



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 11:25 AM UTC

PDB ID : 6U7L / pdb_00006u7l
Title : 2.75 Angstrom Crystal Structure of Galactarate Dehydratase from Escherichia coli.
Authors : Minasov, G.; Shuvalova, L.; Wawrzak, Z.; Dubrovskaya, I.; Kiryukhina, O.; Endres, M.; Satchell, K.J.F.; Center for Structural Genomics of Infectious Diseases (CSGID)
Deposited on : 2019-09-03
Resolution : 2.75 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

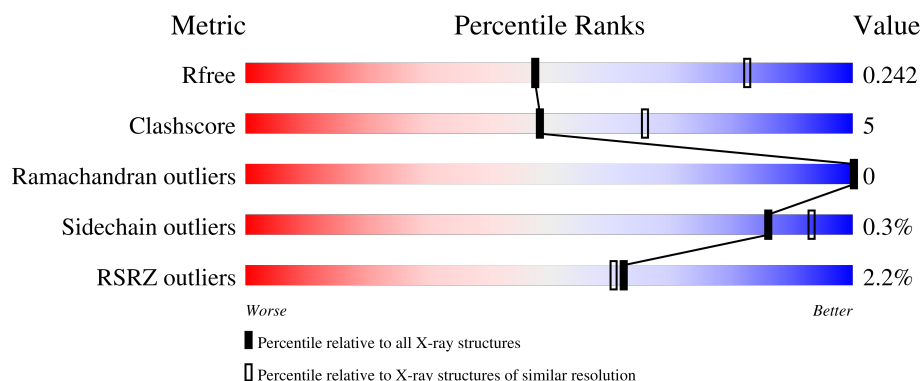
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.75 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	1009 (2.76-2.76)
Clashscore	190562	1044 (2.76-2.76)
Ramachandran outliers	187476	1024 (2.76-2.76)
Sidechain outliers	187428	1024 (2.76-2.76)
RSRZ outliers	180081	1009 (2.76-2.76)

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	526	% 78% 9% 13%
1	B	526	2% 74% 12% 14%
2	C	526	% 79% 7% 14%
2	D	526	3% 76% 10% 14%

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 14035 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 Galactarate dehydratase (L-threo-forming).

Mol	Chain	Residues	Atoms							ZeroOcc	AltConf	Trace
1	A	460	Total	C	N	O	P	S	Se	0	1	0
			3506	2229	600	661	1	6	9			
1	B	451	Total	C	N	O	P	S	Se	0	0	0
			3453	2197	590	651	1	5	9			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	SER	-	expression tag	UNP P39829
A	-1	ASN	-	expression tag	UNP P39829
A	0	ALA	-	expression tag	UNP P39829
B	-2	SER	-	expression tag	UNP P39829
B	-1	ASN	-	expression tag	UNP P39829
B	0	ALA	-	expression tag	UNP P39829

- Molecule 2 is a protein called Galactarate dehydratase (L-threo-forming).

Mol	Chain	Residues	Atoms							ZeroOcc	AltConf	Trace
2	C	453	Total	C	N	O	S	Se		0	0	0
			3455	2200	592	649	5	9				
2	D	452	Total	C	N	O	S	Se		0	1	0
			3455	2201	592	648	5	9				

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	-2	SER	-	expression tag	UNP P39829
C	-1	ASN	-	expression tag	UNP P39829
C	0	ALA	-	expression tag	UNP P39829
D	-2	SER	-	expression tag	UNP P39829
D	-1	ASN	-	expression tag	UNP P39829
D	0	ALA	-	expression tag	UNP P39829

- Molecule 3 is CALCIUM ION (CCD ID: CA) (formula: Ca) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	3	Total Ca 3 3	0	0
3	B	2	Total Ca 2 2	0	0
3	C	2	Total Ca 2 2	0	0
3	D	2	Total Ca 2 2	0	0

- Molecule 4 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

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

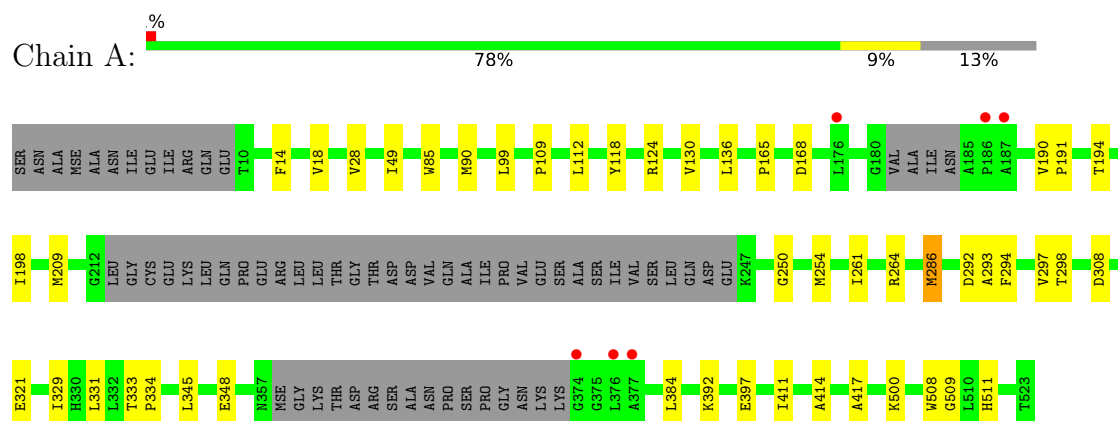
- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	41	Total O 41 41	0	0
5	B	25	Total O 25 25	0	0
5	C	52	Total O 52 52	0	0
5	D	32	Total O 32 32	0	0

