



Full wwPDB X-ray Structure Validation Report ⓘ

Jul 6, 2026 – 12:15 PM EDT

PDB ID : 9PH8 / pdb_00009ph8
Title : Crystal structure of SARS-CoV-2 nsp14 with SAM bound
Authors : Caputo, A.T.; Dennis, M.L.; Blackman, L.D.
Deposited on : 2025-07-08
Resolution : 2.79 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 : 9.0.015 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.50

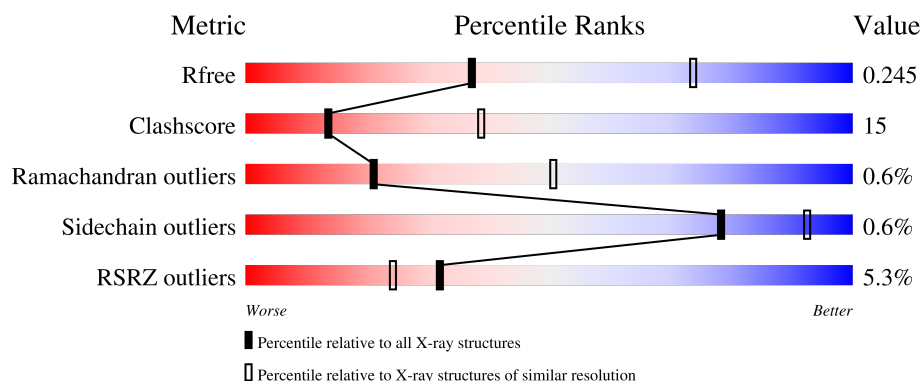
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.79 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	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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	531	
1	B	531	

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 12957 atoms, of which 6338 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 Guanine-N7 methyltransferase nsp14.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	428	Total	C	H	N	O	S	3047	0	0
			6793	2226	3336	584	615	32			
1	B	380	Total	C	H	N	O	S	2955	0	0
			6038	1982	2960	521	547	28			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	21	SER	-	expression tag	UNP P0DTD1
A	22	ASN	-	expression tag	UNP P0DTD1
A	23	GLY	-	expression tag	UNP P0DTD1
A	24	SER	-	expression tag	UNP P0DTD1
B	21	SER	-	expression tag	UNP P0DTD1
B	22	ASN	-	expression tag	UNP P0DTD1
B	23	GLY	-	expression tag	UNP P0DTD1
B	24	SER	-	expression tag	UNP P0DTD1

- Molecule 2 is S-ADENOSYLMETHIONINE (CCD ID: SAM) (formula: C₁₅H₂₂N₆O₅S) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	H	N	O	21	0
			48	15	21	6	5		
2	B	1	Total	C	H	N	O	21	0
			48	15	21	6	5		

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

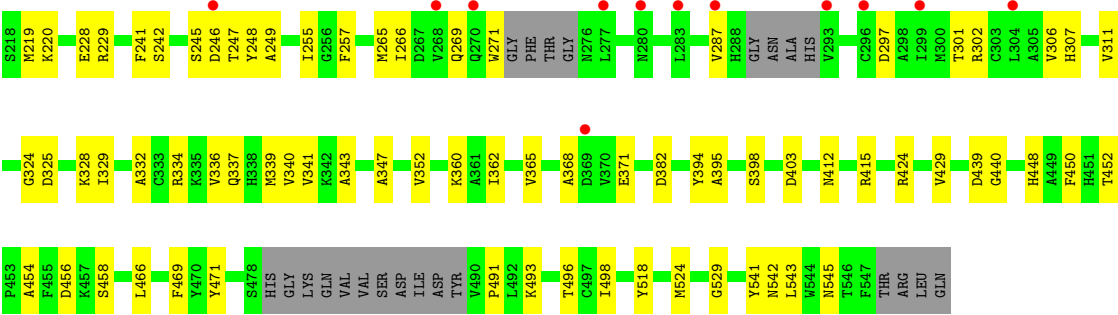
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	4	Total	Zn	0	0
			4	4		
3	B	4	Total	Zn	0	0
			4	4		

- Molecule 4 is MANGANESE (II) ION (CCD ID: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Mn	0	0
			1	1		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	9	Total	O	0	0
			9	9		
5	B	12	Total	O	0	0
			12	12		



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	67.22Å 100.56Å 90.68Å 90.00° 109.35° 90.00°	Depositor
Resolution (Å)	65.17 – 2.79 65.17 – 2.79	Depositor EDS
% Data completeness (in resolution range)	69.3 (65.17-2.79) 64.3 (65.17-2.79)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	-1.34 (at 2.77Å)	Xtriage
Refinement program	PHENIX 1.21_5207	Depositor
R, R_{free}	0.220 , 0.248 0.217 , 0.245	Depositor DCC
R_{free} test set	907 reflections (3.18%)	wwPDB-VP
Wilson B-factor (Å ²)	66.3	Xtriage
Anisotropy	0.050	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 39.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.033 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	12957	wwPDB-VP
Average B, all atoms (Å ²)	87.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.11% 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: ZN, MN, SAM

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	0.37	1/3553 (0.0%)	0.58	1/4819 (0.0%)
1	B	0.34	0/3164	0.57	1/4294 (0.0%)
All	All	0.36	1/6717 (0.0%)	0.58	2/9113 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	262	ASN	C-O	6.90	1.27	1.23

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	262	ASN	O-C-N	-7.95	117.87	121.53
1	B	185	VAL	N-CA-C	-5.85	106.78	111.81

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	3457	3336	3334	95	2
1	B	3078	2960	2959	93	2
2	A	27	21	20	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	27	21	20	1	0
3	A	4	0	0	0	0
3	B	4	0	0	0	0
4	A	1	0	0	0	0
5	A	9	0	0	0	0
5	B	12	0	0	1	0
All	All	6619	6338	6333	187	2

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 15.

