



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 05:47 PM UTC

PDB ID : 8RIP / pdb_00008rip
Title : Beta-keto acid cleavage enzyme from *Paracoccus denitrificans* with bound malonate and Coenzyme A
Authors : Marchal, D.G.; Zarzycki, J.; Erb, T.J.
Deposited on : 2023-12-19
Resolution : 1.81 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

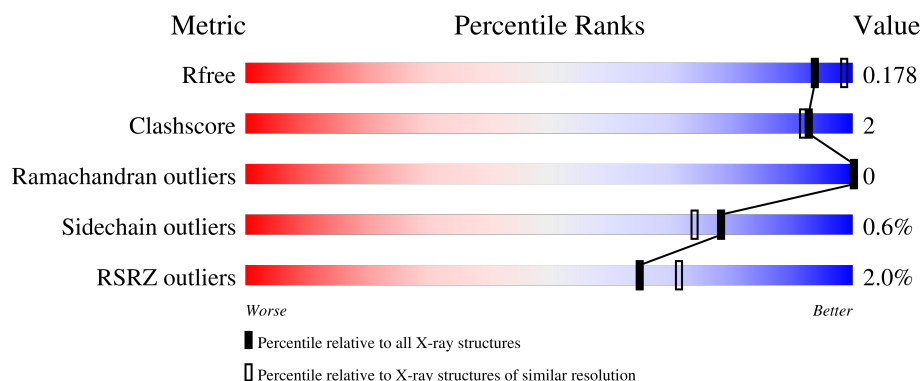
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.81 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	1112 (1.82-1.82)
Clashscore	190562	1148 (1.82-1.82)
Ramachandran outliers	187476	1140 (1.82-1.82)
Sidechain outliers	187428	1140 (1.82-1.82)
RSRZ outliers	180081	1112 (1.82-1.82)

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	334	% 90% 7%
1	B	334	2% 90% 7%
1	C	334	3% 89% 8%
1	D	334	% 90% 8%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 10467 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 3-keto-5-aminohexanoate cleavage protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	311	Total	C	N	O	S	0	0	0
			2371	1488	429	441	13			
1	B	310	Total	C	N	O	S	0	0	0
			2360	1482	425	440	13			
1	C	308	Total	C	N	O	S	0	0	0
			2346	1474	423	437	12			
1	D	308	Total	C	N	O	S	0	0	0
			2346	1474	423	437	12			

There are 96 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-23	MET	-	initiating methionine	UNP A1B802
A	-22	GLY	-	expression tag	UNP A1B802
A	-21	HIS	-	expression tag	UNP A1B802
A	-20	HIS	-	expression tag	UNP A1B802
A	-19	HIS	-	expression tag	UNP A1B802
A	-18	HIS	-	expression tag	UNP A1B802
A	-17	HIS	-	expression tag	UNP A1B802
A	-16	HIS	-	expression tag	UNP A1B802
A	-15	HIS	-	expression tag	UNP A1B802
A	-14	HIS	-	expression tag	UNP A1B802
A	-13	HIS	-	expression tag	UNP A1B802
A	-12	HIS	-	expression tag	UNP A1B802
A	-11	SER	-	expression tag	UNP A1B802
A	-10	SER	-	expression tag	UNP A1B802
A	-9	GLY	-	expression tag	UNP A1B802
A	-8	HIS	-	expression tag	UNP A1B802
A	-7	ILE	-	expression tag	UNP A1B802
A	-6	GLU	-	expression tag	UNP A1B802
A	-5	GLY	-	expression tag	UNP A1B802
A	-4	ARG	-	expression tag	UNP A1B802
A	-3	HIS	-	expression tag	UNP A1B802

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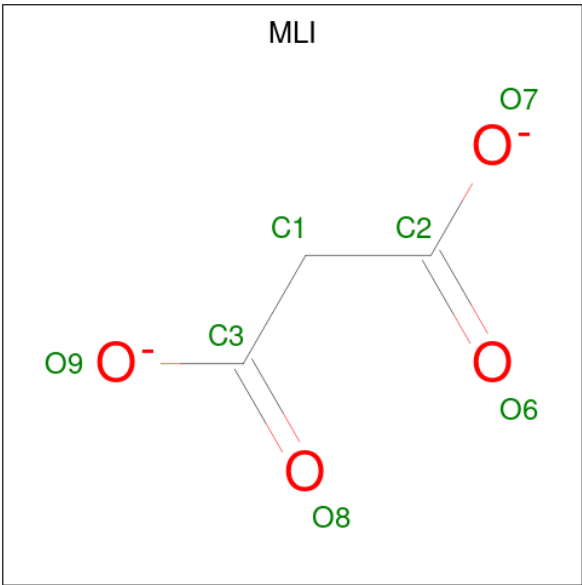
Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	expression tag	UNP A1B802
A	-1	SER	-	expression tag	UNP A1B802
A	0	ARG	-	expression tag	UNP A1B802
B	-23	MET	-	initiating methionine	UNP A1B802
B	-22	GLY	-	expression tag	UNP A1B802
B	-21	HIS	-	expression tag	UNP A1B802
B	-20	HIS	-	expression tag	UNP A1B802
B	-19	HIS	-	expression tag	UNP A1B802
B	-18	HIS	-	expression tag	UNP A1B802
B	-17	HIS	-	expression tag	UNP A1B802
B	-16	HIS	-	expression tag	UNP A1B802
B	-15	HIS	-	expression tag	UNP A1B802
B	-14	HIS	-	expression tag	UNP A1B802
B	-13	HIS	-	expression tag	UNP A1B802
B	-12	HIS	-	expression tag	UNP A1B802
B	-11	SER	-	expression tag	UNP A1B802
B	-10	SER	-	expression tag	UNP A1B802
B	-9	GLY	-	expression tag	UNP A1B802
B	-8	HIS	-	expression tag	UNP A1B802
B	-7	ILE	-	expression tag	UNP A1B802
B	-6	GLU	-	expression tag	UNP A1B802
B	-5	GLY	-	expression tag	UNP A1B802
B	-4	ARG	-	expression tag	UNP A1B802
B	-3	HIS	-	expression tag	UNP A1B802
B	-2	GLY	-	expression tag	UNP A1B802
B	-1	SER	-	expression tag	UNP A1B802
B	0	ARG	-	expression tag	UNP A1B802
C	-23	MET	-	initiating methionine	UNP A1B802
C	-22	GLY	-	expression tag	UNP A1B802
C	-21	HIS	-	expression tag	UNP A1B802
C	-20	HIS	-	expression tag	UNP A1B802
C	-19	HIS	-	expression tag	UNP A1B802
C	-18	HIS	-	expression tag	UNP A1B802
C	-17	HIS	-	expression tag	UNP A1B802
C	-16	HIS	-	expression tag	UNP A1B802
C	-15	HIS	-	expression tag	UNP A1B802
C	-14	HIS	-	expression tag	UNP A1B802
C	-13	HIS	-	expression tag	UNP A1B802
C	-12	HIS	-	expression tag	UNP A1B802
C	-11	SER	-	expression tag	UNP A1B802
C	-10	SER	-	expression tag	UNP A1B802
C	-9	GLY	-	expression tag	UNP A1B802

