



Full wwPDB X-ray Structure Validation Report ⓘ

Jul 7, 2025 – 04:11 PM EDT

PDB ID : 9NY8 / pdb_00009ny8
Title : Crystal structure of the ribose operon repressor, RbsR, bound to ribose operon
Authors : Wells, M.L.; Lu, C.; Sultanov, D.; Weber, K.C.; Gong, Z.; Glasgow, A.
Deposited on : 2025-03-26
Resolution : 2.10 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 : 4-5-2 with Phenix2.0rc1
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0rc1
EDS : 3.0
buster-report : 1.1.7 (2018)
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4 : 9.0.006 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.44

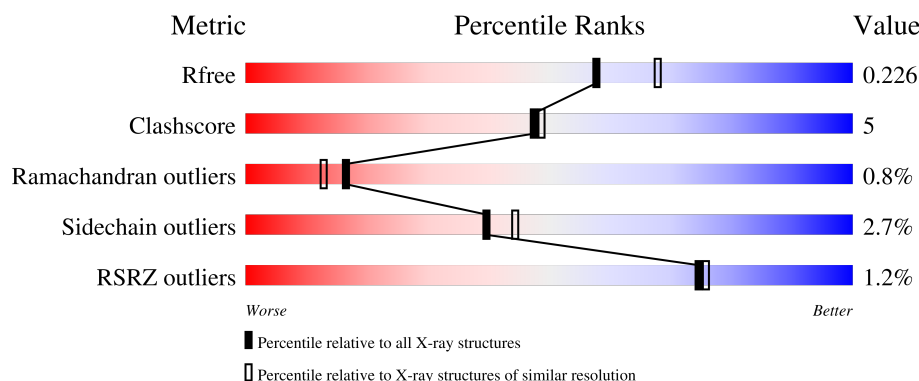
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	164625	6234 (2.10-2.10)
Clashscore	180529	6893 (2.10-2.10)
Ramachandran outliers	177936	6839 (2.10-2.10)
Sidechain outliers	177891	6840 (2.10-2.10)
RSRZ outliers	164620	6234 (2.10-2.10)

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

Mol	Chain	Length	Quality of chain
1	A	330	<div> <div>2%</div> <div> <div></div> <div>87%</div> <div>12%</div> <div>.</div> </div> </div>
1	B	330	<div> <div>%</div> <div> <div></div> <div>85%</div> <div>13%</div> <div>.</div> </div> </div>
2	C	30	<div> <div></div> <div> <div>50%</div> <div>40%</div> <div>10%</div> </div> </div>
3	D	30	<div> <div></div> <div> <div>47%</div> <div>53%</div> </div> </div>

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 6439 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 Ribose operon repressor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	330	Total	C	N	O	S	0	0	0
			2563	1605	448	494	16			
1	B	330	Total	C	N	O	S	0	1	0
			2571	1609	450	496	16			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	SER	-	expression tag	UNP P0ACQ0
B	1	SER	-	expression tag	UNP P0ACQ0

- Molecule 2 is a DNA chain called ribose operon.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	30	Total	C	N	O	P	0	0	0
			626	295	119	182	30			

- Molecule 3 is a DNA chain called ribose operon.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	D	30	Total	C	N	O	P	0	0	0
			606	288	108	180	30			

- Molecule 4 is SULFATE ION (CCD ID: SO4) (formula: O₄S) (labeled as "Ligand of Interest" by depositor).



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

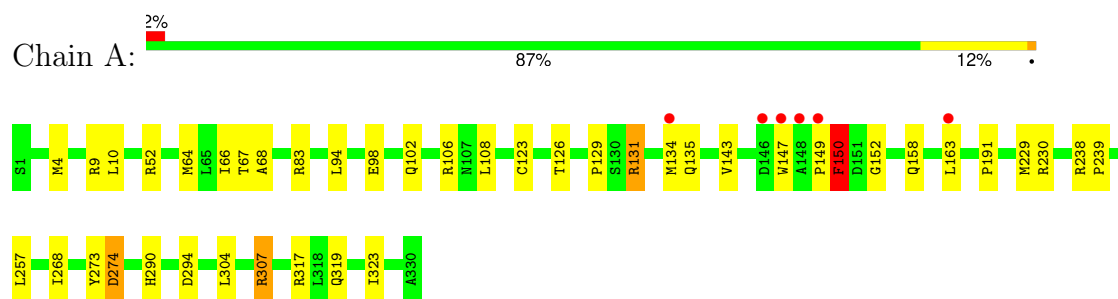
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	24	Total	O	0	0
			24	24		
5	B	29	Total	O	0	0
			29	29		
5	C	3	Total	O	0	0
			3	3		
5	D	2	Total	O	0	0
			2	2		

