



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 08:50 PM UTC

PDB ID : 8ZJ8 / pdb_00008zj8
Title : Acinetobacter baumannii ModA with molybdate H₂O₂
Authors : Wen, Y.; Jiao, M.
Deposited on : 2024-05-14
Resolution : 1.44 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

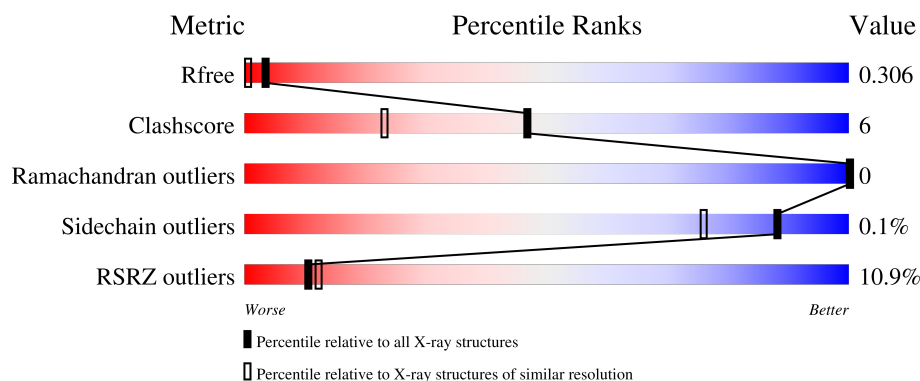
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.44 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	3234 (1.46-1.42)
Clashscore	190562	3289 (1.46-1.42)
Ramachandran outliers	187476	3248 (1.46-1.42)
Sidechain outliers	187428	3248 (1.46-1.42)
RSRZ outliers	180081	3234 (1.46-1.42)

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	241	12% 85% 8% 6%
1	B	241	9% 90% 5% 5%
1	C	241	10% 85% 10% 6%
1	D	241	10% 86% 8% 5%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard

residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	MOO	A	301	-	-	X	-
2	MOO	B	301	-	-	X	-
2	MOO	C	301	-	-	X	-

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 8035 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 Molybdate ABC transporter substrate-binding protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	226	Total	C	N	O	S	0	1	0
			1740	1115	292	329	4			
1	B	229	Total	C	N	O	S	0	2	0
			1765	1130	296	335	4			
1	C	227	Total	C	N	O	S	0	1	0
			1752	1121	295	332	4			
1	D	228	Total	C	N	O	S	0	2	0
			1759	1126	297	332	4			

There are 36 discrepancies between the modelled and reference sequences:

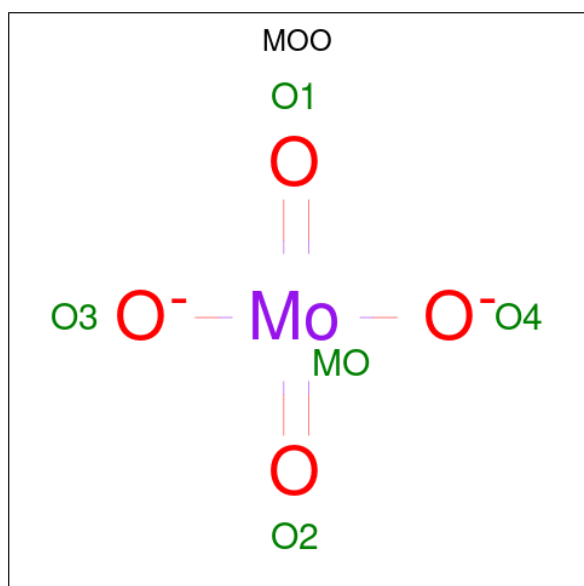
Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	initiating methionine	UNP A0A6F8TFE1
A	2	GLY	-	expression tag	UNP A0A6F8TFE1
A	3	HIS	-	expression tag	UNP A0A6F8TFE1
A	4	HIS	-	expression tag	UNP A0A6F8TFE1
A	5	HIS	-	expression tag	UNP A0A6F8TFE1
A	6	HIS	-	expression tag	UNP A0A6F8TFE1
A	7	HIS	-	expression tag	UNP A0A6F8TFE1
A	8	HIS	-	expression tag	UNP A0A6F8TFE1
A	9	MET	-	expression tag	UNP A0A6F8TFE1
B	1	MET	-	initiating methionine	UNP A0A6F8TFE1
B	2	GLY	-	expression tag	UNP A0A6F8TFE1
B	3	HIS	-	expression tag	UNP A0A6F8TFE1
B	4	HIS	-	expression tag	UNP A0A6F8TFE1
B	5	HIS	-	expression tag	UNP A0A6F8TFE1
B	6	HIS	-	expression tag	UNP A0A6F8TFE1
B	7	HIS	-	expression tag	UNP A0A6F8TFE1
B	8	HIS	-	expression tag	UNP A0A6F8TFE1
B	9	MET	-	expression tag	UNP A0A6F8TFE1
C	1	MET	-	initiating methionine	UNP A0A6F8TFE1
C	2	GLY	-	expression tag	UNP A0A6F8TFE1
C	3	HIS	-	expression tag	UNP A0A6F8TFE1

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Chain	Residue	Modelled	Actual	Comment	Reference
C	4	HIS	-	expression tag	UNP A0A6F8TFE1
C	5	HIS	-	expression tag	UNP A0A6F8TFE1
C	6	HIS	-	expression tag	UNP A0A6F8TFE1
C	7	HIS	-	expression tag	UNP A0A6F8TFE1
C	8	HIS	-	expression tag	UNP A0A6F8TFE1
C	9	MET	-	expression tag	UNP A0A6F8TFE1
D	1	MET	-	initiating methionine	UNP A0A6F8TFE1
D	2	GLY	-	expression tag	UNP A0A6F8TFE1
D	3	HIS	-	expression tag	UNP A0A6F8TFE1
D	4	HIS	-	expression tag	UNP A0A6F8TFE1
D	5	HIS	-	expression tag	UNP A0A6F8TFE1
D	6	HIS	-	expression tag	UNP A0A6F8TFE1
D	7	HIS	-	expression tag	UNP A0A6F8TFE1
D	8	HIS	-	expression tag	UNP A0A6F8TFE1
D	9	MET	-	expression tag	UNP A0A6F8TFE1

- Molecule 2 is MOLYBDATE ION (CCD ID: MOO) (formula: MoO_4) (labeled as "Ligand of Interest" by depositor).