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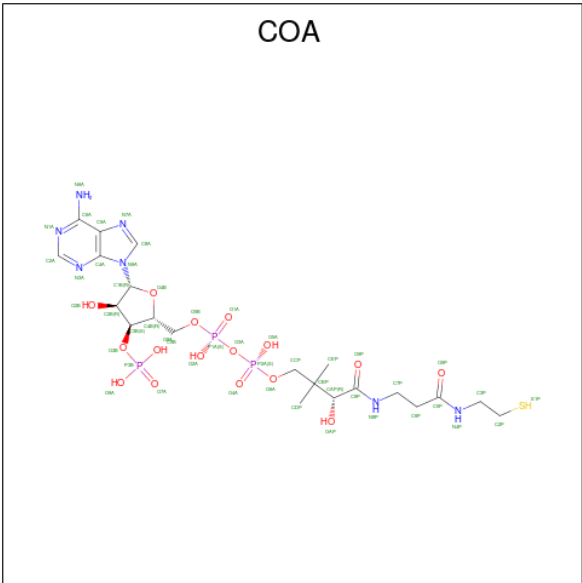
Chain	Residue	Modelled	Actual	Comment	Reference
C	-8	HIS	-	expression tag	UNP A1B802
C	-7	ILE	-	expression tag	UNP A1B802
C	-6	GLU	-	expression tag	UNP A1B802
C	-5	GLY	-	expression tag	UNP A1B802
C	-4	ARG	-	expression tag	UNP A1B802
C	-3	HIS	-	expression tag	UNP A1B802
C	-2	GLY	-	expression tag	UNP A1B802
C	-1	SER	-	expression tag	UNP A1B802
C	0	ARG	-	expression tag	UNP A1B802
D	-23	MET	-	initiating methionine	UNP A1B802
D	-22	GLY	-	expression tag	UNP A1B802
D	-21	HIS	-	expression tag	UNP A1B802
D	-20	HIS	-	expression tag	UNP A1B802
D	-19	HIS	-	expression tag	UNP A1B802
D	-18	HIS	-	expression tag	UNP A1B802
D	-17	HIS	-	expression tag	UNP A1B802
D	-16	HIS	-	expression tag	UNP A1B802
D	-15	HIS	-	expression tag	UNP A1B802
D	-14	HIS	-	expression tag	UNP A1B802
D	-13	HIS	-	expression tag	UNP A1B802
D	-12	HIS	-	expression tag	UNP A1B802
D	-11	SER	-	expression tag	UNP A1B802
D	-10	SER	-	expression tag	UNP A1B802
D	-9	GLY	-	expression tag	UNP A1B802
D	-8	HIS	-	expression tag	UNP A1B802
D	-7	ILE	-	expression tag	UNP A1B802
D	-6	GLU	-	expression tag	UNP A1B802
D	-5	GLY	-	expression tag	UNP A1B802
D	-4	ARG	-	expression tag	UNP A1B802
D	-3	HIS	-	expression tag	UNP A1B802
D	-2	GLY	-	expression tag	UNP A1B802
D	-1	SER	-	expression tag	UNP A1B802
D	0	ARG	-	expression tag	UNP A1B802

- Molecule 2 is MALONATE ION (CCD ID: MLI) (formula: $C_3H_2O_4$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			7	3	4		
2	B	1	Total	C	O	0	0
			7	3	4		
2	C	1	Total	C	O	0	0
			7	3	4		
2	D	1	Total	C	O	0	0
			7	3	4		

- Molecule 3 is COENZYME A (CCD ID: COA) (formula: C₂₁H₃₆N₇O₁₆P₃S) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	S	
			48	21	7	16	3	1	
3	B	1	Total	C	N	O	P	S	
			48	21	7	16	3	1	

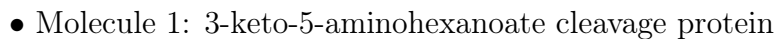
- Molecule 4 is ZINC ION (CCD ID: ZN) (formula: Zn) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Zn		
			1	1	0	0
4	B	1	Total	Zn		
			1	1	0	0
4	C	1	Total	Zn		
			1	1	0	0
4	D	1	Total	Zn		
			1	1	0	0

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	249	Total	O		
			249	249	0	0
5	B	244	Total	O		
			244	244	0	0
5	C	215	Total	O		
			215	215	0	0
5	D	208	Total	O		
			208	208	0	0

- Molecule 1: 3-keto-5-aminohexanoate cleavage protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	79.16Å 134.35Å 140.85Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.99 – 1.81 29.99 – 1.81	Depositor EDS
% Data completeness (in resolution range)	99.8 (29.99-1.81) 99.8 (29.99-1.81)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.24 (at 1.80Å)	Xtriage
Refinement program	PHENIX 1.20.1_4487	Depositor
R, R_{free}	0.158 , 0.178 0.158 , 0.178	Depositor DCC
R_{free} test set	2008 reflections (1.46%)	wwPDB-VP
Wilson B-factor (Å ²)	28.1	Xtriage
Anisotropy	0.310	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 35.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.002 for -h,l,k	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	10467	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.90% 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: MLI, COA, ZN

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	0.32	0/2420	0.51	0/3275
1	B	0.31	0/2409	0.50	0/3261
1	C	0.31	0/2395	0.50	0/3243
1	D	0.32	0/2395	0.49	0/3243
All	All	0.31	0/9619	0.50	0/13022

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	2371	0	2335	8	0
1	B	2360	0	2322	5	0
1	C	2346	0	2305	10	0
1	D	2346	0	2305	6	0
2	A	7	0	2	0	0
2	B	7	0	2	0	0
2	C	7	0	2	0	0
2	D	7	0	2	0	0
3	A	48	0	32	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	48	0	32	2	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
4	C	1	0	0	0	0
4	D	1	0	0	0	0
5	A	249	0	0	1	0
5	B	244	0	0	2	0
5	C	215	0	0	2	0
5	D	208	0	0	1	0
All	All	10467	0	9339	29	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 2.

All (29) 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:C:121:PHE:N	1:C:122:PRO:HD2	2.00	0.77
1:B:202:LEU:HD11	3:B:402:COA:H21	1.72	0.71
1:C:121:PHE:CZ	1:C:203:MET:HE1	2.29	0.67
1:C:120:LEU:C	1:C:122:PRO:HD2	2.27	0.58
1:A:233:LEU:C	1:A:233:LEU:HD23	2.30	0.56
1:B:121:PHE:N	1:B:122:PRO:HD2	2.20	0.56
1:B:222:ARG:NH2	5:B:502:HOH:O	2.25	0.53
1:D:121:PHE:N	1:D:122:PRO:CD	2.72	0.52
1:C:121:PHE:N	1:C:122:PRO:CD	2.72	0.52
1:C:76:GLN:NE2	5:C:505:HOH:O	2.42	0.52
1:D:120:LEU:C	1:D:122:PRO:HD2	2.34	0.52
1:D:121:PHE:O	1:D:124:LEU:HB2	2.10	0.51
1:A:184:PHE:HD1	1:A:189:LEU:HD12	1.75	0.51
1:C:121:PHE:CZ	1:C:203:MET:CE	2.95	0.50
1:B:263:ALA:HA	5:B:501:HOH:O	2.10	0.50
1:B:15:GLY:HA2	1:B:51:HIS:CD2	2.48	0.49
1:D:75:LYS:NZ	5:D:503:HOH:O	2.43	0.45
1:D:121:PHE:CZ	1:D:203:MET:HE1	2.51	0.44
1:A:15:GLY:HA2	1:A:51:HIS:CD2	2.52	0.44
1:A:131:LEU:HD11	1:A:140:LEU:HD12	2.00	0.42
1:A:233:LEU:C	1:A:233:LEU:CD2	2.93	0.41
3:B:402:COA:H51A	3:B:402:COA:H8A	2.03	0.41
1:A:281:GLN:HG3	1:C:281:GLN:HG3	2.02	0.41
1:A:126:ARG:HD2	1:A:127:TYR:CZ	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:267:ARG:NH2	5:A:509:HOH:O	2.49	0.41
1:C:277:ARG:NH1	5:C:502:HOH:O	2.31	0.41
1:C:15:GLY:HA2	1:C:51:HIS:CD2	2.56	0.41
1:D:123:MET:HE3	1:D:123:MET:HB2	2.00	0.40
1:C:184:PHE:HD1	1:C:189:LEU:HD12	1.87	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	309/334 (92%)	303 (98%)	6 (2%)	0	100	100
1	B	308/334 (92%)	302 (98%)	6 (2%)	0	100	100
1	C	306/334 (92%)	299 (98%)	7 (2%)	0	100	100
1	D	306/334 (92%)	298 (97%)	8 (3%)	0	100	100
All	All	1229/1336 (92%)	1202 (98%)	27 (2%)	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	A	242/261 (93%)	240 (99%)	2 (1%)	73	67

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	241/261 (92%)	240 (100%)	1 (0%)	84	81
1	C	239/261 (92%)	238 (100%)	1 (0%)	84	81
1	D	239/261 (92%)	237 (99%)	2 (1%)	73	67
All	All	961/1044 (92%)	955 (99%)	6 (1%)	78	74

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	175	THR
1	A	233	LEU
1	B	175	THR
1	C	175	THR
1	D	175	THR
1	D	214	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	180	ASN
1	B	180	ASN
1	C	68	ASN
1	C	134	GLN
1	C	180	ASN
1	D	102	GLN
1	D	130	GLN
1	D	180	ASN

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

Of 10 ligands modelled in this entry, 4 are monoatomic - leaving 6 for Mogul analysis.