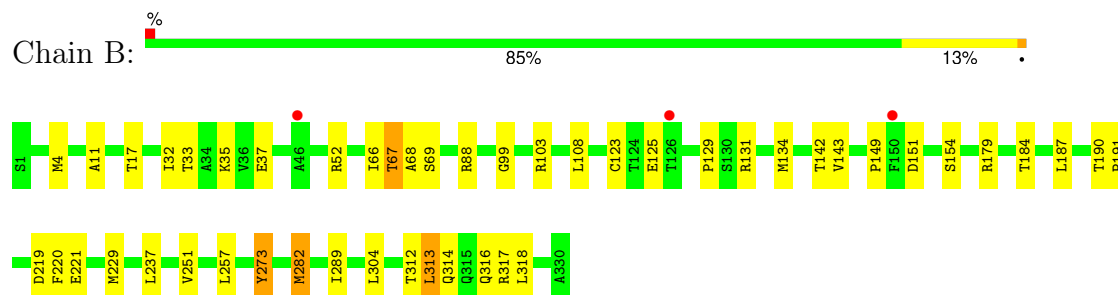
3 Residue-property plots [i](#)

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: Ribose operon repressor



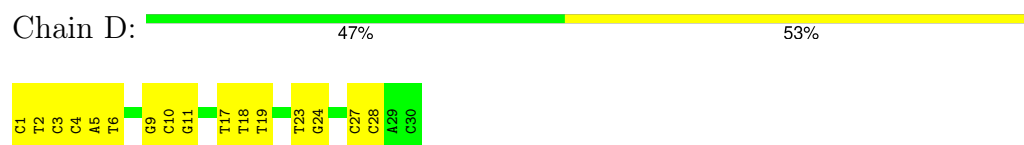
• Molecule 1: Ribose operon repressor



• Molecule 2: ribose operon



• Molecule 3: ribose operon



4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	125.53Å 125.53Å 122.31Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	43.84 – 2.10 43.84 – 2.10	Depositor EDS
% Data completeness (in resolution range)	73.5 (43.84-2.10) 73.5 (43.84-2.10)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.38 (at 2.10Å)	Xtriage
Refinement program	REFMAC 5.8	Depositor
R, R_{free}	0.192 , 0.223 0.196 , 0.226	Depositor DCC
R_{free} test set	3168 reflections (4.85%)	wwPDB-VP
Wilson B-factor (Å ²)	53.1	Xtriage
Anisotropy	0.135	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 52.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.029 for -h,-k,l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	6439	wwPDB-VP
Average B, all atoms (Å ²)	75.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.08% 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

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	0.51	1/2607 (0.0%)	0.83	1/3536 (0.0%)
1	B	0.53	2/2615 (0.1%)	0.85	1/3547 (0.0%)
2	C	0.28	0/703	1.02	3/1084 (0.3%)
3	D	0.29	0/677	0.97	1/1038 (0.1%)
All	All	0.48	3/6602 (0.0%)	0.88	6/9205 (0.1%)

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	A	0	3
1	B	0	2
All	All	0	5

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	4	MET	SD-CE	-6.13	1.64	1.79
1	A	4	MET	SD-CE	-6.11	1.64	1.79
1	B	282	MET	SD-CE	-5.76	1.65	1.79

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	20	DC	O3'-P-O5'	-8.60	91.09	104.00
1	B	273	TYR	N-CA-C	7.21	121.95	112.72
1	A	150	PHE	CA-CB-CG	6.56	120.36	113.80
2	C	27	DG	O3'-P-O5'	-6.45	94.33	104.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	9	DG	O3'-P-O5'	-6.13	94.80	104.00
2	C	28	DG	O3'-P-O5'	-5.80	95.30	104.00

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	131	ARG	Sidechain
1	A	317	ARG	Sidechain
1	A	52	ARG	Sidechain
1	B	103	ARG	Sidechain
1	B	88	ARG	Sidechain

5.2 Too-close contacts

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	2563	0	2569	24	0
1	B	2571	0	2574	22	0
2	C	626	0	338	10	0
3	D	606	0	337	12	0
4	A	5	0	0	0	0
4	B	10	0	0	0	0
5	A	24	0	0	1	1
5	B	29	0	0	0	0
5	C	3	0	0	0	0
5	D	2	0	0	0	0
All	All	6439	0	5818	67	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 5.

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:B:273:TYR:O	1:B:289:ILE:O	1.61	1.17