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

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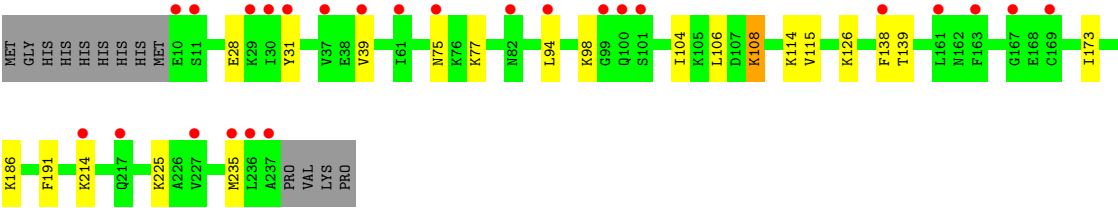
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	D	1	Total 5	Mo 1	O 4	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	244	Total 244	O 244	0	0
3	B	260	Total 260	O 260	0	0
3	C	244	Total 244	O 244	0	0
3	D	251	Total 251	O 251	0	0

- Molecule 1: Molybdate ABC transporter substrate-binding protein





4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	47.31Å 73.77Å 79.15Å 94.72° 107.43° 100.58°	Depositor
Resolution (Å)	33.52 – 1.44 33.52 – 1.44	Depositor EDS
% Data completeness (in resolution range)	92.1 (33.52-1.44) 92.4 (33.52-1.44)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.12 (at 1.44Å)	Xtriage
Refinement program	PHENIX 1.19_4092	Depositor
R, R_{free}	0.276 , 0.305 0.276 , 0.306	Depositor DCC
R_{free} test set	1848 reflections (1.03%)	wwPDB-VP
Wilson B-factor (Å ²)	16.3	Xtriage
Anisotropy	0.455	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 48.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.54$, $\langle L^2 \rangle = 0.39$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	8035	wwPDB-VP
Average B, all atoms (Å ²)	25.0	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 90.72 % 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 2.0418e-08. 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 [i](#)

5.1 Standard geometry [i](#)

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

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	0.31	0/1772	0.56	0/2405
1	B	0.29	0/1801	0.57	0/2445
1	C	0.31	0/1784	0.59	0/2420
1	D	0.29	0/1794	0.56	0/2433
All	All	0.30	0/7151	0.57	0/9703

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1740	0	1749	20	0
1	B	1765	0	1772	14	0
1	C	1752	0	1770	22	0
1	D	1759	0	1774	22	0
2	A	5	0	0	2	0
2	B	5	0	0	2	0
2	C	5	0	0	3	0
2	D	5	0	0	1	0
3	A	244	0	0	6	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	260	0	0	7	0
3	C	244	0	0	9	1
3	D	251	0	0	9	1
All	All	8035	0	7065	80	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 6.

