



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 03:39 PM UTC

PDB ID : 8IIB / pdb_00008iib
Title : Crystal structure of Israeli acute paralysis virus RNA-dependent RNA polymerase delta85 mutant (residues 86-546)
Authors : Fang, X.; Lu, G.; Hou, C.; Gong, P.
Deposited on : 2023-02-24
Resolution : 2.69 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
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

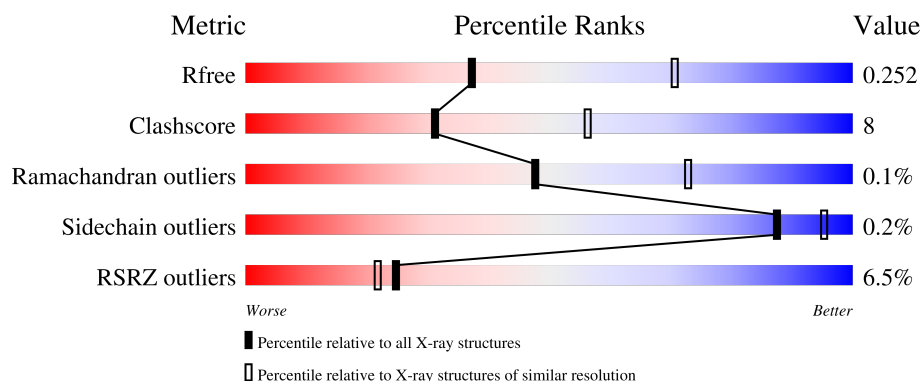
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.69 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	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

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	475	 4% 60% 15% 25%
1	B	475	 5% 65% 9% 26%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 5527 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Polymerase polypeptide.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	355	Total	C	N	O	S	0	1	0
			2719	1730	455	508	26			
1	B	351	Total	C	N	O	S	0	0	0
			2629	1673	446	485	25			

- Molecule 2 is CADMIUM ION (CCD ID: CD) (formula: Cd) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	3	Total	Cd	0	0
			3	3		
2	B	1	Total	Cd	0	0
			1	1		

- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	2	Total	Mg	0	0
			2	2		
3	B	6	Total	Mg	0	0
			6	6		

- Molecule 4 is CHLORIDE ION (CCD ID: CL) (formula: Cl) (labeled as "Ligand of Interest" by depositor).

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

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	81	Total 81	O 81	0	0
5	B	79	Total 79	O 79	0	0



4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	150.51Å 150.51Å 104.63Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	23.41 – 2.69 23.41 – 2.69	Depositor EDS
% Data completeness (in resolution range)	97.4 (23.41-2.69) 97.2 (23.41-2.69)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.71 (at 2.67Å)	Xtriage
Refinement program	PHENIX 1.17	Depositor
R, R_{free}	0.213 , 0.254 0.214 , 0.252	Depositor DCC
R_{free} test set	1819 reflections (4.77%)	wwPDB-VP
Wilson B-factor (Å ²)	37.1	Xtriage
Anisotropy	0.105	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 57.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.024 for -h,-k,l	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	5527	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.30% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: CL, MG, CD

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	0/2770	0.61	0/3754
1	B	0.34	0/2679	0.54	0/3631
All	All	0.36	0/5449	0.58	0/7385

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

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	2719	0	2458	53	0
1	B	2629	0	2330	31	0
2	A	3	0	0	0	0
2	B	1	0	0	0	0
3	A	2	0	0	0	0
3	B	6	0	0	0	0
4	A	4	0	0	0	0
4	B	3	0	0	0	0
5	A	81	0	0	2	0
5	B	79	0	0	2	0
All	All	5527	0	4788	83	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 8.

