



Full wwPDB X-ray Structure Validation Report ⓘ

May 26, 2022 – 04:03 PM EDT

PDB ID : 7U02
Title : Structure of the *C. crescentus* DriD C-domain bound to ssDNA
Authors : Schumacher, M.A.
Deposited on : 2022-02-17
Resolution : 2.48 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 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 : 2.28.1
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) : 2.28.1

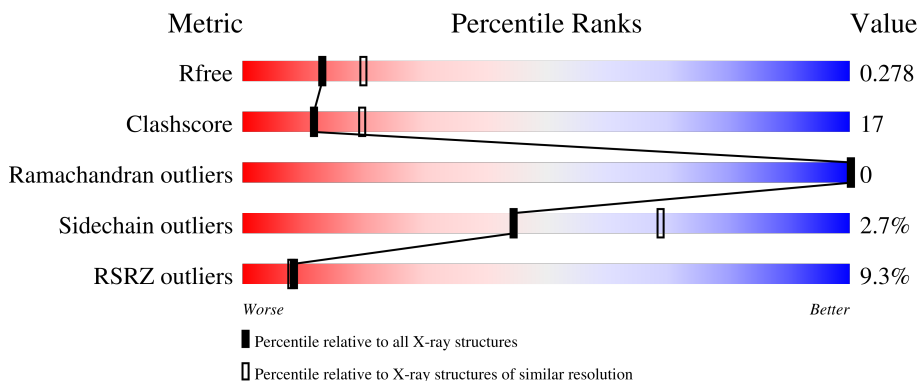
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.48 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	5857 (2.50-2.46)
Clashscore	141614	6594 (2.50-2.46)
Ramachandran outliers	138981	6469 (2.50-2.46)
Sidechain outliers	138945	6471 (2.50-2.46)
RSRZ outliers	127900	5738 (2.50-2.46)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	B	195	 17% 65% 31% ...
1	M	195	 % 70% 29% .
2	H	3	 67% 33%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 3033 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 WYL domain-containing protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	B	193	1478	941	267	266	4	0	0	0
1	M	192	1476	940	265	267	4	0	0	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	133	MET	-	initiating methionine	UNP Q9A999
M	133	MET	-	initiating methionine	UNP Q9A999

- Molecule 2 is a DNA chain called DNA (5'-D(P*AP*CP*G)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
2	H	3	62	29	13	17	3	0	0	0

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O₄S) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	1	Total O S 5 4 1	0	0
3	B	1	Total O S 5 4 1	0	0
3	M	1	Total O S 5 4 1	0	0

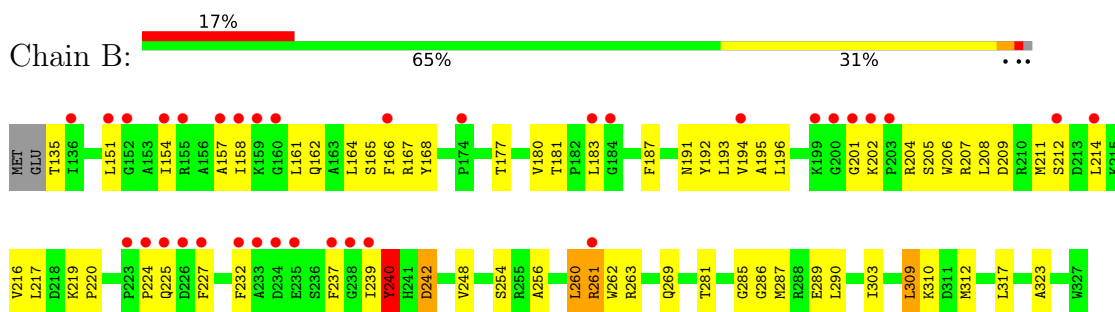
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	M	2	Total O 2 2	0	0

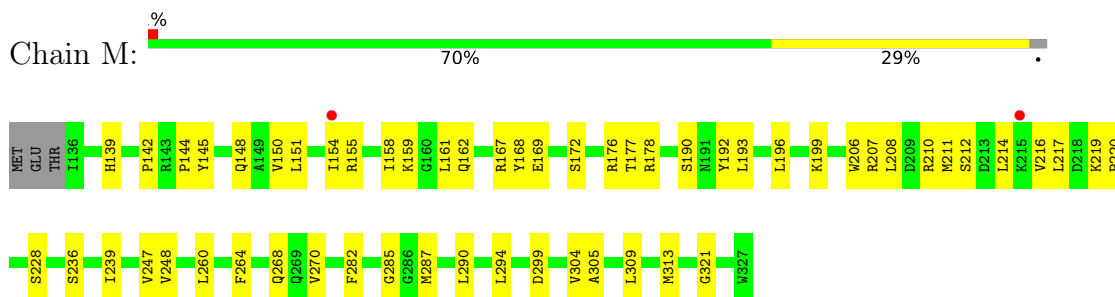
3 Residue-property plots [i](#)

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.