All (80) 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:D:186:LYS:NZ	3:D:401:HOH:O	1.83	1.08
1:B:210:SER:OG	3:B:401:HOH:O	1.83	0.95
1:A:162:ASN:ND2	3:A:402:HOH:O	2.02	0.93
1:A:34:GLN:HE22	1:A:223:GLN:HE22	1.11	0.92
1:A:211:ASN:ND2	3:A:403:HOH:O	2.03	0.92
1:C:211:ASN:ND2	3:C:401:HOH:O	2.00	0.90
1:D:98:LYS:O	3:D:402:HOH:O	1.91	0.87
1:B:75:ASN:O	3:B:402:HOH:O	1.99	0.79
1:C:104:ILE:HD12	1:C:115:VAL:HB	1.66	0.77
1:C:237:ALA:O	3:C:402:HOH:O	2.05	0.73
1:A:225:LYS:HG3	1:A:235:MET:HG3	1.71	0.72
1:D:75:ASN:HB2	3:D:586:HOH:O	1.90	0.70
1:C:119:LYS:NZ	3:C:403:HOH:O	2.11	0.70
1:D:104:ILE:HD12	1:D:115:VAL:HB	1.73	0.69
1:C:138:PHE:HZ	1:C:173:ILE:HD11	1.58	0.68
1:C:225:LYS:HG3	1:C:235:MET:HG3	1.76	0.68
1:C:132:LYS:NZ	3:C:405:HOH:O	2.28	0.67
1:A:135:LYS:NZ	3:A:401:HOH:O	1.96	0.66
1:C:48:THR:O	1:C:52:GLN:HG3	1.96	0.66
1:B:162:ASN:HB3	3:B:420:HOH:O	1.96	0.66
1:A:124:ASP:OD2	1:D:126:LYS:NZ	2.26	0.65
1:D:139:THR:HG21	3:D:620:HOH:O	1.96	0.65
2:D:301:MOO:MO	2:D:301:MOO:O1	1.68	0.64
1:B:101:SER:HB3	1:B:186:LYS:HE3	1.80	0.63
2:A:301:MOO:O1	2:A:301:MOO:MO	1.69	0.63
1:D:94:LEU:HD11	1:D:106:LEU:HD21	1.80	0.63
1:B:39:VAL:O	3:B:403:HOH:O	2.16	0.61
1:A:111:ASP:HB3	1:A:114:LYS:HG2	1.81	0.61
2:B:301:MOO:MO	2:B:301:MOO:O3	1.71	0.61
1:A:101:SER:OG	1:A:186:LYS:HE3	2.00	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:31:TYR:HE2	1:A:37:VAL:HG21	1.67	0.60
2:C:301:MOO:O2	2:C:301:MOO:MO	1.73	0.60
1:A:138:PHE:HZ	1:A:173:ILE:HD11	1.67	0.60
2:A:301:MOO:MO	2:A:301:MOO:O4	1.73	0.59
1:B:163:PHE:HE2	1:C:163:PHE:HE2	1.49	0.59
2:B:301:MOO:MO	2:B:301:MOO:O1	1.73	0.59
2:C:301:MOO:MO	2:C:301:MOO:O4	1.71	0.59
1:C:229:LYS:HE2	3:C:603:HOH:O	2.02	0.59
1:B:124:ASP:OD2	1:C:126:LYS:NZ	2.31	0.58
2:C:301:MOO:MO	2:C:301:MOO:O3	1.73	0.58
1:D:114:LYS:NZ	3:D:409:HOH:O	2.37	0.56
1:A:31:TYR:CE2	1:A:37:VAL:HG21	2.41	0.55
1:D:214:LYS:NZ	3:D:405:HOH:O	2.28	0.55
1:A:111:ASP:H	1:A:114:LYS:NZ	2.05	0.54
1:A:35:ASN:HB2	1:A:37:VAL:HG22	1.89	0.54
1:D:31:TYR:OH	1:D:214:LYS:HG2	2.08	0.54
1:D:225:LYS:HG2	1:D:235:MET:HG3	1.89	0.54
1:D:94:LEU:CD1	1:D:106:LEU:HD21	2.39	0.53
1:C:75:ASN:ND2	3:C:406:HOH:O	2.34	0.53
1:A:67:THR:HG22	3:A:483:HOH:O	2.09	0.53
1:C:11:SER:N	3:C:409:HOH:O	2.42	0.52
1:A:67:THR:HG21	3:A:598:HOH:O	2.09	0.52
1:A:34:GLN:NE2	1:A:223:GLN:HE22	1.94	0.52
1:B:67:THR:HG21	3:B:607:HOH:O	2.09	0.51
1:D:94:LEU:HD12	1:D:106:LEU:HD11	1.94	0.50
1:D:126:LYS:HB2	3:D:479:HOH:O	2.11	0.50
1:A:162:ASN:ND2	3:A:416:HOH:O	2.44	0.50
1:B:28:GLU:HG3	1:B:39:VAL:HB	1.94	0.50
1:B:163:PHE:HE2	1:C:163:PHE:CE2	2.30	0.49
1:D:108:LYS:N	1:D:108:LYS:HD3	2.27	0.48
1:D:138:PHE:HZ	1:D:173:ILE:CD1	2.26	0.47
1:B:163:PHE:CE2	1:C:163:PHE:HE2	2.32	0.47
1:D:75:ASN:HB2	3:D:571:HOH:O	2.15	0.46
1:D:138:PHE:HZ	1:D:173:ILE:HD11	1.80	0.46
1:D:104:ILE:CD1	1:D:115:VAL:HB	2.42	0.46
1:C:138:PHE:CZ	1:C:173:ILE:CD1	3.00	0.45
1:A:111:ASP:H	1:A:114:LYS:HZ2	1.64	0.45
1:C:138:PHE:CZ	1:C:173:ILE:HD11	2.44	0.45
1:C:96:THR:HG22	1:C:171:VAL:HG12	1.98	0.44
1:D:77:LYS:NZ	3:D:410:HOH:O	2.40	0.44
1:D:28:GLU:HG3	1:D:39:VAL:HB	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:225:LYS:HG3	1:A:235:MET:CG	2.43	0.43
1:B:126:LYS:HB3	1:C:126:LYS:HE2	2.01	0.43
1:C:28:GLU:HG3	1:C:39:VAL:HB	2.01	0.43
1:D:106:LEU:HD22	1:D:191:PHE:CE2	2.53	0.43
1:C:105:LYS:HG3	3:C:610:HOH:O	2.19	0.43
1:B:150:LYS:HD3	3:B:473:HOH:O	2.18	0.42
1:A:72:TYR:O	1:A:76:LYS:HG2	2.20	0.42
1:B:40:LYS:NZ	3:B:416:HOH:O	2.49	0.41
1:C:150:LYS:HD3	3:C:458:HOH:O	2.20	0.41

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)
3:C:526:HOH:O	3:D:600:HOH:O[1_545]	2.01	0.19

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	225/241 (93%)	220 (98%)	5 (2%)	0	100	100
1	B	229/241 (95%)	227 (99%)	2 (1%)	0	100	100
1	C	226/241 (94%)	221 (98%)	5 (2%)	0	100	100
1	D	228/241 (95%)	224 (98%)	4 (2%)	0	100	100
All	All	908/964 (94%)	892 (98%)	16 (2%)	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	184/202 (91%)	184 (100%)	0	100	100
1	B	187/202 (93%)	187 (100%)	0	100	100
1	C	187/202 (93%)	187 (100%)	0	100	100
1	D	187/202 (93%)	186 (100%)	1 (0%)	81	63
All	All	745/808 (92%)	744 (100%)	1 (0%)	88	76

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	108	LYS

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	34	GLN
1	A	35	ASN
1	A	103	ASN
1	A	162	ASN
1	A	211	ASN
1	B	34	GLN
1	B	100	GLN
1	B	103	ASN
1	B	211	ASN
1	B	223	GLN
1	C	140	ASN
1	C	211	ASN
1	C	217	GLN
1	D	34	GLN
1	D	82	ASN
1	D	103	ASN
1	D	223	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](#)

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	MOO	A	301	-	2,4,4	1.93	1 (50%)	-		
2	MOO	B	301	-	2,4,4	0.83	0	-		
2	MOO	C	301	-	2,4,4	0.72	0	-		
2	MOO	D	301	-	2,4,4	2.28	1 (50%)	-		

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	301	MOO	O1-MO	-3.18	1.68	1.75
2	A	301	MOO	O1-MO	-2.72	1.69	1.75

There are no bond angle outliers.

There are no chirality outliers.

