



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 17, 2026 – 11:56 PM UTC

PDB ID : 7SRV / pdb_00007srv
Title : Metal dependent activation of Plasmodium falciparum M17 aminopeptidase (inactive form), spacegroup P22121
Authors : Webb, C.T.; McGowan, S.
Deposited on : 2021-11-08
Resolution : 2.03 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
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

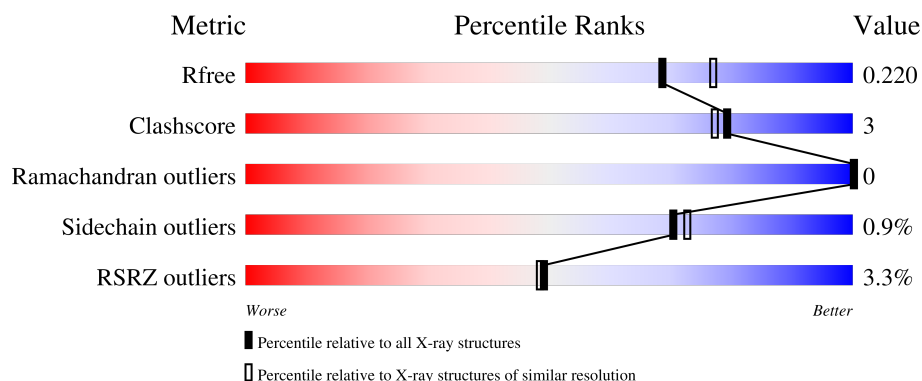
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.03 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	13299 (2.04-2.00)
Clashscore	190562	1022 (2.02-2.02)
Ramachandran outliers	187476	1014 (2.02-2.02)
Sidechain outliers	187428	1014 (2.02-2.02)
RSRZ outliers	180081	13314 (2.04-2.00)

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	527	0% 92% 6% .
1	B	527	6% 91% 6% .
1	C	527	3% 93% 6% .
1	D	527	2% 87% 9% .
1	E	527	4% 91% 6% .

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Mol	Chain	Length	Quality of chain
1	F	527	4% 88% 6% • 5%

2 Entry composition

There are 9 unique types of molecules in this entry. The entry contains 25540 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 M17 leucyl aminopeptidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	519	Total	C	N	O	S	0	3	0
			3964	2546	638	759	21			
1	B	518	Total	C	N	O	S	0	4	0
			3898	2506	630	743	19			
1	C	524	Total	C	N	O	S	0	3	0
			4021	2579	654	769	19			
1	D	506	Total	C	N	O	S	0	2	0
			3867	2488	623	736	20			
1	E	513	Total	C	N	O	S	0	2	0
			3944	2533	641	751	19			
1	F	502	Total	C	N	O	S	0	2	0
			3820	2456	610	736	18			

There are 54 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	152	GLN	ASN	conflict	UNP Q8IL11
A	515	GLN	ASN	conflict	UNP Q8IL11
A	546	GLN	ASN	conflict	UNP Q8IL11
A	606	HIS	-	expression tag	UNP Q8IL11
A	607	HIS	-	expression tag	UNP Q8IL11
A	608	HIS	-	expression tag	UNP Q8IL11
A	609	HIS	-	expression tag	UNP Q8IL11
A	610	HIS	-	expression tag	UNP Q8IL11
A	611	HIS	-	expression tag	UNP Q8IL11
B	152	GLN	ASN	conflict	UNP Q8IL11
B	515	GLN	ASN	conflict	UNP Q8IL11
B	546	GLN	ASN	conflict	UNP Q8IL11
B	606	HIS	-	expression tag	UNP Q8IL11
B	607	HIS	-	expression tag	UNP Q8IL11
B	608	HIS	-	expression tag	UNP Q8IL11
B	609	HIS	-	expression tag	UNP Q8IL11
B	610	HIS	-	expression tag	UNP Q8IL11

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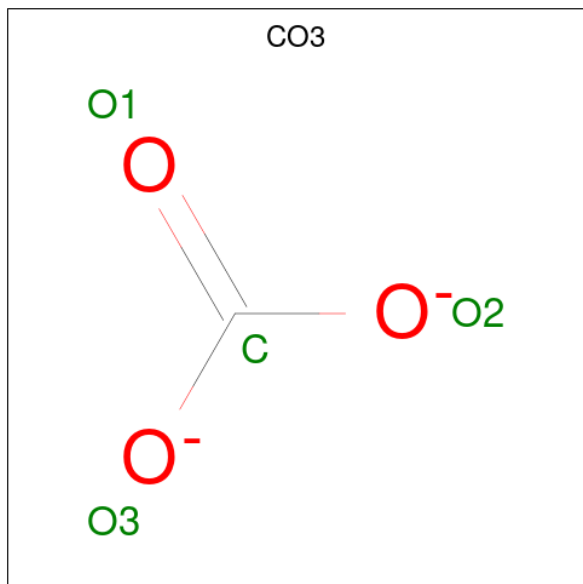
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Chain	Residue	Modelled	Actual	Comment	Reference
B	611	HIS	-	expression tag	UNP Q8IL11
C	152	GLN	ASN	conflict	UNP Q8IL11
C	515	GLN	ASN	conflict	UNP Q8IL11
C	546	GLN	ASN	conflict	UNP Q8IL11
C	606	HIS	-	expression tag	UNP Q8IL11
C	607	HIS	-	expression tag	UNP Q8IL11
C	608	HIS	-	expression tag	UNP Q8IL11
C	609	HIS	-	expression tag	UNP Q8IL11
C	610	HIS	-	expression tag	UNP Q8IL11
C	611	HIS	-	expression tag	UNP Q8IL11
D	152	GLN	ASN	conflict	UNP Q8IL11
D	515	GLN	ASN	conflict	UNP Q8IL11
D	546	GLN	ASN	conflict	UNP Q8IL11
D	606	HIS	-	expression tag	UNP Q8IL11
D	607	HIS	-	expression tag	UNP Q8IL11
D	608	HIS	-	expression tag	UNP Q8IL11
D	609	HIS	-	expression tag	UNP Q8IL11
D	610	HIS	-	expression tag	UNP Q8IL11
D	611	HIS	-	expression tag	UNP Q8IL11
E	152	GLN	ASN	conflict	UNP Q8IL11
E	515	GLN	ASN	conflict	UNP Q8IL11
E	546	GLN	ASN	conflict	UNP Q8IL11
E	606	HIS	-	expression tag	UNP Q8IL11
E	607	HIS	-	expression tag	UNP Q8IL11
E	608	HIS	-	expression tag	UNP Q8IL11
E	609	HIS	-	expression tag	UNP Q8IL11
E	610	HIS	-	expression tag	UNP Q8IL11
E	611	HIS	-	expression tag	UNP Q8IL11
F	152	GLN	ASN	conflict	UNP Q8IL11
F	515	GLN	ASN	conflict	UNP Q8IL11
F	546	GLN	ASN	conflict	UNP Q8IL11
F	606	HIS	-	expression tag	UNP Q8IL11
F	607	HIS	-	expression tag	UNP Q8IL11
F	608	HIS	-	expression tag	UNP Q8IL11
F	609	HIS	-	expression tag	UNP Q8IL11
F	610	HIS	-	expression tag	UNP Q8IL11
F	611	HIS	-	expression tag	UNP Q8IL11

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn) (labeled as "Ligand of Interest" by depositor).

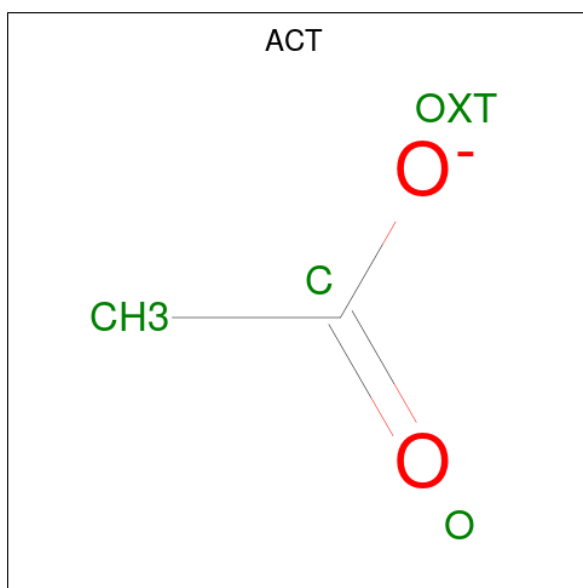
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total 2	Zn 2	0	0
2	B	2	Total 2	Zn 2	0	0
2	C	2	Total 2	Zn 2	0	0
2	D	2	Total 2	Zn 2	0	0
2	E	2	Total 2	Zn 2	0	0
2	F	2	Total 2	Zn 2	0	0

- Molecule 3 is CARBONATE ION (CCD ID: CO3) (formula: CO₃).



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

- Molecule 4 is ACETATE ION (CCD ID: ACT) (formula: $\text{C}_2\text{H}_3\text{O}_2$).



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

- Molecule 5 is CALCIUM ION (CCD ID: CA) (formula: Ca).

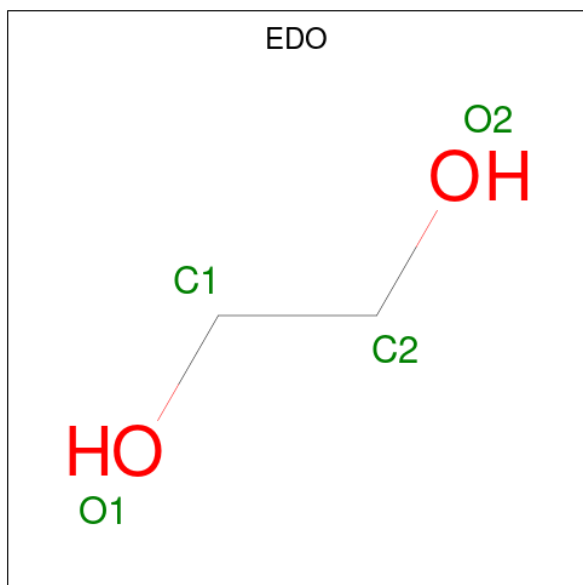
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total Ca 1 1	0	0
5	B	2	Total Ca 2 2	0	0
5	C	4	Total Ca 4 4	0	0

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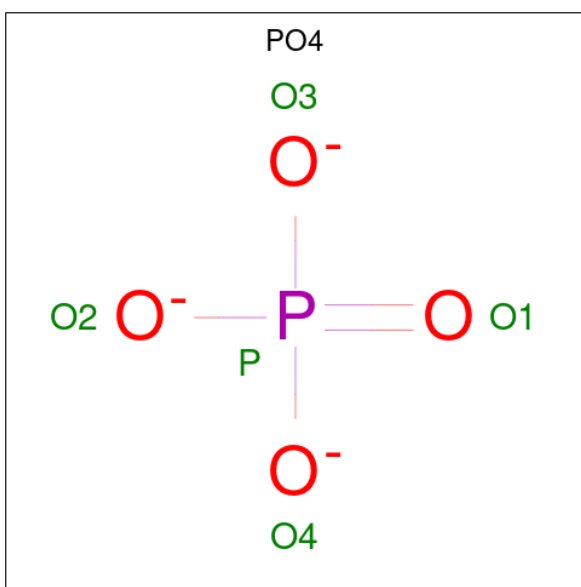
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	D	1	Total	Ca	0	0
			1	1		
5	E	3	Total	Ca	0	0
			3	3		
5	F	1	Total	Ca	0	0
			1	1		

- Molecule 6 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: $C_2H_6O_2$).



