



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 10:40 PM UTC

PDB ID : 8FBF / pdb_00008fbf
Title : Crystal structure of OrfX2 from Clostridium botulinum E1
Authors : Lam, K.H.; Gao, L.
Deposited on : 2022-11-29
Resolution : 1.85 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 Phenix2.0
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

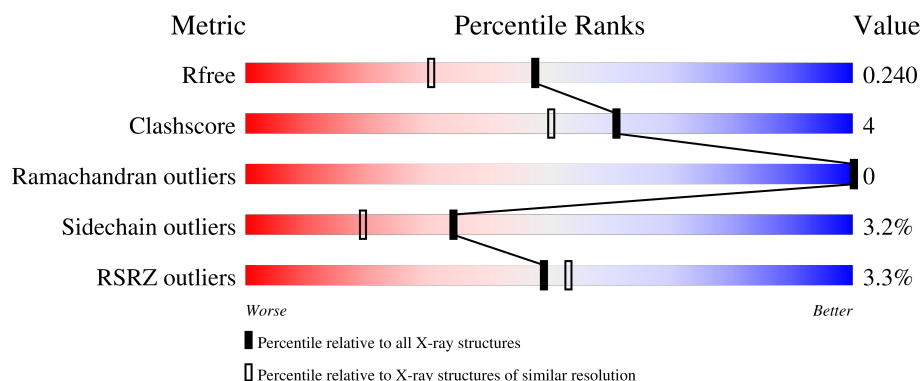
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

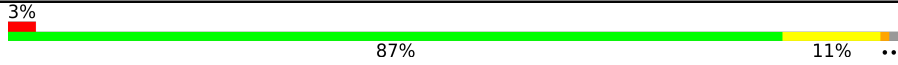
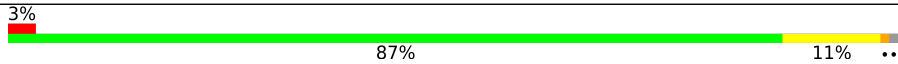
The reported resolution of this entry is 1.85 Å.

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	3428 (1.86-1.86)
Clashscore	190562	3579 (1.86-1.86)
Ramachandran outliers	187476	3553 (1.86-1.86)
Sidechain outliers	187428	3553 (1.86-1.86)
RSRZ outliers	180081	3429 (1.86-1.86)

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	748	
1	B	748	

2 Entry composition [i](#)

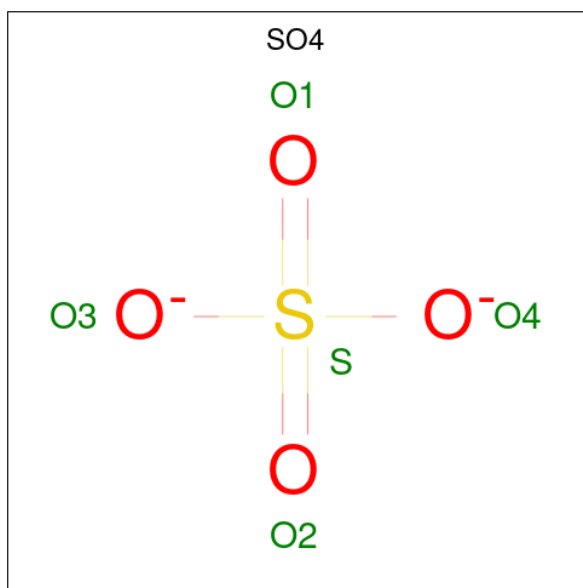
There are 3 unique types of molecules in this entry. The entry contains 12723 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Neurotoxin complex component Orf-X2.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	739	Total	C	N	O	S	Se	0	11	0
			5971	3830	982	1140	7	12			
1	B	738	Total	C	N	O	S	Se	0	3	0
			5884	3770	969	1126	7	12			

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S) (labeled as "Ligand of Interest" by depositor).



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

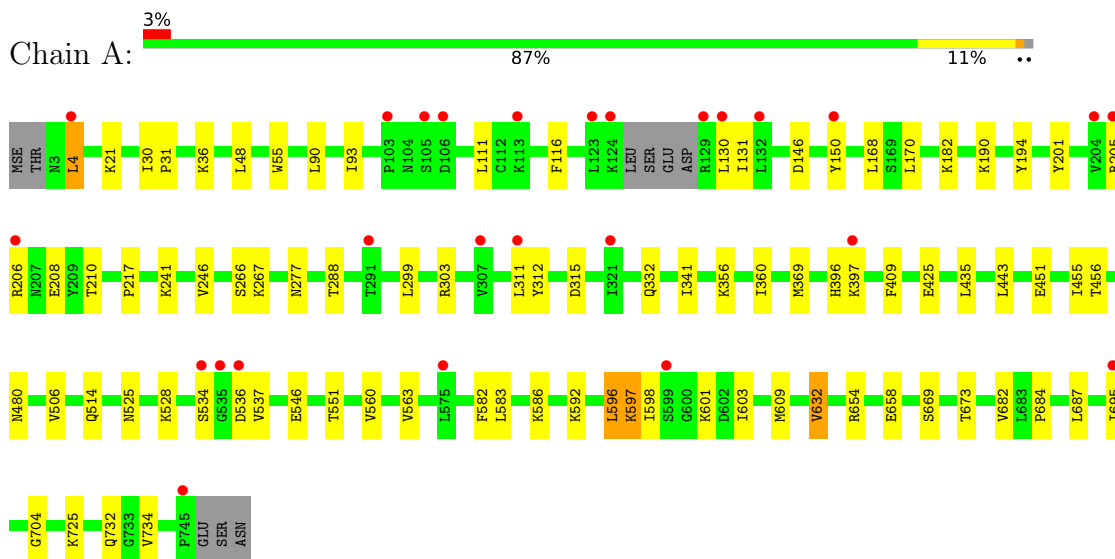
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	440	Total 440	O 440	0	0
3	B	408	Total 408	O 408	0	0

