



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 03:32 PM UTC

PDB ID : 9IYK / pdb_00009iyk
Title : Crystal structure of hSOD1 in C121 space group
Authors : Yapici, I.; DeMirci, H.
Deposited on : 2024-07-30
Resolution : 2.34 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:

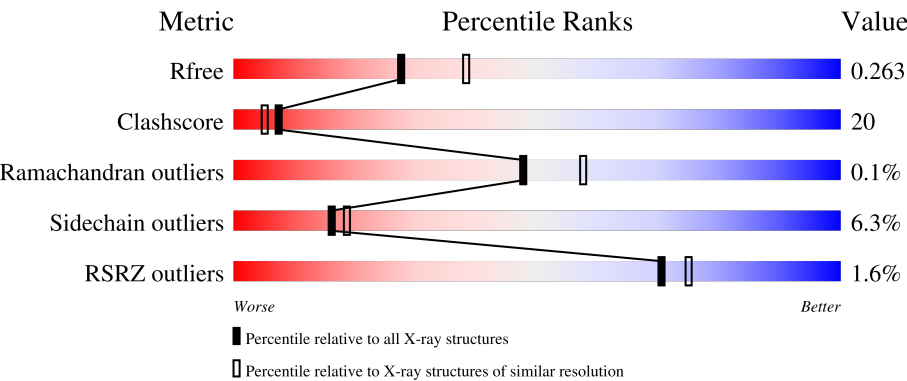
MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal 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 i

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.34 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.



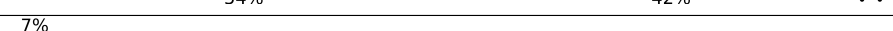
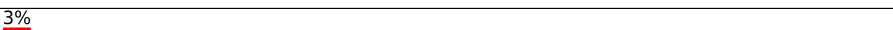
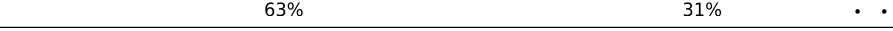
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	3031 (2.36-2.32)
Clashscore	190562	3127 (2.36-2.32)
Ramachandran outliers	187476	3095 (2.36-2.32)
Sidechain outliers	187428	3095 (2.36-2.32)
RSRZ outliers	180081	3033 (2.36-2.32)

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	156	65%27%6%.
1	B	156	75%22%.
1	C	156	76%19%..
1	D	156	3% 60%37%.
1	E	156	% 78%19%..

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Mol	Chain	Length	Quality of chain
1	F	156	 70% 26% ..
1	G	156	 54% 42% ..
1	H	156	 44% 51% ..
1	I	156	 % 60% 37% ..
1	J	156	 3% 63% 31% ..
1	K	156	 68% 28% ..
1	L	156	 % 65% 30% ..

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
4	ACT	C	201	-	-	X	-
4	ACT	C	204	-	-	X	-
5	SO4	C	205	-	-	X	-

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 14391 atoms, of which 12 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 Superoxide dismutase [Cu-Zn].

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	153	Total	C	N	O	S	0	6	0
			1149	703	209	232	5			
1	C	153	Total	C	N	O	S	0	5	0
			1140	699	206	230	5			
1	K	152	Total	C	N	O	S	0	7	0
			1146	702	207	232	5			
1	F	153	Total	C	N	O	S	0	6	0
			1145	702	207	232	4			
1	E	153	Total	C	N	O	S	0	9	0
			1165	714	210	236	5			
1	D	156	Total	C	N	O	S	0	6	0
			1171	717	212	236	6			
1	B	156	Total	C	N	O	S	0	7	0
			1172	718	212	236	6			
1	G	152	Total	C	N	O	S	0	5	0
			1137	697	206	230	4			
1	I	153	Total	C	N	O	S	0	6	0
			1144	702	206	231	5			
1	L	153	Total	C	N	O	S	0	7	0
			1156	707	210	234	5			
1	J	153	Total	C	N	O	S	0	3	0
			1132	696	206	226	4			
1	H	153	Total	C	N	O	S	0	8	0
			1158	709	207	238	4			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	SER	-	expression tag	UNP P00441
A	-1	HIS	-	expression tag	UNP P00441
C	-2	SER	-	expression tag	UNP P00441
C	-1	HIS	-	expression tag	UNP P00441
K	-2	SER	-	expression tag	UNP P00441

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Chain	Residue	Modelled	Actual	Comment	Reference
K	-1	HIS	-	expression tag	UNP P00441
F	-2	SER	-	expression tag	UNP P00441
F	-1	HIS	-	expression tag	UNP P00441
E	-2	SER	-	expression tag	UNP P00441
E	-1	HIS	-	expression tag	UNP P00441
D	-2	SER	-	expression tag	UNP P00441
D	-1	HIS	-	expression tag	UNP P00441
B	-2	SER	-	expression tag	UNP P00441
B	-1	HIS	-	expression tag	UNP P00441
G	-2	SER	-	expression tag	UNP P00441
G	-1	HIS	-	expression tag	UNP P00441
I	-2	SER	-	expression tag	UNP P00441
I	-1	HIS	-	expression tag	UNP P00441
L	-2	SER	-	expression tag	UNP P00441
L	-1	HIS	-	expression tag	UNP P00441
J	-2	SER	-	expression tag	UNP P00441
J	-1	HIS	-	expression tag	UNP P00441
H	-2	SER	-	expression tag	UNP P00441
H	-1	HIS	-	expression tag	UNP P00441

- Molecule 2 is COPPER (II) ION (CCD ID: CU) (formula: Cu) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Cu 1 1	0	0
2	C	1	Total Cu 1 1	0	0
2	K	1	Total Cu 1 1	0	0
2	F	1	Total Cu 1 1	0	0
2	E	1	Total Cu 1 1	0	0
2	D	1	Total Cu 1 1	0	0
2	B	1	Total Cu 1 1	0	0
2	G	1	Total Cu 1 1	0	0
2	I	1	Total Cu 1 1	0	0

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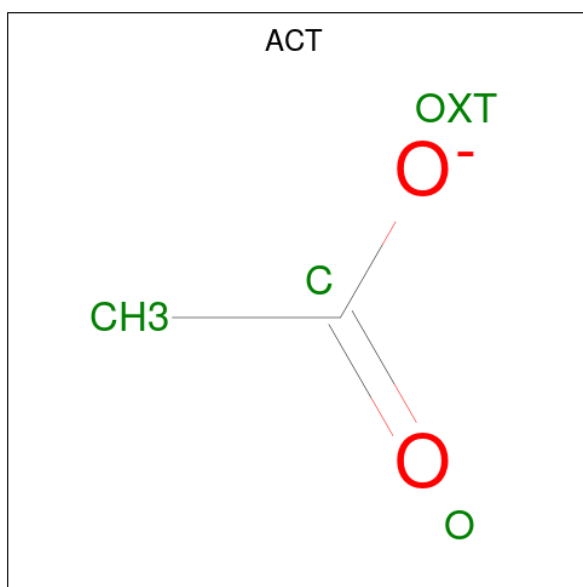
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	L	1	Total	Cu	0	0
			1	1		
2	J	1	Total	Cu	0	0
			1	1		
2	H	1	Total	Cu	0	0
			1	1		

- Molecule 3 is ZINC ION (CCD ID: ZN) (formula: Zn) (labeled as "Ligand of Interest" by depositor).

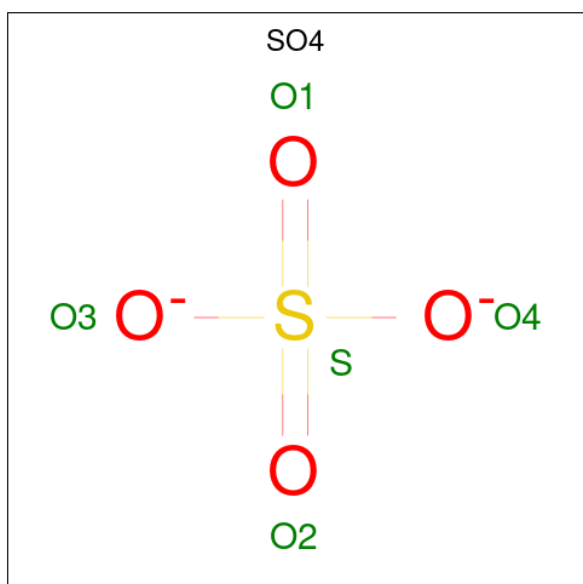
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	2	Total	Zn	0	0
			2	2		
3	C	1	Total	Zn	0	0
			1	1		
3	K	1	Total	Zn	0	0
			1	1		
3	F	1	Total	Zn	0	0
			1	1		
3	E	1	Total	Zn	0	0
			1	1		
3	D	1	Total	Zn	0	0
			1	1		
3	B	1	Total	Zn	0	0
			1	1		
3	G	1	Total	Zn	0	0
			1	1		
3	I	1	Total	Zn	0	0
			1	1		
3	L	1	Total	Zn	0	0
			1	1		
3	J	1	Total	Zn	0	0
			1	1		
3	H	1	Total	Zn	0	0
			1	1		

- Molecule 4 is ACETATE ION (CCD ID: ACT) (formula: C₂H₃O₂) (labeled as "Ligand of Interest" by depositor).



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

- Molecule 5 is SULFATE ION (CCD ID: SO4) (formula: O₄S) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total O S 5 4 1	0	0
5	C	1	Total O S 5 4 1	0	0
5	E	1	Total O S 5 4 1	0	0
5	I	1	Total O S 5 4 1	0	0

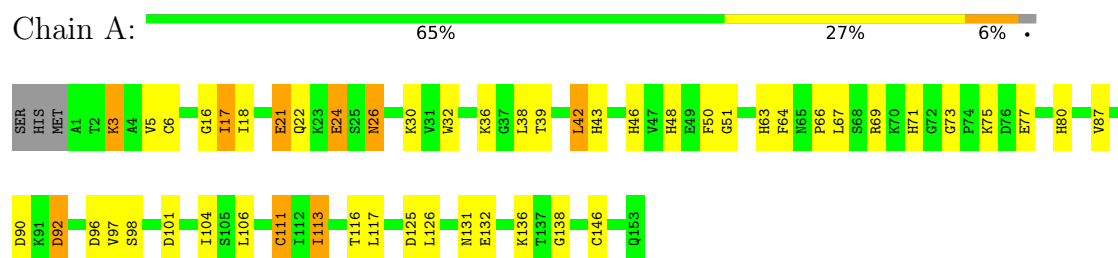
- Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	68	Total O 70 70	0	2
6	C	51	Total O 51 51	0	0
6	K	75	Total O 76 76	0	1
6	F	39	Total O 40 40	0	1
6	E	43	Total O 46 46	0	3
6	D	31	Total O 33 33	0	2
6	B	50	Total O 50 50	0	0
6	G	35	Total O 36 36	0	1
6	I	25	Total O 25 25	0	0
6	L	37	Total O 37 37	0	0
6	J	21	Total O 21 21	0	0
6	H	18	Total O 18 18	0	0

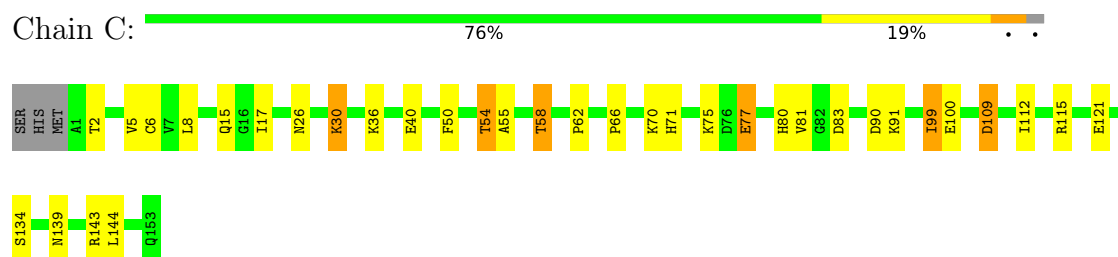
3 Residue-property plots [i](#)

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.

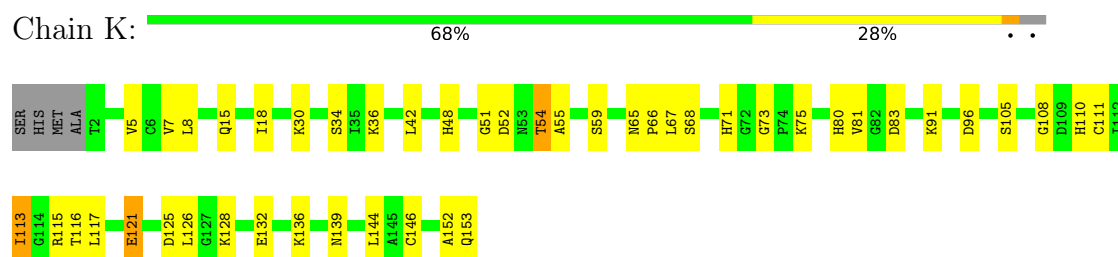
• Molecule 1: Superoxide dismutase [Cu-Zn]



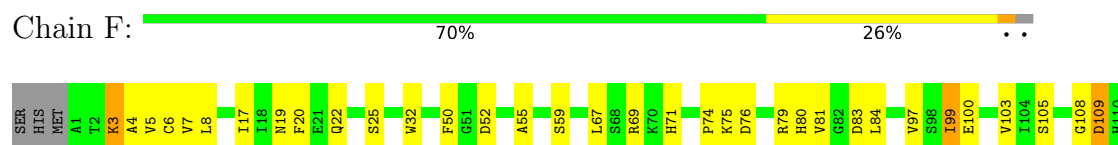
• Molecule 1: Superoxide dismutase [Cu-Zn]



• Molecule 1: Superoxide dismutase [Cu-Zn]

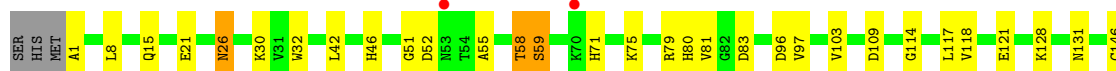
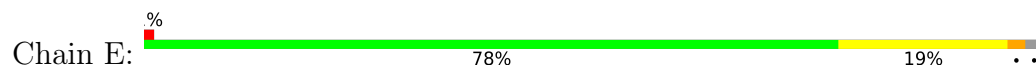


• Molecule 1: Superoxide dismutase [Cu-Zn]





- Molecule 1: Superoxide dismutase [Cu-Zn]



- Molecule 1: Superoxide dismutase [Cu-Zn]



- Molecule 1: Superoxide dismutase [Cu-Zn]

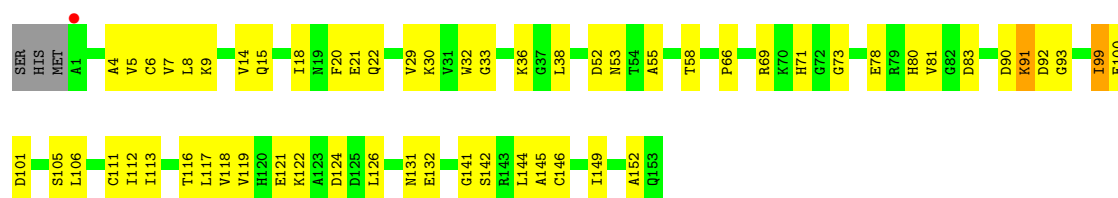


- Molecule 1: Superoxide dismutase [Cu-Zn]

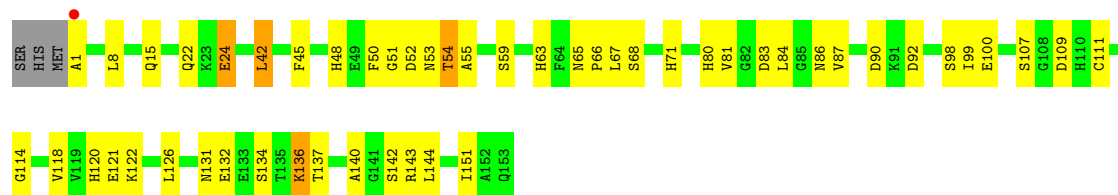


- Molecule 1: Superoxide dismutase [Cu-Zn]

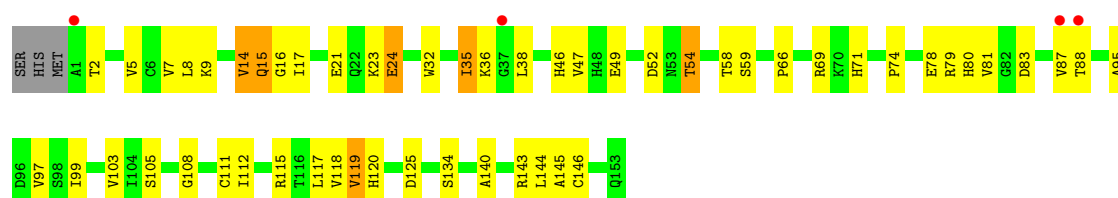




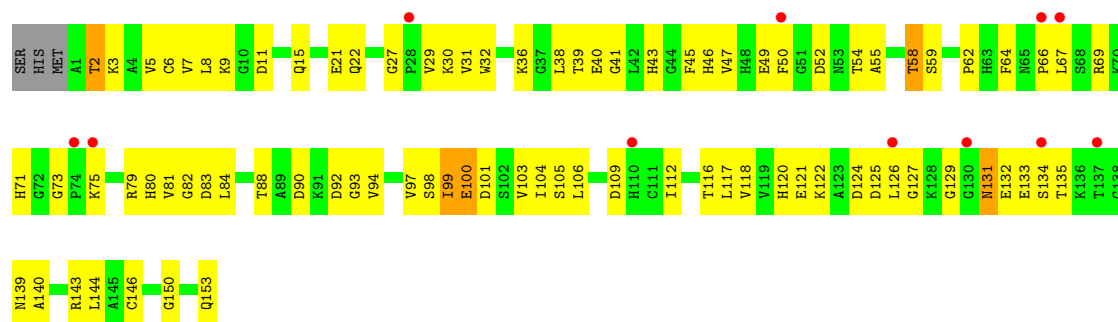
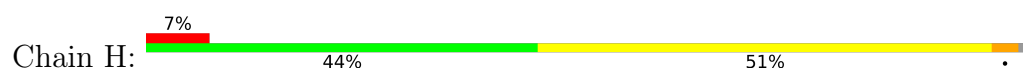
• Molecule 1: Superoxide dismutase [Cu-Zn]



• Molecule 1: Superoxide dismutase [Cu-Zn]



• Molecule 1: Superoxide dismutase [Cu-Zn]



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	178.19Å 138.25Å 112.93Å 90.00° 129.20° 90.00°	Depositor
Resolution (Å)	48.92 – 2.34 48.92 – 2.34	Depositor EDS
% Data completeness (in resolution range)	86.3 (48.92-2.34) 86.3 (48.92-2.34)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.20 (at 2.34Å)	Xtriage
Refinement program	PHENIX (1.20.1_4487: ???)	Depositor
R, R_{free}	0.198 , 0.263 0.199 , 0.263	Depositor DCC
R_{free} test set	3833 reflections (4.29%)	wwPDB-VP
Wilson B-factor (Å ²)	41.9	Xtriage
Anisotropy	0.271	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 70.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.026 for k+l,h+l,-l 0.027 for -k+l,-h-l,-l 0.075 for -h-2*l,-k,l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	14391	wwPDB-VP
Average B, all atoms (Å ²)	69.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.11% 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: ACT, CU, SO4, 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	0.22	0/1178	0.42	0/1589
1	B	0.21	0/1211	0.40	0/1633
1	C	0.21	0/1172	0.38	0/1581
1	D	0.18	0/1207	0.34	0/1628
1	E	0.18	0/1203	0.36	0/1622
1	F	0.17	0/1180	0.35	0/1592
1	G	0.16	0/1169	0.35	0/1577
1	H	0.15	0/1196	0.32	0/1614
1	I	0.16	0/1179	0.31	0/1591
1	J	0.16	0/1161	0.34	0/1565
1	K	0.19	0/1181	0.38	0/1593
1	L	0.21	0/1188	0.40	0/1603
All	All	0.18	0/14225	0.36	0/19188

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	1149	0	1094	45	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	1172	0	1121	28	1
1	C	1140	0	1089	32	0
1	D	1171	0	1115	66	0
1	E	1165	0	1115	33	0
1	F	1145	0	1095	34	0
1	G	1137	0	1082	51	0
1	H	1158	0	1100	98	1
1	I	1144	0	1096	59	0
1	J	1132	0	1089	48	0
1	K	1146	0	1091	45	0
1	L	1156	0	1099	43	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
2	E	1	0	0	0	0
2	F	1	0	0	0	0
2	G	1	0	0	0	0
2	H	1	0	0	0	0
2	I	1	0	0	0	0
2	J	1	0	0	0	0
2	K	1	0	0	0	0
2	L	1	0	0	0	0
3	A	2	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
3	E	1	0	0	0	0
3	F	1	0	0	0	0
3	G	1	0	0	0	0
3	H	1	0	0	0	0
3	I	1	0	0	0	0
3	J	1	0	0	0	0
3	K	1	0	0	0	0
3	L	1	0	0	0	0
4	A	4	3	3	0	0
4	B	4	3	3	0	0
4	C	8	6	6	6	0
5	A	5	0	0	0	0
5	C	5	0	0	2	0
5	E	5	0	0	0	0
5	I	5	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	A	70	0	0	5	0
6	B	50	0	0	1	0
6	C	51	0	0	3	0
6	D	33	0	0	7	0
6	E	46	0	0	3	0
6	F	40	0	0	1	0
6	G	36	0	0	2	0
6	H	18	0	0	4	0
6	I	25	0	0	1	0
6	J	21	0	0	5	0
6	K	76	0	0	3	0
6	L	37	0	0	4	0
All	All	14379	12	13198	552	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 20.