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	MLI	A	401	4	6,6,6	1.23	0	7,7,7	1.27	0
2	MLI	C	401	4	6,6,6	1.19	0	7,7,7	1.33	0
3	COA	A	402	-	47,50,50	1.71	6 (12%)	69,75,75	1.41	10 (14%)
2	MLI	B	401	4	6,6,6	1.34	0	7,7,7	1.27	0
2	MLI	D	401	4	6,6,6	1.26	0	7,7,7	1.28	0
3	COA	B	402	-	47,50,50	1.78	7 (14%)	69,75,75	1.44	10 (14%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	MLI	A	401	4	-	2/4/4/4	-
2	MLI	C	401	4	-	2/4/4/4	-
3	COA	A	402	-	-	9/48/64/64	0/3/3/3
2	MLI	B	401	4	-	2/4/4/4	-
2	MLI	D	401	4	-	2/4/4/4	-
3	COA	B	402	-	-	11/48/64/64	0/3/3/3

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	402	COA	P1A-O3A	7.39	1.67	1.59
3	A	402	COA	P1A-O3A	6.98	1.67	1.59
3	B	402	COA	P2A-O3A	5.98	1.66	1.59
3	A	402	COA	P2A-O3A	5.31	1.65	1.59
3	A	402	COA	C5A-N7A	-3.27	1.33	1.39
3	B	402	COA	C5A-N7A	-3.20	1.33	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	402	COA	C8A-N9A	-2.32	1.33	1.37
3	A	402	COA	CDP-CBP	-2.27	1.49	1.53
3	B	402	COA	P1A-O5B	2.16	1.67	1.59
3	B	402	COA	CDP-CBP	-2.12	1.49	1.53
3	B	402	COA	C2B-C3B	-2.09	1.48	1.53
3	B	402	COA	C8A-N9A	-2.09	1.34	1.37
3	A	402	COA	C2B-C3B	-2.05	1.48	1.53

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	402	COA	C5A-C4A-N3A	-5.60	119.01	126.72
3	B	402	COA	C5A-C4A-N3A	-5.49	119.16	126.72
3	A	402	COA	C4A-C5A-N7A	-3.45	106.63	110.58
3	B	402	COA	C4A-C5A-N7A	-3.28	106.83	110.58
3	B	402	COA	C1B-N9A-C8A	3.10	133.99	127.09
3	A	402	COA	C1B-N9A-C8A	3.02	133.79	127.09
3	A	402	COA	N3A-C4A-N9A	2.97	132.22	127.17
3	B	402	COA	N3A-C4A-N9A	2.96	132.21	127.17
3	B	402	COA	C4A-N9A-C1B	-2.81	120.05	126.63
3	A	402	COA	C5A-C4A-N9A	2.72	108.77	105.81
3	A	402	COA	C4A-N9A-C1B	-2.67	120.38	126.63
3	B	402	COA	C5A-N7A-C8A	2.63	107.58	103.45
3	B	402	COA	C5A-C4A-N9A	2.59	108.63	105.81
3	A	402	COA	C5A-N7A-C8A	2.59	107.51	103.45
3	B	402	COA	C2A-N3A-C4A	2.34	117.54	111.83
3	A	402	COA	C2A-N3A-C4A	2.34	117.54	111.83
3	B	402	COA	C6A-C5A-C4A	2.19	120.17	117.18
3	A	402	COA	C6A-C5A-C4A	2.19	120.17	117.18
3	B	402	COA	N9A-C8A-N7A	-2.08	110.99	113.94
3	A	402	COA	C6P-C5P-N4P	2.00	120.00	116.34

There are no chirality outliers.

All (28) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	402	COA	CCP-O6A-P2A-O3A
3	A	402	COA	CCP-O6A-P2A-O4A
3	B	402	COA	CCP-O6A-P2A-O3A
3	B	402	COA	CCP-O6A-P2A-O4A
3	B	402	COA	CCP-O6A-P2A-O5A
3	B	402	COA	OAP-CAP-CBP-CCP

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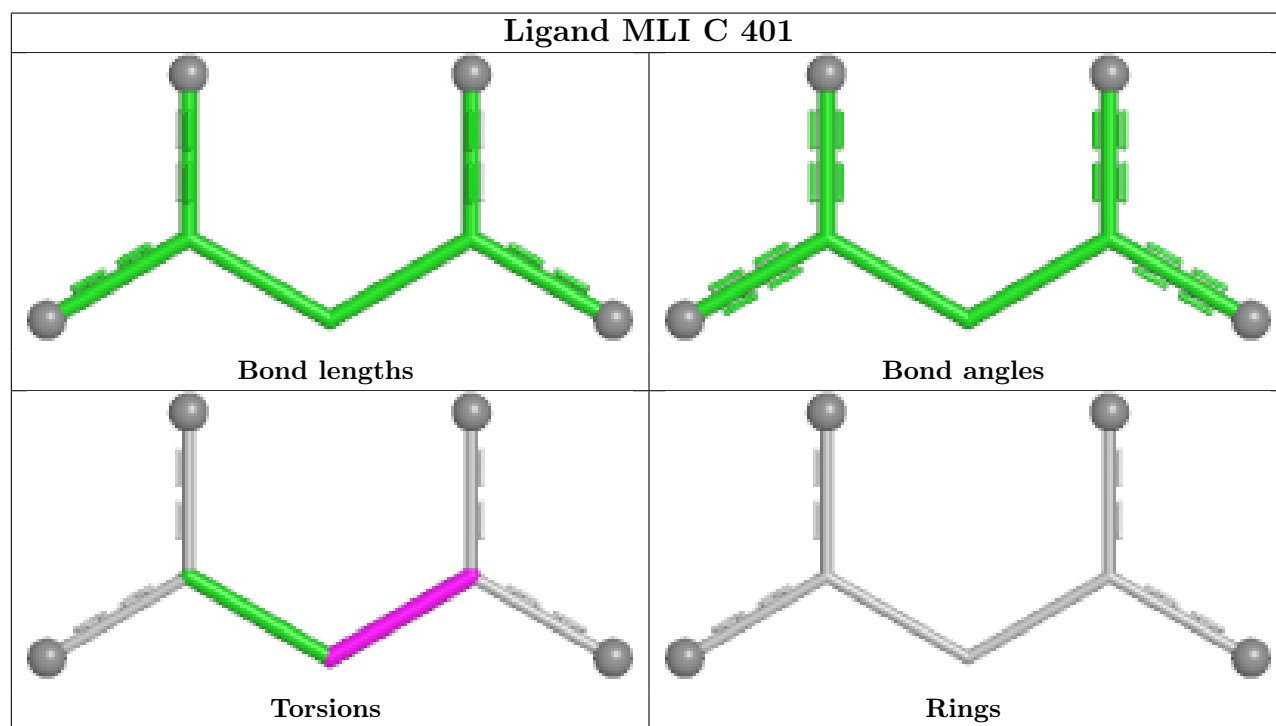
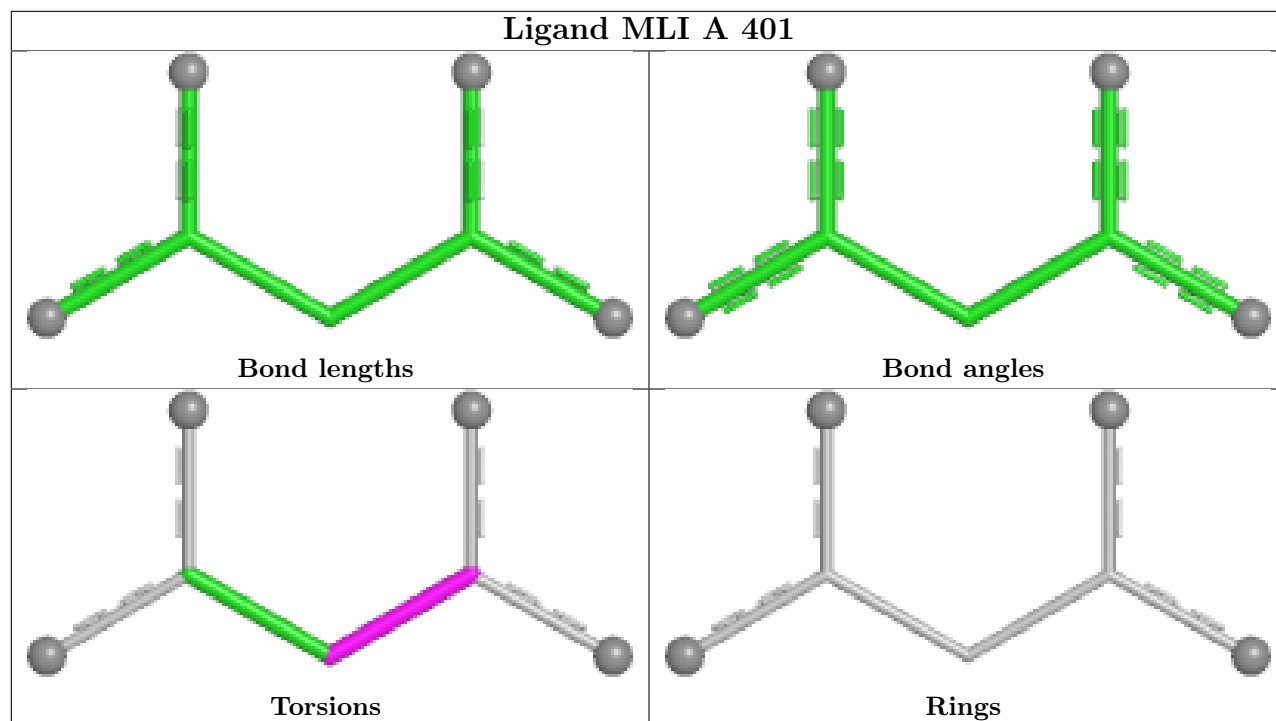
Mol	Chain	Res	Type	Atoms
3	B	402	COA	C9P-CAP-CBP-CCP
3	B	402	COA	OAP-CAP-CBP-CDP
3	B	402	COA	C9P-CAP-CBP-CDP
3	B	402	COA	C9P-CAP-CBP-CEP
3	A	402	COA	OAP-CAP-CBP-CEP
3	B	402	COA	OAP-CAP-CBP-CEP
2	A	401	MLI	C3-C1-C2-O7
2	C	401	MLI	C3-C1-C2-O7
2	C	401	MLI	C3-C1-C2-O6
3	A	402	COA	C9P-CAP-CBP-CDP
2	A	401	MLI	C3-C1-C2-O6
3	A	402	COA	C9P-CAP-CBP-CCP
3	A	402	COA	OAP-CAP-CBP-CCP
2	B	401	MLI	C3-C1-C2-O6
3	B	402	COA	C4B-C5B-O5B-P1A
2	D	401	MLI	C3-C1-C2-O6
2	B	401	MLI	C3-C1-C2-O7
3	A	402	COA	C4B-C5B-O5B-P1A
2	D	401	MLI	C3-C1-C2-O7
3	A	402	COA	C9P-CAP-CBP-CEP
3	B	402	COA	C3B-C4B-C5B-O5B
3	A	402	COA	P1A-O3A-P2A-O5A

