.57
1:M:155:MET:HE1	1:M:184:ALA:HA	1.87	0.57
1:P:46:GLN:HB2	1:P:104:LEU:HD11	1.86	0.57
1:B:40:LEU:HD22	1:B:60:VAL:HG23	1.87	0.57
1:P:164:ILE:HD13	1:P:170:ILE:HD12	1.87	0.56
2:E:201:DG:H5'	2:E:201:DG:H8	1.70	0.56
1:A:194:GLU:HG2	1:A:195:PHE:CD2	2.41	0.56
1:M:164:ILE:HG21	1:M:192:LEU:HD11	1.87	0.56
1:B:154:ARG:NH2	2:C:205:DT:H5''	2.21	0.56

*Continued on next page...*



*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:183:LYS:O	1:A:186:GLU:N	2.39	0.56
2:C:208:DA:N7	2:C:209:DA:N6	2.55	0.55
1:A:34:LEU:HA	1:A:109:LYS:O	2.05	0.55
1:G:143:TYR:HB2	1:G:151:LEU:CD1	2.36	0.55
1:B:123:ASP:HB2	1:B:169:LYS:HD3	1.87	0.55
1:G:151:LEU:HD23	1:G:164:ILE:HD13	1.88	0.55
2:H:201:DG:H2'	2:H:202:DC:H6	1.72	0.55
1:P:87:GLU:OE2	1:P:178:ARG:HD2	2.05	0.55
1:M:153:TYR:HE1	1:M:155:MET:HG3	1.72	0.55
1:G:152:ILE:HD12	2:H:205:DT:H5'	1.88	0.55
1:J:143:TYR:HE2	1:J:145:PRO:HB3	1.71	0.55
1:M:63:ARG:HG2	1:M:70:THR:HB	1.89	0.55
1:B:64:ILE:HG23	1:B:97:TYR:CE1	2.42	0.55
1:A:117:ASN:C	1:A:118:ILE:HD12	2.32	0.55
1:P:23:PRO:HD2	1:P:182:TYR:HD1	1.72	0.55
1:B:194:GLU:HB3	1:B:195:PHE:HD1	1.72	0.55
1:J:40:LEU:HD13	1:J:60:VAL:HG23	1.89	0.55
1:M:96:LYS:O	1:M:100:ILE:HG13	2.07	0.54
1:A:74:PHE:CZ	2:E:209:DA:H1'	2.42	0.54
1:G:62:MET:HE2	1:G:73:ILE:HD12	1.89	0.54
1:G:161:VAL:HG11	2:H:205:DT:H4'	1.89	0.54
1:A:62:MET:HE3	1:A:73:ILE:HD13	1.90	0.54
1:M:41:LYS:O	1:M:45:LEU:HD12	2.07	0.54
1:B:143:TYR:CZ	1:B:145:PRO:HD3	2.43	0.54
1:J:154:ARG:NH2	2:K:206:DA:OP1	2.37	0.54
1:M:114:LYS:HG2	1:M:116:GLN:CG	2.37	0.54
1:P:62:MET:HE2	1:P:97:TYR:HD1	1.72	0.54
1:B:27:ASN:ND2	3:D:222:DT:O2	2.41	0.54
1:G:102:GLN:O	1:G:105:GLY:N	2.37	0.54
1:B:193:SER:HA	1:B:196:ARG:NE	2.22	0.54
1:P:163:LEU:HD21	2:Q:205:DT:H1'	1.90	0.53
1:A:86:SER:HB3	1:A:89:PHE:CD2	2.43	0.53
1:M:63:ARG:HG2	1:M:70:THR:HG22	1.90	0.53
1:M:133:LEU:HD11	1:M:191:VAL:HG12	1.90	0.53
2:N:204:DA:N1	3:O:224:DA:N6	2.57	0.53
1:B:74:PHE:CE1	2:C:209:DA:H1'	2.44	0.53
1:J:23:PRO:HA	1:J:122:CYS:HB3	1.90	0.53
1:J:178:ARG:HD3	1:J:182:TYR:CE1	2.44	0.53
2:Q:207:DA:H2'	2:Q:208:DA:C8	2.43	0.53
1:B:102:GLN:HA	1:B:106:PHE:O	2.09	0.53
1:G:34:LEU:HA	1:G:109:LYS:O	2.09	0.53

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:165:PHE:CD2	3:I:226:DA:H5'	2.44	0.53
1:M:174:GLY:O	1:M:176:LYS:NZ	2.41	0.53
1:P:27:ASN:HB2	3:R:223:DT:O4'	2.09	0.53
1:A:26:GLN:HG3	3:F:224:DA:H4'	1.91	0.53
1:J:52:TYR:CZ	1:J:54:PRO:HB3	2.44	0.53
1:B:89:PHE:CZ	1:M:88:ASP:HB3	2.44	0.53
1:G:127:PRO:HB2	1:G:197:LYS:HB3	1.90	0.52
2:Q:211:DC:H2''	2:Q:212:DG:C8	2.44	0.52
1:P:41:LYS:HD3	1:P:41:LYS:N	2.20	0.52
1:A:73:ILE:HG12	1:A:79:MET:HE2	1.91	0.52
1:B:99:ARG:HE	1:B:103:LYS:CE	2.23	0.52
1:P:23:PRO:HD2	1:P:182:TYR:CD1	2.45	0.52
1:B:67:PRO:O	1:B:69:THR:HG23	2.09	0.52
1:J:100:ILE:O	1:J:104:LEU:HD13	2.10	0.52
1:A:124:VAL:HG11	1:A:192:LEU:HD12	1.91	0.52
2:K:201:DG:C8	2:K:201:DG:H5'	2.45	0.52
1:B:26:GLN:HG3	3:D:224:DA:H4'	1.92	0.51
1:B:128:ILE:HD13	1:B:168:GLY:HA2	1.92	0.51
1:J:131:GLU:N	1:J:131:GLU:OE1	2.42	0.51
1:M:40:LEU:HD21	1:M:75:ALA:HA	1.92	0.51
1:B:162:LEU:O	1:B:163:LEU:HD23	2.10	0.51
1:B:117:ASN:C	1:B:118:ILE:HD12	2.35	0.51
1:M:59:ALA:HB2	1:M:74:PHE:CE1	2.46	0.51
1:M:164:ILE:HG23	1:M:170:ILE:HD12	1.93	0.51
1:M:169:LYS:C	1:M:170:ILE:HD13	2.35	0.51
1:G:64:ILE:HG23	1:G:97:TYR:CE1	2.46	0.51
1:G:95:ARG:HG2	1:G:110:PHE:CZ	2.45	0.51
2:K:201:DG:H5'	2:K:201:DG:H8	1.76	0.51
1:A:165:PHE:CD2	3:F:226:DA:H5'	2.46	0.51
1:B:143:TYR:O	1:B:144:GLU:HG3	2.10	0.51
1:P:18:PRO:HD2	1:P:186:GLU:HG3	1.92	0.51
1:P:123:ASP:OD2	1:P:125:LYS:HD3	2.10	0.51
1:B:174:GLY:HA2	2:C:206:DA:H2''	1.93	0.51
1:P:40:LEU:HD21	1:P:59:ALA:HA	1.93	0.51
1:B:159:LYS:O	1:B:160:ILE:HD13	2.11	0.51
1:A:67:PRO:O	1:A:69:THR:HG23	2.11	0.51
1:M:153:TYR:CE1	1:M:155:MET:HG3	2.46	0.51
1:A:46:GLN:HB3	1:A:104:LEU:HD22	1.93	0.50
1:G:25:LEU:HD22	1:G:118:ILE:HG22	1.93	0.50
1:J:130:LEU:HD22	1:J:151:LEU:HB2	1.92	0.50
1:A:19:SER:HB3	1:A:189:TYR:CD2	2.46	0.50

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:116:GLN:HE21	2:C:208:DA:H4'	1.76	0.50
1:P:130:LEU:HD12	1:P:166:VAL:HG23	1.94	0.50
1:B:86:SER:HB3	1:B:89:PHE:CD2	2.47	0.50
2:E:212:DG:H2''	2:E:213:DC:OP2	2.11	0.50
1:M:40:LEU:CD2	1:M:75:ALA:HA	2.42	0.50
1:G:23:PRO:HG3	1:G:185:PHE:CD2	2.47	0.50
2:K:207:DA:N7	2:K:208:DA:N6	2.59	0.50
1:M:61:ILE:HD13	3:O:221:DT:H5''	1.94	0.50
1:A:40:LEU:HD22	1:A:60:VAL:HG23	1.94	0.50
1:B:34:LEU:HA	1:B:109:LYS:O	2.12	0.49
1:G:165:PHE:CE2	3:I:225:DT:H2''	2.47	0.49
1:J:74:PHE:CE2	2:K:209:DA:H1'	2.47	0.49
1:J:153:TYR:CE1	1:J:155:MET:HB2	2.30	0.49
1:P:133:LEU:HA	1:P:195:PHE:CD2	2.47	0.49
1:B:127:PRO:C	1:B:128:ILE:HD12	2.38	0.49
1:B:109:LYS:HE3	1:B:109:LYS:HA	1.94	0.49
1:M:62:MET:HE2	1:M:97:TYR:HD1	1.77	0.49
1:P:173:THR:OG1	1:P:174:GLY:N	2.45	0.49
1:G:149:PRO:HG3	3:I:227:DG:O4'	2.13	0.49
1:J:161:VAL:HG11	2:K:205:DT:H4'	1.95	0.49
1:G:26:GLN:HG3	3:I:224:DA:H4'	1.95	0.49
1:M:34:LEU:HA	1:M:109:LYS:O	2.13	0.49
1:B:99:ARG:HE	1:B:103:LYS:HE2	1.78	0.49
1:J:44:ALA:CB	1:J:52:TYR:HB3	2.42	0.49
1:P:133:LEU:HD13	1:P:195:PHE:HD2	1.76	0.49
1:B:123:ASP:CG	1:B:169:LYS:HD3	2.37	0.48
1:M:17:HIS:CD2	1:M:22:VAL:HG12	2.48	0.48
1:P:130:LEU:HD22	1:P:151:LEU:HD12	1.95	0.48
1:M:42:ALA:O	1:M:46:GLN:HG3	2.13	0.48
2:C:201:DG:H2'	2:C:202:DC:C6	2.47	0.48
1:G:117:ASN:HD22	1:G:118:ILE:H	1.58	0.48
1:M:146:GLU:HG2	1:M:147:LEU:N	2.29	0.48
1:J:185:PHE:HA	1:J:188:ILE:HG22	1.95	0.48
1:J:59:ALA:HB2	1:J:74:PHE:CD1	2.48	0.48
1:J:163:LEU:HD21	2:K:205:DT:H1'	1.94	0.48
1:P:148:PHE:HD1	1:P:149:PRO:CD	2.23	0.48
1:B:141:SER:HB2	1:B:153:TYR:HD1	1.78	0.48
1:A:38:LEU:HB2	1:A:75:ALA:HA	1.96	0.48
1:A:124:VAL:HG11	1:A:192:LEU:CD1	2.44	0.48
2:H:201:DG:H5'	2:H:201:DG:H8	1.79	0.48
1:M:63:ARG:HG2	1:M:70:THR:CB	2.43	0.48

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:169:LYS:C	1:J:170:ILE:HD13	2.39	0.48
1:B:18:PRO:HG2	1:B:19:SER:N	2.28	0.47
2:C:208:DA:C5	2:C:209:DA:C6	3.02	0.47
1:B:158:PRO:O	1:B:160:ILE:N	2.47	0.47
1:J:62:MET:HE2	1:J:97:TYR:HD1	1.79	0.47
1:M:130:LEU:HB3	1:M:143:TYR:CE2	2.50	0.47
1:A:149:PRO:HG3	3:F:227:DG:O4'	2.14	0.47
1:P:127:PRO:HB2	1:P:197:LYS:HB3	1.96	0.47
1:P:144:GLU:HG2	1:P:147:LEU:CB	2.43	0.47
1:B:118:ILE:HD12	1:B:118:ILE:N	2.28	0.47
1:A:173:THR:HG21	2:E:205:DT:O2	2.15	0.47
1:G:19:SER:OG	1:G:186:GLU:HA	2.14	0.47
1:B:18:PRO:HD2	1:B:186:GLU:HG2	1.96	0.47
1:G:56:ARG:NH2	3:I:220:DT:H5''	2.30	0.47
1:G:159:LYS:O	1:G:160:ILE:HD13	2.15	0.47
1:B:74:PHE:CZ	2:C:209:DA:H1'	2.50	0.47
1:G:67:PRO:O	1:G:69:THR:HG23	2.15	0.47
1:J:130:LEU:HD12	1:J:166:VAL:HG23	1.96	0.47
1:M:151:LEU:CD2	1:M:153:TYR:HB2	2.44	0.47
1:M:161:VAL:HG11	2:N:205:DT:H4'	1.97	0.47
1:B:152:ILE:HB	1:B:154:ARG:HH12	1.79	0.47
1:A:27:ASN:OD1	1:A:28:ILE:N	2.48	0.47
1:A:91:LYS:HG3	1:A:113:PHE:CE2	2.50	0.47
1:A:102:GLN:HA	1:A:106:PHE:O	2.15	0.47
1:A:151:LEU:O	1:A:163:LEU:HA	2.15	0.47
1:P:46:GLN:HB2	1:P:104:LEU:CD1	2.44	0.47
1:B:59:ALA:HB2	1:B:74:PHE:CE2	2.50	0.46
2:C:208:DA:N7	2:C:209:DA:C6	2.83	0.46
1:J:178:ARG:HD3	1:J:182:TYR:HE1	1.78	0.46
1:A:116:GLN:O	2:E:207:DA:H4'	2.16	0.46
3:F:226:DA:H2''	3:F:227:DG:C8	2.50	0.46
1:M:54:PRO:HA	1:M:57:PHE:O	2.14	0.46
1:M:130:LEU:HG	1:M:143:TYR:HE2	1.80	0.46
1:P:148:PHE:CD1	1:P:149:PRO:HD2	2.45	0.46
1:B:37:LYS:H	1:B:37:LYS:HG2	1.42	0.46
1:M:23:PRO:HB3	1:M:122:CYS:SG	2.55	0.46
1:A:159:LYS:O	1:A:160:ILE:HD13	2.16	0.46
2:H:206:DA:N7	2:H:207:DA:C6	2.83	0.46
1:P:133:LEU:HA	1:P:195:PHE:CE2	2.50	0.46
1:B:173:THR:HG21	2:C:205:DT:O2	2.15	0.46
1:P:154:ARG:HG3	1:P:154:ARG:HH11	1.80	0.46