All (552) 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:121:GLU:HB2	1:D:144:LEU:HD21	1.34	1.07
1:L:132:GLU:HG2	1:L:136:LYS:HE2	1.36	1.05
1:H:121:GLU:HB2	1:H:144:LEU:HD21	1.37	1.04
1:H:121:GLU:HA	1:H:144:LEU:HD11	1.43	0.99
1:I:121:GLU:HA	1:I:144:LEU:HD11	1.40	0.99
1:H:22:GLN:HB2	1:H:29:VAL:HG22	1.46	0.98
1:J:14:VAL:HG13	1:J:145:ALA:HB2	1.49	0.94
1:C:62:PRO:HD3	4:C:201:ACT:H2	1.50	0.94
1:I:73:GLY:HA2	1:I:126:LEU:HD12	1.50	0.92
1:C:62:PRO:CD	4:C:201:ACT:H2	2.02	0.89
1:A:90:ASP:HB2	1:A:92:ASP:OD1	1.74	0.88
1:I:5:VAL:HG12	1:I:152:ALA:HB2	1.55	0.87
1:D:73:GLY:HA3	1:D:126:LEU:HD13	1.58	0.86
1:A:111[A]:CYS:SG	6:A:332:HOH:O	2.33	0.85
1:I:66:PRO:HG2	1:I:81:VAL:HG21	1.57	0.84
1:C:15:GLN:HE21	1:C:36:LYS:HE3	1.41	0.83
1:H:79:ARG:NE	1:H:80:HIS:O	2.10	0.83
1:I:112:ILE:HD12	1:I:149:ILE:HD13	1.61	0.82
1:C:109:ASP:OD2	1:D:-2:SER:HB3	1.79	0.82
1:B:30:LYS:NZ	1:B:100:GLU:OE1	2.11	0.82
1:D:70:LYS:O	1:D:79:ARG:HA	1.81	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:80:HIS:HB2	1:G:83:ASP:OD1	1.80	0.81
1:F:5:VAL:HG23	1:F:152:ALA:HB2	1.62	0.81
1:C:77:GLU:HB2	4:C:204:ACT:H2	1.64	0.80
1:D:121:GLU:HA	1:D:144:LEU:HD11	1.64	0.80
1:D:101:ASP:OD1	1:D:103:VAL:N	2.14	0.80
1:J:15:GLN:O	6:J:301:HOH:O	1.99	0.79
1:A:111[A]:CYS:SG	1:A:113:ILE:HB	2.23	0.79
1:G:66:PRO:HD2	1:G:81:VAL:CG2	2.14	0.78
1:I:80:HIS:HB2	1:I:83:ASP:OD2	1.84	0.78
1:F:52:ASP:O	1:F:59:SER:HB2	1.84	0.78
1:H:81:VAL:O	1:H:104:ILE:HG23	1.85	0.75
1:C:40:GLU:OE2	6:C:301:HOH:O	2.04	0.75
1:D:86:ASN:O	6:D:301:HOH:O	2.03	0.75
1:L:1:ALA:HB1	6:L:318:HOH:O	1.85	0.75
1:H:81:VAL:HG13	1:H:103:VAL:HG12	1.70	0.74
1:H:129:GLY:O	6:H:301:HOH:O	2.05	0.74
1:B:55:ALA:HB3	1:B:59:SER:OG	1.88	0.74
1:G:40:GLU:HG3	1:G:91:LYS:HD2	1.69	0.73
1:H:121:GLU:CB	1:H:144:LEU:HD21	2.18	0.73
1:K:113:ILE:CD1	1:L:114:GLY:HA3	2.18	0.73
1:H:80:HIS:HB2	1:H:83:ASP:OD1	1.89	0.72
1:A:3:LYS:HD3	1:A:21:GLU:HG3	1.70	0.72
1:K:52:ASP:O	1:K:59:SER:HB2	1.90	0.72
1:H:135:THR:HA	6:H:303:HOH:O	1.89	0.72
1:H:73:GLY:HA2	1:H:126:LEU:HD22	1.69	0.71
1:H:81:VAL:HG12	1:H:104:ILE:HG22	1.72	0.71
1:J:81:VAL:HG13	1:J:103:VAL:HG12	1.72	0.71
1:H:36:LYS:HB3	1:H:94:VAL:HG22	1.73	0.71
1:H:66:PRO:HG2	1:H:81:VAL:HG21	1.71	0.71
1:D:121:GLU:CB	1:D:144:LEU:HD21	2.15	0.70
1:D:79:ARG:HD3	1:D:83:ASP:HB2	1.73	0.70
1:D:69:ARG:O	6:D:303:HOH:O	2.09	0.70
1:F:84:LEU:HD22	1:F:99:ILE:HD11	1.72	0.70
1:K:71:HIS:CE1	1:K:83:ASP:OD1	2.44	0.69
1:D:130:GLY:O	6:D:304:HOH:O	2.11	0.69
1:G:105:SER:O	1:G:111:CYS:HA	1.92	0.69
1:G:50:PHE:CZ	1:H:153:GLN:HB2	2.27	0.69
1:H:29:VAL:HB	6:H:312:HOH:O	1.92	0.69
1:H:8:LEU:HD11	1:H:117:LEU:HD23	1.73	0.69
1:A:16:GLY:C	1:A:17:ILE:HD13	2.18	0.69
1:D:40:GLU:OE2	1:D:91:LYS:NZ	2.16	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:38:LEU:O	1:B:93:GLY:HA2	1.91	0.69
1:J:35:ILE:HA	6:J:301:HOH:O	1.93	0.68
1:G:5:VAL:HG12	1:G:152:ALA:HB2	1.76	0.68
1:I:121:GLU:HB3	1:I:144:LEU:HD21	1.75	0.68
1:H:64:PHE:CZ	1:H:66:PRO:HG3	2.29	0.68
1:A:111[B]:CYS:SG	6:A:332:HOH:O	2.52	0.68
1:I:66:PRO:HD2	1:I:81:VAL:HB	1.74	0.68
1:F:109:ASP:OD1	1:F:109:ASP:N	2.25	0.67
1:J:47:VAL:HG11	1:J:112:ILE:HG22	1.76	0.67
1:J:118:VAL:HG11	1:J:143:ARG:HG2	1.75	0.67
1:D:65:ASN:HB2	1:D:80:HIS:HB3	1.76	0.67
1:F:5:VAL:HG23	1:F:152:ALA:CB	2.25	0.66
1:L:52:ASP:O	1:L:59:SER:HB2	1.94	0.66
1:K:121:GLU:HB2	1:K:144:LEU:HD21	1.78	0.66
1:I:73:GLY:CA	1:I:126:LEU:HD12	2.26	0.66
1:B:96[A]:ASP:OD1	6:B:302:HOH:O	2.14	0.65
1:H:45:PHE:O	1:H:84:LEU:HB2	1.97	0.65
1:A:3:LYS:HD3	1:A:21:GLU:CG	2.27	0.65
1:A:30:LYS:HD3	6:A:338:HOH:O	1.97	0.65
1:K:80:HIS:HB2	1:K:83:ASP:OD2	1.96	0.65
1:C:62:PRO:HD2	4:C:201:ACT:H2	1.78	0.64
1:H:31:VAL:HB	1:H:99:ILE:HG23	1.79	0.64
1:H:66:PRO:HG2	1:H:81:VAL:CG2	2.27	0.64
1:F:7:VAL:HG22	1:F:17:ILE:CD1	2.28	0.64
1:D:66:PRO:HD2	1:D:81:VAL:CG2	2.28	0.64
1:H:52:ASP:O	1:H:59:SER:HB2	1.97	0.64
1:L:53[A]:ASN:O	1:L:55:ALA:N	2.31	0.64
1:G:64:PHE:CE1	1:G:66:PRO:HG3	2.33	0.63
1:J:52:ASP:OD1	1:J:54:THR:OG1	2.16	0.63
1:H:118:VAL:HG11	1:H:143:ARG:HG2	1.81	0.63
1:A:90:ASP:HB2	1:A:92:ASP:CG	2.22	0.63
1:J:38:LEU:HD23	1:J:144:LEU:HD12	1.80	0.63
1:H:2:THR:HG21	1:H:106:LEU:HB2	1.79	0.63
1:L:121:GLU:HG2	1:L:142:SER:H	1.63	0.63
1:G:80:HIS:HB2	1:G:83:ASP:CG	2.23	0.62
1:L:65:ASN:ND2	1:L:68:SER:HA	2.13	0.62
1:A:77:GLU:OE2	1:D:2:THR:HG21	1.99	0.62
1:H:64:PHE:O	1:H:80:HIS:HB3	2.00	0.62
1:I:4:ALA:HB3	1:I:20:PHE:HB2	1.81	0.62
1:H:104:ILE:HG13	6:H:312:HOH:O	1.98	0.62
1:G:21:GLU:OE2	1:G:23:LYS:HG3	1.99	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:125:ASP:OD1	1:G:138:GLY:HA2	2.00	0.62
1:I:66:PRO:CG	1:I:81:VAL:HG21	2.29	0.62
1:A:39:THR:O	1:A:43:HIS:HE1	1.82	0.62
1:L:53[B]:ASN:O	1:L:55:ALA:N	2.32	0.62
1:K:111[B]:CYS:SG	1:K:113:ILE:HG23	2.40	0.62
1:I:30:LYS:NZ	1:I:30:LYS:HB3	2.15	0.62
1:K:111[A]:CYS:SG	1:K:113:ILE:HG23	2.40	0.61
1:D:72:GLY:HA3	1:D:79:ARG:HB3	1.82	0.61
1:B:80:HIS:HB2	1:B:83:ASP:OD2	1.99	0.61
1:I:90:ASP:HB2	1:I:92:ASP:OD1	2.00	0.61
1:H:120:HIS:CE1	1:H:143:ARG:HG2	2.34	0.61
1:E:1:ALA:N	6:E:304:HOH:O	2.32	0.61
1:I:121:GLU:HG2	1:I:142:SER:N	2.16	0.61
1:J:21:GLU:HB2	1:J:32[A]:TRP:HH2	1.66	0.61
1:J:14:VAL:CG1	1:J:145:ALA:HB2	2.26	0.61
1:D:79:ARG:HD2	1:D:80:HIS:O	2.01	0.61
1:J:5:VAL:HG22	1:J:17:ILE:HD11	1.82	0.61
1:J:5:VAL:CG2	1:J:17:ILE:HD11	2.30	0.61
1:K:67:LEU:HD12	6:K:332:HOH:O	2.01	0.60
1:C:80:HIS:HB2	1:C:83:ASP:CG	2.26	0.60
1:I:5:VAL:CG1	1:I:152:ALA:HB2	2.28	0.60
1:I:38:LEU:O	1:I:93:GLY:HA2	2.01	0.60
1:B:92:ASP:OD1	1:B:92:ASP:N	2.33	0.60
1:I:99:ILE:HD11	1:I:101:ASP:HB2	1.84	0.60
1:F:50:PHE:CZ	1:E:153:GLN:HB2	2.37	0.60
1:H:2:THR:HB	1:H:106:LEU:HD12	1.82	0.60
1:G:14:VAL:HG21	1:G:144:LEU:HB3	1.84	0.60
1:L:121:GLU:HG2	1:L:142:SER:N	2.17	0.59
1:H:122:LYS:HB2	1:H:140:ALA:HA	1.84	0.59
1:H:3:LYS:HD2	1:H:153:GLN:OXT	2.03	0.59
1:K:5:VAL:HG12	1:K:152:ALA:HB2	1.85	0.59
1:K:80:HIS:CG	1:K:83:ASP:OD2	2.55	0.59
1:E:71:HIS:HB2	1:E:80:HIS:CE1	2.37	0.59
1:I:121:GLU:CA	1:I:144:LEU:HD11	2.25	0.59
1:H:22:GLN:CB	1:H:29:VAL:HG22	2.26	0.59
1:D:118:VAL:HG12	1:D:120:HIS:HD2	1.67	0.59
1:B:109:ASP:OD1	1:B:109:ASP:N	2.32	0.59
1:J:21:GLU:HB2	1:J:32[A]:TRP:CH2	2.38	0.59
1:B:55:ALA:O	1:B:58:THR:HG23	2.03	0.58
1:B:19:ASN:HB2	1:B:32[B]:TRP:NE1	2.19	0.58
1:H:21:GLU:O	1:H:29:VAL:HG13	2.04	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:50:PHE:CZ	1:D:153:GLN:HB2	2.38	0.58
1:F:55:ALA:N	1:F:59:SER:OG	2.34	0.58
1:D:61:GLY:O	1:D:143:ARG:NH2	2.35	0.58
1:A:16:GLY:O	1:A:17:ILE:HD13	2.03	0.58
1:E:96[B]:ASP:OD1	6:E:301:HOH:O	2.17	0.58
1:J:23:LYS:O	1:J:24:GLU:HG3	2.04	0.58
1:F:22:GLN:NE2	1:F:25:SER:O	2.37	0.58
1:D:65:ASN:HB2	1:D:80:HIS:CB	2.33	0.58
1:I:71:HIS:HB2	1:I:80:HIS:CE1	2.39	0.58
1:G:120:HIS:HB3	1:G:140:ALA:O	2.04	0.57
1:H:64:PHE:CE1	1:H:66:PRO:HG3	2.40	0.57
1:I:80:HIS:HB2	1:I:83:ASP:CG	2.29	0.57
1:L:118:VAL:HG11	1:L:143:ARG:HG2	1.85	0.57
1:H:105:SER:HG	1:H:106:LEU:H	1.52	0.57
1:J:16:GLY:HA2	6:J:301:HOH:O	2.04	0.57
1:H:122:LYS:HB2	1:H:140:ALA:C	2.29	0.57
1:H:81:VAL:HG12	1:H:104:ILE:CG2	2.34	0.57
1:K:7:VAL:HG21	1:L:53[A]:ASN:HB3	1.86	0.57
1:F:5:VAL:HG12	1:F:6:CYS:N	2.20	0.57
1:C:75:LYS:HD3	1:K:128:LYS:NZ	2.20	0.57
1:K:110:HIS:HB2	6:K:305:HOH:O	2.05	0.57
1:K:132:GLU:HG2	1:K:136:LYS:HE2	1.87	0.57
1:A:36:LYS:HE2	6:A:333:HOH:O	2.05	0.56
1:J:8:LEU:O	1:J:15:GLN:NE2	2.38	0.56
1:F:80:HIS:N	1:F:83:ASP:OD2	2.33	0.56
1:E:52:ASP:O	1:E:59:SER:HB2	2.05	0.56
1:G:71:HIS:HB2	1:G:80:HIS:CE1	2.40	0.56
1:J:120:HIS:HB3	1:J:140:ALA:O	2.06	0.56
1:A:39:THR:HG22	1:A:43:HIS:NE2	2.20	0.56
1:K:80:HIS:CB	1:K:83:ASP:OD2	2.54	0.56
1:J:23:LYS:HD3	6:J:316:HOH:O	2.05	0.56
1:H:32[B]:TRP:HA	1:H:97:VAL:O	2.06	0.56
1:D:101:ASP:OD1	1:D:102:SER:N	2.39	0.56
1:I:55:ALA:O	1:I:58[A]:THR:HG22	2.06	0.56
1:D:52:ASP:O	1:D:59:SER:HB2	2.06	0.56
1:F:7:VAL:HG22	1:F:17:ILE:HD13	1.88	0.56
1:G:49:GLU:O	1:G:115:ARG:NH1	2.37	0.55
1:A:21:GLU:HB2	1:A:32[B]:TRP:CZ3	2.41	0.55
1:A:64:PHE:CD1	1:A:66:PRO:HD3	2.41	0.55
1:C:90:ASP:HB3	6:C:304:HOH:O	2.05	0.55
1:H:131:ASN:OD1	1:H:133:GLU:HB3	2.06	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:80:HIS:N	1:D:83:ASP:OD2	2.37	0.55
1:J:47:VAL:HG11	1:J:112:ILE:CG2	2.36	0.55
1:H:134:SER:HB2	1:H:139:ASN:ND2	2.21	0.55
1:H:105:SER:O	1:H:112:ILE:HG12	2.06	0.55
1:A:117:LEU:O	1:A:146:CYS:HA	2.07	0.55
1:E:79:ARG:NH2	1:E:103:VAL:HB	2.22	0.55
1:D:45:PHE:O	1:D:84:LEU:HB2	2.07	0.55
1:L:107:SER:HA	1:L:111[C]:CYS:SG	2.47	0.55
1:G:91:LYS:O	6:G:301:HOH:O	2.18	0.55
1:H:32[A]:TRP:HA	1:H:97:VAL:O	2.06	0.55
1:G:73:GLY:HA3	1:G:126:LEU:HD22	1.88	0.54
1:H:47:VAL:HG11	1:H:112:ILE:HG22	1.87	0.54
1:D:81:VAL:HG13	1:D:103:VAL:HG12	1.88	0.54
1:A:42[A]:LEU:HA	1:A:87:VAL:O	2.08	0.54
1:K:113:ILE:HD11	1:L:114:GLY:HA3	1.89	0.54
1:E:80:HIS:HB2	1:E:83:ASP:CG	2.32	0.54
1:H:38:LEU:HD22	1:H:43:HIS:ND1	2.22	0.54
1:C:54:THR:O	1:G:24:GLU:HG3	2.06	0.54
1:C:30:LYS:NZ	1:C:30:LYS:HB3	2.23	0.54
1:I:8:LEU:O	1:I:15:GLN:HA	2.07	0.54
1:D:65:ASN:HB2	1:D:80:HIS:HA	1.90	0.54
1:J:14:VAL:HG11	1:J:144:LEU:HB3	1.89	0.54
1:F:32[A]:TRP:HA	1:F:97:VAL:O	2.08	0.54
1:C:77:GLU:OE1	1:B:2:THR:HG21	2.07	0.54
1:F:32[B]:TRP:HA	1:F:97:VAL:O	2.08	0.54
1:L:54:THR:HG22	1:L:55:ALA:N	2.23	0.54
1:H:79:ARG:O	1:H:79:ARG:HD3	2.08	0.54
1:H:131:ASN:O	1:H:134:SER:HB3	2.08	0.54
1:F:99:ILE:HG13	1:F:100:GLU:N	2.23	0.53
1:I:9:LYS:NZ	6:I:301:HOH:O	2.20	0.53
1:F:50:PHE:CE2	1:E:153:GLN:HB2	2.43	0.53
1:K:54:THR:HG21	6:K:340:HOH:O	2.06	0.53
1:D:118:VAL:HG12	1:D:120:HIS:CD2	2.44	0.53
1:A:69:ARG:HH22	1:D:2:THR:HB	1.73	0.53
1:D:22:GLN:HE22	1:D:105:SER:HA	1.73	0.53
1:D:125:ASP:OD2	1:D:134:SER:OG	2.22	0.53
1:H:22:GLN:HE22	1:H:105:SER:CB	2.22	0.53
1:I:121:GLU:HG2	1:I:142:SER:H	1.72	0.53
1:K:117:LEU:O	1:K:146:CYS:HA	2.09	0.53
1:D:131[A]:ASN:OD1	1:D:133:GLU:HB3	2.09	0.52
1:G:45:PHE:O	1:G:84:LEU:HB2	2.09	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:132:GLU:OE2	1:A:136:LYS:NZ	2.42	0.52
1:K:52:ASP:OD1	1:K:54:THR:HB	2.10	0.52
1:L:42[A]:LEU:HA	1:L:87:VAL:O	2.08	0.52
1:L:55:ALA:H	1:L:59:SER:HG	1.57	0.52
1:L:42[B]:LEU:HA	1:L:87:VAL:O	2.09	0.52
1:H:40:GLU:HG3	1:H:90[A]:ASP:C	2.33	0.52
1:C:77:GLU:CB	4:C:204:ACT:H2	2.37	0.52
1:K:125:ASP:OD2	1:K:139:ASN:ND2	2.36	0.52
1:A:64:PHE:CE1	1:A:66:PRO:HD3	2.45	0.52
1:B:38:LEU:HD22	1:B:43:HIS:CE1	2.45	0.52
1:G:51:GLY:O	1:H:150:GLY:HA3	2.10	0.52
1:D:26:ASN:OD1	1:D:26:ASN:N	2.38	0.52
1:C:112:ILE:O	1:C:115:ARG:HG3	2.10	0.51
1:G:117:LEU:O	1:G:146:CYS:HA	2.10	0.51
1:H:118:VAL:HG22	1:H:146:CYS:HB3	1.92	0.51
1:D:5:VAL:HG22	1:D:6:CYS:N	2.25	0.51
1:I:99:ILE:HD12	1:I:100:GLU:N	2.26	0.51
1:C:143:ARG:N	5:C:205:SO4:O1	2.39	0.51
1:F:81:VAL:HG13	1:F:103:VAL:HG12	1.93	0.51
1:K:66:PRO:CG	1:K:81:VAL:HG21	2.40	0.51
1:G:69:ARG:HB2	1:G:79:ARG:O	2.10	0.51
1:G:133:GLU:HB3	1:G:139:ASN:OD1	2.10	0.51
1:J:9:LYS:HA	1:J:15:GLN:NE2	2.25	0.51
1:J:32[A]:TRP:HA	1:J:97:VAL:O	2.11	0.51
1:J:46:HIS:HD1	1:J:120:HIS:CD2	2.29	0.51
1:H:47:VAL:HG11	1:H:112:ILE:CG2	2.41	0.51
1:A:132:GLU:HG2	1:A:136:LYS:NZ	2.26	0.51
1:I:118:VAL:HG22	1:I:146:CYS:HB3	1.92	0.51
1:D:66:PRO:HD2	1:D:81:VAL:HG21	1.92	0.51
1:B:28:PRO:HB3	1:B:102[A]:SER:OG	2.11	0.51
1:G:22:GLN:HG3	1:G:28:PRO:O	2.11	0.51
1:A:132:GLU:N	1:G:98[B]:SER:OG	2.43	0.50
1:G:24:GLU:HB2	1:G:27:GLY:HA3	1.93	0.50
1:H:40:GLU:HG3	1:H:90[B]:ASP:C	2.35	0.50
1:J:69:ARG:HB3	1:J:78:GLU:OE2	2.11	0.50
1:K:108:GLY:O	1:K:111[A]:CYS:HB2	2.11	0.50
1:D:66:PRO:HD2	1:D:81:VAL:HG23	1.93	0.50
1:H:2:THR:CG2	1:H:106:LEU:HB2	2.42	0.50
1:D:22:GLN:NE2	1:D:105:SER:HA	2.26	0.50
1:H:67:LEU:HB2	1:H:69:ARG:HG2	1.93	0.50
1:D:72:GLY:CA	1:D:79:ARG:HB3	2.42	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:55:ALA:HB3	1:K:59:SER:OG	2.11	0.50
1:E:8:LEU:HD11	1:E:117:LEU:HD23	1.94	0.50
1:D:121:GLU:HB2	1:D:144:LEU:CD2	2.24	0.50
1:F:105:SER:O	1:F:111:CYS:HA	2.12	0.50
1:D:1:ALA:CB	1:D:113:ILE:HD11	2.42	0.50
1:L:66:PRO:HG2	1:L:81:VAL:HG21	1.93	0.50
1:J:32[B]:TRP:HA	1:J:97:VAL:O	2.11	0.50
1:E:121:GLU:O	1:E:121:GLU:HG2	2.10	0.49
1:G:73:GLY:CA	1:G:126:LEU:HD22	2.42	0.49
1:H:43:HIS:HA	1:H:122:LYS:O	2.11	0.49
1:B:112:ILE:HD12	1:B:149:ILE:HD13	1.93	0.49
1:G:66:PRO:HD2	1:G:81:VAL:HG21	1.91	0.49
1:J:38:LEU:CD2	1:J:144:LEU:HD12	2.42	0.49
1:L:121:GLU:HG3	1:L:122:LYS:HG3	1.93	0.49
1:J:15:GLN:HE21	1:J:15:GLN:HA	1.76	0.49
1:A:71:HIS:HB2	1:A:80:HIS:CE1	2.47	0.49
1:G:32[B]:TRP:HA	1:G:97:VAL:O	2.12	0.49
1:I:105:SER:O	1:I:111[B]:CYS:HA	2.12	0.49
1:I:105:SER:O	1:I:111[A]:CYS:HA	2.12	0.49
1:H:22:GLN:NE2	1:H:27:GLY:O	2.44	0.49
1:H:30:LYS:NZ	1:H:30:LYS:HB3	2.26	0.49
1:C:134:SER:HB2	1:C:139:ASN:ND2	2.28	0.49
1:I:7:VAL:HG13	1:I:7:VAL:O	2.12	0.49
1:B:108:GLY:O	1:B:111[A]:CYS:HB2	2.12	0.49
1:H:2:THR:O	1:H:106:LEU:HD12	2.13	0.49
1:D:3:LYS:NZ	6:D:306:HOH:O	2.45	0.49
1:I:18:ILE:HD13	1:I:33:GLY:HA3	1.93	0.49
1:A:48:HIS:CE1	1:A:63:HIS:HD2	2.31	0.49
1:G:66:PRO:HD2	1:G:81:VAL:HG23	1.94	0.49
1:E:80:HIS:HB2	1:E:83:ASP:OD1	2.13	0.49
1:C:5:VAL:HG22	1:C:6:CYS:H	1.78	0.49
1:I:117:LEU:O	1:I:146:CYS:HA	2.13	0.49
1:H:132:GLU:O	1:H:135:THR:OG1	2.30	0.49
1:K:108:GLY:O	1:K:111[B]:CYS:HB2	2.11	0.48
1:B:66:PRO:HD2	1:B:81:VAL:HB	1.96	0.48
1:G:99:ILE:HG22	1:G:100:GLU:H	1.78	0.48
1:J:66:PRO:HD2	1:J:81:VAL:HG23	1.95	0.48
1:F:134:SER:HB2	1:F:139:ASN:OD1	2.13	0.48
1:H:73:GLY:CA	1:H:126:LEU:HD22	2.40	0.48
1:B:81:VAL:HG22	1:B:103:VAL:HG12	1.94	0.48
1:H:134:SER:HB2	1:H:139:ASN:HD21	1.78	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:5:VAL:HG22	1:I:6:CYS:N	2.29	0.48
1:G:5:VAL:HG22	1:G:6:CYS:N	2.29	0.48
1:H:81:VAL:HA	1:H:103:VAL:HG11	1.96	0.48
1:A:26:ASN:O	1:A:26:ASN:ND2	2.47	0.48
1:E:71:HIS:HB2	1:E:80:HIS:HE1	1.79	0.48
1:D:65:ASN:HB2	1:D:80:HIS:CA	2.44	0.47
1:G:32[A]:TRP:HA	1:G:97:VAL:O	2.12	0.47
1:G:52:ASP:O	1:G:59:SER:HB2	2.14	0.47
1:H:38:LEU:HD22	1:H:43:HIS:CG	2.49	0.47
1:C:99:ILE:HD12	1:C:100:GLU:N	2.29	0.47
1:C:143:ARG:HG3	5:C:205:SO4:O1	2.15	0.47
1:K:65:ASN:ND2	1:K:68:SER:HA	2.29	0.47
1:K:75:LYS:HD3	1:E:128:LYS:NZ	2.30	0.47
1:E:26:ASN:OD1	1:E:26:ASN:N	2.31	0.47
1:G:5:VAL:CG1	1:G:152:ALA:HB2	2.43	0.47
1:E:109[A]:ASP:HB2	6:E:315:HOH:O	2.14	0.47
1:B:80:HIS:CB	1:B:83:ASP:OD2	2.62	0.47
1:K:7:VAL:HG21	1:L:53[B]:ASN:HB3	1.95	0.47
1:F:4:ALA:HB3	1:F:20:PHE:HB2	1.96	0.47
1:J:87:VAL:CG1	1:J:95:ALA:HB1	2.44	0.47
1:H:39:THR:O	1:H:43:HIS:NE2	2.34	0.47
1:F:108:GLY:O	1:F:111:CYS:HB2	2.14	0.47
1:E:30:LYS:HE3	1:E:32[B]:TRP:CD1	2.50	0.47
1:E:32[A]:TRP:HA	1:E:97:VAL:O	2.15	0.47
1:D:38:LEU:O	1:D:93:GLY:HA2	2.14	0.47
1:D:71:HIS:HB3	1:D:135:THR:C	2.40	0.47
1:I:106:LEU:HD22	1:I:113:ILE:HD11	1.96	0.47
1:L:65:ASN:HD21	1:L:68:SER:HA	1.77	0.47
1:L:90:ASP:HB2	6:L:312:HOH:O	2.13	0.47
1:H:8:LEU:HD12	1:H:8:LEU:N	2.29	0.47
1:H:81:VAL:HA	1:H:103:VAL:CG1	2.44	0.47
1:C:5:VAL:HG22	1:C:6:CYS:N	2.29	0.47
1:B:52:ASP:OD1	1:B:54:THR:HB	2.14	0.47
1:G:134:SER:HB2	1:G:139:ASN:ND2	2.29	0.47
1:L:8:LEU:O	1:L:15:GLN:HA	2.14	0.47
1:C:26:ASN:OD1	1:C:26:ASN:N	2.44	0.47
1:E:75:LYS:HB3	1:E:75:LYS:HE3	1.69	0.47
1:I:21:GLU:HB3	1:I:32[A]:TRP:CZ3	2.50	0.47
1:H:122:LYS:HB2	1:H:140:ALA:CA	2.45	0.47
1:J:14:VAL:HG11	1:J:144:LEU:CB	2.45	0.47
1:I:30:LYS:HG2	1:I:100:GLU:OE2	2.14	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:66:PRO:HD2	1:I:81:VAL:CG2	2.46	0.46
1:E:32[B]:TRP:HA	1:E:97:VAL:O	2.15	0.46
1:H:79:ARG:HH11	1:H:103:VAL:HG21	1.79	0.46
1:B:8:LEU:HD23	1:B:146:CYS:HA	1.97	0.46
1:I:52:ASP:OD1	1:I:53:ASN:N	2.48	0.46
1:H:21:GLU:O	1:H:29:VAL:HA	2.16	0.46
1:K:73:GLY:HA2	1:K:126:LEU:HD22	1.97	0.46
1:G:29:VAL:HG21	1:G:104:ILE:HG13	1.98	0.46
1:I:21:GLU:HB2	1:I:32[A]:TRP:HH2	1.81	0.46
1:A:39:THR:HG22	6:A:320:HOH:O	2.15	0.46
1:K:105:SER:O	1:K:111[A]:CYS:HA	2.16	0.46
1:K:105:SER:O	1:K:111[B]:CYS:HA	2.16	0.46
1:L:120:HIS:HB3	1:L:140:ALA:O	2.16	0.46
1:E:55:ALA:O	1:E:58:THR:HG23	2.16	0.46
1:I:121:GLU:HG3	1:I:122:LYS:N	2.31	0.46
1:K:15:GLN:OE1	1:K:36:LYS:HE2	2.16	0.46
1:B:5:VAL:HG22	1:B:6:CYS:N	2.29	0.46
1:F:80:HIS:HB2	1:F:83:ASP:CG	2.40	0.46
1:L:80:HIS:HB2	1:L:83:ASP:OD1	2.16	0.46
1:L:121:GLU:HA	1:L:144:LEU:HD11	1.98	0.46
1:H:40:GLU:HG3	1:H:90[A]:ASP:O	2.15	0.46
1:H:105:SER:HG	1:H:106:LEU:N	2.12	0.46
1:B:118:VAL:HG22	1:B:146:CYS:HB3	1.98	0.45
1:J:66:PRO:HD2	1:J:81:VAL:CG2	2.45	0.45
1:D:73:GLY:O	1:D:76:ASP:HB2	2.16	0.45
1:I:36:LYS:HA	1:I:93:GLY:O	2.17	0.45
1:J:38:LEU:HD11	1:J:119:VAL:HG21	1.98	0.45
1:D:1:ALA:HB3	1:D:113:ILE:HD11	1.98	0.45
1:D:119:VAL:O	1:D:119:VAL:HG13	2.16	0.45
1:I:80:HIS:CB	1:I:83:ASP:OD2	2.57	0.45
1:H:125[A]:ASP:OD1	1:H:134:SER:OG	2.35	0.45
1:D:22:GLN:HE22	1:D:105:SER:CB	2.29	0.45
1:J:105:SER:O	1:J:111:CYS:HA	2.16	0.45
1:A:50:PHE:CZ	1:B:153:GLN:HB2	2.51	0.45
1:D:1:ALA:CB	1:D:113:ILE:CD1	2.94	0.45
1:G:123:ALA:N	6:G:307:HOH:O	2.50	0.45
1:G:153:GLN:HB2	1:H:50:PHE:CZ	2.51	0.45
1:I:22:GLN:HB2	1:I:29:VAL:HG22	1.97	0.45
1:H:120:HIS:CB	1:H:140:ALA:HB1	2.46	0.45
1:D:140:ALA:N	6:D:302:HOH:O	2.06	0.45
1:G:76:ASP:O	1:G:79:ARG:HG2	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:5:VAL:HG22	1:H:6:CYS:N	2.32	0.45
1:A:3:LYS:HE3	1:A:3:LYS:HB2	1.76	0.45
1:H:41:GLY:O	1:H:88:THR:HA	2.17	0.45
1:G:118:VAL:HG11	1:G:143:ARG:HG2	1.98	0.45
1:H:79:ARG:HH21	1:H:83:ASP:N	2.15	0.45
1:D:80:HIS:CB	1:D:83:ASP:OD2	2.64	0.44
1:L:137:THR:HG22	1:L:137:THR:O	2.17	0.44
1:C:77:GLU:CD	1:B:2:THR:HG21	2.42	0.44
1:G:50:PHE:CE1	1:H:153:GLN:HB2	2.52	0.44
1:G:66:PRO:HD2	1:G:81:VAL:HB	1.99	0.44
1:I:69:ARG:HD2	1:I:78:GLU:OE1	2.16	0.44
1:L:55:ALA:HB3	1:L:59:SER:OG	2.17	0.44
1:C:80:HIS:HB2	1:C:83:ASP:OD1	2.18	0.44
1:K:8:LEU:HD21	1:K:117:LEU:HD23	2.00	0.44
1:F:3:LYS:NZ	1:F:153:GLN:O	2.48	0.44
1:E:21:GLU:HB2	1:E:32[A]:TRP:HH2	1.81	0.44
1:J:87:VAL:HG13	1:J:95:ALA:HB1	1.99	0.44
1:E:8:LEU:O	1:E:15:GLN:HA	2.17	0.44
1:E:81:VAL:HA	1:E:103:VAL:HG11	1.99	0.44
1:B:80:HIS:HB2	1:B:83:ASP:CG	2.42	0.44
1:C:70:LYS:NZ	6:C:305:HOH:O	2.51	0.44
1:K:55:ALA:N	1:K:59:SER:OG	2.49	0.44
1:G:55:ALA:HB3	1:G:59:SER:OG	2.17	0.44
1:J:118:VAL:HG22	1:J:146:CYS:HB3	2.00	0.44
1:C:71:HIS:HB2	1:C:80:HIS:CE1	2.53	0.44
1:I:66:PRO:HD2	1:I:81:VAL:CB	2.43	0.44
1:A:21:GLU:HB2	1:A:32[B]:TRP:CH2	2.53	0.44
1:D:22:GLN:HE22	1:D:105:SER:CA	2.31	0.44
1:L:71:HIS:HB2	1:L:80:HIS:CE1	2.53	0.44
1:I:116:THR:HG22	1:I:118:VAL:HG23	2.00	0.44
1:J:23:LYS:C	1:J:24:GLU:HG3	2.43	0.44
1:B:55:ALA:CB	1:B:59:SER:OG	2.62	0.43
1:L:126:LEU:HD11	6:L:326:HOH:O	2.17	0.43
1:A:131[A]:ASN:HB3	1:G:96[A]:ASP:O	2.17	0.43
1:C:121:GLU:HB2	1:C:144:LEU:HD21	2.00	0.43
1:G:66:PRO:HD2	1:G:81:VAL:CB	2.48	0.43
1:J:117:LEU:O	1:J:146:CYS:HA	2.18	0.43
1:A:98[B]:SER:HB3	1:I:132:GLU:N	2.32	0.43
1:J:71:HIS:CE1	1:J:83:ASP:OD1	2.71	0.43
1:J:74:PRO:O	1:J:79:ARG:NH1	2.51	0.43
1:H:21:GLU:HB2	1:H:32[A]:TRP:CH2	2.54	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:144:LEU:N	1:H:144:LEU:HD23	2.33	0.43
1:L:22:GLN:HG2	1:L:24:GLU:O	2.19	0.43
1:L:45:PHE:O	1:L:84:LEU:HB2	2.18	0.43
1:J:8:LEU:O	1:J:15:GLN:HA	2.17	0.43
1:E:26:ASN:HD21	1:B:16:GLY:HA2	1.83	0.43
1:J:9:LYS:HA	1:J:15:GLN:HE22	1.83	0.43
1:J:36:LYS:N	6:J:301:HOH:O	2.51	0.43
1:H:3:LYS:HB2	1:H:3:LYS:HE3	1.82	0.43
1:H:21:GLU:HB2	1:H:32[A]:TRP:HH2	1.83	0.43
1:A:98[A]:SER:HB3	1:I:132:GLU:N	2.32	0.43
1:A:21:GLU:HB2	1:A:32[B]:TRP:HZ3	1.83	0.43
1:A:101:ASP:HB3	1:A:104:ILE:HG12	1.99	0.43
1:K:51:GLY:N	1:L:151:ILE:O	2.52	0.43
1:D:47:VAL:O	1:D:64:PHE:HB3	2.18	0.43
1:D:131[A]:ASN:HB2	6:D:304:HOH:O	2.18	0.43
1:A:73:GLY:CA	1:A:126:LEU:HD22	2.49	0.43
1:F:5:VAL:CG1	1:F:6:CYS:N	2.82	0.43
1:I:14:VAL:HG12	1:I:145:ALA:HB2	1.99	0.43
1:I:121:GLU:HG3	1:I:141:GLY:HA3	2.00	0.43
1:C:8:LEU:O	1:C:15:GLN:HA	2.19	0.43
1:F:71:HIS:HB2	1:F:80:HIS:CE1	2.54	0.43
1:F:76:ASP:O	1:F:79:ARG:HG2	2.19	0.43
1:F:120:HIS:HE1	6:F:319[A]:HOH:O	2.01	0.43
1:D:46:HIS:HA	1:D:82:GLY:O	2.19	0.43
1:H:79:ARG:NH1	1:H:103:VAL:HG21	2.34	0.43
1:H:71:HIS:HE2	1:H:124:ASP:CG	2.27	0.42
1:A:125:ASP:OD1	1:A:138:GLY:HA2	2.19	0.42
1:C:62:PRO:HD3	4:C:201:ACT:CH3	2.34	0.42
1:G:38:LEU:O	1:G:93:GLY:HA2	2.19	0.42
1:H:7:VAL:CG1	1:H:9:LYS:HE3	2.49	0.42
1:H:99:ILE:HG12	1:H:100:GLU:N	2.34	0.42
1:F:80:HIS:HB2	1:F:83:ASP:OD1	2.18	0.42
1:D:79:ARG:NH2	1:D:103:VAL:HG21	2.34	0.42
1:J:80:HIS:HB2	1:J:83:ASP:OD2	2.19	0.42
1:H:55:ALA:O	1:H:58:THR:HG23	2.19	0.42
1:K:153:GLN:HB2	1:L:50:PHE:CZ	2.55	0.42
1:F:5:VAL:HG22	1:F:19:ASN:OD1	2.20	0.42
1:J:108:GLY:O	1:J:111:CYS:HB2	2.20	0.42
1:H:71:HIS:HB2	1:H:80:HIS:CE1	2.54	0.42
1:K:71:HIS:HB2	1:K:80:HIS:CE1	2.55	0.42
1:E:81:VAL:HG13	1:E:103:VAL:HG12	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:99:ILE:CG2	1:L:100:GLU:N	2.83	0.42
1:K:116:THR:CG2	1:K:146:CYS:HB2	2.50	0.42
1:I:121:GLU:HG3	1:I:122:LYS:H	1.84	0.42
1:G:64:PHE:CZ	1:G:66:PRO:HG3	2.54	0.42
1:I:113:ILE:N	1:I:113:ILE:HD12	2.34	0.42
1:I:121:GLU:CG	1:I:141:GLY:HA3	2.48	0.42
1:L:109:ASP:OD1	1:L:109:ASP:N	2.52	0.42
1:L:131:ASN:OD1	1:L:134:SER:N	2.48	0.42
1:L:121:GLU:HA	1:L:144:LEU:CD1	2.50	0.42
1:H:38:LEU:O	1:H:93:GLY:HA2	2.19	0.42
1:A:22:GLN:OE1	1:A:106:LEU:N	2.45	0.42
1:K:48:HIS:O	1:K:115:ARG:HB3	2.20	0.42
1:F:151:ILE:O	1:E:51:GLY:N	2.47	0.42
1:E:118:VAL:HG22	1:E:146:CYS:HB3	2.02	0.42
1:D:44:GLY:O	1:D:119:VAL:HA	2.20	0.42
1:G:65:ASN:ND2	1:G:68:SER:HA	2.34	0.42
1:I:124:ASP:O	1:I:126:LEU:HD22	2.20	0.42
1:L:52:ASP:OD1	1:L:54:THR:HB	2.20	0.42
1:A:22:GLN:HG2	1:A:24:GLU:O	2.20	0.42
1:K:18:ILE:HD13	1:K:18:ILE:HA	1.91	0.42
1:K:30:LYS:HB3	1:K:30:LYS:NZ	2.35	0.42
1:I:21:GLU:HB3	1:I:32[A]:TRP:HZ3	1.85	0.42
1:H:2:THR:C	1:H:106:LEU:HD12	2.45	0.42
1:H:67:LEU:HG	1:H:81:VAL:CG2	2.50	0.42
1:H:125[A]:ASP:OD1	1:H:127:GLY:N	2.53	0.42
1:F:74:PRO:HA	1:F:79:ARG:NH1	2.34	0.41
1:D:79:ARG:CD	1:D:80:HIS:O	2.68	0.41
1:D:80:HIS:ND1	1:D:83:ASP:OD2	2.53	0.41
1:G:3:LYS:HG2	1:G:21:GLU:OE1	2.20	0.41
1:G:48:HIS:CE1	1:G:63:HIS:HD2	2.38	0.41
1:A:51:GLY:HA2	1:A:116:THR:OG1	2.20	0.41
1:J:125:ASP:OD1	1:J:134:SER:OG	2.36	0.41
1:A:96:ASP:O	1:I:131:ASN:HB3	2.20	0.41
1:D:80:HIS:HB2	1:D:83:ASP:OD2	2.20	0.41
1:D:2:THR:HG22	1:D:3:LYS:HG3	2.02	0.41
1:G:112:ILE:HA	1:G:115:ARG:HD2	2.03	0.41
1:I:119:VAL:CG1	1:I:145:ALA:HB3	2.51	0.41
1:L:51:GLY:HA3	1:L:114:GLY:O	2.20	0.41
1:H:49:GLU:HB2	1:H:62:PRO:O	2.21	0.41
1:H:120:HIS:HB2	1:H:140:ALA:HB1	2.02	0.41
1:A:18:ILE:HD11	1:A:97:VAL:HG21	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:81:VAL:HG22	1:E:103:VAL:CG1	2.50	0.41
1:J:74:PRO:HA	1:J:79:ARG:NH1	2.35	0.41
1:H:118:VAL:HG11	1:H:143:ARG:CG	2.50	0.41
1:K:75:LYS:HD3	1:E:128:LYS:HZ3	1.85	0.41
1:J:80:HIS:CG	1:J:83:ASP:OD2	2.73	0.41
1:H:46:HIS:HA	1:H:82:GLY:O	2.21	0.41
1:K:66:PRO:HD2	1:K:81:VAL:CG2	2.51	0.41
1:K:96[B]:ASP:O	1:E:131:ASN:HB2	2.20	0.41
1:F:75:LYS:H	1:F:75:LYS:HG2	1.69	0.41
1:E:21:GLU:HB2	1:E:32[A]:TRP:CH2	2.56	0.41
1:B:19:ASN:HB2	1:B:32[B]:TRP:CD1	2.56	0.41
1:B:74:PRO:O	1:B:79:ARG:NH1	2.53	0.41
1:I:8:LEU:HD12	1:I:8:LEU:N	2.35	0.41
1:I:14:VAL:HG12	1:I:145:ALA:CB	2.51	0.41
1:H:8:LEU:N	1:H:8:LEU:CD1	2.84	0.41
1:A:21:GLU:HG3	1:A:32[B]:TRP:HH2	1.86	0.41
1:C:66:PRO:HD2	1:C:81:VAL:HB	2.03	0.41
1:D:22:GLN:CD	1:D:106:LEU:H	2.29	0.41
1:L:121:GLU:HG3	1:L:122:LYS:N	2.35	0.41
1:D:134:SER:O	1:D:138:GLY:HA2	2.20	0.40
1:I:91:LYS:O	1:I:91:LYS:HD3	2.21	0.40
1:C:55:ALA:O	1:C:58:THR:HG23	2.21	0.40
1:E:51:GLY:HA3	1:E:114:GLY:O	2.20	0.40
1:D:15:GLN:HE21	1:D:36:LYS:CE	2.34	0.40
1:D:70:LYS:HA	6:D:303:HOH:O	2.21	0.40
1:J:49:GLU:O	1:J:115:ARG:NH1	2.51	0.40
1:A:5:VAL:HG22	1:A:6:CYS:N	2.36	0.40
1:L:48:HIS:CE1	1:L:63:HIS:HD2	2.40	0.40
1:A:5:VAL:HG22	1:A:6:CYS:H	1.87	0.40
1:F:67:LEU:HD13	1:F:69:ARG:NH1	2.36	0.40
1:F:119:VAL:HG12	1:F:145:ALA:O	2.22	0.40
1:D:121:GLU:CA	1:D:144:LEU:HD21	2.51	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:132:GLU:N	1:H:98[B]:SER:OG[3_555]	2.10	0.10