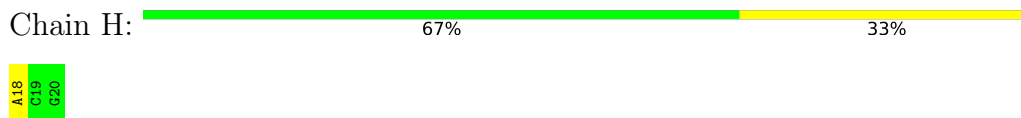
- Molecule 1: WYL domain-containing protein



- Molecule 1: WYL domain-containing protein



- Molecule 2: DNA (5'-D(P*AP*CP*G)-3')



4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	87.40Å 87.40Å 139.60Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	66.54 – 2.48 75.69 – 2.30	Depositor EDS
% Data completeness (in resolution range)	99.5 (66.54-2.48) 97.3 (75.69-2.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.95 (at 2.29Å)	Xtrriage
Refinement program	PHENIX 1.17.1_3660	Depositor
R, R_{free}	0.230 , 0.279 0.231 , 0.278	Depositor DCC
R_{free} test set	1996 reflections (7.32%)	wwPDB-VP
Wilson B-factor (Å ²)	70.0	Xtrriage
Anisotropy	0.372	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 77.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtrriage
Estimated twinning fraction	0.026 for -h,-k,l	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	3033	wwPDB-VP
Average B, all atoms (Å ²)	117.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.10% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: SO4

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# $ Z > 5$	RMSZ	# $ Z > 5$
1	B	0.54	1/1516 (0.1%)	0.80	4/2062 (0.2%)
1	M	0.48	0/1514	0.60	0/2056
2	H	0.73	0/69	0.93	0/104
All	All	0.52	1/3099 (0.0%)	0.71	4/4222 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	261	ARG	CB-CG	5.32	1.66	1.52

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	261	ARG	NE-CZ-NH1	-9.24	115.68	120.30
1	B	261	ARG	CG-CD-NE	-7.90	95.22	111.80
1	B	261	ARG	CA-CB-CG	7.21	129.26	113.40
1	B	240	TYR	CA-CB-CG	-5.86	102.27	113.40

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	B	1478	0	1429	60	0
1	M	1476	0	1437	49	1
2	H	62	0	34	3	0
3	B	10	0	0	1	0
3	M	5	0	0	0	0
4	M	2	0	0	0	0
All	All	3033	0	2900	103	1

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 17.