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:P:57:PHE:HE1	1:P:74:PHE:CE1	2.33	0.46
1:B:18:PRO:HG2	1:B:19:SER:H	1.80	0.46
1:B:61:ILE:HD13	3:D:221:DT:H5''	1.98	0.46
1:A:133:LEU:HD12	1:A:137:HIS:HD2	1.81	0.46
1:G:165:PHE:CZ	3:I:225:DT:H1'	2.51	0.46
1:J:98:ALA:O	1:J:102:GLN:HG3	2.16	0.46
1:M:178:ARG:HD3	1:M:182:TYR:OH	2.15	0.46
1:B:143:TYR:CB	1:B:151:LEU:HD12	2.47	0.45
1:A:162:LEU:O	1:A:163:LEU:HD23	2.16	0.45
3:I:227:DG:H2''	3:I:228:DC:O5'	2.16	0.45
1:J:175:ALA:HB1	1:J:177:MET:O	2.17	0.45
1:A:135:TYR:HA	1:A:138:ALA:HB2	1.99	0.45
1:J:87:GLU:OE1	1:J:87:GLU:N	2.28	0.45
1:M:23:PRO:HA	1:M:122:CYS:HB3	1.97	0.45
1:G:96:LYS:HE2	1:J:66:GLU:OE1	2.16	0.45
1:G:163:LEU:HB3	1:G:165:PHE:HE1	1.81	0.45
1:P:127:PRO:C	1:P:128:ILE:HD12	2.41	0.45
1:P:144:GLU:HG2	1:P:147:LEU:HB2	1.97	0.45
1:B:64:ILE:HG12	1:B:97:TYR:OH	2.17	0.45
1:B:116:GLN:OE1	1:B:116:GLN:HA	2.16	0.45
1:B:140:PHE:HB3	1:B:153:TYR:CE1	2.50	0.45
1:A:55:LYS:HD2	1:A:55:LYS:HA	1.68	0.45
3:L:224:DA:H2'	3:L:225:DT:O4'	2.17	0.45
1:P:34:LEU:HB2	1:P:77:GLY:O	2.17	0.45
1:B:79:MET:HE1	1:B:97:TYR:HB2	1.99	0.45
1:A:27:ASN:HB2	3:F:223:DT:O4'	2.16	0.45
1:B:89:PHE:CE1	1:M:88:ASP:HB3	2.51	0.45
1:A:74:PHE:CE1	2:E:209:DA:H1'	2.52	0.45
1:G:124:VAL:HG11	1:G:192:LEU:HD23	1.99	0.45
1:P:40:LEU:HG	1:P:60:VAL:CG2	2.45	0.45
1:B:66:GLU:HA	1:B:67:PRO:C	2.42	0.45
3:L:225:DT:H2''	3:L:226:DA:H5'	1.98	0.45
1:B:99:ARG:O	1:B:103:LYS:HG2	2.17	0.45
1:G:59:ALA:HB2	1:G:74:PHE:CD1	2.52	0.45
1:M:93:ALA:O	1:M:97:TYR:HD2	1.99	0.45
1:M:117:ASN:HD21	2:N:206:DA:H1'	1.82	0.45
1:B:143:TYR:CG	1:B:151:LEU:HD12	2.52	0.45
1:G:18:PRO:HG2	1:G:186:GLU:HB3	1.99	0.45
2:H:204:DA:H2''	2:H:205:DT:H5'	1.97	0.45
1:J:26:GLN:HG3	3:L:224:DA:H4'	2.00	0.45
1:B:165:PHE:CD1	3:D:226:DA:H5'	2.52	0.44

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:173:THR:OG1	2:C:206:DA:N3	2.41	0.44
1:J:167:SER:OG	1:J:169:LYS:HG2	2.16	0.44
1:M:130:LEU:HD22	1:M:166:VAL:HG23	1.98	0.44
1:P:18:PRO:HD2	1:P:186:GLU:CD	2.42	0.44
1:P:42:ALA:O	1:P:46:GLN:HG3	2.16	0.44
1:B:182:TYR:N	1:B:182:TYR:CD1	2.85	0.44
1:B:183:LYS:O	1:B:186:GLU:N	2.46	0.44
2:H:202:DC:H2'	2:H:203:DT:C6	2.52	0.44
3:O:224:DA:H2'	3:O:225:DT:O4'	2.18	0.44
3:D:222:DT:C4	3:D:223:DT:O4	2.71	0.44
1:G:59:ALA:HB2	1:G:74:PHE:CE1	2.52	0.44
1:J:111:LYS:HD3	1:J:111:LYS:HA	1.68	0.44
1:P:97:TYR:O	1:P:101:VAL:HG22	2.18	0.44
1:P:130:LEU:HD22	1:P:151:LEU:HB2	2.00	0.44
1:P:163:LEU:HB2	1:P:171:VAL:HB	1.98	0.44
1:M:130:LEU:CG	1:M:143:TYR:HE2	2.31	0.44
1:P:121:SER:HA	1:P:170:ILE:O	2.18	0.44
1:A:118:ILE:HD12	1:A:118:ILE:N	2.33	0.44
1:P:73:ILE:HG12	1:P:79:MET:CG	2.48	0.44
1:P:160:ILE:HG21	1:P:184:ALA:CB	2.48	0.44
1:B:189:TYR:HB3	1:B:190:PRO:HD3	2.00	0.44
1:G:25:LEU:HD23	1:G:25:LEU:HA	1.84	0.44
1:M:95:ARG:O	1:M:98:ALA:N	2.51	0.44
1:P:143:TYR:CD1	1:P:145:PRO:HD3	2.53	0.44
1:B:169:LYS:H	1:B:169:LYS:HG2	1.68	0.43
3:D:224:DA:H2'	3:D:225:DT:O4'	2.17	0.43
1:M:109:LYS:HD2	1:M:109:LYS:HA	1.72	0.43
1:A:53:ASN:HD22	1:A:56:ARG:HD3	1.83	0.43
1:A:111:LYS:O	1:A:112:ASP:HB2	2.17	0.43
1:B:47:ALA:HB2	1:B:104:LEU:HD11	2.00	0.43
1:B:122:CYS:HB3	1:B:170:ILE:HD11	2.01	0.43
1:A:64:ILE:HG23	1:A:97:TYR:CZ	2.53	0.43
1:G:52:TYR:CZ	1:G:54:PRO:HB3	2.53	0.43
1:G:87:GLU:OE1	1:G:178:ARG:NH2	2.39	0.43
1:G:163:LEU:HB3	1:G:165:PHE:CE1	2.53	0.43
1:J:118:ILE:HB	1:J:175:ALA:HB3	1.99	0.43
1:J:126:PHE:CZ	1:J:193:SER:HA	2.53	0.43
1:P:160:ILE:HG21	1:P:184:ALA:HB1	2.01	0.43
1:A:35:ASP:HB3	1:A:109:LYS:HB2	2.01	0.43
1:G:118:ILE:HB	1:G:175:ALA:HB3	2.00	0.43
1:P:40:LEU:HD23	1:P:52:TYR:CE2	2.54	0.43

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:P:148:PHE:CD2	2:Q:204:DA:O4'	2.72	0.43
1:G:31:THR:HG21	1:G:116:GLN:HE21	1.83	0.43
1:G:89:PHE:CZ	1:J:88:ASP:HB3	2.54	0.43
1:G:87:GLU:OE2	1:G:178:ARG:HB3	2.18	0.43
1:M:133:LEU:HA	1:M:195:PHE:HD1	1.84	0.43
1:B:123:ASP:CB	1:B:169:LYS:HD3	2.48	0.43
1:G:61:ILE:HD13	3:I:221:DT:H5''	2.00	0.43
1:M:133:LEU:HA	1:M:195:PHE:CD1	2.54	0.43
1:B:87:GLU:OE1	1:B:178:ARG:NH2	2.35	0.42
1:B:170:ILE:HD12	1:B:170:ILE:O	2.19	0.42
1:J:173:THR:OG1	1:J:174:GLY:N	2.52	0.42
1:M:179:ASP:HA	1:M:182:TYR:HD1	1.83	0.42
1:B:54:PRO:HA	1:B:57:PHE:O	2.20	0.42
1:B:165:PHE:CZ	3:D:225:DT:H1'	2.54	0.42
1:A:101:VAL:HG12	1:A:106:PHE:HB2	2.00	0.42
1:A:152:ILE:HA	1:A:162:LEU:O	2.19	0.42
1:J:156:LYS:O	1:J:159:LYS:HE3	2.20	0.42
3:L:222:DT:C4	3:L:223:DT:O4	2.72	0.42
1:B:158:PRO:HG2	1:B:160:ILE:HG12	2.01	0.42
1:A:46:GLN:HB3	1:A:104:LEU:CD2	2.49	0.42
1:J:129:ARG:HG3	1:J:129:ARG:O	2.19	0.42
1:M:91:LYS:HD2	1:M:113:PHE:CD2	2.54	0.42
1:B:149:PRO:HG3	3:D:227:DG:O4'	2.20	0.42
1:A:40:LEU:HD23	1:A:40:LEU:HA	1.78	0.42
1:A:127:PRO:HA	1:A:166:VAL:O	2.19	0.42
1:G:64:ILE:HG12	1:G:97:TYR:OH	2.19	0.42
2:H:206:DA:N7	2:H:207:DA:N6	2.67	0.42
1:J:73:ILE:HG23	1:J:73:ILE:HD12	1.78	0.42
1:J:99:ARG:HA	1:J:102:GLN:HG3	2.01	0.42
1:A:162:LEU:HD23	1:A:163:LEU:N	2.35	0.42
1:M:156:LYS:HA	1:M:156:LYS:HD3	1.78	0.42
1:B:143:TYR:C	1:B:144:GLU:HG3	2.45	0.42
1:M:158:PRO:HB2	1:M:160:ILE:HG12	2.01	0.42
1:P:128:ILE:HD13	1:P:168:GLY:HA2	2.01	0.42
1:P:173:THR:OG1	2:Q:206:DA:H1'	2.20	0.42
1:A:49:ASN:O	1:A:62:MET:HA	2.20	0.42
1:A:162:LEU:C	1:A:163:LEU:HD23	2.44	0.42
1:M:52:TYR:HB2	1:M:60:VAL:HG13	2.02	0.42
1:B:26:GLN:HG3	3:D:224:DA:C4'	2.49	0.42
2:C:201:DG:H2'	2:C:202:DC:H6	1.84	0.42
3:F:224:DA:N7	3:F:225:DT:C4	2.88	0.42

*Continued on next page...*



*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:143:TYR:HE2	1:J:145:PRO:CB	2.33	0.42
1:M:126:PHE:CE2	1:M:193:SER:HA	2.55	0.42
1:P:117:ASN:OD1	1:P:118:ILE:N	2.53	0.42
1:B:137:HIS:CB	1:B:140:PHE:HB2	2.43	0.42
1:B:164:ILE:HG12	1:B:170:ILE:HG22	2.01	0.42
1:P:34:LEU:HA	1:P:109:LYS:O	2.20	0.42
1:P:169:LYS:C	1:P:170:ILE:HD13	2.45	0.41
1:B:181:THR:HG22	1:B:182:TYR:HD1	1.83	0.41
1:A:25:LEU:HD23	1:A:25:LEU:HA	1.74	0.41
1:M:93:ALA:O	1:M:97:TYR:CD2	2.73	0.41
1:B:73:ILE:HD11	1:B:79:MET:SD	2.61	0.41
2:C:201:DG:H5'	2:C:201:DG:H8	1.85	0.41
1:G:34:LEU:HB2	1:G:77:GLY:O	2.19	0.41
1:G:143:TYR:HB2	1:G:151:LEU:HD12	2.02	0.41
1:J:25:LEU:H	1:J:25:LEU:HG	1.75	0.41
1:P:95:ARG:HG2	1:P:110:PHE:CZ	2.55	0.41
1:G:167:SER:HB2	1:G:169:LYS:HG2	2.02	0.41
1:P:61:ILE:HD13	3:R:221:DT:H5''	2.02	0.41
1:J:153:TYR:CD1	1:J:153:TYR:C	2.98	0.41
1:M:169:LYS:HD3	3:O:226:DA:P	2.60	0.41
1:P:28:ILE:HD12	1:P:86:SER:C	2.45	0.41
1:G:189:TYR:HD1	1:G:189:TYR:HA	1.68	0.41
1:B:31:THR:HG22	1:B:80:VAL:HG22	2.03	0.41
1:A:65:ARG:N	4:A:301:SO4:O4	2.48	0.41
1:A:140:PHE:HB2	1:A:153:TYR:CE1	2.55	0.41
1:J:27:ASN:C	1:J:28:ILE:HD12	2.46	0.41
1:G:152:ILE:HD12	2:H:205:DT:C5'	2.49	0.41
1:J:115:ILE:H	1:J:115:ILE:HG13	1.74	0.41
1:J:129:ARG:HG2	1:J:195:PHE:HB3	2.02	0.41
1:P:195:PHE:HD1	1:P:195:PHE:HA	1.74	0.41
1:B:60:VAL:HB	1:B:73:ILE:HB	2.02	0.41
1:B:190:PRO:HA	1:B:193:SER:HB3	2.02	0.41
1:A:85:LYS:HE3	1:A:85:LYS:HB2	1.87	0.41
1:J:40:LEU:HD21	1:J:75:ALA:HA	2.03	0.41
1:J:47:ALA:HB2	1:J:104:LEU:HD21	2.03	0.41
1:J:128:ILE:HD13	1:J:192:LEU:HD22	2.02	0.41
1:J:141:SER:CB	1:J:151:LEU:HD21	2.49	0.41
1:P:153:TYR:CD1	1:P:153:TYR:C	2.99	0.41
1:B:99:ARG:NH2	4:B:302:SO4:O2	2.54	0.41
1:G:104:LEU:HD13	1:G:104:LEU:HA	1.81	0.41
1:J:109:LYS:HE3	1:J:110:PHE:N	2.23	0.41

*Continued on next page...*



Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:155:MET:HG2	1:J:158:PRO:O	2.20	0.41
1:B:34:LEU:HG	1:B:108:ALA:HB1	2.02	0.40
1:B:91:LYS:HE3	1:B:95:ARG:HH22	1.86	0.40
2:C:211:DC:H2''	2:C:212:DG:C8	2.56	0.40
2:H:206:DA:H62	2:H:207:DA:H61	1.69	0.40
1:J:40:LEU:HG	1:J:75:ALA:HB2	2.01	0.40
3:O:222:DT:C4	3:O:223:DT:O4	2.74	0.40
1:B:159:LYS:C	1:B:160:ILE:HD13	2.47	0.40
1:A:161:VAL:CG1	2:E:205:DT:H4'	2.48	0.40
1:P:96:LYS:O	1:P:100:ILE:HG13	2.21	0.40
1:B:109:LYS:HA	1:B:109:LYS:CE	2.51	0.40
1:J:42:ALA:O	1:J:46:GLN:HG3	2.21	0.40
1:M:153:TYR:HE1	1:M:155:MET:CG	2.35	0.40
1:B:52:TYR:CZ	1:B:54:PRO:HB3	2.56	0.40
1:B:155:MET:HB2	1:B:155:MET:HE3	1.77	0.40
1:P:18:PRO:HD2	1:P:186:GLU:CG	2.52	0.40
1:P:141:SER:CB	1:P:151:LEU:HD21	2.51	0.40
2:C:206:DA:N6	2:C:207:DA:H61	2.18	0.40
1:G:27:ASN:HB2	3:I:223:DT:O4'	2.21	0.40
1:J:128:ILE:HD12	1:J:168:GLY:HA2	2.03	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	185/203 (91%)	166 (90%)	19 (10%)	0	100	100
1	B	185/203 (91%)	164 (89%)	21 (11%)	0	100	100
1	G	183/203 (90%)	167 (91%)	16 (9%)	0	100	100
1	J	185/203 (91%)	174 (94%)	11 (6%)	0	100	100

Continued on next page...