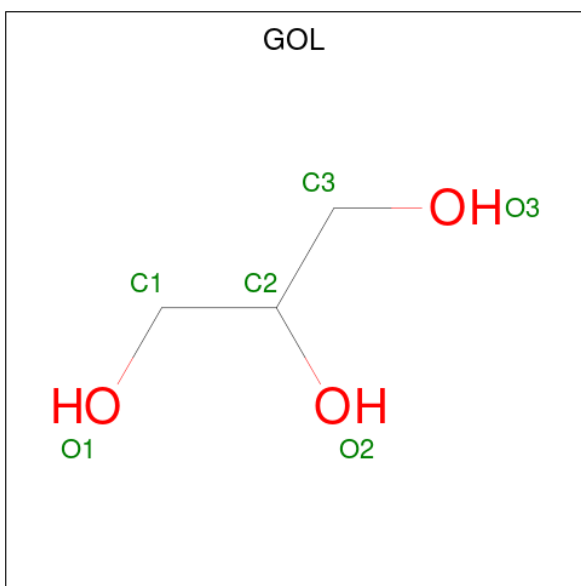
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			4	2	2		
6	C	1	Total	C	O	0	0
			4	2	2		
6	C	1	Total	C	O	0	0
			4	2	2		
6	C	1	Total	C	O	0	0
			4	2	2		
6	D	1	Total	C	O	0	0
			4	2	2		
6	F	1	Total	C	O	0	0
			4	2	2		

- Molecule 7 is PHOSPHATE ION (CCD ID: PO4) (formula: O_4P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	B	1	Total	O	P	0	0
			5	4	1		
7	F	1	Total	O	P	0	0
			5	4	1		

- Molecule 8 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	E	1	Total	C	O	0	0
			6	3	3		

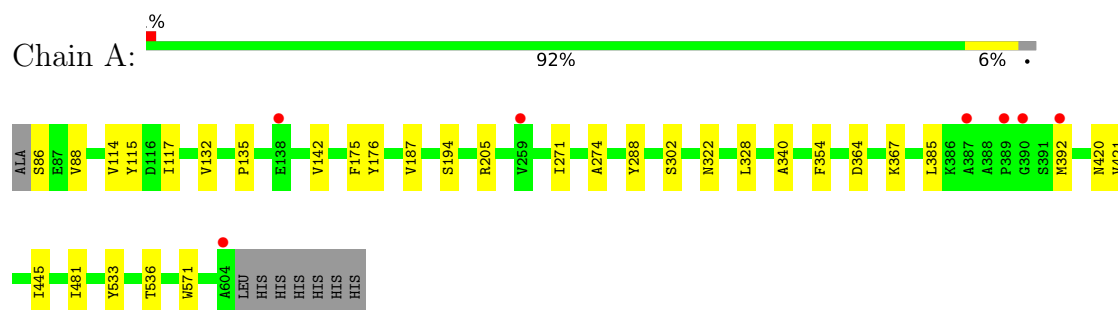
- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	343	Total 343	O 343	0	0
9	B	236	Total 236	O 236	0	0
9	C	370	Total 370	O 370	0	0
9	D	312	Total 312	O 312	0	0
9	E	387	Total 387	O 387	0	0
9	F	258	Total 258	O 258	0	0

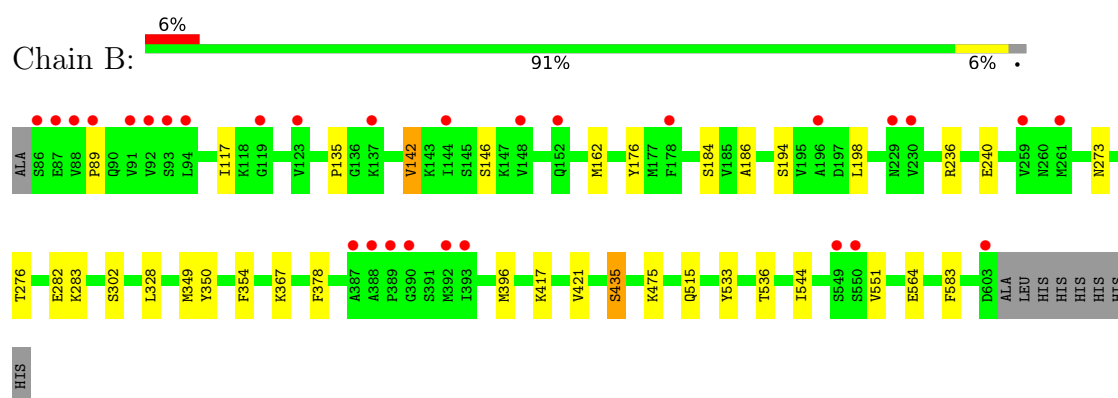
3 Residue-property plots [i](#)

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.

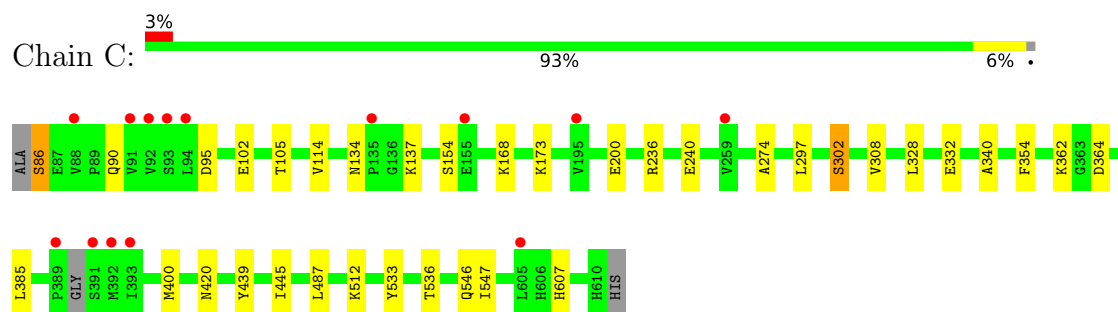
- Molecule 1: M17 leucyl aminopeptidase



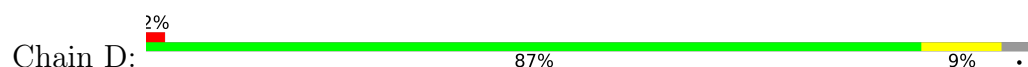
- Molecule 1: M17 leucyl aminopeptidase

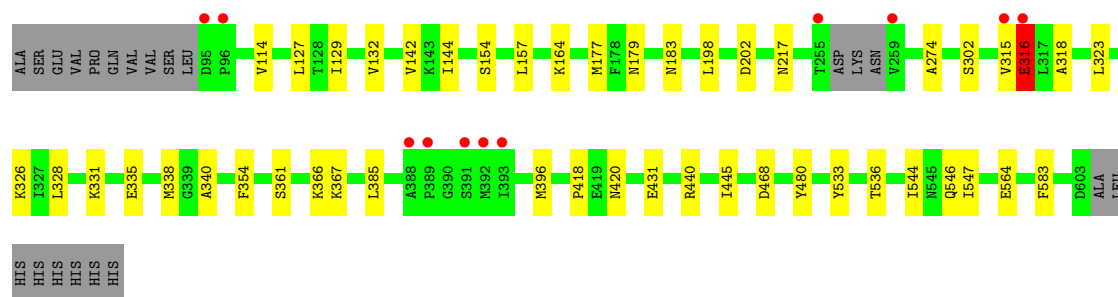


- Molecule 1: M17 leucyl aminopeptidase

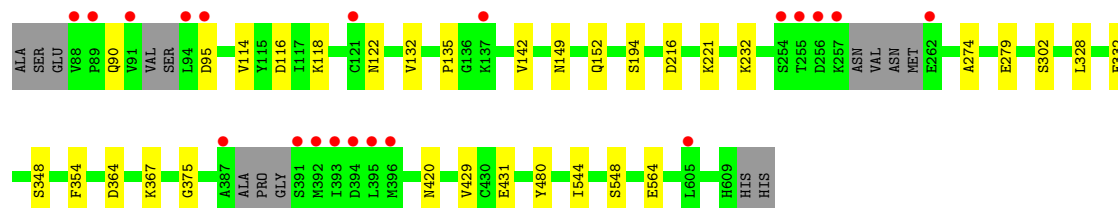
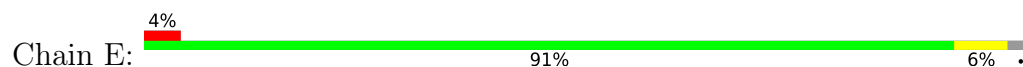


- Molecule 1: M17 leucyl aminopeptidase

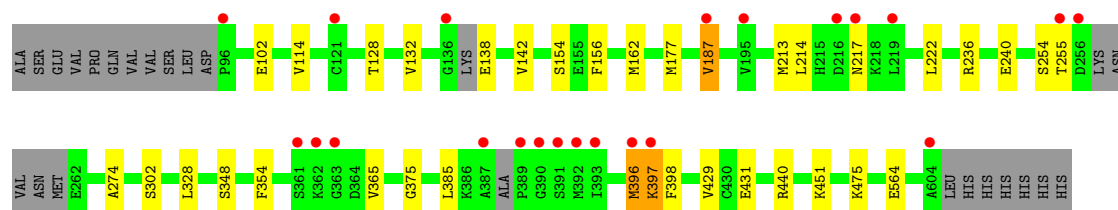
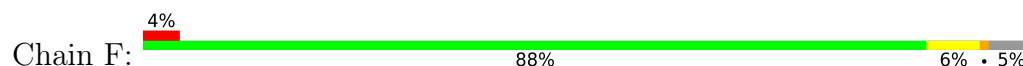




- Molecule 1: M17 leucyl aminopeptidase



- Molecule 1: M17 leucyl aminopeptidase



4 Data and refinement statistics

Property	Value	Source
Space group	P 2 21 21	Depositor
Cell constants a, b, c, α , β , γ	111.54Å 172.67Å 179.40Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.19 – 2.03 49.19 – 2.03	Depositor EDS
% Data completeness (in resolution range)	100.0 (49.19-2.03) 100.0 (49.19-2.03)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.12 (at 2.00Å)	Xtriage
Refinement program	PHENIX 1.19.2_4158	Depositor
R, R_{free}	0.189 , 0.219 0.190 , 0.220	Depositor DCC
R_{free} test set	11550 reflections (4.96%)	wwPDB-VP
Wilson B-factor (Å ²)	36.5	Xtriage
Anisotropy	0.068	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 35.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.000 for -h,l,k	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	25540	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.48% 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: PO4, ACT, CO3, GOL, CA, EDO, 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	0.13	1/4048 (0.0%)	0.28	0/5497
1	B	0.09	0/3982	0.29	0/5417
1	C	0.15	0/4109	0.40	3/5577 (0.1%)
1	D	0.12	0/3949	0.37	2/5358 (0.0%)
1	E	0.10	0/4028	0.32	0/5462
1	F	0.14	0/3900	0.45	7/5296 (0.1%)
All	All	0.12	1/24016 (0.0%)	0.36	12/32607 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	88	VAL	CA-C	-5.17	1.49	1.53