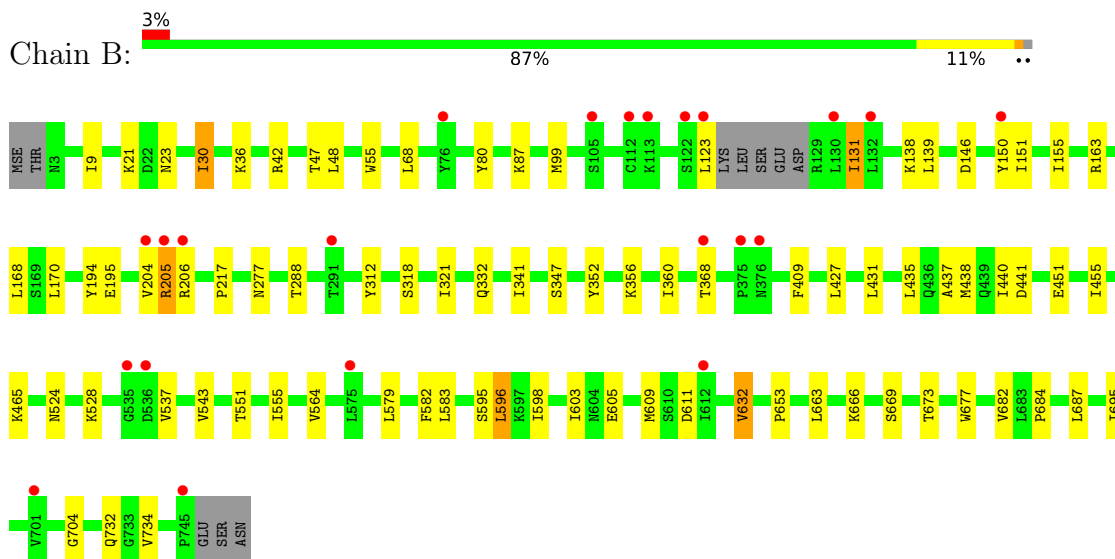
3 Residue-property plots [i](#)

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: Neurotoxin complex component Orf-X2



- Molecule 1: Neurotoxin complex component Orf-X2



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	106.32Å 83.19Å 108.42Å 90.00° 105.77° 90.00°	Depositor
Resolution (Å)	19.89 – 1.85 19.89 – 1.85	Depositor EDS
% Data completeness (in resolution range)	99.6 (19.89-1.85) 99.5 (19.89-1.85)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.91 (at 1.85Å)	Xtriage
Refinement program	PHENIX 1.15	Depositor
R, R_{free}	0.191 , 0.239 0.194 , 0.240	Depositor DCC
R_{free} test set	7661 reflections (4.94%)	wwPDB-VP
Wilson B-factor (Å ²)	23.3	Xtriage
Anisotropy	0.981	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 45.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.56$, $\langle L^2 \rangle = 0.40$	Xtriage
Estimated twinning fraction	0.000 for l,-k,h	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	12723	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 45.63 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.2863e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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	0.44	0/6078	0.72	1/8212 (0.0%)
1	B	0.47	0/5990	0.72	0/8094
All	All	0.45	0/12068	0.72	1/16306 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	205	ARG	CB-CA-C	-5.08	110.73	116.63

There are no chirality outliers.

There are no planarity outliers.

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	5971	0	6003	48	0
1	B	5884	0	5895	53	0
2	A	10	0	0	1	0
2	B	10	0	0	1	0
3	A	440	0	0	5	0
3	B	408	0	0	2	0
All	All	12723	0	11898	100	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 4.