All (83) 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:188:LEU:HD13	1:A:188:LEU:O	1.74	0.88
1:A:421:PHE:CD1	1:A:426:PHE:HB2	2.09	0.87
1:B:197:ALA:O	1:B:321:LEU:CD1	2.24	0.86
1:A:148:SER:O	1:A:149:ILE:HG13	1.79	0.82
1:B:238:ILE:HG22	1:B:242:MET:HE2	1.60	0.82
1:A:107:MET:HE1	1:A:372:ALA:HB3	1.62	0.82
1:A:107:MET:CE	1:A:372:ALA:HB3	2.15	0.77
1:A:421:PHE:HD1	1:A:426:PHE:HB2	1.51	0.75
1:B:134:GLU:OE2	1:B:187:ARG:NH2	2.21	0.73
1:B:190:VAL:HG11	1:B:242:MET:HE1	1.72	0.71
1:A:421:PHE:HE1	1:A:426:PHE:CD2	2.13	0.67
1:B:197:ALA:O	1:B:321:LEU:HD12	1.96	0.66
1:A:421:PHE:CE1	1:A:426:PHE:CD2	2.86	0.64
1:B:541:PHE:CD1	1:B:545:TYR:HB2	2.35	0.62
1:A:500:GLU:HG3	5:A:727:HOH:O	2.00	0.60
1:B:137:ILE:HD13	1:B:181:ILE:HD13	1.84	0.59
1:A:257:THR:HG23	1:A:273:LYS:HE2	1.83	0.59
1:A:107:MET:CE	1:A:372:ALA:CB	2.81	0.59
1:A:107:MET:HE1	1:A:372:ALA:CB	2.32	0.58
1:A:382:MET:HE1	1:A:390:HIS:NE2	2.18	0.58
1:A:327:LEU:HD12	1:A:330:VAL:HG21	1.84	0.57
1:A:421:PHE:HE1	1:A:426:PHE:CG	2.22	0.57
1:A:330:VAL:HG13	1:A:354:ALA:HB1	1.86	0.56
1:A:188:LEU:O	1:A:192:GLN:HG2	2.06	0.55
1:B:239:ALA:HA	1:B:242:MET:HE3	1.88	0.55
1:A:315:ASP:CG	1:A:319:ASN:HD22	2.13	0.55
1:B:383:THR:HG23	1:B:385:LYS:H	1.71	0.54
1:A:367:LEU:HD12	1:A:393:MET:HE1	1.87	0.54
1:A:136:ALA:HB2	1:A:242:MET:HG3	1.91	0.53
1:B:502:SER:OG	1:B:540:ARG:NH2	2.42	0.52
1:A:395:SER:HB3	5:A:736:HOH:O	2.10	0.52
1:A:491:LYS:NZ	1:A:529:ASP:OD2	2.43	0.51
1:A:541:PHE:HD1	1:A:542:LEU:HD12	1.75	0.51
1:A:124:ARG:NH2	1:A:253:GLU:OE1	2.43	0.51
1:A:198:ARG:HG3	1:A:322:ILE:HD11	1.93	0.51
1:B:190:VAL:CG1	1:B:242:MET:HE1	2.37	0.51
1:B:540:ARG:HD2	5:B:717:HOH:O	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:330:VAL:HG12	1:A:355:THR:HG22	1.94	0.51
1:B:301:ILE:HD11	1:B:369:MET:HE1	1.92	0.51
1:A:421:PHE:CE1	1:A:426:PHE:CG	3.00	0.50
1:B:538:GLN:OE1	1:B:542:LEU:HD11	2.12	0.50
1:A:145:TYR:CD1	1:A:252:MET:HE2	2.47	0.49
1:A:182:LEU:O	1:A:186:VAL:HG22	2.12	0.49
1:B:454:LYS:O	1:B:467:PRO:HD2	2.11	0.49
1:A:369:MET:O	1:A:373:ILE:HG22	2.12	0.49
1:A:468:LEU:HD23	1:A:500:GLU:HG2	1.93	0.49
1:B:109:ASP:O	1:B:113:GLN:HG2	2.12	0.49
1:B:281:PHE:O	1:B:405:SER:HA	2.12	0.49
1:A:486:ILE:O	1:A:486:ILE:HG13	2.12	0.48
1:A:281:PHE:O	1:A:405:SER:HA	2.13	0.48
1:A:131:LEU:HD11	1:A:245:LEU:HB3	1.96	0.48
1:A:107:MET:HE2	1:A:372:ALA:HB3	1.95	0.47
1:B:197:ALA:O	1:B:321:LEU:HD13	2.10	0.47
1:A:133:TYR:OH	1:A:186:VAL:HG23	2.13	0.47
1:A:479:TRP:CG	1:B:474:LEU:HD12	2.50	0.47
1:B:320:ASN:N	1:B:320:ASN:HD22	2.12	0.47
1:A:321:LEU:O	1:A:325:VAL:HG23	2.15	0.47
1:A:107:MET:HE2	1:A:384:MET:CE	2.44	0.46
1:B:123:MET:HG2	1:B:312:PHE:CE2	2.51	0.46
1:A:202:ARG:HB3	1:A:325:VAL:HG13	1.97	0.45
1:A:298:ASN:ND2	1:A:301:ILE:HD12	2.32	0.45
1:B:123:MET:HE3	1:B:123:MET:HB2	1.77	0.45
1:A:278:LEU:HD13	1:A:401:VAL:HG11	2.00	0.44
1:A:298:ASN:CG	1:A:301:ILE:HD12	2.42	0.44
1:B:520:TYR:CE1	1:B:526:ASP:HB3	2.52	0.44
1:B:310:ASN:OD1	1:B:323:ARG:HD2	2.17	0.44
1:B:248:ILE:HG22	1:B:252:MET:HE3	1.99	0.44
1:A:145:TYR:CZ	1:A:255:ARG:NH2	2.87	0.43
1:B:142:VAL:CG2	1:B:146:ILE:HD12	2.49	0.43
1:A:193:ARG:NE	1:A:235:ASP:OD1	2.48	0.42
1:A:188:LEU:HD13	1:A:188:LEU:C	2.43	0.42
1:B:413:ASN:HB2	1:B:441:LYS:O	2.20	0.42
1:B:505:GLU:HG3	5:B:751:HOH:O	2.19	0.42
1:A:474:LEU:HD23	1:A:474:LEU:HA	1.88	0.42
1:A:137:ILE:HD11	1:A:238:ILE:HD13	2.02	0.41
1:A:236:PHE:CE1	1:A:329:ASP:HB2	2.56	0.41
1:A:526:ASP:CG	1:A:527:HIS:H	2.29	0.41
1:B:521:ALA:HA	1:B:526:ASP:O	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:248:ILE:HG22	1:B:252:MET:CE	2.51	0.41
1:A:431[B]:GLU:H	1:A:431[B]:GLU:HG3	1.27	0.41
1:B:393:MET:HE3	1:B:393:MET:HB2	1.82	0.41
1:A:465:GLU:OE2	1:A:540:ARG:NH2	2.48	0.40
1:A:421:PHE:CE1	1:A:426:PHE:HB2	2.53	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	344/475 (72%)	330 (96%)	13 (4%)	1 (0%)	36	60
1	B	337/475 (71%)	326 (97%)	11 (3%)	0	100	100
All	All	681/950 (72%)	656 (96%)	24 (4%)	1 (0%)	48	73

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	203	LEU

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.

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	262/420 (62%)	261 (100%)	1 (0%)	84	93
1	B	242/420 (58%)	242 (100%)	0	100	100
All	All	504/840 (60%)	503 (100%)	1 (0%)	87	95

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

Mol	Chain	Res	Type
1	A	183	ASN

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

Mol	Chain	Res	Type
1	A	191	GLN
1	A	192	GLN
1	A	195	GLN
1	A	319	ASN
1	A	324	HIS
1	A	377	ASN
1	A	449	GLN
1	B	320	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 19 ligands modelled in this entry, 19 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	355/475 (74%)	0.19	20 (5%)	30 26	15, 37, 61, 81	1 (0%)
1	B	351/475 (73%)	0.33	26 (7%)	20 18	16, 45, 68, 79	0
All	All	706/950 (74%)	0.26	46 (6%)	25 22	15, 40, 65, 81	1 (0%)

All (46) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	198	ARG	4.4
1	B	314	ASP	4.1
1	B	296	SER	4.0
1	B	200	GLY	4.0
1	B	205	VAL	3.7
1	B	149	ILE	3.6
1	A	421	PHE	3.5
1	A	376	LYS	3.5
1	B	107	MET	3.5
1	A	431[A]	GLU	3.2
1	B	292	THR	3.2
1	A	314	ASP	3.1
1	A	149	ILE	3.0
1	B	110	GLU	3.0
1	A	373	ILE	2.8
1	B	297	LEU	2.8
1	B	283	ASN	2.8
1	A	107	MET	2.8
1	A	442	TRP	2.6
1	A	297	LEU	2.6
1	A	204	PRO	2.5
1	B	353	PRO	2.5
1	A	331	TYR	2.5
1	B	106	ASP	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	375	ALA	2.4
1	B	183	ASN	2.3
1	B	315	ASP	2.3
1	A	353	PRO	2.3
1	B	487	GLN	2.3
1	A	106	ASP	2.3
1	A	233	PRO	2.2
1	B	378	ALA	2.2
1	B	233	PRO	2.2
1	A	423	THR	2.2
1	B	204	PRO	2.2
1	B	232	GLY	2.1
1	A	292	THR	2.1
1	B	538	GLN	2.1
1	A	432	LEU	2.1
1	B	322	ILE	2.1
1	B	430	ASP	2.1
1	B	352	ASN	2.1
1	A	485	ASP	2.1
1	A	105	LYS	2.0
1	A	144	GLU	2.0
1	B	423	THR	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	MG	A	605	1/1	0.81	0.19	54,54,54,54	0