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	397	LYS	CB-CG-CD	13.70	142.81	111.30
1	C	154	SER	CA-C-N	-10.52	106.17	122.37
1	C	154	SER	C-N-CA	-10.52	106.17	122.37
1	D	316	GLU	CB-CA-C	-9.19	96.31	110.92
1	F	396	MET	CA-C-N	-8.94	108.46	122.67
1	F	396	MET	C-N-CA	-8.94	108.46	122.67
1	F	397	LYS	CG-CD-CE	-6.02	97.45	111.30
1	F	187	VAL	CG1-CB-CG2	5.92	123.82	110.80
1	D	316	GLU	CA-CB-CG	5.85	125.80	114.10
1	F	397	LYS	CD-CE-NZ	5.84	130.58	111.90
1	C	168	LYS	CB-CG-CD	5.70	124.41	111.30
1	F	187	VAL	CA-CB-CG1	5.25	119.32	110.40

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	3964	0	3861	16	0
1	B	3898	0	3737	24	0
1	C	4021	0	3910	19	0
1	D	3867	0	3784	27	0
1	E	3944	0	3846	22	0
1	F	3820	0	3682	24	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
2	C	2	0	0	0	0
2	D	2	0	0	0	0
2	E	2	0	0	0	0
2	F	2	0	0	0	0
3	A	4	0	0	0	0
3	B	4	0	0	0	0
3	C	4	0	0	0	0
3	D	4	0	0	0	0
3	E	4	0	0	0	0
3	F	4	0	0	0	0
4	A	4	0	3	0	0
4	B	4	0	3	0	0
4	C	8	0	6	0	0
4	D	8	0	6	0	0
4	E	4	0	3	1	0
4	F	4	0	3	0	0
5	A	1	0	0	0	0
5	B	2	0	0	0	0
5	C	4	0	0	0	0
5	D	1	0	0	0	0
5	E	3	0	0	0	0
5	F	1	0	0	0	0
6	A	4	0	6	0	0
6	C	12	0	18	1	0
6	D	4	0	6	0	0
6	F	4	0	6	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
7	B	5	0	0	1	0
7	F	5	0	0	0	0
8	E	6	0	8	0	0
9	A	343	0	0	4	0
9	B	236	0	0	5	0
9	C	370	0	0	2	0
9	D	312	0	0	2	0
9	E	387	0	0	5	0
9	F	258	0	0	6	0
All	All	25540	0	22888	127	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 3.

All (127) 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:440:ARG:NH1	1:E:431:GLU:OE2	2.19	0.70
1:D:366:LYS:HG3	1:D:420:ASN:HB3	1.75	0.69
1:E:90:GLN:NE2	1:E:95:ASP:O	2.26	0.68
1:A:205:ARG:NE	9:A:1106:HOH:O	2.28	0.67
1:D:144:ILE:HG13	1:D:157:LEU:HD22	1.78	0.66
1:F:132:VAL:HG21	1:F:142:VAL:HG13	1.77	0.64
1:F:128:THR:HB	1:F:187:VAL:HG22	1.79	0.64
1:E:328:LEU:HB2	1:E:354:PHE:HB3	1.78	0.64
1:B:367:LYS:HB3	1:B:421:VAL:HG23	1.79	0.64
1:C:236:ARG:NE	1:C:240:GLU:OE2	2.22	0.64
1:E:364:ASP:O	1:E:420:ASN:HA	1.99	0.62
1:D:326:LYS:NZ	9:D:1105:HOH:O	2.29	0.62
1:C:364:ASP:O	1:C:420:ASN:HA	1.99	0.62
1:F:236:ARG:NE	1:F:240:GLU:OE2	2.34	0.61
1:C:173:LYS:NZ	1:E:216:ASP:OD2	2.35	0.60
1:B:282:GLU:OE2	9:B:1101:HOH:O	2.15	0.60
1:E:364:ASP:OD1	9:E:1101:HOH:O	2.16	0.60
1:F:397:LYS:HB2	9:F:1199:HOH:O	2.02	0.60
1:B:176:TYR:HB3	1:F:177:MET:HE2	1.84	0.59
1:E:221:LYS:NZ	9:E:1109:HOH:O	2.36	0.59
1:B:515:GLN:NE2	9:B:1111:HOH:O	2.35	0.59
1:B:328:LEU:HB2	1:B:354:PHE:HB3	1.85	0.59
1:E:122:ASN:OD1	1:E:149:ASN:ND2	2.27	0.58
1:E:232:LYS:NZ	1:E:279:GLU:OE2	2.32	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:328:LEU:HB2	1:A:354:PHE:HB3	1.84	0.58
1:A:533:TYR:O	1:A:536:THR:HG22	2.03	0.58
1:D:396:MET:HE1	1:D:583:PHE:HZ	1.67	0.57
1:C:328:LEU:HD23	1:C:332:GLU:HG2	1.86	0.57
1:F:213:MET:O	1:F:217:ASN:HB2	2.05	0.56
1:D:154:SER:OG	9:D:1101:HOH:O	2.18	0.56
1:C:86:SER:HB3	1:C:308:VAL:HG13	1.86	0.56
1:F:138:GLU:N	9:F:1110:HOH:O	2.39	0.55
1:C:328:LEU:HB2	1:C:354:PHE:HB3	1.87	0.55
1:C:297:LEU:HB3	1:C:400:MET:HE1	1.89	0.54
1:A:322:ASN:ND2	9:A:1116:HOH:O	2.36	0.54
1:D:328:LEU:HB2	1:D:354:PHE:HB3	1.89	0.53
1:B:533:TYR:O	1:B:536[A]:THR:HG22	2.08	0.53
1:E:118:LYS:NZ	9:E:1104:HOH:O	2.40	0.53
1:D:179:ASN:HD21	1:D:183:ASN:HB2	1.74	0.53
1:E:132:VAL:HG21	1:E:142:VAL:HG13	1.91	0.52
1:F:328:LEU:HB2	1:F:354:PHE:HB3	1.91	0.52
1:B:273:ASN:O	1:B:276:THR:OG1	2.28	0.51
1:C:114:VAL:HG12	1:C:274:ALA:HB1	1.93	0.51
1:C:90:GLN:HB3	1:C:95:ASP:HB2	1.92	0.50
1:A:114:VAL:HG12	1:A:274:ALA:HB1	1.93	0.50
1:F:114:VAL:HG12	1:F:274:ALA:HB1	1.94	0.50
1:D:331:LYS:O	1:D:335:GLU:HG3	2.12	0.49
1:F:348:SER:OG	1:F:431:GLU:OE1	2.29	0.49
1:B:396:MET:HE1	1:B:583:PHE:HZ	1.75	0.49
1:E:332:GLU:OE1	9:E:1102:HOH:O	2.20	0.49
1:B:417:LYS:NZ	9:B:1116:HOH:O	2.45	0.48
1:D:367:LYS:HD3	1:D:480:TYR:CE2	2.48	0.48
1:F:397:LYS:HD3	1:F:398:PHE:CE1	2.49	0.48
1:F:236:ARG:NH1	9:F:1114:HOH:O	2.44	0.48
1:F:396:MET:HA	9:F:1171:HOH:O	2.13	0.48
1:B:176:TYR:CE1	1:B:184:SER:HB2	2.49	0.48
1:B:435:SER:HB2	7:B:1004:PO4:O4	2.14	0.47
1:D:361:SER:OG	1:D:418:PRO:O	2.28	0.47
1:C:607:HIS:CD2	6:C:1010:EDO:H12	2.49	0.47
1:B:135:PRO:HA	1:B:194:SER:O	2.15	0.47
1:D:544:ILE:HD12	1:D:564:GLU:HG3	1.97	0.46
1:E:348:SER:OG	1:E:431:GLU:OE1	2.31	0.46
1:A:364:ASP:O	1:A:420:ASN:HA	2.16	0.46
1:E:367:LYS:HD2	1:E:480:TYR:CE2	2.51	0.46
1:B:302:SER:OG	1:B:378:PHE:HB2	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:132:VAL:HG21	1:A:142:VAL:HG13	1.98	0.46
1:D:132:VAL:HG21	1:D:142:VAL:HG13	1.97	0.46
1:B:186:ALA:O	9:B:1102:HOH:O	2.21	0.45
1:E:114:VAL:HG12	1:E:274:ALA:HB1	1.98	0.45
1:F:156:PHE:O	1:F:162:MET:HE3	2.17	0.45
1:F:451:LYS:NZ	1:F:564:GLU:O	2.46	0.45
1:D:338:MET:HG2	1:D:468:ASP:HB3	1.99	0.45
1:C:134:ASN:OD1	1:C:137:LYS:HB2	2.16	0.45
1:D:315:VAL:O	1:D:316:GLU:C	2.60	0.45
1:E:152:GLN:OE1	9:E:1103:HOH:O	2.21	0.44
1:F:102:GLU:HG3	9:F:1317:HOH:O	2.17	0.44
1:B:117:ILE:HD11	1:B:146:SER:OG	2.18	0.44
1:C:533:TYR:O	1:C:536:THR:HG22	2.18	0.44
1:B:236:ARG:NE	1:B:240:GLU:OE2	2.43	0.44
1:D:318:ALA:HA	1:D:323:LEU:HB2	1.99	0.44
1:E:367:LYS:HB2	1:E:367:LYS:HE2	1.61	0.44
1:E:544:ILE:HD12	1:E:564:GLU:HG3	2.00	0.44
1:A:86:SER:N	9:A:1135:HOH:O	2.51	0.43
1:B:198:LEU:HD23	1:B:198:LEU:HA	1.85	0.43
1:A:135:PRO:HA	1:A:194:SER:O	2.18	0.43
1:B:89:PRO:HD2	1:B:350:TYR:HD1	1.82	0.43
1:D:367:LYS:HD3	1:D:480:TYR:HE2	1.84	0.43
1:F:475:LYS:NZ	9:F:1128:HOH:O	2.51	0.43
1:B:349:MET:HE3	1:B:349:MET:HB2	1.90	0.43
1:D:431:GLU:OE2	1:F:440:ARG:NH1	2.48	0.43
1:A:115:TYR:OH	9:A:1101:HOH:O	2.21	0.43
1:D:546:GLN:HG2	1:D:547:ILE:HG23	2.00	0.43
1:F:214:LEU:HD21	1:F:222:LEU:HD22	2.00	0.43
1:B:536[B]:THR:HG21	1:B:551:VAL:HG23	2.00	0.42
1:D:198:LEU:HD22	1:D:202:ASP:HB3	2.01	0.42
1:C:512:LYS:NZ	9:C:1127:HOH:O	2.52	0.42
1:F:156:PHE:CD1	1:F:177:MET:HE1	2.54	0.42
1:E:375:GLY:O	1:E:429:VAL:HA	2.20	0.42
1:F:156:PHE:CE2	1:F:187:VAL:CG1	3.03	0.42
1:B:544:ILE:HD12	1:B:564:GLU:HG3	2.02	0.42
1:E:548:SER:HB3	4:E:1003:ACT:H2	2.01	0.42
1:E:135:PRO:HA	1:E:194:SER:O	2.20	0.42
1:F:375:GLY:O	1:F:429:VAL:HA	2.19	0.41
1:A:176:TYR:HB3	1:D:177:MET:HE2	2.02	0.41
1:C:340:ALA:HA	1:C:445:ILE:HD12	2.02	0.41
1:B:475:LYS:NZ	9:B:1120:HOH:O	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:142:VAL:HG22	1:B:162:MET:O	2.20	0.41
1:D:127:LEU:HD11	1:D:129:ILE:HD11	2.01	0.41
1:A:117:ILE:HG13	1:A:271:ILE:C	2.45	0.41
1:A:481:ILE:O	1:A:571:TRP:HA	2.21	0.41
1:C:546:GLN:HG2	1:C:547:ILE:HG23	2.02	0.41
1:A:340:ALA:HA	1:A:445:ILE:HD12	2.03	0.41
1:C:487:LEU:HD12	1:C:487:LEU:HA	1.91	0.41
1:D:114:VAL:HG12	1:D:274:ALA:HB1	2.02	0.41
1:D:533:TYR:O	1:D:536:THR:HG22	2.21	0.41
1:B:236:ARG:HD2	1:B:283:LYS:HD3	2.02	0.41
1:C:362:LYS:O	9:C:1101:HOH:O	2.22	0.41
1:D:340:ALA:HA	1:D:445:ILE:HD12	2.03	0.41
1:F:177:MET:HE2	1:F:177:MET:HB3	1.91	0.41
1:F:254:SER:OG	1:F:255:THR:N	2.52	0.41
1:A:175:PHE:N	1:A:187:VAL:O	2.46	0.40
1:D:396:MET:HE1	1:D:583:PHE:CZ	2.52	0.40
1:E:116:ASP:OD1	1:E:118:LYS:HE2	2.20	0.40
1:C:102:GLU:HG2	1:C:105:THR:HG22	2.03	0.40
1:D:164:LYS:HE3	1:D:164:LYS:HB2	1.77	0.40
1:A:392:MET:HE3	1:A:392:MET:HB2	1.76	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	519/527 (98%)	506 (98%)	13 (2%)	0	100	100
1	B	518/527 (98%)	504 (97%)	14 (3%)	0	100	100
1	C	522/527 (99%)	509 (98%)	13 (2%)	0	100	100
1	D	504/527 (96%)	492 (98%)	12 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	E	507/527 (96%)	493 (97%)	14 (3%)	0	100	100
1	F	496/527 (94%)	484 (98%)	12 (2%)	0	100	100
All	All	3066/3162 (97%)	2988 (98%)	78 (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	420/454 (92%)	415 (99%)	5 (1%)	63	65
1	B	402/454 (88%)	400 (100%)	2 (0%)	81	81
1	C	428/454 (94%)	423 (99%)	5 (1%)	63	65
1	D	409/454 (90%)	405 (99%)	4 (1%)	68	71
1	E	419/454 (92%)	418 (100%)	1 (0%)	87	88
1	F	399/454 (88%)	394 (99%)	5 (1%)	61	63
All	All	2477/2724 (91%)	2455 (99%)	22 (1%)	70	73