All (103) 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:M:148:GLN:O	1:M:148:GLN:OE1	1.81	0.97
1:B:232:PHE:O	1:B:240:TYR:OH	1.85	0.95
1:B:167:ARG:HB2	1:B:212:SER:HB2	1.63	0.78
1:M:248:VAL:HG13	1:M:304:VAL:HG13	1.66	0.75
1:B:287:MET:SD	1:B:309:LEU:HG	2.28	0.73
1:M:150:VAL:HG12	1:M:214:LEU:CD2	2.19	0.72
1:B:237:PHE:HB3	1:B:286:GLY:HA3	1.72	0.70
1:B:180:VAL:HG22	1:B:195:ALA:HB1	1.73	0.69
1:M:150:VAL:HG12	1:M:214:LEU:HD22	1.73	0.69
2:H:18:DA:P	1:M:178:ARG:HH22	2.15	0.69
1:M:208:LEU:HA	1:M:211:MET:HE2	1.78	0.66
1:M:176:ARG:HG3	1:M:177:THR:N	2.10	0.66
1:M:162:GLN:NE2	1:M:217:LEU:O	2.26	0.65
1:B:323:ALA:HB2	1:M:299:ASP:HB2	1.81	0.63
1:M:167:ARG:HG3	1:M:212:SER:OG	1.99	0.63
1:B:167:ARG:HG2	1:B:177:THR:HG22	1.81	0.63
1:B:154:ILE:HD11	1:B:214:LEU:HD13	1.82	0.62
1:B:192:TYR:HB3	1:B:205:SER:HB3	1.81	0.61
1:B:289:GLU:N	1:B:289:GLU:OE1	2.33	0.60
1:M:219:LYS:HG2	1:M:220:PRO:HD2	1.84	0.59
1:B:166:PHE:HB3	1:B:214:LEU:HG	1.85	0.58
1:B:217:LEU:HD13	1:B:219:LYS:HD3	1.85	0.58
1:M:248:VAL:HG12	1:M:305:ALA:HB3	1.84	0.58
1:B:183:LEU:HD12	1:B:195:ALA:HA	1.86	0.56
1:B:135:THR:HG22	1:M:155:ARG:HH12	1.69	0.56
1:B:256:ALA:O	1:B:260:LEU:HD12	2.06	0.56
1:M:150:VAL:O	1:M:154:ILE:HG13	2.06	0.55
1:B:196:LEU:HD22	1:B:202:LYS:C	2.27	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:M:148:GLN:HG2	1:M:151:LEU:HD23	1.88	0.55
1:M:247:VAL:HG21	1:M:282:PHE:CE1	2.41	0.55
1:M:199:LYS:HA	1:M:199:LYS:HE2	1.89	0.55
1:B:239:ILE:HG13	1:B:286:GLY:HA2	1.88	0.54
1:B:196:LEU:HB3	1:B:201:GLY:HA3	1.89	0.54
1:B:206:TRP:HB2	1:B:211:MET:CE	2.37	0.54
1:M:148:GLN:OE1	1:M:148:GLN:C	2.47	0.53
1:M:287:MET:CE	1:M:290:LEU:HD23	2.38	0.53
1:B:157:ALA:HB2	1:B:216:VAL:HG13	1.90	0.53
2:H:18:DA:H5 ⁷	1:M:206:TRP:CZ3	2.44	0.52
1:B:206:TRP:HB2	1:B:211:MET:HE2	1.91	0.52
1:B:217:LEU:HB3	1:B:219:LYS:HD2	1.90	0.52
1:B:239:ILE:HB	1:B:242:ASP:OD1	2.10	0.52
1:M:144:PRO:HG2	1:M:190:SER:HA	1.92	0.52
1:B:196:LEU:HD13	1:B:201:GLY:C	2.31	0.51
1:B:290:LEU:HD23	1:B:309:LEU:HD11	1.92	0.51
1:B:191:ASN:HB3	1:B:208:LEU:HD12	1.92	0.51
1:B:208:LEU:HD23	1:B:211:MET:HE3	1.92	0.50
1:B:187:PHE:HB2	1:M:139:HIS:ND1	2.25	0.50
1:B:183:LEU:HD21	1:B:224:PRO:HD2	1.94	0.50
1:B:183:LEU:HB3	1:B:227:PHE:CE2	2.46	0.50
1:B:261:ARG:CZ	1:B:261:ARG:HA	2.42	0.50
1:M:264:PHE:N	1:M:268:GLN:OE1	2.44	0.50
1:M:150:VAL:HG12	1:M:214:LEU:CD1	2.42	0.49
1:M:208:LEU:HA	1:M:211:MET:CE	2.43	0.49
1:M:287:MET:HE2	1:M:309:LEU:HB2	1.93	0.49
1:B:181:THR:HG22	1:B:196:LEU:O	2.12	0.49
1:B:269:GLN:OE1	1:B:281:THR:OG1	2.30	0.49
1:B:193:LEU:HB2	1:B:208:LEU:HD21	1.93	0.49
1:B:164:LEU:HB3	1:B:214:LEU:HD22	1.94	0.48
1:B:262:TRP:HE1	1:M:142:PRO:HB2	1.77	0.48
1:B:217:LEU:C	1:B:219:LYS:H	2.16	0.48
1:B:162:GLN:HA	1:B:220:PRO:HA	1.96	0.47
1:B:208:LEU:HD23	1:B:211:MET:CE	2.45	0.47
1:M:287:MET:HE3	1:M:290:LEU:HD23	1.97	0.46
1:M:247:VAL:HG21	1:M:282:PHE:CZ	2.50	0.46
1:B:196:LEU:HD13	1:B:201:GLY:CA	2.46	0.46
1:M:294:LEU:HB2	1:M:313:MET:HE1	1.98	0.46
1:B:239:ILE:CG1	1:B:286:GLY:HA2	2.46	0.45
1:M:248:VAL:CG1	1:M:305:ALA:HB3	2.47	0.45
1:B:254:SER:OG	3:B:401:SO4:S	2.62	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:158:ILE:O	1:B:161:LEU:HD23	2.18	0.44
1:B:194:VAL:HG22	1:B:205:SER:OG	2.17	0.44
1:M:216:VAL:O	1:M:216:VAL:HG23	2.17	0.44
1:B:237:PHE:HB2	1:B:285:GLY:O	2.17	0.44
1:B:191:ASN:CB	1:B:208:LEU:HD12	2.48	0.44
1:B:263:ARG:HB2	1:M:145:TYR:CD2	2.53	0.44
1:B:239:ILE:HG22	1:B:240:TYR:N	2.32	0.43
1:M:192:TYR:CE1	1:M:239:ILE:HG13	2.53	0.43
1:M:158:ILE:O	1:M:161:LEU:HD12	2.18	0.43
1:B:303:ILE:O	1:B:310:LYS:HE2	2.19	0.43
1:B:193:LEU:O	1:B:205:SER:HA	2.19	0.42
1:B:151:LEU:HD22	1:B:151:LEU:O	2.20	0.41
1:B:248:VAL:HG13	1:B:281:THR:HG22	2.02	0.41
1:M:148:GLN:HA	1:M:151:LEU:HB3	2.01	0.41
1:M:236:SER:HA	1:M:285:GLY:H	1.86	0.41
1:M:247:VAL:HG11	1:M:287:MET:HE1	2.03	0.41
1:B:196:LEU:HD13	1:B:201:GLY:HA3	2.02	0.41
1:B:207:ARG:HB3	1:B:209:ASP:OD1	2.19	0.41
1:M:260:LEU:HD23	1:M:270:VAL:HG21	2.02	0.41
1:B:165:SER:O	1:B:214:LEU:HA	2.19	0.41
1:B:237:PHE:CB	1:B:286:GLY:HA3	2.47	0.41
1:M:287:MET:HE1	1:M:290:LEU:HD23	2.02	0.41
1:M:219:LYS:HG2	1:M:220:PRO:CD	2.49	0.41
1:B:202:LYS:HD3	1:B:204:ARG:HD3	2.03	0.41
1:B:217:LEU:HB3	1:B:219:LYS:CD	2.52	0.40
1:M:168:TYR:OH	1:M:207:ARG:N	2.50	0.40
1:M:193:LEU:O	1:M:206:TRP:HD1	2.05	0.40
1:M:247:VAL:HG22	1:M:282:PHE:O	2.21	0.40
1:M:169:GLU:N	1:M:210:ARG:O	2.37	0.40
1:B:181:THR:CG2	1:B:196:LEU:HB2	2.52	0.40
1:B:317:LEU:HD13	1:M:321:GLY:HA2	2.04	0.40
2:H:18:DA:OP1	1:M:172:SER:N	2.49	0.40
1:M:150:VAL:CG1	1:M:214:LEU:HD22	2.49	0.40
1:M:196:LEU:HD23	1:M:196:LEU:HA	1.90	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:M:159:LYS:NZ	1:M:299:ASP:OD1[2_665]	1.84	0.36