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:52:ARG:NH1	2:C:18:DT:OP1	1.99	0.94
1:A:9:ARG:NH2	5:A:501:HOH:O	2.08	0.85
1:A:66:ILE:HD12	1:A:68:ALA:HB2	1.67	0.75
2:C:23:DT:H2''	2:C:24:DG:C8	2.26	0.70
3:D:23:DT:H2''	3:D:24:DG:C8	2.27	0.69
1:B:33:THR:O	1:B:37:GLU:HG2	1.94	0.67
3:D:1:DC:O2	3:D:1:DC:H2'	1.96	0.66
1:A:147:TRP:CD1	1:A:191:PRO:HG3	2.32	0.64
1:A:102:GLN:HE21	1:A:106:ARG:HD2	1.64	0.62
1:B:184:THR:HG21	1:B:187:LEU:HD21	1.82	0.61
1:B:66:ILE:HD12	1:B:68:ALA:HB2	1.83	0.60
1:B:229:MET:HE2	1:B:257:LEU:HD21	1.83	0.59
1:A:108:LEU:HD21	1:A:134:MET:HE3	1.86	0.58
2:C:10:DC:H2'	2:C:11:DG:C8	2.40	0.57
1:A:229:MET:HE2	1:A:257:LEU:HD21	1.87	0.57
1:A:83:ARG:NH2	1:A:294:ASP:OD1	2.25	0.55
1:A:307:ARG:HH11	1:A:307:ARG:HB2	1.72	0.55
3:D:4:DC:H2''	3:D:5:DA:OP2	2.06	0.54
1:B:143:VAL:HG23	1:B:304:LEU:HD13	1.89	0.54
1:A:238:ARG:HD3	1:A:239:PRO:HD2	1.89	0.54
1:B:190:THR:N	1:B:191:PRO:HD2	2.23	0.54
3:D:10:DC:H2''	3:D:11:DG:C8	2.43	0.53
3:D:17:DT:H2''	3:D:18:DT:O5'	2.09	0.53
1:B:313:LEU:HD13	1:B:314:GLN:N	2.25	0.52
3:D:2:DT:H1'	3:D:3:DC:H5'	1.91	0.51
1:A:134:MET:HE2	1:A:134:MET:HA	1.91	0.51
1:A:108:LEU:CD2	1:A:134:MET:HE3	2.41	0.50
2:C:18:DT:H2''	2:C:19:DT:H5'	1.94	0.50
2:C:26:DT:H2''	2:C:27:DG:OP2	2.12	0.50
2:C:29:DA:H2''	2:C:30:DG:O5'	2.13	0.49
3:D:18:DT:H2'	3:D:19:DT:H72	1.94	0.49
2:C:17:DT:H2'	2:C:18:DT:C6	2.48	0.49
2:C:20:DC:H2''	2:C:21:DG:C8	2.48	0.48
2:C:27:DG:H2''	2:C:28:DG:OP2	2.12	0.48
1:A:143:VAL:HG23	1:A:304:LEU:HD13	1.97	0.47
1:A:108:LEU:CD2	1:A:129:PRO:HB3	2.45	0.46
3:D:5:DA:H2''	3:D:6:DT:OP2	2.16	0.46
3:D:17:DT:C2'	3:D:18:DT:O5'	2.64	0.46
1:B:134:MET:HA	1:B:134:MET:HE2	1.98	0.46
2:C:25:DA:C2'	2:C:26:DT:H71	2.46	0.45
1:A:290:HIS:HB3	1:A:323:ILE:HG13	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:251:VAL:HG22	1:B:282:MET:HE2	1.98	0.45
1:B:11:ALA:O	1:B:35:LYS:HD2	2.16	0.45
1:B:179:ARG:HD2	1:B:237:LEU:HG	1.97	0.45
3:D:3:DC:H2''	3:D:4:DC:C5	2.52	0.44
1:A:163:LEU:HD22	1:A:163:LEU:N	2.32	0.44
1:B:17:THR:HG21	1:B:32:ILE:HG21	1.99	0.44
3:D:5:DA:H2''	3:D:6:DT:H71	1.98	0.44
1:B:108:LEU:CD1	1:B:129:PRO:HB3	2.49	0.43
1:B:67:THR:OG1	1:B:125:GLU:OE1	2.28	0.43
1:B:219:ASP:O	1:B:221:GLU:N	2.44	0.43
1:A:158:GLN:HB2	1:A:319:GLN:HG2	2.01	0.42
1:B:108:LEU:HD13	1:B:134:MET:HE3	2.01	0.42
1:B:67:THR:HG23	1:B:99:GLY:H	1.84	0.42
1:A:64:MET:HE3	1:A:94:LEU:HD13	2.02	0.42
3:D:27:DC:H1'	3:D:28:DC:H5'	2.00	0.42
1:A:147:TRP:O	1:A:149:PRO:HD3	2.20	0.42
1:A:150:PHE:CE1	1:A:152:GLY:C	2.98	0.41
1:A:10:LEU:HD23	1:A:10:LEU:C	2.46	0.41
1:A:131:ARG:HB3	1:A:135:GLN:HE21	1.86	0.41
1:A:108:LEU:HD21	1:A:134:MET:CE	2.50	0.41
1:B:314:GLN:H	1:B:314:GLN:CD	2.28	0.41
1:A:290:HIS:HB3	1:A:323:ILE:CG1	2.51	0.41
1:B:142:THR:HB	1:B:154:SER:HB2	2.03	0.40
1:A:273:TYR:O	1:A:274:ASP:CB	2.69	0.40
1:B:316:GLN:NE2	1:B:316:GLN:HA	2.35	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)
5:A:524:HOH:O	5:A:524:HOH:O[5_554]	2.15	0.05