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	157/156 (101%)	152 (97%)	5 (3%)	0	100	100
1	B	161/156 (103%)	159 (99%)	2 (1%)	0	100	100
1	C	156/156 (100%)	150 (96%)	6 (4%)	0	100	100
1	D	160/156 (103%)	151 (94%)	9 (6%)	0	100	100
1	E	160/156 (103%)	157 (98%)	3 (2%)	0	100	100
1	F	157/156 (101%)	153 (98%)	4 (2%)	0	100	100
1	G	155/156 (99%)	150 (97%)	5 (3%)	0	100	100
1	H	159/156 (102%)	151 (95%)	8 (5%)	0	100	100
1	I	157/156 (101%)	153 (98%)	4 (2%)	0	100	100
1	J	154/156 (99%)	149 (97%)	5 (3%)	0	100	100
1	K	157/156 (101%)	155 (99%)	2 (1%)	0	100	100
1	L	158/156 (101%)	151 (96%)	6 (4%)	1 (1%)	21	22
All	All	1891/1872 (101%)	1831 (97%)	59 (3%)	1 (0%)	48	57

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	L	54	THR

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	124/121 (102%)	109 (88%)	15 (12%)	5	4
1	B	128/121 (106%)	120 (94%)	8 (6%)	16	19
1	C	123/121 (102%)	114 (93%)	9 (7%)	13	14
1	D	127/121 (105%)	119 (94%)	8 (6%)	16	19
1	E	127/121 (105%)	121 (95%)	6 (5%)	23	30
1	F	124/121 (102%)	119 (96%)	5 (4%)	28	36
1	G	123/121 (102%)	116 (94%)	7 (6%)	18	23
1	H	126/121 (104%)	115 (91%)	11 (9%)	9	10
1	I	124/121 (102%)	122 (98%)	2 (2%)	55	67
1	J	121/121 (100%)	109 (90%)	12 (10%)	7	7
1	K	125/121 (103%)	118 (94%)	7 (6%)	19	23
1	L	125/121 (103%)	118 (94%)	7 (6%)	19	23
All	All	1497/1452 (103%)	1400 (94%)	97 (6%)	16	17

All (97) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	LYS
1	A	17	ILE
1	A	21	GLU
1	A	24	GLU
1	A	26	ASN
1	A	38	LEU
1	A	42[A]	LEU
1	A	42[B]	LEU
1	A	46	HIS
1	A	67	LEU
1	A	75	LYS
1	A	92	ASP
1	A	111[A]	CYS
1	A	111[B]	CYS
1	A	113	ILE
1	C	2	THR
1	C	17	ILE
1	C	30	LYS
1	C	54	THR
1	C	58	THR
1	C	77	GLU
1	C	91	LYS

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Mol	Chain	Res	Type
1	C	99	ILE
1	C	109	ASP
1	K	34	SER
1	K	42[A]	LEU
1	K	42[B]	LEU
1	K	54	THR
1	K	91	LYS
1	K	113	ILE
1	K	121	GLU
1	F	3	LYS
1	F	8	LEU
1	F	99	ILE
1	F	109	ASP
1	F	136	LYS
1	E	26	ASN
1	E	42[A]	LEU
1	E	42[B]	LEU
1	E	46	HIS
1	E	58	THR
1	E	59	SER
1	D	2	THR
1	D	25	SER
1	D	26	ASN
1	D	46	HIS
1	D	58	THR
1	D	65	ASN
1	D	98	SER
1	D	144	LEU
1	B	42[A]	LEU
1	B	42[B]	LEU
1	B	54	THR
1	B	58	THR
1	B	81	VAL
1	B	92	ASP
1	B	99	ILE
1	B	109	ASP
1	G	2	THR
1	G	34	SER
1	G	42[A]	LEU
1	G	42[B]	LEU
1	G	84	LEU
1	G	100	GLU

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Mol	Chain	Res	Type
1	G	135	THR
1	I	91	LYS
1	I	99	ILE
1	L	24	GLU
1	L	42[A]	LEU
1	L	42[B]	LEU
1	L	67	LEU
1	L	92	ASP
1	L	98	SER
1	L	136	LYS
1	J	2	THR
1	J	7	VAL
1	J	14	VAL
1	J	15	GLN
1	J	24	GLU
1	J	35	ILE
1	J	54	THR
1	J	58	THR
1	J	59	SER
1	J	88	THR
1	J	99	ILE
1	J	119	VAL
1	H	2	THR
1	H	11	ASP
1	H	15	GLN
1	H	54	THR
1	H	58	THR
1	H	75	LYS
1	H	92	ASP
1	H	99	ILE
1	H	100	GLU
1	H	116	THR
1	H	131	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	43	HIS
1	A	53	ASN
1	C	15	GLN
1	K	19	ASN
1	F	15	GLN

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Mol	Chain	Res	Type
1	E	15	GLN
1	D	15	GLN
1	B	15	GLN
1	B	19	ASN
1	B	153	GLN
1	G	153	GLN
1	I	153	GLN
1	L	15	GLN
1	J	15	GLN
1	J	19	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 33 ligands modelled in this entry, 25 are monoatomic - leaving 8 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	$\# Z > 2$
4	ACT	C	204	-	3,3,3	1.38	0	3,3,3	1.41	0
4	ACT	B	203	-	3,3,3	1.40	1 (33%)	3,3,3	1.34	0
5	SO4	A	205	-	4,4,4	0.30	0	6,6,6	0.18	0
5	SO4	C	205	-	4,4,4	0.23	0	6,6,6	0.12	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	SO4	I	203	-	4,4,4	0.26	0	6,6,6	0.11	0
4	ACT	A	204	-	3,3,3	1.48	1 (33%)	3,3,3	1.38	0
4	ACT	C	201	-	3,3,3	1.18	0	3,3,3	1.58	1 (33%)
5	SO4	E	203	-	4,4,4	0.24	0	6,6,6	0.13	0

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	203	ACT	CH3-C	2.04	1.57	1.49
4	A	204	ACT	CH3-C	2.02	1.57	1.49

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	C	201	ACT	OXT-C-O	2.02	129.53	122.03

There are no chirality outliers.

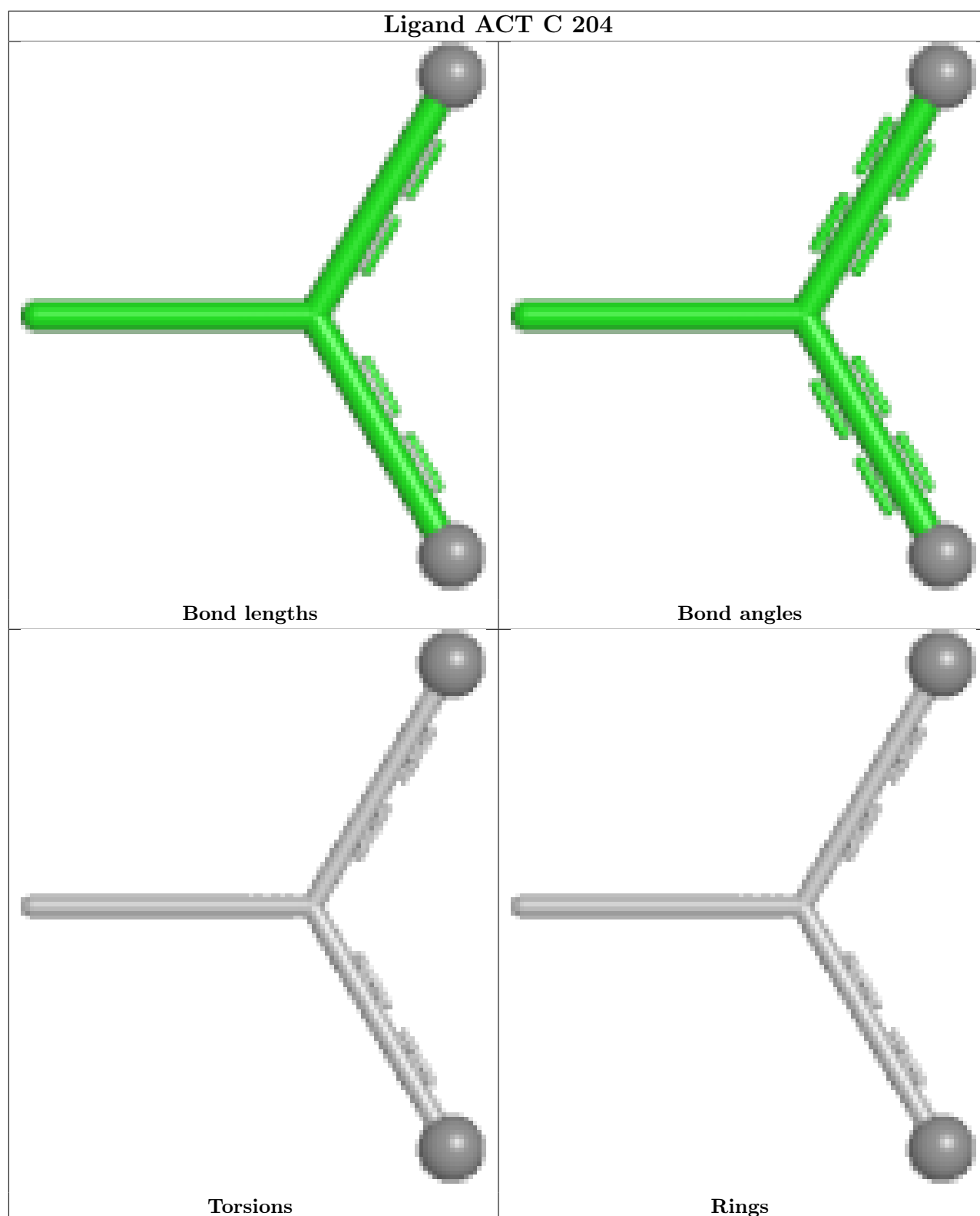
There are no torsion outliers.

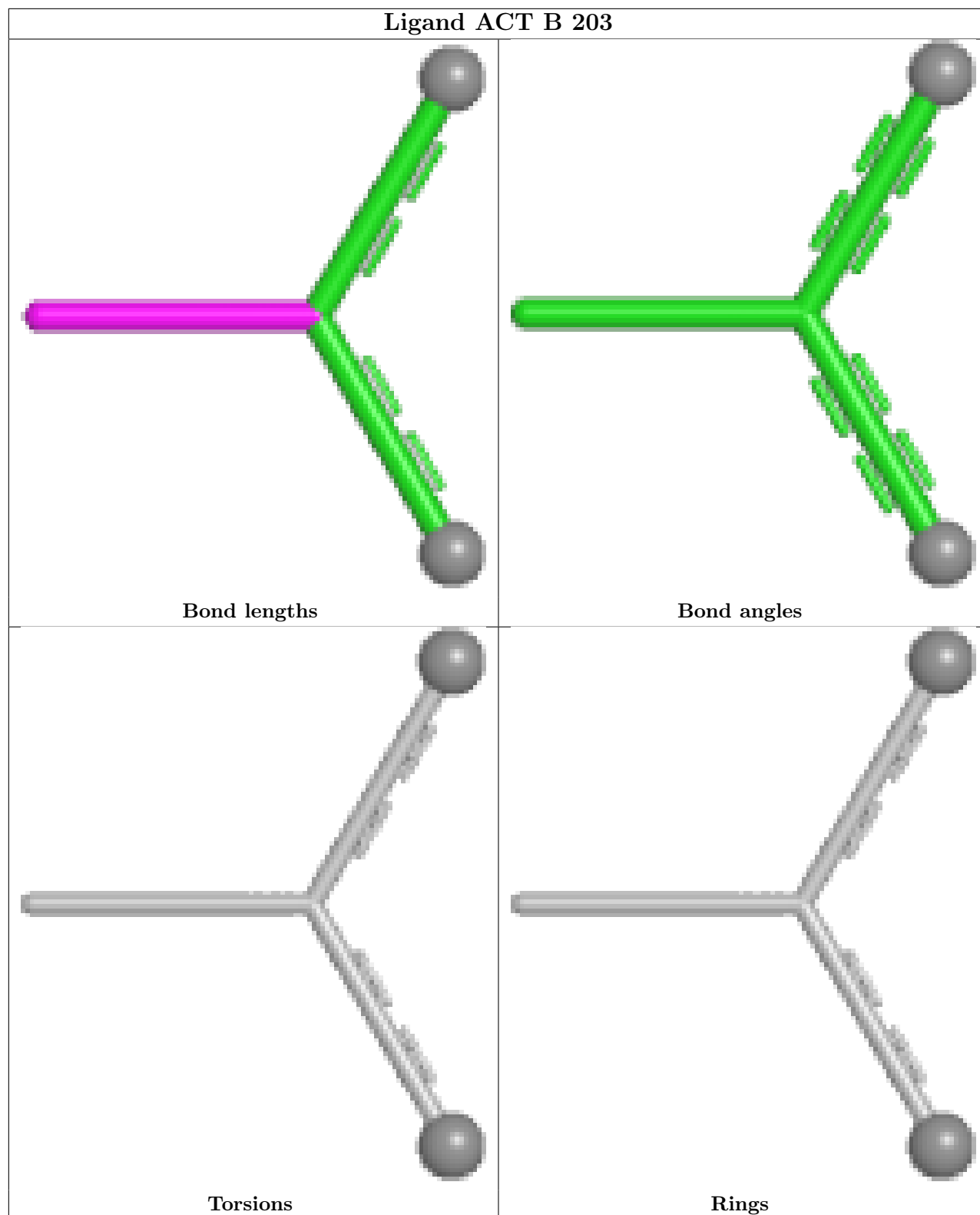
There are no ring outliers.