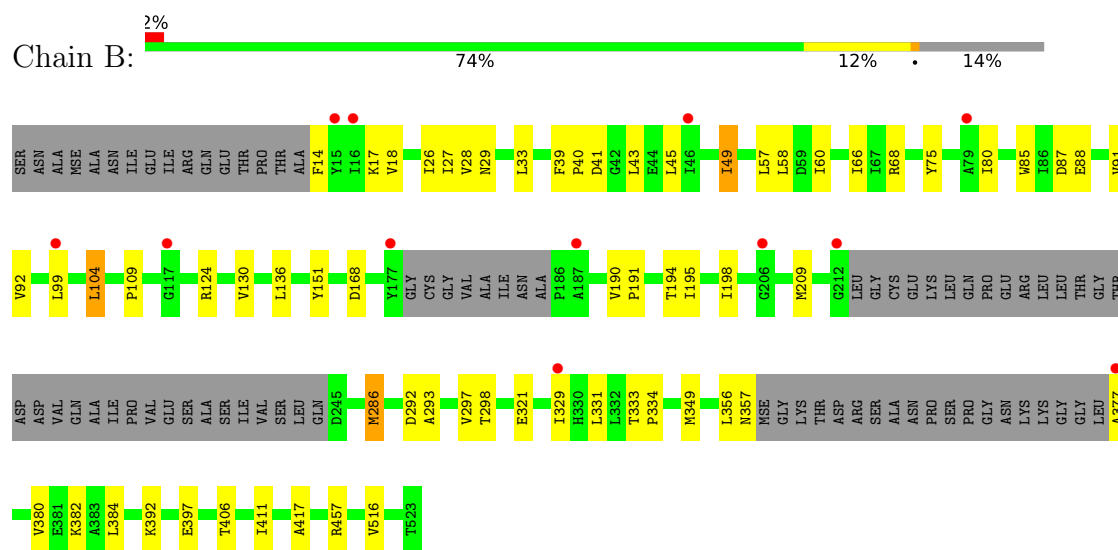
3 Residue-property plots [i](#)

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.

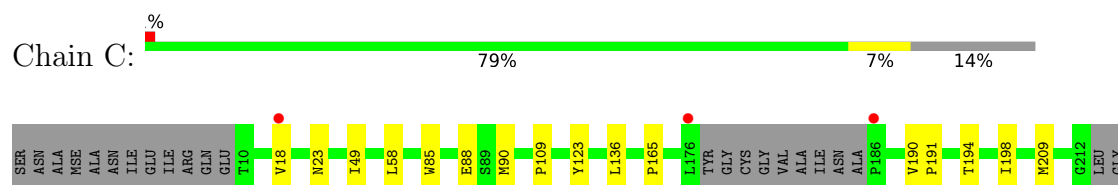
- Molecule 1: Galactarate dehydratase (L-threo-forming)

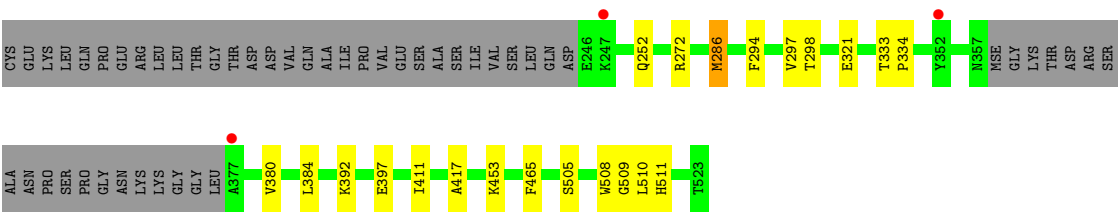


- Molecule 1: Galactarate dehydratase (L-threo-forming)

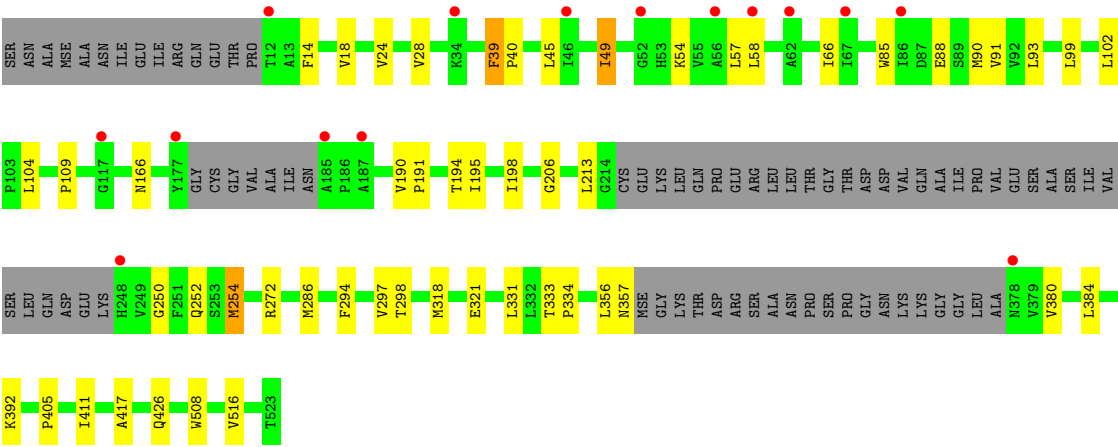
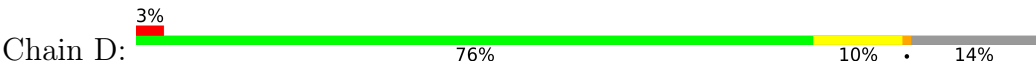


- Molecule 2: Galactarate dehydratase (L-threo-forming)





● Molecule 2: Galactarate dehydratase (L-threo-forming)