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	B	191/195 (98%)	172 (90%)	19 (10%)	0	100	100
1	M	190/195 (97%)	182 (96%)	8 (4%)	0	100	100
All	All	381/390 (98%)	354 (93%)	27 (7%)	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	B	145/156 (93%)	138 (95%)	7 (5%)	25	45
1	M	147/156 (94%)	146 (99%)	1 (1%)	84	93
All	All	292/312 (94%)	284 (97%)	8 (3%)	44	69

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	168	TYR
1	B	225	GLN
1	B	240	TYR
1	B	242	ASP
1	B	260	LEU
1	B	309	LEU
1	B	312	MET
1	M	228	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	324	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](#)

3 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	$\# Z > 2$
3	SO4	B	401	-	4,4,4	0.16	0	6,6,6	0.11	0
3	SO4	B	402	-	4,4,4	0.13	0	6,6,6	0.15	0
3	SO4	M	401	-	4,4,4	0.19	0	6,6,6	0.13	0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

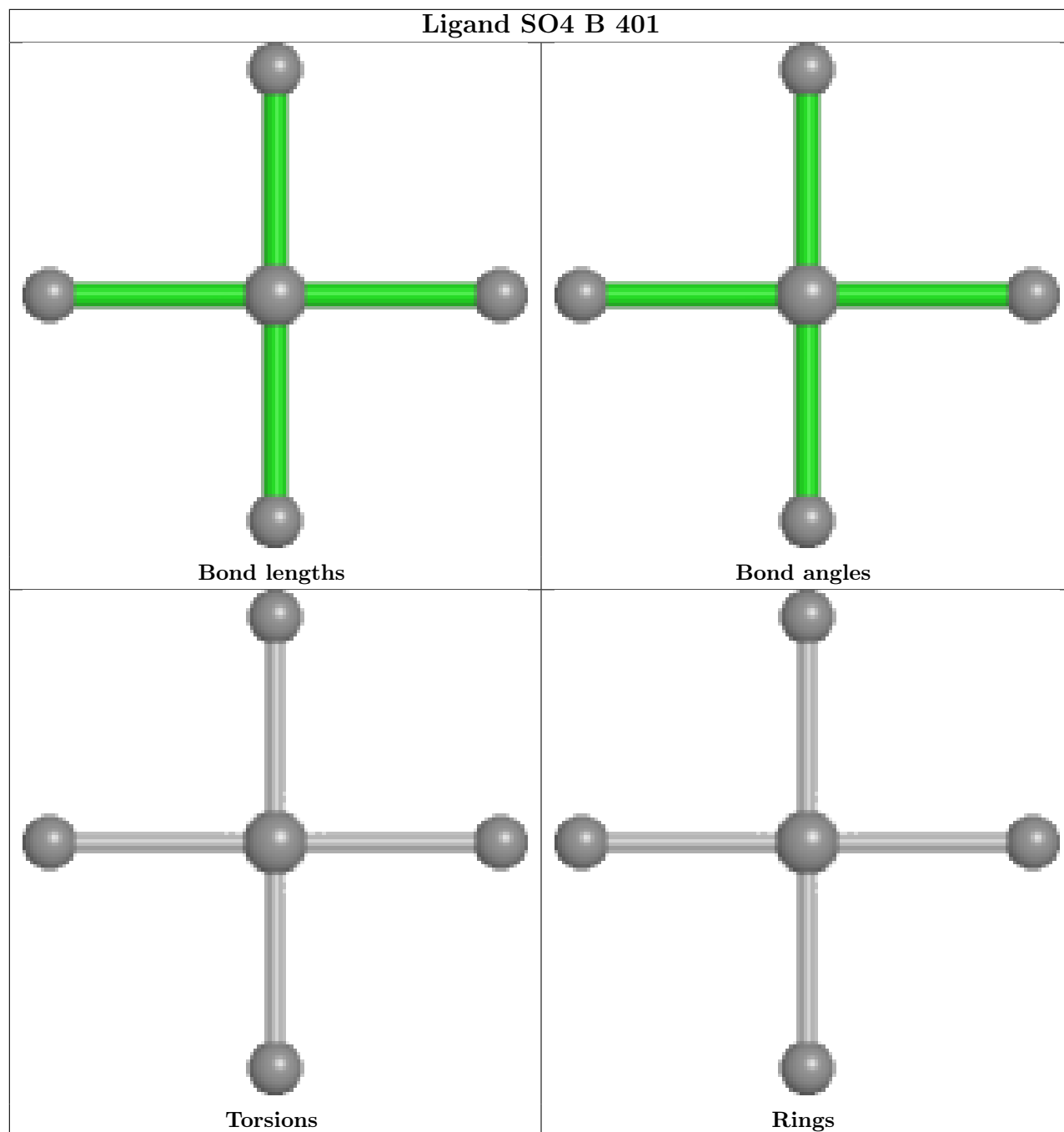
There are no torsion outliers.