*Continued from previous page...*

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	M	186/203 (92%)	174 (94%)	12 (6%)	0	100	100
1	P	185/203 (91%)	171 (92%)	14 (8%)	0	100	100
All	All	1109/1218 (91%)	1016 (92%)	93 (8%)	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	149/173 (86%)	145 (97%)	4 (3%)	39	71
1	B	147/173 (85%)	146 (99%)	1 (1%)	76	91
1	G	145/173 (84%)	140 (97%)	5 (3%)	32	65
1	J	156/173 (90%)	152 (97%)	4 (3%)	40	72
1	M	154/173 (89%)	153 (99%)	1 (1%)	78	92
1	P	149/173 (86%)	147 (99%)	2 (1%)	61	84
All	All	900/1038 (87%)	883 (98%)	17 (2%)	50	78

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	18	PRO
1	A	103	LYS
1	A	151	LEU
1	A	167	SER
1	A	176	LYS
1	G	104	LEU
1	G	142	SER
1	G	163	LEU
1	G	167	SER
1	G	188	ILE
1	J	22	VAL
1	J	41	LYS

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type
1	J	121	SER
1	J	166	VAL
1	M	103	LYS
1	P	22	VAL
1	P	109	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	49	ASN
1	A	137	HIS
1	G	102	GLN
1	G	117	ASN
1	M	137	HIS
1	P	26	GLN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

12 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
4	SO4	J	301	-	4,4,4	0.24	0	6,6,6	0.12	0
4	SO4	G	302	-	4,4,4	0.26	0	6,6,6	0.11	0
4	SO4	M	301	-	4,4,4	0.29	0	6,6,6	0.21	0
4	SO4	P	302	-	4,4,4	0.30	0	6,6,6	0.28	0
4	SO4	P	301	-	4,4,4	0.20	0	6,6,6	0.13	0
4	SO4	B	301	-	4,4,4	0.30	0	6,6,6	0.40	0
4	SO4	A	301	-	4,4,4	0.27	0	6,6,6	0.28	0
4	SO4	J	302	-	4,4,4	0.29	0	6,6,6	0.08	0
4	SO4	M	302	-	4,4,4	0.20	0	6,6,6	0.20	0
4	SO4	B	302	-	4,4,4	0.25	0	6,6,6	0.25	0
4	SO4	P	303	-	4,4,4	0.29	0	6,6,6	0.30	0
4	SO4	G	301	-	4,4,4	0.25	0	6,6,6	0.23	0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

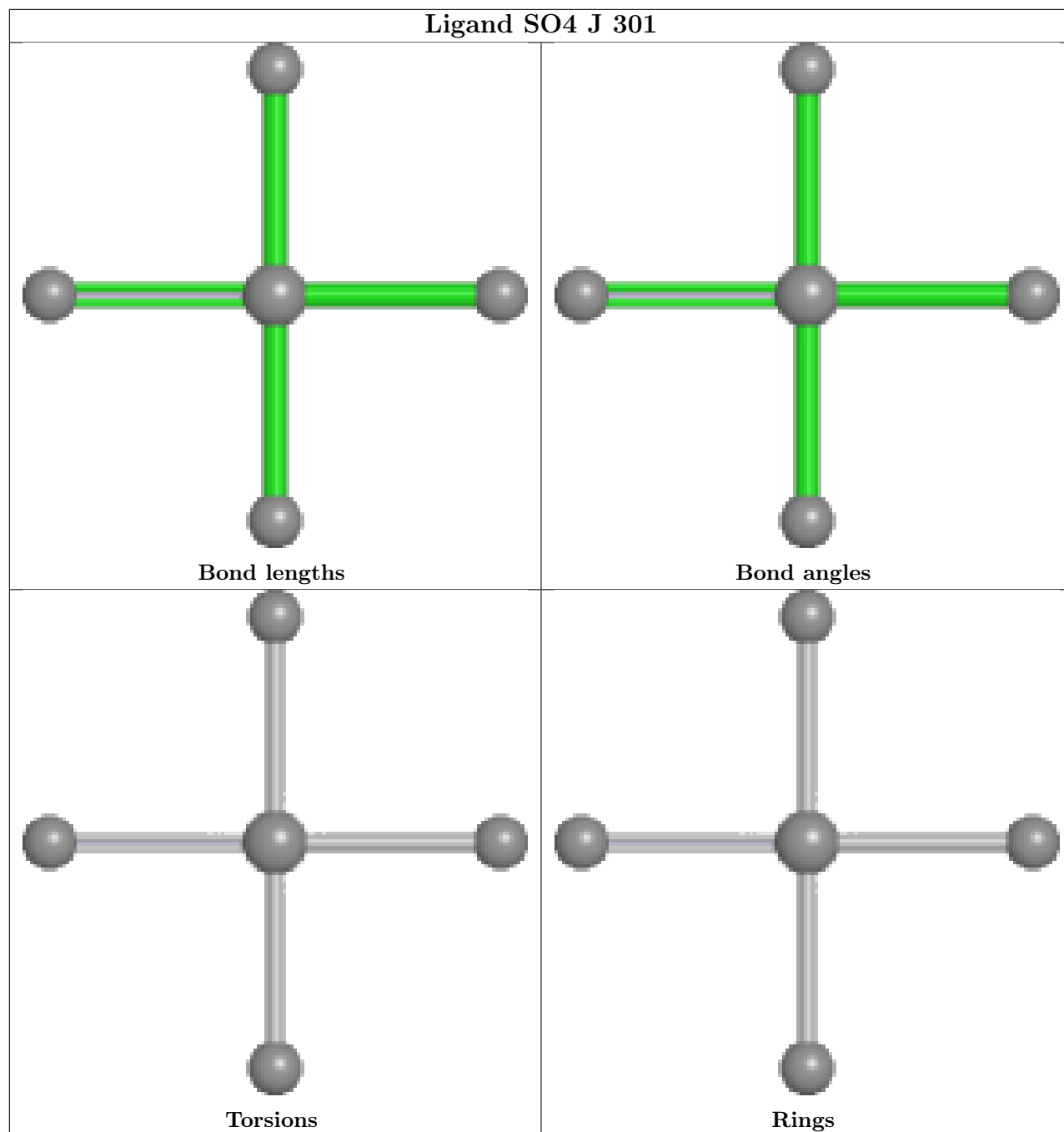
There are no torsion outliers.

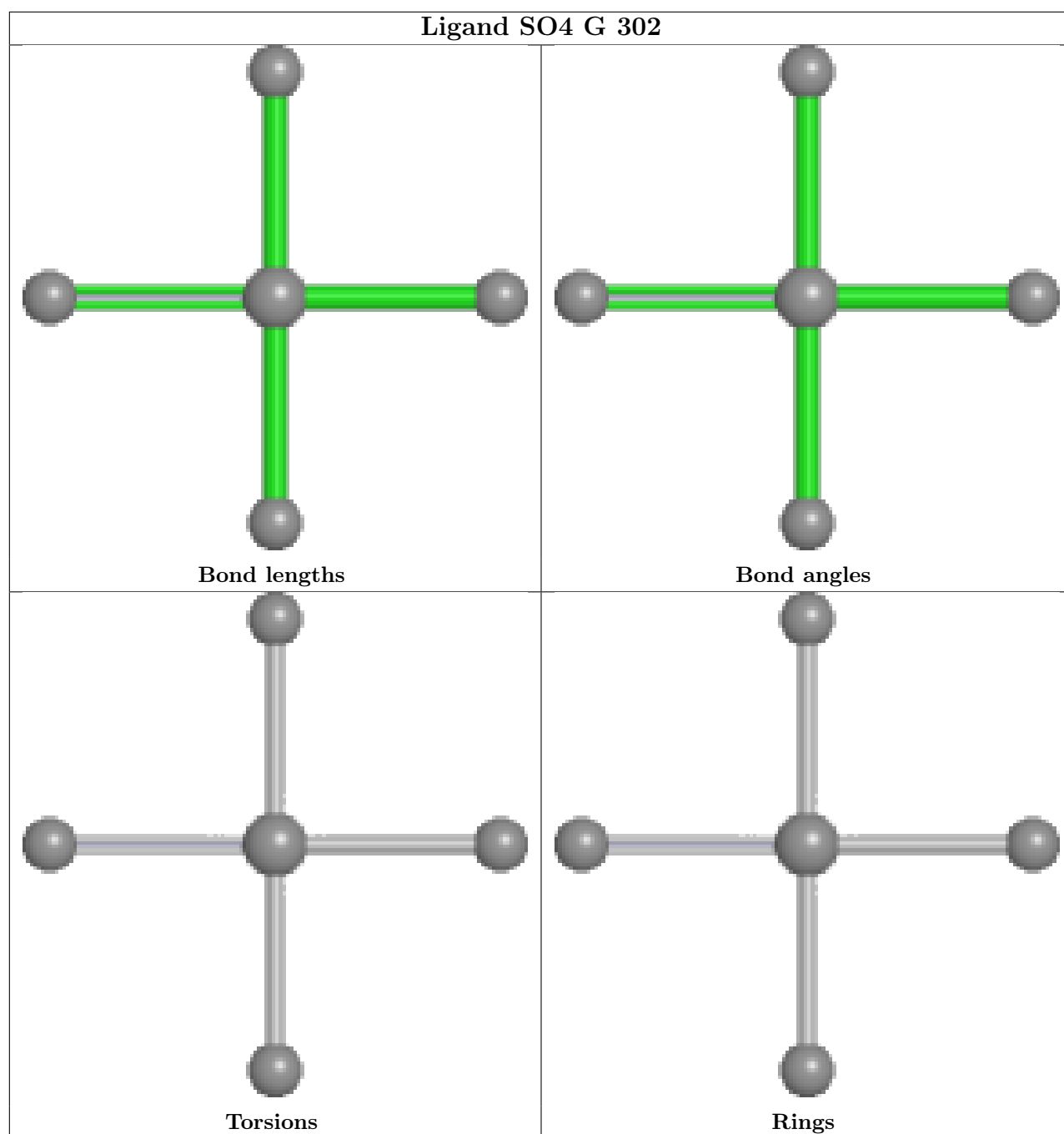
There are no ring outliers.

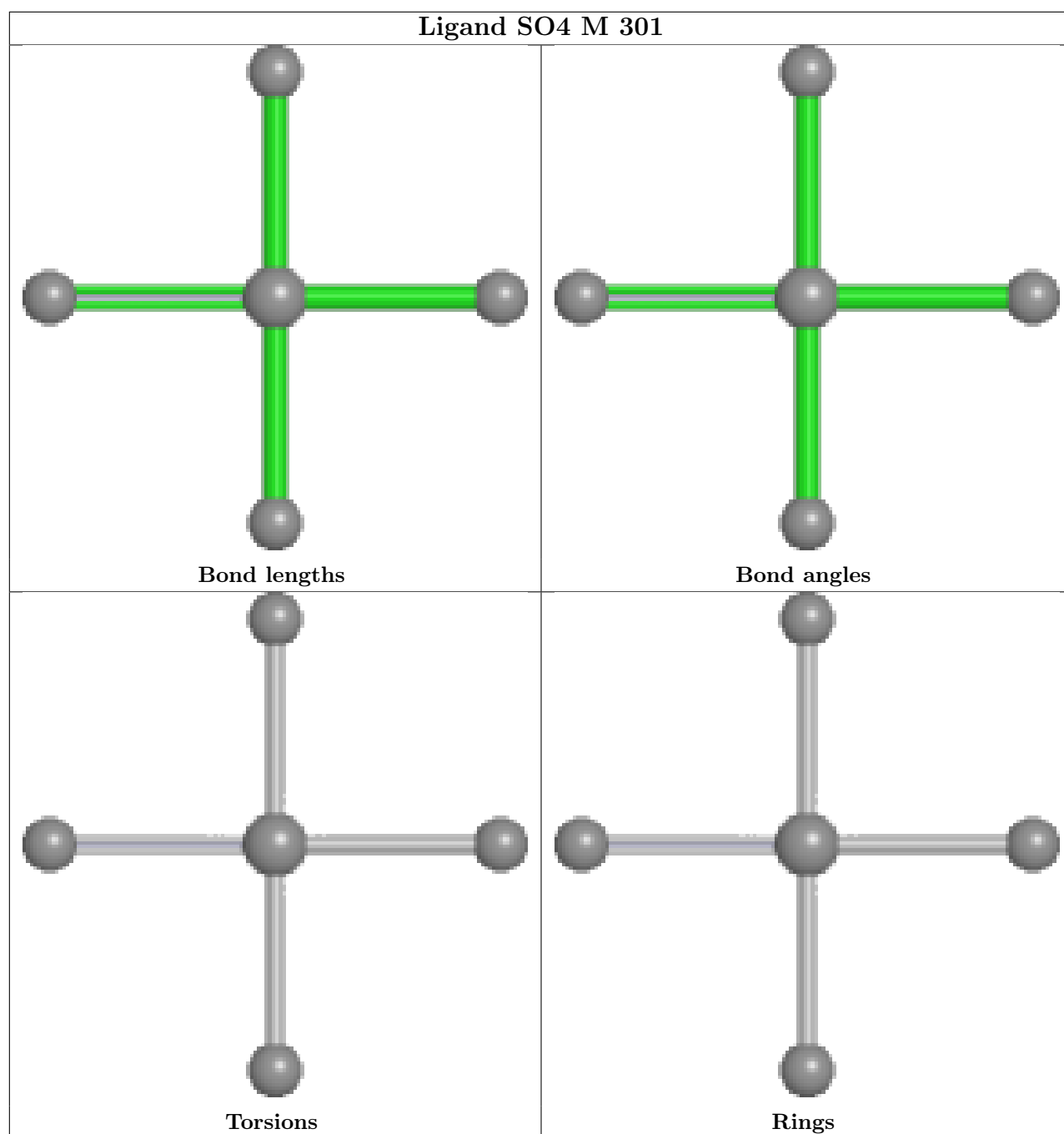
4 monomers are involved in 4 short contacts:

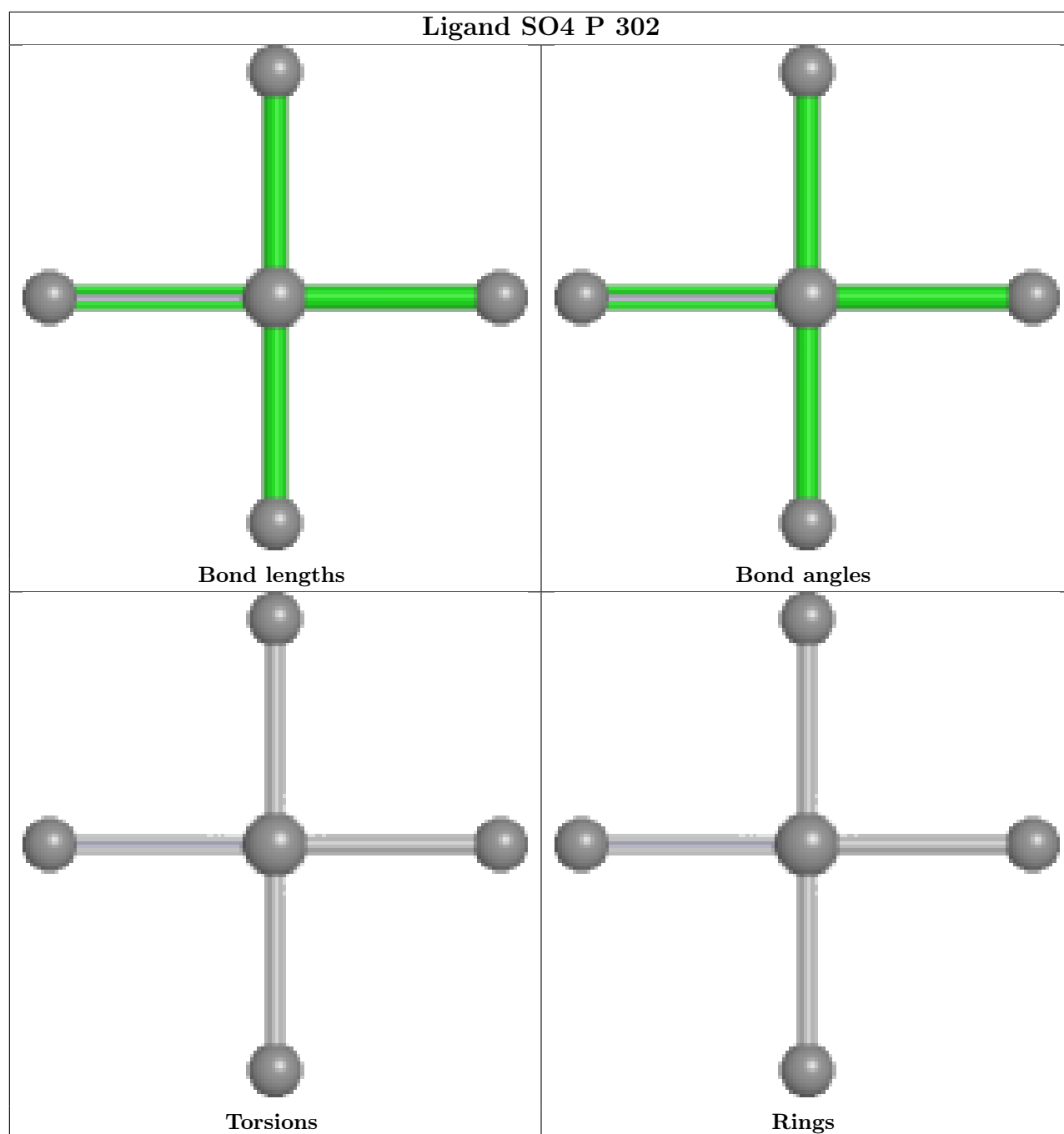
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	J	301	SO4	1	0
4	A	301	SO4	1	0
4	B	302	SO4	1	0
4	G	301	SO4	1	0