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	328/330 (99%)	320 (98%)	7 (2%)	1 (0%)	37	37
1	B	329/330 (100%)	318 (97%)	7 (2%)	4 (1%)	11	7
All	All	657/660 (100%)	638 (97%)	14 (2%)	5 (1%)	16	13

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	274	ASP
1	B	220	PHE
1	B	69	SER
1	B	151	ASP
1	B	149	PRO

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	280/280 (100%)	272 (97%)	8 (3%)	37	41
1	B	281/280 (100%)	274 (98%)	7 (2%)	42	47
All	All	561/560 (100%)	546 (97%)	15 (3%)	40	44

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	67	THR
1	A	98	GLU
1	A	123	CYS
1	A	126	THR
1	A	150	PHE
1	A	230	ARG
1	A	268	ILE
1	A	307	ARG

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Mol	Chain	Res	Type
1	B	67	THR
1	B	123	CYS
1	B	131	ARG
1	B	312	THR
1	B	313	LEU
1	B	317	ARG
1	B	318	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	58	GLN
1	A	60	HIS
1	A	102	GLN
1	A	128	GLN
1	A	135	GLN
1	A	160	ASN
1	A	177	HIS
1	A	223	ASN
1	A	291	GLN
1	A	316	GLN
1	B	23	ASN
1	B	127	HIS
1	B	160	ASN
1	B	177	HIS
1	B	306	HIS
1	B	316	GLN
1	B	319	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

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	$\# Z > 2$
4	SO4	B	401	-	4,4,4	0.31	0	6,6,6	0.12	0
4	SO4	A	401	-	4,4,4	0.34	0	6,6,6	0.08	0
4	SO4	B	402	-	4,4,4	0.36	0	6,6,6	0.06	0

There are no bond length outliers.

There are no bond angle outliers.

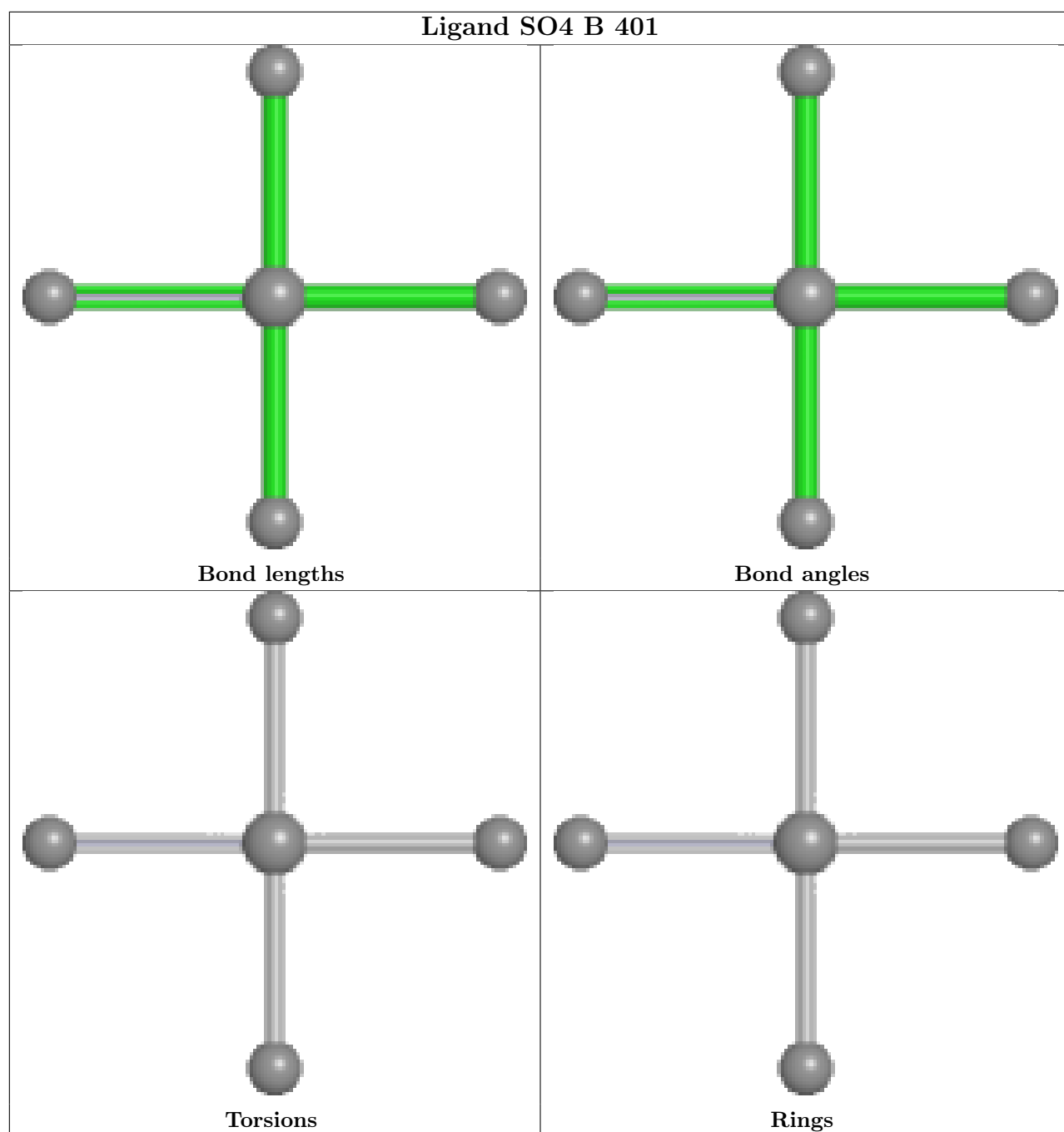
There are no chirality outliers.