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	288	TYR
1	A	302	SER
1	A	367	LYS
1	A	385	LEU
1	A	421	VAL
1	B	142	VAL
1	B	435	SER
1	C	86	SER
1	C	200	GLU
1	C	302	SER
1	C	385	LEU
1	C	439	TYR

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Mol	Chain	Res	Type
1	D	217	ASN
1	D	302	SER
1	D	316	GLU
1	D	385	LEU
1	E	302	SER
1	F	154[A]	SER
1	F	154[B]	SER
1	F	302	SER
1	F	365	VAL
1	F	385	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	215	HIS
1	A	420	ASN
1	A	521	ASN
1	B	113	GLN
1	B	272	ASN
1	B	515	GLN
1	C	602	ASN
1	D	104	ASN
1	D	139	ASN
1	D	161	ASN
1	D	183	ASN
1	D	266	HIS
1	D	521	ASN
1	D	568	ASN
1	E	161	ASN
1	E	174	HIS
1	E	266	HIS
1	E	437	ASN
1	E	606	HIS
1	F	139	ASN
1	F	266	HIS

5.3.3 RNA ⓘ

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](#)

Of 47 ligands modelled in this entry, 24 are monoatomic - leaving 23 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	$\# Z > 2$
6	EDO	D	1005	-	3,3,3	0.42	0	2,2,2	0.41	0
4	ACT	B	1003	-	3,3,3	1.41	1 (33%)	3,3,3	1.49	0
4	ACT	A	1003	-	3,3,3	1.36	0	3,3,3	1.36	0
4	ACT	E	1003	-	3,3,3	1.36	0	3,3,3	1.36	0
6	EDO	A	1005	-	3,3,3	0.43	0	2,2,2	0.39	0
4	ACT	D	1003	-	3,3,3	1.36	0	3,3,3	1.36	0
3	CO3	E	1002	-	3,3,3	0.85	0	2,3,3	0.17	0
6	EDO	C	1008	-	3,3,3	0.43	0	2,2,2	0.38	0
3	CO3	F	1002	-	3,3,3	0.85	0	2,3,3	0.18	0
4	ACT	F	1006	-	3,3,3	1.39	0	3,3,3	1.34	0
4	ACT	D	1006	-	3,3,3	1.40	1 (33%)	3,3,3	1.50	0
3	CO3	C	1002	-	3,3,3	0.79	0	2,3,3	0.07	0
6	EDO	C	1009	-	3,3,3	0.43	0	2,2,2	0.39	0
4	ACT	C	1003	-	3,3,3	1.36	0	3,3,3	1.51	0
3	CO3	A	1002	-	3,3,3	0.80	0	2,3,3	0.10	0
3	CO3	B	1002	-	3,3,3	0.83	0	2,3,3	0.15	0
4	ACT	C	1011	-	3,3,3	1.34	0	3,3,3	1.36	0
7	PO4	B	1004	-	4,4,4	0.99	0	6,6,6	0.51	0
3	CO3	D	1002	-	3,3,3	0.85	0	2,3,3	0.18	0
7	PO4	F	1003	-	4,4,4	0.97	0	6,6,6	0.44	0
8	GOL	E	1007	-	5,5,5	0.96	0	5,5,5	1.04	0
6	EDO	C	1010	-	3,3,3	0.41	0	2,2,2	0.44	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	EDO	F	1005	-	3,3,3	0.43	0	2,2,2	0.36	0

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
6	EDO	A	1005	-	-	0/1/1/1	-
6	EDO	C	1008	-	-	0/1/1/1	-
6	EDO	D	1005	-	-	0/1/1/1	-
8	GOL	E	1007	-	-	1/4/4/4	-
6	EDO	C	1010	-	-	0/1/1/1	-
6	EDO	C	1009	-	-	1/1/1/1	-
6	EDO	F	1005	-	-	0/1/1/1	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	D	1006	ACT	CH3-C	2.01	1.57	1.49
4	B	1003	ACT	CH3-C	2.00	1.57	1.49

There are no bond angle outliers.

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
8	E	1007	GOL	O1-C1-C2-C3
6	C	1009	EDO	O1-C1-C2-O2

There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	E	1003	ACT	1	0
7	B	1004	PO4	1	0
6	C	1010	EDO	1	0

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	519/527 (98%)	-0.12	7 (1%)	75 75	17, 35, 56, 80	7 (1%)
1	B	518/527 (98%)	0.27	29 (5%)	30 29	24, 40, 69, 83	11 (2%)
1	C	524/527 (99%)	-0.12	14 (2%)	56 56	22, 33, 58, 80	11 (2%)
1	D	506/527 (96%)	-0.07	11 (2%)	62 62	18, 36, 54, 87	13 (2%)
1	E	513/527 (97%)	-0.12	20 (3%)	43 43	22, 34, 54, 70	14 (2%)
1	F	502/527 (95%)	0.19	22 (4%)	39 38	25, 40, 68, 88	11 (2%)
All	All	3082/3162 (97%)	0.01	103 (3%)	49 48	17, 36, 63, 88	67 (2%)

All (103) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	96	PRO	4.8
1	D	95	ASP	4.4
1	A	604	ALA	4.2
1	C	393	ILE	4.1
1	F	255	THR	4.1
1	F	389	PRO	4.1
1	B	137	LYS	4.0
1	E	395	LEU	3.9
1	E	254	SER	3.9
1	F	362	LYS	3.9
1	F	393	ILE	3.8
1	B	88	VAL	3.7
1	D	316	GLU	3.7
1	C	392	MET	3.5
1	E	396	MET	3.5
1	E	262	GLU	3.5
1	C	391	SER	3.4
1	B	389	PRO	3.4
1	F	391	SER	3.4

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Mol	Chain	Res	Type	RSRZ
1	A	390	GLY	3.3
1	F	216	ASP	3.3
1	C	92	VAL	3.3
1	B	603	ASP	3.3
1	E	255	THR	3.2
1	F	387	ALA	3.2
1	F	256	ASP	3.2
1	E	88	VAL	3.1
1	F	363	GLY	3.1
1	E	89	PRO	3.1
1	B	123	VAL	3.0
1	D	392	MET	3.0
1	E	394	ASP	3.0
1	E	393	ILE	3.0
1	D	255	THR	3.0
1	B	89	PRO	3.0
1	D	391	SER	2.9
1	B	259	VAL	2.9
1	B	92	VAL	2.8
1	C	88	VAL	2.8
1	E	257	LYS	2.8
1	C	93	SER	2.8
1	A	392	MET	2.8
1	B	94	LEU	2.8
1	A	389	PRO	2.8
1	D	393	ILE	2.8
1	B	388	ALA	2.8
1	F	604	ALA	2.8
1	E	91	VAL	2.8
1	C	259	VAL	2.8
1	D	388	ALA	2.7
1	F	397	LYS	2.7
1	B	178	PHE	2.7
1	B	93	SER	2.7
1	F	217	ASN	2.6
1	F	219	LEU	2.6
1	F	392	MET	2.6
1	E	137	LYS	2.6
1	B	148	VAL	2.6
1	F	187	VAL	2.6
1	B	87	GLU	2.6
1	B	390	GLY	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	393	ILE	2.5
1	D	259	VAL	2.5
1	A	387	ALA	2.5
1	D	389	PRO	2.5
1	C	155	GLU	2.5
1	F	136	GLY	2.5
1	E	94	LEU	2.5
1	E	121	CYS	2.5
1	E	391	SER	2.5
1	F	96	PRO	2.5
1	E	392	MET	2.5
1	B	392	MET	2.4
1	B	550	SER	2.4
1	C	195	VAL	2.4
1	B	86	SER	2.4
1	B	144	ILE	2.4
1	F	121	CYS	2.4
1	E	387	ALA	2.3
1	C	91	VAL	2.3
1	C	135	PRO	2.3
1	B	196	ALA	2.3
1	E	95	ASP	2.3
1	B	261	MET	2.3
1	C	389	PRO	2.3
1	B	91	VAL	2.3
1	B	230	VAL	2.3
1	E	256	ASP	2.2
1	A	138	GLU	2.2
1	B	229	ASN	2.2
1	A	259	VAL	2.2
1	F	195	VAL	2.2
1	B	387	ALA	2.2
1	B	549	SER	2.2
1	F	361	SER	2.2
1	C	605	LEU	2.2
1	E	605	LEU	2.1
1	D	315	VAL	2.0
1	F	396	MET	2.0
1	B	152	GLN	2.0
1	C	94	LEU	2.0
1	B	119	GLY	2.0
1	F	390	GLY	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
6	EDO	C	1009	4/4	0.63	0.22	40,54,57,59	0
6	EDO	A	1005	4/4	0.73	0.12	51,54,58,60	0
5	CA	B	1006	1/1	0.73	0.14	83,83,83,83	0
8	GOL	E	1007	6/6	0.76	0.17	43,49,63,66	0
4	ACT	D	1006	4/4	0.79	0.11	43,50,53,54	0
5	CA	C	1006	1/1	0.83	0.19	60,60,60,60	0
6	EDO	D	1005	4/4	0.84	0.13	47,48,52,53	0
5	CA	C	1005	1/1	0.84	0.08	84,84,84,84	0
5	CA	E	1006	1/1	0.85	0.10	79,79,79,79	0
6	EDO	C	1008	4/4	0.87	0.16	48,51,56,61	0
6	EDO	F	1005	4/4	0.88	0.10	46,50,53,54	0
6	EDO	C	1010	4/4	0.88	0.12	41,43,51,53	0
4	ACT	C	1011	4/4	0.89	0.12	34,45,46,49	0
5	CA	C	1007	1/1	0.89	0.07	75,75,75,75	0
4	ACT	E	1003	4/4	0.91	0.10	31,40,42,45	0
4	ACT	D	1003	4/4	0.92	0.12	35,37,39,43	0
3	CO3	F	1002	4/4	0.93	0.10	28,32,37,47	0
4	ACT	B	1003	4/4	0.93	0.08	30,34,34,38	0
3	CO3	D	1002	4/4	0.93	0.09	29,31,36,41	0
3	CO3	A	1002	4/4	0.94	0.09	26,29,33,45	0
3	CO3	C	1002	4/4	0.95	0.11	25,28,34,42	0
4	ACT	C	1003	4/4	0.95	0.07	31,32,33,37	0
7	PO4	F	1003	5/5	0.95	0.13	48,51,52,57	0
4	ACT	F	1006	4/4	0.95	0.08	35,35,39,43	0
3	CO3	E	1002	4/4	0.96	0.08	28,31,40,44	0
4	ACT	A	1003	4/4	0.96	0.07	32,36,38,42	0
2	ZN	D	1007	1/1	0.97	0.04	37,37,37,37	1