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

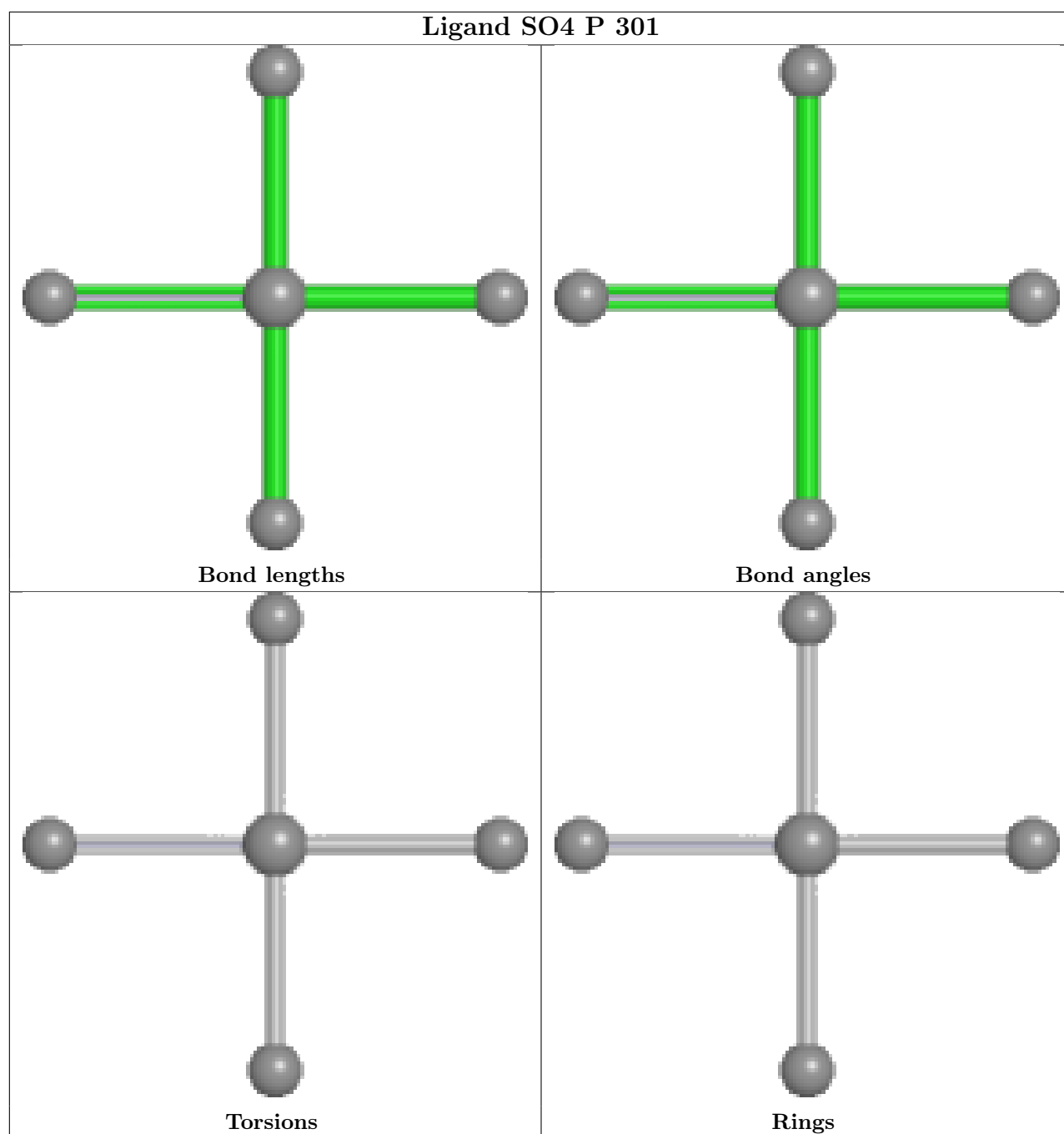


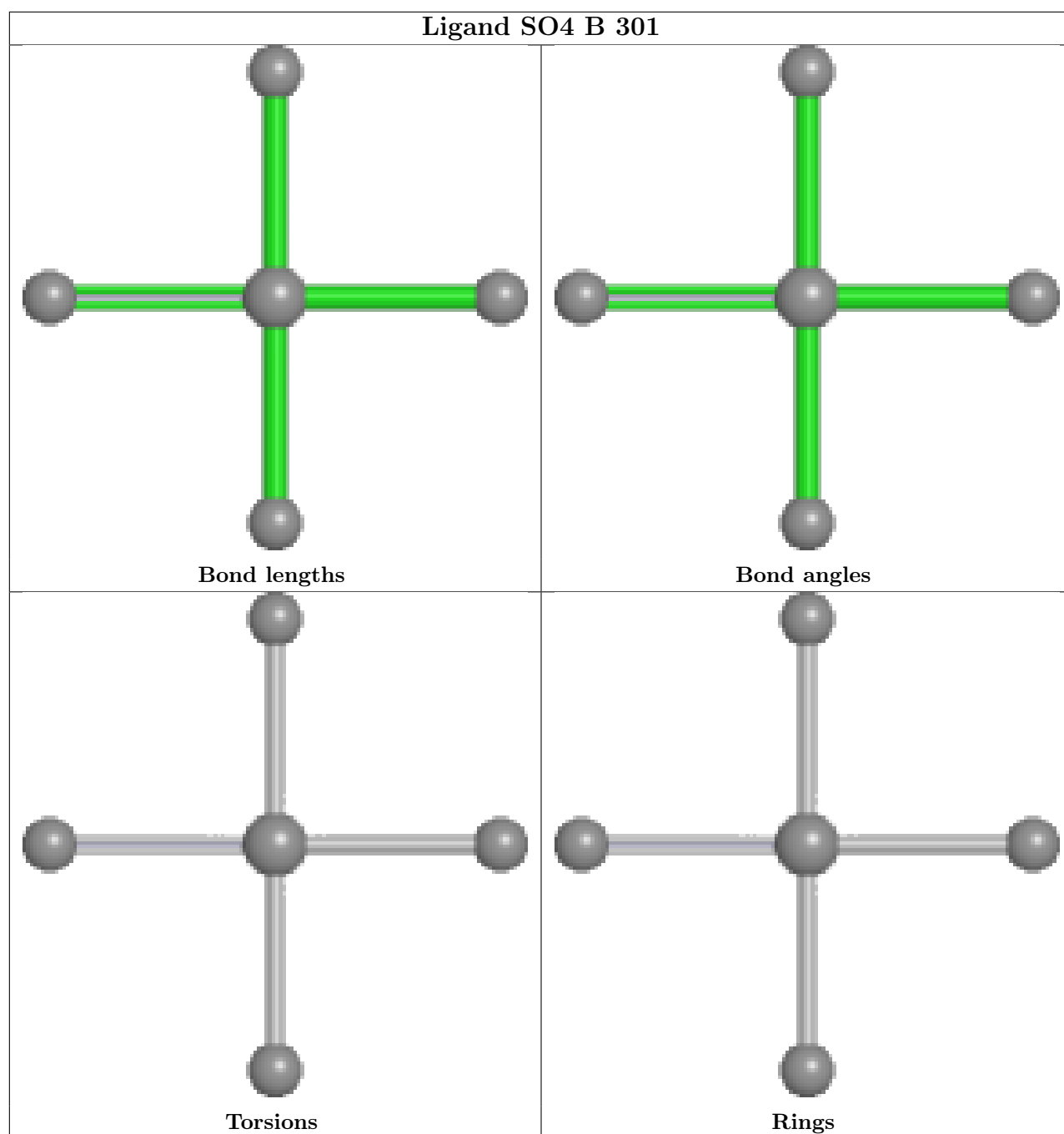


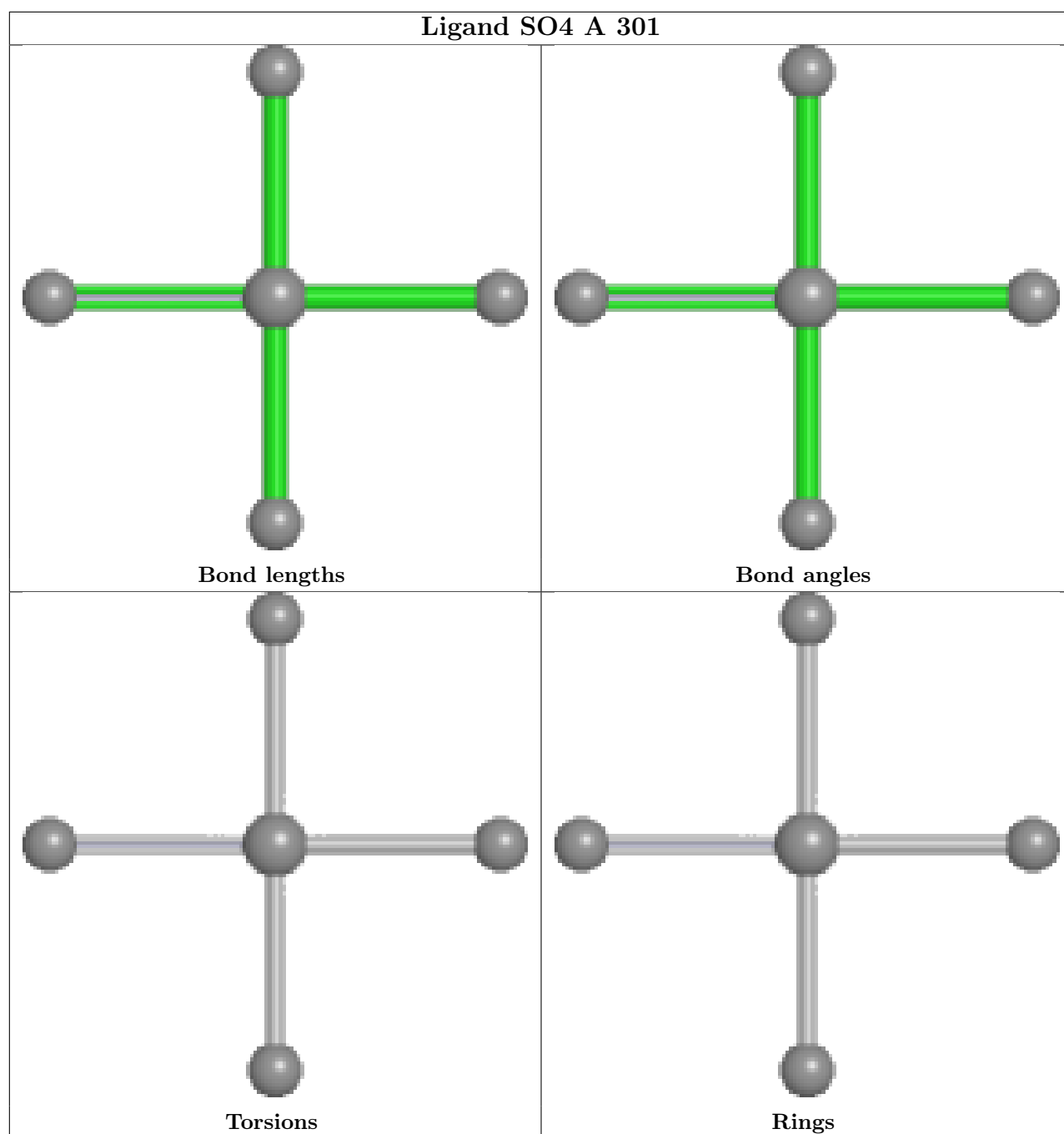


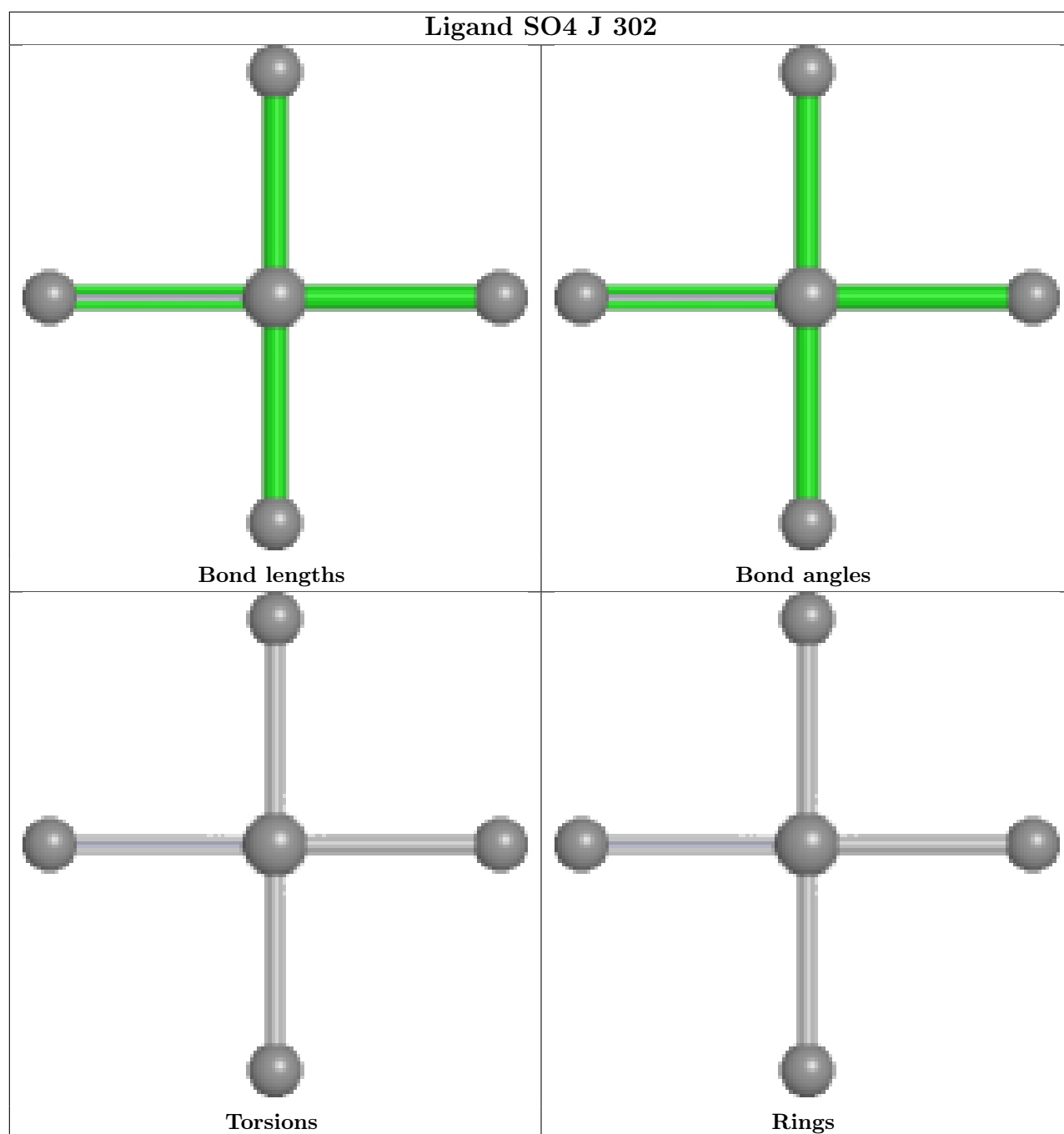


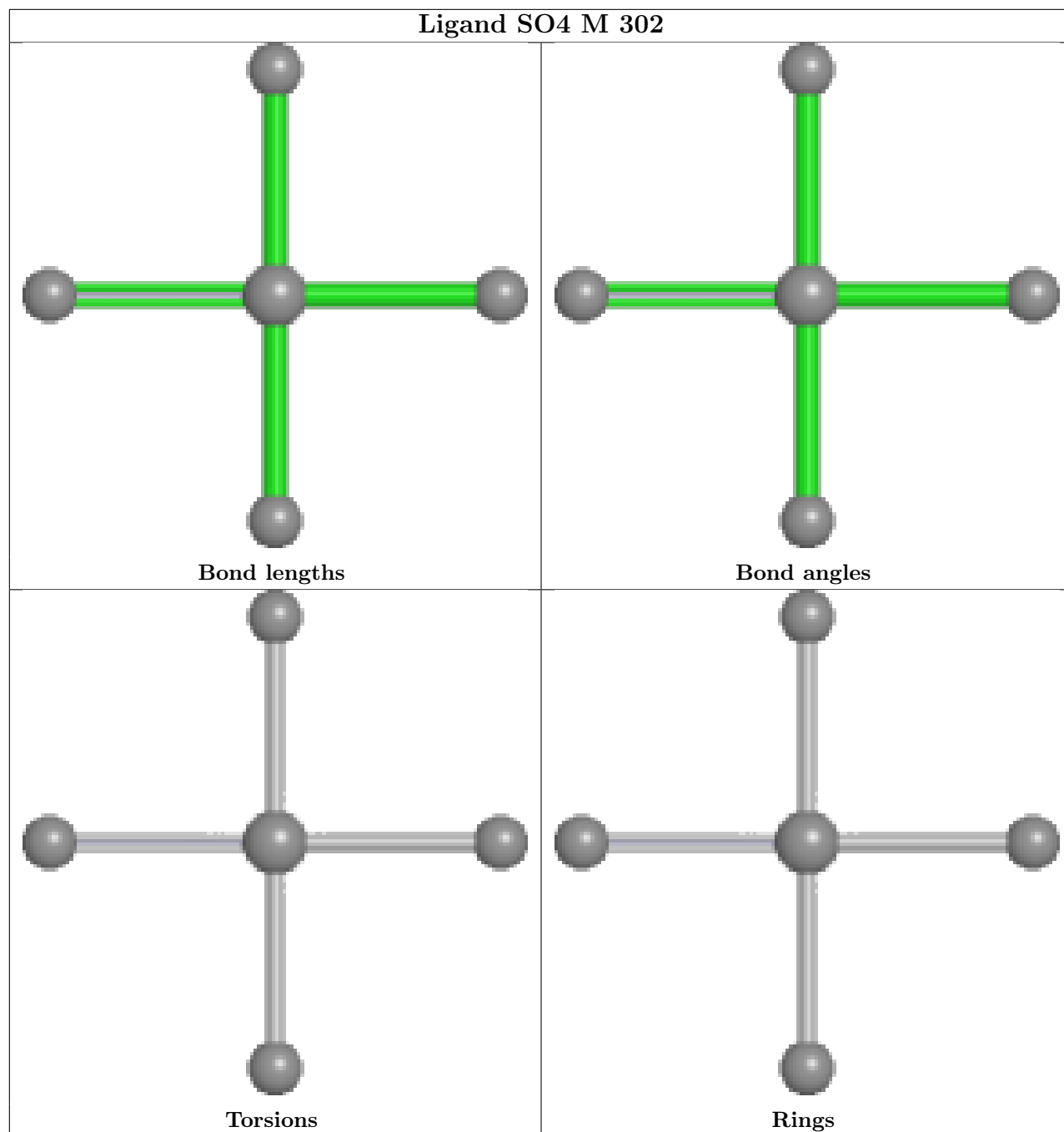


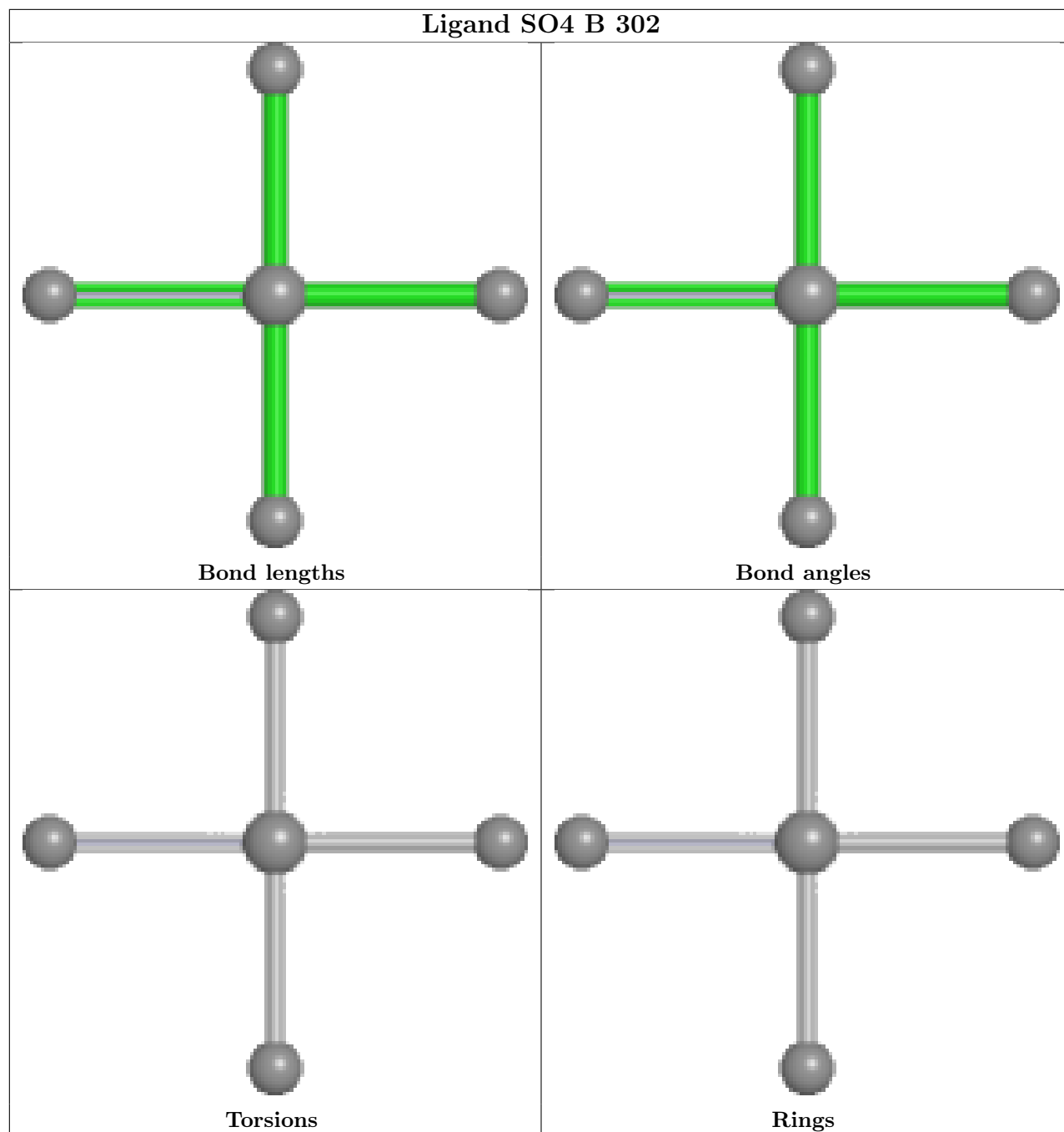


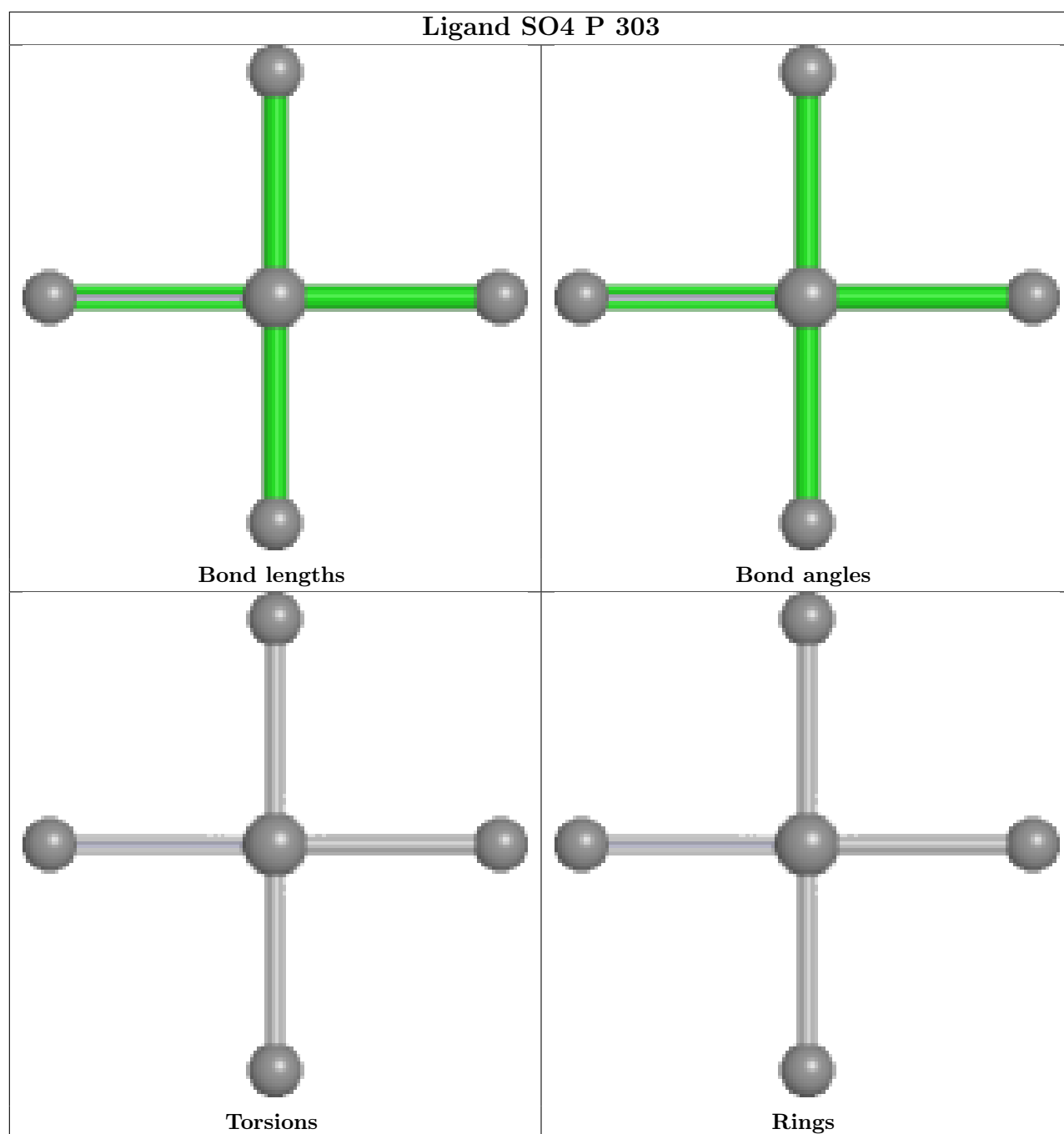


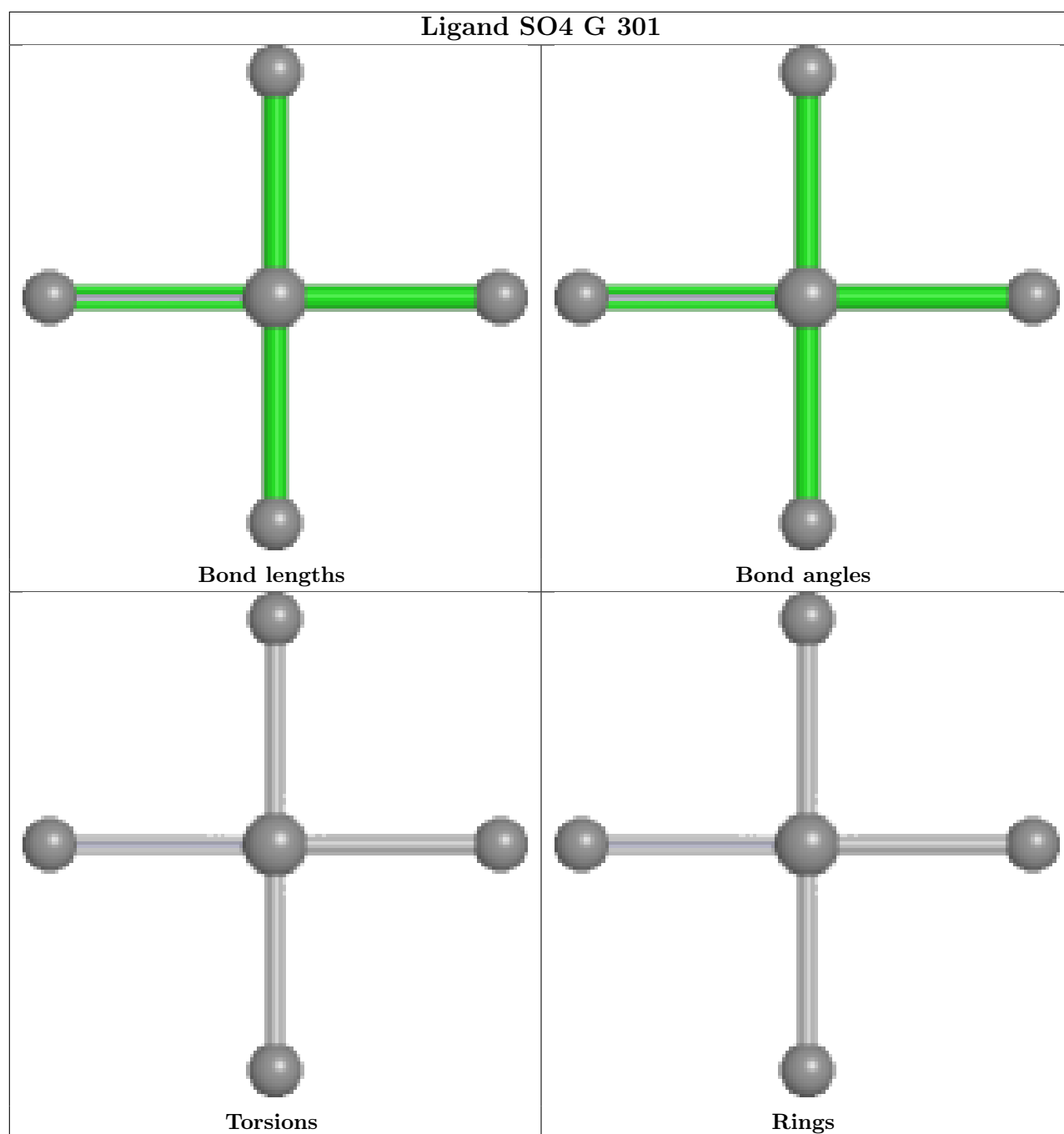












## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.



## 6 Fit of model and data

### 6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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	187/203 (92%)	-0.90	1 (0%) 87 83	56, 77, 119, 147	0
1	B	187/203 (92%)	-0.96	0 100 100	48, 73, 111, 117	0
1	G	185/203 (91%)	-0.92	0 100 100	51, 74, 119, 141	0
1	J	187/203 (92%)	-1.06	0 100 100	64, 86, 105, 116	0
1	M	188/203 (92%)	-0.90	0 100 100	59, 86, 109, 130	0
1	P	187/203 (92%)	-1.01	0 100 100	63, 86, 105, 122	0
2	C	13/13 (100%)	-1.49	0 100 100	66, 79, 99, 103	0
2	E	13/13 (100%)	-1.40	0 100 100	64, 79, 106, 107	0
2	H	13/13 (100%)	-1.35	0 100 100	63, 80, 103, 104	0
2	K	13/13 (100%)	-1.22	0 100 100	79, 86, 128, 133	0
2	N	13/13 (100%)	-1.16	0 100 100	80, 90, 123, 124	0
2	Q	13/13 (100%)	-1.23	0 100 100	81, 90, 127, 130	0
3	D	9/9 (100%)	-1.48	0 100 100	58, 64, 94, 103	0
3	F	9/9 (100%)	-1.49	0 100 100	61, 68, 109, 110	0