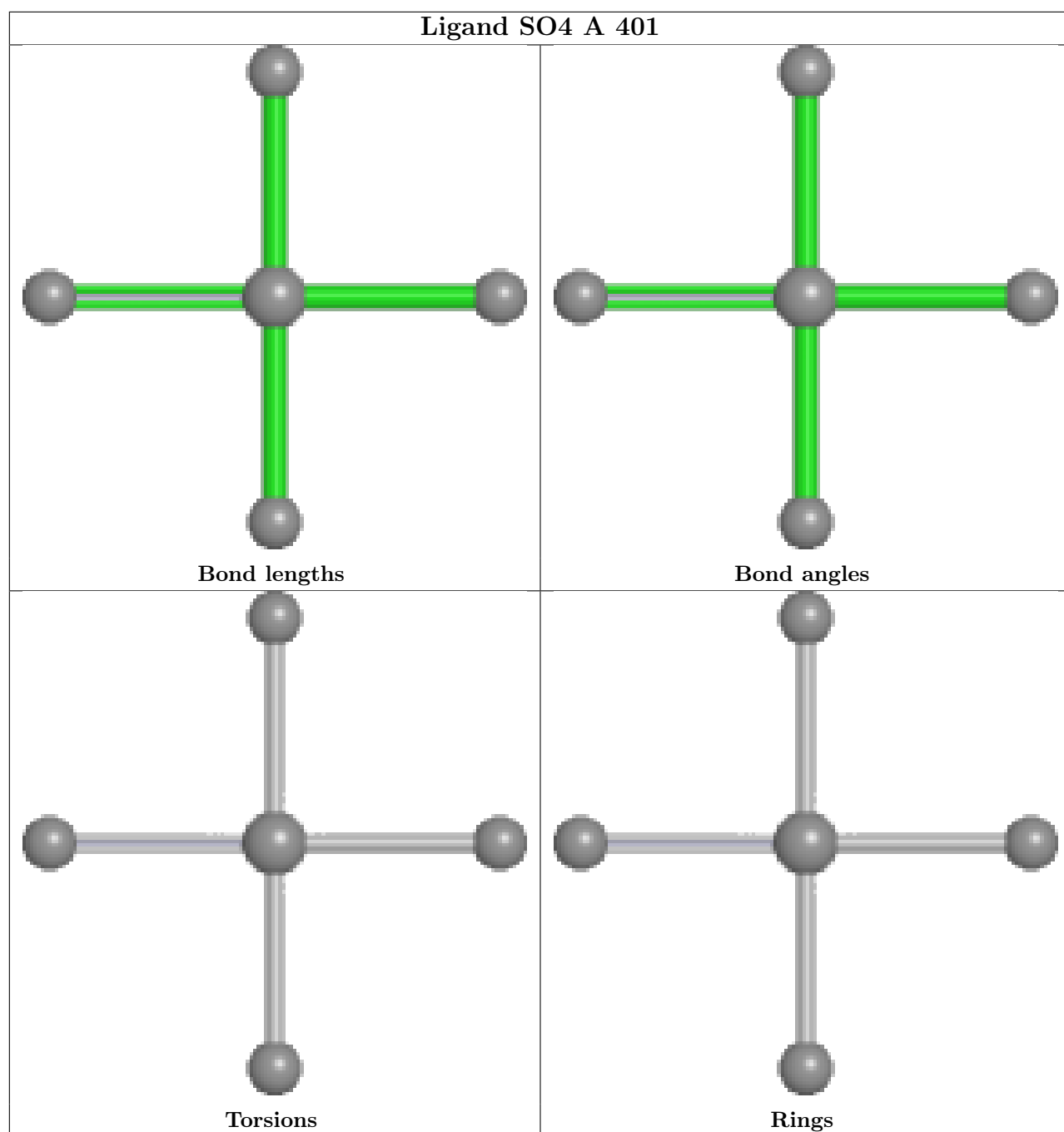
There are no torsion outliers.