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	58.46Å 167.57Å 117.06Å 90.00° 103.97° 90.00°	Depositor
Resolution (Å)	29.23 – 2.75 29.23 – 2.75	Depositor EDS
% Data completeness (in resolution range)	98.8 (29.23-2.75) 98.8 (29.23-2.75)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.33 (at 2.76Å)	Xtriage
Refinement program	REFMAC 5.8.0253	Depositor
R, R_{free}	0.206 , 0.241 0.208 , 0.242	Depositor DCC
R_{free} test set	2734 reflections (4.82%)	wwPDB-VP
Wilson B-factor (Å ²)	59.2	Xtriage
Anisotropy	0.345	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 40.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.036 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	14035	wwPDB-VP
Average B, all atoms (Å ²)	77.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.41% 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: CL, PHD, CA

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	1.01	0/3557	1.19	4/4826 (0.1%)
1	B	1.01	1/3503 (0.0%)	1.21	1/4751 (0.0%)
2	C	1.01	0/3518	1.20	3/4775 (0.1%)
2	D	1.02	2/3518 (0.1%)	1.22	3/4776 (0.1%)
All	All	1.01	3/14096 (0.0%)	1.20	11/19128 (0.1%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	39	PHE	N-CA	5.28	1.49	1.46
1	B	49	ILE	N-CA	5.13	1.49	1.45
2	D	49	ILE	N-CA	5.12	1.49	1.45

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	250	GLY	CA-C-O	-7.15	116.87	122.52
2	C	286	MSE	CG-SE-CE	5.94	112.00	98.92
1	A	254	MSE	CG-SE-CE	5.86	111.81	98.92
2	C	90	MSE	CG-SE-CE	5.67	111.39	98.92
1	A	286	MSE	CG-SE-CE	5.64	111.33	98.92
2	D	90	MSE	CG-SE-CE	5.48	110.97	98.92
1	A	90	MSE	CG-SE-CE	5.46	110.93	98.92
1	B	286	MSE	CG-SE-CE	5.44	110.89	98.92
2	D	254	MSE	CG-SE-CE	5.26	110.49	98.92
1	A	250	GLY	CA-C-O	-5.12	118.48	122.52
2	C	453	LYS	CB-CA-C	5.03	118.33	110.19

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	3506	0	3508	35	0
1	B	3453	0	3455	52	0
2	C	3455	0	3469	27	0
2	D	3455	0	3463	46	0
3	A	3	0	0	0	0
3	B	2	0	0	0	0
3	C	2	0	0	0	0
3	D	2	0	0	0	0
4	A	4	0	0	0	0
4	D	3	0	0	0	0
5	A	41	0	0	0	0
5	B	25	0	0	0	0
5	C	52	0	0	0	0
5	D	32	0	0	2	0
All	All	14035	0	13895	152	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 5.