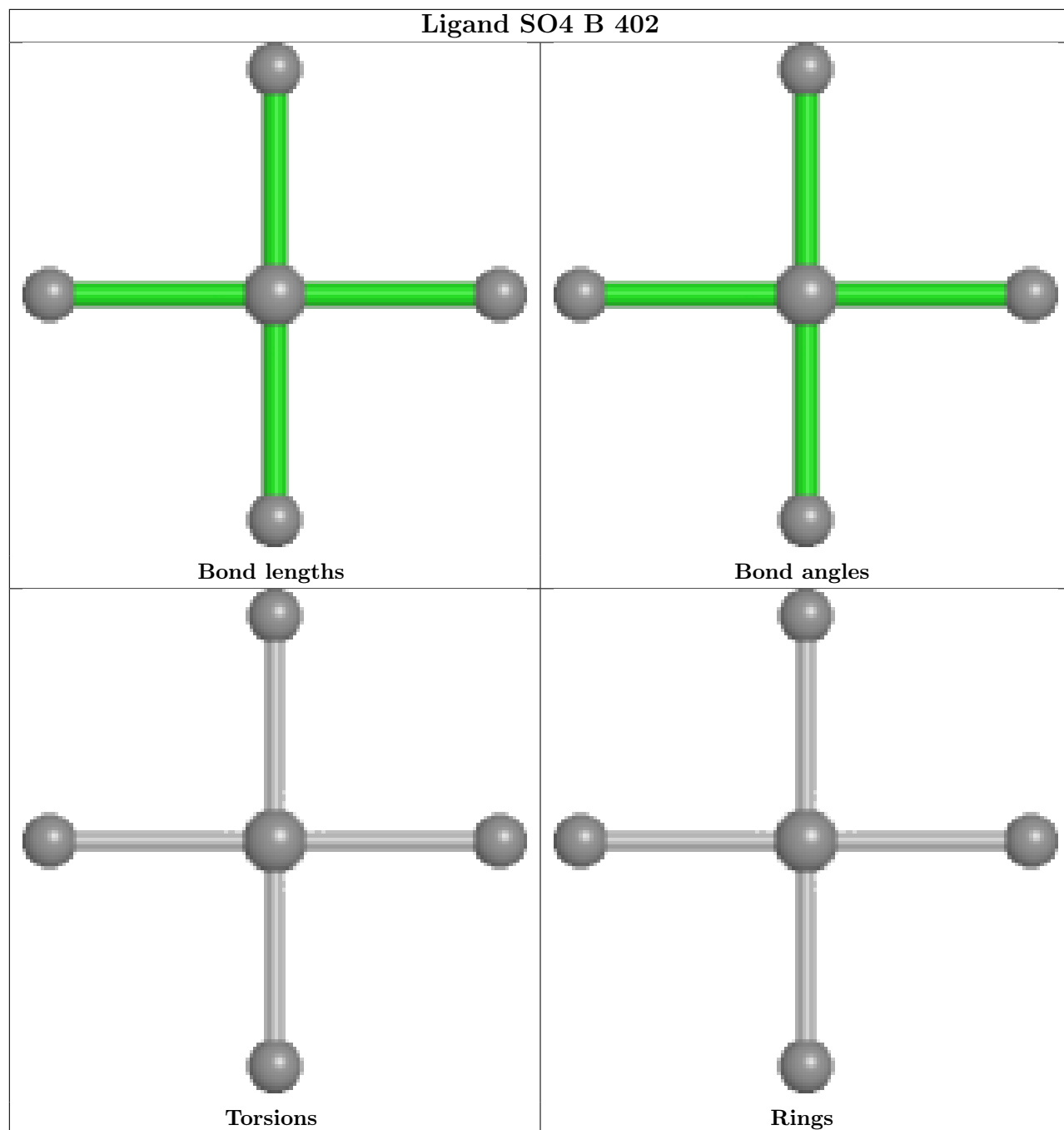
There are no ring outliers.