All (100) 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:435:LEU:HA	1:B:438:MSE:HE2	1.66	0.78
1:B:596:LEU:C	1:B:596:LEU:HD23	2.15	0.72
1:A:36:LYS:HE3	1:A:146:ASP:OD2	1.92	0.69
1:B:205:ARG:CG	1:B:205:ARG:HH11	2.07	0.67
1:A:36:LYS:HE2	2:A:801:SO4:O2	1.96	0.66
1:B:36:LYS:HE2	2:B:801:SO4:O3	1.96	0.66
1:A:303:ARG:NH1	3:A:904:HOH:O	2.28	0.65
1:A:182:LYS:NZ	3:A:908:HOH:O	2.31	0.64
1:A:210[B]:THR:OG1	1:A:241:LYS:HB3	2.00	0.62
1:B:537:VAL:HG13	1:B:603:ILE:HD13	1.81	0.61
1:A:673:THR:HG21	1:A:687:LEU:HD23	1.86	0.58
1:A:725:LYS:NZ	3:A:903:HOH:O	2.27	0.57
1:A:528:LYS:HG2	1:A:704:GLY:HA2	1.86	0.56
1:B:288:THR:HG21	1:B:332:GLN:HA	1.86	0.56
1:A:30[A]:ILE:HG22	1:A:150[A]:TYR:HE2	1.71	0.56
1:B:528:LYS:HG2	1:B:704:GLY:HA2	1.86	0.56
1:A:669:SER:O	1:A:673:THR:HG23	2.06	0.56
1:B:605:GLU:HG2	1:B:609:MSE:HE2	1.88	0.55
1:B:451:GLU:H	1:B:451:GLU:CD	2.15	0.54
1:B:682:VAL:HG12	1:B:684:PRO:HD2	1.89	0.53
1:B:669:SER:O	1:B:673:THR:HG23	2.08	0.53
1:A:312:TYR:HB2	1:A:315:ASP:OD2	2.09	0.53
1:A:682:VAL:HG12	1:A:684:PRO:HD2	1.90	0.52
1:A:30[A]:ILE:HG21	1:A:55:TRP:CD2	2.44	0.52
1:A:210[B]:THR:HG1	1:A:241:LYS:HB3	1.72	0.52
1:A:597:LYS:HB2	1:A:597:LYS:NZ	2.25	0.52
1:B:435:LEU:HD22	1:B:455:ILE:HG21	1.91	0.52
1:B:131:ILE:HG12	1:B:204:VAL:HG13	1.92	0.52
1:B:48:LEU:HB2	1:B:139:LEU:HD13	1.93	0.51
1:B:30:ILE:HG22	1:B:150:TYR:HE2	1.76	0.51
1:B:205:ARG:CG	1:B:205:ARG:NH1	2.73	0.50
1:A:451:GLU:H	1:A:451:GLU:CD	2.18	0.50
1:A:31:PRO:HD2	1:A:150[A]:TYR:CE2	2.47	0.50
1:A:288:THR:HG21	1:A:332:GLN:HA	1.94	0.49
1:B:318:SER:O	1:B:321:ILE:HG22	2.12	0.49
1:A:514:GLN:OE1	1:A:537:VAL:HG22	2.12	0.48
1:A:396:HIS:HE1	1:A:425:GLU:OE2	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:266:SER:OG	3:A:901:HOH:O	2.20	0.48
1:A:4:LEU:HD21	1:A:168[A]:LEU:HD13	1.95	0.48
1:B:543:VAL:HG21	1:B:555:ILE:HD13	1.96	0.47
1:A:546:GLU:OE2	1:B:163:ARG:NH2	2.46	0.47
1:B:352:TYR:CE2	1:B:677:TRP:HB2	2.49	0.47
1:A:194:TYR:CZ	1:A:217:PRO:HA	2.50	0.47
1:B:465:LYS:HB3	1:B:653:PRO:HA	1.97	0.47
1:B:596:LEU:C	1:B:596:LEU:CD2	2.86	0.47
1:B:673:THR:HG21	1:B:687:LEU:HD23	1.96	0.47
1:B:9:ILE:HB	1:B:163:ARG:HB2	1.97	0.47
1:B:194:TYR:CZ	1:B:217:PRO:HA	2.50	0.47
1:A:201:TYR:HD1	1:A:210[B]:THR:HG22	1.80	0.47
1:A:168[A]:LEU:HB2	1:A:732:GLN:O	2.14	0.46
1:B:347:SER:HB3	1:B:356:LYS:HE2	1.97	0.46
1:A:435:LEU:HD22	1:A:455:ILE:HG21	1.97	0.46
1:B:30:ILE:HG21	1:B:55:TRP:CE3	2.50	0.46
1:B:36:LYS:HE3	1:B:146:ASP:OD2	2.15	0.46
1:B:30:ILE:HG22	1:B:150:TYR:CE2	2.51	0.45
1:B:427:LEU:HD12	1:B:431:LEU:HD22	1.98	0.45
1:B:564:VAL:HG21	1:B:579:LEU:HD13	1.97	0.45
1:B:30:ILE:HG21	1:B:55:TRP:CD2	2.51	0.45
1:B:341:ILE:HG21	1:B:360:ILE:HG21	1.99	0.45
1:B:168[A]:LEU:HB2	1:B:732:GLN:O	2.17	0.44
1:B:87:LYS:NZ	3:B:916:HOH:O	2.46	0.44
1:A:341:ILE:HG21	1:A:360:ILE:HG21	1.98	0.44
1:B:42:ARG:HG3	1:B:312:TYR:CG	2.53	0.44
1:B:583:LEU:HD22	1:B:598:ILE:HD11	1.98	0.44
1:B:68:LEU:HD12	1:B:99:MSE:HE1	2.00	0.44
1:A:592:LYS:NZ	1:A:592:LYS:HB3	2.33	0.44
1:B:47[B]:THR:HG23	1:B:80:TYR:HE2	1.83	0.44
1:B:170:LEU:HG	1:B:734:VAL:HG11	2.00	0.44
1:B:582:PHE:CG	1:B:609:MSE:HE1	2.52	0.43
1:B:205:ARG:HB2	3:B:989:HOH:O	2.18	0.43
1:A:30[A]:ILE:HG22	1:A:150[A]:TYR:CE2	2.52	0.43
1:A:456:THR:OG1	1:A:480:ASN:HA	2.18	0.43
1:A:111:LEU:HD21	1:A:116:PHE:CZ	2.53	0.43
1:B:437:ALA:O	1:B:666:LYS:HD3	2.19	0.43
1:A:560:VAL:HA	1:A:583:LEU:HD12	2.01	0.43
1:A:168[A]:LEU:HG	1:A:409:PHE:CE1	2.54	0.43
1:A:563:VAL:HG11	1:A:598:ILE:HD11	2.01	0.43
1:B:440:ILE:HD11	1:B:663:LEU:HD13	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:356:LYS:NZ	3:A:905:HOH:O	2.53	0.42
1:A:632:VAL:HG22	1:A:687:LEU:HD13	2.02	0.42
1:B:205:ARG:HB3	1:B:206:ARG:H	1.67	0.42
1:B:596:LEU:HD23	1:B:596:LEU:O	2.18	0.42
1:A:170:LEU:HG	1:A:734:VAL:HG11	2.02	0.42
1:A:654:ARG:HD3	1:A:658:GLU:OE2	2.20	0.42
1:A:534:SER:O	1:A:534:SER:OG	2.38	0.42
1:A:596[A]:LEU:HD22	1:A:597:LYS:N	2.34	0.42
1:A:90:LEU:HA	1:A:93:ILE:HD12	2.01	0.41
1:B:131:ILE:HD12	1:B:131:ILE:HA	1.92	0.41
1:B:205:ARG:NH1	1:B:205:ARG:HG2	2.35	0.41
1:A:201:TYR:HD1	1:A:210[B]:THR:CG2	2.33	0.41
1:B:205:ARG:HH11	1:B:205:ARG:HG3	1.83	0.41
1:A:563:VAL:CG1	1:A:598:ILE:HD11	2.50	0.41
1:A:111:LEU:HD21	1:A:116:PHE:HZ	1.85	0.41
1:B:168[A]:LEU:HG	1:B:409:PHE:CE1	2.55	0.41
1:B:632:VAL:HG22	1:B:687:LEU:HD13	2.02	0.41
1:A:586:LYS:HB2	1:A:596[A]:LEU:HD23	2.02	0.41
1:A:201:TYR:HE1	1:A:208:GLU:HB3	1.86	0.40
1:B:582:PHE:CD2	1:B:609:MSE:HE1	2.56	0.40
1:B:151:ILE:O	1:B:155:ILE:HG12	2.21	0.40
1:A:582:PHE:CG	1:A:609:MSE:HE1	2.57	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	746/748 (100%)	736 (99%)	10 (1%)	0	100	100
1	B	737/748 (98%)	726 (98%)	11 (2%)	0	100	100
All	All	1483/1496 (99%)	1462 (99%)	21 (1%)	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	677/664 (102%)	649 (96%)	28 (4%)	27	12
1	B	665/664 (100%)	647 (97%)	18 (3%)	39	24
All	All	1342/1328 (101%)	1296 (97%)	46 (3%)	34	17

All (46) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	LEU
1	A	21	LYS
1	A	48[A]	LEU
1	A	48[B]	LEU
1	A	130	LEU
1	A	131	ILE
1	A	190	LYS
1	A	206	ARG
1	A	246	VAL
1	A	267	LYS
1	A	277	ASN
1	A	299	LEU
1	A	311[A]	LEU
1	A	311[B]	LEU
1	A	369	MSE
1	A	397	LYS
1	A	443	LEU
1	A	506	VAL
1	A	525	ASN
1	A	536	ASP
1	A	551	THR
1	A	596[A]	LEU
1	A	596[B]	LEU
1	A	597	LYS
1	A	601	LYS