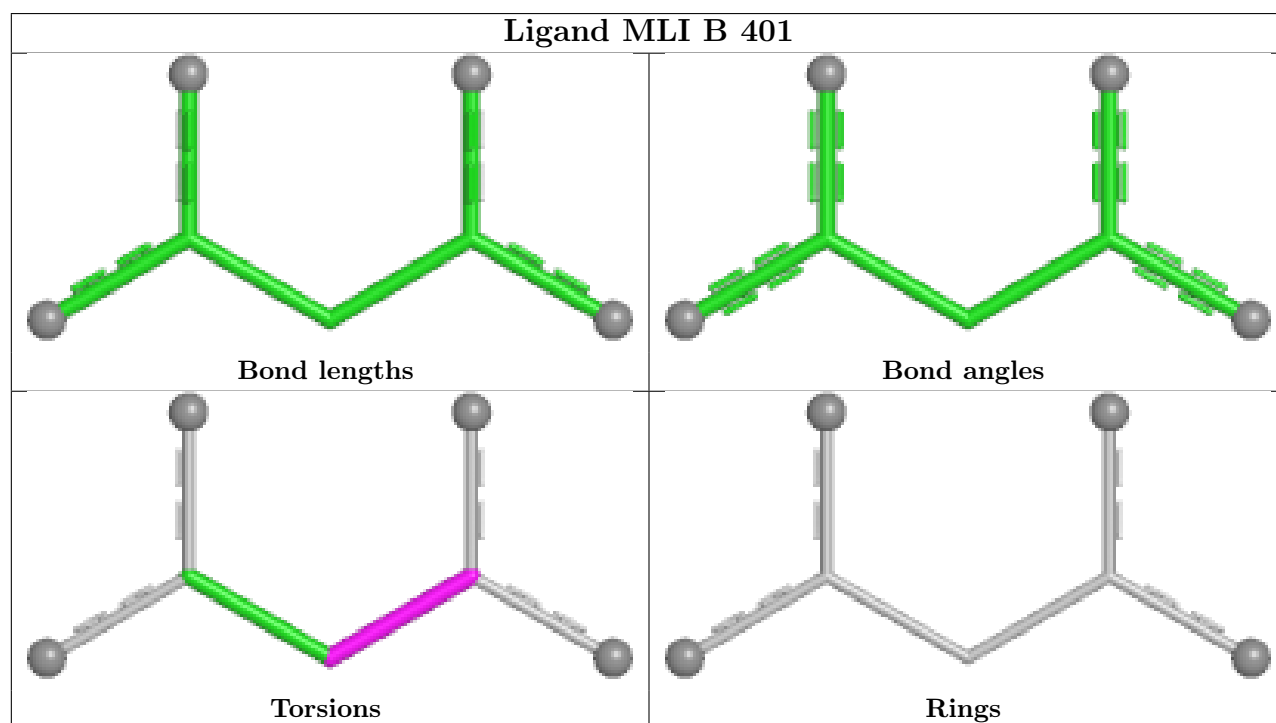
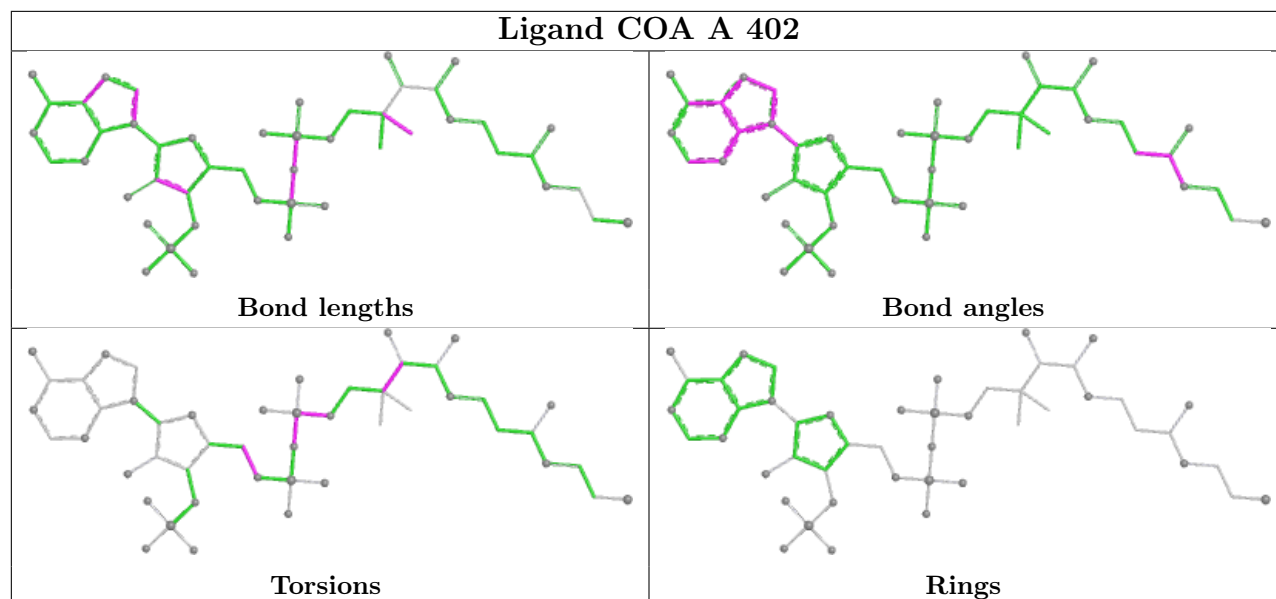
There are no ring outliers.