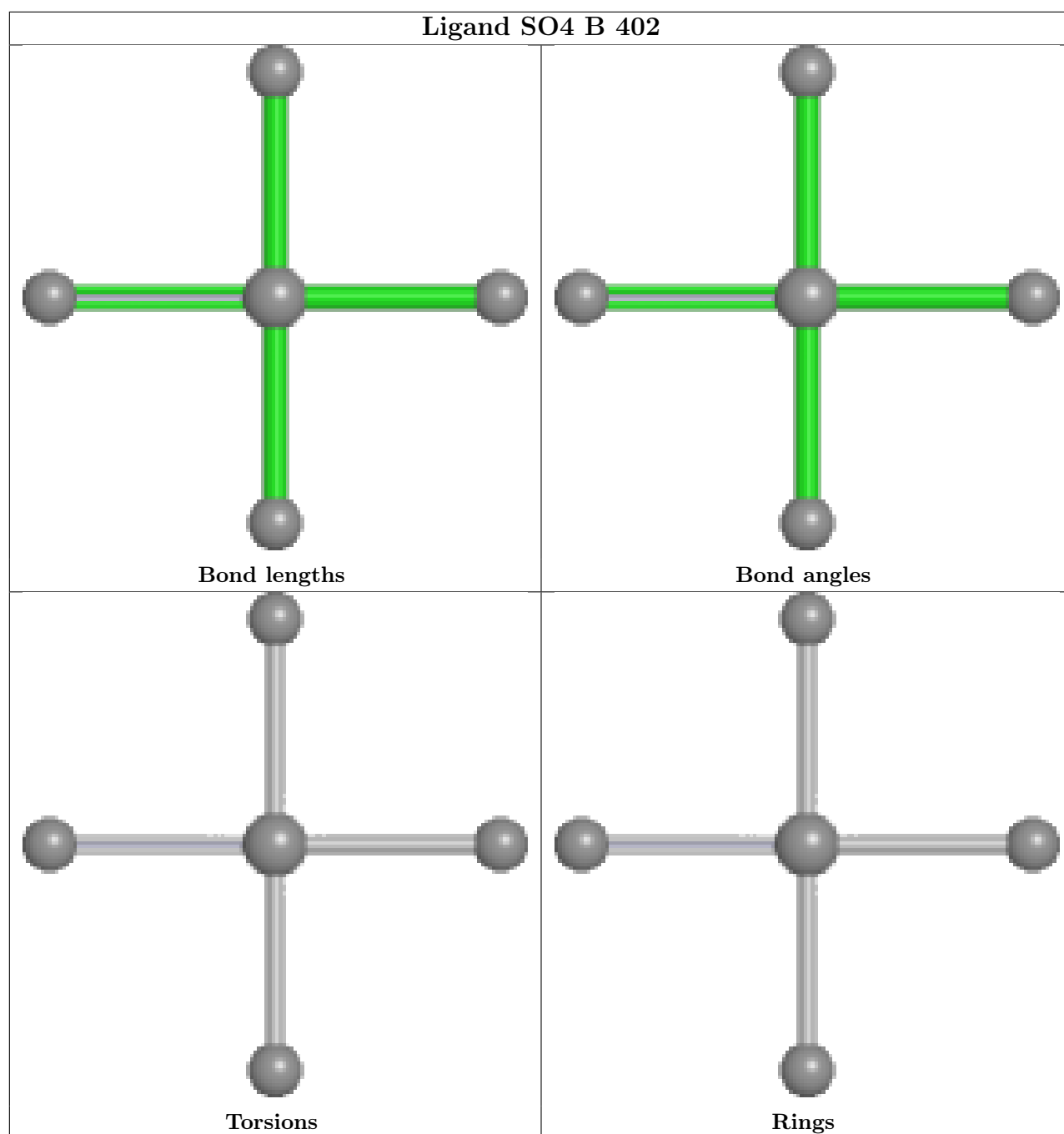
There are no ring outliers.

No monomer is involved in short contacts.

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







5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	330/330 (100%)	0.15	6 (1%) 67 69	45, 61, 97, 130	0
1	B	330/330 (100%)	0.12	3 (0%) 81 82	28, 60, 94, 157	1 (0%)
2	C	30/30 (100%)	-0.24	0 100 100	55, 101, 199, 228	0
3	D	30/30 (100%)	-0.21	0 100 100	53, 104, 203, 245	0
All	All	720/720 (100%)	0.11	9 (1%) 74 76	28, 62, 119, 245	1 (0%)

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	148	ALA	4.6
1	A	147	TRP	4.5
1	A	149	PRO	2.9
1	B	126	THR	2.7
1	A	146	ASP	2.6
1	B	46	ALA	2.4
1	B	150	PHE	2.4
1	A	163	LEU	2.3
1	A	134	MET	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 oligosaccharides in this entry.