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Mol	Chain	Res	Type
1	A	603	ILE
1	A	632	VAL
1	A	695	ILE
1	B	21	LYS
1	B	23	ASN
1	B	30	ILE
1	B	123	LEU
1	B	131	ILE
1	B	138	LYS
1	B	195	GLU
1	B	205	ARG
1	B	277	ASN
1	B	368	THR
1	B	441	ASP
1	B	524	ASN
1	B	551	THR
1	B	595	SER
1	B	596	LEU
1	B	611	ASP
1	B	632	VAL
1	B	695	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	228	ASN
1	A	280	GLN
1	A	344	ASN
1	A	452	ASN
1	A	480	ASN
1	A	711	GLN
1	A	740	ASN
1	B	51	ASN
1	B	104	ASN
1	B	148	HIS
1	B	227	GLN
1	B	295	ASN
1	B	395	ASN
1	B	436	GLN
1	B	470	ASN
1	B	525	ASN
1	B	711	GLN

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Mol	Chain	Res	Type
1	B	740	ASN

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](#)

4 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	SO4	A	801	-	4,4,4	0.13	0	6,6,6	0.29	0
2	SO4	B	802	-	4,4,4	0.24	0	6,6,6	0.32	0
2	SO4	B	801	-	4,4,4	0.31	0	6,6,6	0.23	0
2	SO4	A	802	-	4,4,4	0.31	0	6,6,6	0.43	0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

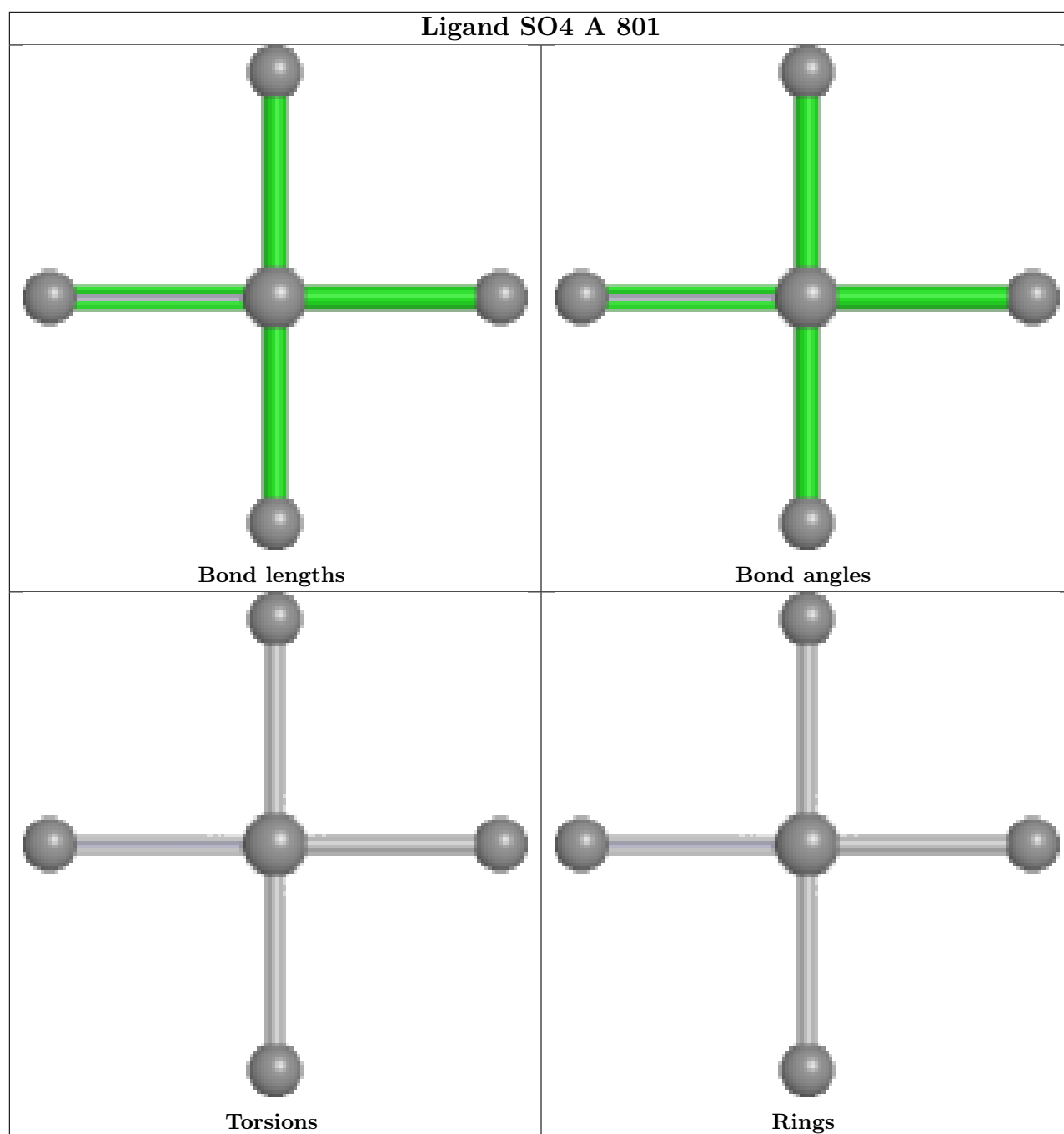
There are no torsion outliers.