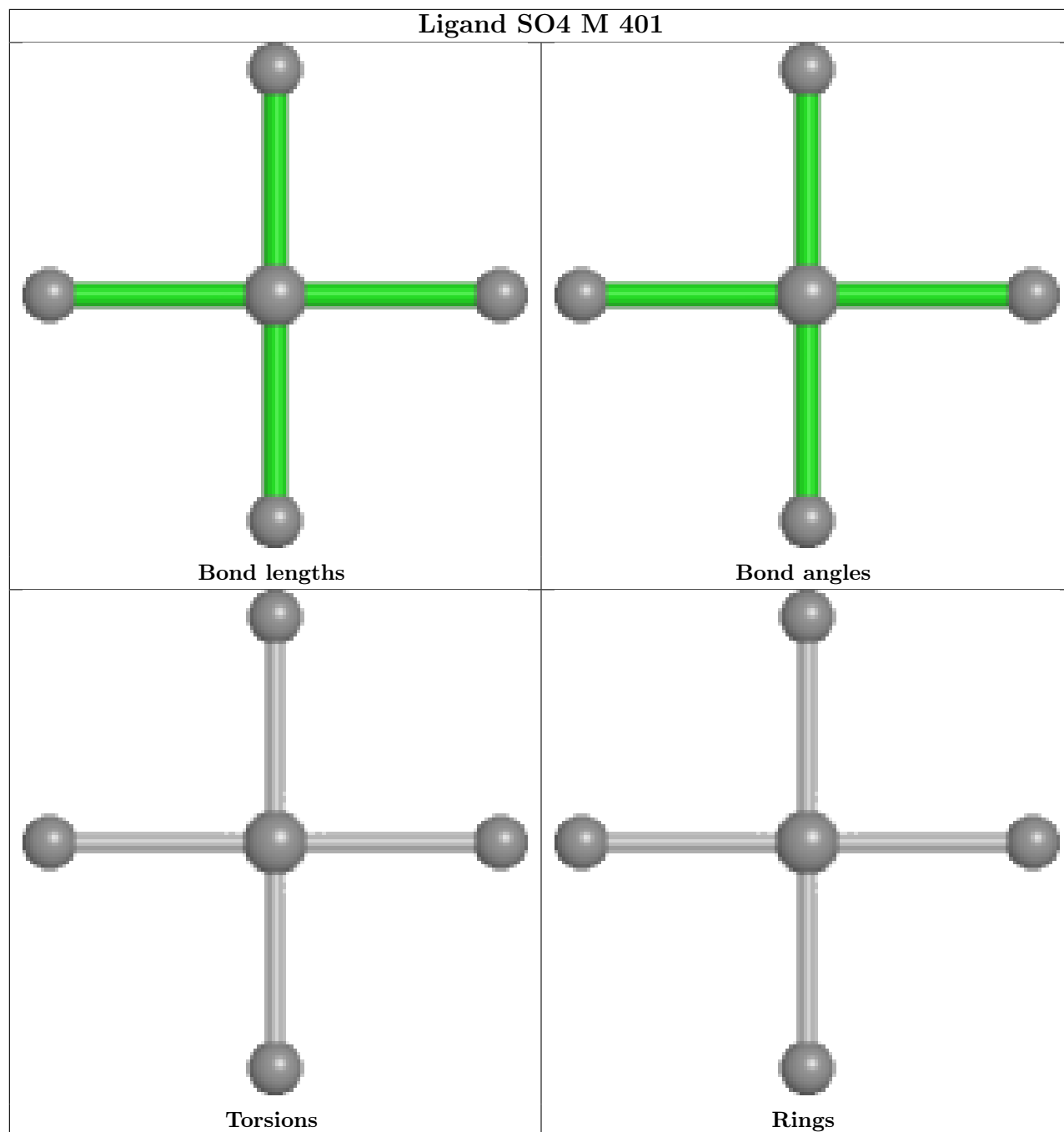
1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	401	SO4	1	0

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







5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	B	193/195 (98%)	1.03	34 (17%) 1 1	65, 129, 208, 234	0
1	M	192/195 (98%)	0.12	2 (1%) 82 84	61, 94, 159, 181	0
2	H	3/3 (100%)	-0.21	0 100 100	95, 95, 117, 121	0
All	All	388/393 (98%)	0.57	36 (9%) 8 8	61, 107, 196, 234	0

All (36) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	184	GLY	12.3
1	B	214	LEU	10.2
1	B	183	LEU	7.8
1	B	238	GLY	7.6
1	B	239	ILE	6.2
1	B	154	ILE	6.0
1	B	237	PHE	5.3
1	B	201	GLY	5.3
1	B	226	ASP	5.2
1	B	151	LEU	4.6
1	B	159	LYS	4.0
1	B	235	GLU	3.9
1	B	174	PRO	3.7
1	B	200	GLY	3.6
1	B	227	PHE	3.5
1	B	224	PRO	3.3
1	B	158	ILE	3.3
1	B	166	PHE	3.2
1	B	223	PRO	3.2
1	B	233	ALA	3.0
1	B	225	GLN	3.0
1	B	136	ILE	2.9
1	B	234	ASP	2.9