6.4 Ligands

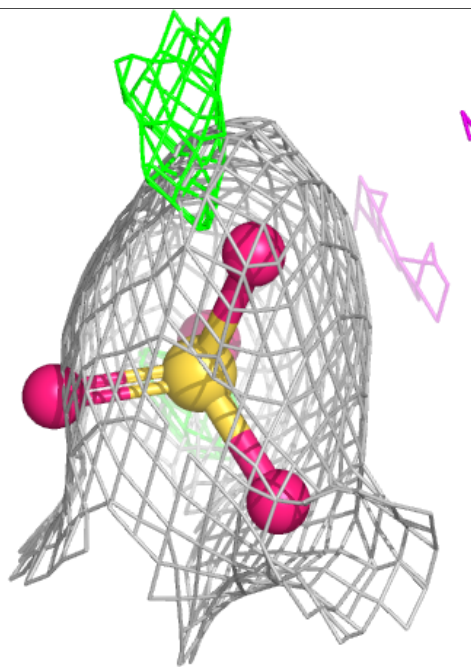
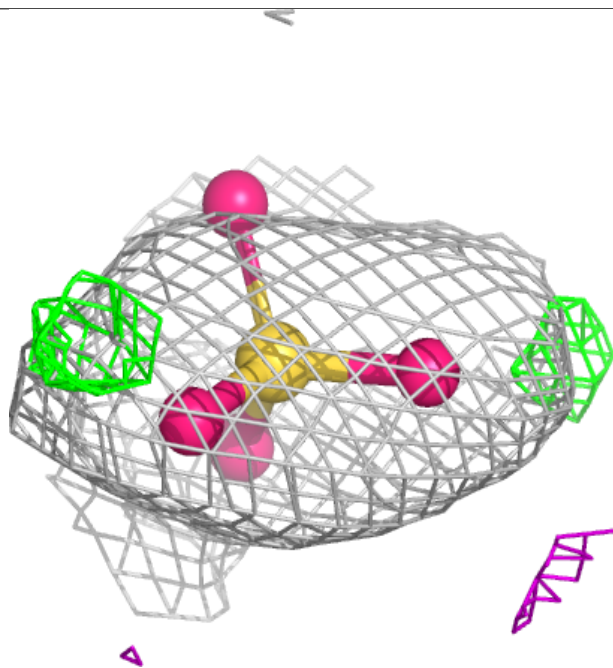
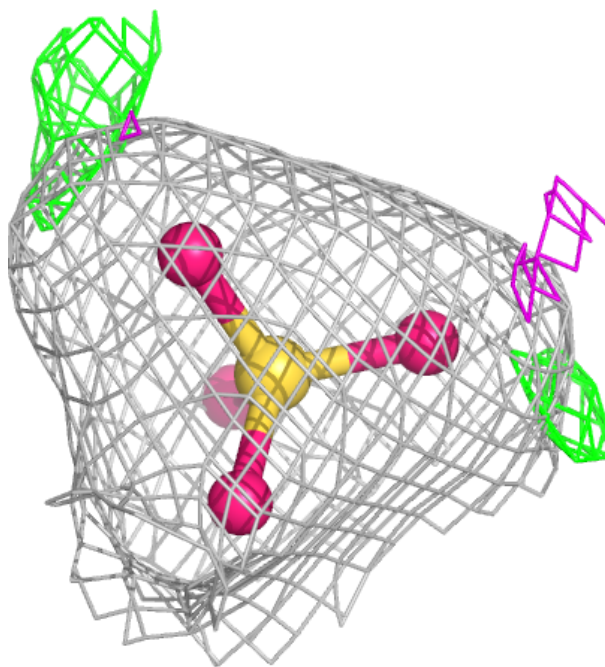
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	SO4	A	401	5/5	0.89	0.11	52,73,93,102	5
4	SO4	B	402	5/5	0.90	0.10	82,84,103,126	0
4	SO4	B	401	5/5	0.91	0.11	52,68,87,92	5

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.