All (152) 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:B:27:ILE:CD1	1:B:49:ILE:HB	2.16	0.74
2:D:66:ILE:CD1	2:D:91:VAL:HG12	2.19	0.72
2:D:93:LEU:HD12	2:D:93:LEU:O	1.92	0.69
1:B:329:ILE:HG21	1:B:349:MSE:SE	2.43	0.69
2:D:213:LEU:C	2:D:213:LEU:HD23	2.21	0.66
1:A:348:GLU:HG2	1:A:414:ALA:HB2	1.79	0.64
2:D:318:MSE:HE1	2:D:426:GLN:HA	1.80	0.63
1:B:39:PHE:HB3	1:B:40:PRO:HD2	1.81	0.63
2:C:252:GLN:NE2	2:D:508:TRP:CD2	2.68	0.61
1:A:112:LEU:HD12	1:A:112:LEU:N	2.15	0.60
1:A:348:GLU:HG2	1:A:414:ALA:CB	2.31	0.60
1:B:66:ILE:HD12	1:B:91:VAL:HG23	1.81	0.60
1:B:190:VAL:HB	1:B:191:PRO:HD3	1.84	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:104:LEU:HD12	2:D:104:LEU:N	2.17	0.59
2:D:39:PHE:HB3	2:D:40:PRO:HD2	1.85	0.59
2:C:190:VAL:HB	2:C:191:PRO:HD3	1.85	0.59
2:C:252:GLN:NE2	2:D:508:TRP:CE3	2.70	0.59
1:B:27:ILE:HD11	1:B:49:ILE:HB	1.83	0.59
1:B:104:LEU:N	1:B:104:LEU:HD12	2.16	0.59
1:A:190:VAL:HB	1:A:191:PRO:HD3	1.85	0.59
2:D:99:LEU:HD12	2:D:99:LEU:N	2.18	0.58
2:D:88:GLU:OE1	2:D:88:GLU:N	2.35	0.58
2:D:356:LEU:O	2:D:357:ASN:C	2.47	0.58
2:D:190:VAL:HB	2:D:191:PRO:HD3	1.85	0.57
1:B:60:ILE:CG1	1:B:80:ILE:HB	2.35	0.57
2:D:24:VAL:CG1	2:D:54:LYS:HB2	2.35	0.57
1:A:99:LEU:HD12	1:A:99:LEU:N	2.20	0.56
2:C:88:GLU:N	2:C:88:GLU:OE1	2.36	0.56
2:C:397:GLU:HB3	2:C:411:ILE:HD12	1.88	0.56
1:B:88:GLU:OE1	1:B:88:GLU:N	2.34	0.56
1:B:397:GLU:HB3	1:B:411:ILE:HD12	1.87	0.55
1:B:29:ASN:OD1	1:B:33:LEU:HD21	2.06	0.55
1:A:397:GLU:HB3	1:A:411:ILE:HD12	1.88	0.54
2:C:18:VAL:CG2	2:C:297:VAL:HG12	2.37	0.54
1:B:41:ASP:OD1	1:B:43:LEU:HD22	2.08	0.54
1:B:58:LEU:C	1:B:58:LEU:HD12	2.32	0.53
2:D:405:PRO:HB3	2:D:411:ILE:HD13	1.90	0.53
1:B:292:ASP:OD2	1:B:457:ARG:NH1	2.38	0.53
1:B:356:LEU:O	1:B:357:ASN:CB	2.56	0.53
1:A:333:THR:OG1	1:A:334:PRO:HD3	2.09	0.53
1:B:333:THR:OG1	1:B:334:PRO:HD3	2.09	0.53
2:C:297:VAL:HG23	2:C:298:THR:HG23	1.90	0.53
2:D:18:VAL:CG2	2:D:297:VAL:HG12	2.39	0.53
1:B:321:GLU:HG2	1:B:417:ALA:HA	1.91	0.52
2:D:206:GLY:HA2	5:D:727:HOH:O	2.07	0.52
2:D:321:GLU:HG2	2:D:417:ALA:HA	1.91	0.52
1:B:75:TYR:HB2	1:B:92:VAL:CG2	2.39	0.52
2:C:333:THR:OG1	2:C:334:PRO:HD3	2.09	0.52
1:B:377:ALA:HB3	1:B:382:LYS:HE3	1.92	0.52
2:C:509:GLY:HA2	2:C:511:HIS:CE1	2.45	0.52
2:D:206:GLY:N	5:D:701:HOH:O	2.42	0.52
2:C:321:GLU:HG2	2:C:417:ALA:HA	1.92	0.52
1:B:49:ILE:HG12	1:B:85:TRP:HB2	1.92	0.52
2:D:297:VAL:HG23	2:D:298:THR:HG23	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:165:PRO:HB2	2:C:465:PHE:CD1	2.45	0.51
1:B:17:LYS:CE	1:B:57:LEU:HD13	2.41	0.51
2:D:333:THR:OG1	2:D:334:PRO:HD3	2.10	0.51
1:A:165:PRO:HB2	2:C:465:PHE:CG	2.46	0.51
1:A:321:GLU:HG2	1:A:417:ALA:HA	1.93	0.51
2:D:194:THR:O	2:D:198:ILE:HG13	2.11	0.51
1:A:49:ILE:HG12	1:A:85:TRP:HB2	1.93	0.50
1:B:17:LYS:CE	1:B:57:LEU:CD1	2.89	0.50
1:B:194:THR:O	1:B:198:ILE:HG13	2.12	0.50
2:D:66:ILE:HD13	2:D:91:VAL:HG12	1.91	0.50
1:A:18:VAL:CG2	1:A:297:VAL:HG12	2.42	0.49
1:B:41:ASP:OD1	1:B:43:LEU:CD2	2.59	0.49
1:B:60:ILE:HG13	1:B:80:ILE:HB	1.93	0.49
2:C:49:ILE:HG12	2:C:85:TRP:HB2	1.95	0.49
1:A:14:PHE:HB2	1:A:28:VAL:CG2	2.42	0.49
2:D:14:PHE:HB2	2:D:28:VAL:CG2	2.43	0.49
1:A:500:LYS:HA	2:C:508:TRP:CZ3	2.48	0.49
1:B:356:LEU:O	1:B:357:ASN:CG	2.56	0.49
1:A:118:TYR:CE1	2:C:165:PRO:HG3	2.48	0.48
1:B:39:PHE:HE2	1:B:45:LEU:HD11	1.78	0.48
2:D:49:ILE:HG12	2:D:85:TRP:HB2	1.94	0.48
1:B:297:VAL:HG23	1:B:298:THR:HG23	1.96	0.48
1:B:45:LEU:N	1:B:45:LEU:HD12	2.28	0.48
1:A:331:LEU:CD1	1:A:384:LEU:CD2	2.91	0.48
1:B:18:VAL:CG2	1:B:297:VAL:HG12	2.43	0.47
1:B:39:PHE:CE2	1:B:45:LEU:HD11	2.49	0.47
1:B:331:LEU:CD1	1:B:384:LEU:CD2	2.92	0.47
1:A:509:GLY:HA2	1:A:511:HIS:CE1	2.49	0.47