3	I	9/9 (100%)	-1.46	0 100 100	58, 69, 104, 111	0
3	L	9/9 (100%)	-1.57	0 100 100	66, 78, 96, 100	0
3	O	9/9 (100%)	-1.54	0 100 100	69, 81, 99, 100	0
3	R	9/9 (100%)	-1.39	0 100 100	69, 78, 100, 103	0
All	All	1253/1350 (92%)	-1.00	1 (0%) 92 90	48, 83, 114, 147	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	20	GLY	2.4

## 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 ⓘ

There are no oligosaccharides in this entry.

## 6.4 Ligands ⓘ

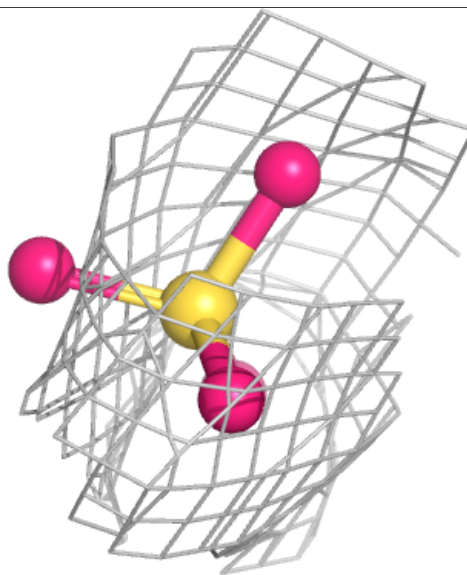
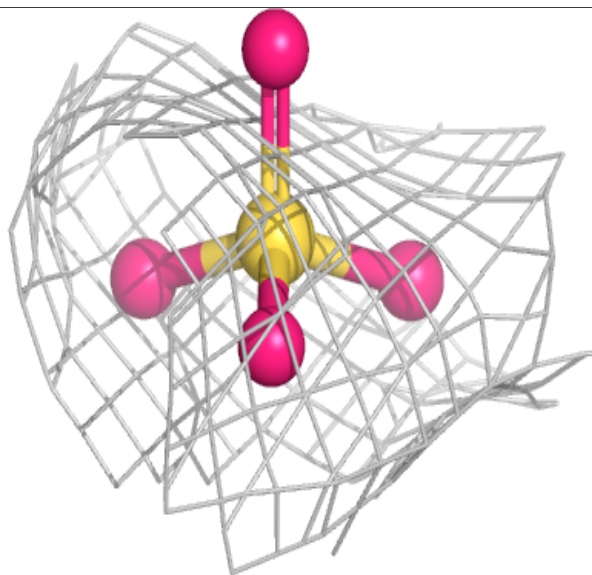
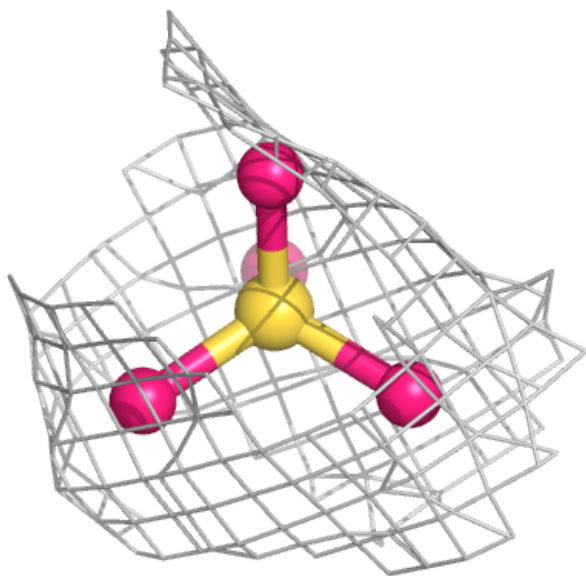
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
4	SO4	G	302	5/5	0.95	0.06	123,123,140,146	0
4	SO4	B	302	5/5	0.97	0.06	95,96,116,122	0
4	SO4	J	301	5/5	0.97	0.05	130,135,156,159	0
4	SO4	P	303	5/5	0.97	0.09	101,104,107,120	0
4	SO4	P	301	5/5	0.98	0.04	141,144,149,155	0
4	SO4	G	301	5/5	0.98	0.16	83,87,96,100	0
4	SO4	J	302	5/5	0.99	0.07	91,93,105,106	0
4	SO4	M	301	5/5	0.99	0.06	88,93,100,101	0
4	SO4	M	302	5/5	0.99	0.04	101,108,122,123	0
4	SO4	A	301	5/5	0.99	0.06	94,97,105,106	0
4	SO4	P	302	5/5	0.99	0.05	86,88,95,99	0
4	SO4	B	301	5/5	0.99	0.13	80,82,96,99	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.

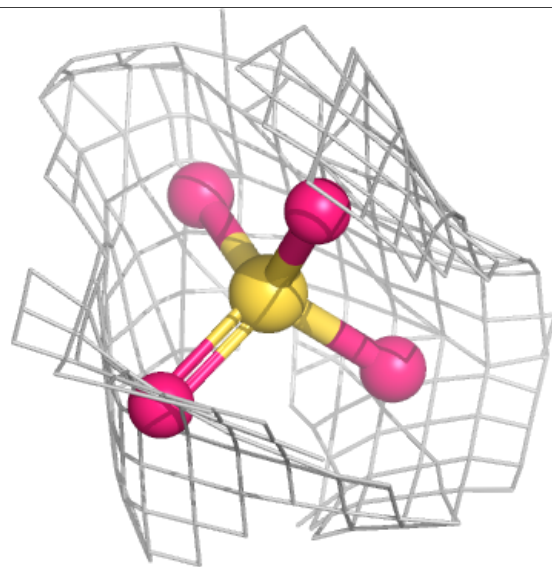
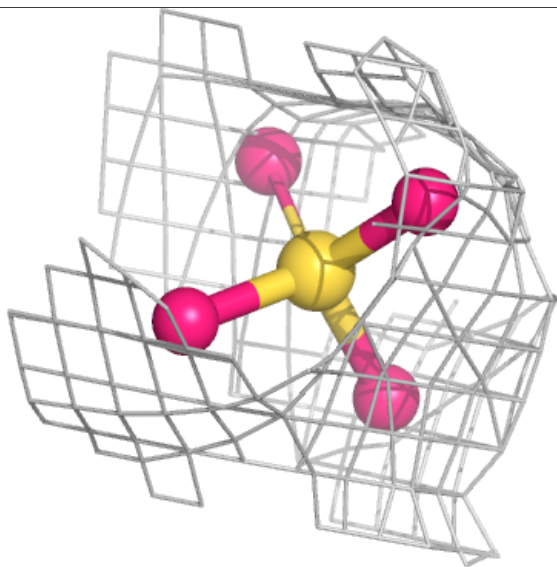
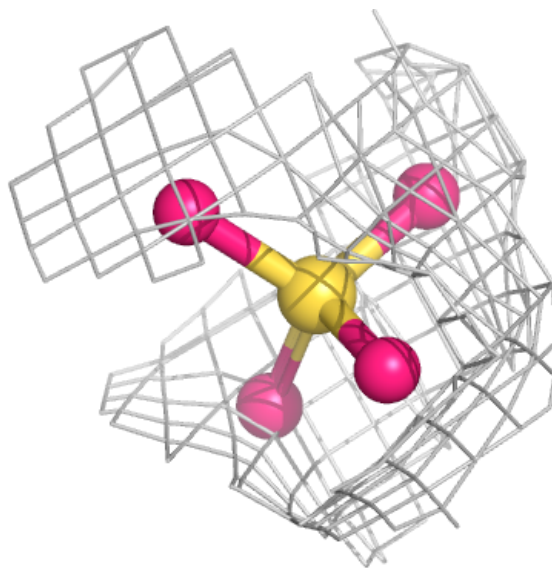
**Electron density around SO4 G 302:**