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Mol	Chain	Res	Type	RSRZ
1	M	215	LYS	2.8
1	B	261	ARG	2.7
1	B	152	GLY	2.7
1	B	160	GLY	2.7
1	B	199	LYS	2.6
1	B	155	ARG	2.4
1	M	154	ILE	2.3
1	B	202	LYS	2.3
1	B	157	ALA	2.3
1	B	203	PRO	2.2
1	B	194	VAL	2.2
1	B	232	PHE	2.2
1	B	212	SER	2.1

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

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

6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

6.4 Ligands [i](#)

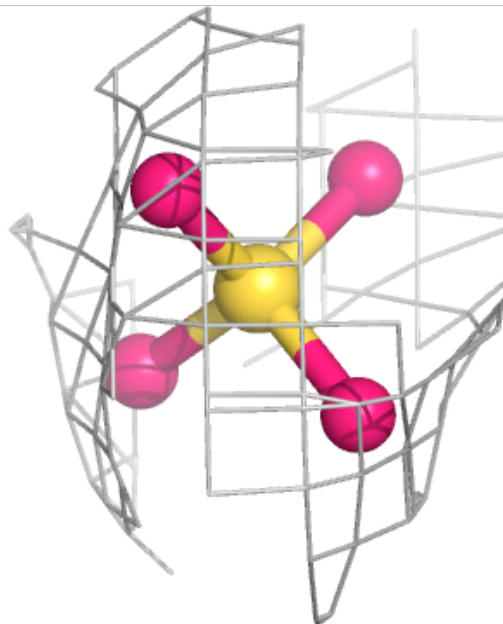
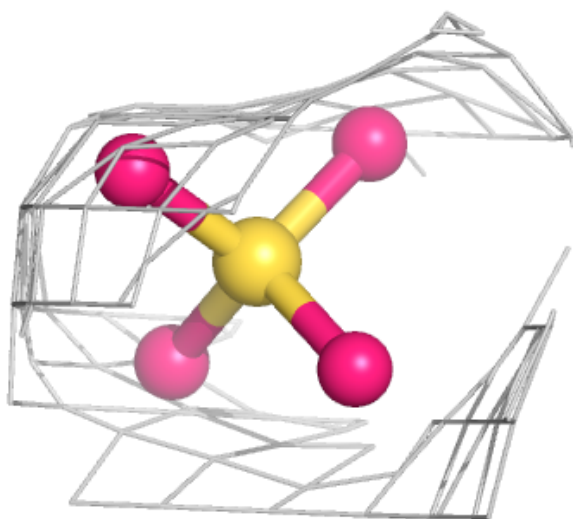
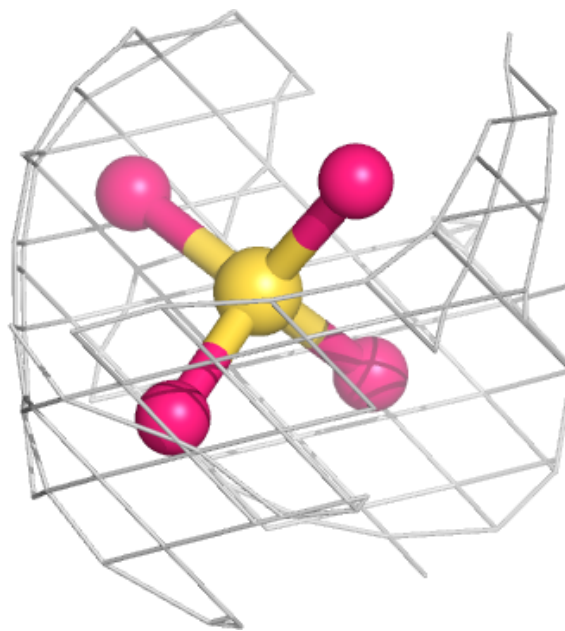
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	SO4	B	402	5/5	0.68	0.17	198,200,204,208	0
3	SO4	M	401	5/5	0.75	0.27	164,164,167,175	0
3	SO4	B	401	5/5	0.91	0.15	158,159,161,164	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