1:B:17:LYS:HE2	1:B:57:LEU:CD1	2.45	0.47
1:B:27:ILE:HD12	1:B:49:ILE:HB	1.95	0.47
2:C:18:VAL:HG22	2:C:297:VAL:HG12	1.96	0.47
2:C:194:THR:O	2:C:198:ILE:HG13	2.13	0.47
2:D:58:LEU:HD12	2:D:58:LEU:N	2.30	0.47
2:C:23:ASN:HB2	2:C:58:LEU:HD21	1.96	0.46
1:A:194:THR:O	1:A:198:ILE:HG13	2.14	0.46
1:B:14:PHE:HB2	1:B:28:VAL:CG2	2.45	0.46
1:B:329:ILE:CG2	1:B:349:MSE:SE	3.13	0.46
2:D:331:LEU:CD1	2:D:384:LEU:CD2	2.94	0.46
1:B:124:ARG:HD2	1:B:130:VAL:HG22	1.97	0.46
1:B:397:GLU:OE2	1:B:406:THR:HG22	2.16	0.46
1:A:112:LEU:N	1:A:112:LEU:CD1	2.79	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:45:LEU:HD12	2:D:45:LEU:N	2.30	0.46
1:B:41:ASP:CG	1:B:43:LEU:HD22	2.41	0.45
1:B:26:ILE:HG13	1:B:68:ARG:NH2	2.32	0.45
2:C:508:TRP:CE3	2:D:252:GLN:NE2	2.85	0.45
2:D:14:PHE:HB2	2:D:28:VAL:HG22	1.99	0.45
2:D:24:VAL:CA	2:D:57:LEU:HD13	2.47	0.44
1:A:14:PHE:HB2	1:A:28:VAL:HG22	1.98	0.44
1:A:297:VAL:HG23	1:A:298:THR:HG23	1.98	0.44
2:D:405:PRO:HB3	2:D:411:ILE:CD1	2.46	0.44
1:A:99:LEU:N	1:A:99:LEU:CD1	2.81	0.44
2:D:166[B]:ASN:ND2	2:D:272:ARG:O	2.50	0.44
1:A:508:TRP:O	1:B:151:TYR:HE1	2.01	0.44
2:D:318:MSE:CE	2:D:426:GLN:HG2	2.48	0.43
1:A:348:GLU:CG	1:A:414:ALA:CB	2.96	0.43
2:D:213:LEU:C	2:D:213:LEU:CD2	2.90	0.43
1:A:109:PRO:HG3	1:A:392:LYS:O	2.18	0.43
1:B:109:PRO:HG2	1:B:392:LYS:O	2.18	0.43
2:D:213:LEU:HD12	2:D:254:MSE:HE1	2.01	0.43
2:D:24:VAL:HG13	2:D:54:LYS:HB2	1.99	0.43
1:B:87:ASP:OD1	1:B:88:GLU:OE1	2.37	0.43
1:A:331:LEU:HD12	1:A:384:LEU:CD2	2.48	0.43
1:B:136:LEU:HD21	1:B:209:MSE:CE	2.49	0.42
2:D:99:LEU:N	2:D:99:LEU:CD1	2.81	0.42
1:A:329:ILE:HD11	1:A:345:LEU:HG	2.01	0.42
2:D:109:PRO:HG3	2:D:392:LYS:O	2.19	0.42
2:C:123:TYR:HB3	2:C:272:ARG:NH2	2.35	0.42
2:D:294:PHE:HA	2:D:297:VAL:HG22	2.00	0.42
1:B:14:PHE:HB2	1:B:28:VAL:HG22	2.02	0.42
2:D:57:LEU:N	2:D:57:LEU:HD12	2.35	0.42
2:D:195:ILE:HD11	2:D:516:VAL:HG21	2.02	0.42
1:B:87:ASP:OD1	1:B:88:GLU:N	2.52	0.42
1:B:356:LEU:O	1:B:357:ASN:HB3	2.20	0.42
2:C:23:ASN:CB	2:C:58:LEU:HD21	2.50	0.42
2:C:294:PHE:HA	2:C:297:VAL:HG22	2.01	0.42
1:A:308:ASP:OD2	1:A:392:LYS:HG2	2.19	0.41
2:D:24:VAL:HA	2:D:57:LEU:HD13	2.02	0.41
2:D:93:LEU:HD12	2:D:93:LEU:C	2.45	0.41
1:A:136:LEU:HD21	1:A:209:MSE:CE	2.51	0.41
1:A:294:PHE:HA	1:A:297:VAL:HG22	2.02	0.41
2:C:136:LEU:HD21	2:C:209:MSE:CE	2.51	0.41
1:B:99:LEU:CD2	1:B:329:ILE:CD1	2.98	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:292:ASP:OD1	1:B:293:ALA:N	2.54	0.41
2:C:109:PRO:HG3	2:C:392:LYS:O	2.21	0.41
2:D:102:LEU:O	2:D:104:LEU:HD12	2.20	0.41
1:A:261:ILE:HD12	1:A:264:ARG:HH22	1.85	0.41
2:D:213:LEU:HA	2:D:254:MSE:SE	2.71	0.41
1:A:124:ARG:HD3	1:A:130:VAL:HG22	2.02	0.41
2:C:505:SER:HA	2:C:510:LEU:HD12	2.02	0.41
1:A:333:THR:N	1:A:334:PRO:CD	2.84	0.41
1:A:292:ASP:OD1	1:A:293:ALA:N	2.54	0.40
2:C:380:VAL:O	2:C:384:LEU:HG	2.22	0.40
1:B:195:ILE:HD11	1:B:516:VAL:HG21	2.04	0.40
1:B:333:THR:N	1:B:334:PRO:CD	2.85	0.40
1:B:380:VAL:O	1:B:384:LEU:HG	2.21	0.40
2:C:333:THR:N	2:C:334:PRO:CD	2.84	0.40
1:A:112:LEU:HD12	1:A:112:LEU:H	1.87	0.40
2:D:380:VAL:O	2:D:384:LEU:HG	2.21	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	452/526 (86%)	430 (95%)	22 (5%)	0	100	100
1	B	442/526 (84%)	425 (96%)	17 (4%)	0	100	100
2	C	445/526 (85%)	425 (96%)	20 (4%)	0	100	100
2	D	445/526 (85%)	425 (96%)	20 (4%)	0	100	100
All	All	1784/2104 (85%)	1705 (96%)	79 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	376/420 (90%)	375 (100%)	1 (0%)	86	93
1	B	372/420 (89%)	370 (100%)	2 (0%)	81	89
2	C	374/421 (89%)	373 (100%)	1 (0%)	86	93
2	D	373/421 (89%)	372 (100%)	1 (0%)	86	93
All	All	1495/1682 (89%)	1490 (100%)	5 (0%)	86	93