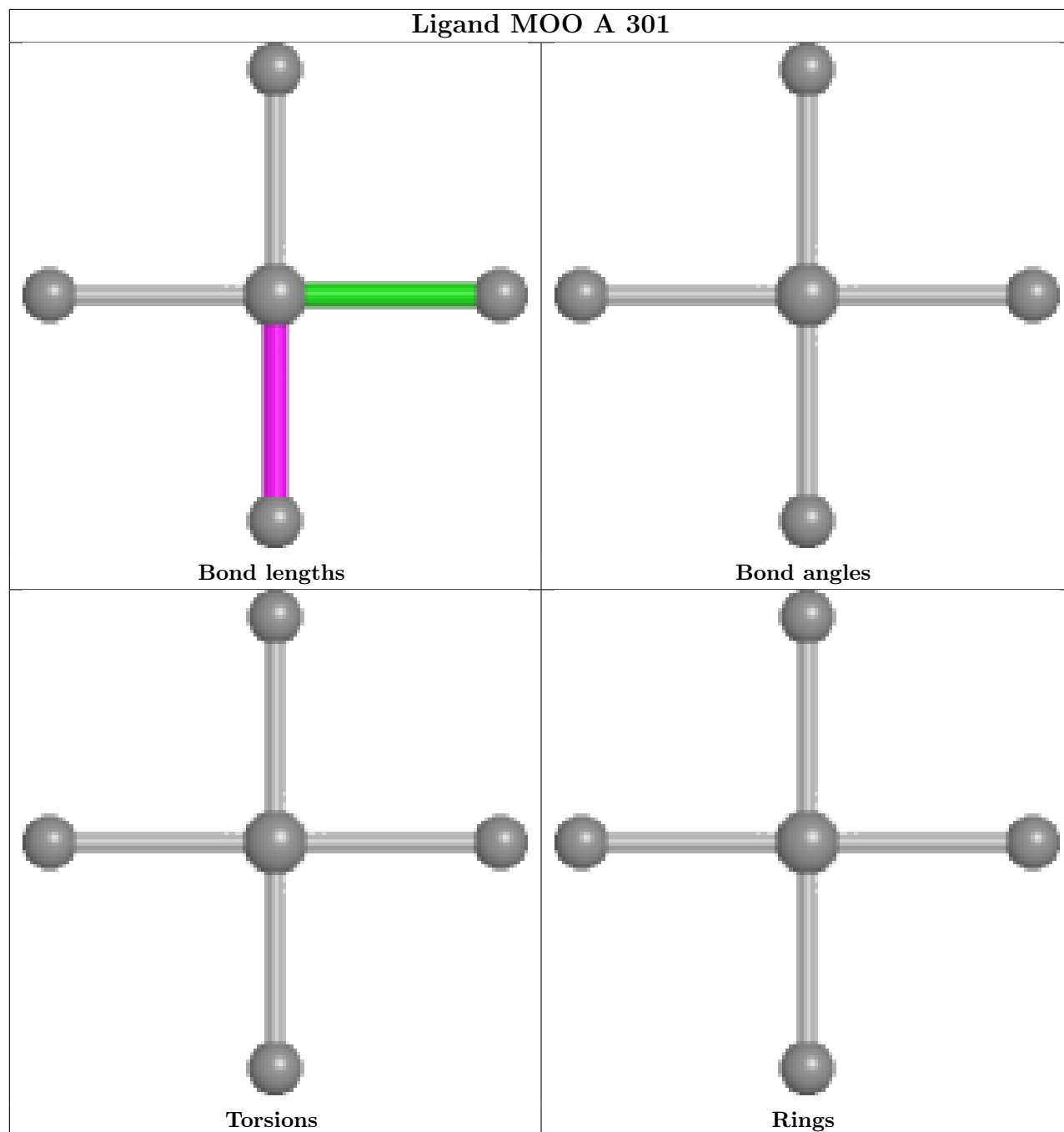
There are no torsion outliers.

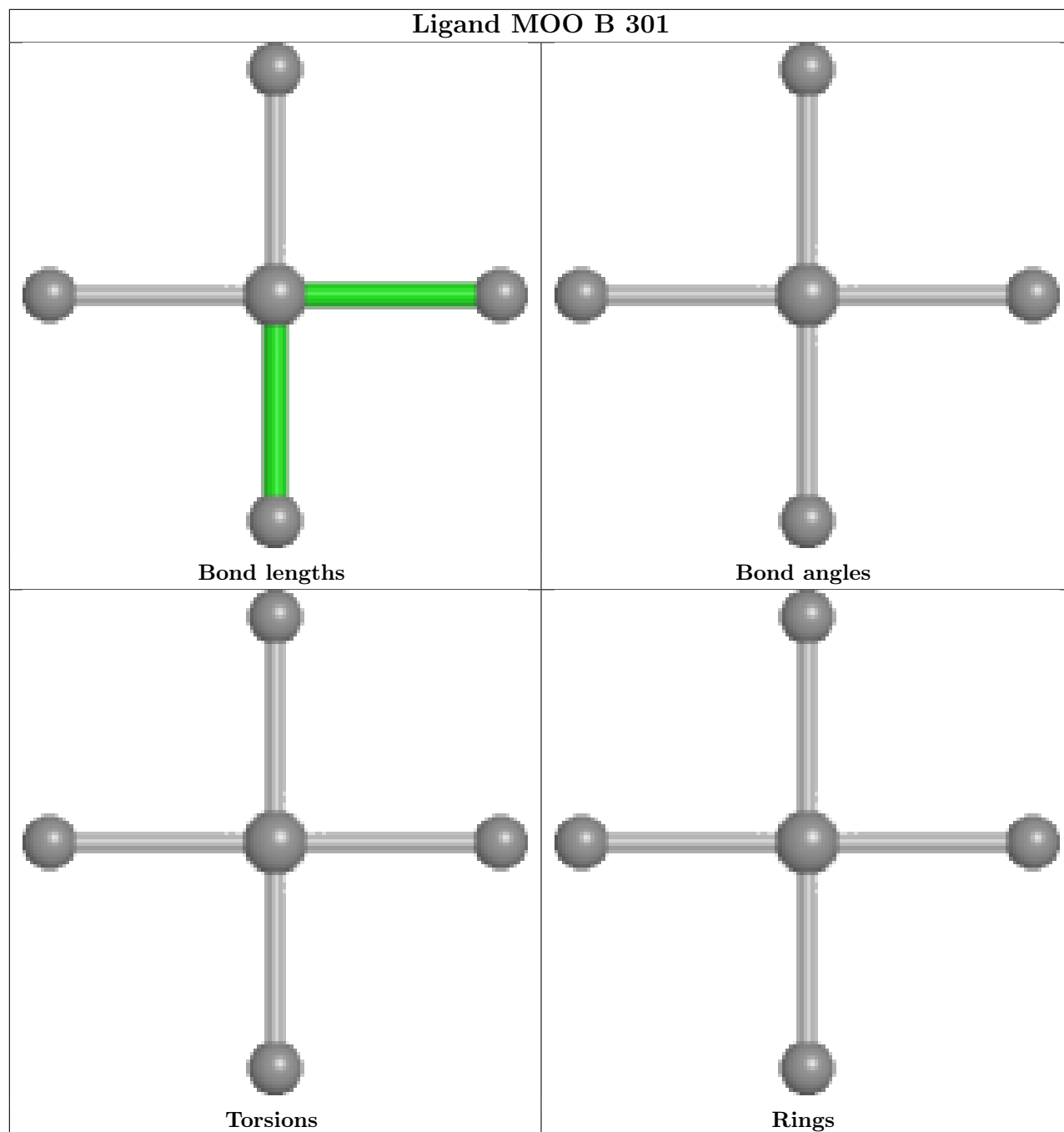
There are no ring outliers.