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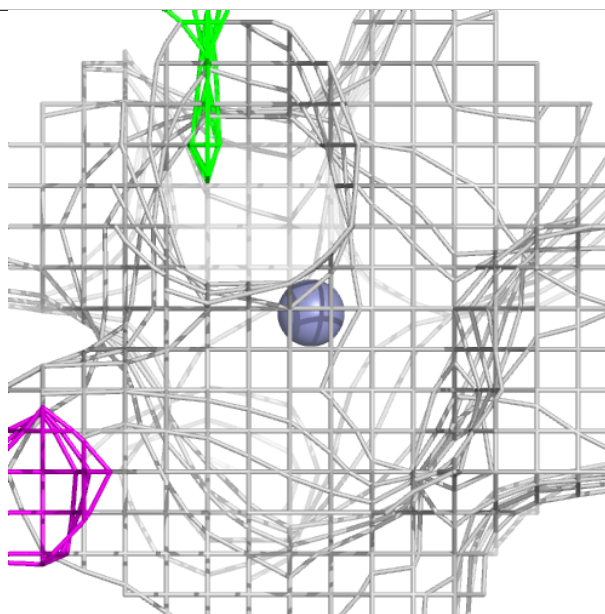
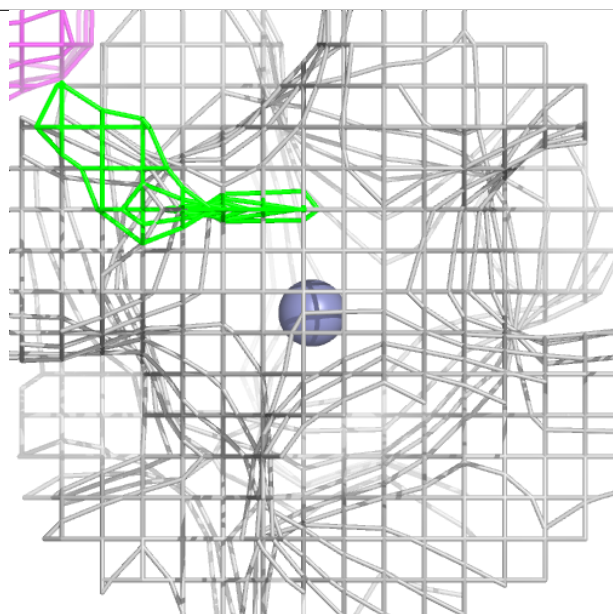
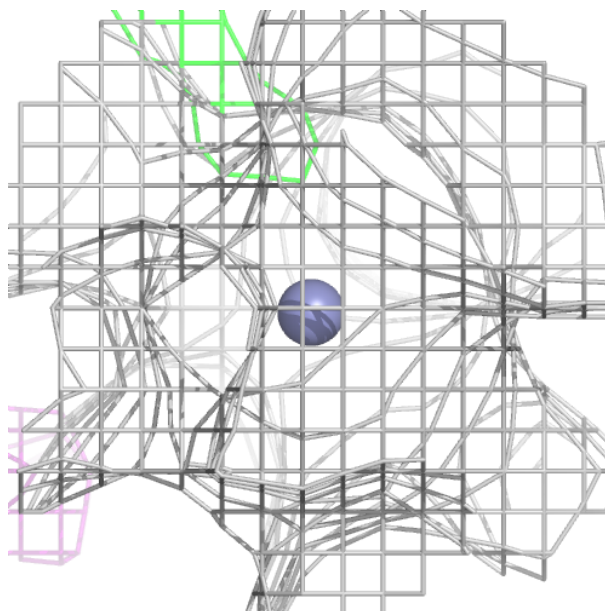
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
7	PO4	B	1004	5/5	0.97	0.08	34,39,44,50	0
3	CO3	B	1002	4/4	0.97	0.08	29,32,33,38	0
2	ZN	F	1007	1/1	0.97	0.04	33,33,33,33	1
2	ZN	C	1012	1/1	0.98	0.03	31,31,31,31	1
5	CA	E	1005	1/1	0.98	0.04	48,48,48,48	0
2	ZN	E	1008	1/1	0.98	0.04	34,34,34,34	1
5	CA	B	1005	1/1	0.99	0.11	36,36,36,36	0
2	ZN	E	1001	1/1	0.99	0.02	32,32,32,32	1
2	ZN	B	1007	1/1	0.99	0.02	39,39,39,39	1
2	ZN	F	1001	1/1	0.99	0.02	30,30,30,30	1
2	ZN	A	1006	1/1	0.99	0.02	34,34,34,34	1
5	CA	D	1004	1/1	0.99	0.08	34,34,34,34	0
2	ZN	D	1001	1/1	0.99	0.02	33,33,33,33	1
2	ZN	B	1001	1/1	0.99	0.01	34,34,34,34	1
5	CA	F	1004	1/1	0.99	0.04	36,36,36,36	0
2	ZN	A	1001	1/1	1.00	0.01	29,29,29,29	1
5	CA	C	1004	1/1	1.00	0.09	33,33,33,33	0
5	CA	E	1004	1/1	1.00	0.12	33,33,33,33	0
5	CA	A	1004	1/1	1.00	0.08	33,33,33,33	0
2	ZN	C	1001	1/1	1.00	0.07	44,44,44,44	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.

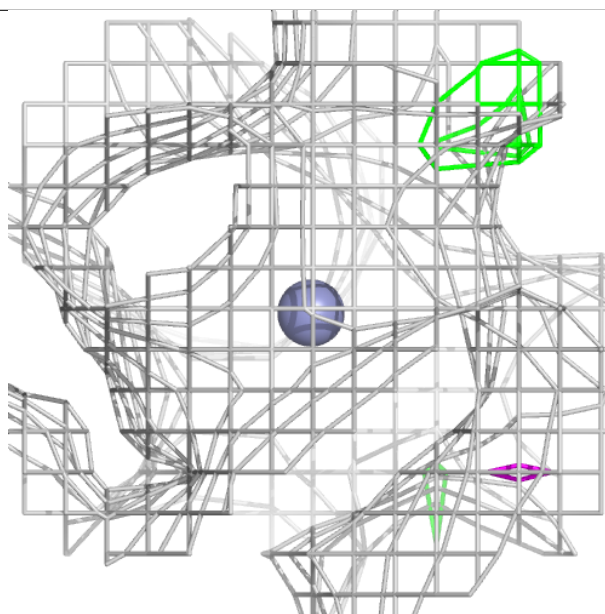
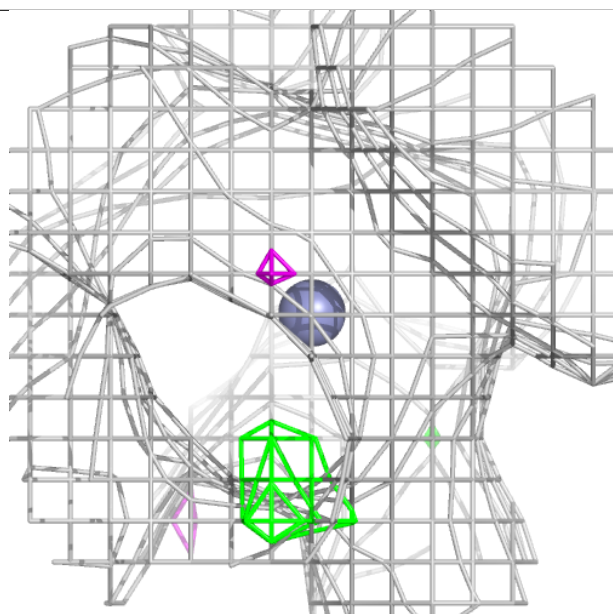
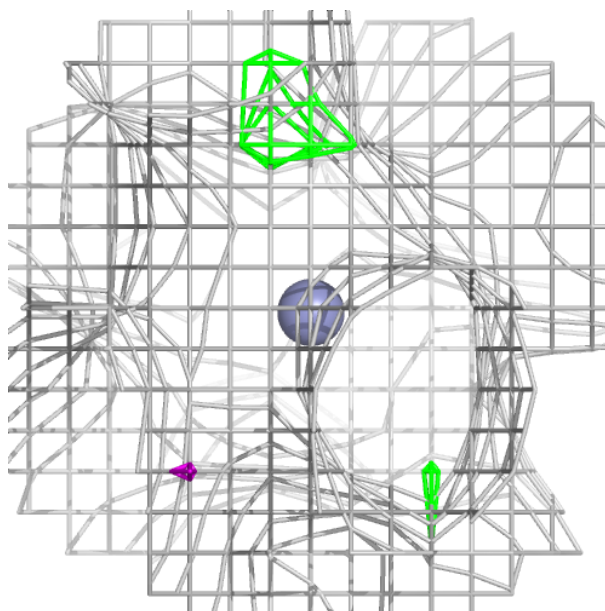
Electron density around ZN D 1007:

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



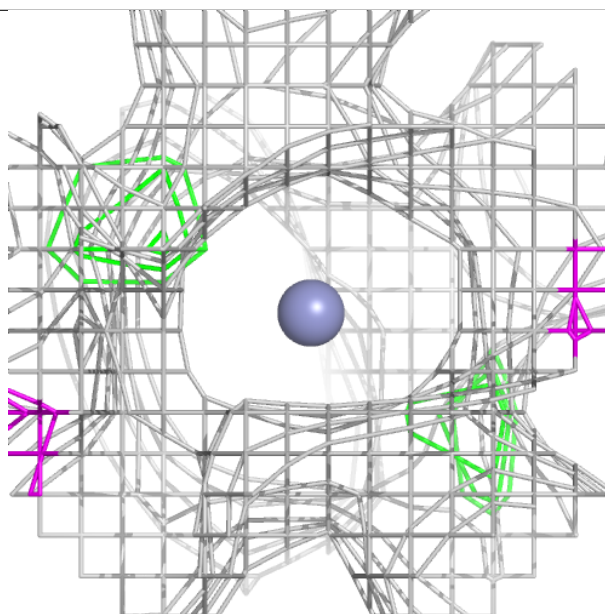
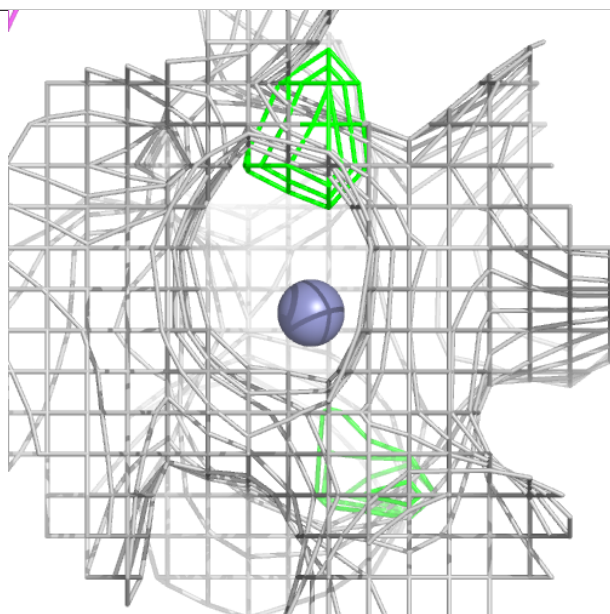
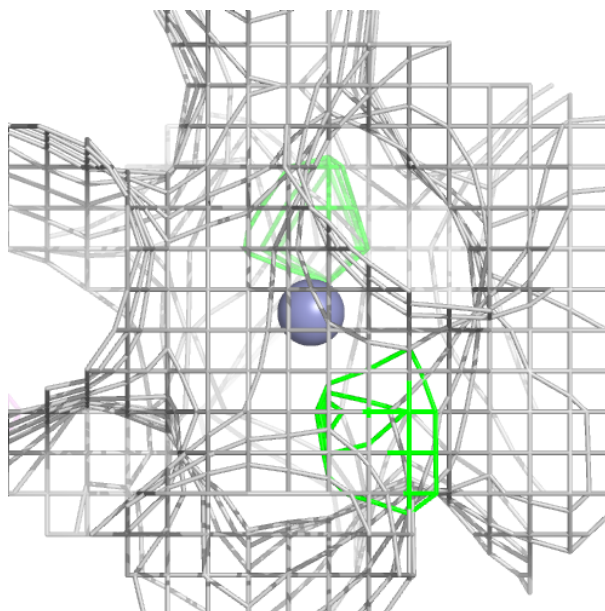
Electron density around ZN F 1007:

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



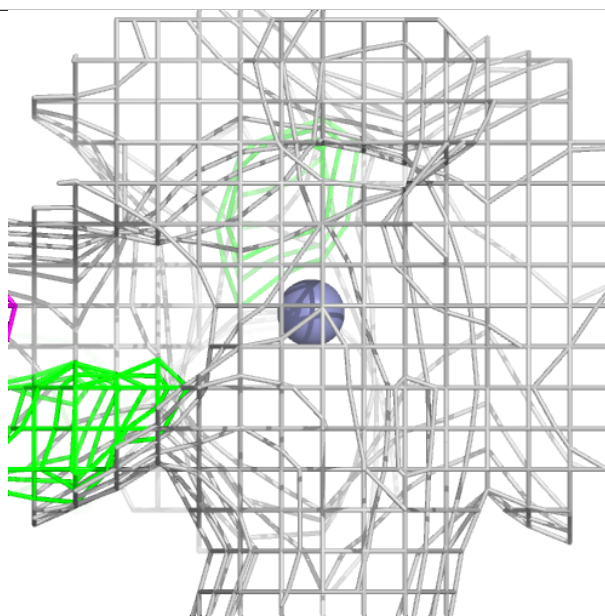
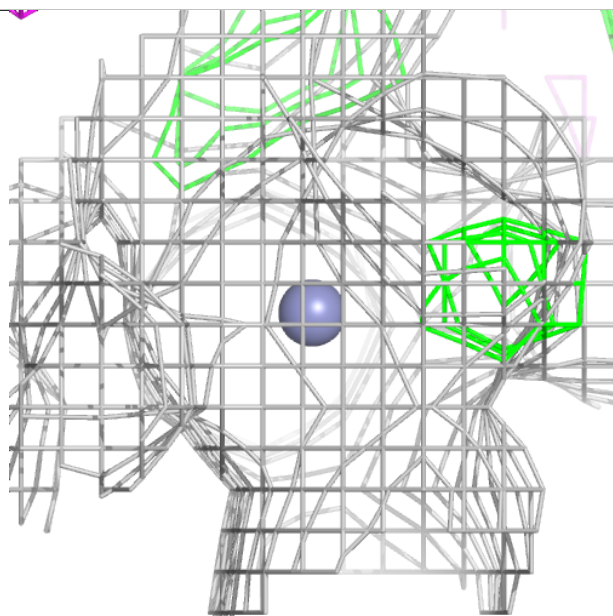
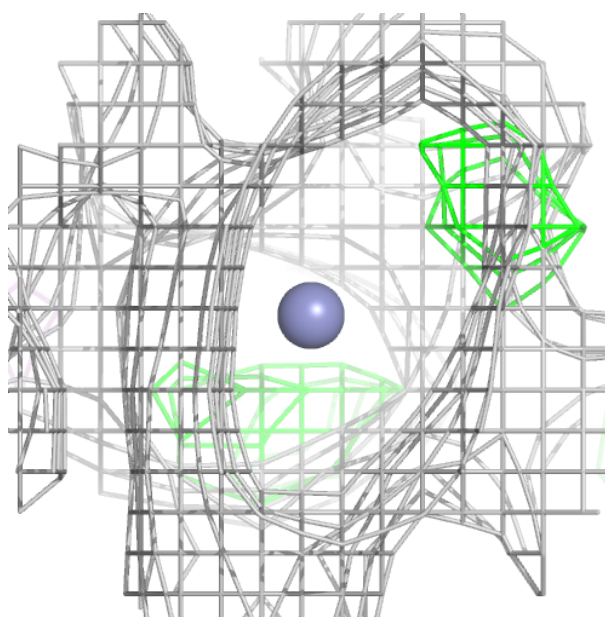
Electron density around ZN C 1012:

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and green (positive)



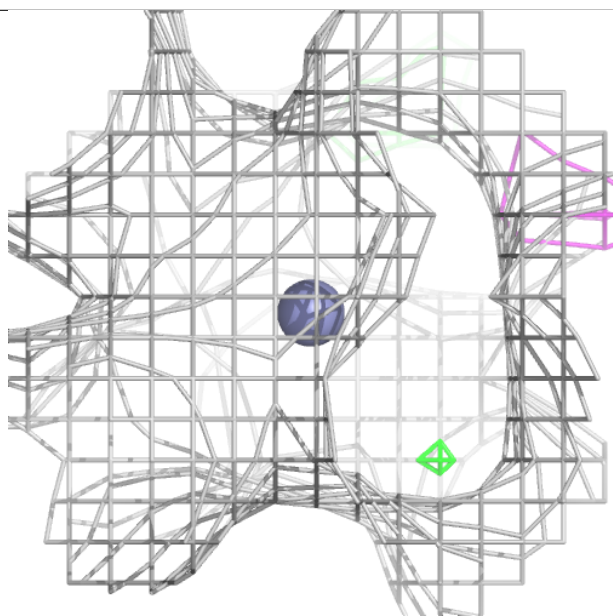
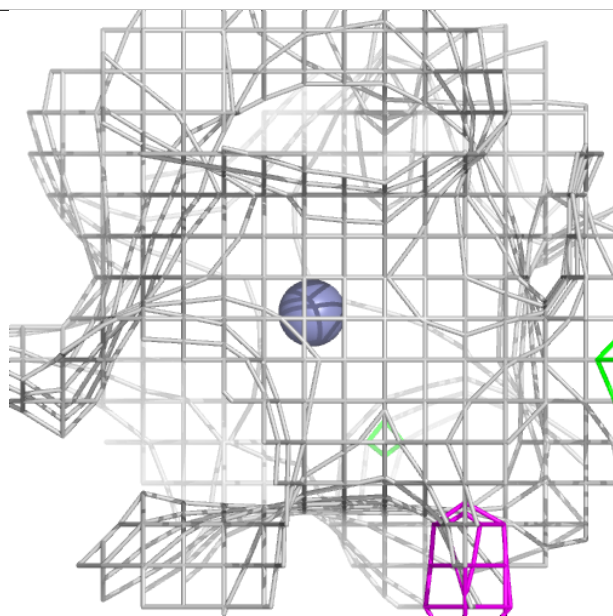
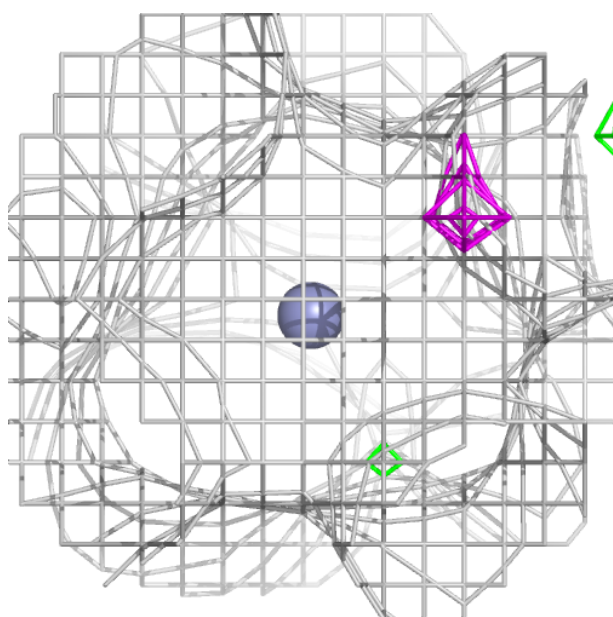
Electron density around ZN E 1008:

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and green (positive)