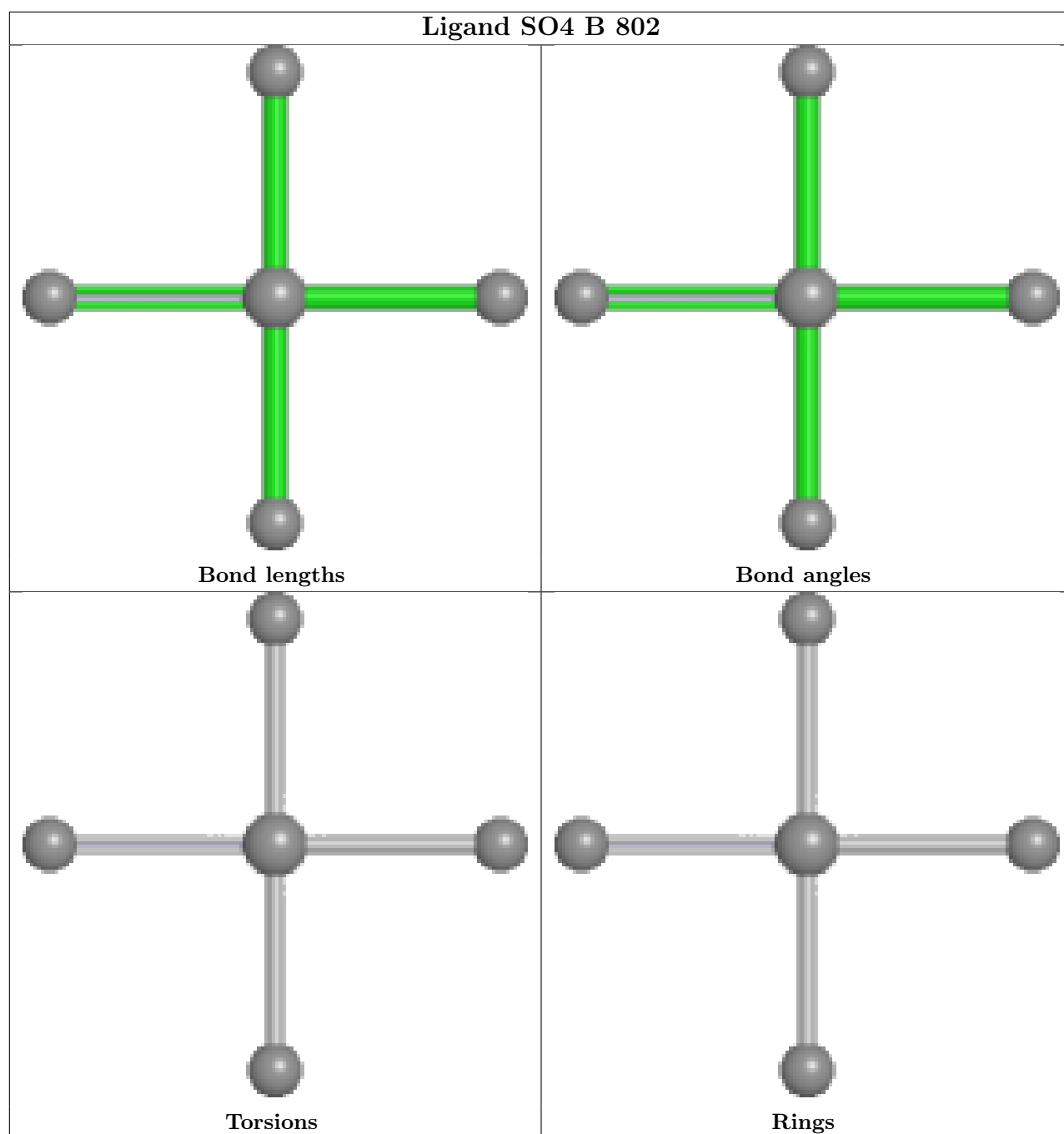
There are no ring outliers.

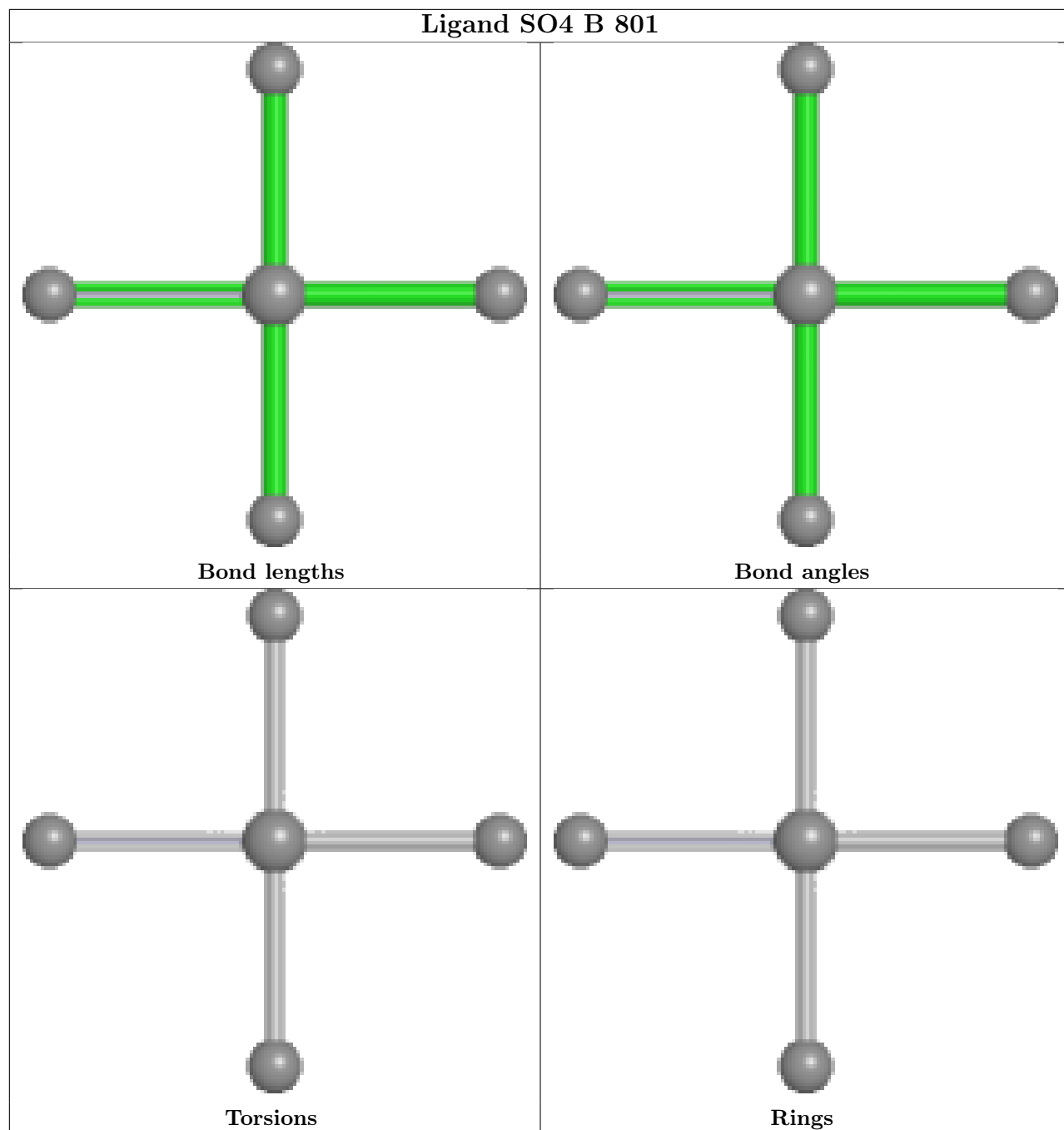
2 monomers are involved in 2 short contacts:

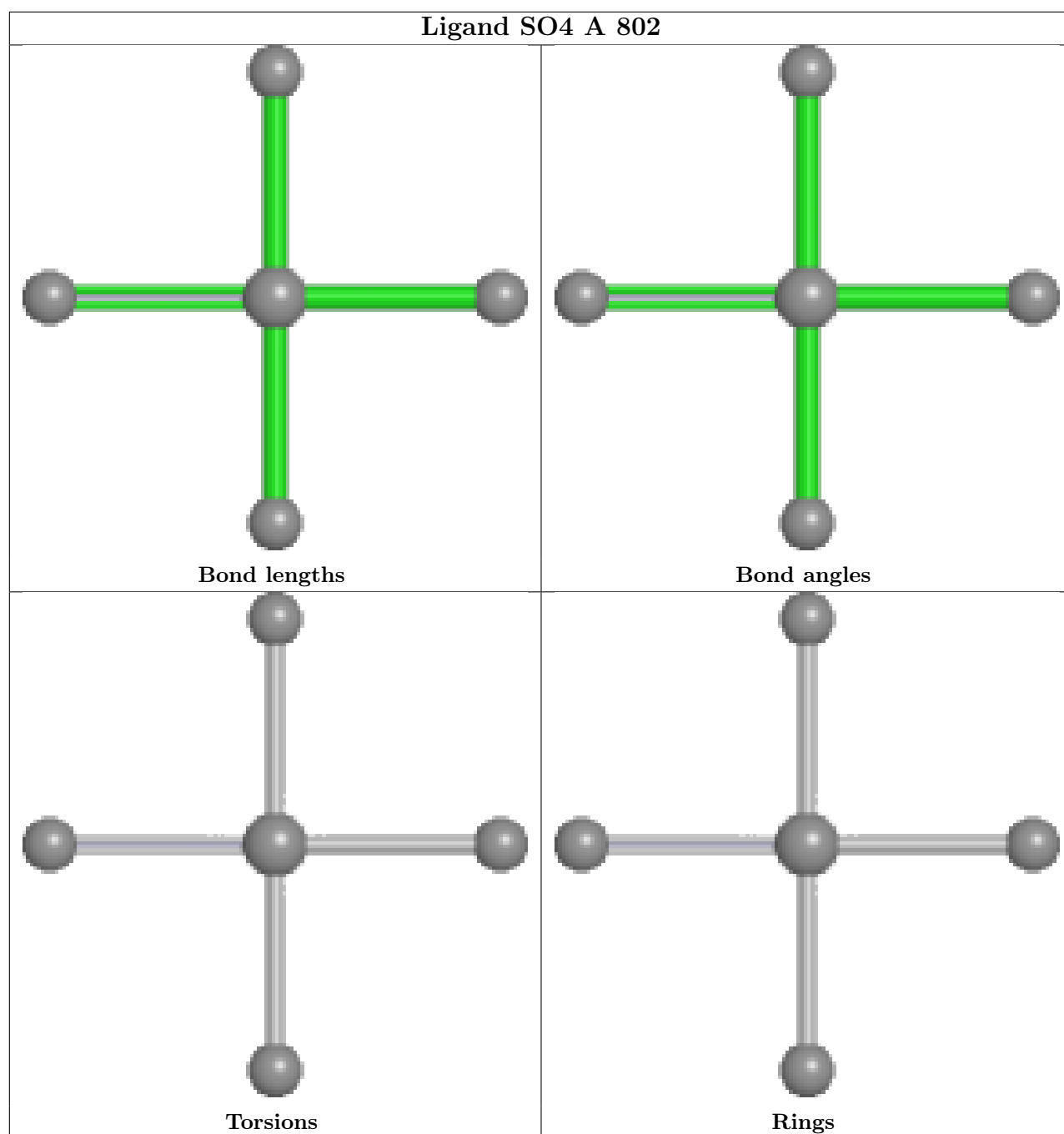
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	801	SO4	1	0
2	B	801	SO4	1	0

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	A	727/748 (97%)	0.16	26 (3%)	46 50	12, 32, 55, 80	12 (1%)
1	B	726/748 (97%)	0.18	22 (3%)	52 56	10, 34, 58, 77	3 (0%)
All	All	1453/1496 (97%)	0.17	48 (3%)	49 53	10, 33, 56, 80	15 (1%)

All (48) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	535	GLY	4.9
1	A	150[A]	TYR	4.6
1	A	536	ASP	4.1
1	B	113	LYS	3.9
1	A	103	PRO	3.2
1	A	206	ARG	3.1
1	B	291	THR	3.1
1	B	123	LEU	3.0
1	B	204	VAL	3.0
1	A	113	LYS	3.0
1	B	536	ASP	3.0
1	A	124	LYS	3.0
1	A	534	SER	2.9
1	B	150	TYR	2.9
1	B	535	GLY	2.9
1	A	204	VAL	2.8
1	B	205	ARG	2.8
1	A	106	ASP	2.7
1	A	307	VAL	2.7
1	A	132	LEU	2.6
1	B	105	SER	2.6
1	B	612	ILE	2.6
1	A	4	LEU	2.6
1	B	575	LEU	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	745	PRO	2.4
1	A	599	SER	2.4
1	A	397	LYS	2.4
1	A	575	LEU	2.4
1	B	122	SER	2.4
1	A	123	LEU	2.3
1	A	129	ARG	2.3
1	B	132	LEU	2.3
1	B	745	PRO	2.3
1	A	105	SER	2.3
1	B	130	LEU	2.3
1	B	375	PRO	2.3
1	A	695	ILE	2.3
1	A	291[A]	THR	2.3
1	B	368	THR	2.3
1	A	205	ARG	2.2
1	B	206	ARG	2.2
1	B	701	VAL	2.2
1	B	76	TYR	2.2
1	A	130	LEU	2.2
1	B	112	CYS	2.1
1	A	321	ILE	2.1
1	B	376	ASN	2.1
1	A	311[A]	LEU	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 [i](#)