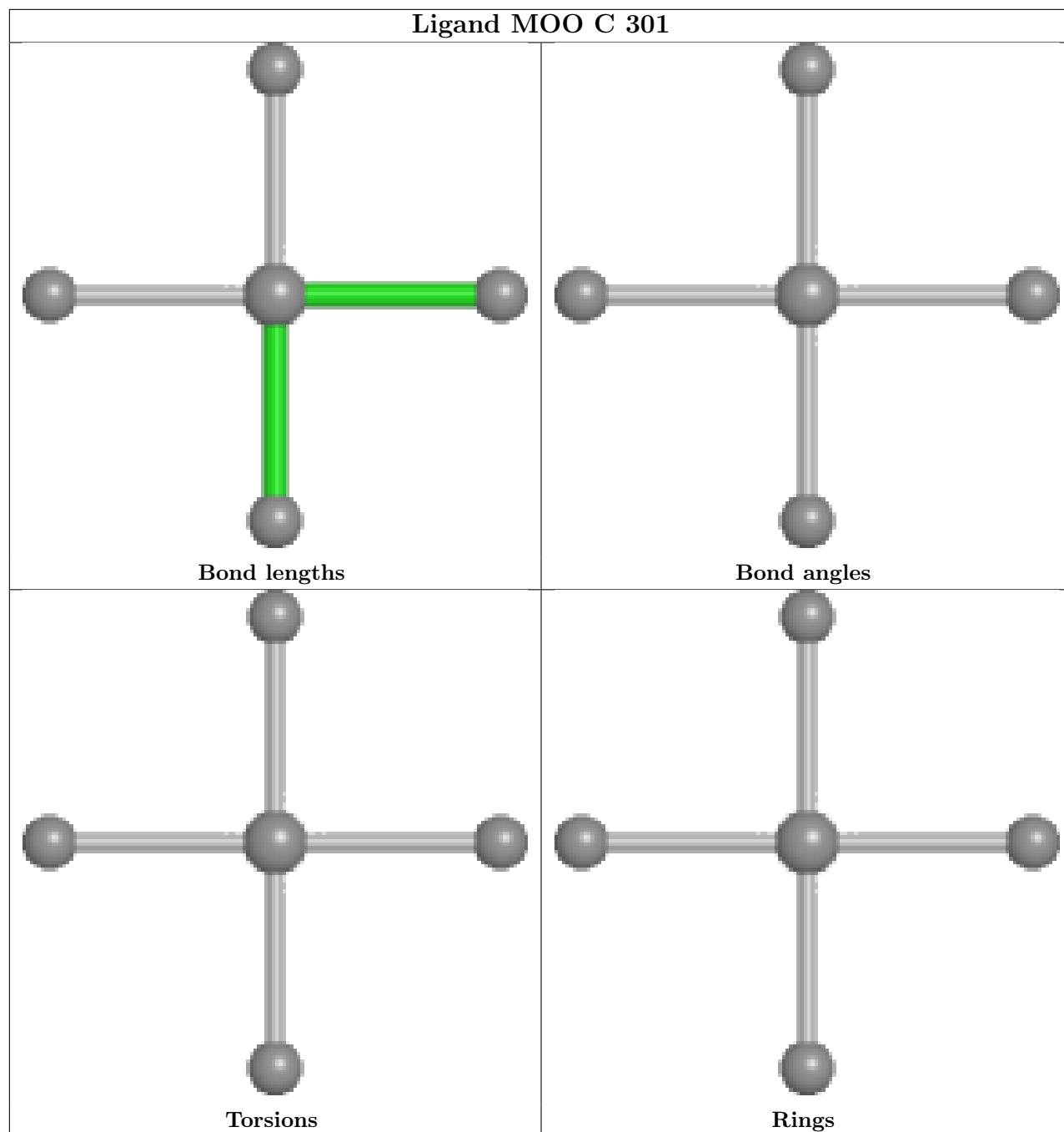
4 monomers are involved in 8 short contacts:

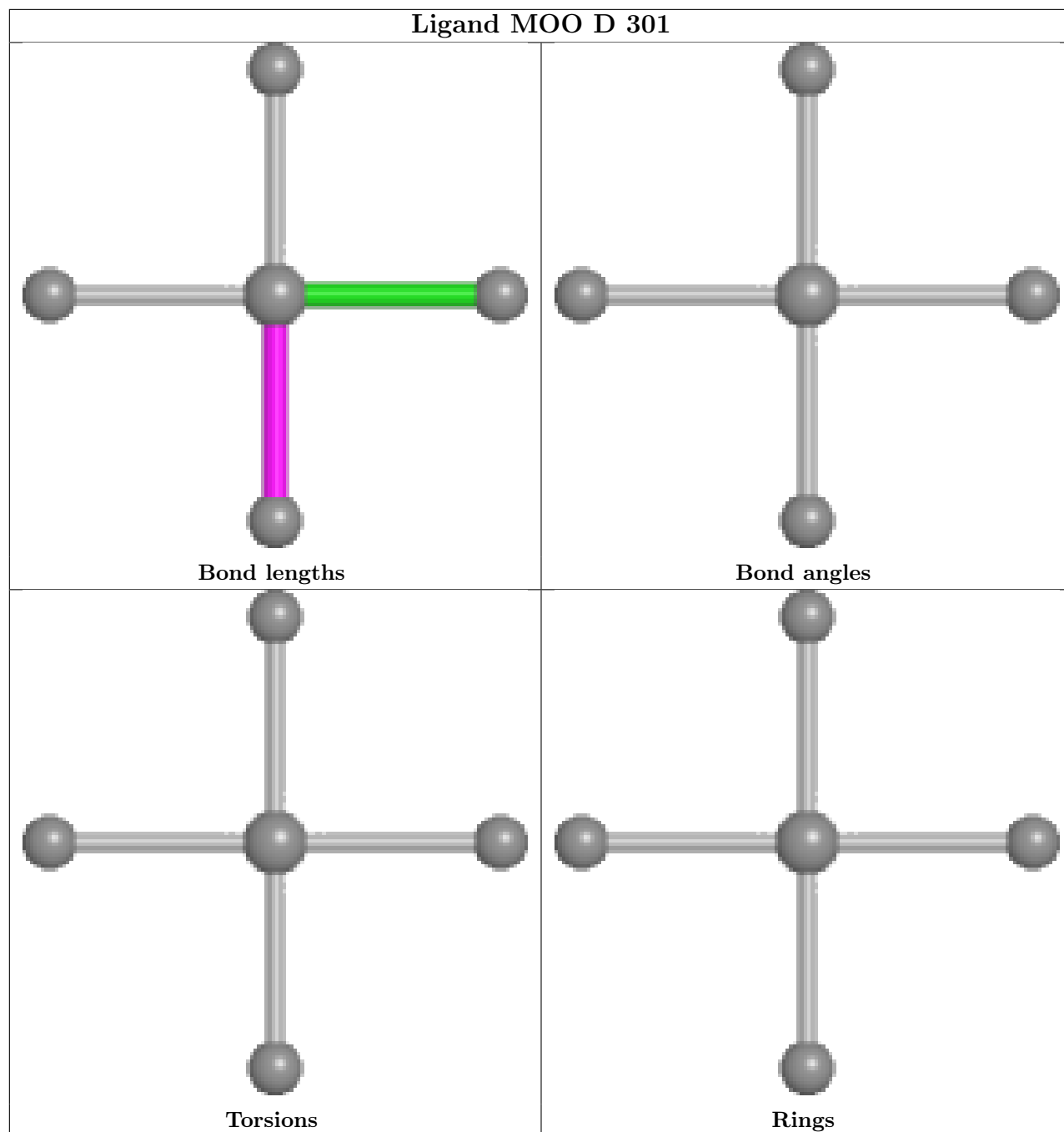
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	301	MOO	2	0
2	B	301	MOO	2	0
2	C	301	MOO	3	0
2	D	301	MOO	1	0

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	226/241 (93%)	1.07	29 (12%) 7 9	11, 24, 36, 41	1 (0%)
1	B	229/241 (95%)	1.07	22 (9%) 13 15	11, 24, 36, 55	2 (0%)
1	C	227/241 (94%)	1.05	23 (10%) 12 14	11, 24, 36, 46	1 (0%)
1	D	228/241 (94%)	1.05	25 (10%) 10 12	11, 24, 36, 48	2 (0%)
All	All	910/964 (94%)	1.06	99 (10%) 10 12	11, 24, 36, 55	6 (0%)

All (99) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	237	ALA	8.1
1	B	237	ALA	6.2
1	A	235	MET	4.3
1	D	237	ALA	4.3
1	B	236	LEU	4.2
1	B	10	GLU	3.8
1	B	238	PRO	3.6
1	C	236	LEU	3.6
1	B	37	VAL	3.5
1	D	236	LEU	3.4
1	A	37	VAL	3.4
1	A	163	PHE	3.4
1	D	235	MET	3.3
1	B	235	MET	3.3
1	D	39	VAL	3.3
1	B	234	SER	3.2
1	D	99	GLY	3.2
1	C	11	SER	3.1
1	A	167	GLY	3.0
1	D	100	GLN	3.0
1	A	39	VAL	3.0

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Mol	Chain	Res	Type	RSRZ
1	D	163	PHE	3.0
1	D	10	GLU	3.0
1	C	229	LYS	2.9
1	B	11	SER	2.9
1	B	180	ALA	2.8
1	A	99	GLY	2.8
1	A	12	VAL	2.8
1	A	210	SER	2.7
1	D	167	GLY	2.7
1	C	109	ALA	2.7
1	B	138	PHE	2.7
1	D	161	LEU	2.6
1	B	227	VAL	2.6
1	D	29	LYS	2.6
1	C	138	PHE	2.6
1	C	37	VAL	2.6
1	B	226	ALA	2.6
1	D	94	LEU	2.6
1	A	114	LYS	2.6
1	B	100	GLN	2.5
1	C	210	SER	2.5
1	B	61	ILE	2.5
1	A	27	LEU	2.5
1	A	67	THR	2.5
1	C	194	ASN	2.5
1	C	166	ARG	2.5
1	A	81	ALA	2.4
1	C	99	GLY	2.4
1	A	82	ASN	2.4
1	D	31	TYR	2.4
1	A	36	LYS	2.4
1	C	167	GLY	2.4
1	D	82	ASN	2.4
1	C	181	ILE	2.4
1	D	138	PHE	2.4
1	A	98	LYS	2.4
1	B	126	LYS	2.4
1	A	224	ALA	2.4
1	D	169	CYS	2.4
1	B	99	GLY	2.4
1	C	168	GLU	2.4
1	A	89	GLY	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	181	ILE	2.3
1	A	100	GLN	2.3
1	A	40	LYS	2.3
1	D	101	SER	2.3
1	C	213	ALA	2.2
1	B	143	TRP	2.2
1	C	143	TRP	2.2
1	A	236	LEU	2.2
1	C	228	PHE	2.2
1	A	11	SER	2.2
1	D	227	VAL	2.2
1	A	145	ASN	2.2
1	A	173	ILE	2.2
1	C	30	ILE	2.2
1	D	30	ILE	2.2
1	D	61	ILE	2.2
1	D	217	GLN	2.2
1	C	108	LYS	2.2
1	D	214	LYS	2.1
1	D	37	VAL	2.1
1	C	100	GLN	2.1
1	D	11	SER	2.1
1	C	112	PRO	2.1
1	B	36	LYS	2.1
1	A	38	GLU	2.1
1	C	163	PHE	2.0
1	B	109	ALA	2.0
1	A	166	ARG	2.0
1	A	94	LEU	2.0
1	B	166	ARG	2.0
1	A	106	LEU	2.0
1	A	162	ASN	2.0
1	D	75	ASN	2.0
1	A	15	TYR	2.0
1	B	31	TYR	2.0
1	C	31	TYR	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