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



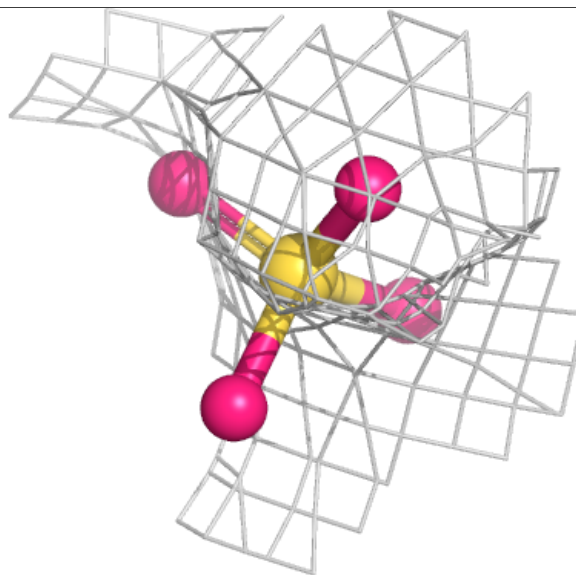
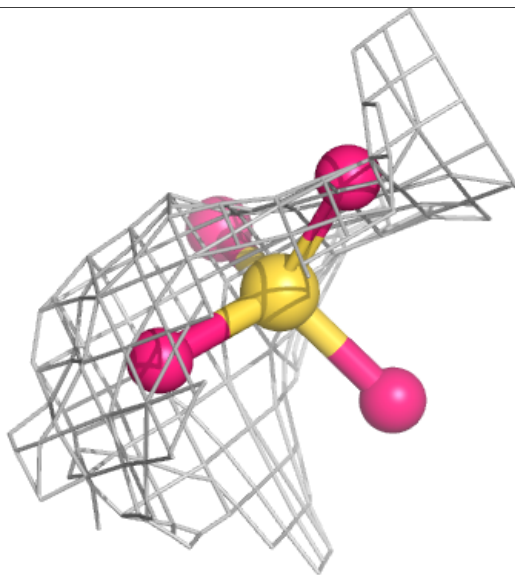
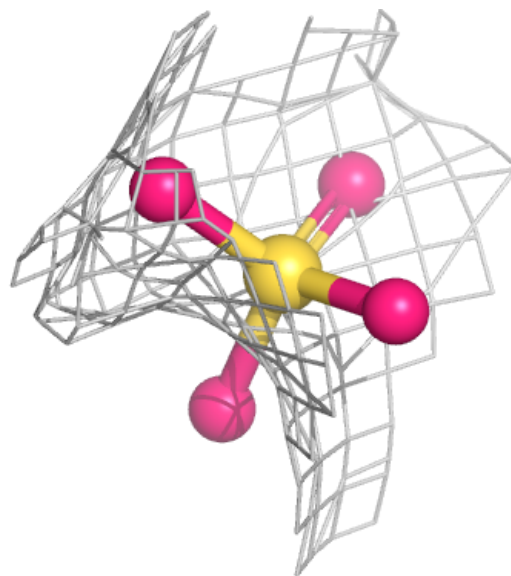
**Electron density around SO4 B 302:**

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



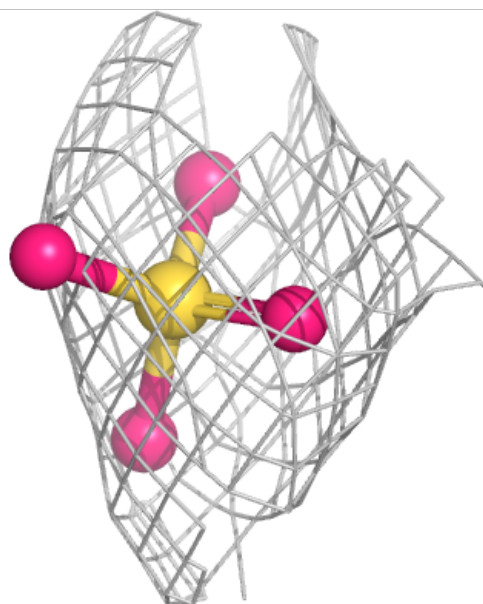
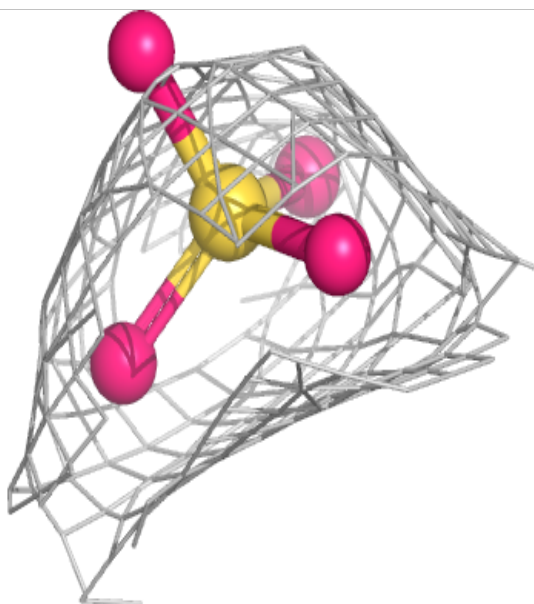
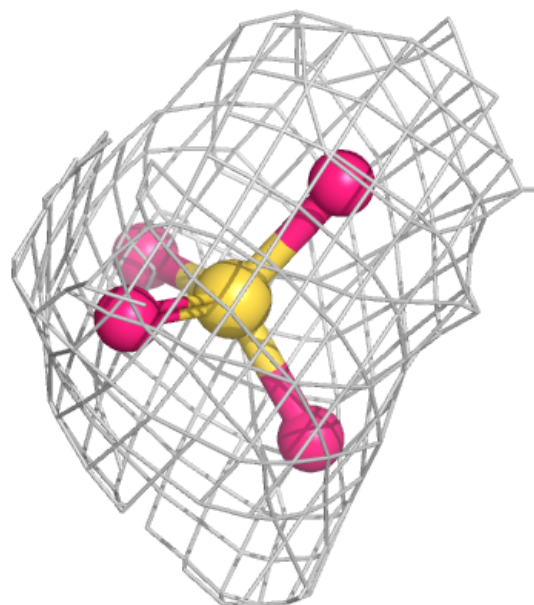
**Electron density around SO4 J 301:**

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



**Electron density around SO4 P 303:**

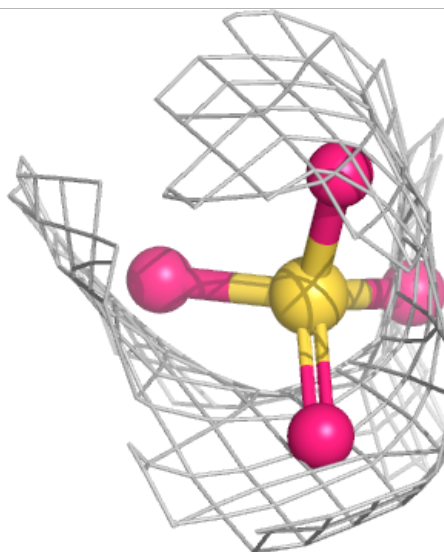
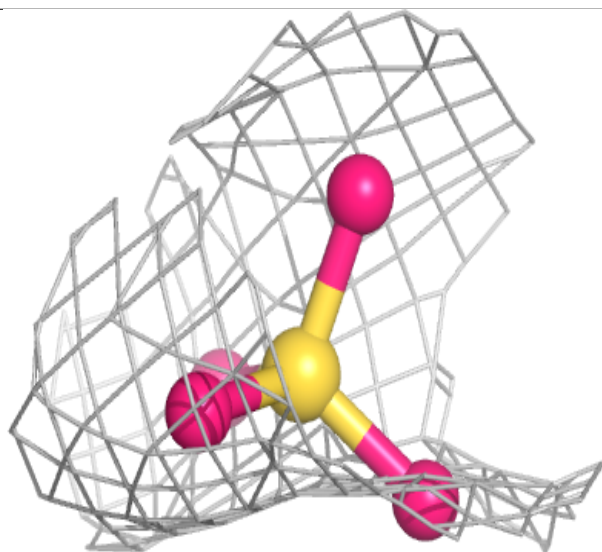
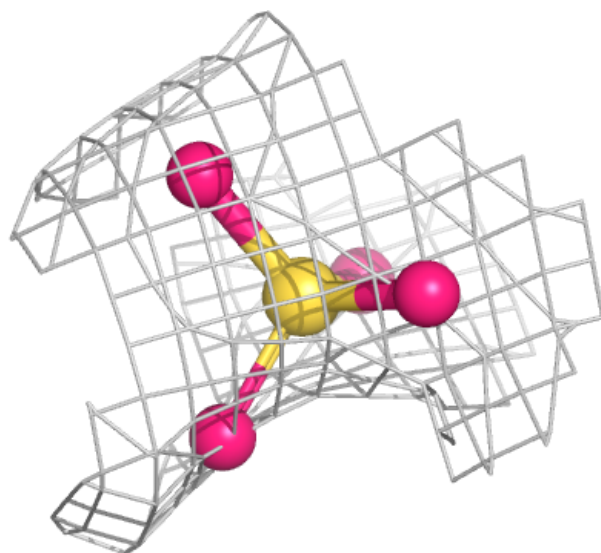
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





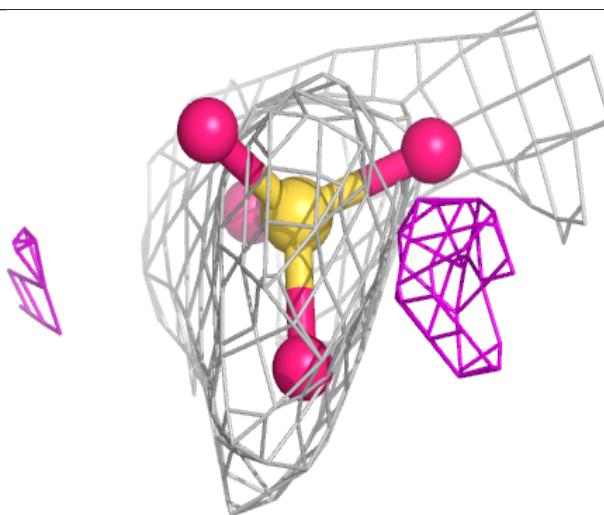
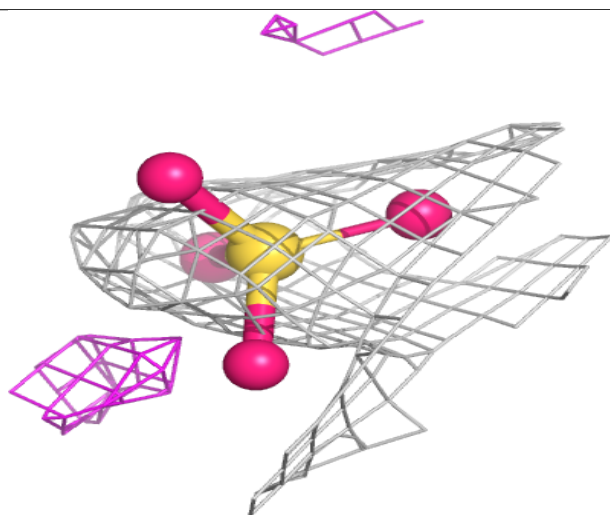
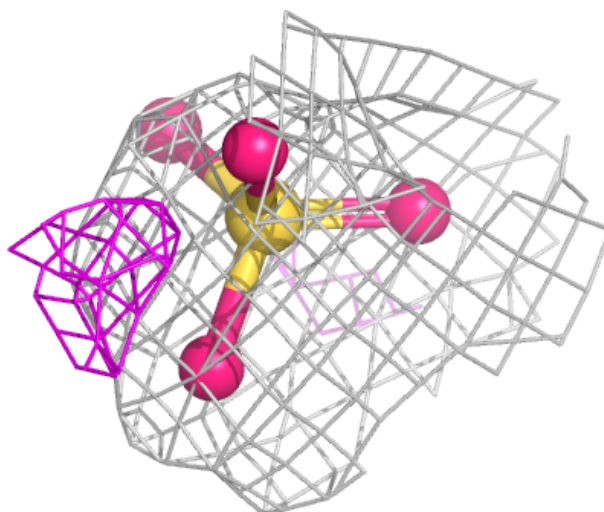
**Electron density around SO4 P 301:**

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



**Electron density around SO4 G 301:**

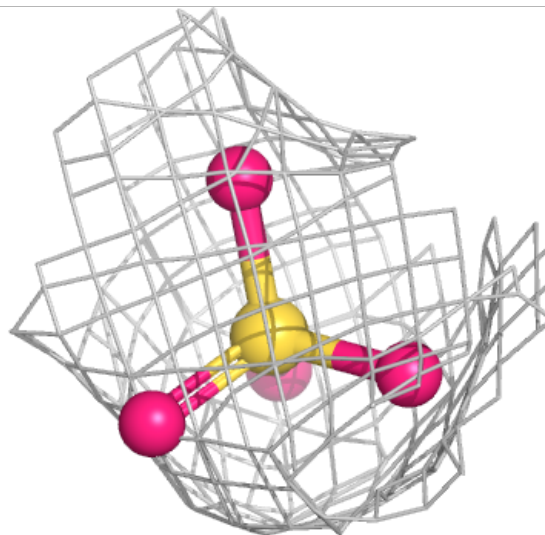
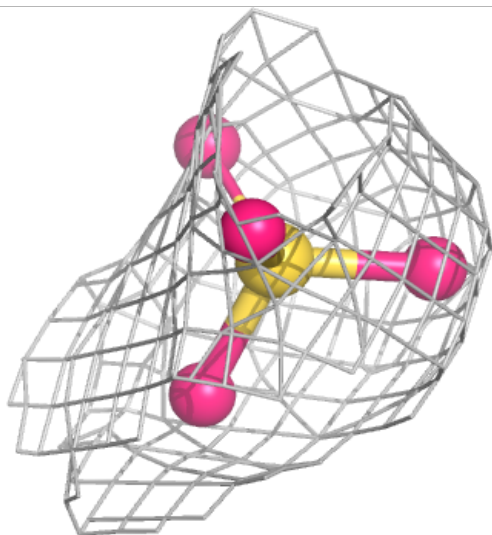
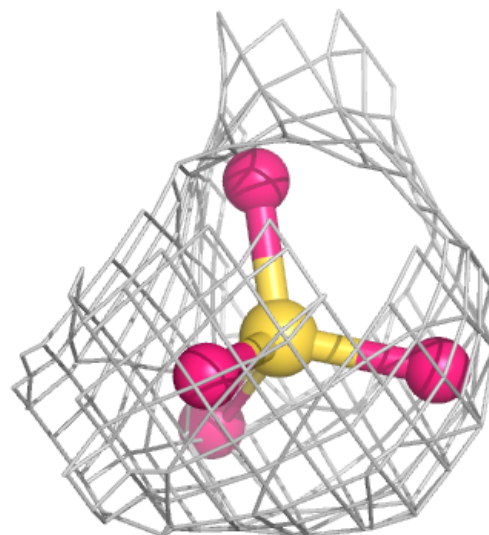
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