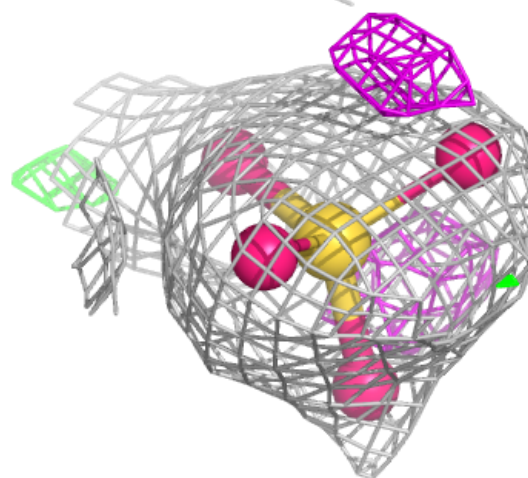
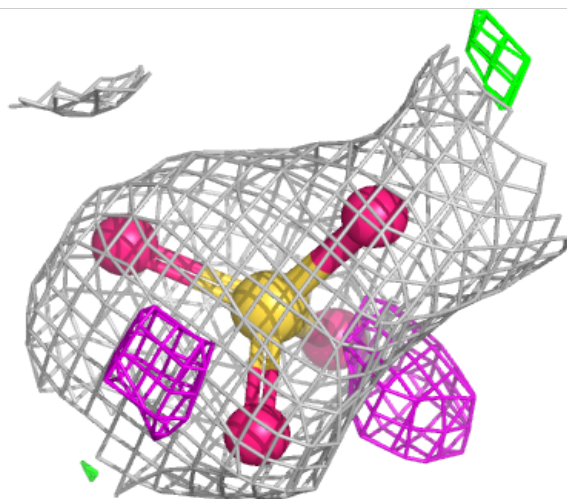
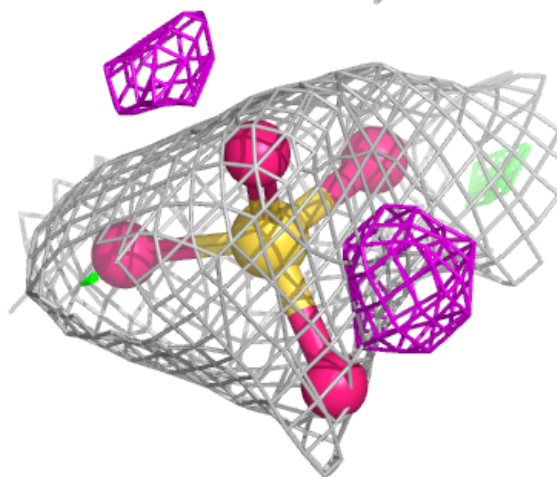
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	SO4	A	801	5/5	0.84	0.16	50,50,50,50	5
2	SO4	B	801	5/5	0.90	0.14	47,47,47,47	5
2	SO4	B	802	5/5	0.92	0.17	49,49,49,50	0
2	SO4	A	802	5/5	0.95	0.13	45,46,46,47	0

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

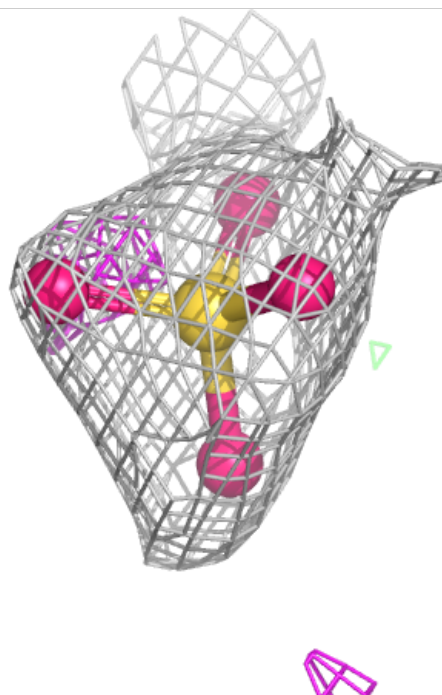
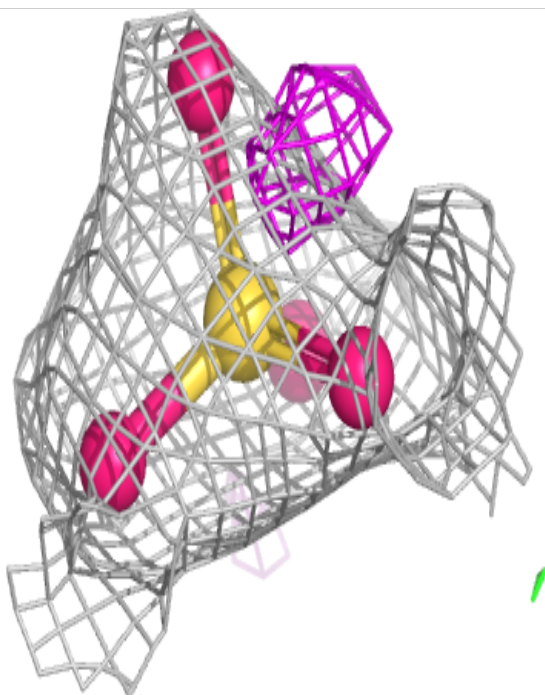
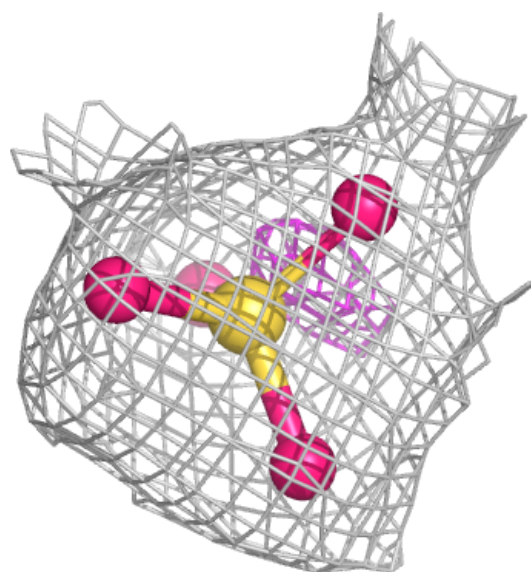
Electron density around SO4 A 801:

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



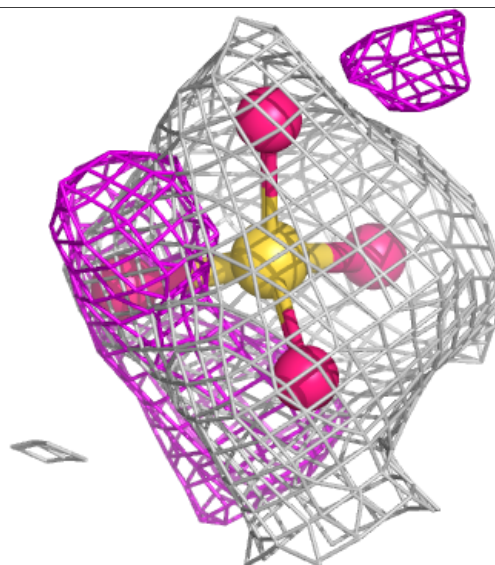
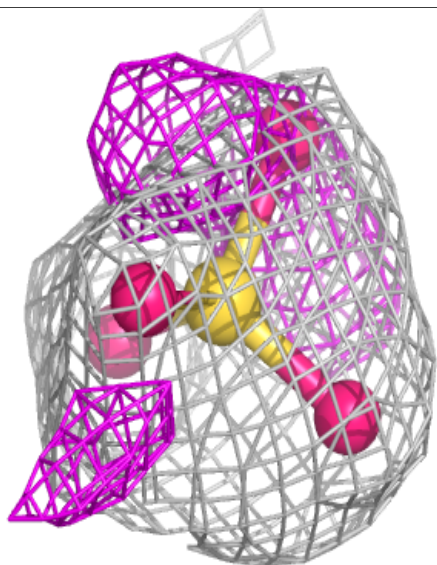
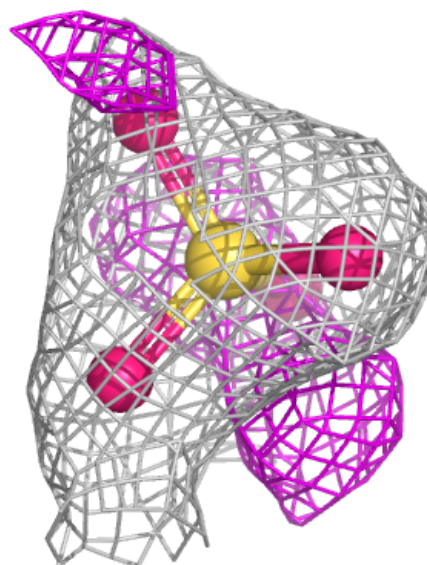
Electron density around SO4 B 801:

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



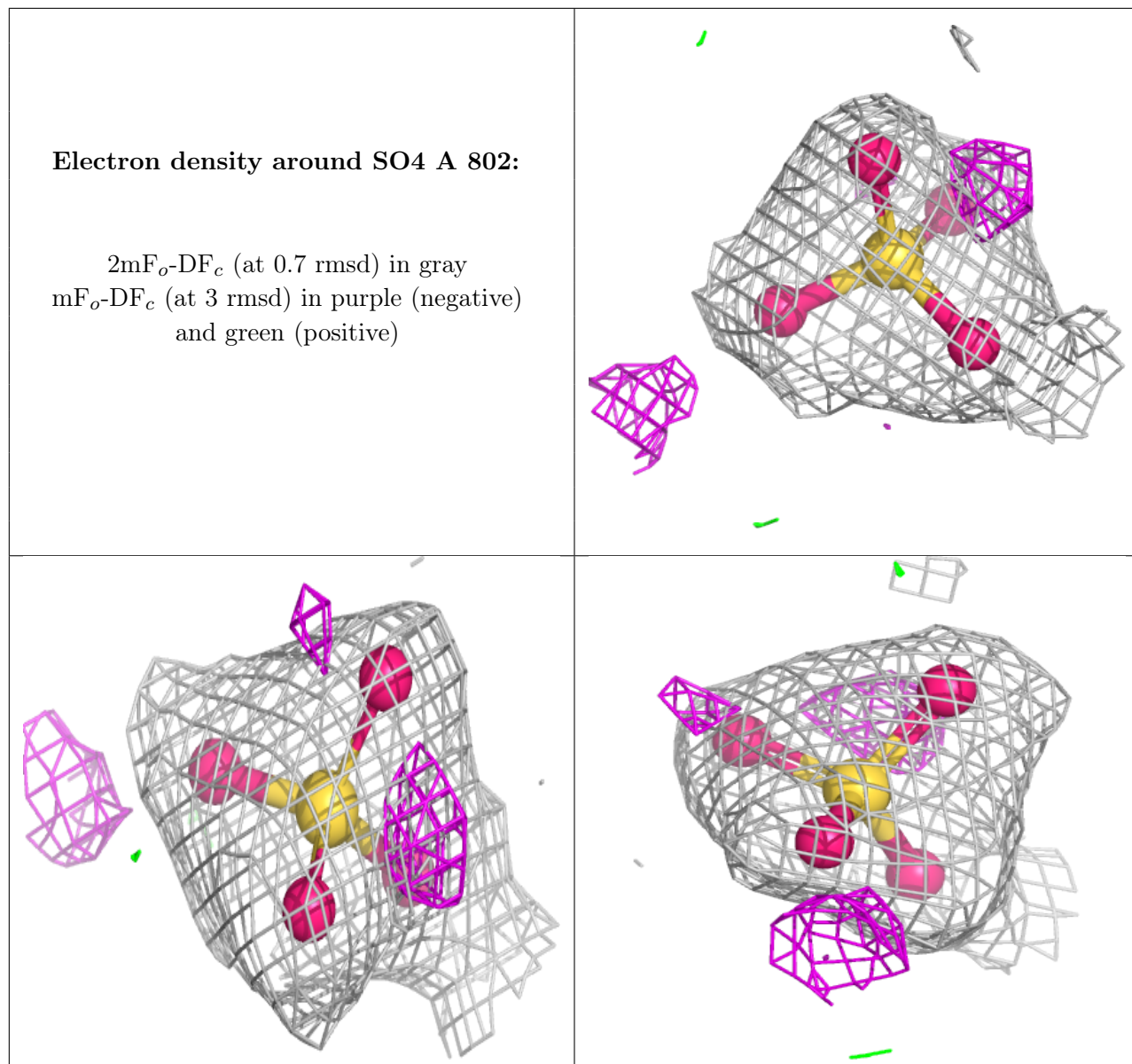
Electron density around SO4 B 802:

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



Electron density around SO4 A 802:

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



6.5 Other polymers ⓘ

There are no such residues in this entry.