All (187) 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:A:290:ASN:OD1	1:A:291:ALA:N	2.09	0.86
1:A:347:ALA:HB2	1:A:543:LEU:HD11	1.61	0.82
1:A:59:THR:HG22	1:A:73:MET:HE2	1.62	0.79
1:B:329:ILE:HG12	1:B:524:MET:HE2	1.68	0.76
1:A:96:MET:HE3	1:A:265:MET:HE3	1.69	0.73
1:A:352:VAL:HG23	1:A:371:GLU:HB3	1.71	0.72
1:B:219:MET:SD	1:B:265:MET:HE1	2.31	0.71
1:B:229:ARG:HD3	1:B:248:TYR:CE1	2.24	0.71
1:A:365:VAL:CG1	1:A:368:ALA:HB2	2.21	0.70
1:B:216:LEU:HA	1:B:219:MET:HE2	1.74	0.69
1:A:219:MET:HE1	1:A:265:MET:SD	2.31	0.69
1:A:93:TYR:OH	1:A:263:PRO:O	2.11	0.68
1:A:365:VAL:HG11	1:A:368:ALA:HB2	1.76	0.68
1:B:217:THR:O	1:B:220:LYS:HE2	1.94	0.68
1:A:96:MET:HE3	1:A:265:MET:CE	2.24	0.67
1:A:499:THR:OG1	1:A:501:CYS:SG	2.52	0.67
1:B:210:TRP:CZ3	1:B:269:GLN:HG3	2.29	0.67
1:A:115:VAL:HG21	1:A:215:GLU:HA	1.77	0.65
1:B:325:ASP:O	1:B:329:ILE:HG13	1.97	0.64
1:B:337:GLN:O	1:B:341:VAL:HG23	1.98	0.64
1:A:347:ALA:HB2	1:A:543:LEU:CD1	2.28	0.64
1:B:255:ILE:HD12	1:B:255:ILE:H	1.63	0.63
1:A:51:LEU:HD23	1:A:52:SER:N	2.12	0.63
1:B:241:PHE:HD1	1:B:242:SER:N	1.96	0.63
1:B:241:PHE:CD1	1:B:242:SER:N	2.66	0.63
1:B:334:ARG:NH2	2:B:601:SAM:OXT	2.32	0.62
1:B:101:GLU:OE1	1:B:101:GLU:N	2.32	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:86:MET:O	1:A:89:GLN:N	2.33	0.62
1:B:211:ALA:CB	1:B:269:GLN:OE1	2.47	0.61
1:B:343:ALA:HB1	1:B:543:LEU:HD21	1.83	0.61
1:A:365:VAL:HG12	1:A:365:VAL:O	2.00	0.61
1:A:98:ILE:O	1:A:266:ILE:HD13	2.00	0.61
1:B:307:HIS:HA	1:B:311:VAL:HG23	1.83	0.60
1:A:74:THR:HG23	1:A:77:ARG:H	1.65	0.60
1:A:212:HIS:HB2	1:A:215:GLU:OE1	2.02	0.60
1:A:57:PHE:HB2	1:A:75:TYR:CE1	2.37	0.60
1:B:429:VAL:HG21	1:B:529:GLY:HA2	1.84	0.59
1:A:115:VAL:CG2	1:A:215:GLU:HA	2.33	0.59
1:A:79:ILE:HA	1:A:82:MET:HE3	1.84	0.58
1:B:242:SER:N	1:B:257:PHE:CE1	2.72	0.58
1:A:75:TYR:O	1:A:79:ILE:HD13	2.03	0.57
1:A:79:ILE:HG13	1:A:96:MET:HE2	1.87	0.57
1:B:241:PHE:C	1:B:257:PHE:HD1	2.12	0.57
1:B:352:VAL:HG23	1:B:371:GLU:HB3	1.85	0.57
1:B:246:ASP:OD1	1:B:246:ASP:O	2.23	0.57
1:A:241:PHE:HE2	1:A:243:THR:HG22	1.69	0.56
1:A:96:MET:HB3	1:A:265:MET:CE	2.36	0.56
1:B:297:ASP:O	1:B:301:THR:OG1	2.23	0.56
1:A:53:VAL:HG21	1:A:76:ARG:CZ	2.36	0.56
1:B:365:VAL:HG12	1:B:365:VAL:O	2.04	0.56
1:A:219:MET:HE1	1:A:265:MET:HE1	1.89	0.55
1:A:99:THR:HB	1:A:101:GLU:OE1	2.07	0.55
1:A:195:SER:O	1:A:199:LYS:HG2	2.06	0.55
1:A:291:ALA:CB	1:A:293:VAL:HG23	2.37	0.55
1:B:241:PHE:CD1	1:B:241:PHE:C	2.84	0.55
1:B:135:PHE:CD2	1:B:193:MET:HE1	2.41	0.55
1:B:210:TRP:CH2	1:B:269:GLN:HG3	2.42	0.55
1:A:212:HIS:O	1:A:215:GLU:HG3	2.06	0.55
1:B:332:ALA:O	1:B:336:VAL:HG23	2.07	0.55
1:B:241:PHE:C	1:B:257:PHE:CD1	2.86	0.54
1:B:245:SER:OG	1:B:247:THR:HG23	2.07	0.54
1:A:429:VAL:HG21	1:A:529:GLY:HA2	1.88	0.54
1:B:347:ALA:HB2	1:B:543:LEU:CD1	2.37	0.54
1:A:115:VAL:CG1	1:A:131:LEU:HD22	2.38	0.54
1:B:439:ASP:OD2	1:B:439:ASP:O	2.25	0.54
1:A:137:THR:OG1	1:A:193:MET:HE1	2.07	0.54
1:B:135:PHE:CE2	1:B:193:MET:HE1	2.44	0.53
1:B:229:ARG:HD3	1:B:248:TYR:CD1	2.44	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:195:SER:O	1:B:199:LYS:HG2	2.08	0.53
1:A:246:ASP:O	1:A:246:ASP:OD1	2.27	0.53
1:B:307:HIS:HA	1:B:311:VAL:CG2	2.37	0.53
1:B:211:ALA:HB2	1:B:269:GLN:OE1	2.09	0.53
1:B:211:ALA:HB3	1:B:269:GLN:OE1	2.09	0.52
1:B:491:PRO:HB2	1:B:493:LYS:NZ	2.23	0.52
1:B:340:VAL:HG13	1:B:466:LEU:HD11	1.91	0.52
1:A:333:CYS:SG	1:A:525:MET:CE	2.98	0.52
1:A:337:GLN:HB2	1:A:409:TRP:CE2	2.44	0.52
1:A:139:VAL:HG21	1:A:193:MET:HE2	1.92	0.52
1:B:98:ILE:O	1:B:266:ILE:HD13	2.09	0.52
1:A:58:LYS:HD3	1:A:270:GLN:NE2	2.25	0.51
1:B:365:VAL:CG1	1:B:368:ALA:HB2	2.40	0.51
1:A:498:ILE:HG13	1:A:544:TRP:CZ2	2.45	0.51
1:A:354:HIS:O	1:A:356:ILE:HD12	2.10	0.50
1:A:513:ASN:OD1	1:A:516:ARG:NH1	2.44	0.50
1:A:133:LEU:HD12	1:A:133:LEU:N	2.26	0.50
1:A:278:GLN:HA	1:A:300:MET:HE3	1.94	0.50
1:B:139:VAL:HG21	1:B:193:MET:SD	2.51	0.50
1:B:242:SER:N	1:B:257:PHE:CD1	2.80	0.50
1:B:210:TRP:HB2	1:B:301:THR:HG23	1.93	0.50
1:B:241:PHE:HD1	1:B:242:SER:H	1.58	0.50
1:B:360:LYS:O	1:B:362:ILE:N	2.44	0.50
1:A:291:ALA:HB3	1:A:293:VAL:HG23	1.92	0.50
1:B:394:TYR:O	1:B:398:SER:HB3	2.12	0.50
1:A:284:TYR:CD2	1:A:307:HIS:CE1	3.00	0.50
1:B:182:PRO:HG2	1:B:185:VAL:HG23	1.92	0.50
1:A:115:VAL:HG11	1:A:131:LEU:HD22	1.93	0.50
1:A:473:ASP:OD1	1:A:473:ASP:C	2.53	0.50
1:B:498:ILE:HD12	1:B:498:ILE:N	2.27	0.49
1:A:219:MET:HE1	1:A:265:MET:CE	2.43	0.49
1:B:210:TRP:HB2	1:B:301:THR:CG2	2.43	0.49
1:A:96:MET:HB3	1:A:265:MET:HE2	1.93	0.49
1:B:448:HIS:HB2	1:B:450:PHE:CZ	2.47	0.49
1:B:412:ASN:OD1	1:B:424:ARG:HD2	2.13	0.49
1:A:365:VAL:HG12	1:A:368:ALA:HB2	1.91	0.48
1:A:59:THR:HG22	1:A:73:MET:CE	2.41	0.48
1:B:182:PRO:O	1:B:186:VAL:HG23	2.14	0.48
1:B:247:THR:O	1:B:257:PHE:HE1	1.96	0.48
1:A:86:MET:O	1:A:89:GLN:HB2	2.13	0.48
1:A:273:PHE:O	1:A:274:THR:HG23	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:139:VAL:HB	1:A:193:MET:CE	2.44	0.48
1:B:185:VAL:O	1:B:189:LYS:HG2	2.14	0.47
1:B:334:ARG:NH1	1:B:337:GLN:OE1	2.37	0.47
1:B:210:TRP:CE3	1:B:269:GLN:HG3	2.49	0.47
1:A:342:LYS:O	1:A:346:LEU:HD13	2.15	0.47
1:B:241:PHE:HE1	1:B:246:ASP:HA	1.80	0.47
1:B:469:PHE:HB2	1:B:498:ILE:HD13	1.95	0.47
1:A:297:ASP:O	1:A:301:THR:OG1	2.27	0.47
1:A:392:TYR:HB3	1:A:413:VAL:HG11	1.97	0.46
1:A:316:TRP:CZ2	1:A:359:PRO:HG3	2.50	0.46
1:A:57:PHE:HB2	1:A:75:TYR:CD1	2.50	0.46
1:A:412:ASN:HD22	1:A:453:PRO:HD2	1.81	0.46
1:B:271:TRP:CD1	1:B:271:TRP:N	2.83	0.46
1:A:100:ARG:O	1:A:104:ILE:HG13	2.16	0.46
1:B:339:MET:HG2	1:B:518:TYR:CZ	2.51	0.46
1:A:59:THR:CG2	1:A:73:MET:HE2	2.39	0.45
1:A:85:LYS:HG3	1:A:87:ASN:HB2	1.98	0.45
1:B:325:ASP:HA	1:B:328:LYS:HZ1	1.81	0.45
1:A:58:LYS:H	1:A:270:GLN:HE22	1.64	0.45
1:B:365:VAL:HG11	1:B:368:ALA:HB2	1.98	0.45
1:A:357:GLY:O	2:A:601:SAM:HG2	2.17	0.45
1:A:333:CYS:HA	1:A:525:MET:CE	2.47	0.45
1:A:360:LYS:O	1:A:362:ILE:N	2.50	0.45
1:B:395:ALA:N	5:B:703:HOH:O	2.43	0.44
1:B:242:SER:HB2	1:B:257:PHE:CD1	2.53	0.44
1:A:89:GLN:C	1:A:91:ASN:N	2.75	0.44
1:A:241:PHE:CE2	1:A:243:THR:HG22	2.51	0.44
1:A:241:PHE:CD1	1:A:248:TYR:CE2	3.05	0.44
1:B:210:TRP:CE2	1:B:269:GLN:NE2	2.86	0.44
1:A:57:PHE:CD2	1:A:75:TYR:HE1	2.36	0.44
1:B:217:THR:HG22	1:B:220:LYS:NZ	2.33	0.44
1:A:456:ASP:CG	1:A:458:SER:HG	2.26	0.43
1:B:212:HIS:HB2	1:B:215:GLU:OE2	2.17	0.43
1:B:498:ILE:HG21	1:B:543:LEU:HB3	1.99	0.43
1:B:219:MET:HE3	1:B:219:MET:HB2	1.86	0.43
1:A:112:GLY:H	1:A:136:SER:HB3	1.83	0.43
1:A:113:PHE:HD2	1:A:209:LEU:HD21	1.84	0.43
1:A:131:LEU:O	1:A:142:VAL:HA	2.19	0.43
1:A:365:VAL:HG12	1:A:368:ALA:CB	2.49	0.43
1:B:302:ARG:O	1:B:306:VAL:HG23	2.19	0.43
1:B:542:ASN:HA	1:B:545:ASN:HB3	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:96:MET:HB3	1:A:265:MET:HE3	2.01	0.43
1:B:456:ASP:OD2	1:B:458:SER:OG	2.37	0.43
1:A:343:ALA:HB1	1:A:543:LEU:HD21	2.01	0.43
1:A:493:LYS:HB3	1:B:541:TYR:CE2	2.53	0.43
1:B:97:PHE:HA	1:B:265:MET:O	2.19	0.43
1:B:496:THR:C	1:B:498:ILE:HD12	2.44	0.43
1:B:100:ARG:HG3	1:B:271:TRP:CZ3	2.54	0.42
1:B:325:ASP:OD1	1:B:328:LYS:HD2	2.19	0.42
1:B:491:PRO:HB2	1:B:493:LYS:HZ3	1.85	0.42
1:B:249:ALA:HB2	1:B:257:PHE:CE2	2.54	0.42
1:B:412:ASN:ND2	1:B:452:THR:HG23	2.33	0.42
1:A:338:HIS:CE1	1:A:365:VAL:HG22	2.54	0.42
1:B:347:ALA:HB2	1:B:543:LEU:HD13	2.01	0.42
1:A:82:MET:SD	1:A:216:LEU:HD23	2.59	0.42
1:B:394:TYR:CD2	1:B:415:ARG:HD2	2.55	0.42
1:A:211:ALA:O	1:A:215:GLU:OE1	2.38	0.42
1:B:542:ASN:OD1	1:B:542:ASN:N	2.52	0.42
1:A:242:SER:HB2	1:A:257:PHE:CD2	2.55	0.41
1:B:287:VAL:HG23	1:B:439:ASP:HA	2.03	0.41
1:A:100:ARG:HG3	1:A:271:TRP:CZ2	2.55	0.41
1:B:471:TYR:HD1	1:B:498:ILE:HG22	1.85	0.41
1:B:100:ARG:HG3	1:B:271:TRP:CH2	2.56	0.41
1:A:337:GLN:HB2	1:A:409:TRP:CZ2	2.55	0.41
1:B:137:THR:OG1	1:B:139:VAL:HG23	2.20	0.41
1:A:290:ASN:OD1	1:A:290:ASN:C	2.62	0.41
1:B:112:GLY:O	1:B:135:PHE:HA	2.21	0.41
1:B:403:ASP:OD1	1:B:403:ASP:N	2.52	0.41
1:A:338:HIS:ND1	1:A:365:VAL:HG22	2.35	0.41
1:A:57:PHE:HA	1:A:270:GLN:NE2	2.36	0.41
1:A:79:ILE:HG23	1:A:84:PHE:HB2	2.02	0.41
1:A:84:PHE:HB3	1:A:225:ILE:HD13	2.01	0.41
1:A:273:PHE:O	1:A:274:THR:CG2	2.69	0.41
1:B:200:ASN:C	1:B:201:LEU:HD22	2.46	0.41
1:A:192:GLN:O	1:A:196:ASP:OD1	2.40	0.40
1:A:312:LYS:H	1:A:312:LYS:HG3	1.68	0.40
1:B:216:LEU:HD12	1:B:216:LEU:N	2.36	0.40
1:B:440:GLY:C	1:B:454:ALA:HB2	2.47	0.40