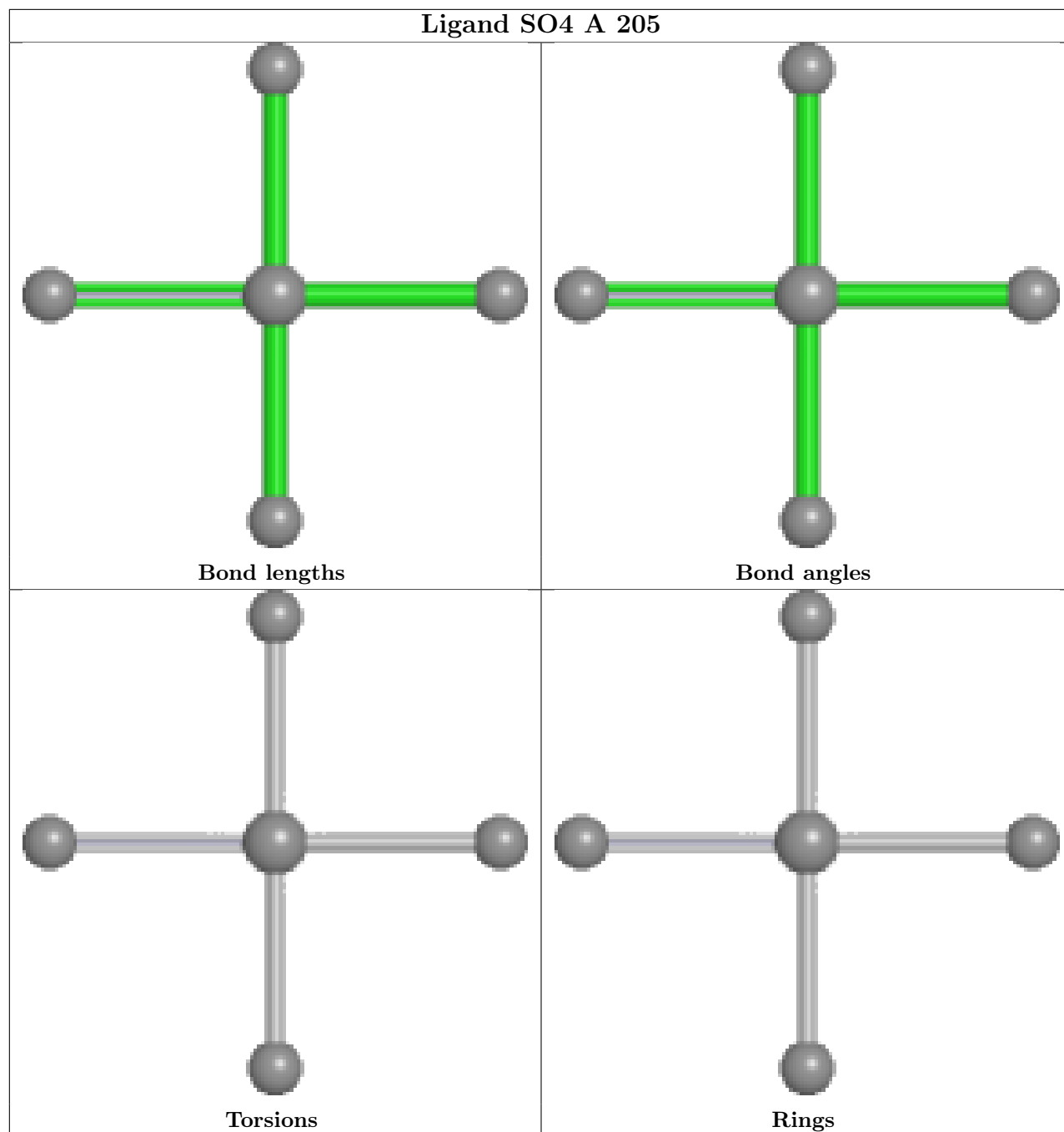
3 monomers are involved in 8 short contacts:

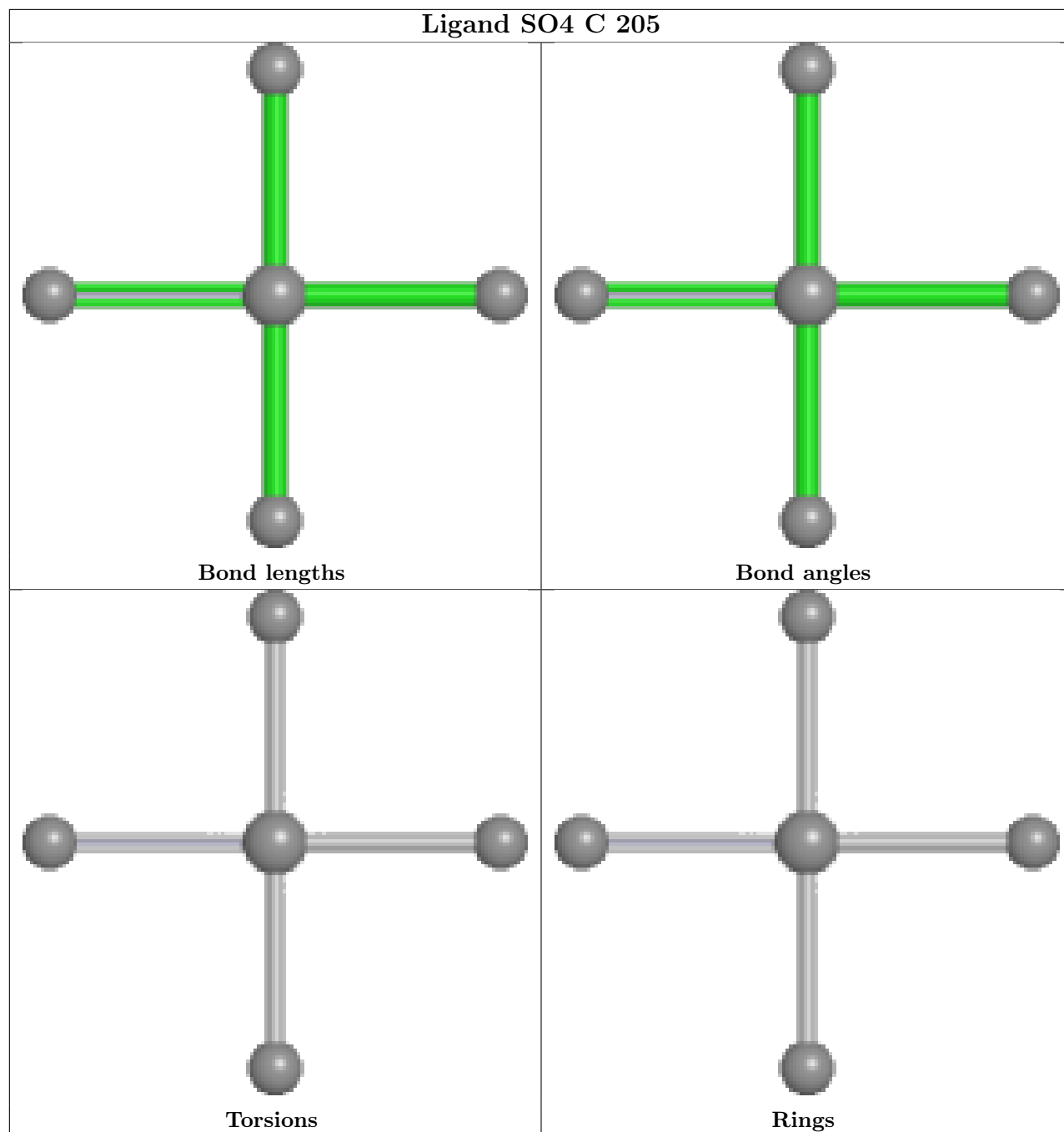
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	C	204	ACT	2	0
5	C	205	SO4	2	0
4	C	201	ACT	4	0

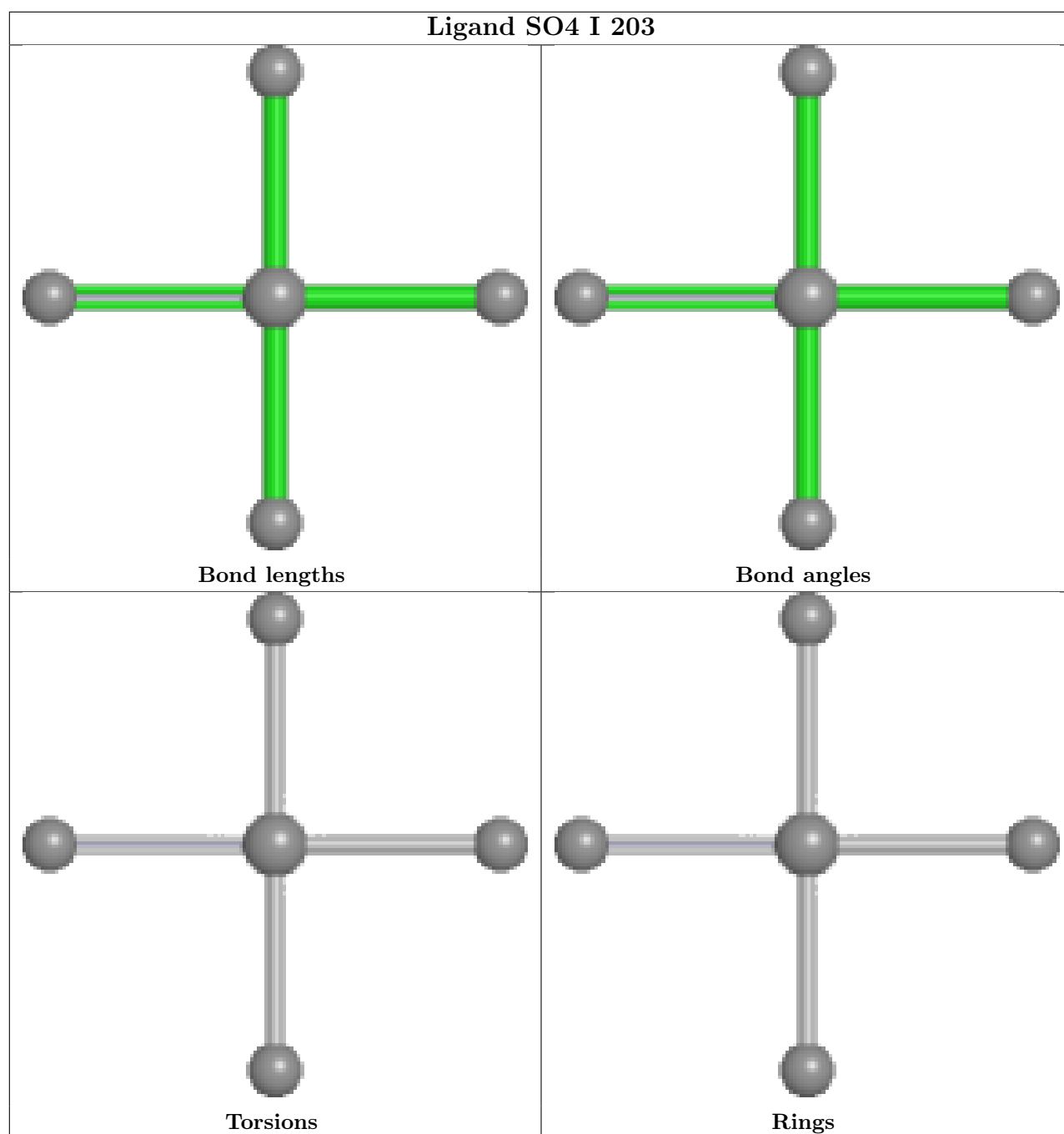
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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.

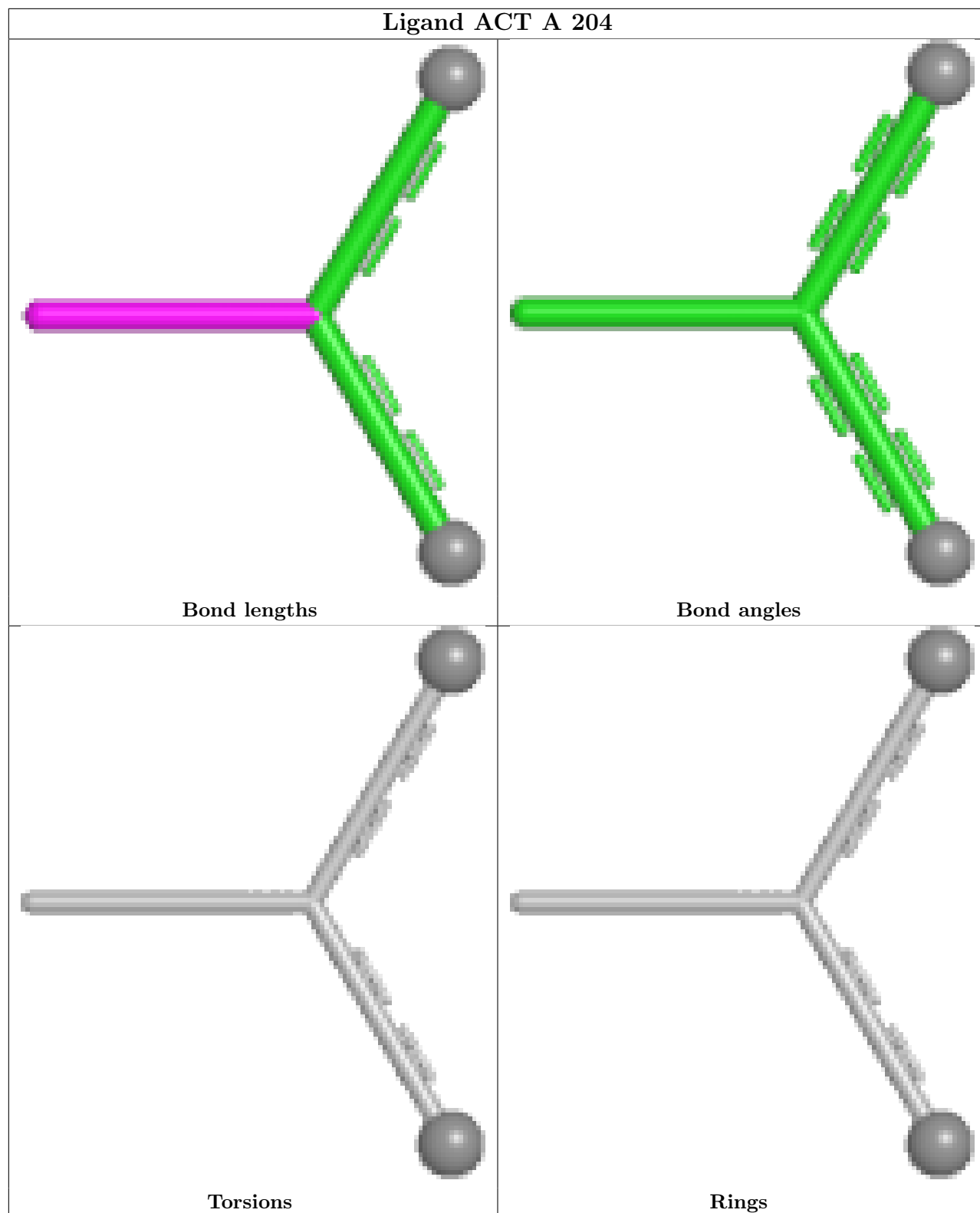


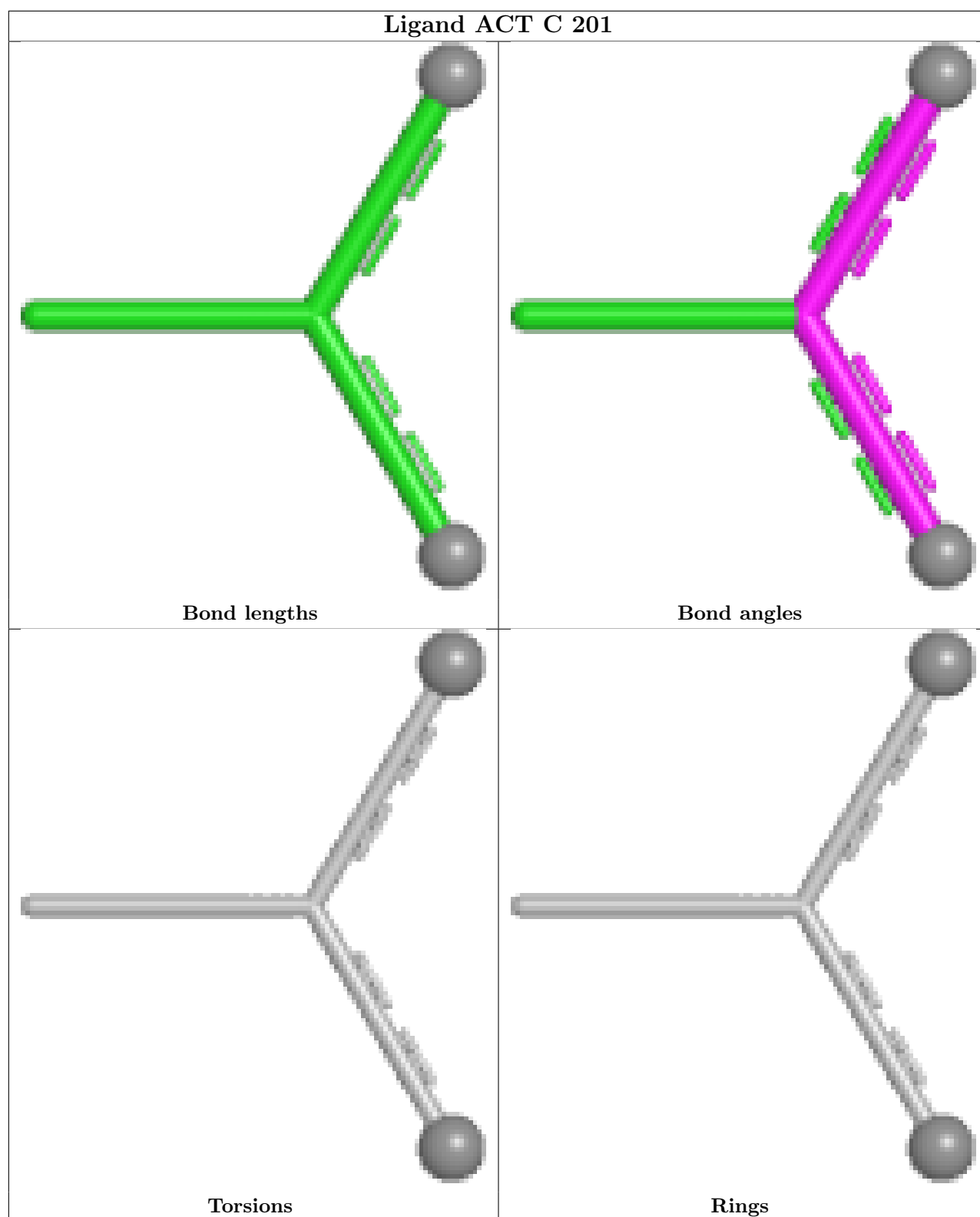


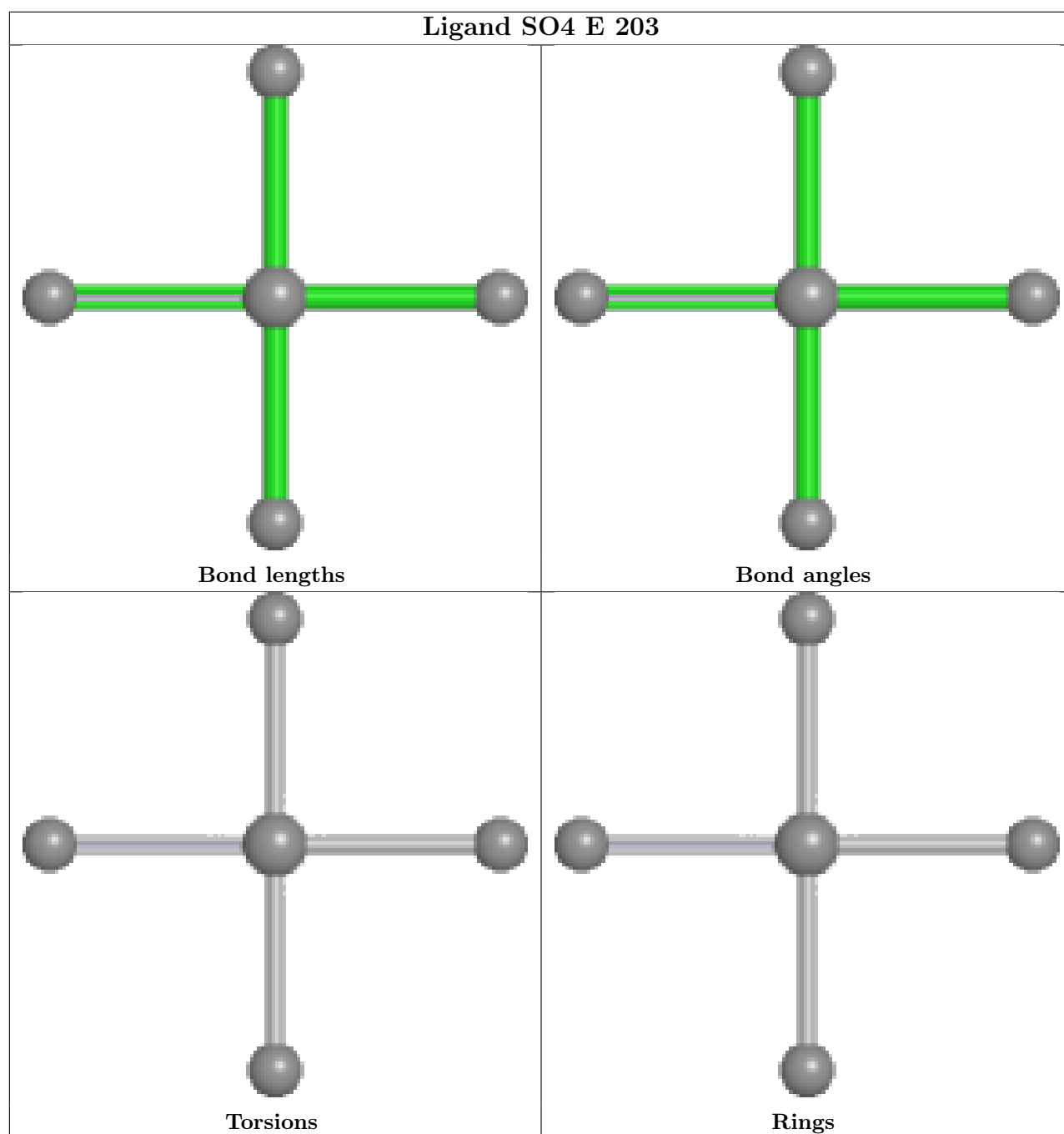












5.7 Other polymers [i](#)

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	153/156 (98%)	-0.19	0	100	100	24, 49, 82, 134	7 (4%)
1	B	156/156 (100%)	-0.14	0	100	100	22, 53, 76, 113	11 (7%)
1	C	153/156 (98%)	-0.14	0	100	100	28, 48, 89, 119	8 (5%)
1	D	156/156 (100%)	0.71	5 (3%)	50	57	30, 86, 120, 165	7 (4%)
1	E	153/156 (98%)	-0.10	2 (1%)	75	79	27, 54, 83, 116	17 (11%)
1	F	153/156 (98%)	0.10	0	100	100	28, 64, 99, 124	6 (3%)
1	G	152/156 (97%)	0.48	6 (3%)	43	49	30, 72, 115, 156	8 (5%)
1	H	153/156 (98%)	0.84	11 (7%)	21	25	30, 89, 152, 188	12 (7%)
1	I	153/156 (98%)	0.39	1 (0%)	84	87	39, 77, 118, 159	8 (5%)
1	J	153/156 (98%)	0.47	4 (2%)	57	63	26, 81, 127, 153	4 (2%)
1	K	152/156 (97%)	-0.22	0	100	100	23, 49, 81, 126	8 (5%)
1	L	153/156 (98%)	0.03	1 (0%)	84	87	28, 56, 86, 113	11 (7%)
All	All	1840/1872 (98%)	0.19	30 (1%)	70	75	22, 64, 113, 188	107 (5%)

All (30) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	G	28	PRO	5.7
1	H	67	LEU	4.5
1	H	130	GLY	4.4
1	D	130	GLY	4.2
1	I	1	ALA	3.4
1	H	28	PRO	3.4
1	H	74	PRO	3.2
1	E	53[A]	ASN	2.8
1	G	152	ALA	2.8
1	J	87	VAL	2.7
1	G	26	ASN	2.6

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Mol	Chain	Res	Type	RSRZ
1	D	128	LYS	2.6
1	D	64	PHE	2.6
1	H	66	PRO	2.6
1	J	37	GLY	2.5
1	D	134	SER	2.4
1	J	1	ALA	2.4
1	H	50	PHE	2.3
1	G	130	GLY	2.3
1	H	134	SER	2.3
1	H	126	LEU	2.2
1	H	137	THR	2.2
1	H	75	LYS	2.2
1	G	4	ALA	2.1
1	L	1	ALA	2.1
1	J	88	THR	2.1
1	D	104	ILE	2.1
1	E	70[A]	LYS	2.0
1	H	110	HIS	2.0
1	G	30	LYS	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(Å ²)	Q<0.9
5	SO4	I	203	5/5	0.74	0.11	87,99,112,117	0
4	ACT	A	204	4/4	0.76	0.13	33,41,49,49	7
4	ACT	C	204	4/4	0.80	0.13	62,72,80,80	0
5	SO4	E	203	5/5	0.81	0.11	92,93,103,114	0

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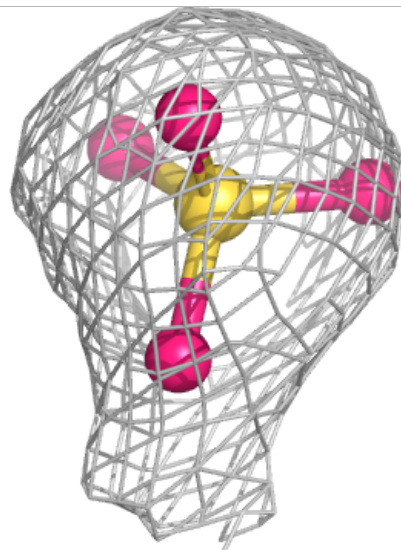
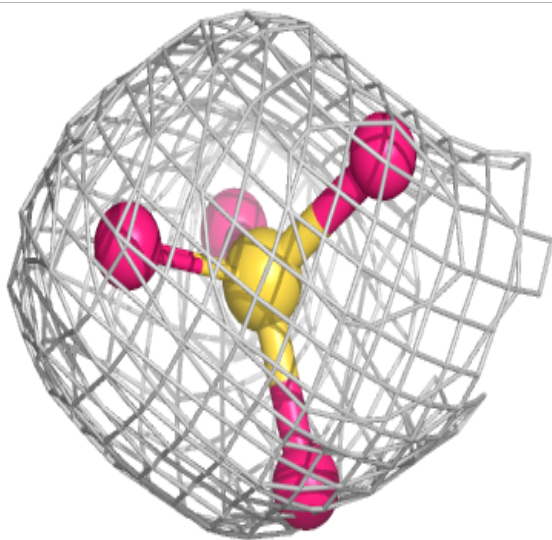
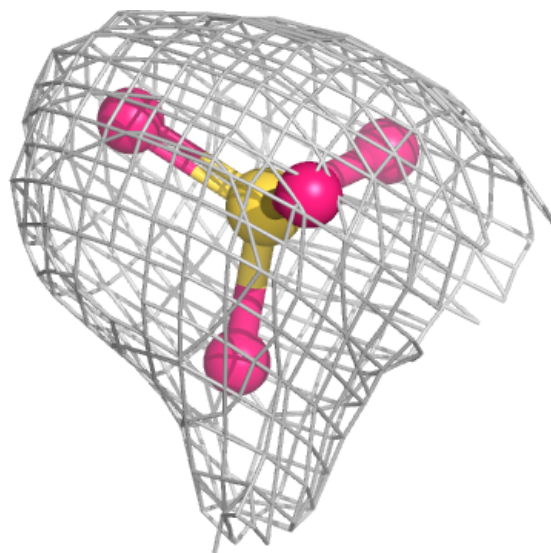
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	ACT	C	201	4/4	0.82	0.16	47,65,74,75	7
4	ACT	B	203	4/4	0.84	0.15	52,64,77,77	7
3	ZN	H	202	1/1	0.91	0.08	120,120,120,120	0
5	SO4	A	205	5/5	0.92	0.10	48,50,60,86	3
5	SO4	C	205	5/5	0.92	0.09	39,43,50,56	5
2	CU	H	201	1/1	0.93	0.08	90,90,90,90	0
2	CU	D	201	1/1	0.95	0.07	85,85,85,85	0
2	CU	G	201	1/1	0.97	0.04	60,60,60,60	0
3	ZN	D	202	1/1	0.97	0.05	84,84,84,84	0
3	ZN	G	202	1/1	0.97	0.05	67,67,67,67	0
3	ZN	J	202	1/1	0.98	0.03	58,58,58,58	0
2	CU	J	201	1/1	0.98	0.04	60,60,60,60	0
2	CU	I	201	1/1	0.98	0.04	56,56,56,56	0
3	ZN	I	202	1/1	0.98	0.04	59,59,59,59	0
3	ZN	L	202	1/1	0.98	0.04	46,46,46,46	0
2	CU	B	201	1/1	0.99	0.03	50,50,50,50	0
2	CU	C	202	1/1	0.99	0.04	43,43,43,43	0
2	CU	K	201	1/1	0.99	0.02	45,45,45,45	0
2	CU	L	201	1/1	0.99	0.03	46,46,46,46	0
2	CU	F	201	1/1	0.99	0.03	55,55,55,55	0
2	CU	E	201	1/1	0.99	0.02	52,52,52,52	0
3	ZN	A	202	1/1	0.99	0.04	32,32,32,32	0
3	ZN	A	203	1/1	0.99	0.03	59,59,59,59	0
3	ZN	K	202	1/1	0.99	0.02	44,44,44,44	0
3	ZN	F	202	1/1	0.99	0.03	52,52,52,52	0
3	ZN	E	202	1/1	0.99	0.03	53,53,53,53	0
2	CU	A	201	1/1	0.99	0.02	42,42,42,42	0
3	ZN	B	202	1/1	0.99	0.02	49,49,49,49	0
3	ZN	C	203	1/1	1.00	0.02	29,29,29,29	0