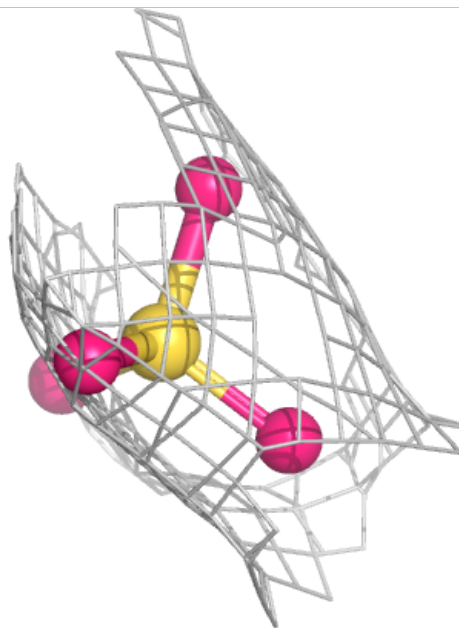
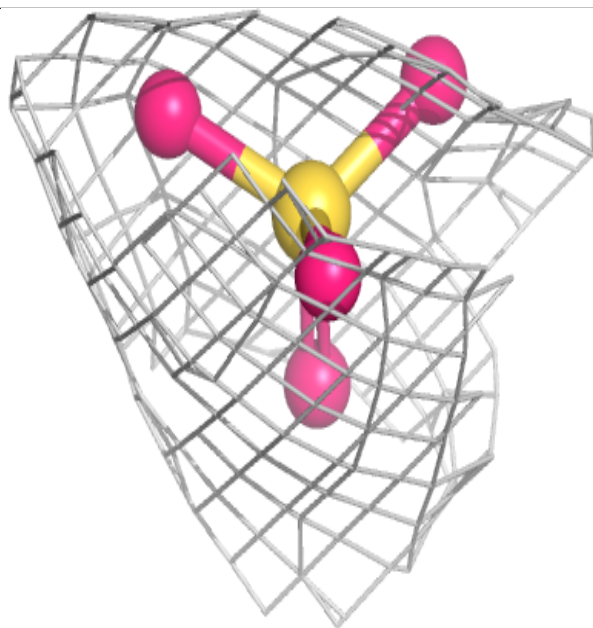
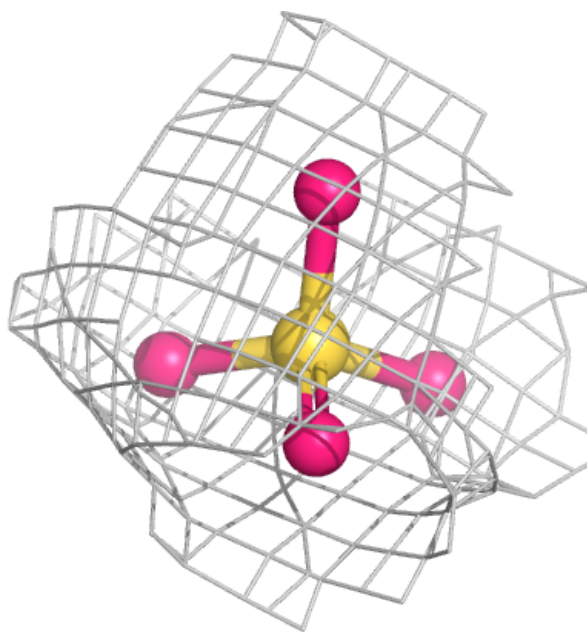
**Electron density around SO4 J 302:**

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



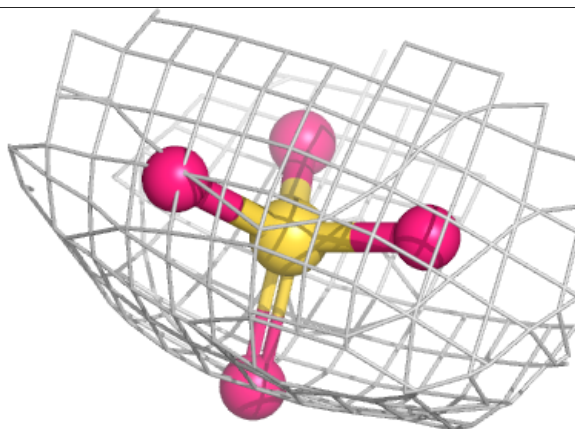
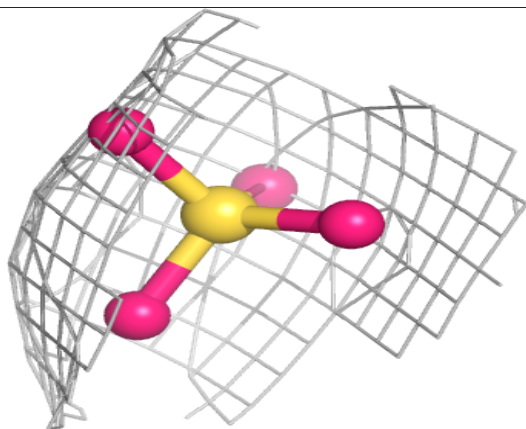
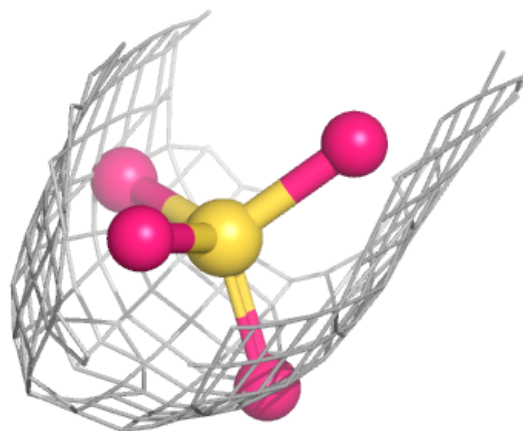
**Electron density around SO4 M 301:**

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



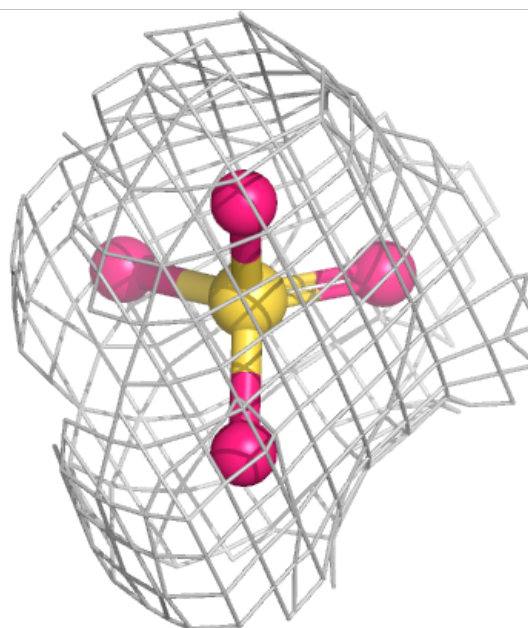
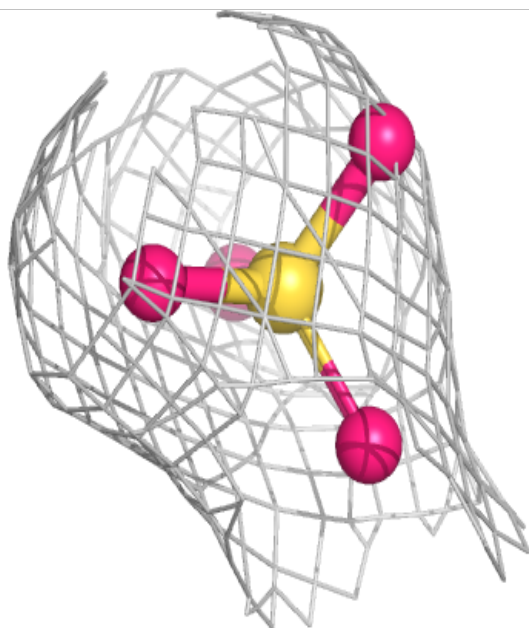
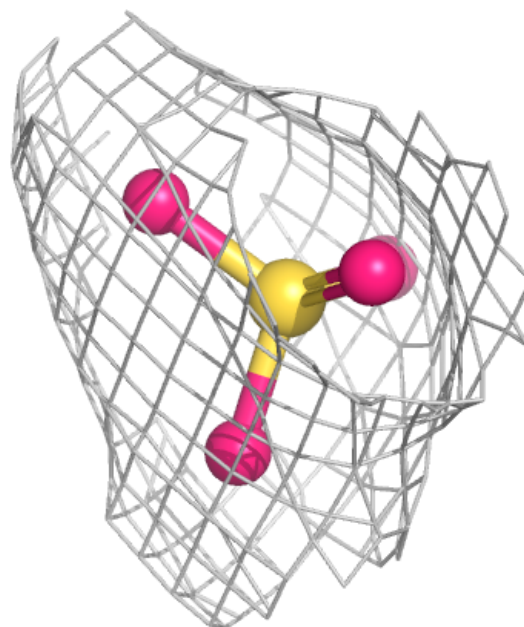
**Electron density around SO4 M 302:**

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



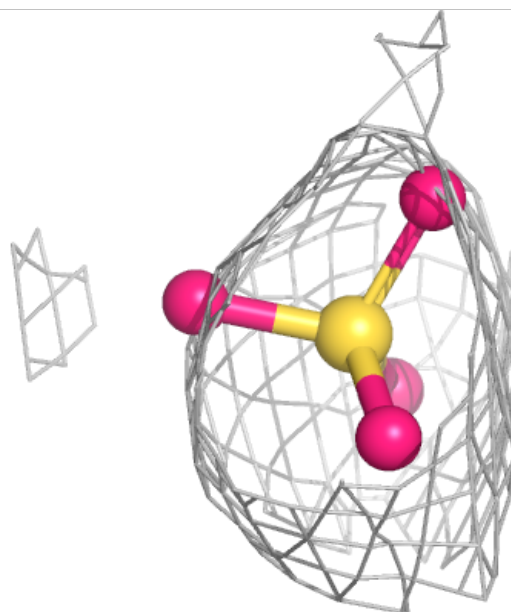
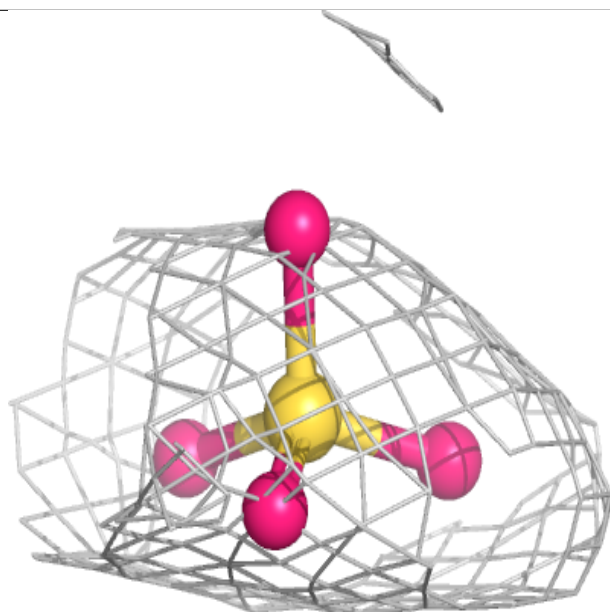
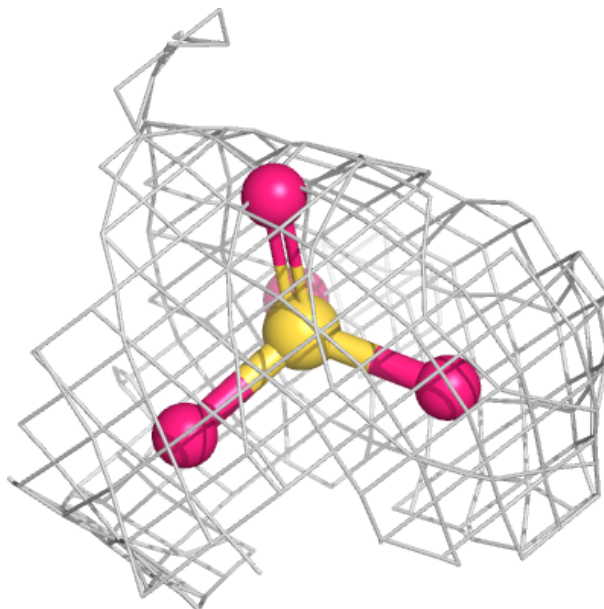
**Electron density around SO4 A 301:**

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



**Electron density around SO4 P 302:**

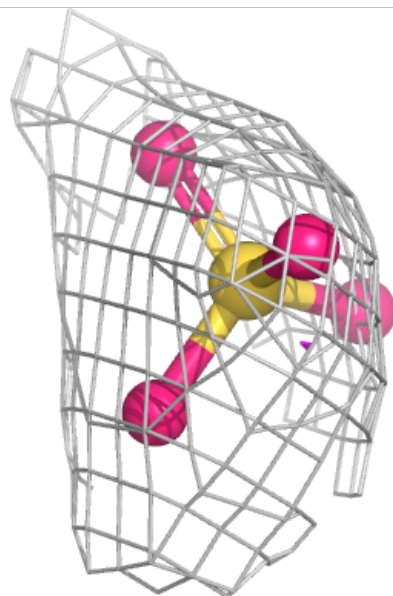
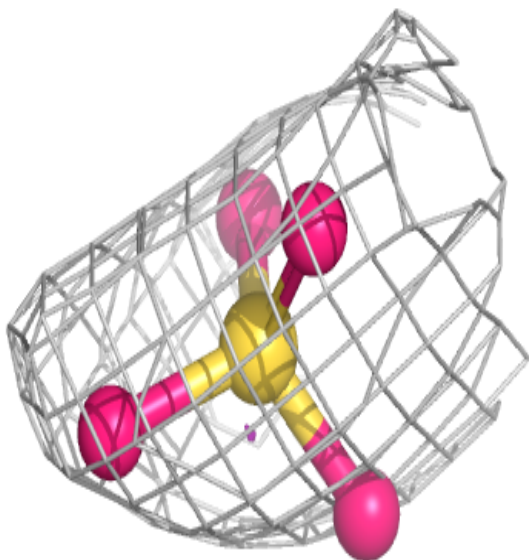
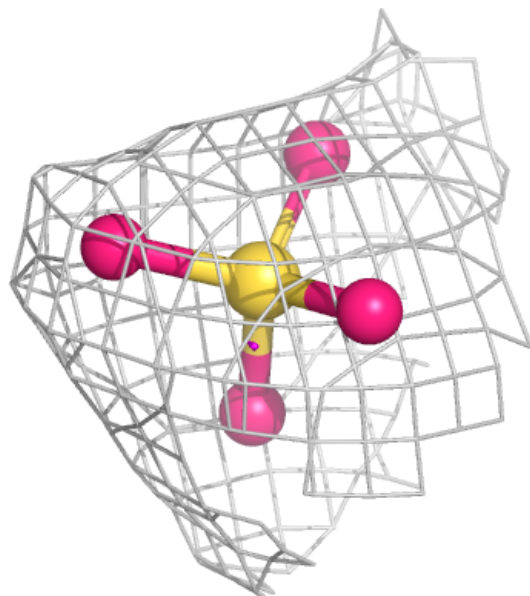
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





**Electron density around SO4 B 301:**

$2mF_o-DF_c$  (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.