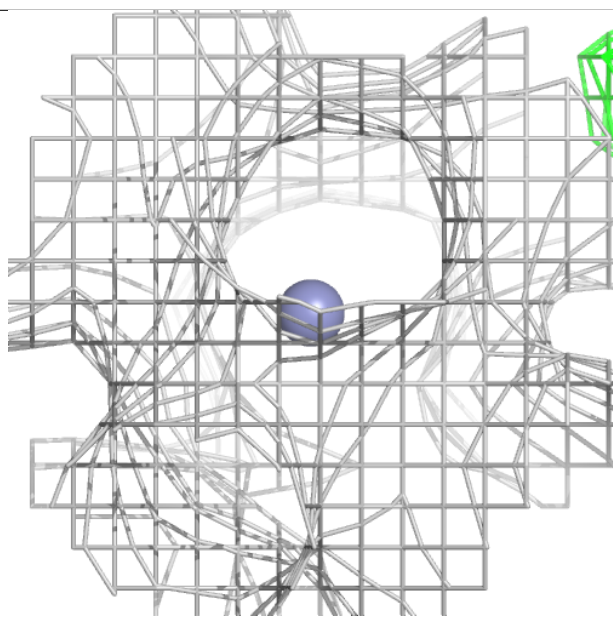
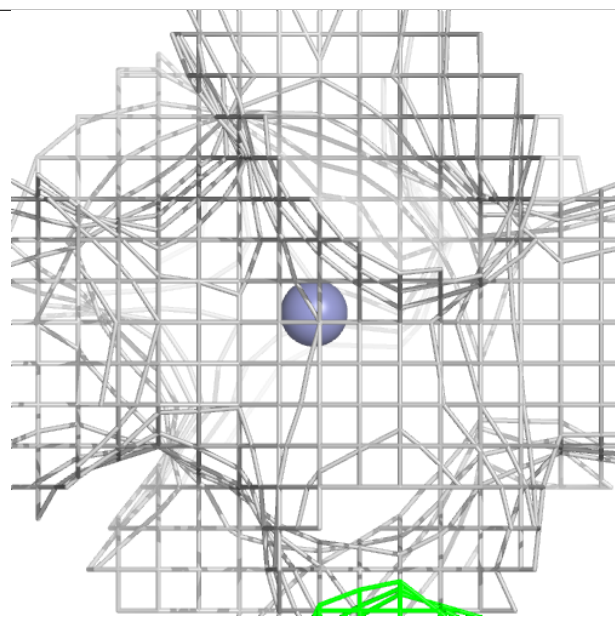
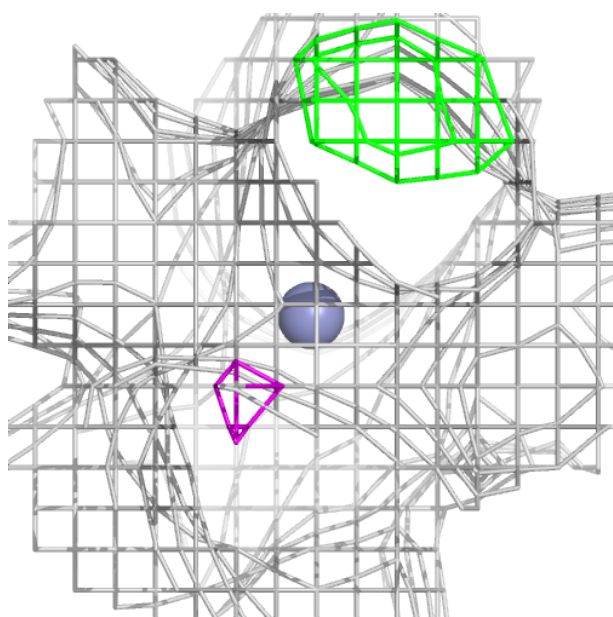
Electron density around ZN E 1001:

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and green (positive)



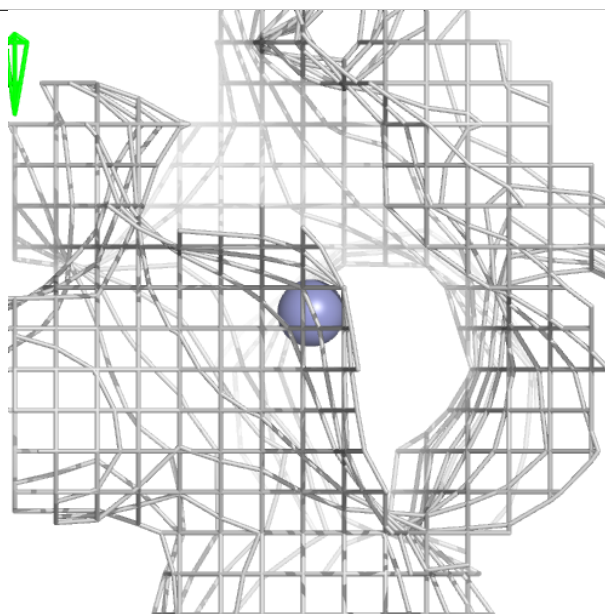
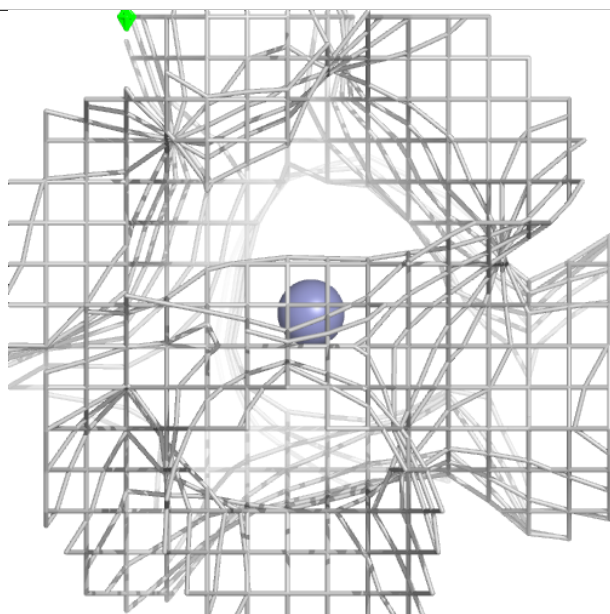
Electron density around ZN B 1007:

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and green (positive)



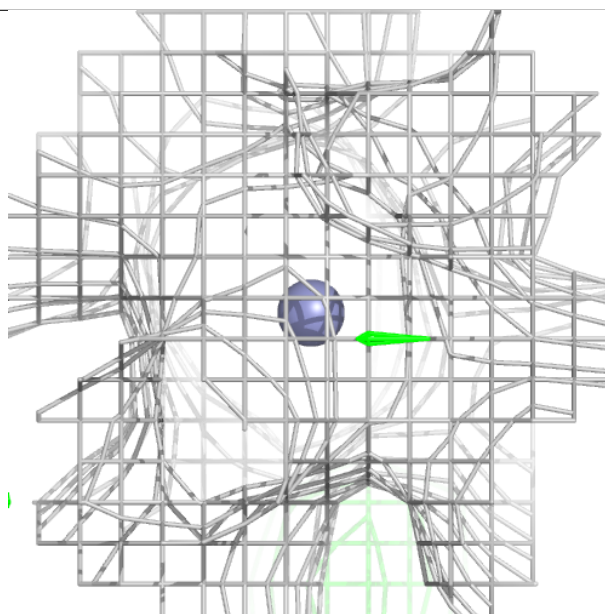
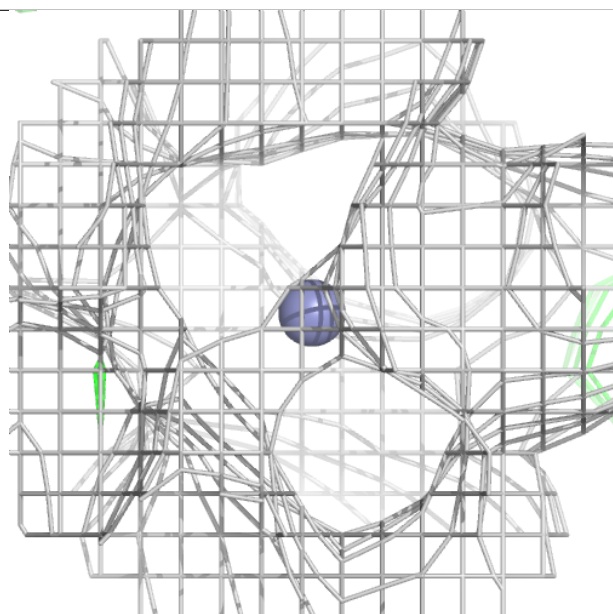
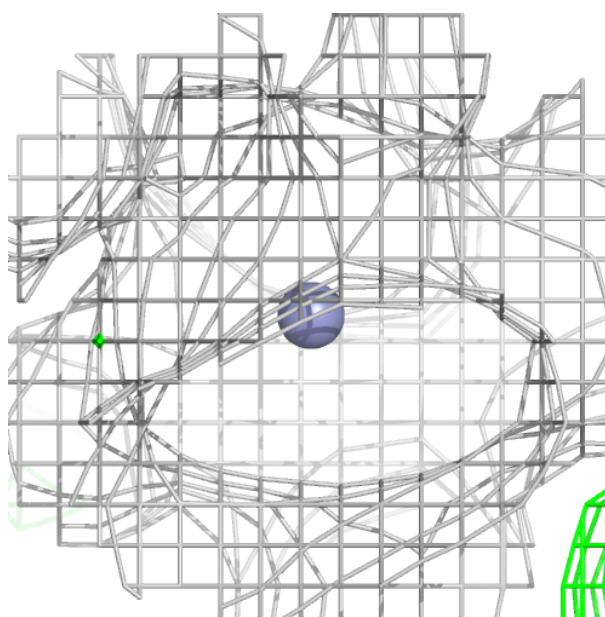
Electron density around ZN F 1001:

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



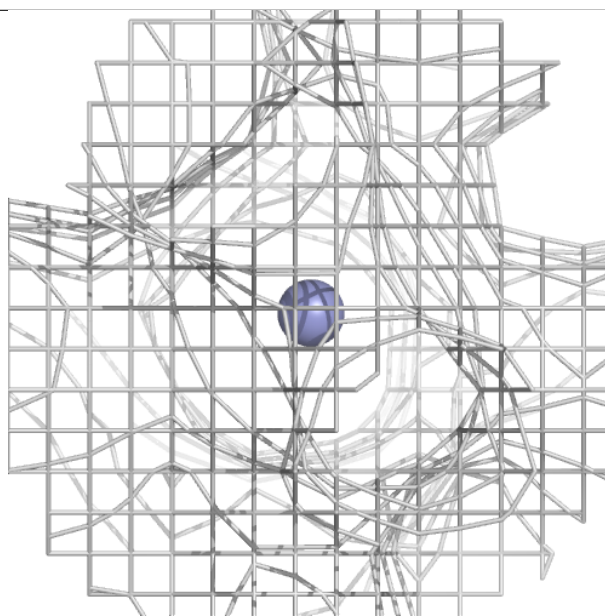
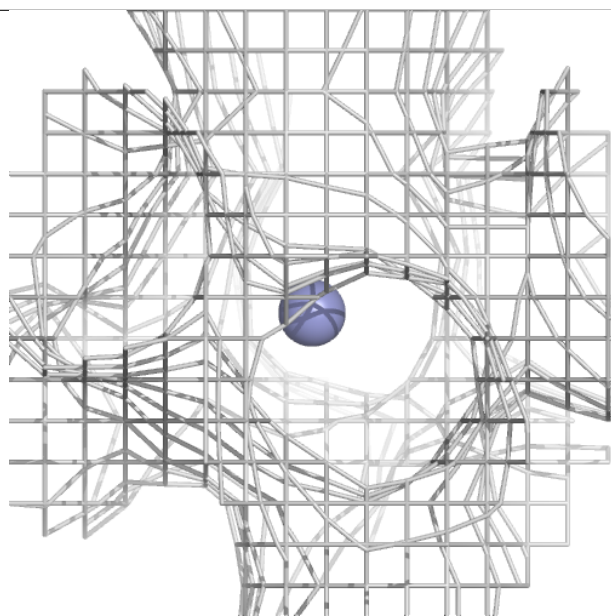
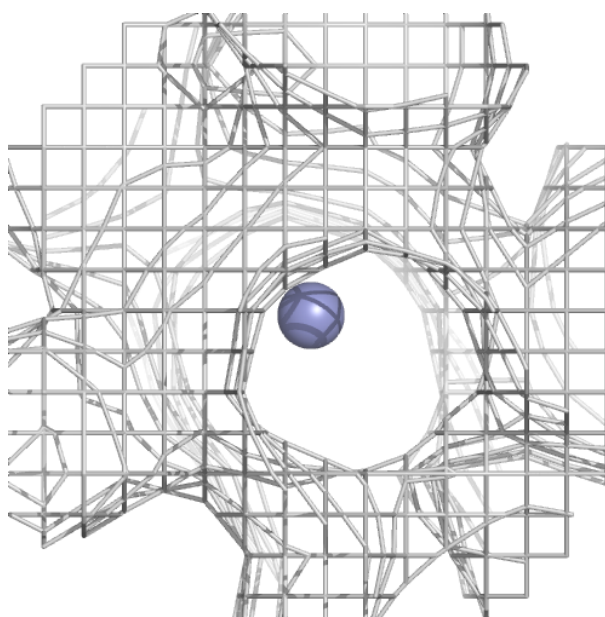
Electron density around ZN A 1006:

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



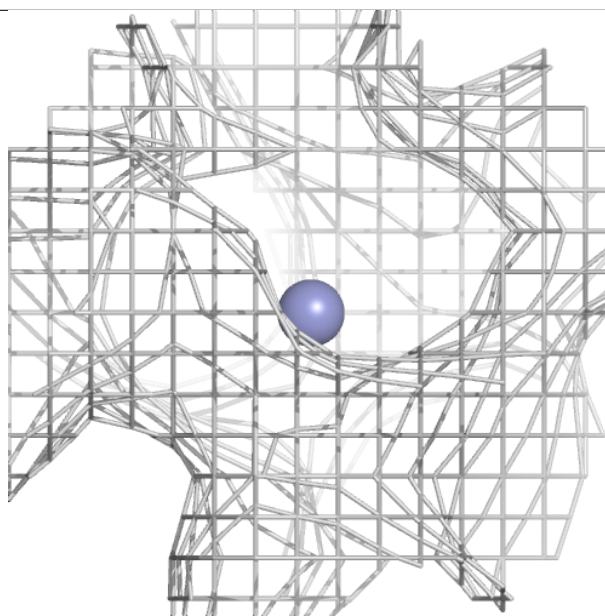
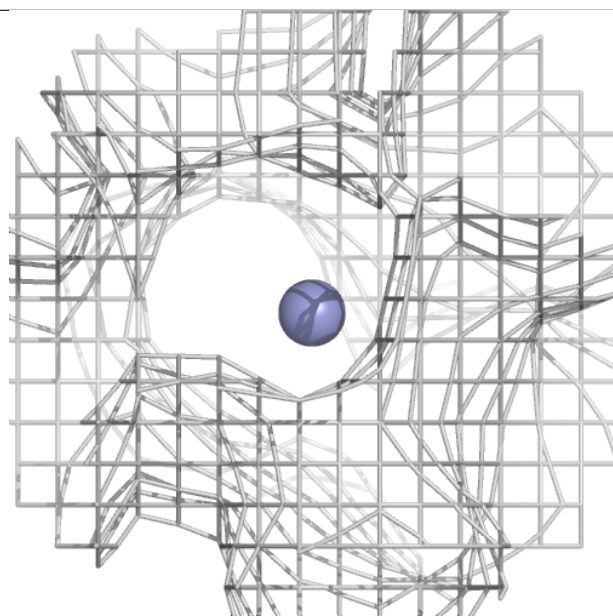
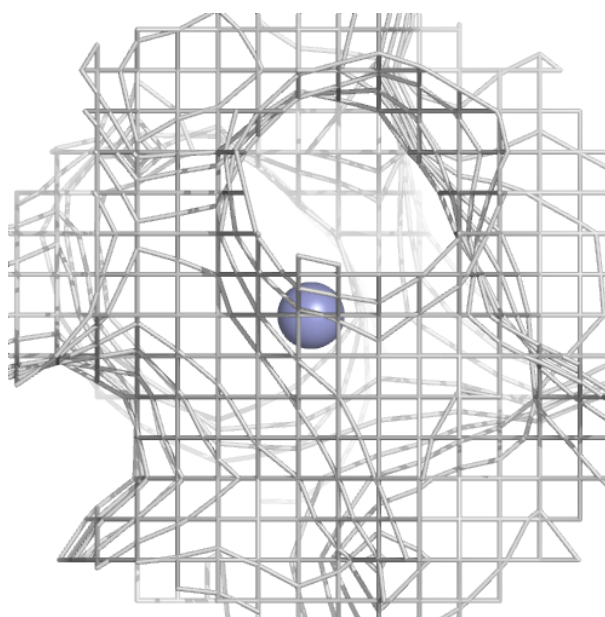
Electron density around ZN D 1001:

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



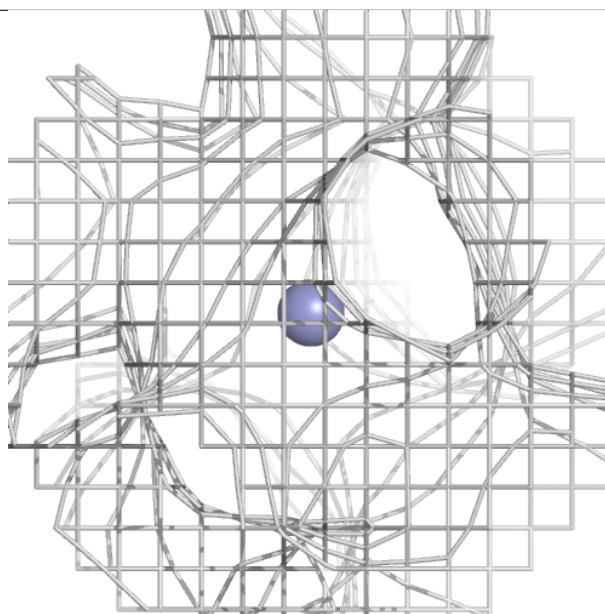
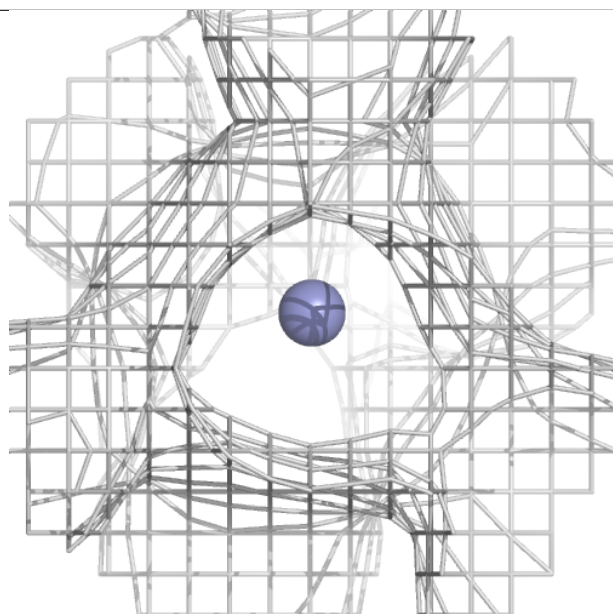
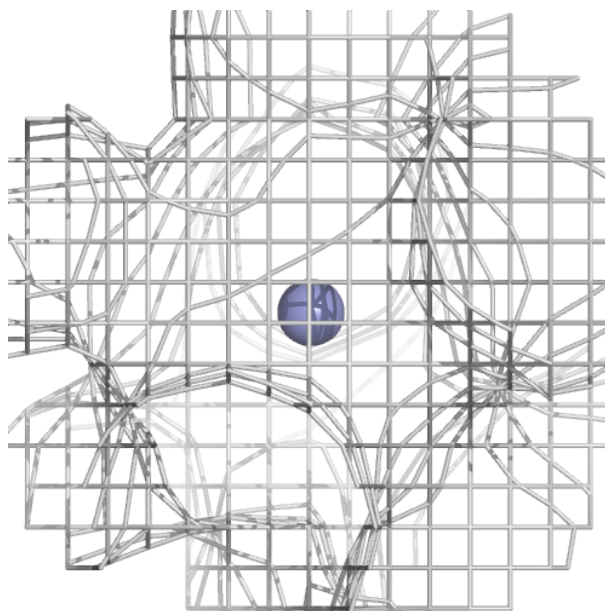
Electron density around ZN B 1001:

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



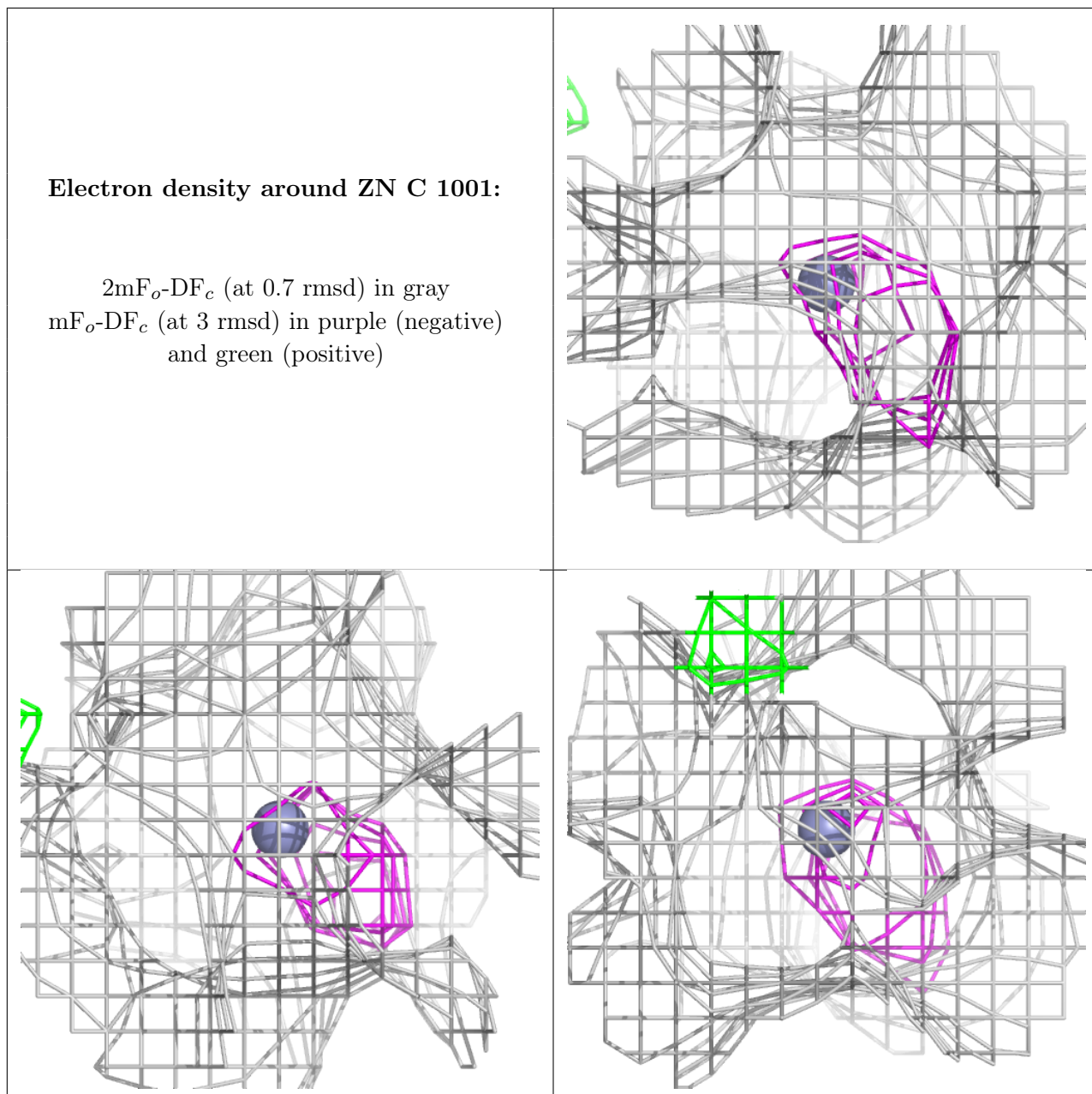
Electron density around ZN A 1001:

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



Electron density around ZN C 1001:

$2mF_o-DF_c$ (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.