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

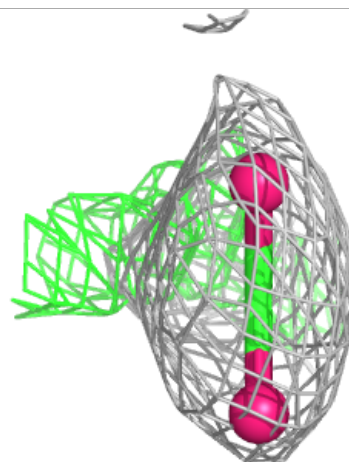
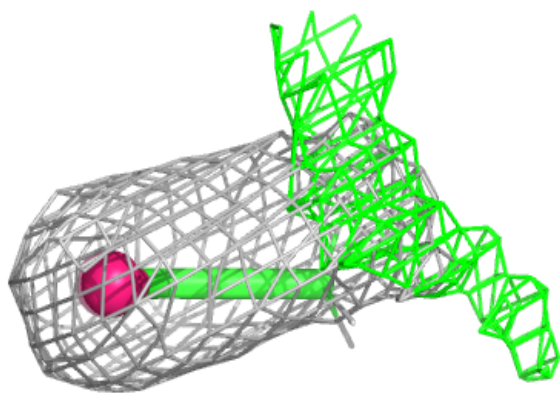
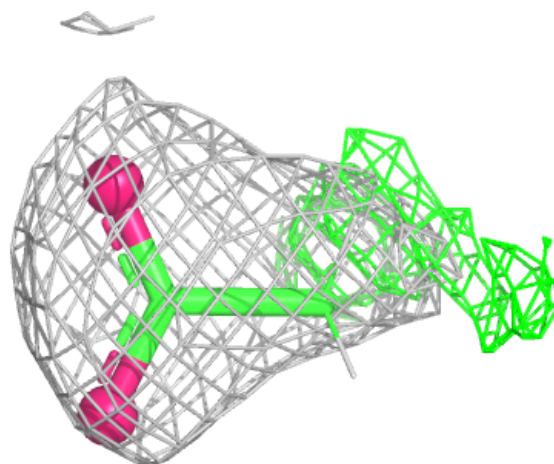
Electron density around SO4 I 203:

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



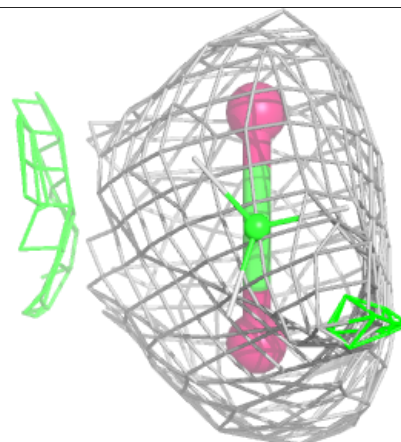
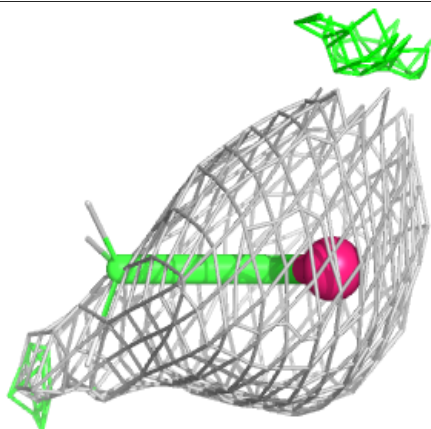
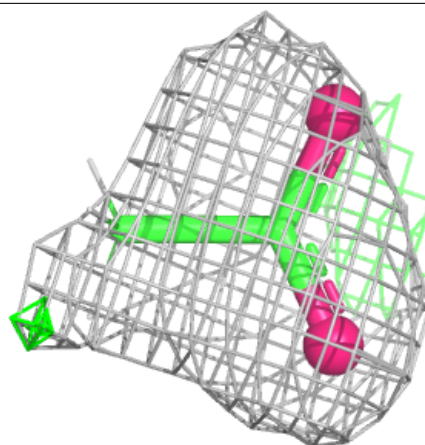
Electron density around ACT A 204:

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



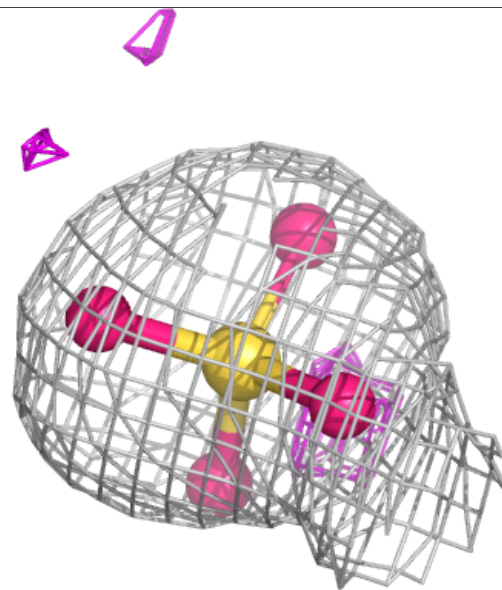
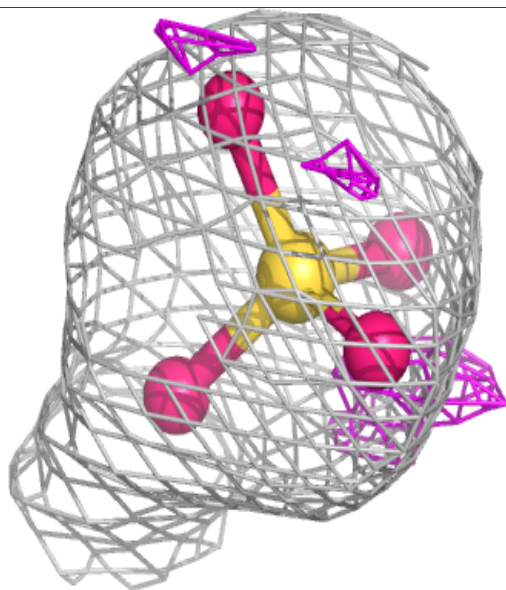
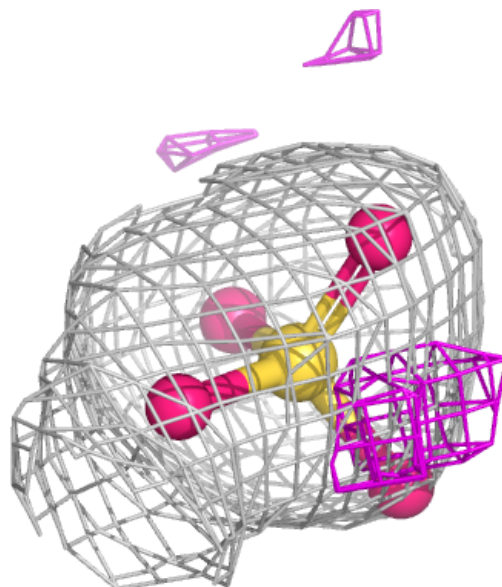
Electron density around ACT C 204:

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



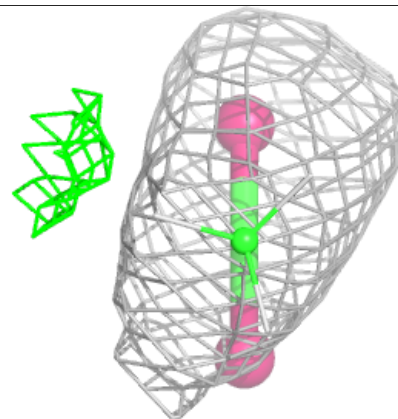
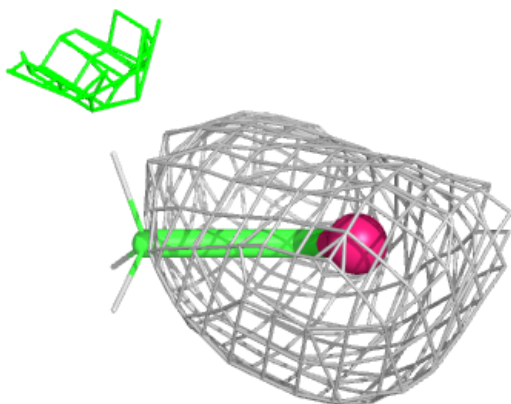
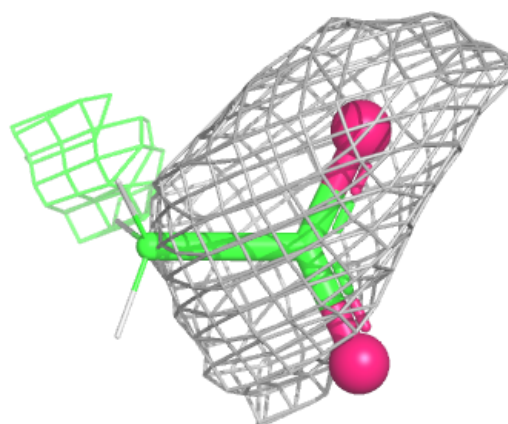
Electron density around SO4 E 203:

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

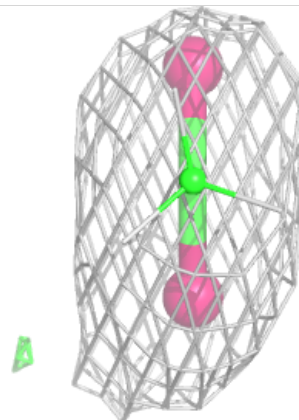
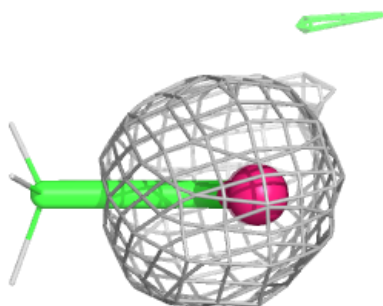
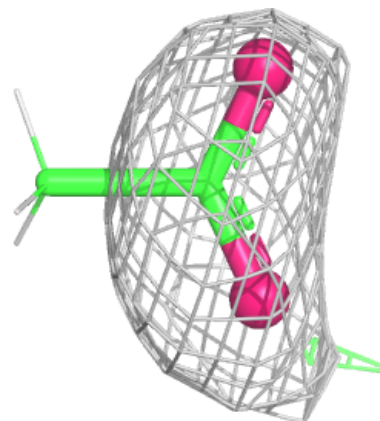


Electron density around ACT C 201:

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

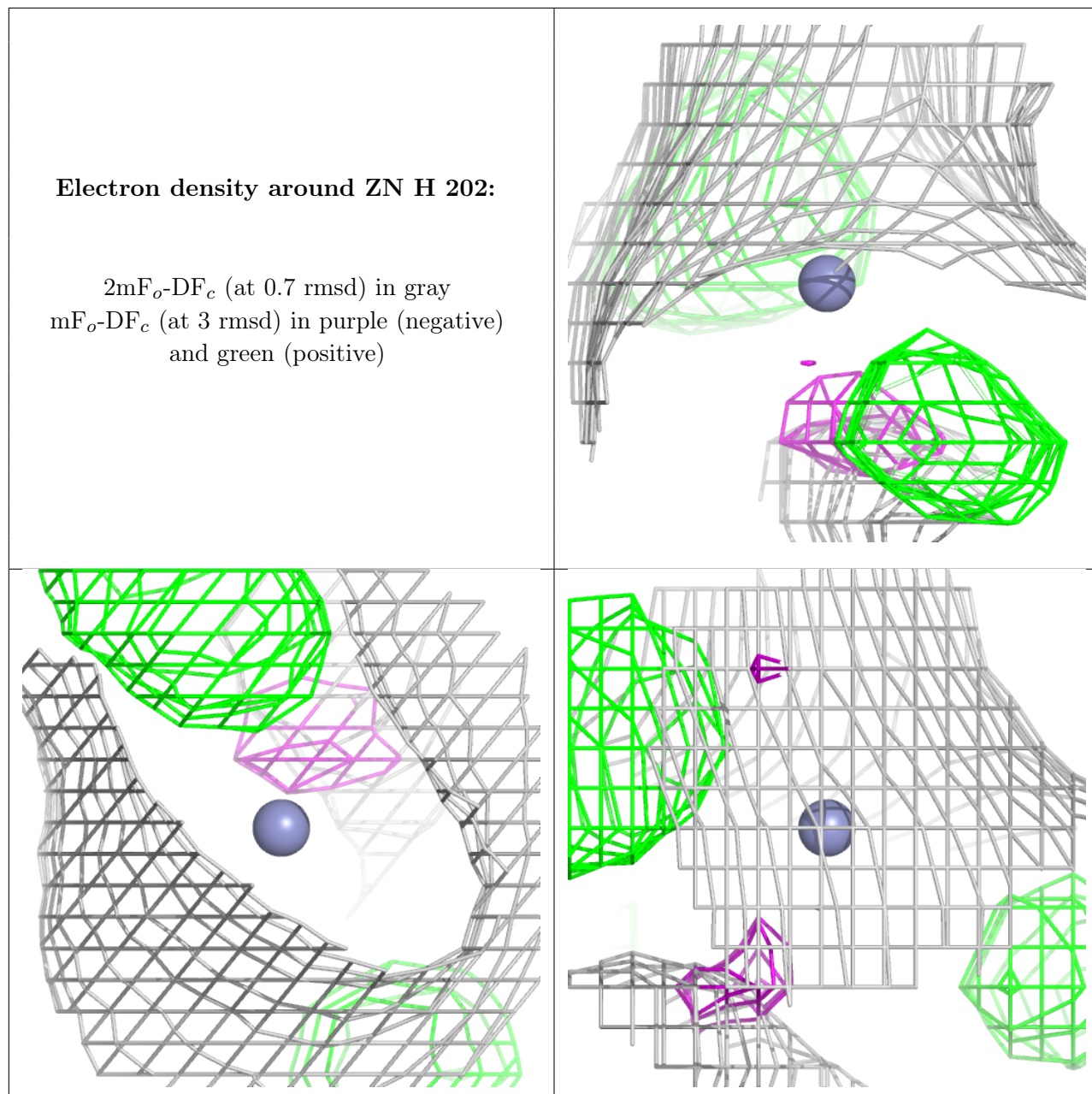
**Electron density around ACT B 203:**

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



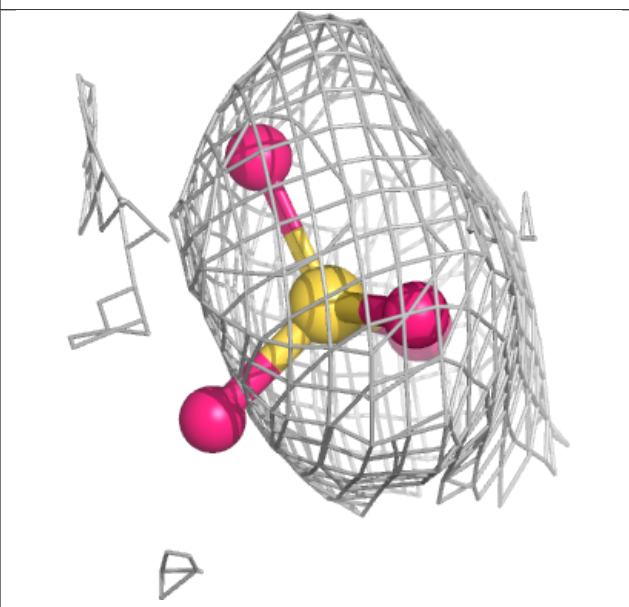
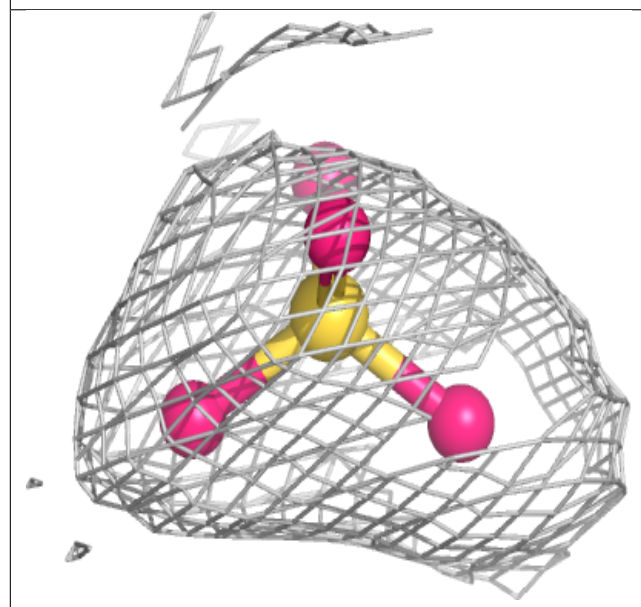
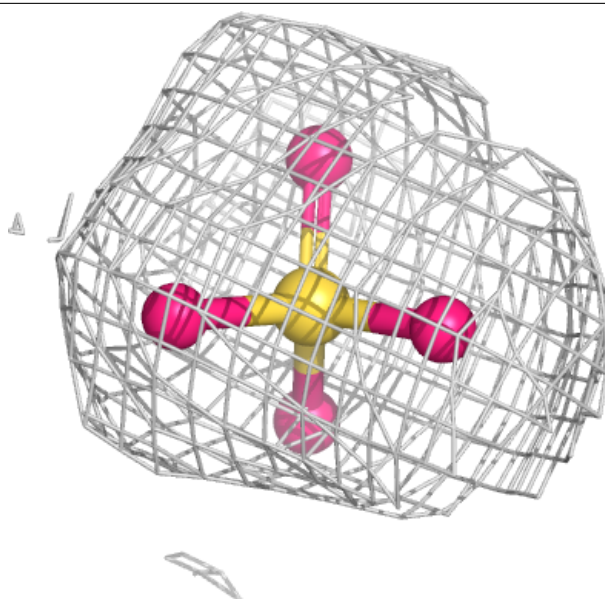
Electron density around ZN H 202:

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



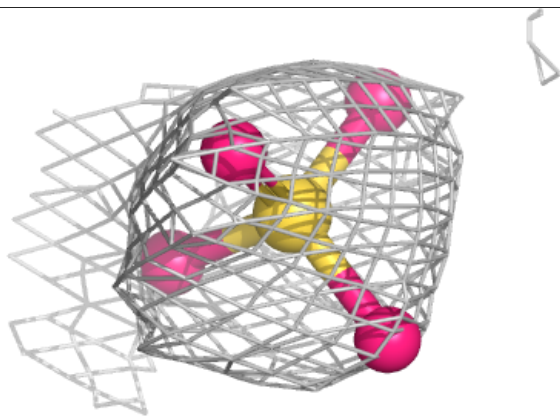
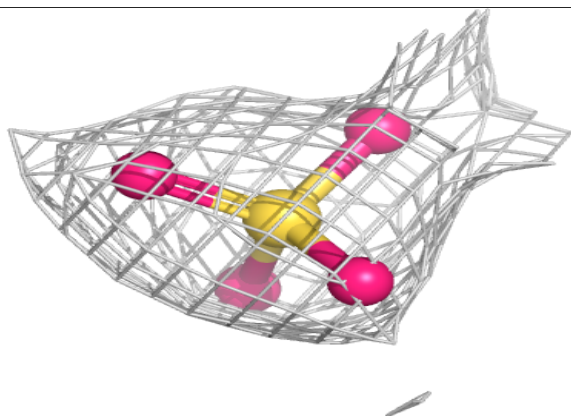
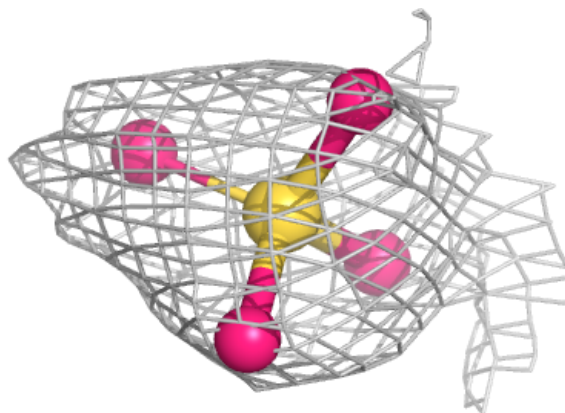
Electron density around SO4 A 205:

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



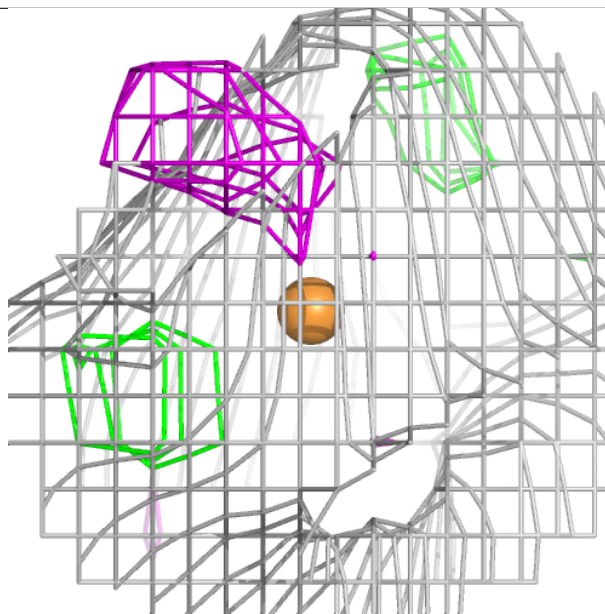
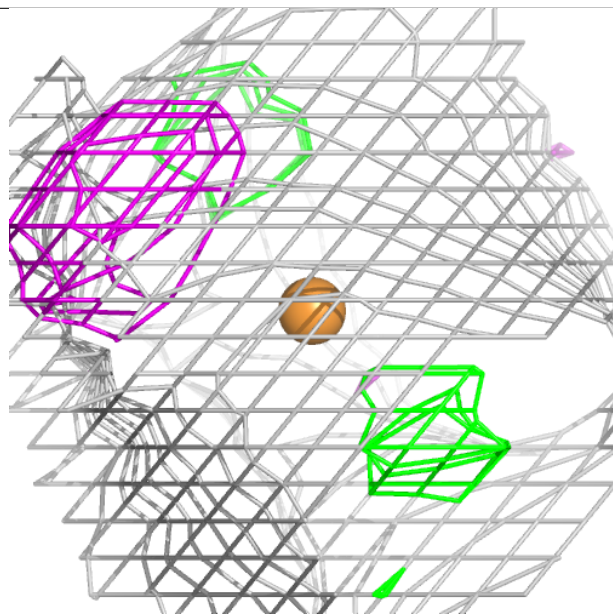
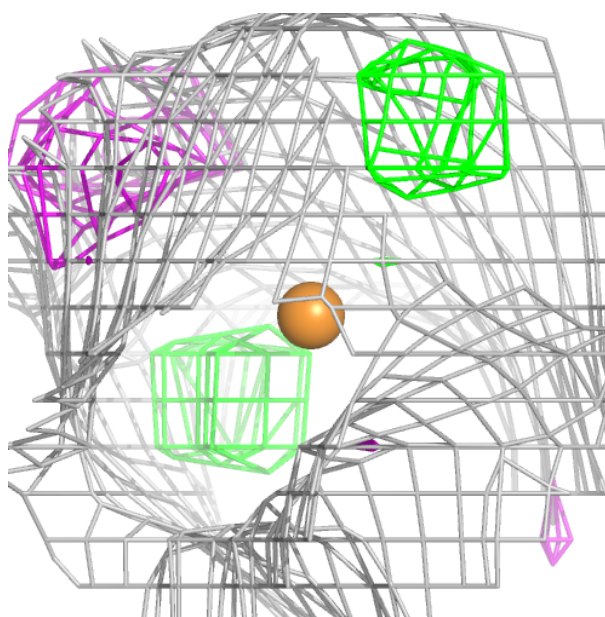
Electron density around SO4 C 205:

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



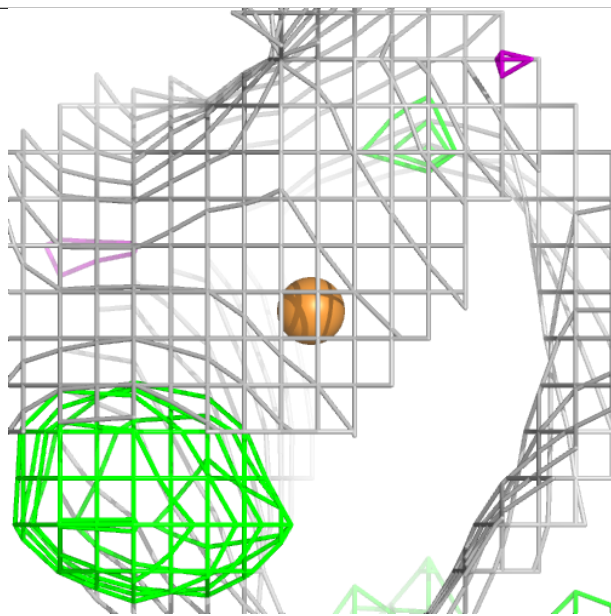
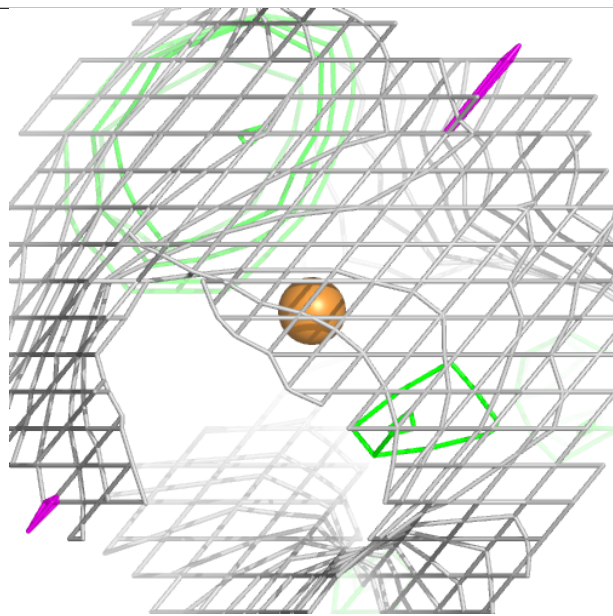
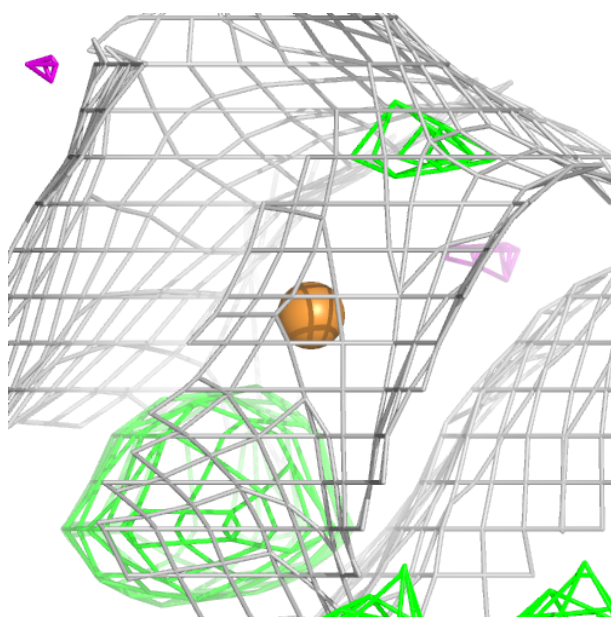
Electron density around CU H 201:

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



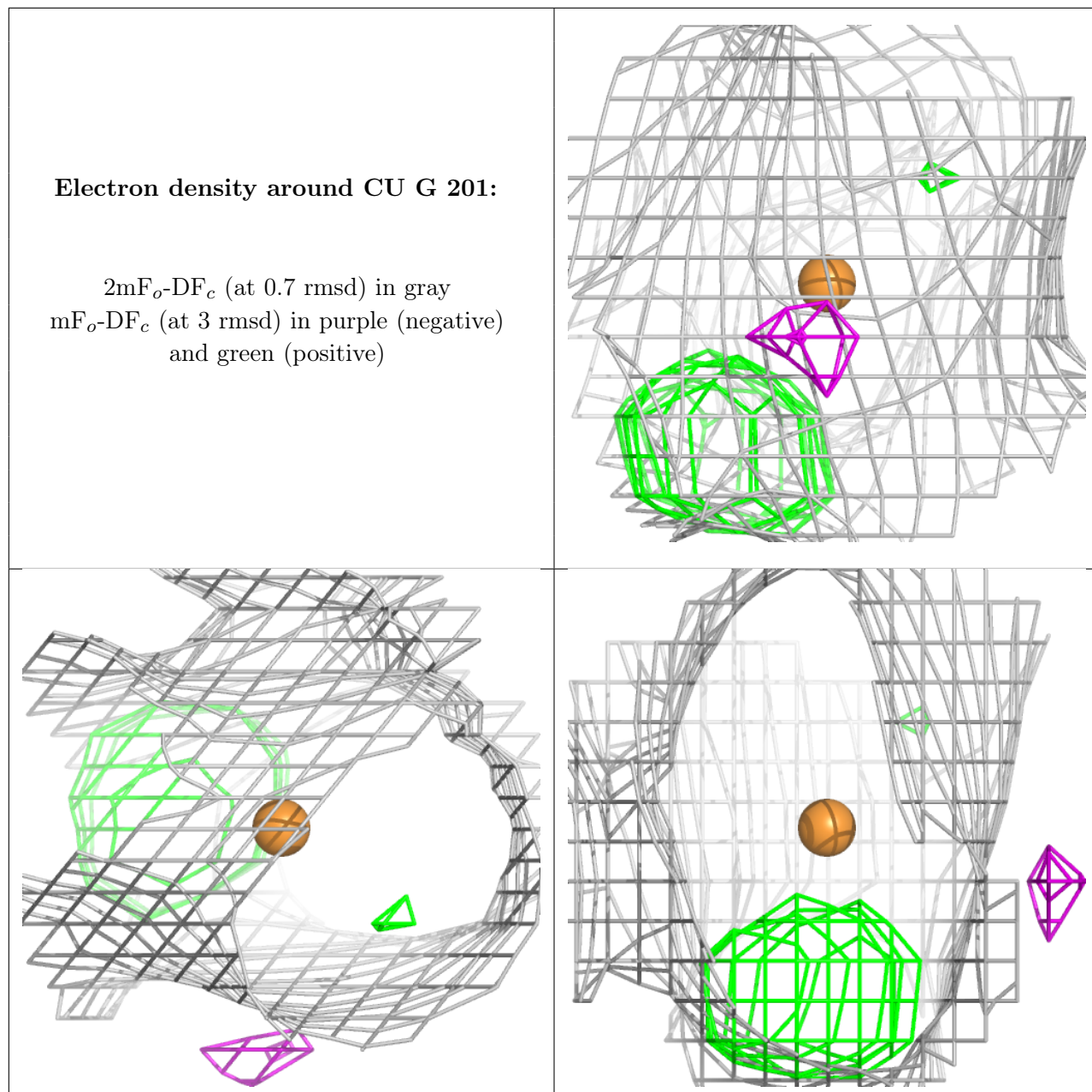
Electron density around CU D 201:

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



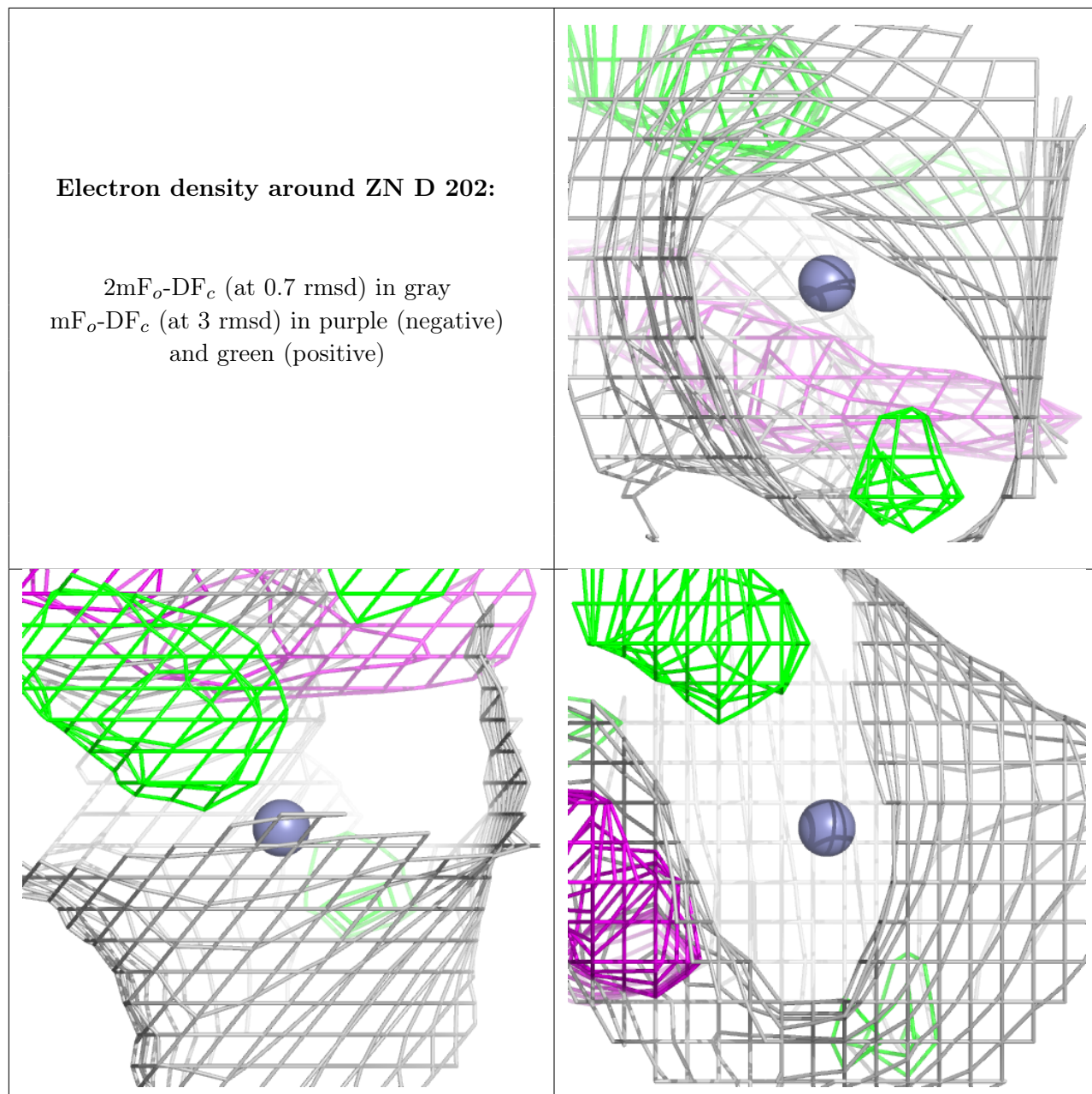
Electron density around CU G 201:

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



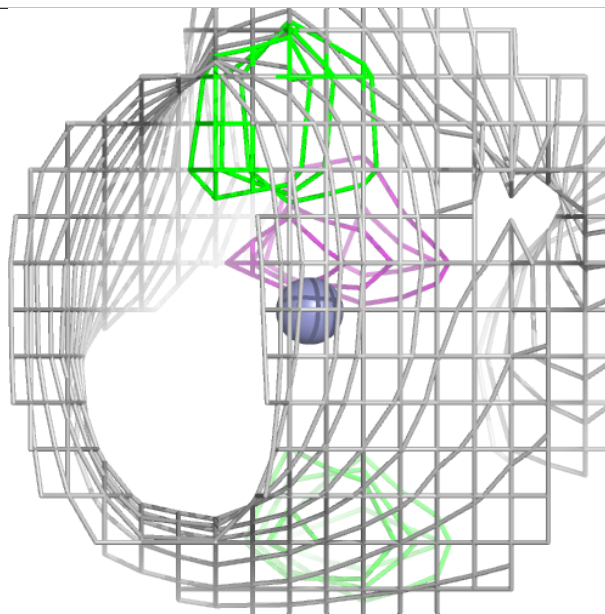
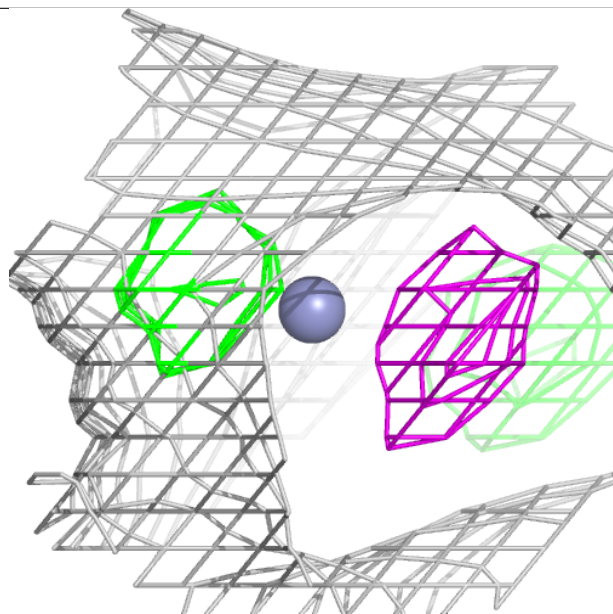
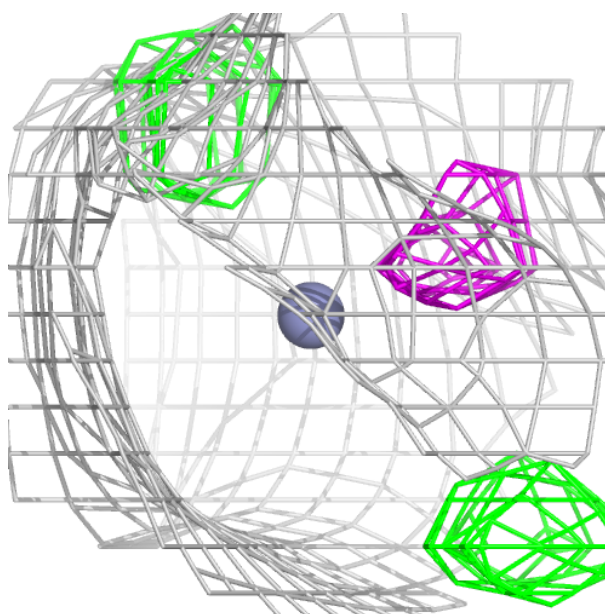
Electron density around ZN D 202:

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



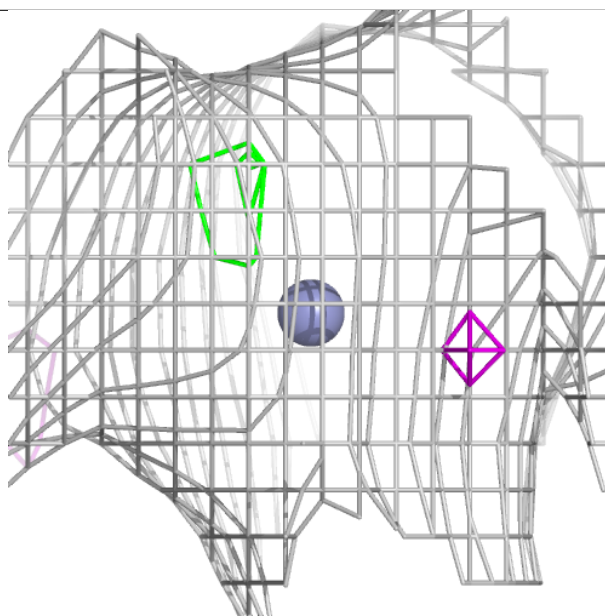
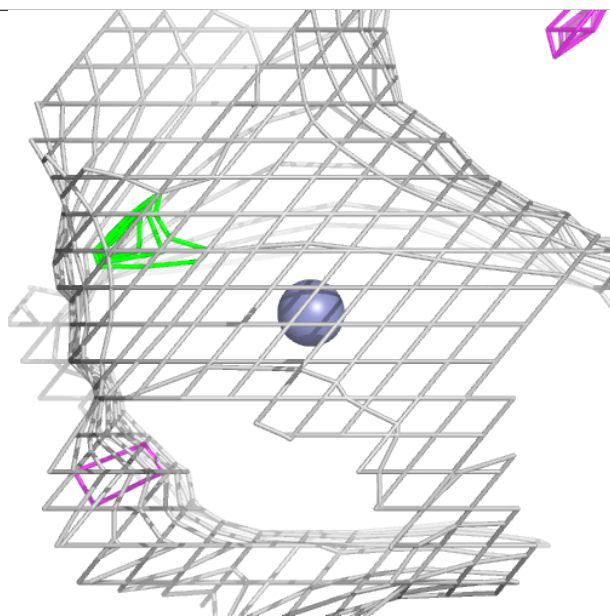
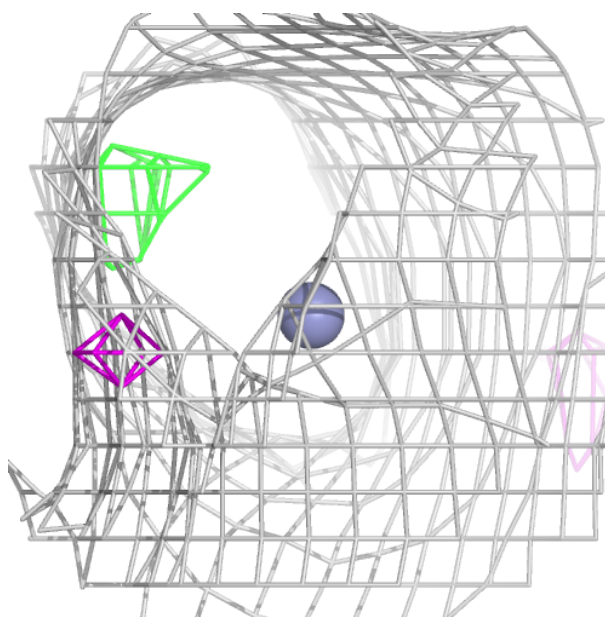
Electron density around ZN G 202:

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



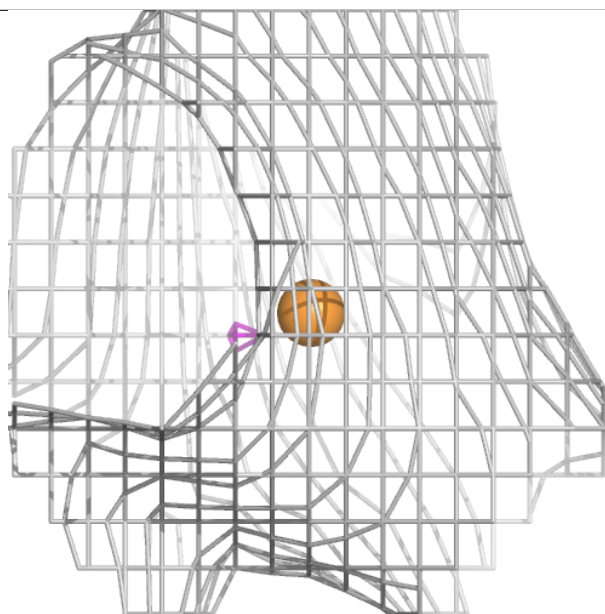
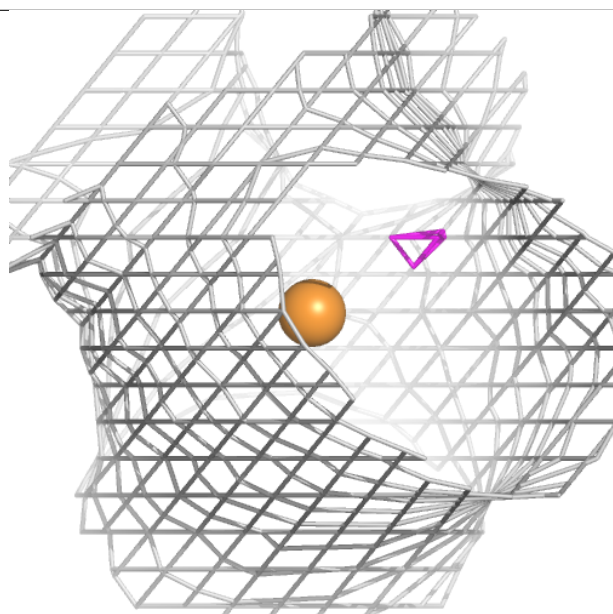
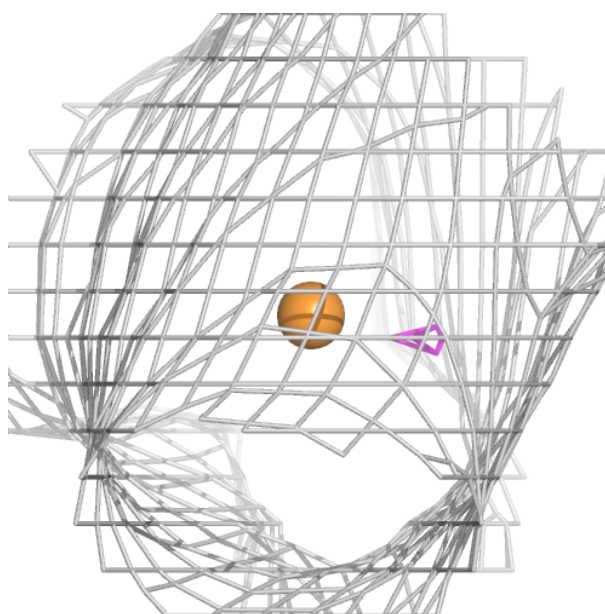
Electron density around ZN J 202:

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



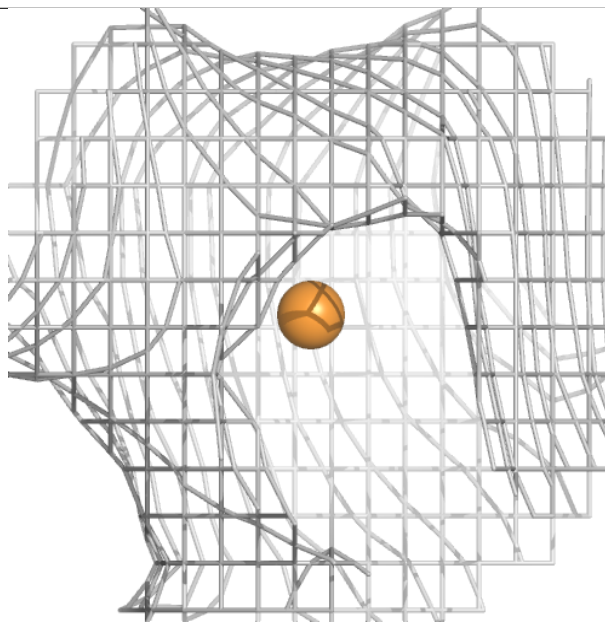
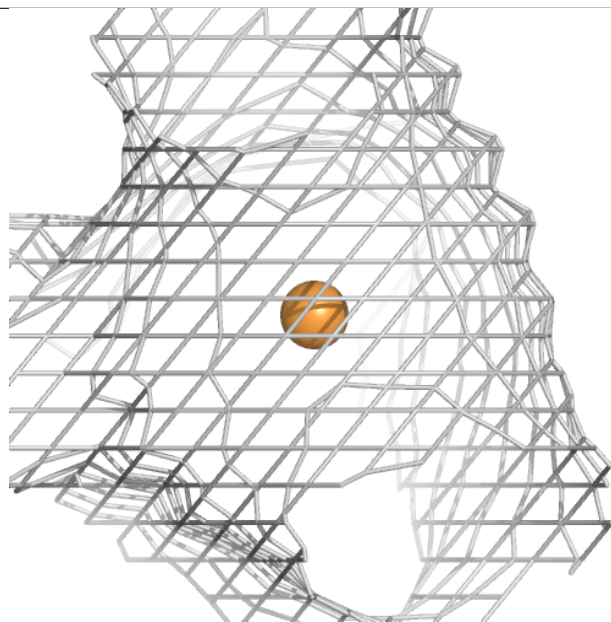
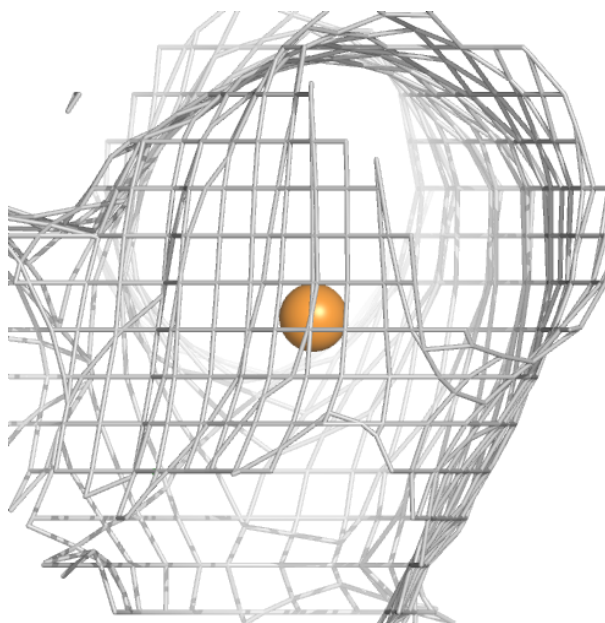
Electron density around CU J 201:

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



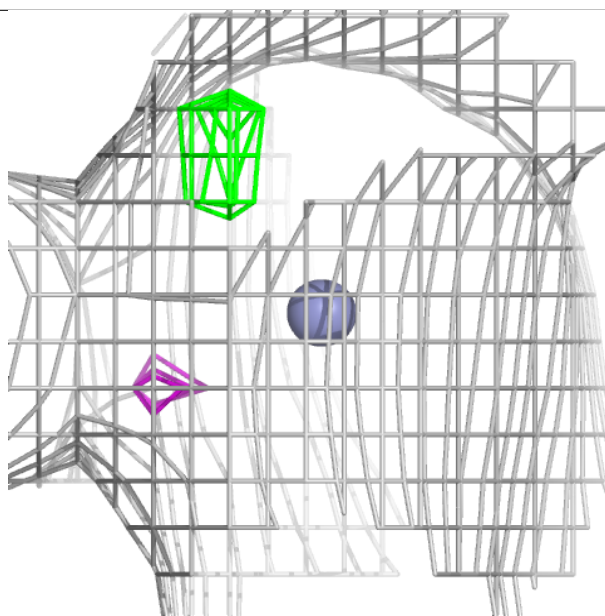
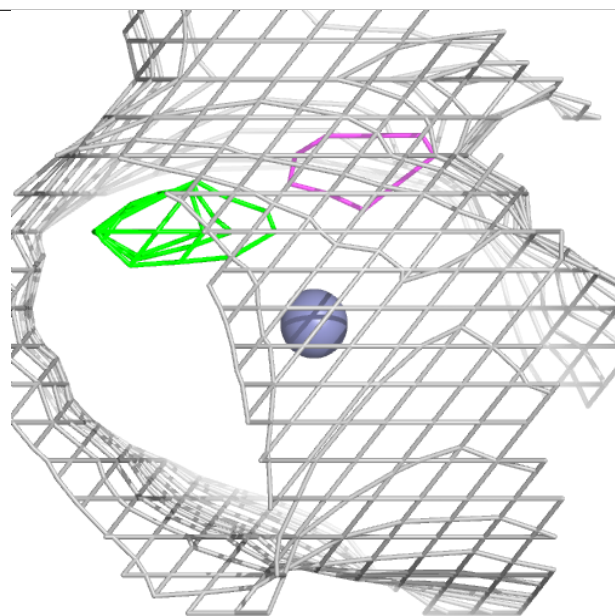
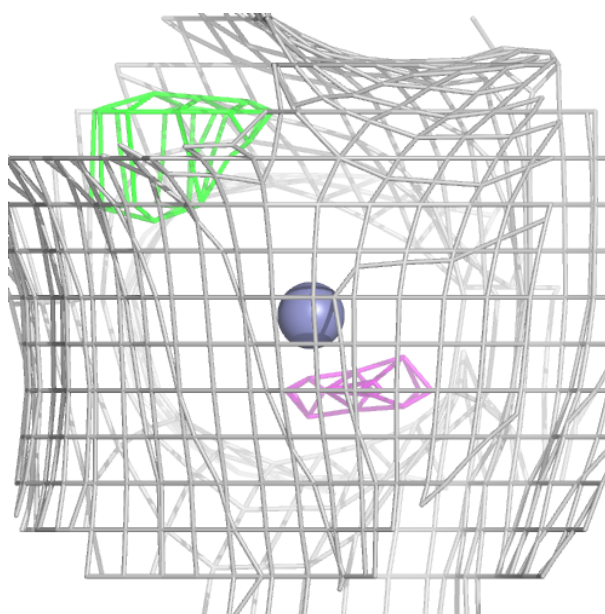
Electron density around CU I 201:

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



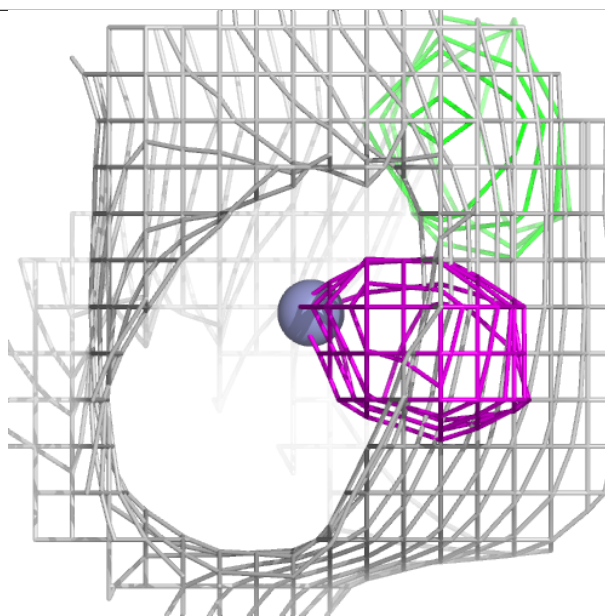
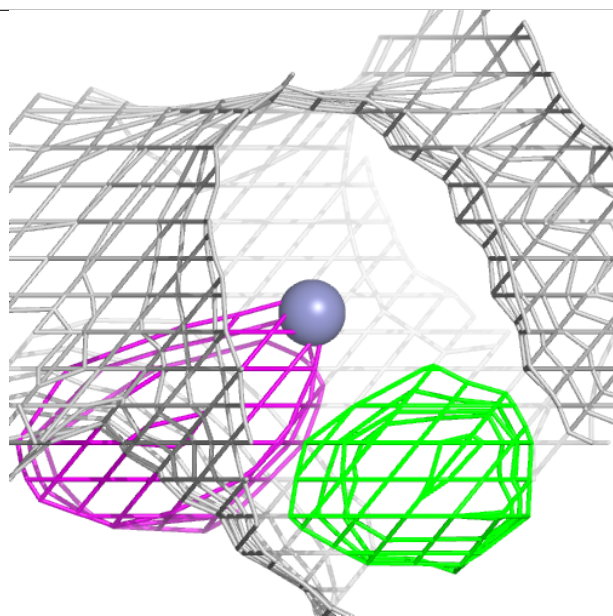
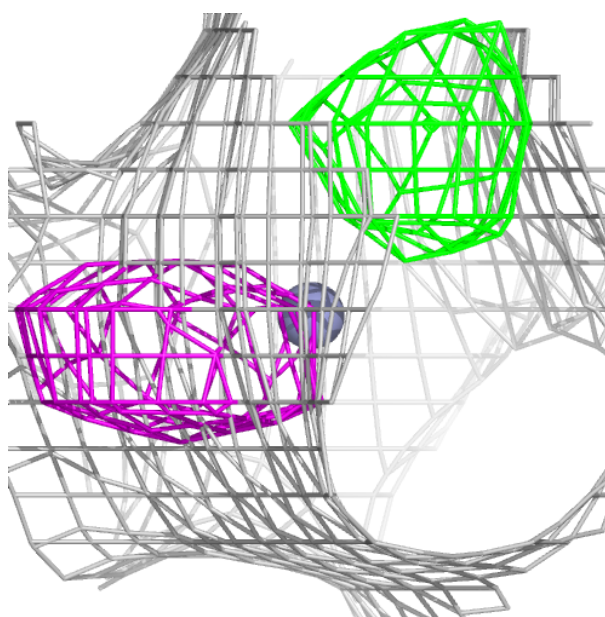
Electron density around ZN I 202:

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



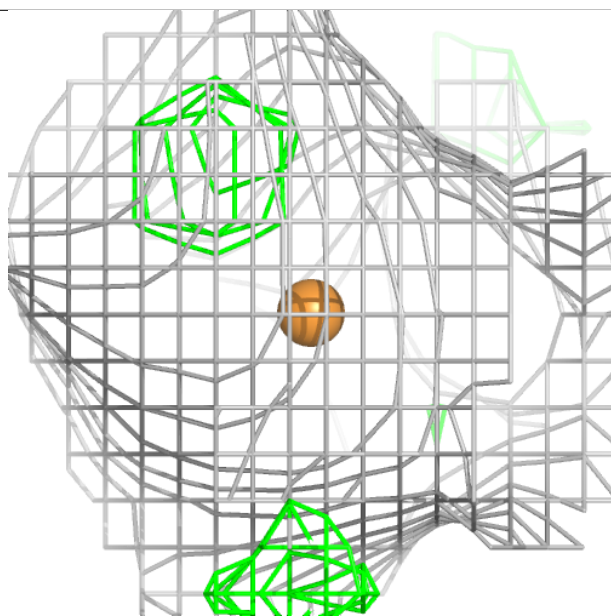
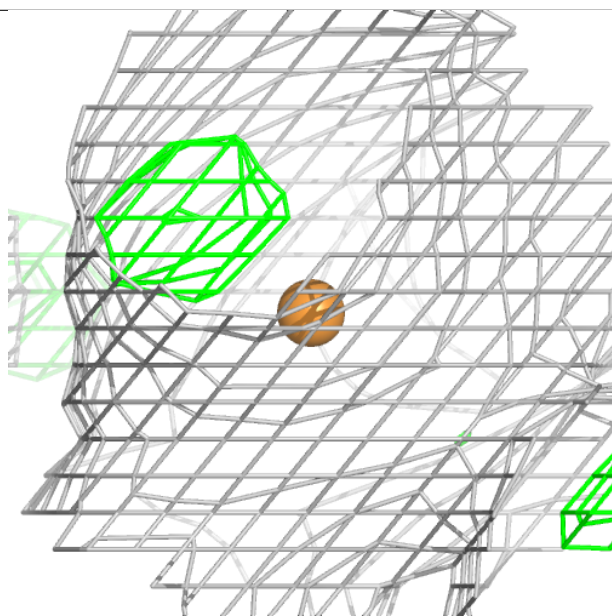
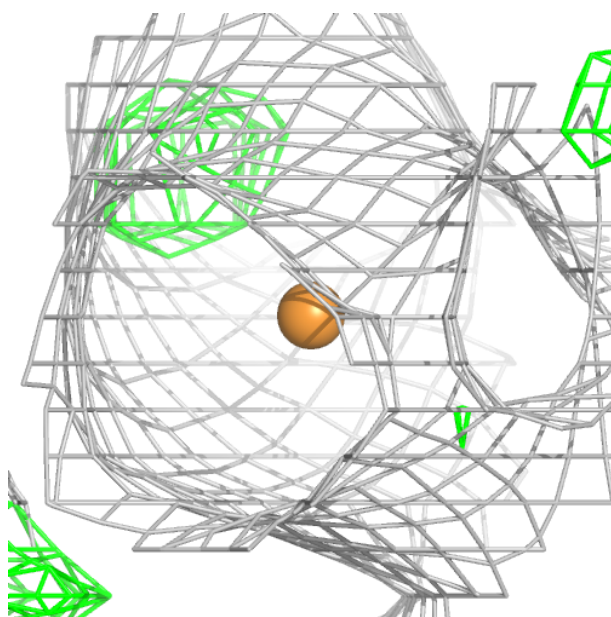
Electron density around ZN L 202:

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



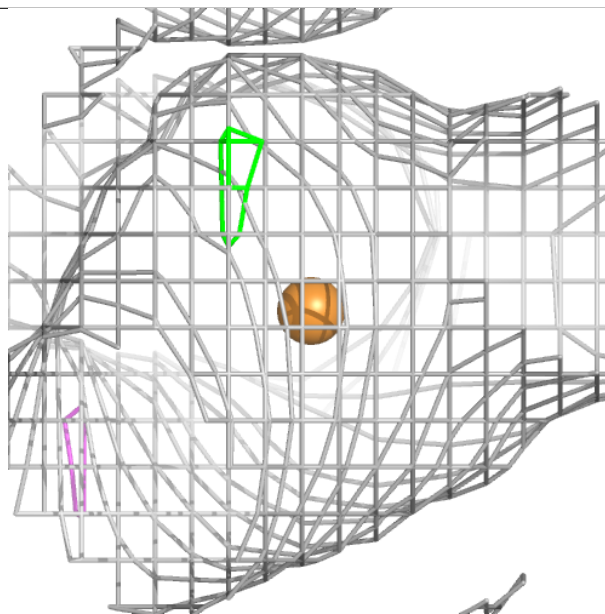
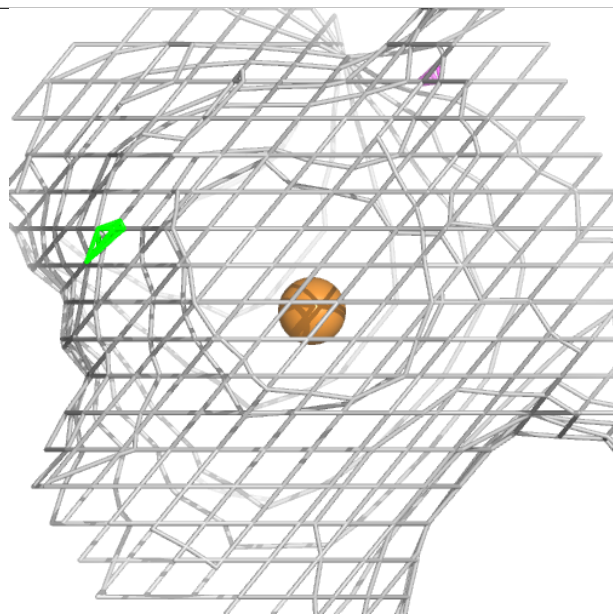
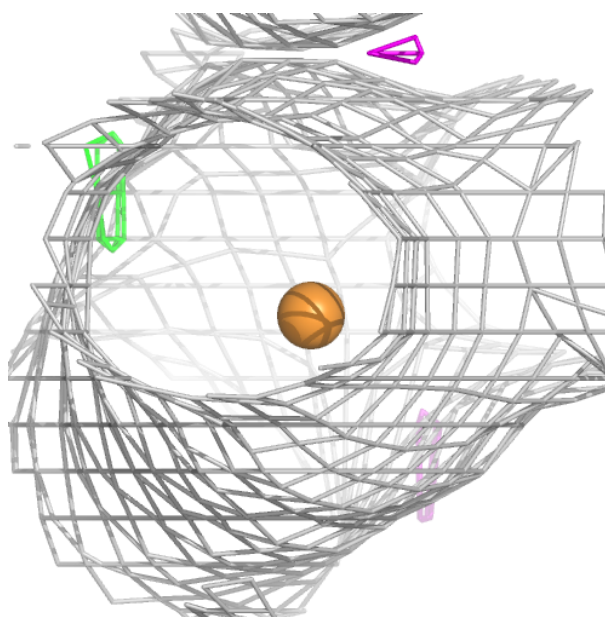
Electron density around CU B 201:

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



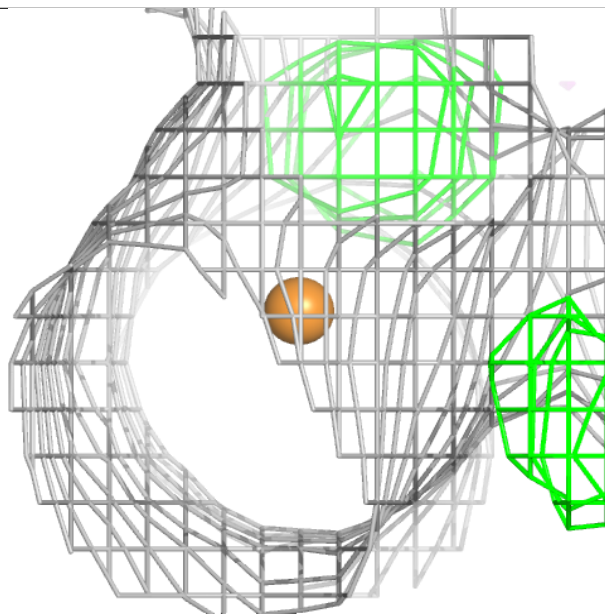
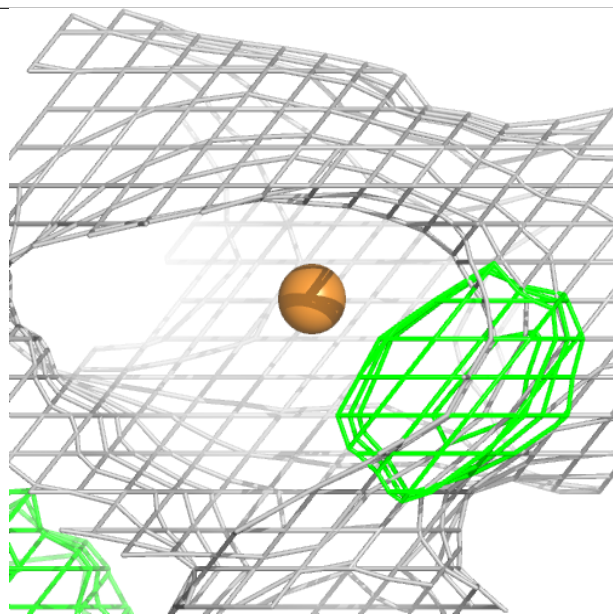
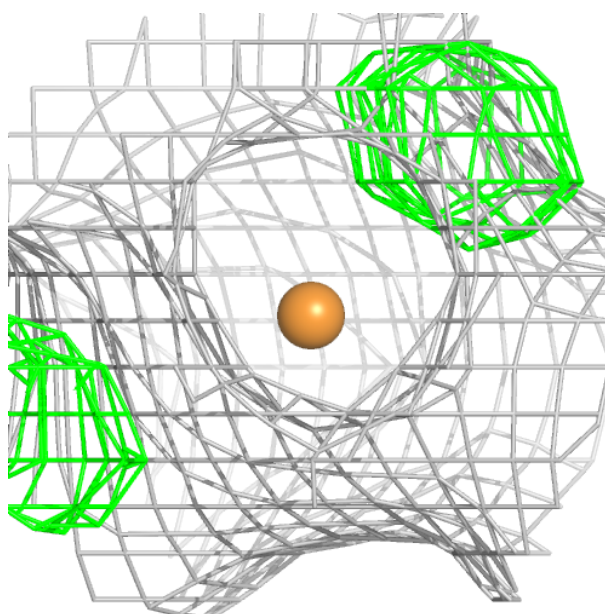
Electron density around CU C 202:

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



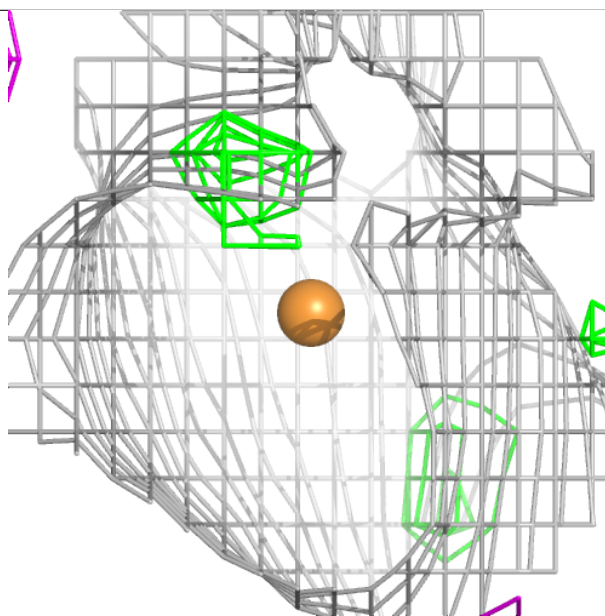
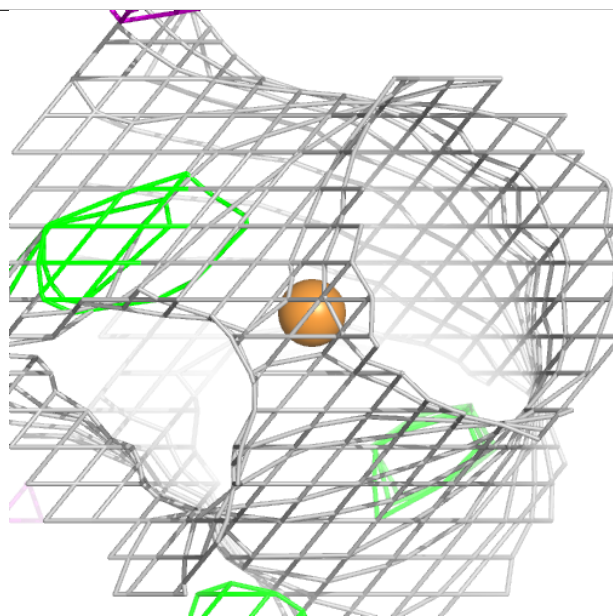
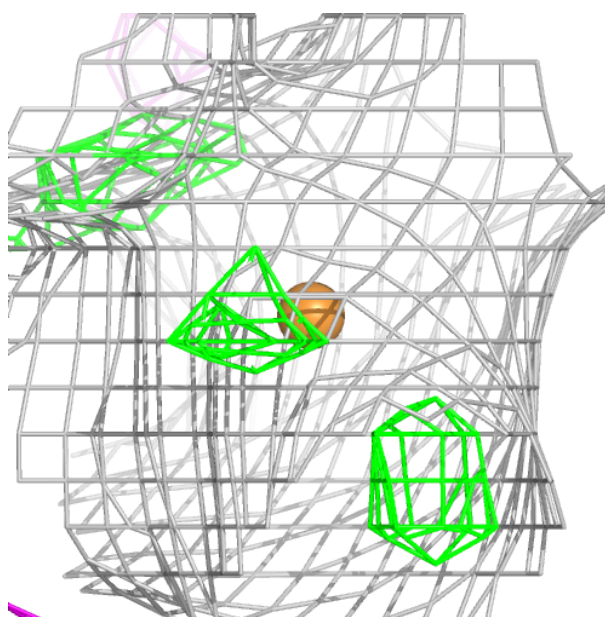
Electron density around CU K 201:

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



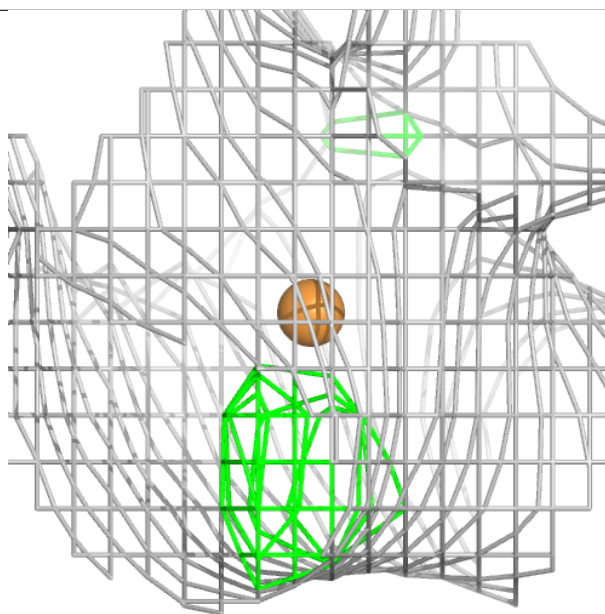
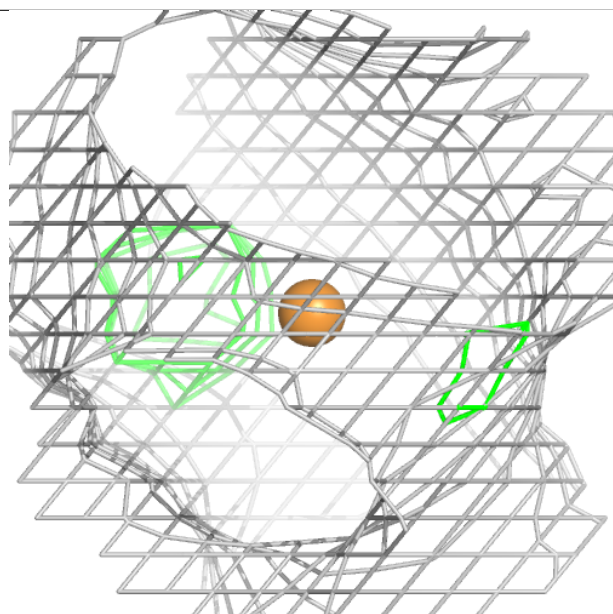
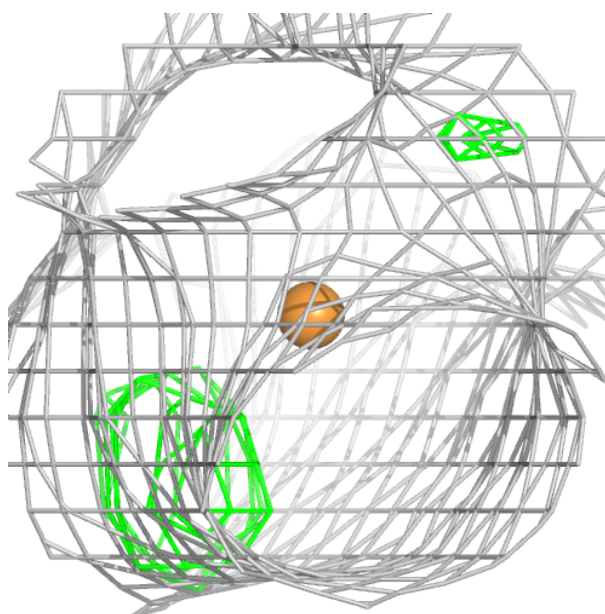
Electron density around CU L 201:

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



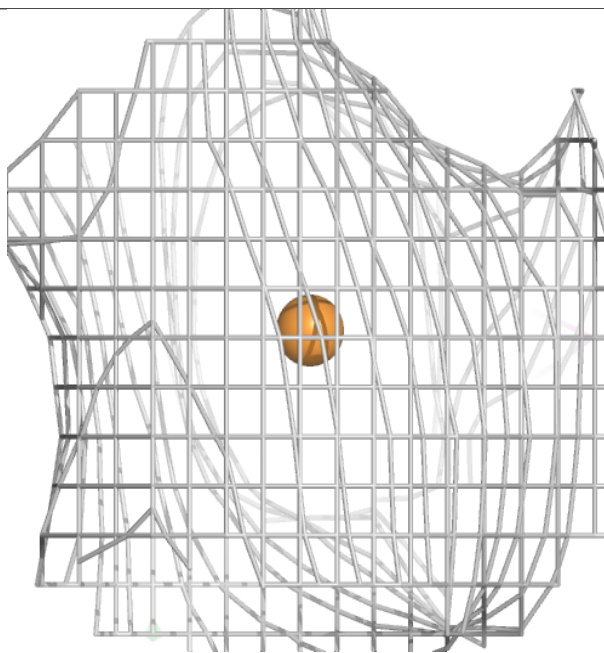
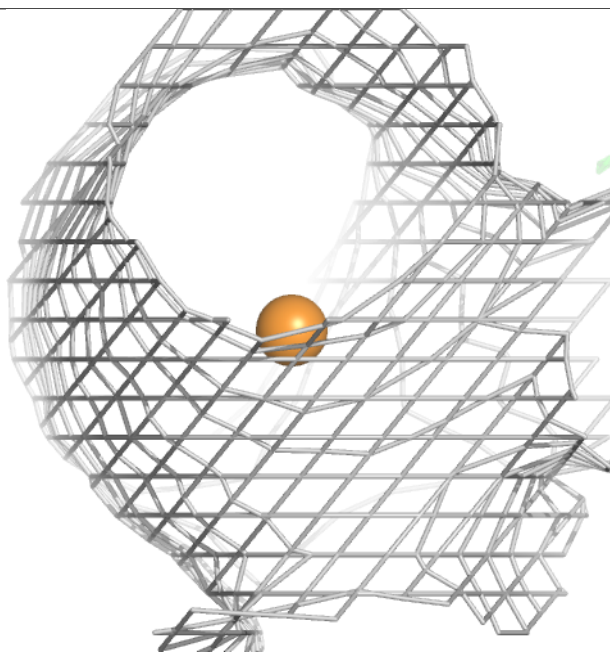
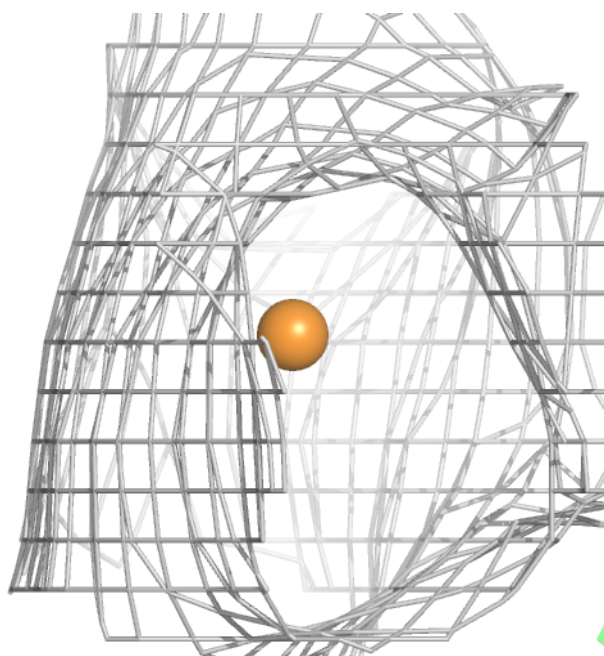
Electron density around CU F 201:

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



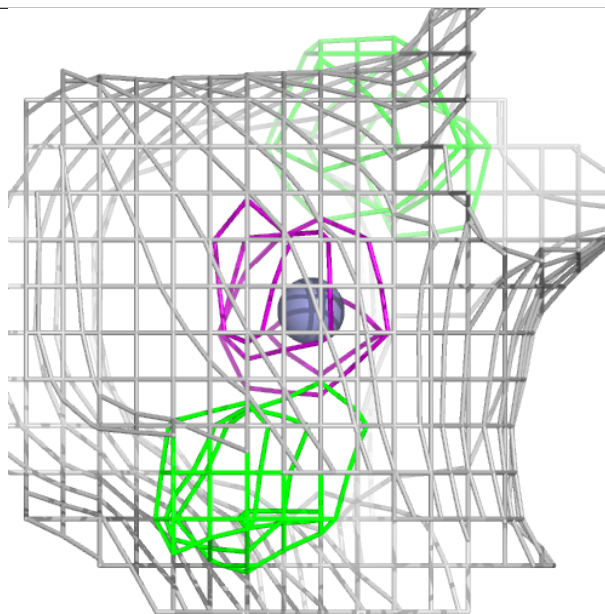
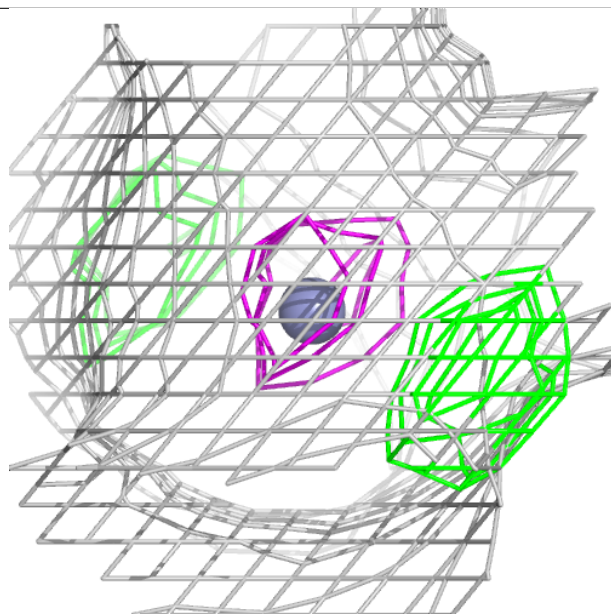
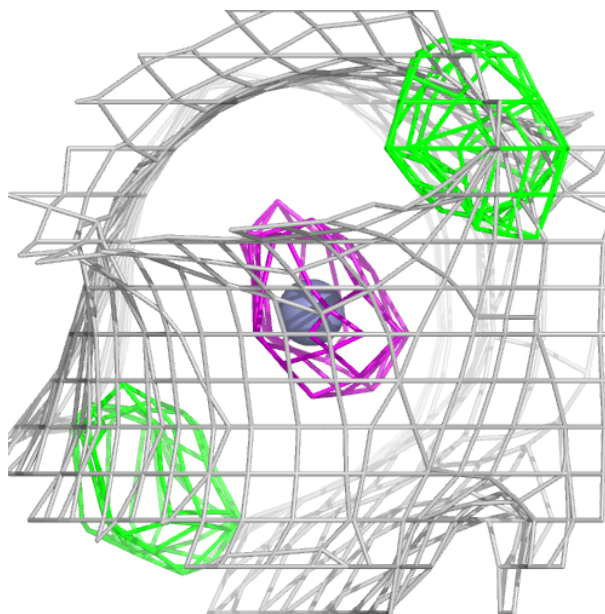
Electron density around CU E 201:

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



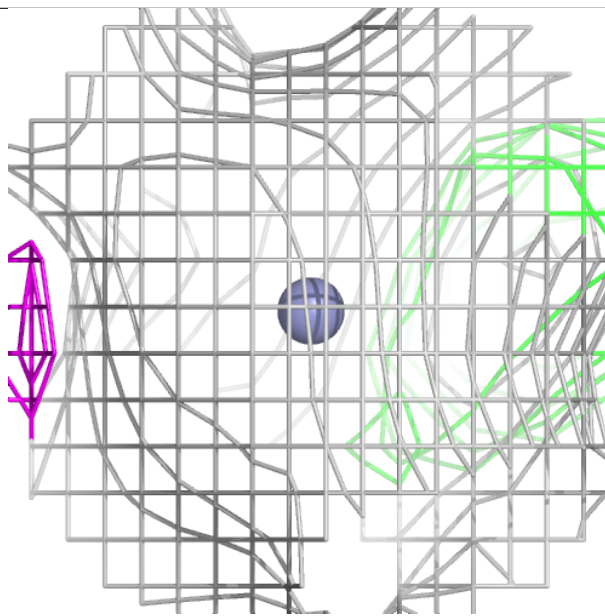
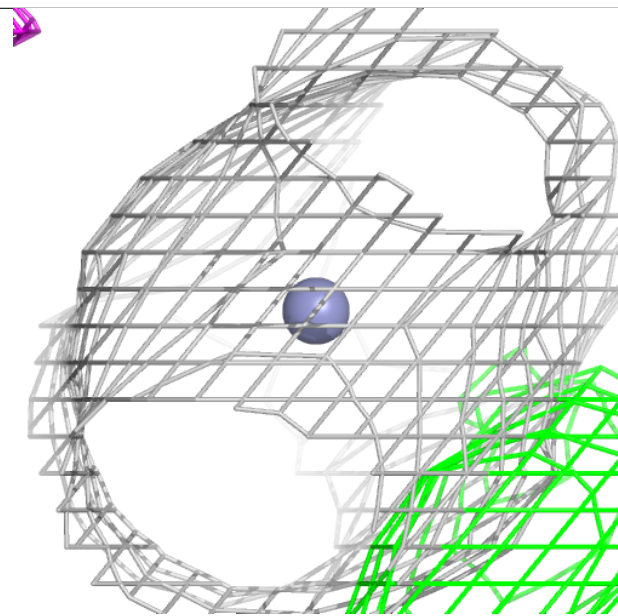
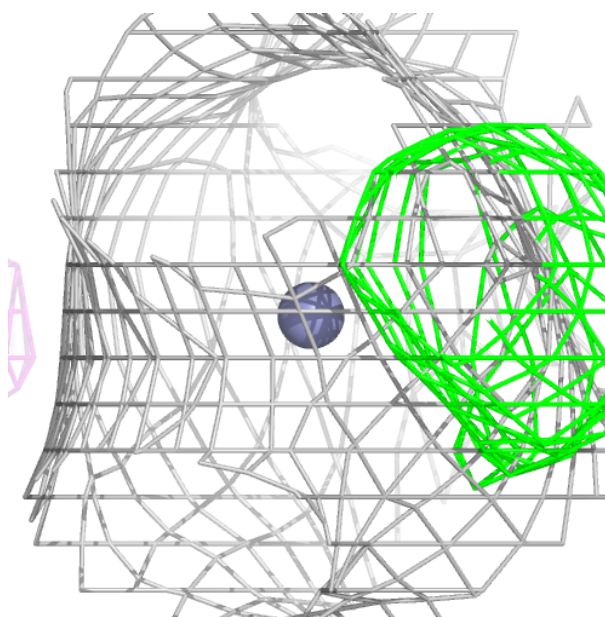
Electron density around ZN A 202:

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



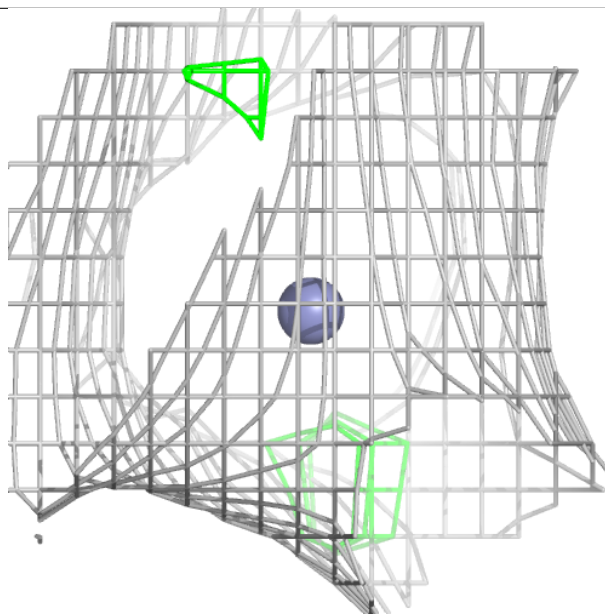
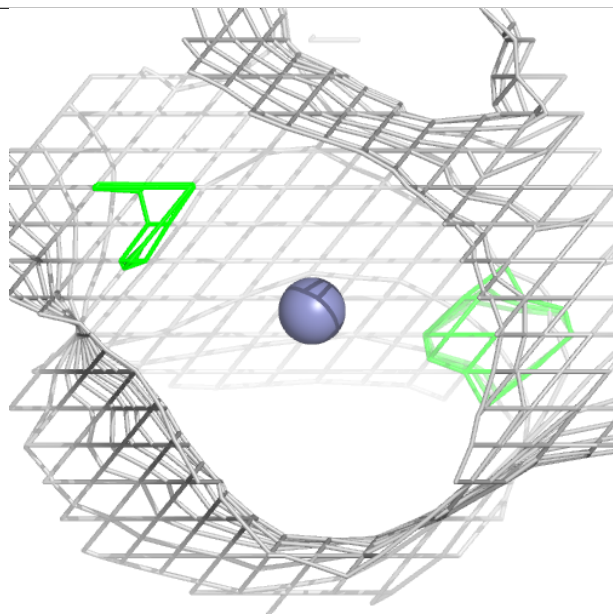
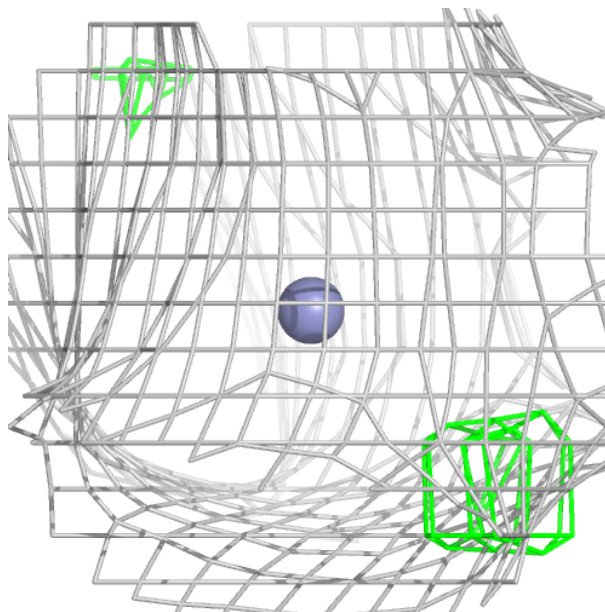
Electron density around ZN A 203:

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



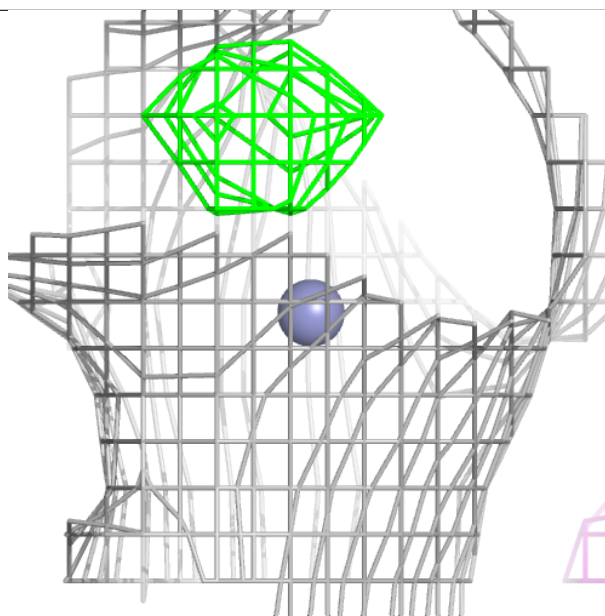
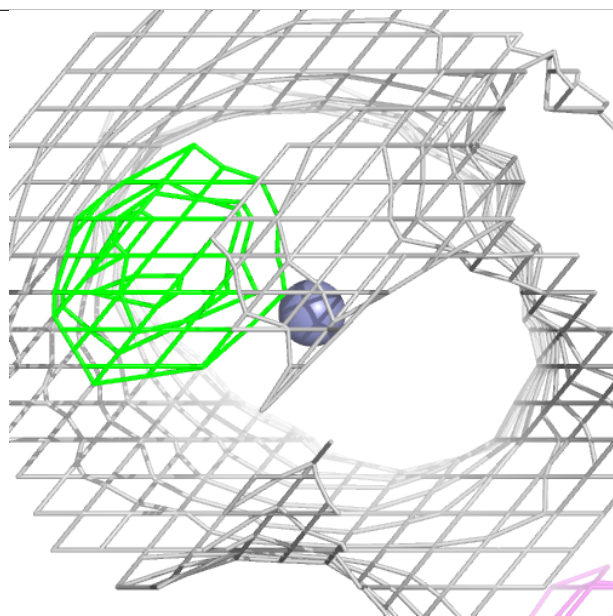
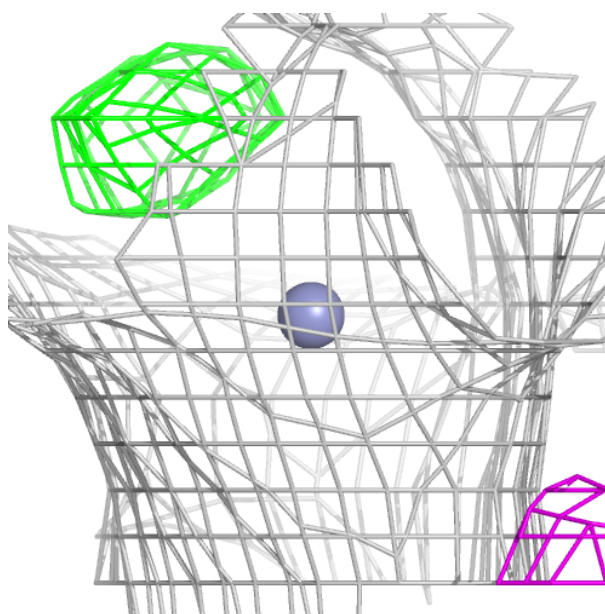
Electron density around ZN K 202:

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



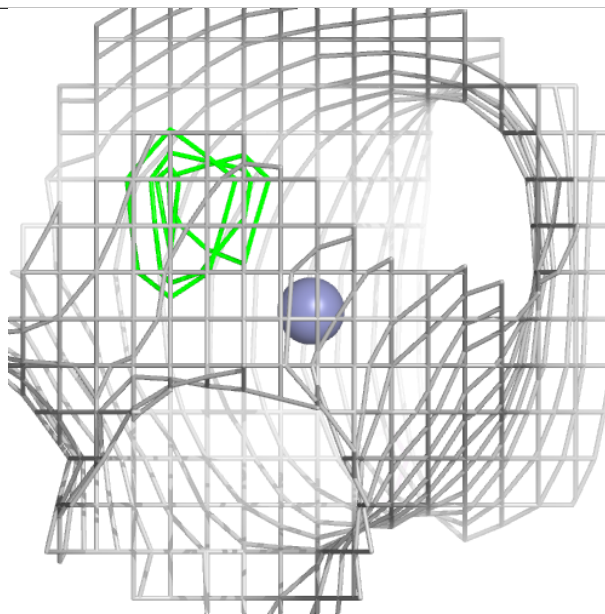
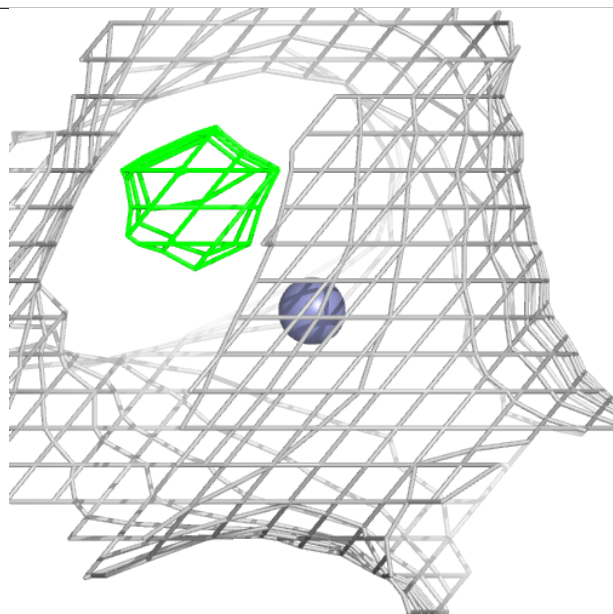
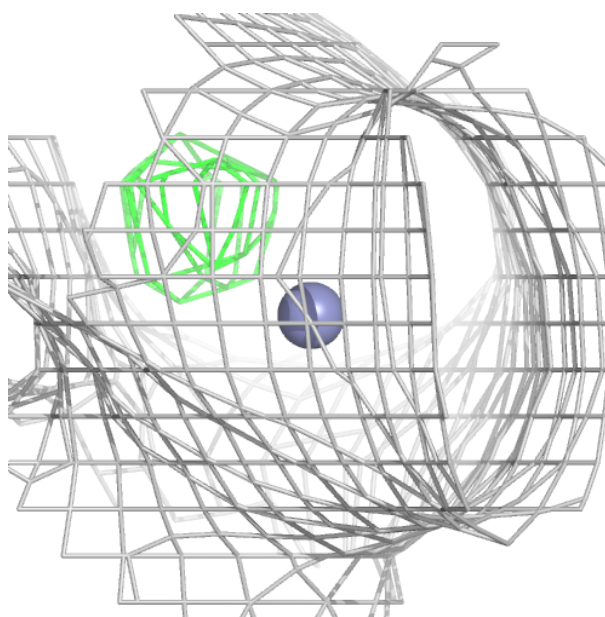
Electron density around ZN F 202:

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



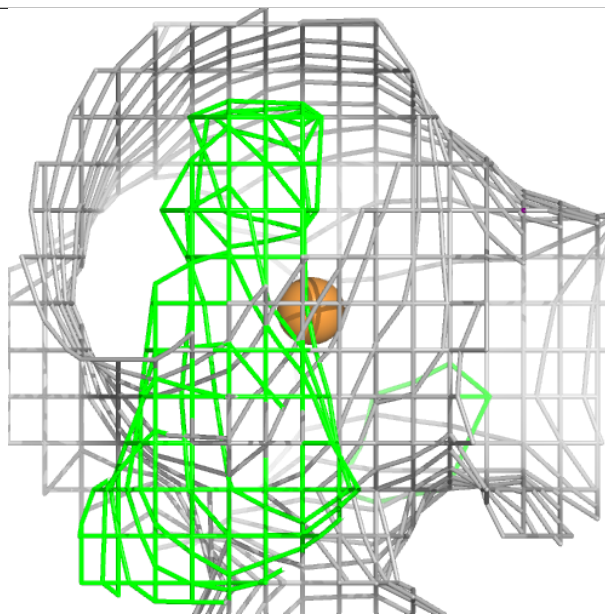
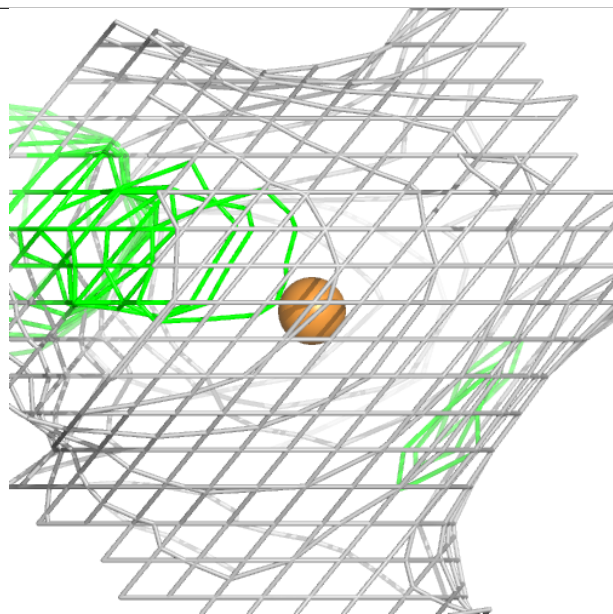
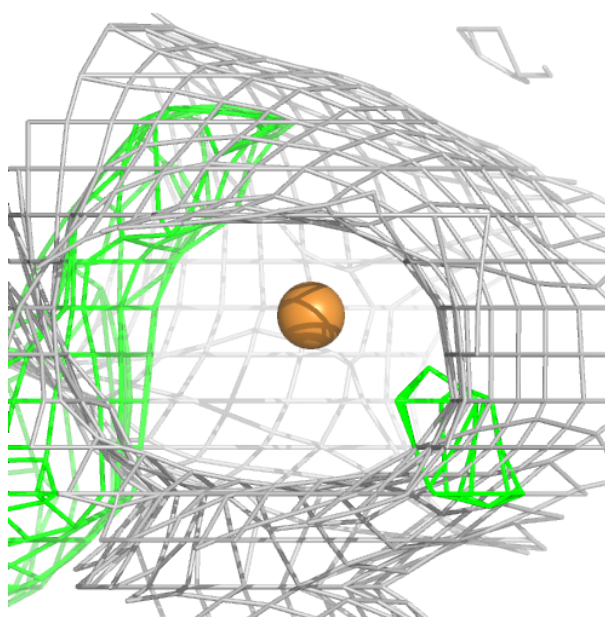
Electron density around ZN E 202:

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



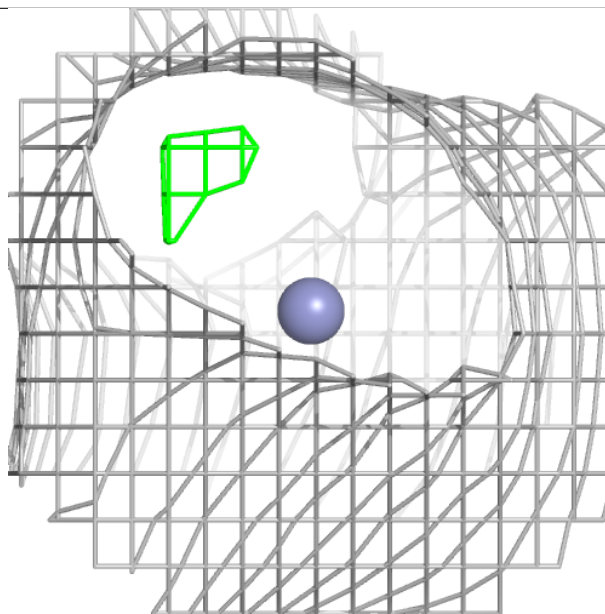
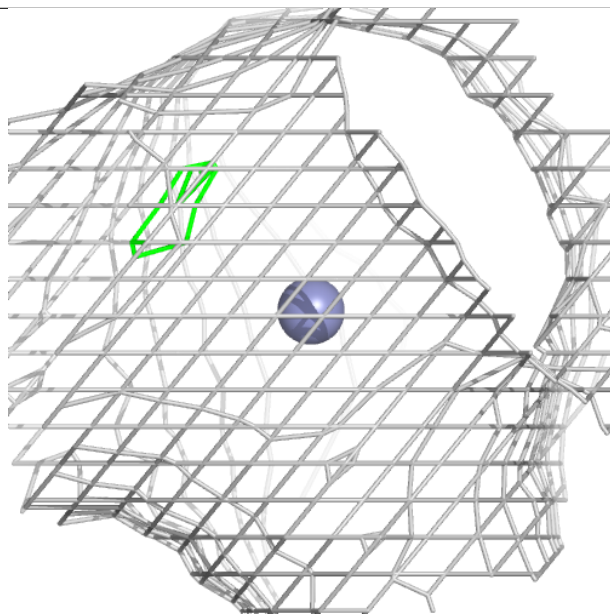
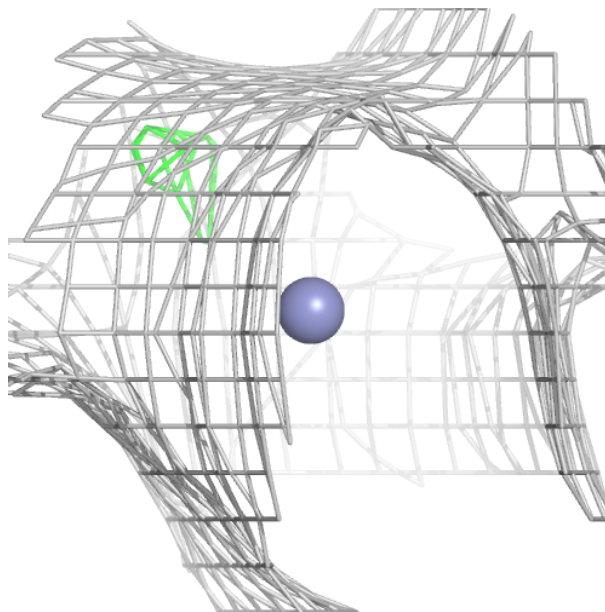
Electron density around CU A 201:

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



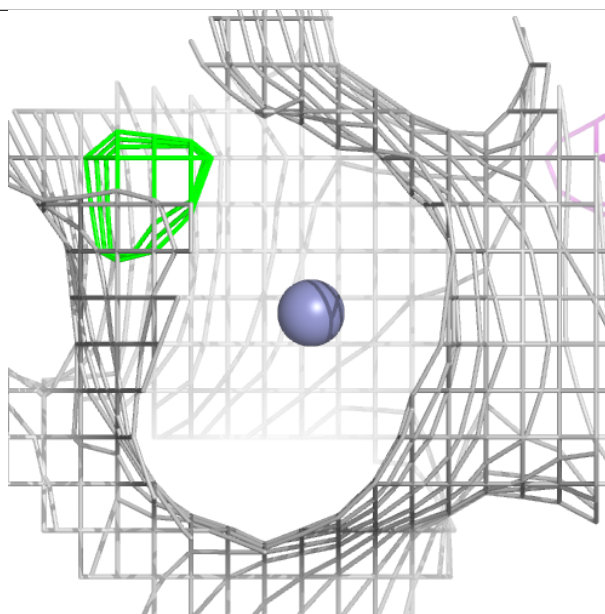
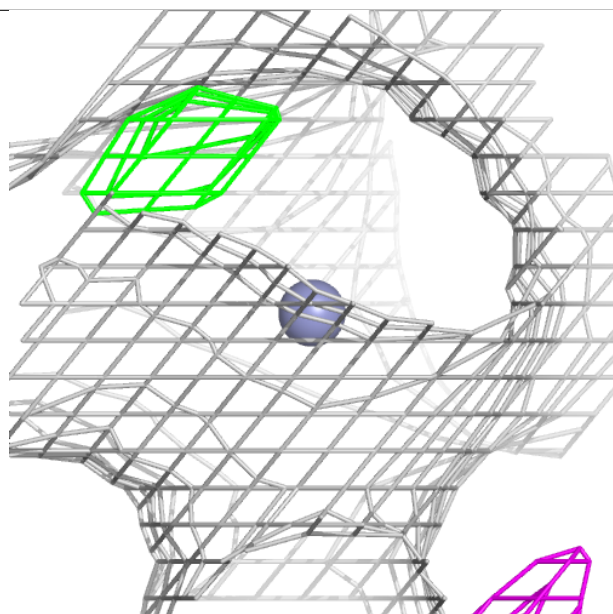
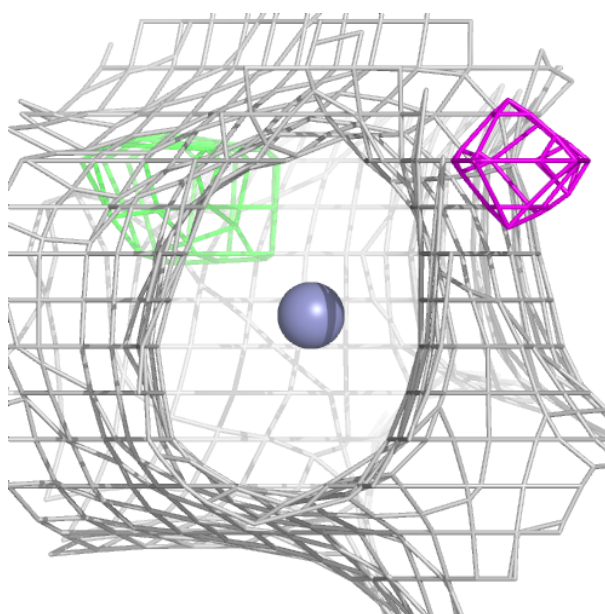
Electron density around ZN B 202:

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



Electron density around ZN C 203:

$2mF_o-DF_c$ (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.