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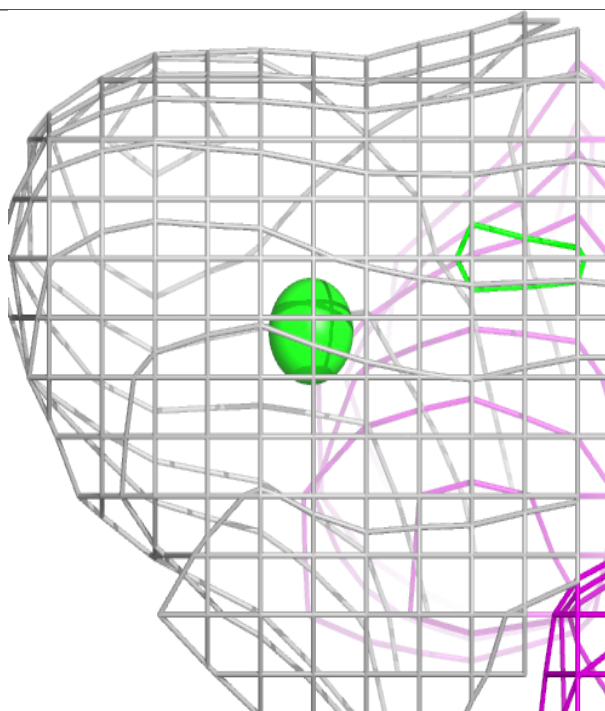
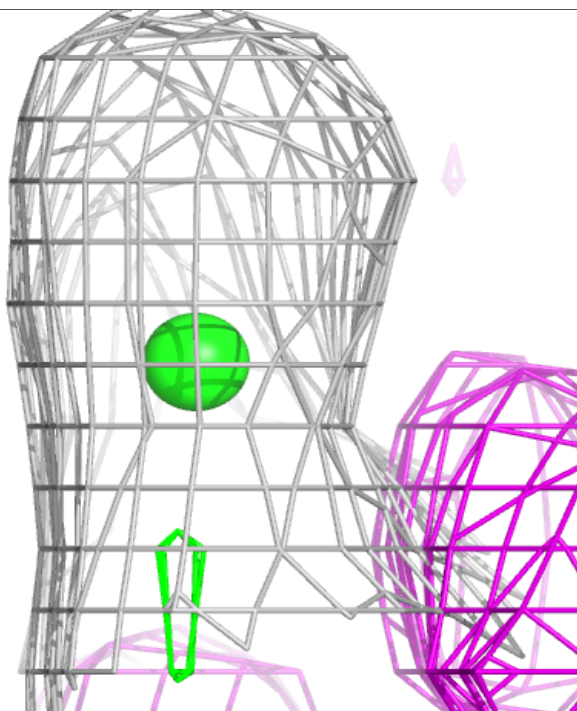
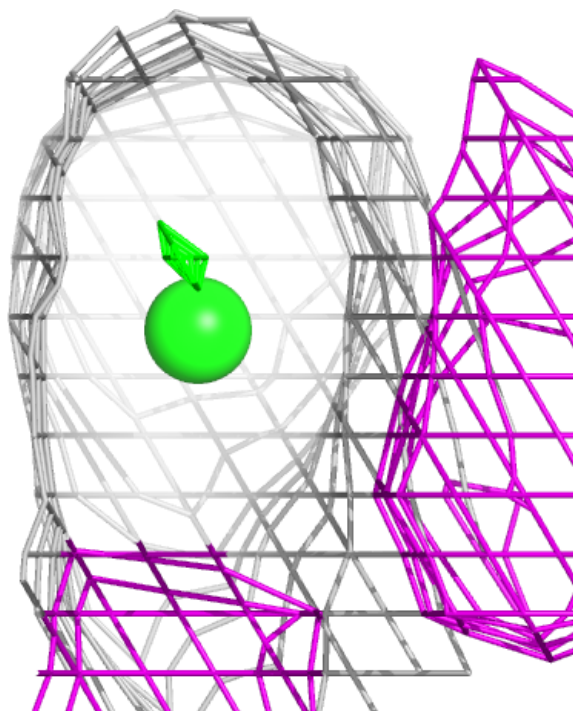
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	MG	B	606	1/1	0.85	0.20	52,52,52,52	0
4	CL	B	610	1/1	0.87	0.15	75,75,75,75	0
3	MG	A	604	1/1	0.89	0.11	49,49,49,49	0
4	CL	A	609	1/1	0.89	0.12	75,75,75,75	0
3	MG	B	604	1/1	0.89	0.19	48,48,48,48	0
3	MG	B	603	1/1	0.90	0.09	57,57,57,57	0
2	CD	B	601	1/1	0.92	0.09	77,77,77,77	0
3	MG	B	605	1/1	0.93	0.06	41,41,41,41	0
3	MG	B	602	1/1	0.95	0.11	50,50,50,50	0
4	CL	B	609	1/1	0.96	0.07	59,59,59,59	0
4	CL	A	607	1/1	0.97	0.08	30,30,30,30	0
2	CD	A	603	1/1	0.97	0.05	76,76,76,76	0
4	CL	A	606	1/1	0.98	0.07	30,30,30,30	0
4	CL	B	608	1/1	0.98	0.07	30,30,30,30	0
3	MG	B	607	1/1	0.99	0.06	52,52,52,52	0
2	CD	A	602	1/1	0.99	0.03	24,24,24,24	0
2	CD	A	601	1/1	0.99	0.04	29,29,29,29	0
4	CL	A	608	1/1	0.99	0.05	24,24,24,24	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.

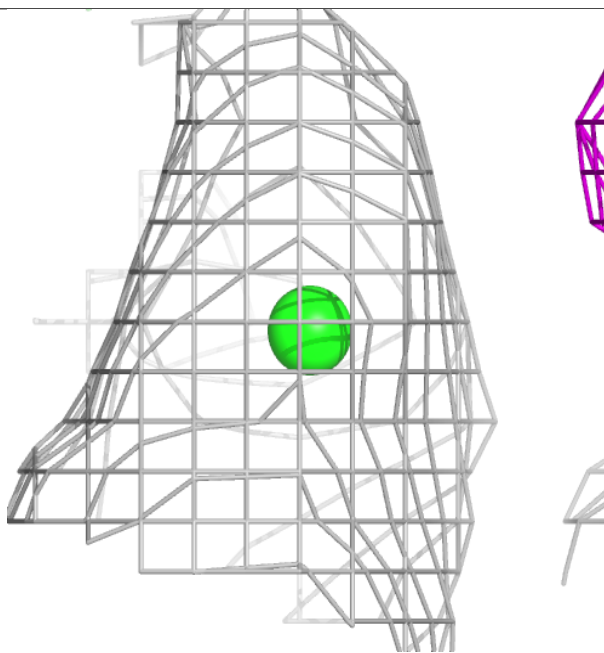
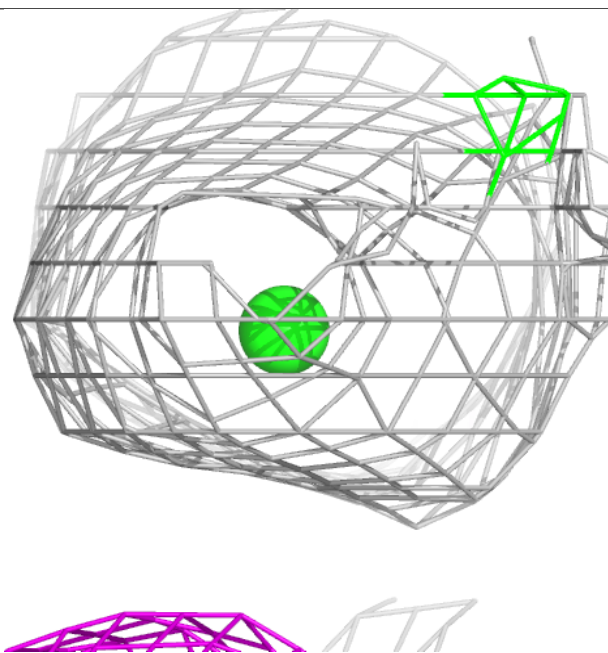
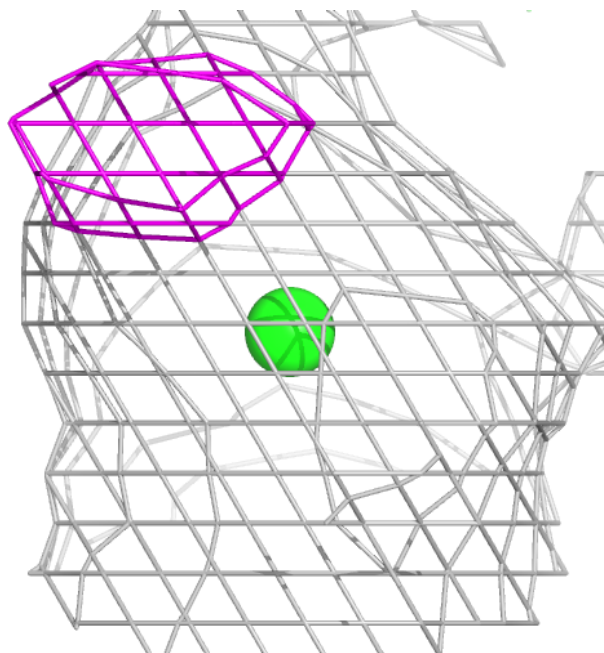
Electron density around CL B 610:

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



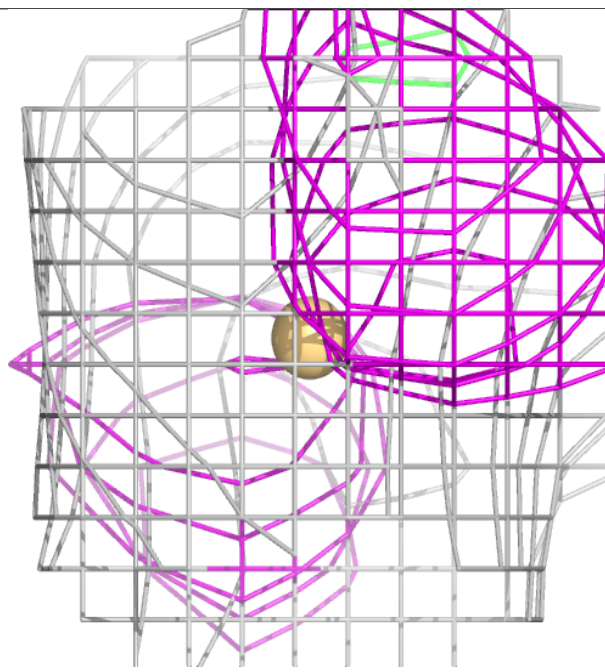
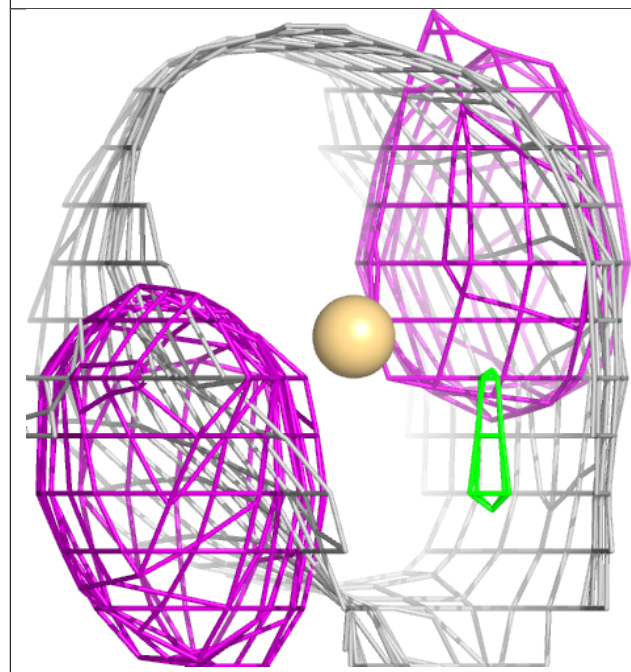
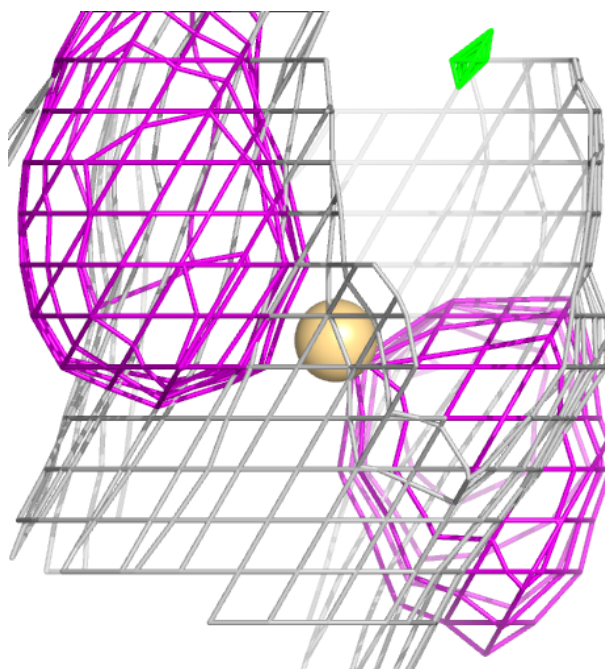
Electron density around CL A 609:

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



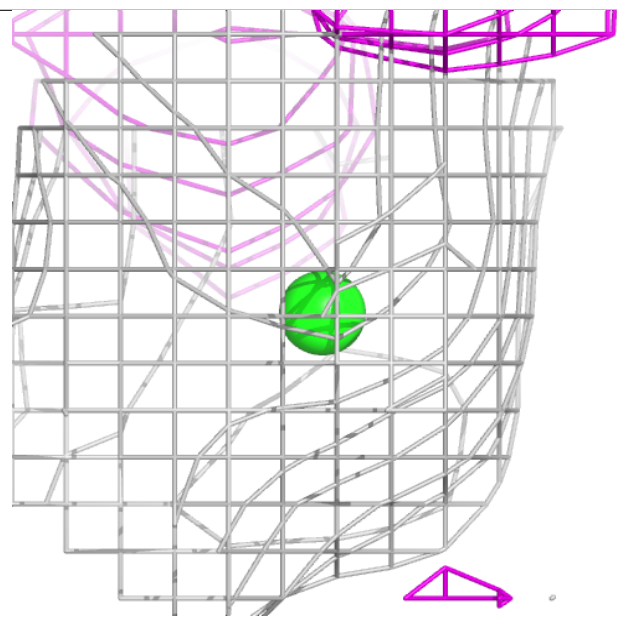
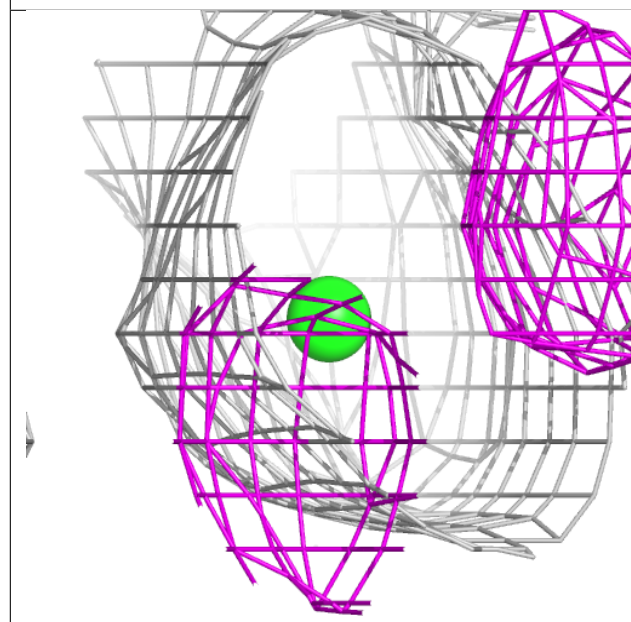
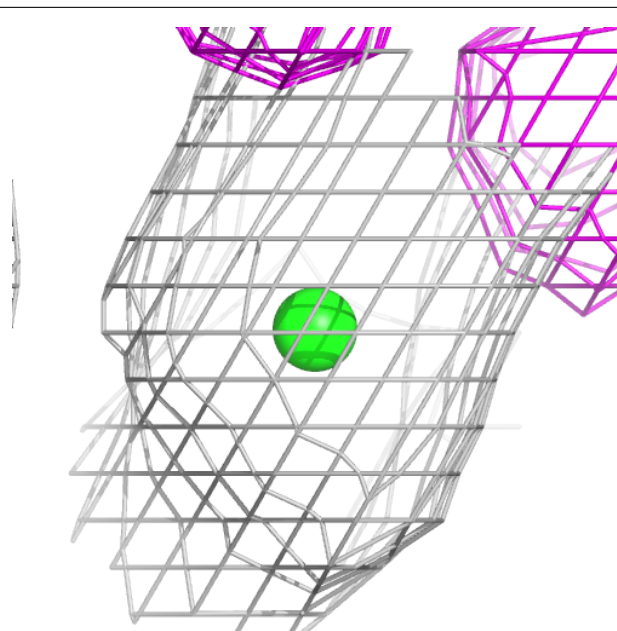
Electron density around CD B 601:

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



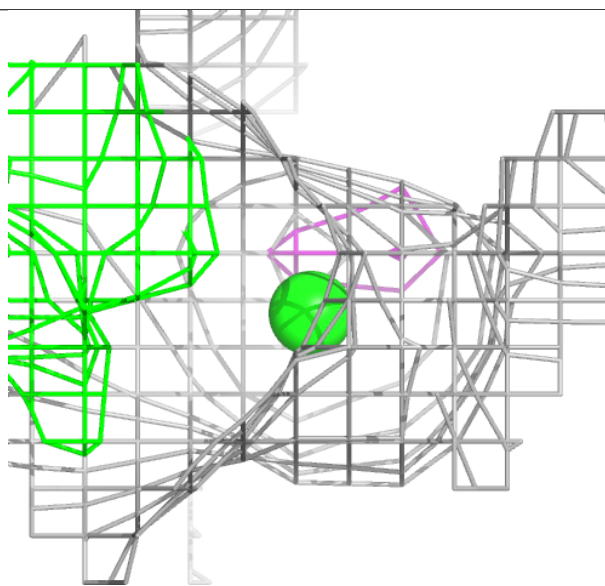
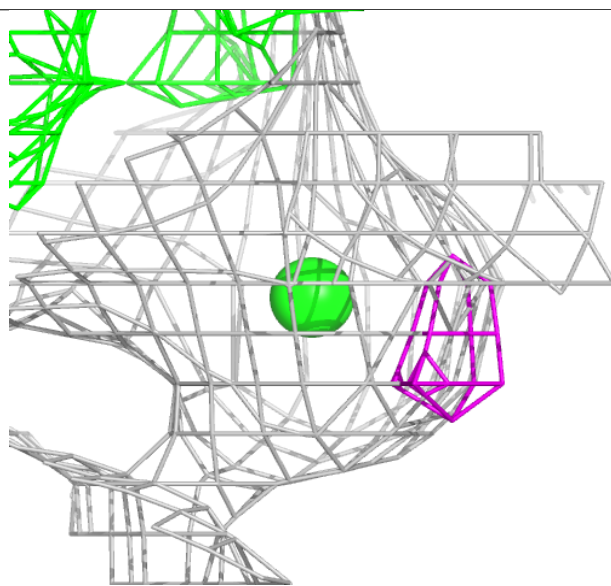
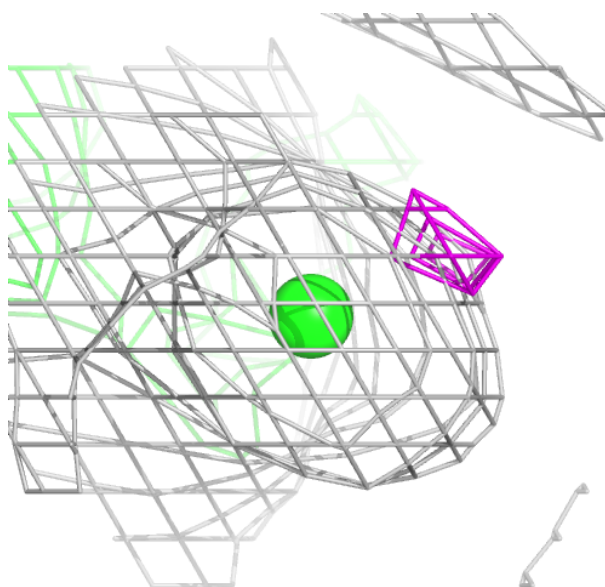
Electron density around CL B 609:

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



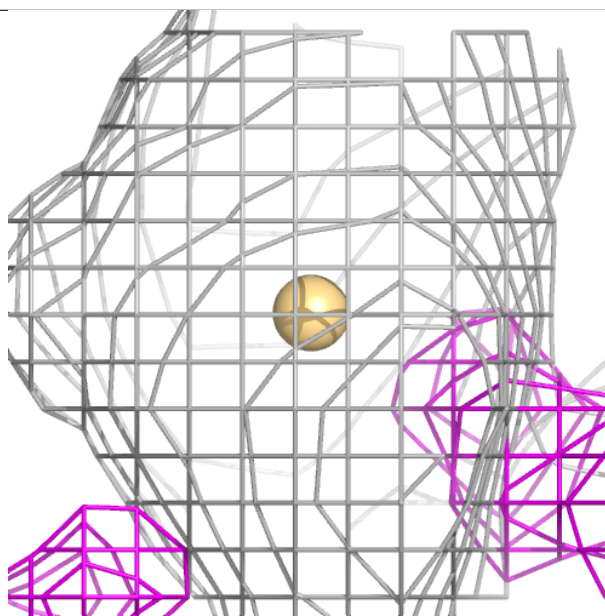
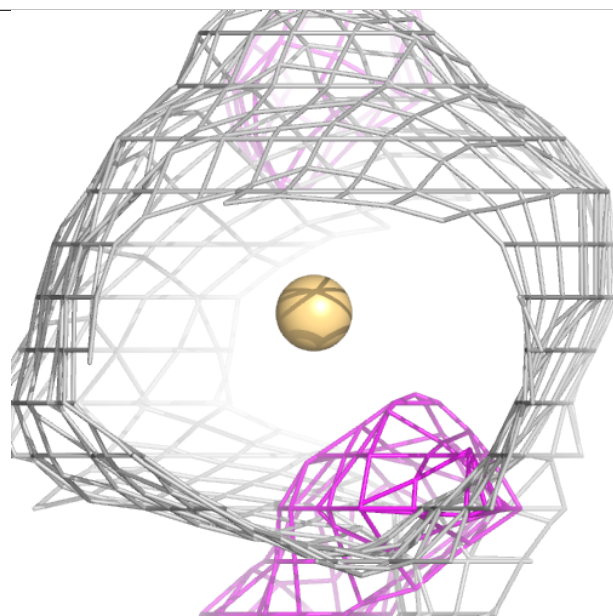
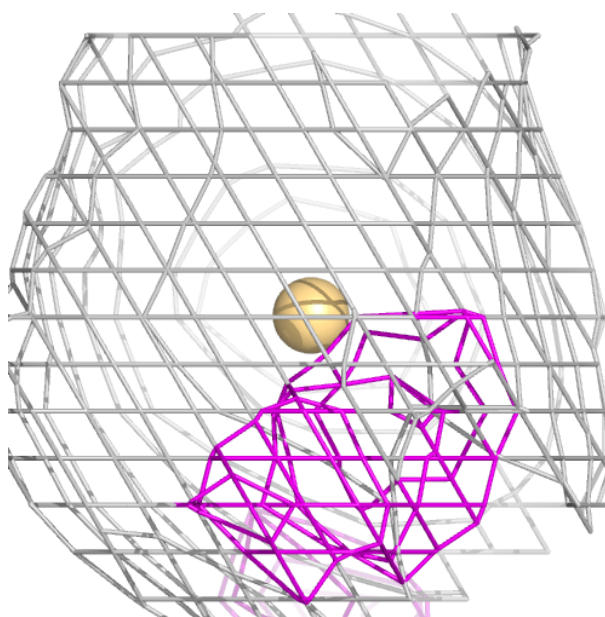
Electron density around CL A 607:

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



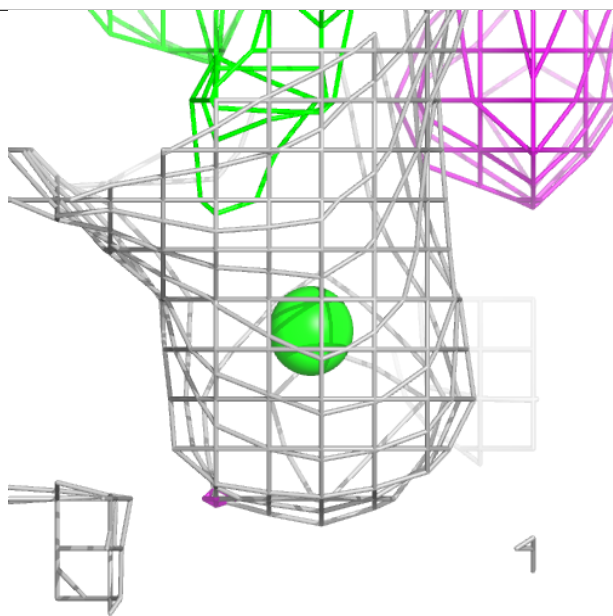
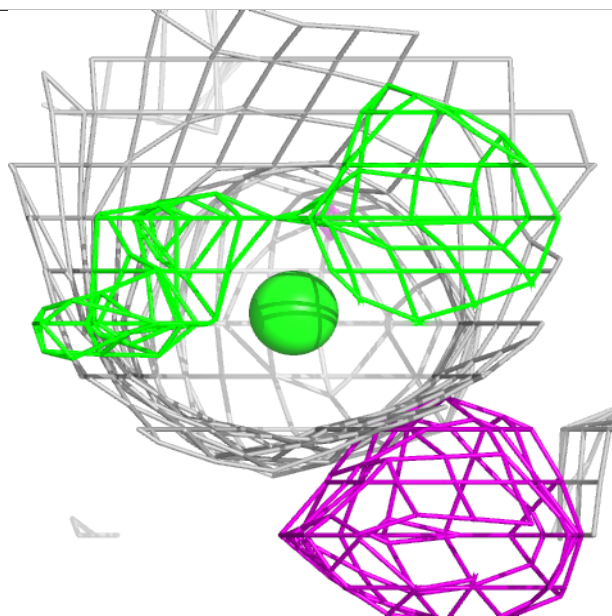
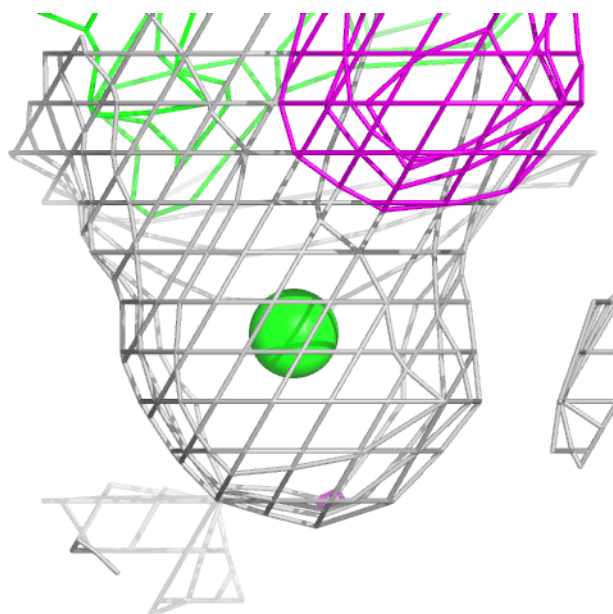
Electron density around CD A 603:

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



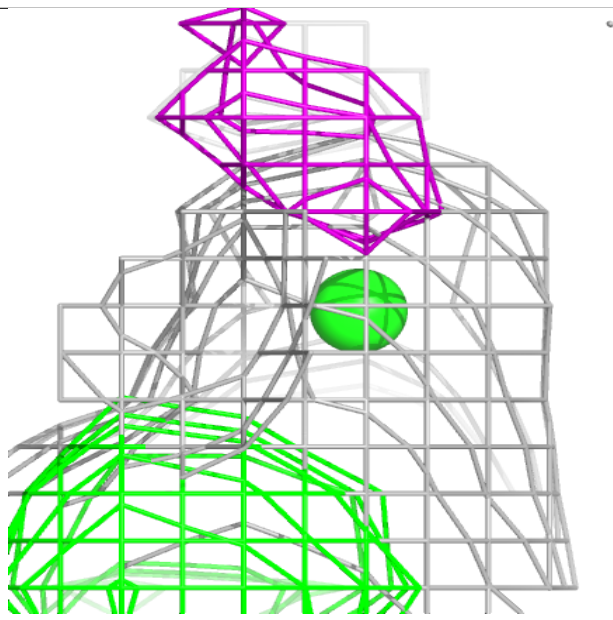
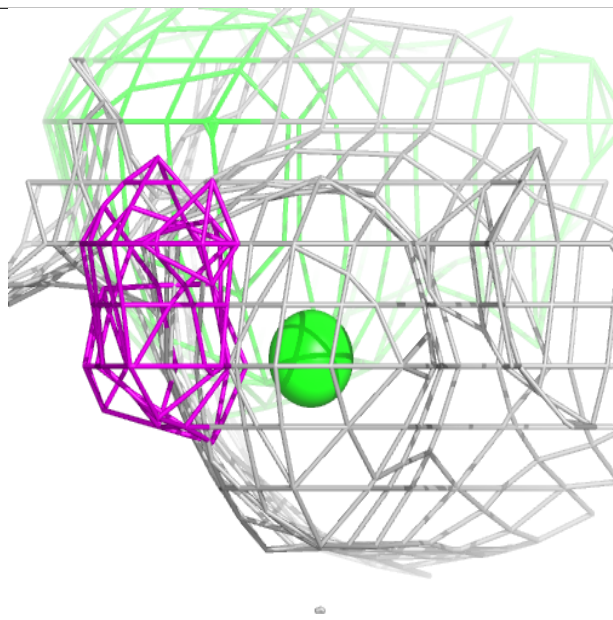
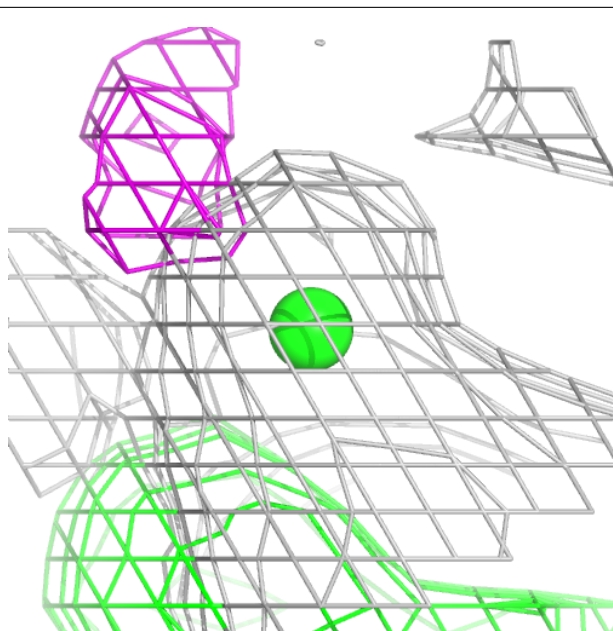
Electron density around CL A 606:

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



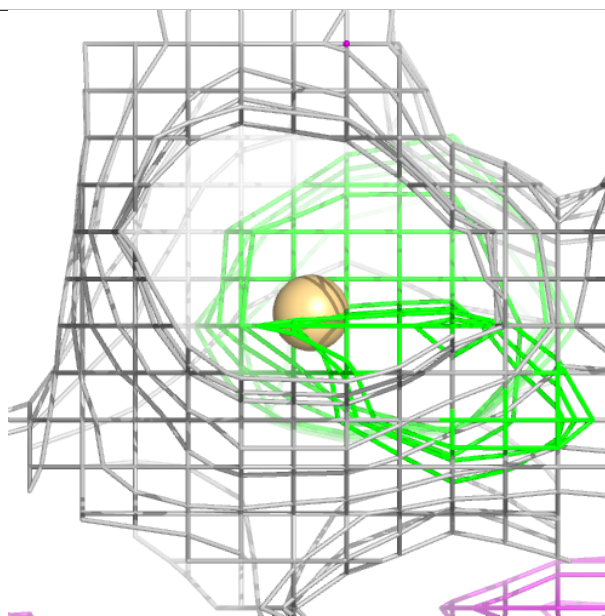
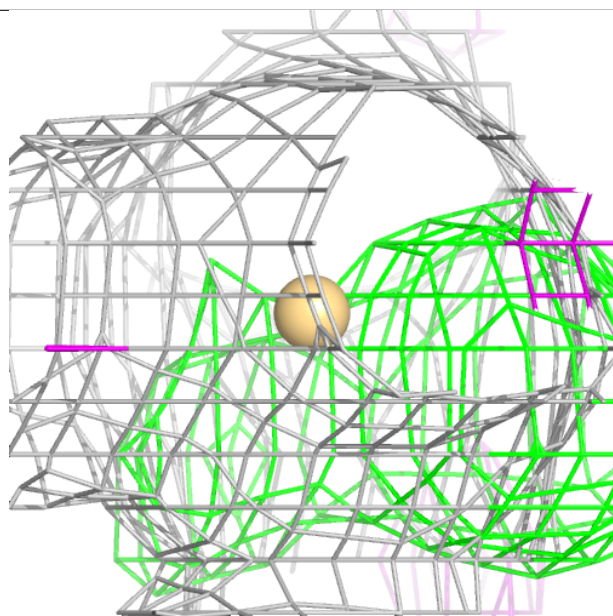
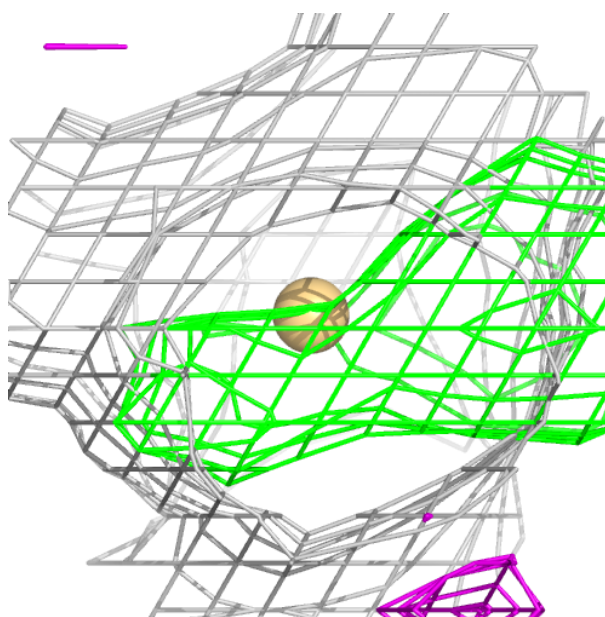
Electron density around CL B 608:

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



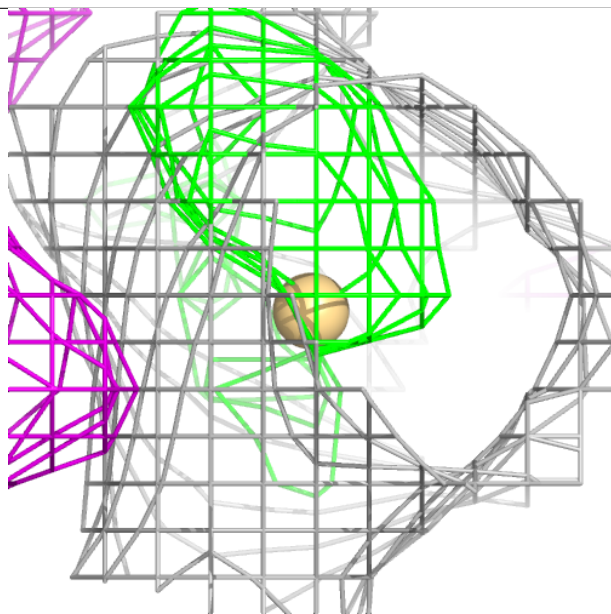
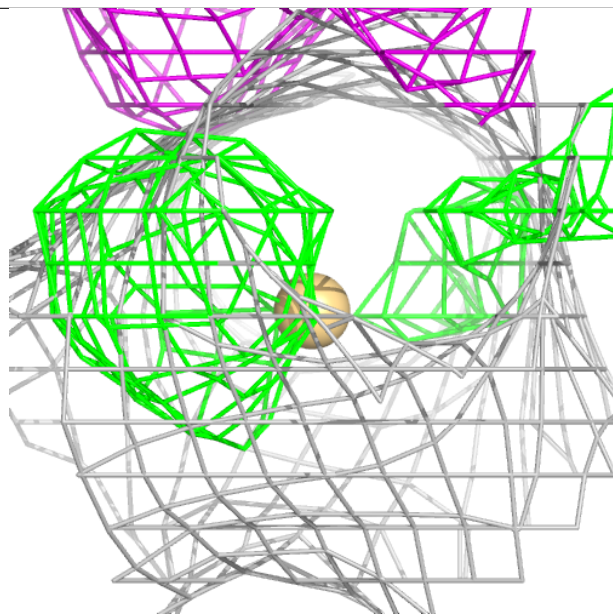
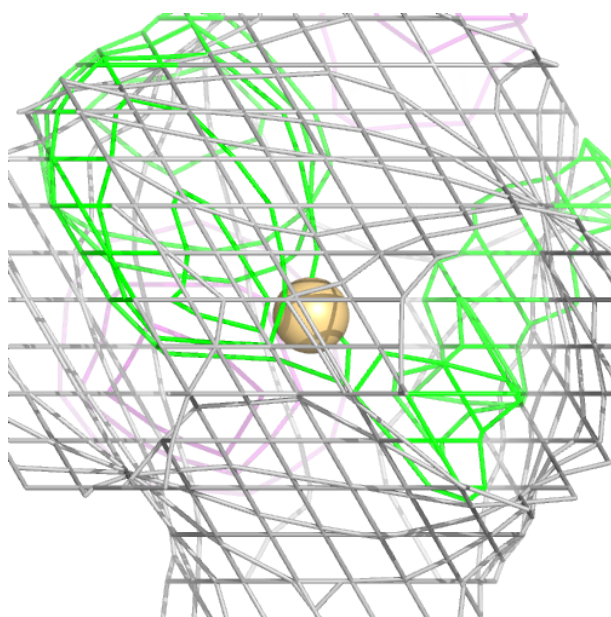
Electron density around CD A 602:

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



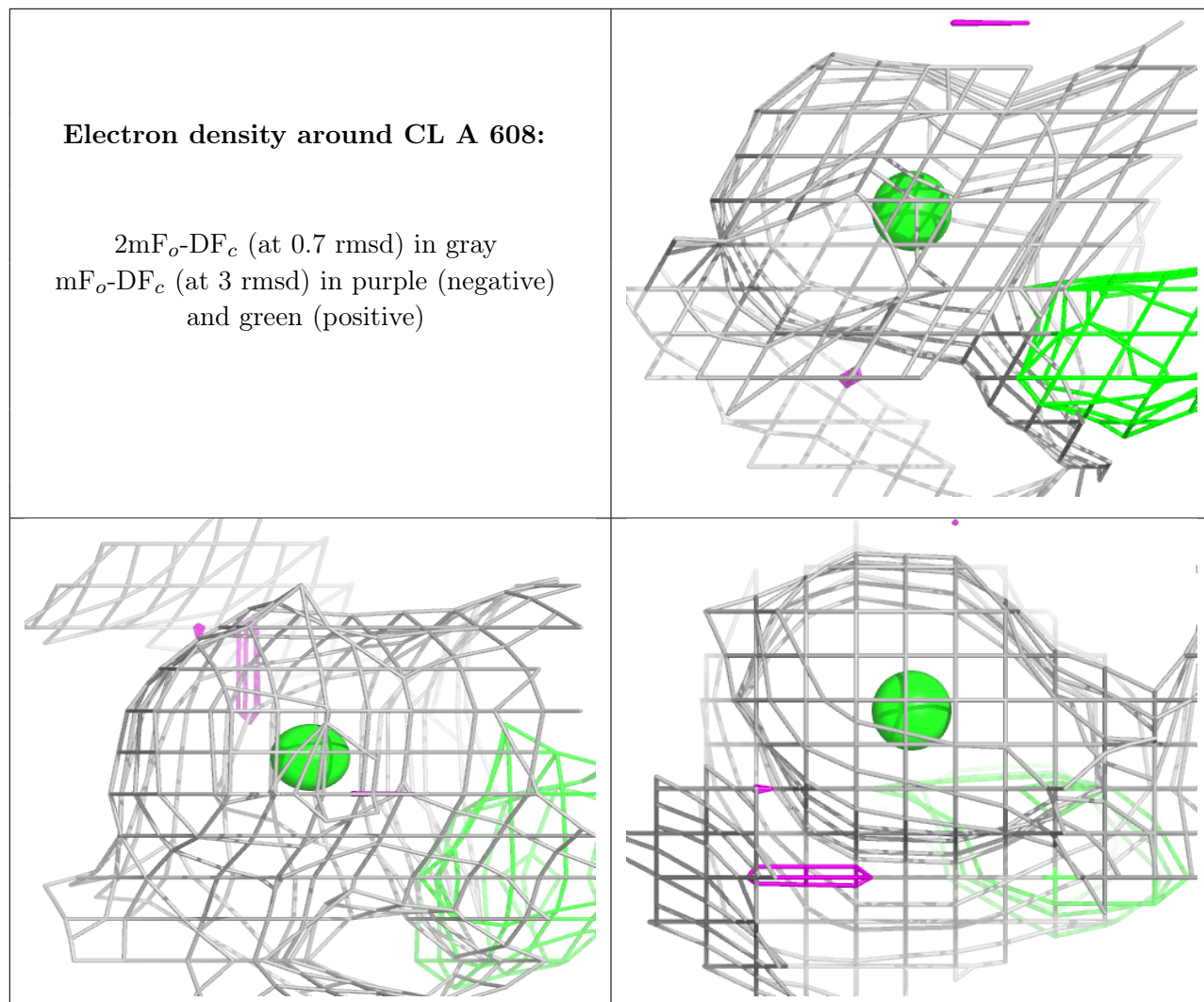
Electron density around CD A 601:

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



Electron density around CL A 608:

$2mF_o-DF_c$ (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.