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	286	MSE
1	B	104	LEU
1	B	286	MSE
2	C	286	MSE
2	D	286	MSE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	175	HIS
1	A	491	HIS
1	A	512	ASN
1	A	518	ASN
1	B	175	HIS
1	B	518	ASN
2	C	513	GLN
2	D	175	HIS
2	D	518	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

2 non-standard protein/DNA/RNA residues are modelled in this entry.

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	$\# Z > 2$
1	PHD	B	168	1	9,11,12	0.94	1 (11%)	9,15,17	0.71	0
1	PHD	A	168	1	9,11,12	1.12	1 (11%)	9,15,17	0.69	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
1	PHD	B	168	1	-	1/8/11/13	-
1	PHD	A	168	1	-	1/8/11/13	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	168	PHD	P-OD1	2.66	1.64	1.59
1	B	168	PHD	P-OD1	2.16	1.63	1.59

There are no bond angle outliers.

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	168	PHD	O-C-CA-CB
1	B	168	PHD	O-C-CA-CB

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 16 ligands modelled in this entry, 16 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	450/526 (85%)	-0.06	6 (1%) 75 75	28, 64, 112, 171	1 (0%)
1	B	441/526 (83%)	0.26	12 (2%) 56 54	42, 81, 146, 178	0
2	C	444/526 (84%)	-0.13	6 (1%) 73 73	30, 61, 101, 162	0
2	D	443/526 (84%)	0.21	15 (3%) 48 46	29, 73, 183, 208	1 (0%)
All	All	1778/2104 (84%)	0.07	39 (2%) 62 60	28, 68, 149, 208	2 (0%)

All (39) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	C	176	LEU	3.9
1	B	187	ALA	3.3
1	A	176	LEU	3.2
2	D	177	TYR	3.2
2	D	248	HIS	3.1
2	C	18	VAL	3.0
2	D	52	GLY	2.8
1	B	15	TYR	2.8
2	D	58	LEU	2.8
2	D	185	ALA	2.6
1	B	117	GLY	2.6
1	A	187	ALA	2.5
1	B	377	ALA	2.5
1	B	16	ILE	2.5
1	A	376	LEU	2.5
1	B	206	GLY	2.5
2	D	46	ILE	2.4
2	D	187	ALA	2.4
1	B	177	TYR	2.4
2	D	62	ALA	2.4
2	D	12	THR	2.4

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Mol	Chain	Res	Type	RSRZ
2	D	67	ILE	2.3
2	D	34	LYS	2.3
1	A	374	GLY	2.3
1	B	46	ILE	2.2
2	D	56	ALA	2.2
1	A	186	PRO	2.2
2	D	117	GLY	2.2
1	B	99	LEU	2.2
2	C	247	LYS	2.2
2	C	377	ALA	2.2
1	A	377	ALA	2.1
1	B	212	GLY	2.1
1	B	329	ILE	2.1
2	C	352	TYR	2.1
2	D	86	ILE	2.1
1	B	79	ALA	2.0
2	C	186	PRO	2.0
2	D	378	ASN	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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
1	PHD	B	168	12/13	0.82	0.11	61,73,90,91	4
1	PHD	A	168	12/13	0.85	0.11	46,56,92,93	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

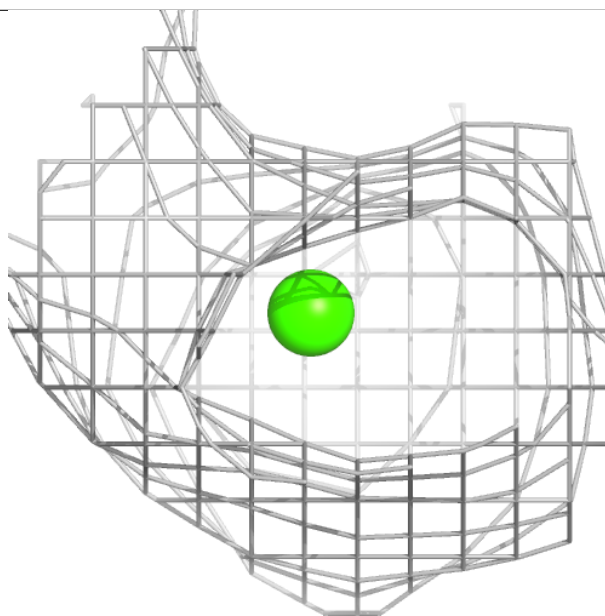
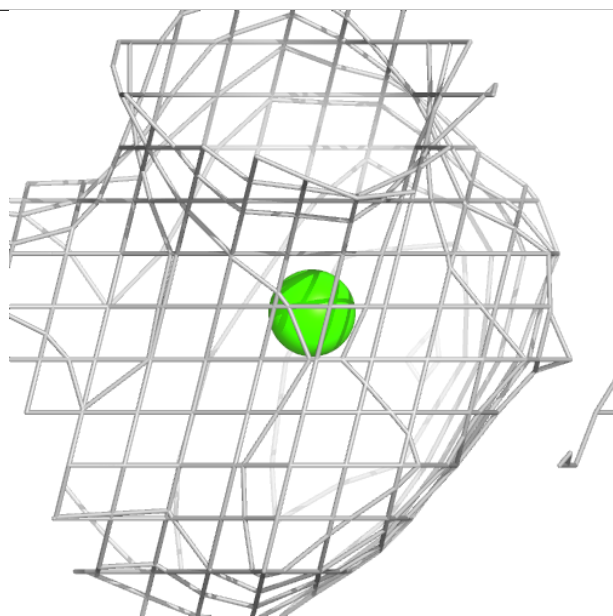
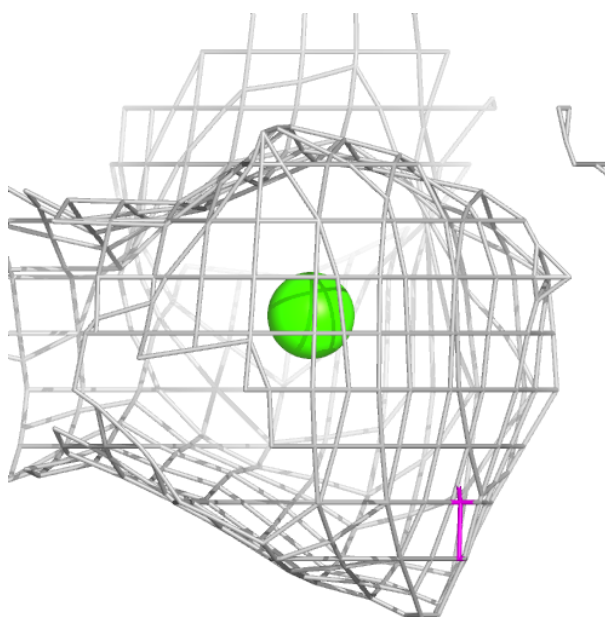
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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
4	CL	A	606	1/1	0.71	0.14	88,88,88,88	0
3	CA	A	602	1/1	0.73	0.10	92,92,92,92	0
3	CA	A	603	1/1	0.79	0.16	63,63,63,63	1
4	CL	A	604	1/1	0.80	0.08	86,86,86,86	0
4	CL	D	605	1/1	0.81	0.12	89,89,89,89	0
4	CL	D	603	1/1	0.84	0.10	85,85,85,85	0
4	CL	D	604	1/1	0.84	0.08	97,97,97,97	0
3	CA	B	602	1/1	0.84	0.10	104,104,104,104	0
3	CA	C	602	1/1	0.90	0.08	99,99,99,99	0
4	CL	A	605	1/1	0.91	0.09	76,76,76,76	0
3	CA	B	601	1/1	0.95	0.06	69,69,69,69	0
4	CL	A	607	1/1	0.96	0.10	77,77,77,77	0
3	CA	D	602	1/1	0.96	0.10	102,102,102,102	0
3	CA	D	601	1/1	0.97	0.07	65,65,65,65	0
3	CA	C	601	1/1	0.97	0.07	51,51,51,51	0
3	CA	A	601	1/1	0.97	0.07	62,62,62,62	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.