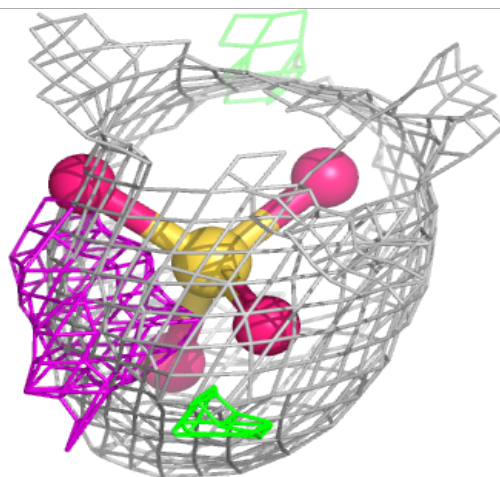
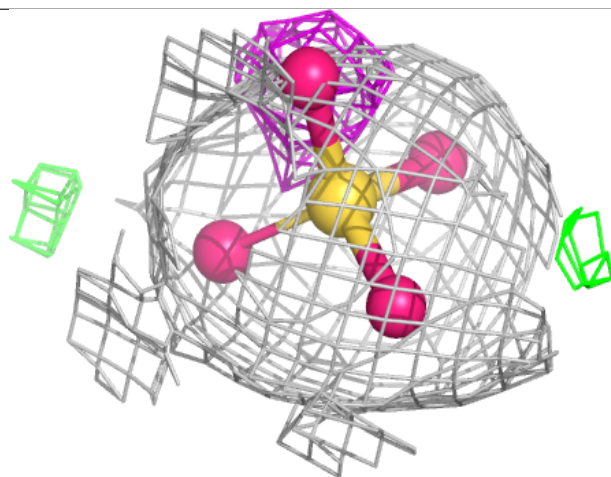
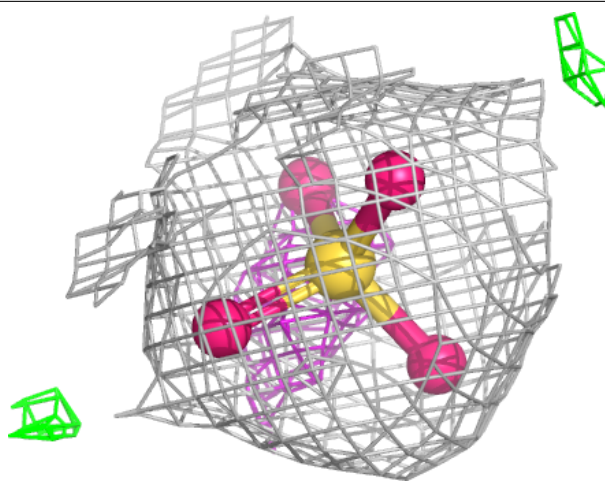
Electron density around SO4 A 401:

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



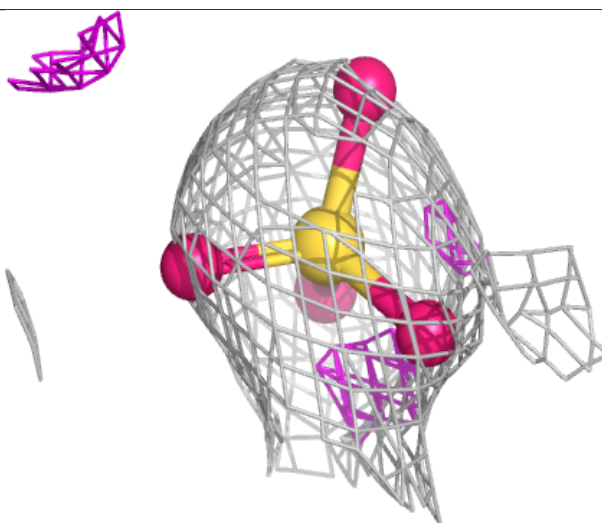
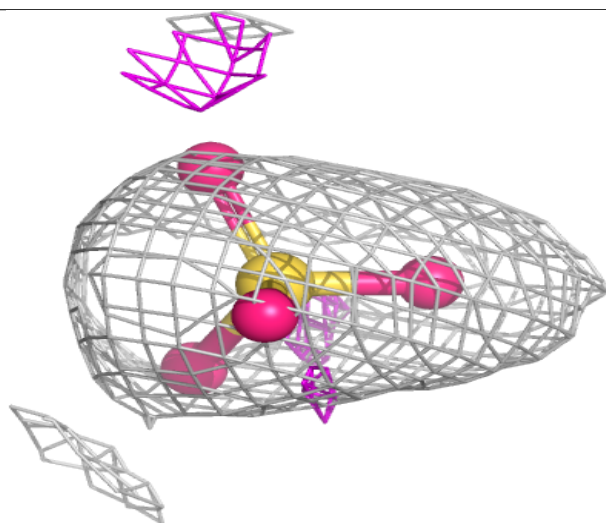
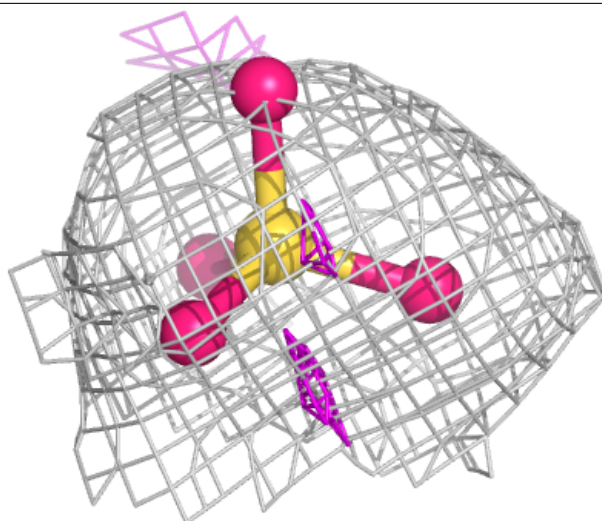
Electron density around SO4 B 402:

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



Electron density around SO4 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.