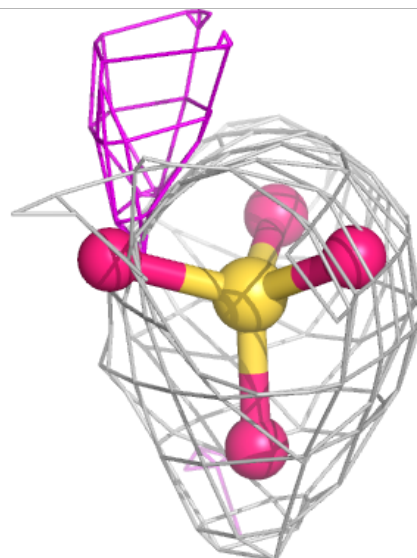
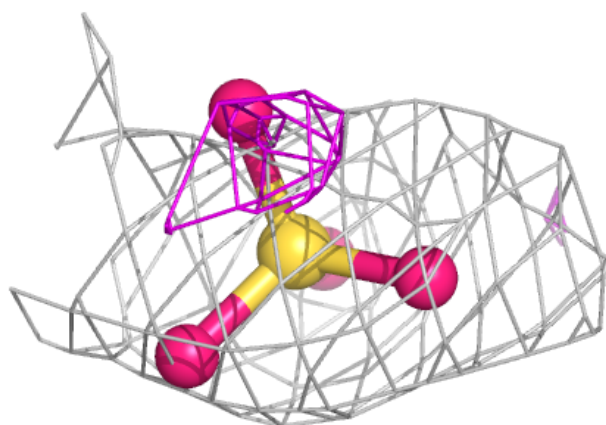
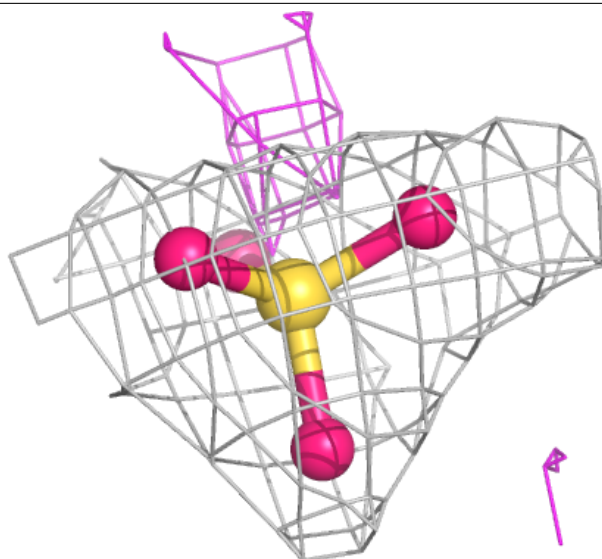
Electron density around SO4 B 402:

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



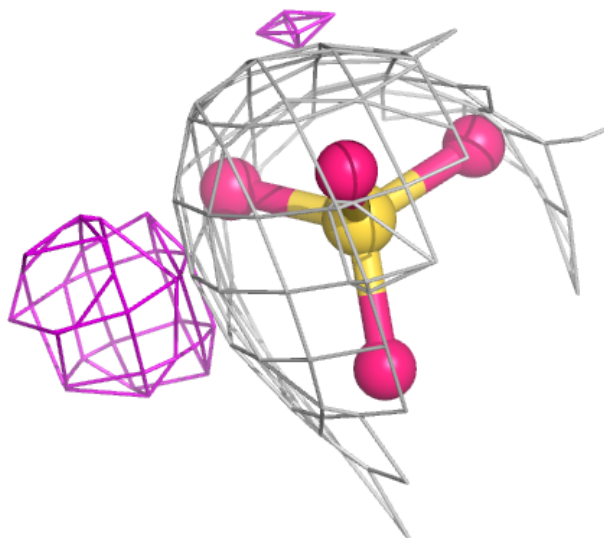
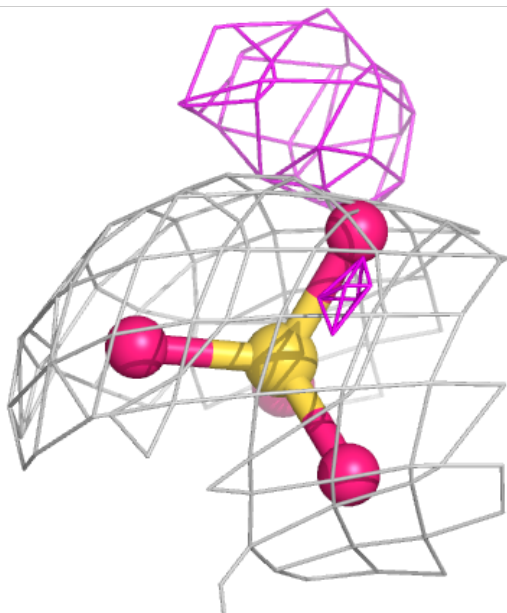
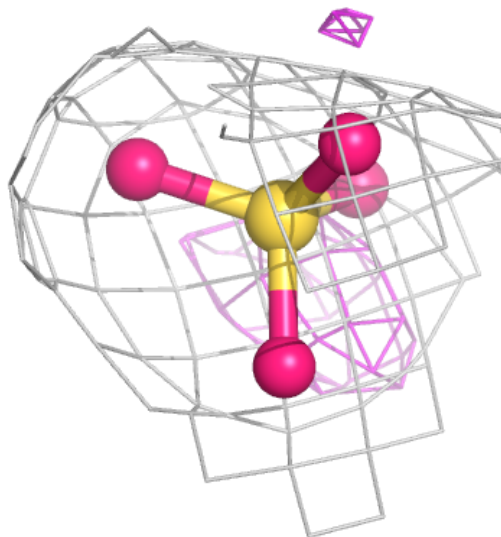
Electron density around SO4 M 401:

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



Electron density around SO4 B 401:

$2mF_o-DF_c$ (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.