All (2) 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:A:274:THR:OG1	1:B:228:GLU:O[2_455]	2.16	0.04
1:A:308:GLU:OE2	1:B:229:ARG:HH22[2_455]	1.56	0.04

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	418/531 (79%)	391 (94%)	25 (6%)	2 (0%)	24	55
1	B	368/531 (69%)	348 (95%)	17 (5%)	3 (1%)	16	44
All	All	786/1062 (74%)	739 (94%)	42 (5%)	5 (1%)	21	51

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	212	HIS
1	A	90	VAL
1	B	212	HIS
1	B	211	ALA
1	B	324	GLY

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	377/465 (81%)	374 (99%)	3 (1%)	73	90
1	B	337/465 (72%)	336 (100%)	1 (0%)	86	95

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
All	All	714/930 (77%)	710 (99%)	4 (1%)	78	92

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	89	GLN
1	A	136	SER
1	A	348	ASP
1	B	382	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	91	ASN
1	A	252	HIS
1	A	412	ASN
1	B	330	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 11 ligands modelled in this entry, 9 are monoatomic - leaving 2 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the

expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	SAM	A	601	-	27,29,29	1.21	1 (3%)	34,42,42	0.95	2 (5%)
2	SAM	B	601	-	27,29,29	1.30	2 (7%)	34,42,42	0.83	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
2	SAM	A	601	-	-	2/17/33/33	0/3/3/3
2	SAM	B	601	-	-	5/17/33/33	0/3/3/3

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	601	SAM	CA-N	-5.11	1.22	1.48
2	B	601	SAM	CA-N	-5.10	1.22	1.48
2	B	601	SAM	CB-CA	-2.15	1.48	1.53

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	601	SAM	CG-SD-C5'	-2.19	98.08	103.43
2	A	601	SAM	OXT-C-O	2.04	128.70	124.08

There are no chirality outliers.