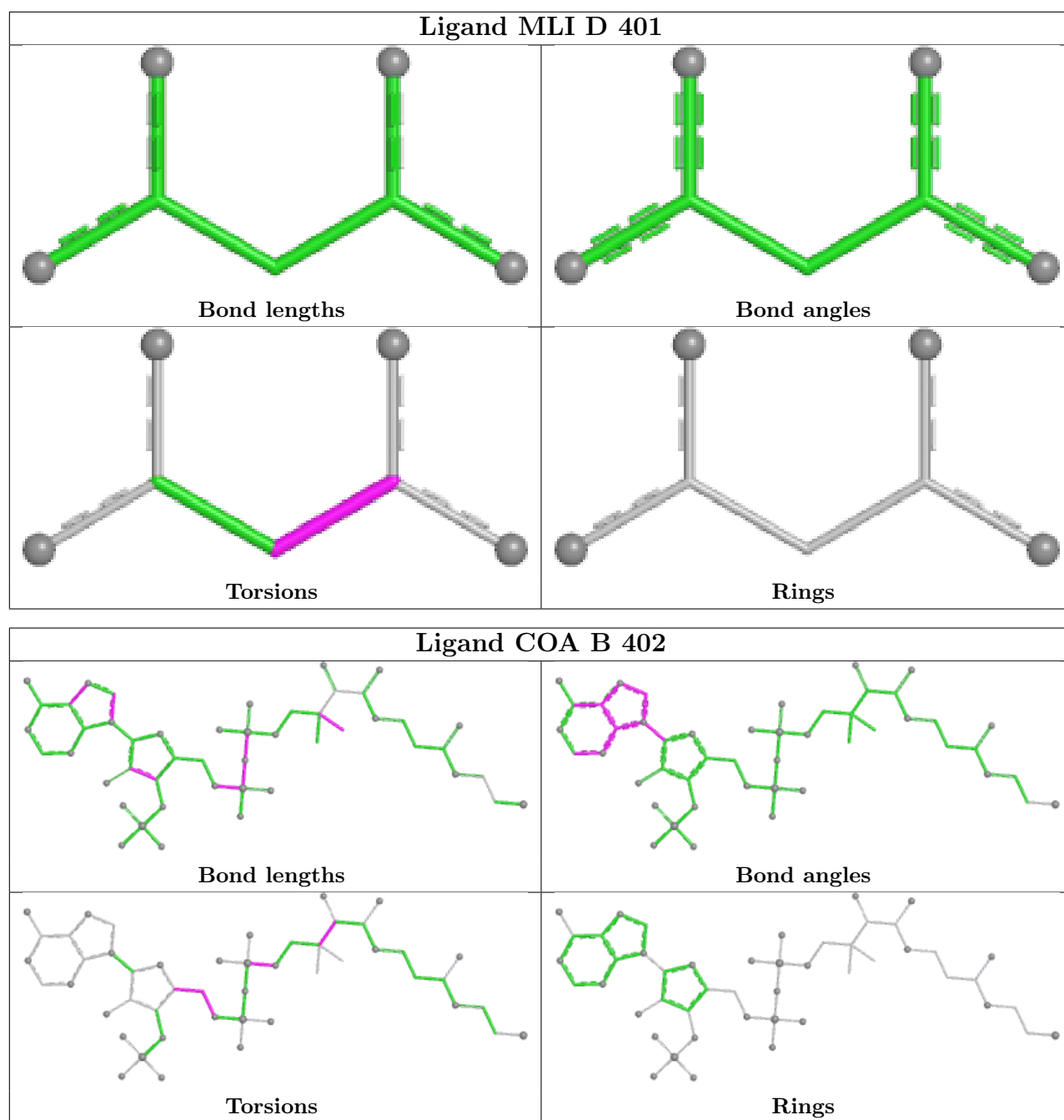
1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	402	COA	2	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	311/334 (93%)	-0.36	3 (0%) 79 84	21, 27, 38, 76	0
1	B	310/334 (92%)	-0.35	6 (1%) 66 72	20, 26, 41, 95	0
1	C	308/334 (92%)	-0.14	11 (3%) 46 53	21, 29, 45, 93	0
1	D	308/334 (92%)	-0.30	5 (1%) 70 77	21, 28, 45, 83	0
All	All	1237/1336 (92%)	-0.29	25 (2%) 65 71	20, 28, 43, 95	0

All (25) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	124	LEU	4.3
1	C	121	PHE	4.2
1	C	122	PRO	3.9
1	B	1	MET	3.7
1	A	0	ARG	3.4
1	B	265	LYS	3.2
1	C	120	LEU	3.2
1	B	264	GLY	3.1
1	A	1	MET	2.9
1	B	4	ASN	2.8
1	C	123	MET	2.8
1	C	266	GLY	2.7
1	C	118	PHE	2.7
1	B	79	ASP	2.5
1	D	132	LYS	2.5
1	C	125	ASN	2.4
1	D	124	LEU	2.4
1	D	266	GLY	2.3
1	C	126	ARG	2.3
1	D	126	ARG	2.3
1	C	267	ARG	2.2

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Mol	Chain	Res	Type	RSRZ
1	C	4	ASN	2.1
1	B	2	SER	2.1
1	A	79	ASP	2.1
1	D	4	ASN	2.0

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](#)

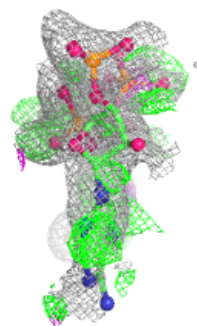
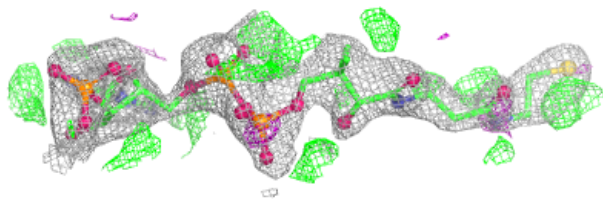
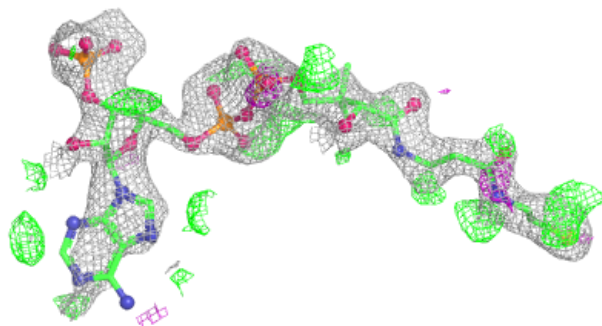
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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
3	COA	B	402	48/48	0.79	0.17	28,50,70,73	48
2	MLI	C	401	7/7	0.94	0.08	27,29,33,36	0
3	COA	A	402	48/48	0.95	0.07	23,30,36,39	48
2	MLI	B	401	7/7	0.96	0.07	24,25,30,32	0
2	MLI	A	401	7/7	0.97	0.06	23,25,29,32	0
2	MLI	D	401	7/7	0.97	0.07	24,26,33,34	0
4	ZN	A	403	1/1	1.00	0.01	25,25,25,25	0
4	ZN	B	403	1/1	1.00	0.01	25,25,25,25	0
4	ZN	C	402	1/1	1.00	0.01	27,27,27,27	0
4	ZN	D	402	1/1	1.00	0.01	25,25,25,25	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.

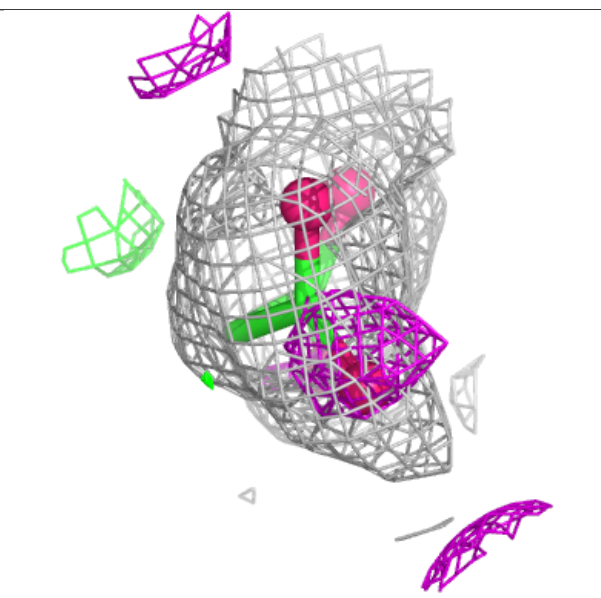
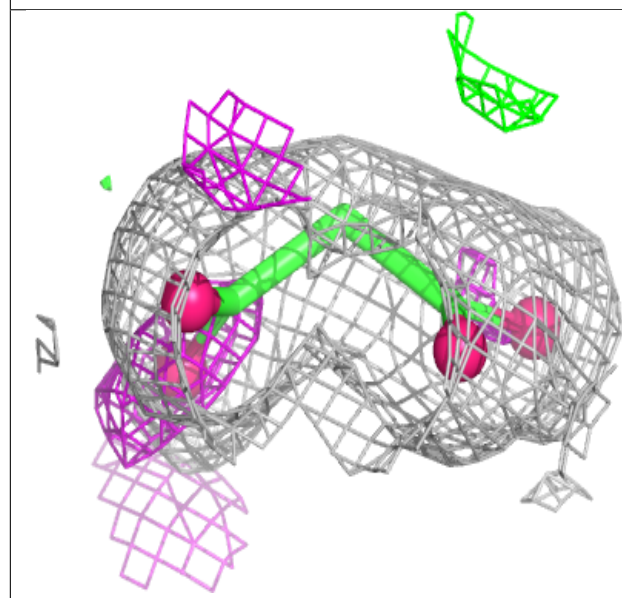
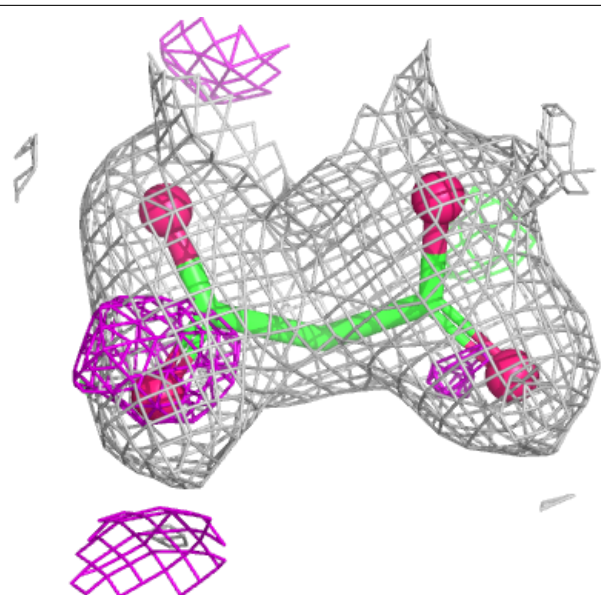
Electron density around COA B 402:

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



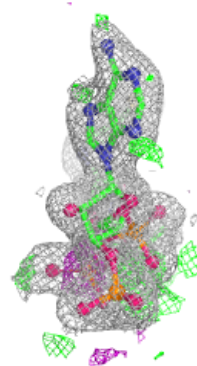
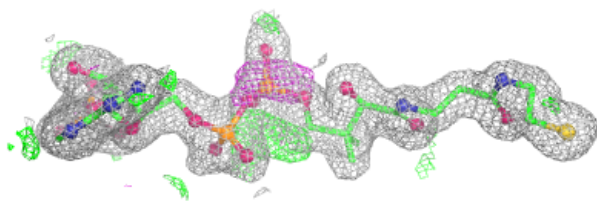
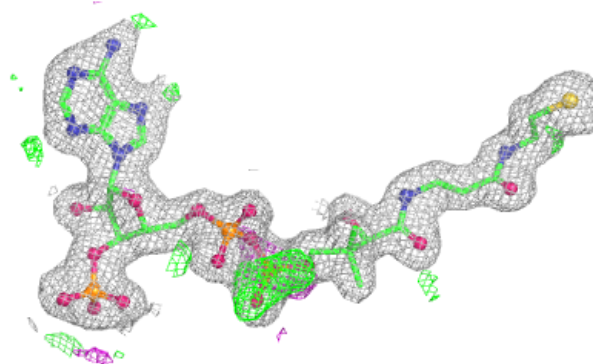
Electron density around MLI C 401:

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and green (positive)



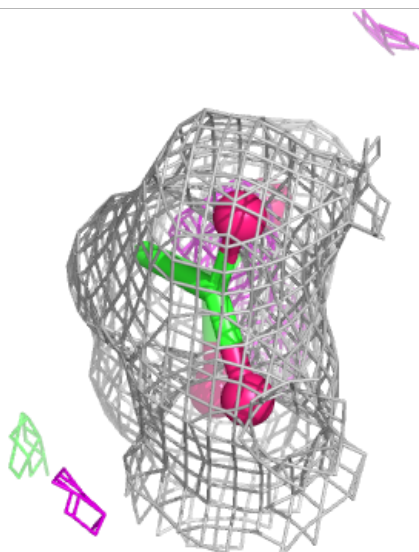
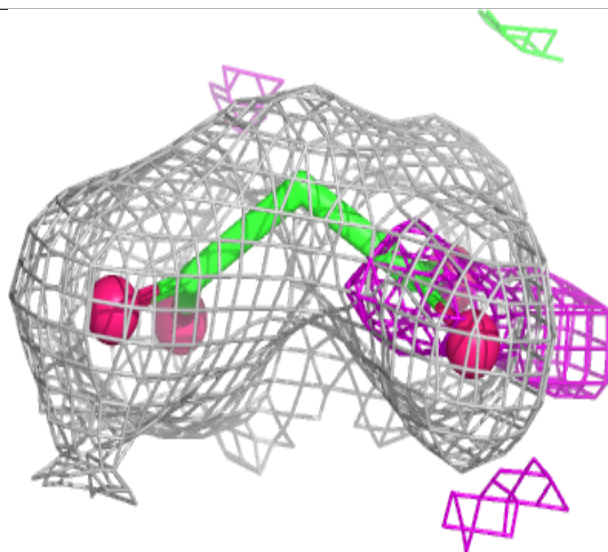
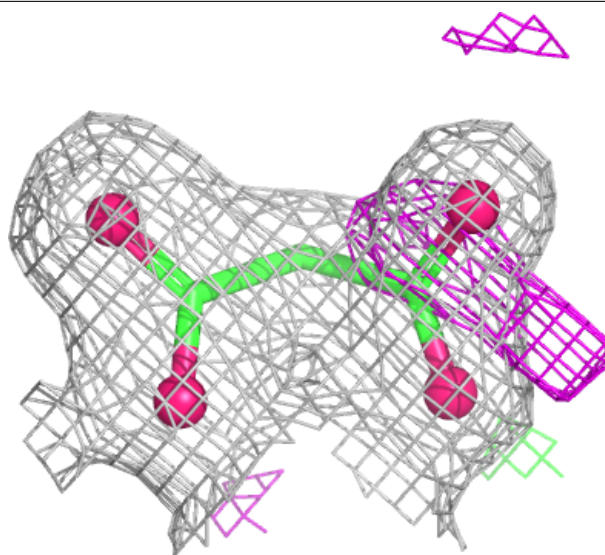
Electron density around COA A 402:

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



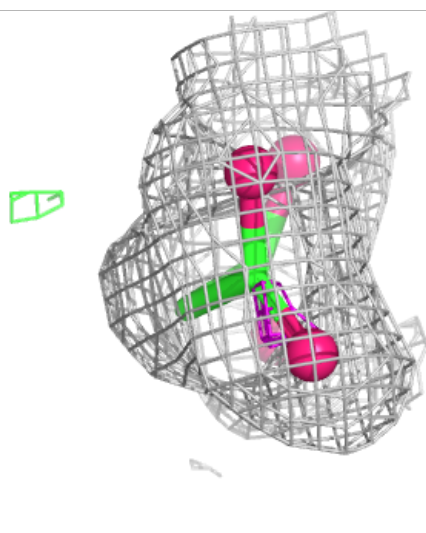
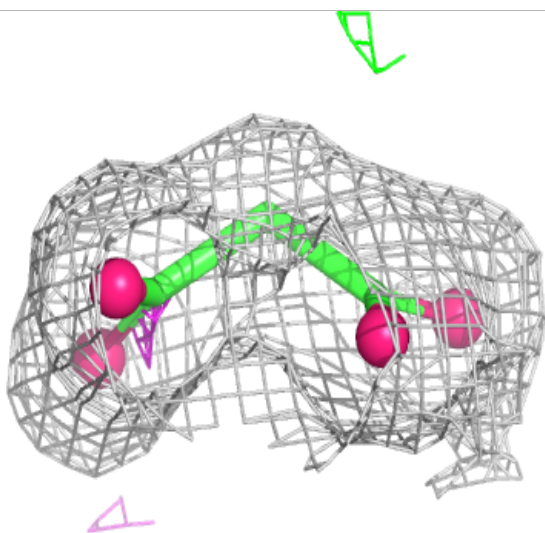
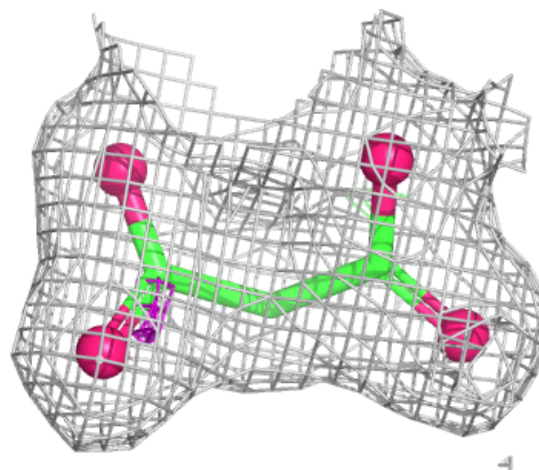
Electron density around MLI B 401:

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and green (positive)



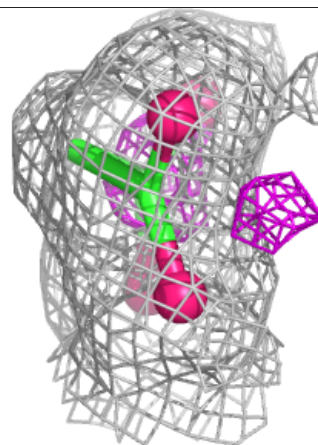
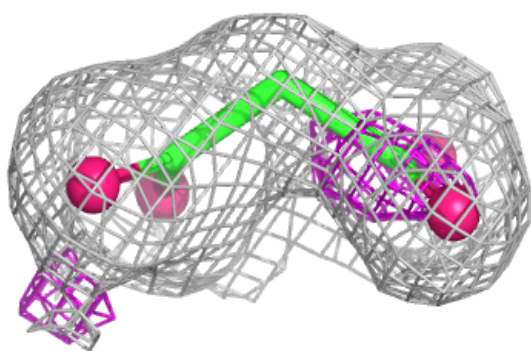
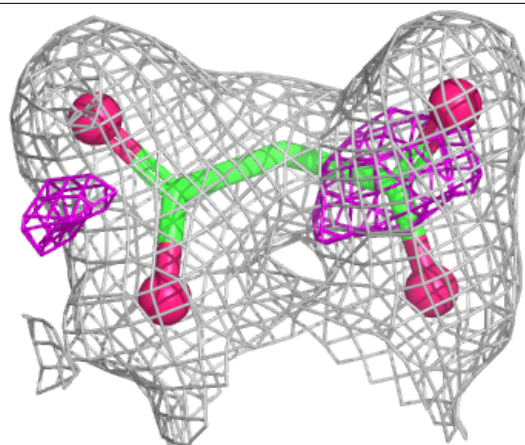
Electron density around MLI A 401:

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



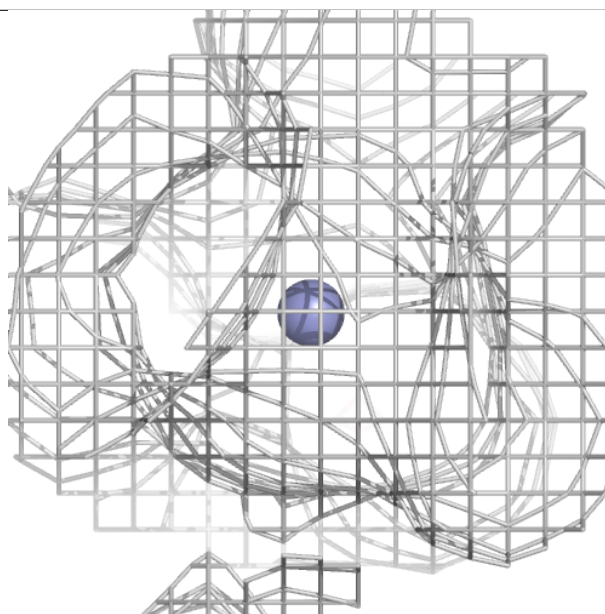
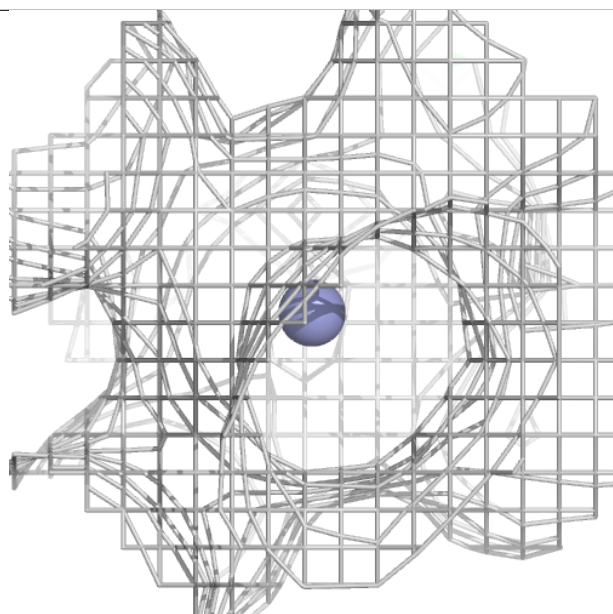
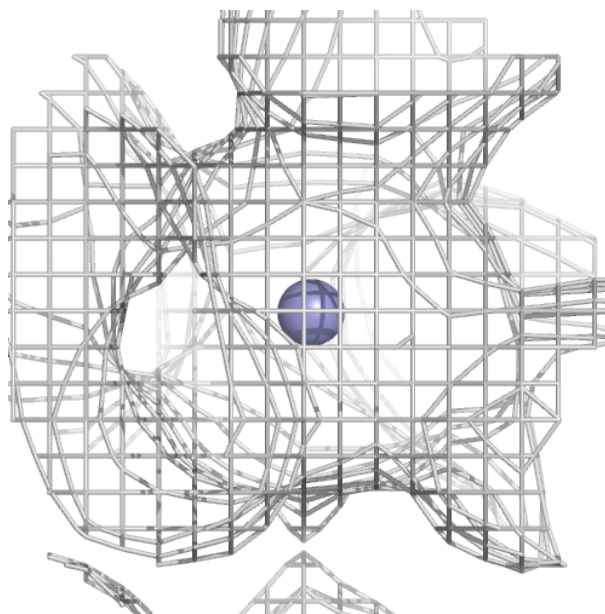
Electron density around MLI D 401:

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



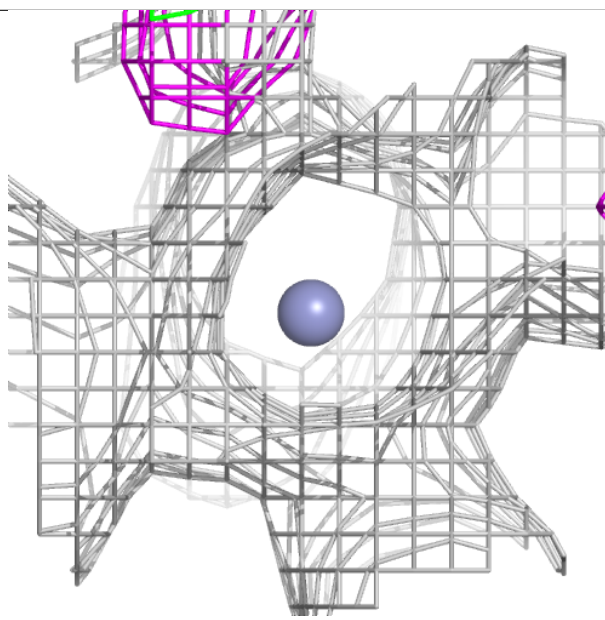
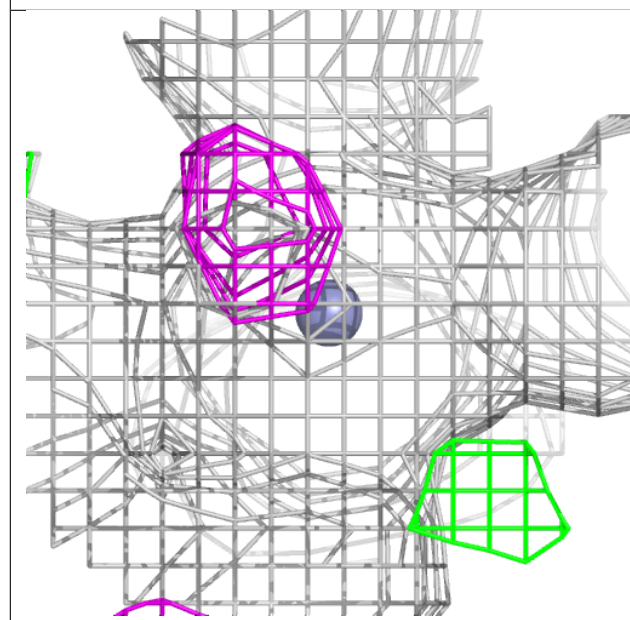
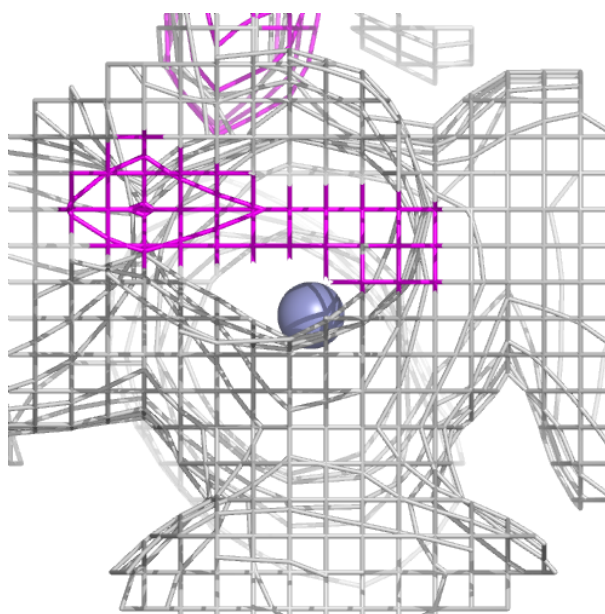
Electron density around ZN A 403:

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and green (positive)



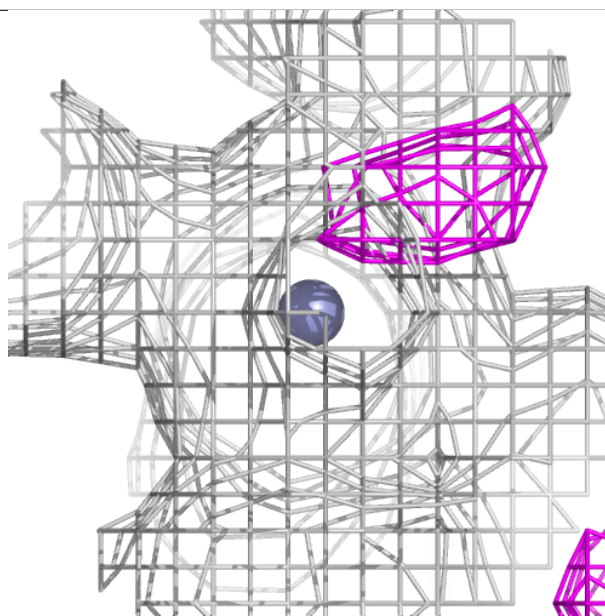
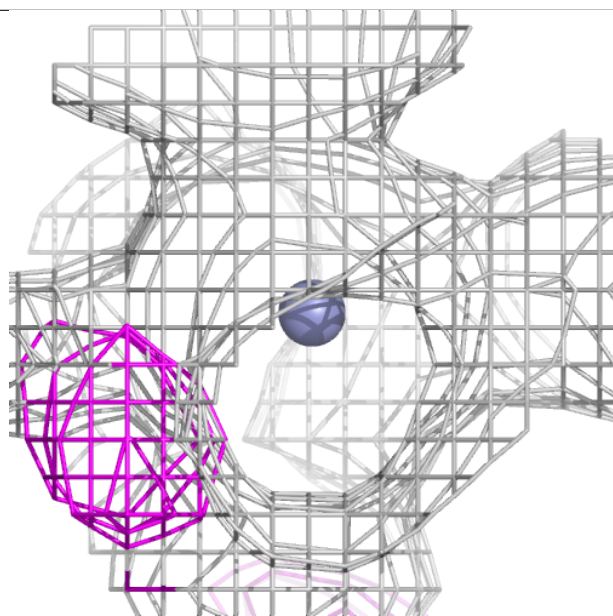
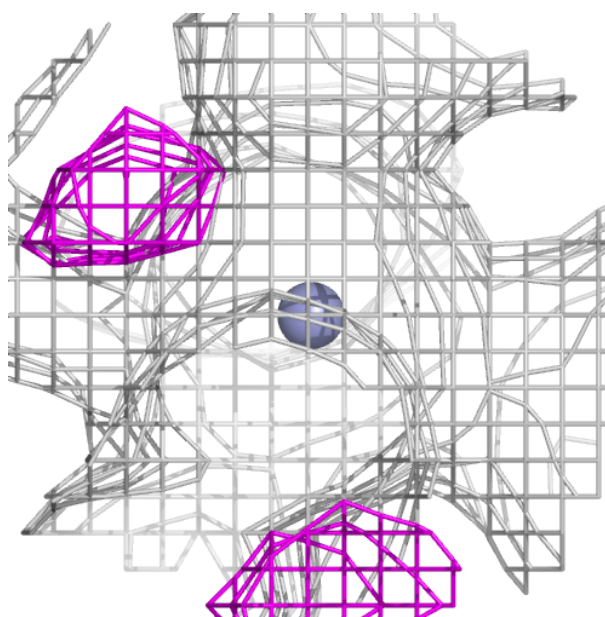
Electron density around ZN B 403:

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and green (positive)



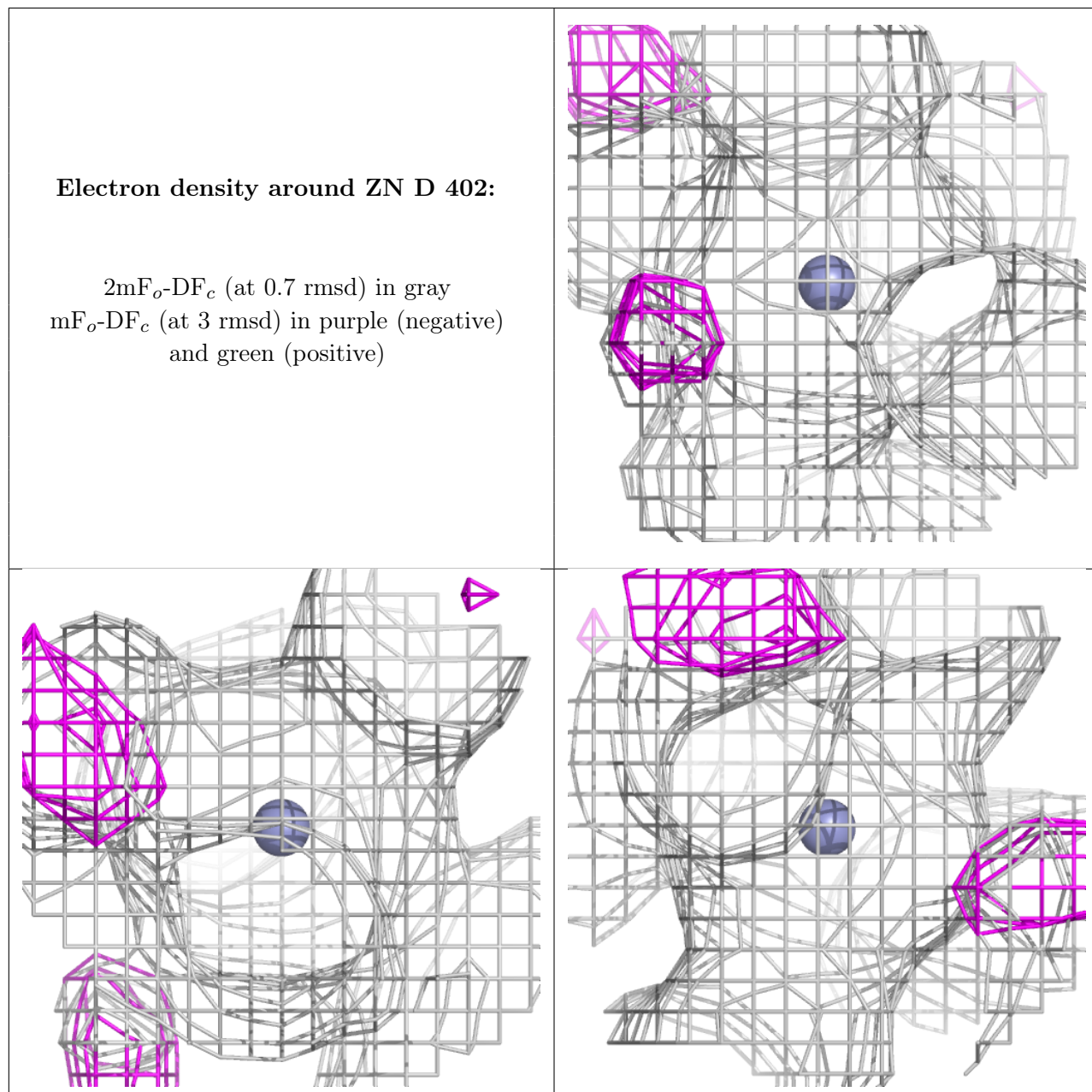
Electron density around ZN C 402:

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



Electron density around ZN D 402:

$2mF_o-DF_c$ (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.