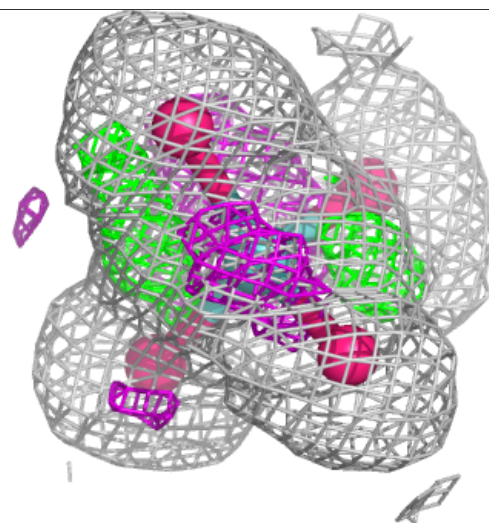
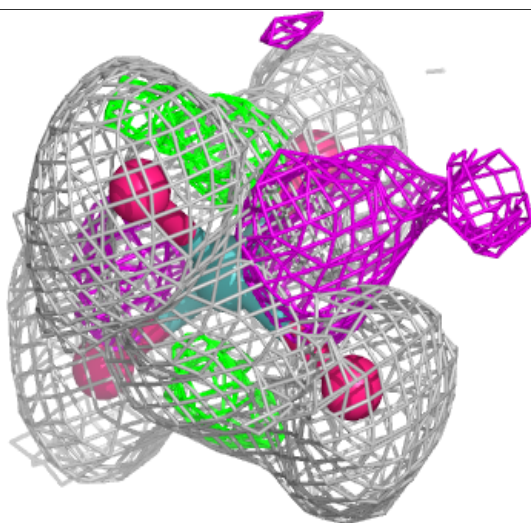
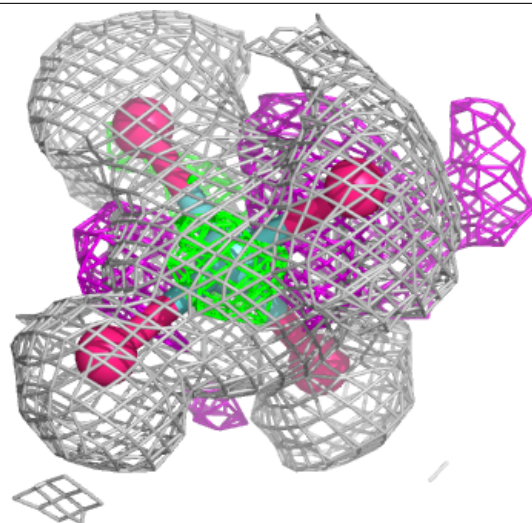
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	MOO	A	301	5/5	0.99	0.06	15,16,18,19	0
2	MOO	C	301	5/5	0.99	0.05	14,14,15,16	0
2	MOO	D	301	5/5	0.99	0.05	15,15,16,16	0
2	MOO	B	301	5/5	1.00	0.04	14,14,15,16	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.

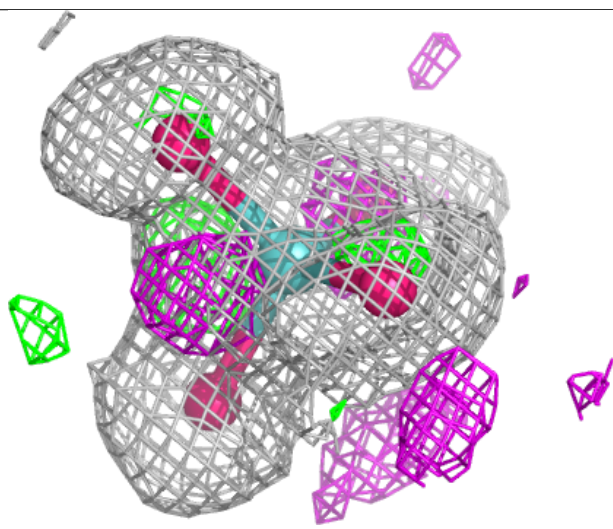
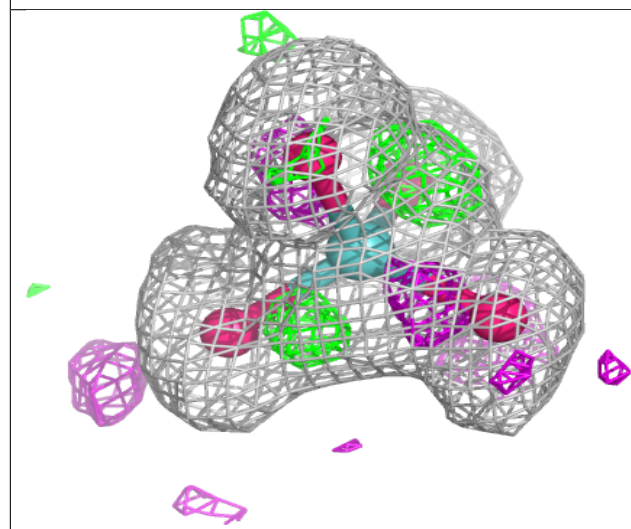
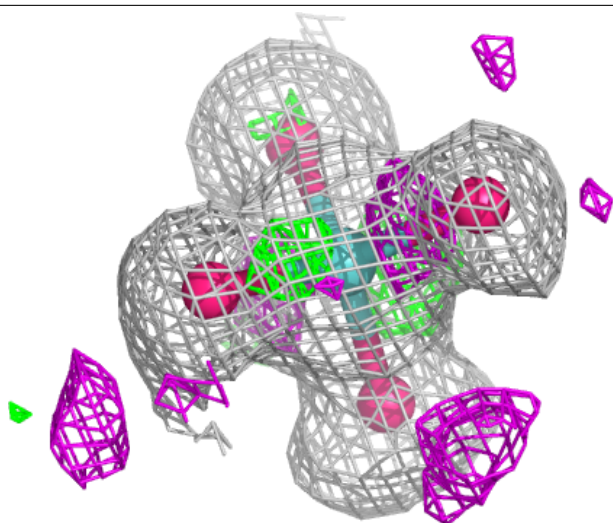
Electron density around MOO A 301:

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



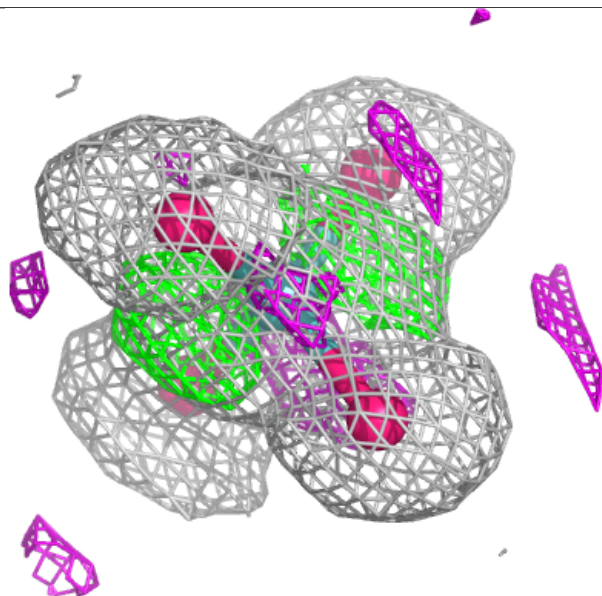
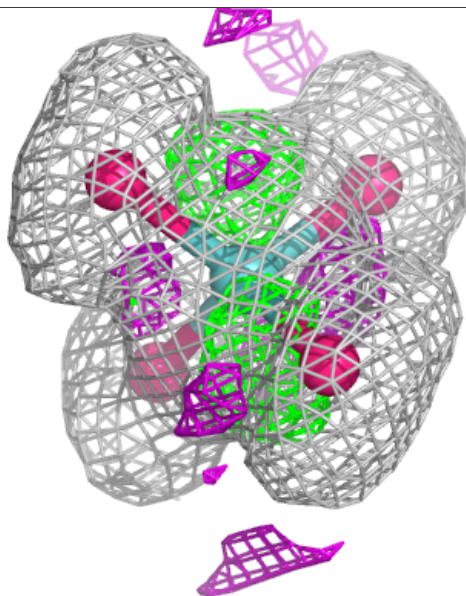
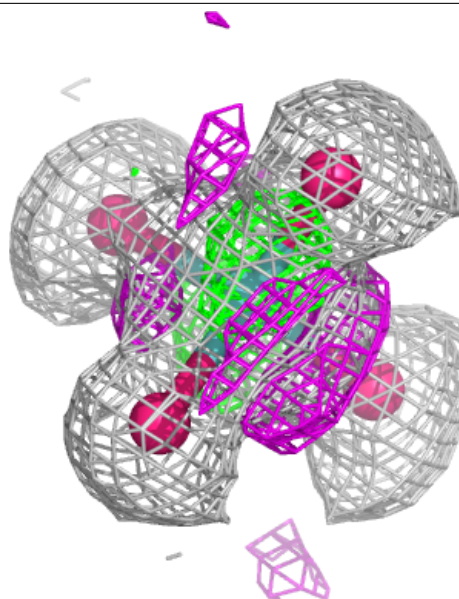
Electron density around MOO C 301:

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



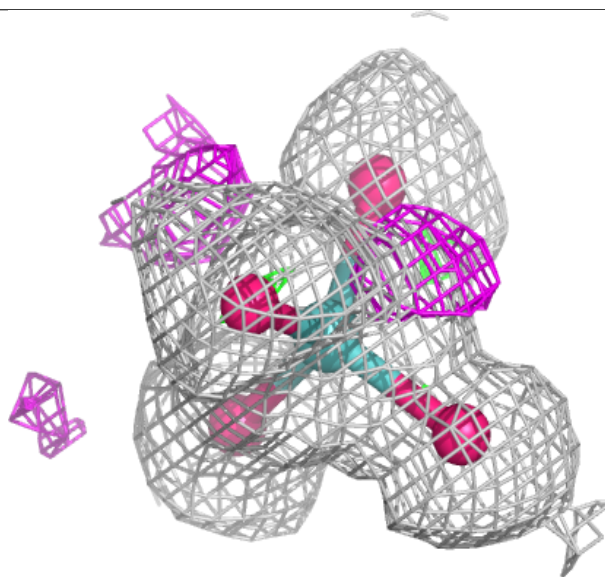
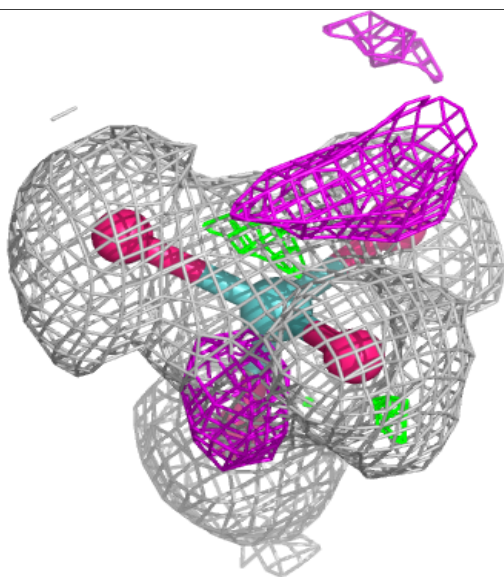
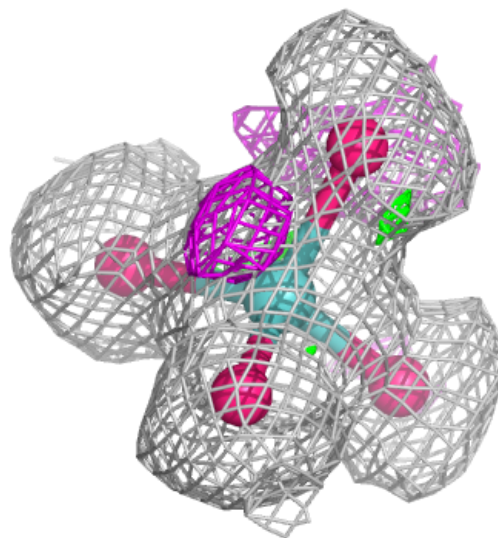
Electron density around MOO D 301:

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



Electron density around MOO 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.