All (7) torsion outliers are listed below:

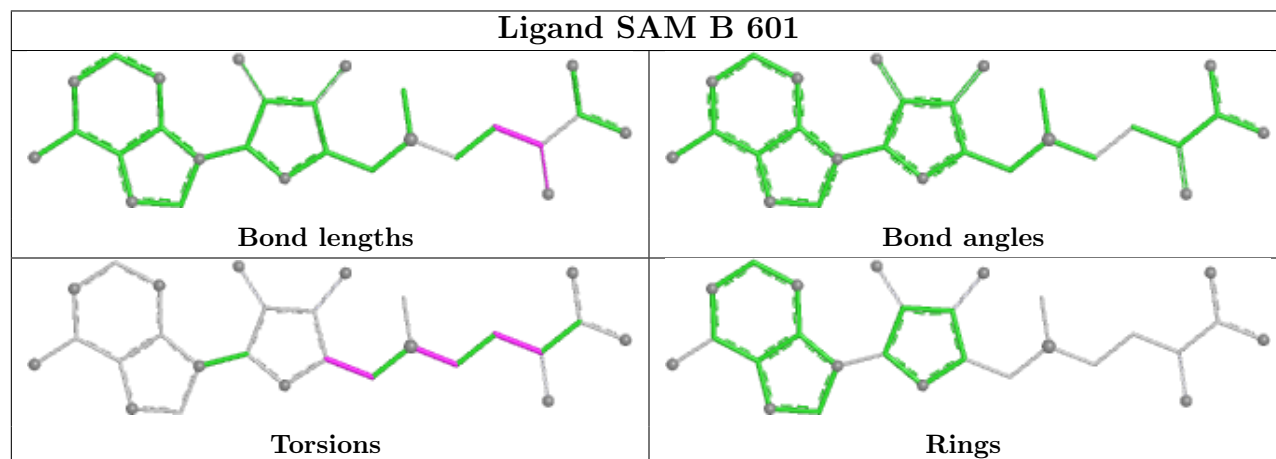
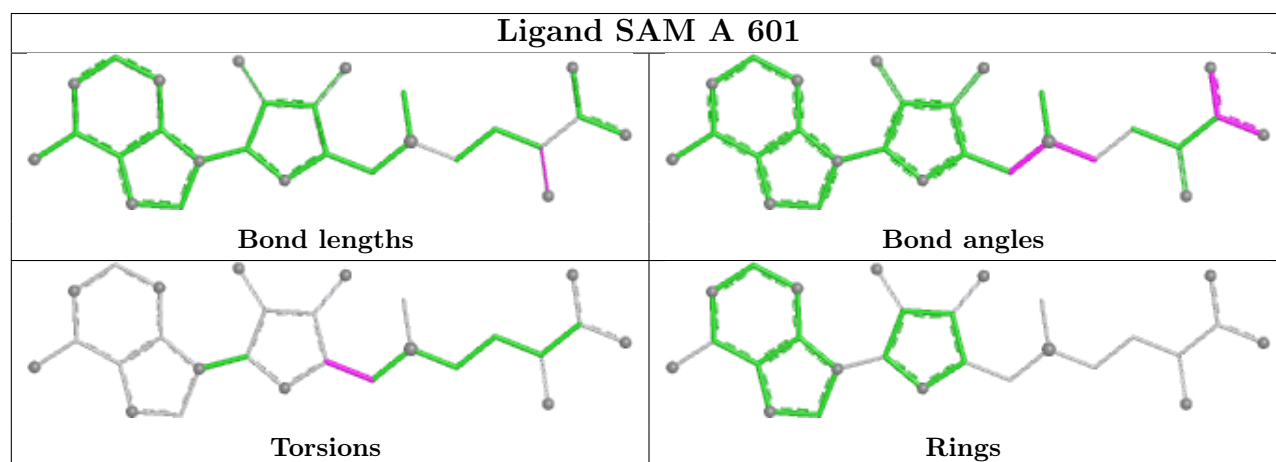
Mol	Chain	Res	Type	Atoms
2	B	601	SAM	CB-CG-SD-CE
2	B	601	SAM	CB-CG-SD-C5'
2	A	601	SAM	O4'-C4'-C5'-SD
2	B	601	SAM	O4'-C4'-C5'-SD
2	A	601	SAM	C3'-C4'-C5'-SD
2	B	601	SAM	C3'-C4'-C5'-SD
2	B	601	SAM	C-CA-CB-CG

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	601	SAM	1	0
2	B	601	SAM	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ > 2			OWAB(Å ²)	Q < 0.9
1	A	428/531 (80%)	0.29	19 (4%)	39	31	15, 35, 150, 196	6 (1%)
1	B	380/531 (71%)	0.45	24 (6%)	26	19	18, 41, 72, 130	10 (2%)
All	All	808/1062 (76%)	0.37	43 (5%)	32	24	15, 38, 91, 196	16 (1%)

All (43) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	304	LEU	4.5
1	A	209	LEU	3.9
1	B	268	VAL	3.9
1	A	92	GLY	3.8
1	A	93	TYR	3.8
1	B	94	PRO	3.6
1	B	216	LEU	3.2
1	A	271	TRP	2.9
1	A	61	GLY	2.8
1	B	369	ASP	2.7
1	B	277	LEU	2.7
1	B	210	TRP	2.7
1	B	183	TRP	2.6
1	B	246	ASP	2.6
1	B	270	GLN	2.6
1	B	209	LEU	2.6
1	B	296	CYS	2.6
1	B	293	VAL	2.6
1	A	268	VAL	2.5
1	A	131	LEU	2.5
1	A	101	GLU	2.5
1	A	214	PHE	2.4
1	A	289	GLY	2.4
1	A	51	LEU	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	133	LEU	2.4
1	A	82	MET	2.4
1	B	283	LEU	2.3
1	B	287	VAL	2.2
1	B	299	ILE	2.2
1	B	106	HIS	2.1
1	A	97	PHE	2.1
1	A	491	PRO	2.1
1	B	208	VAL	2.1
1	A	340	VAL	2.1
1	A	507	VAL	2.1
1	A	216	LEU	2.1
1	A	91	ASN	2.1
1	B	280	ASN	2.1
1	B	190	ILE	2.1
1	B	207	PHE	2.0
1	A	96	MET	2.0
1	B	96	MET	2.0
1	B	213	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
2	SAM	A	601	27/27	0.90	0.09	58,79,98,104	21
2	SAM	B	601	27/27	0.91	0.09	55,68,81,86	21
3	ZN	B	605	1/1	0.93	0.06	139,139,139,139	0
3	ZN	B	603	1/1	0.95	0.06	129,129,129,129	0

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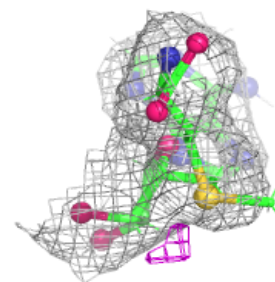
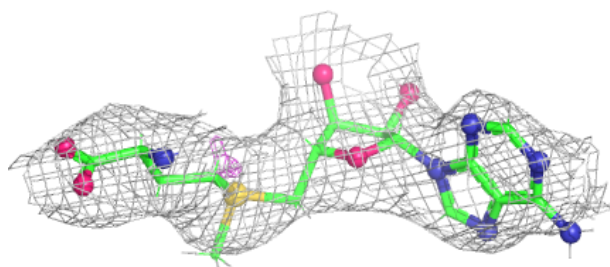
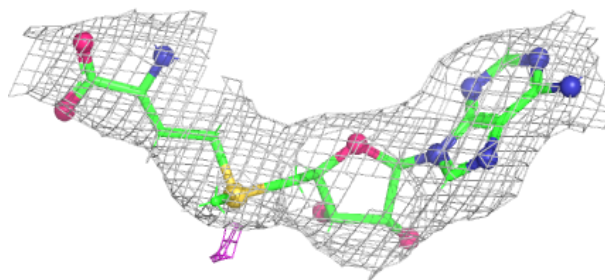
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	MN	A	606	1/1	0.96	0.21	68,68,68,68	0
3	ZN	A	605	1/1	0.99	0.03	100,100,100,100	0
3	ZN	A	602	1/1	0.99	0.04	54,54,54,54	0
3	ZN	B	604	1/1	0.99	0.05	98,98,98,98	0
3	ZN	A	603	1/1	0.99	0.15	87,87,87,87	0
3	ZN	A	604	1/1	0.99	0.06	103,103,103,103	0
3	ZN	B	602	1/1	1.00	0.02	74,74,74,74	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.

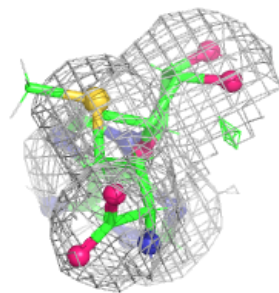
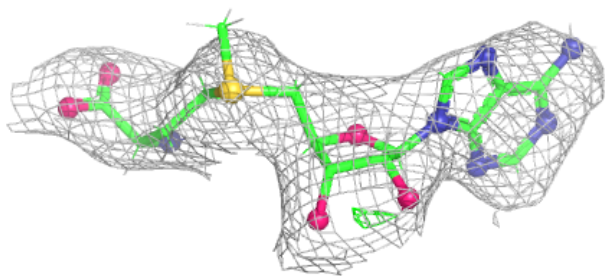
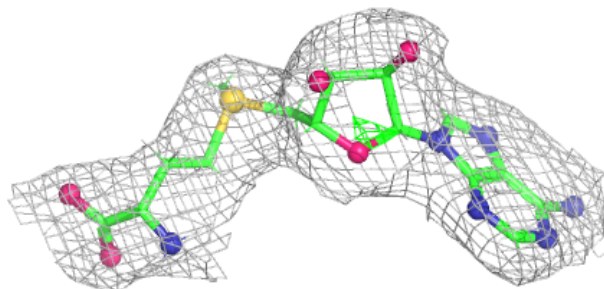
Electron density around SAM A 601:

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



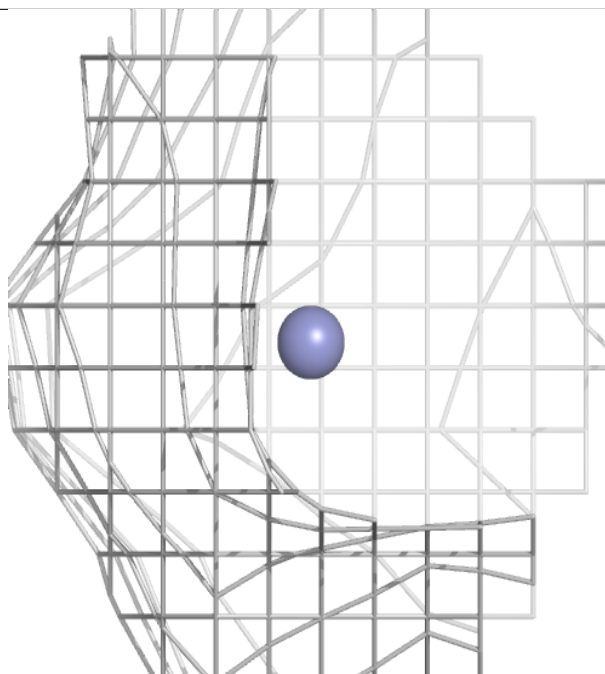
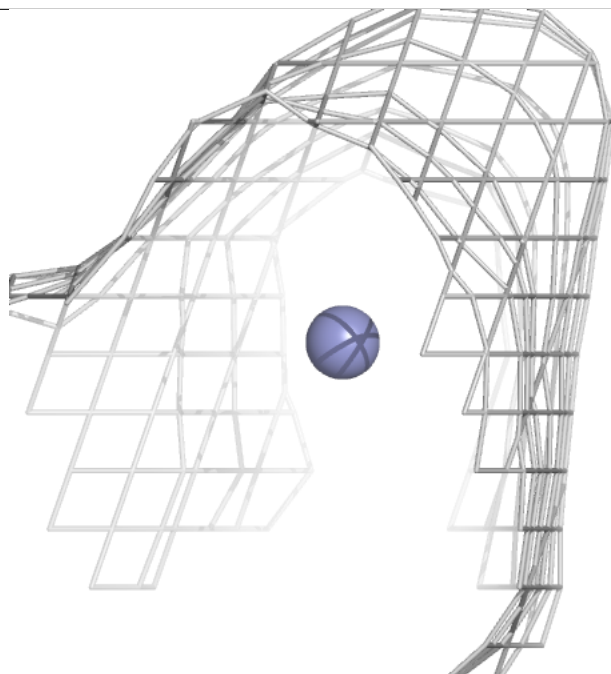
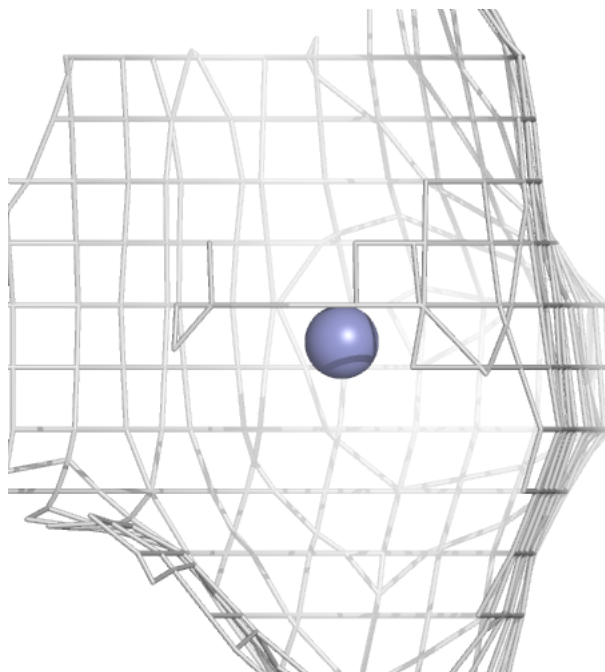
Electron density around SAM B 601:

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



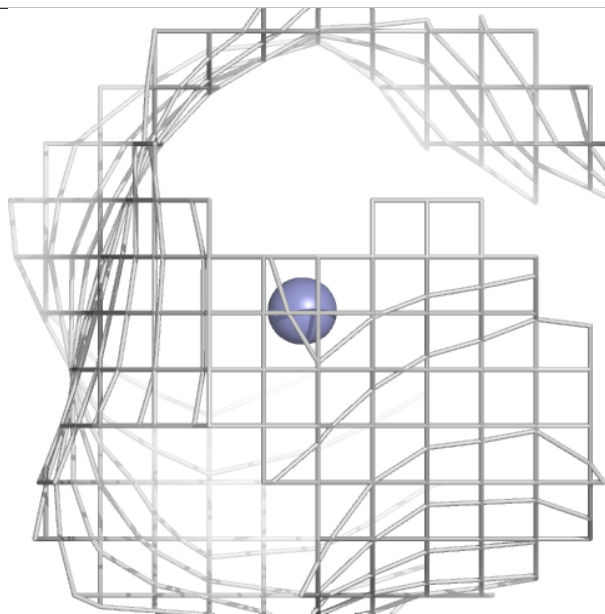
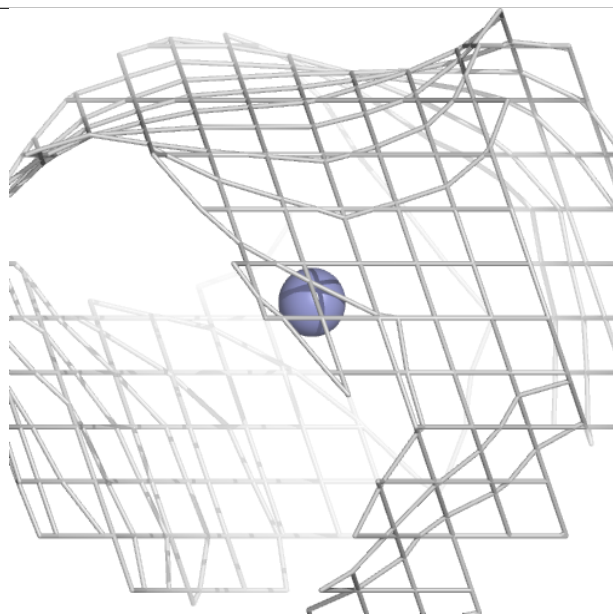
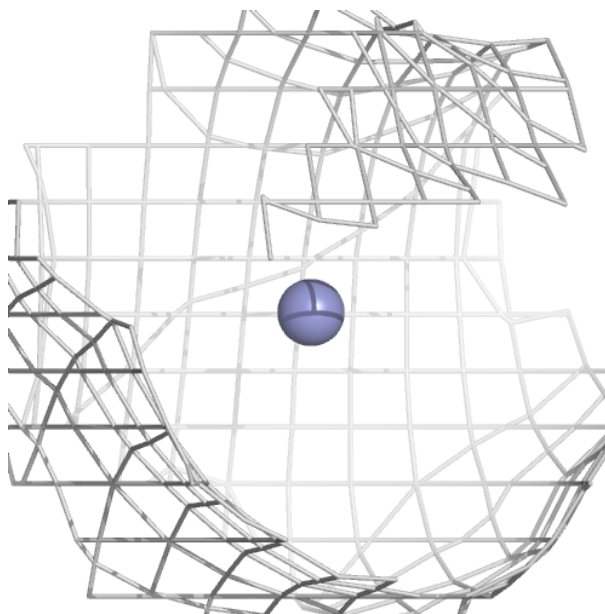
Electron density around ZN B 605:

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



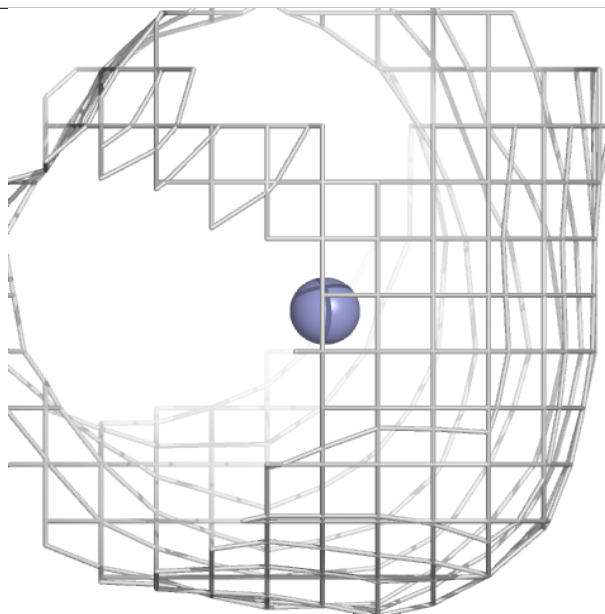
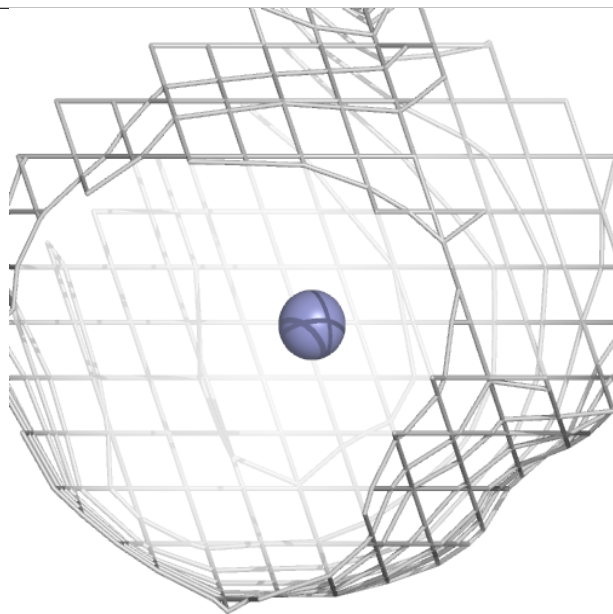
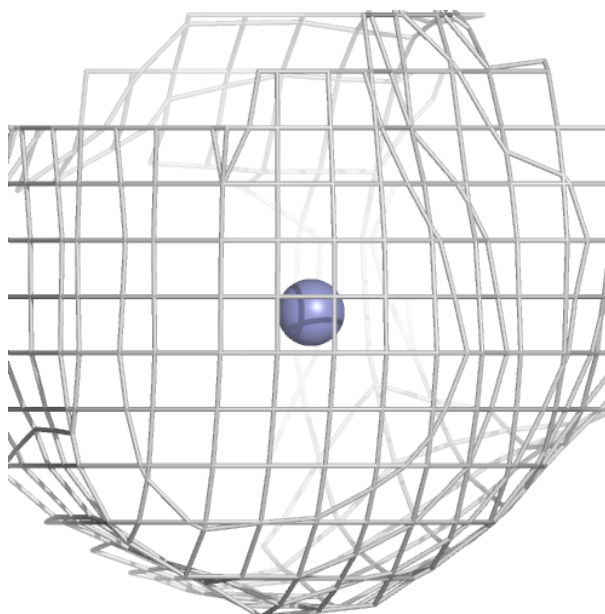
Electron density around ZN B 603:

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



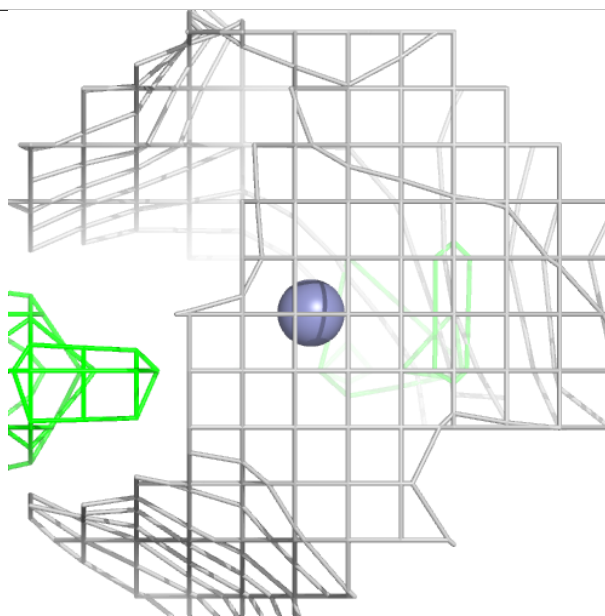
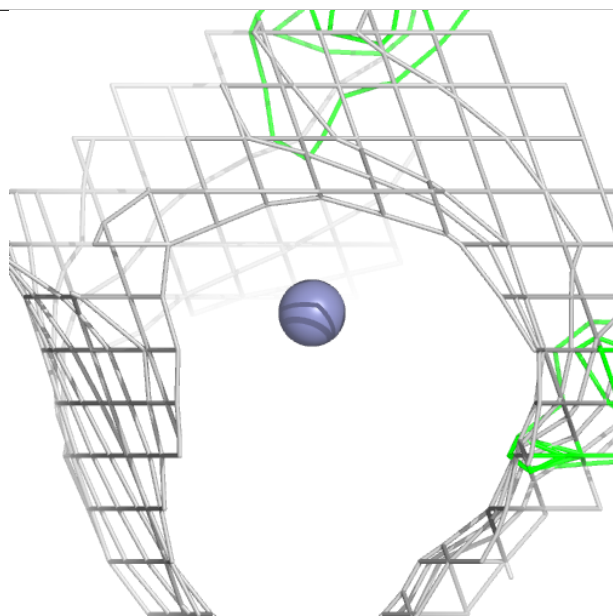
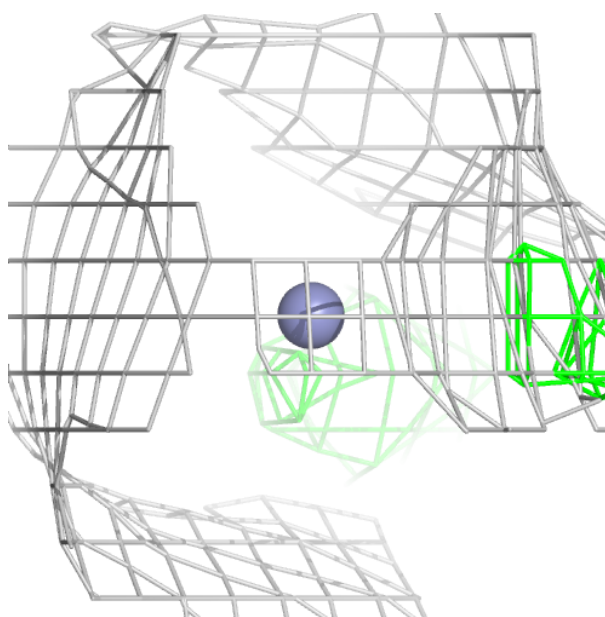
Electron density around ZN A 605:

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



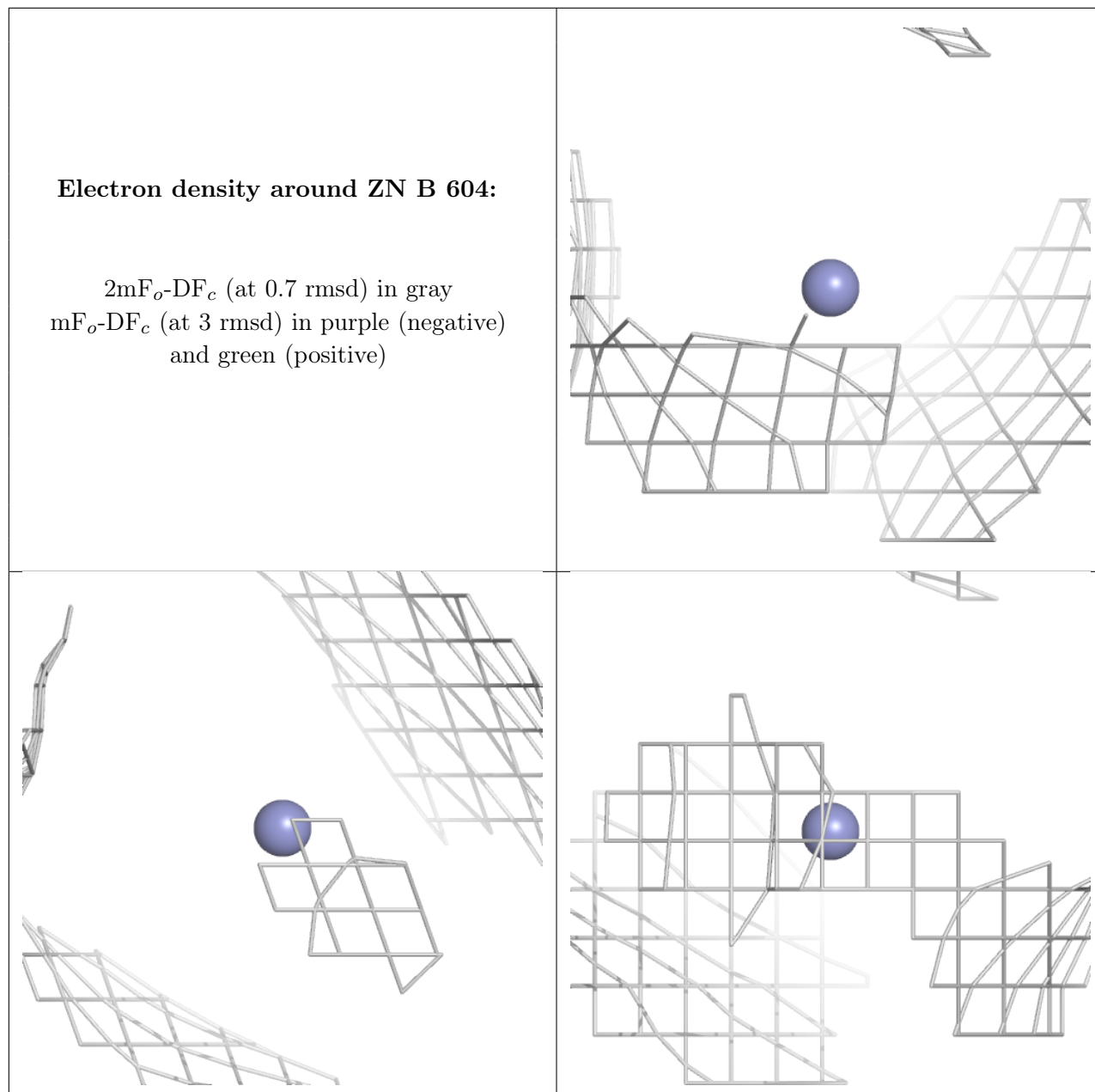
Electron density around ZN A 602:

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and green (positive)



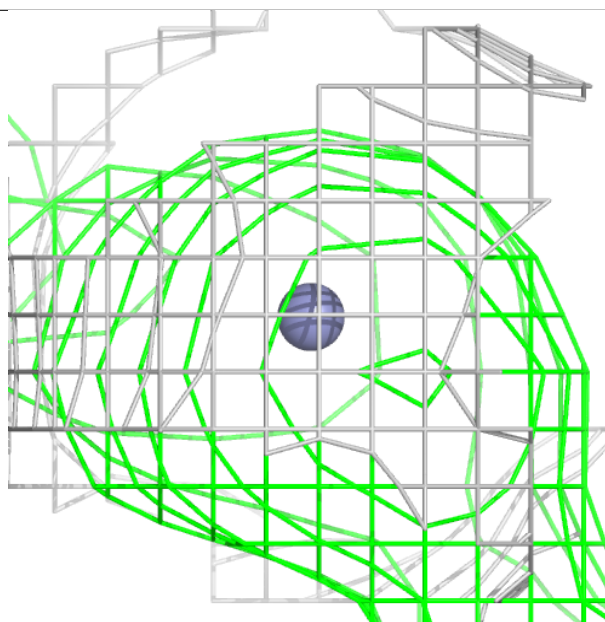
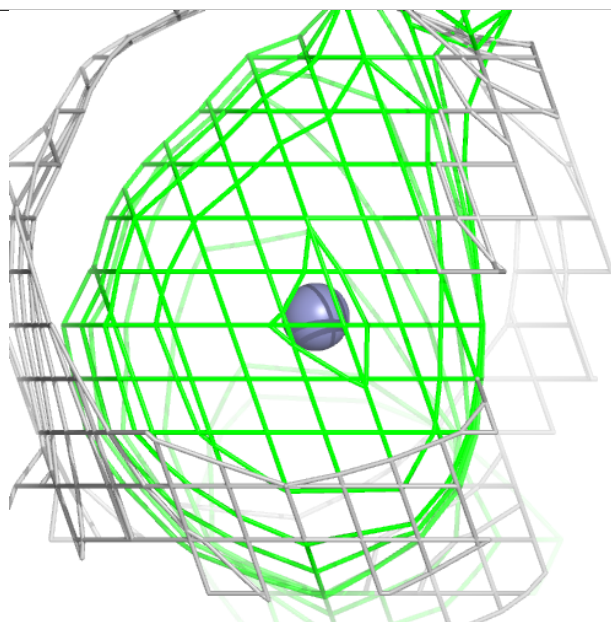
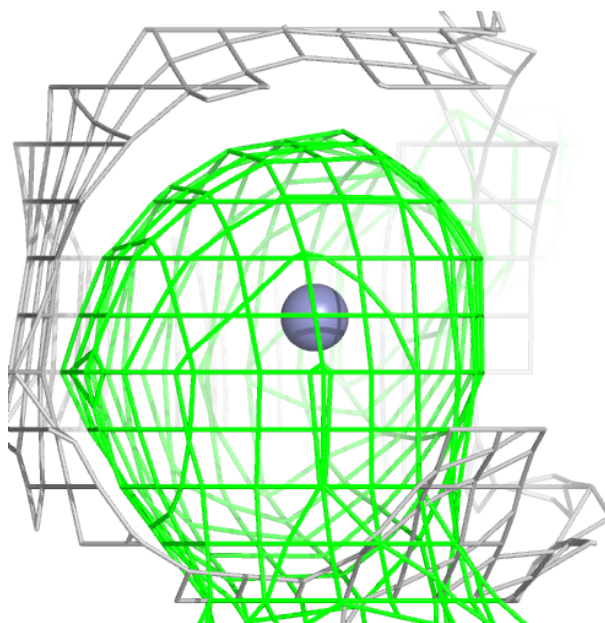
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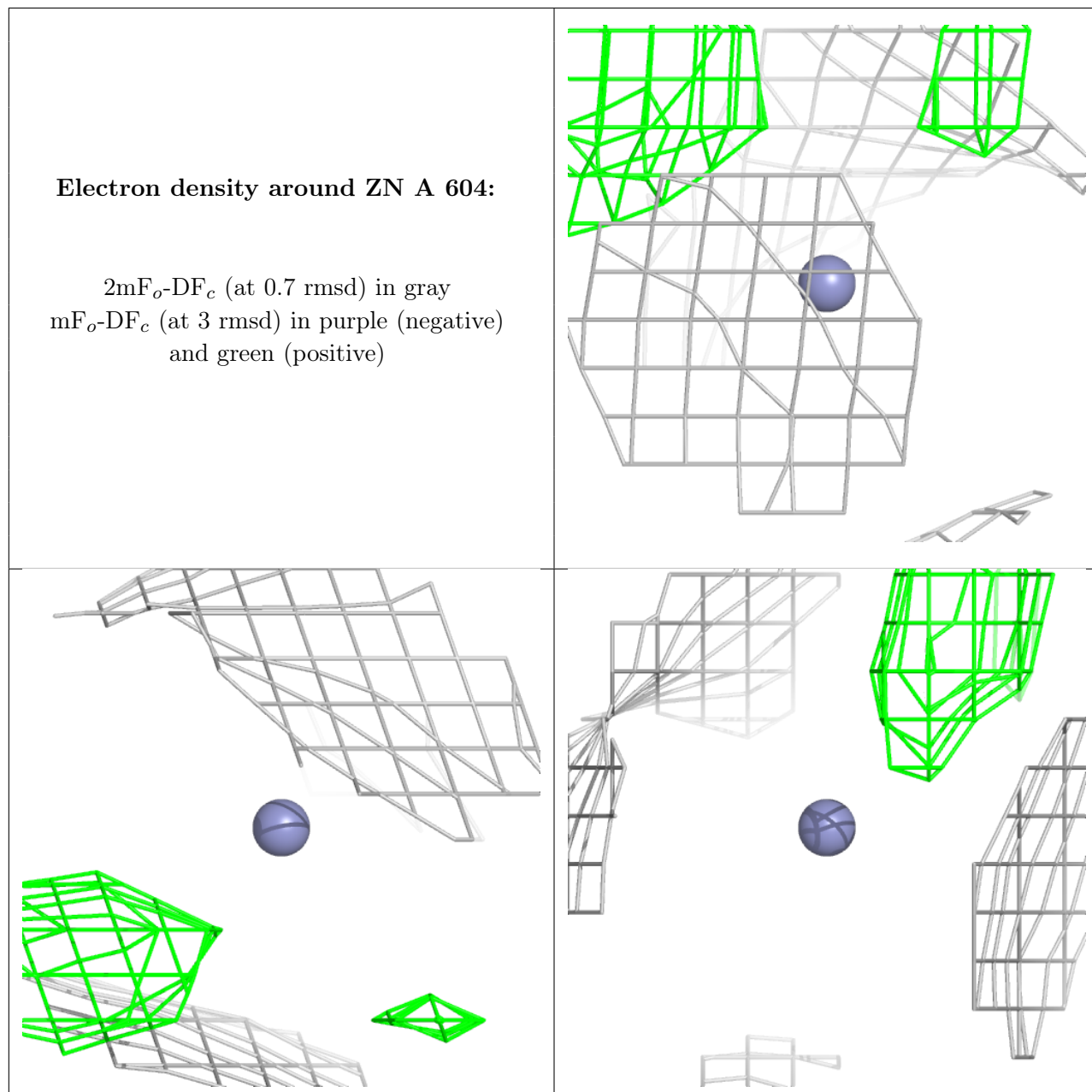
Electron density around ZN A 603:

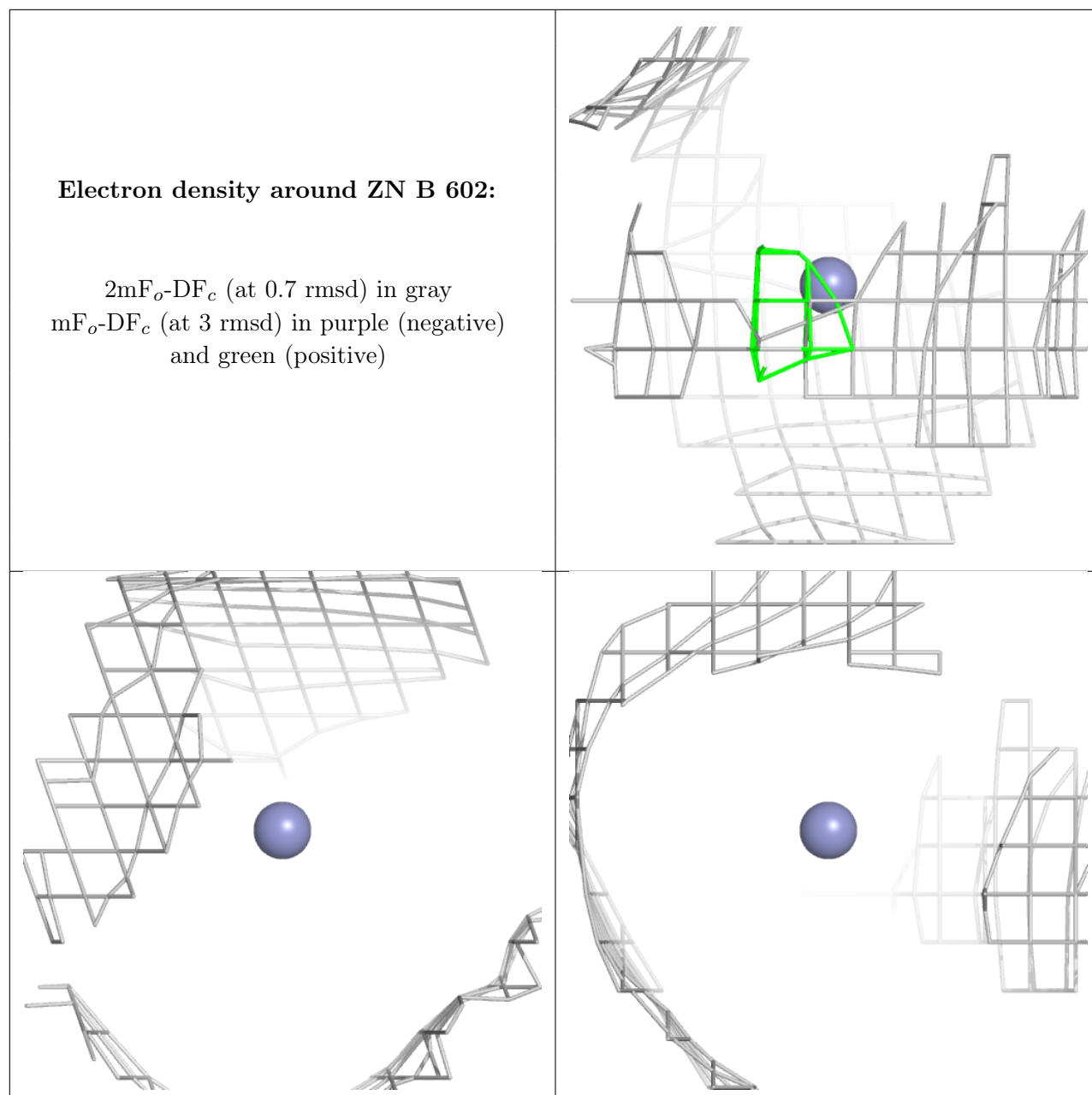
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and green (positive)



Electron density around ZN A 604:

$2mF_o-DF_c$ (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.