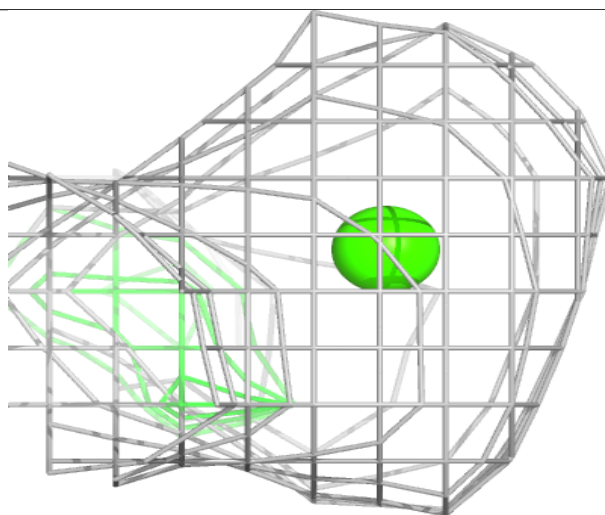
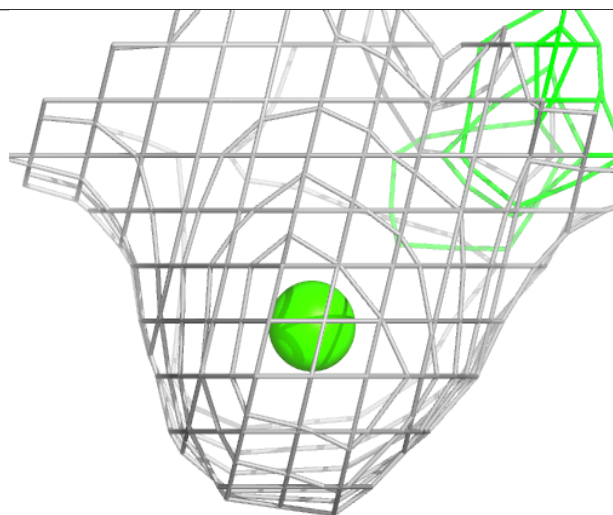
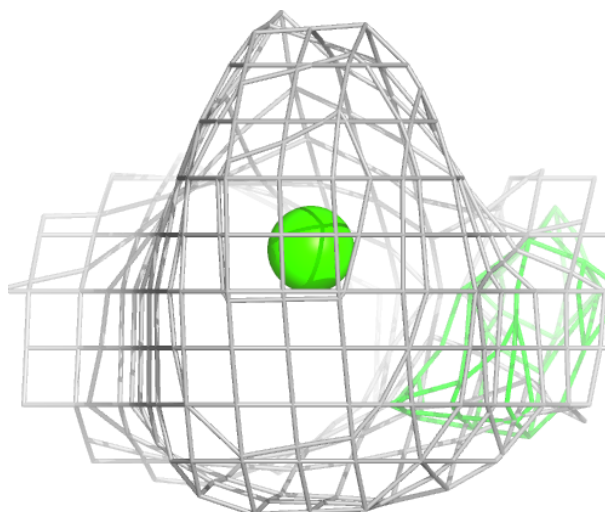
Electron density around CA A 602:

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



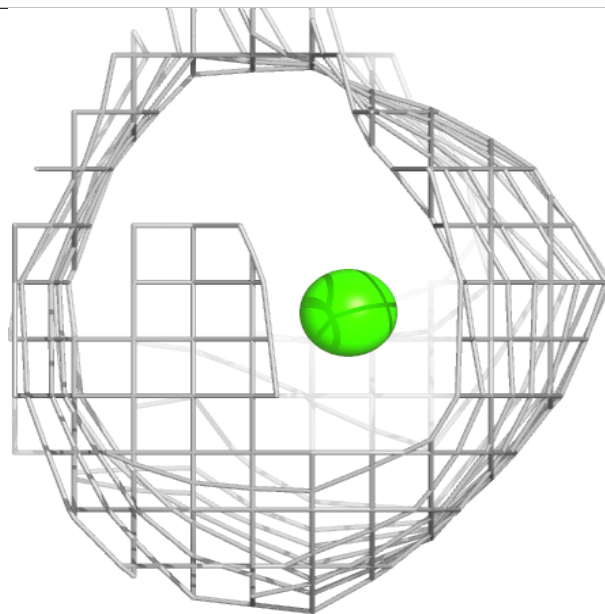
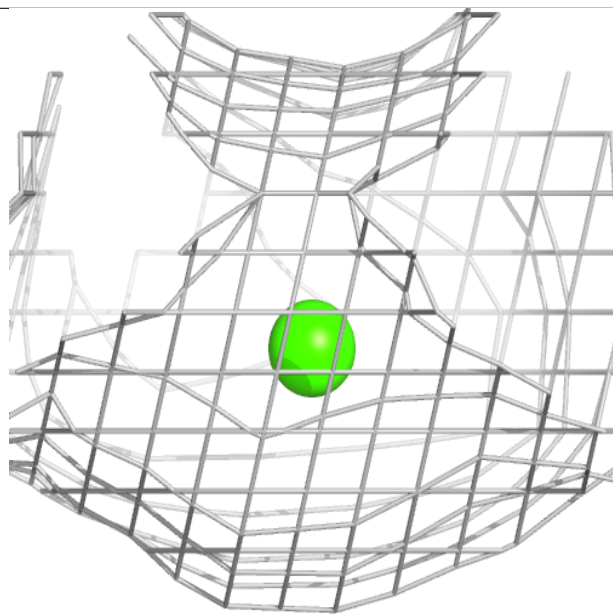
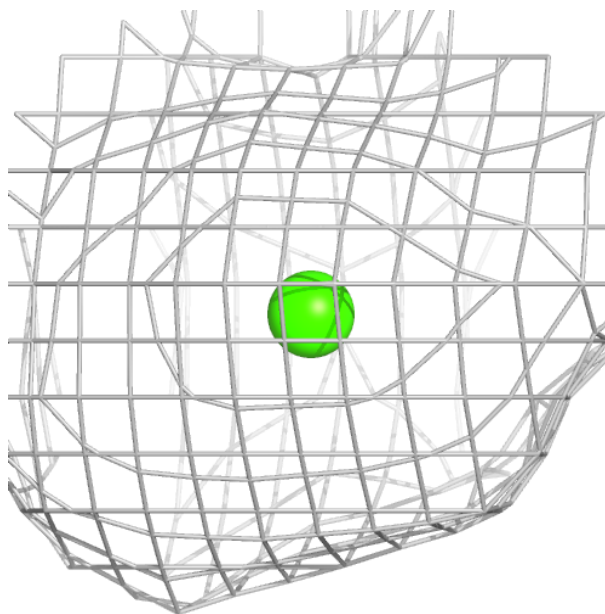
Electron density around CA A 603:

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



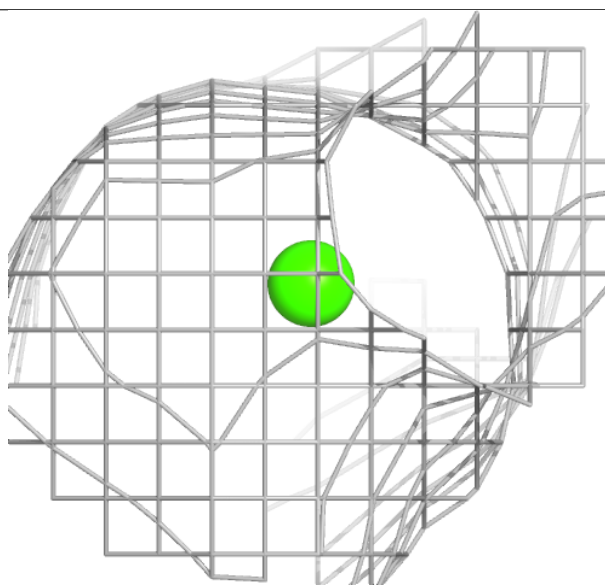
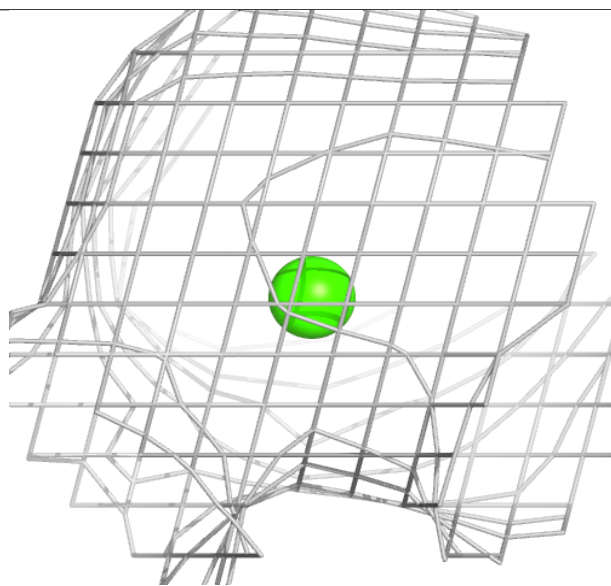
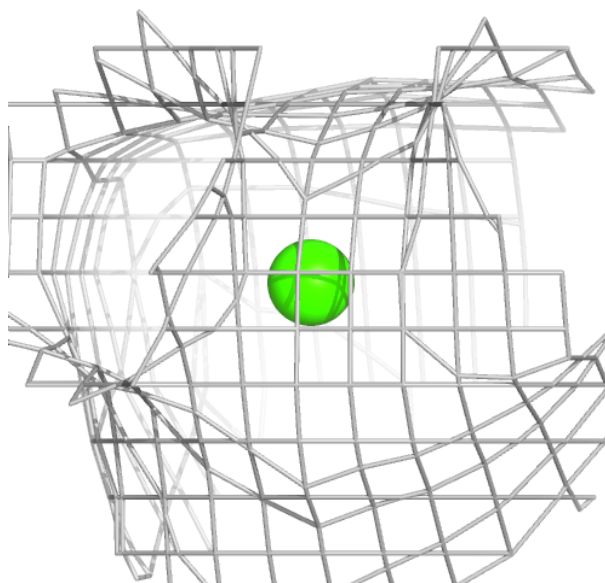
Electron density around CA B 602:

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



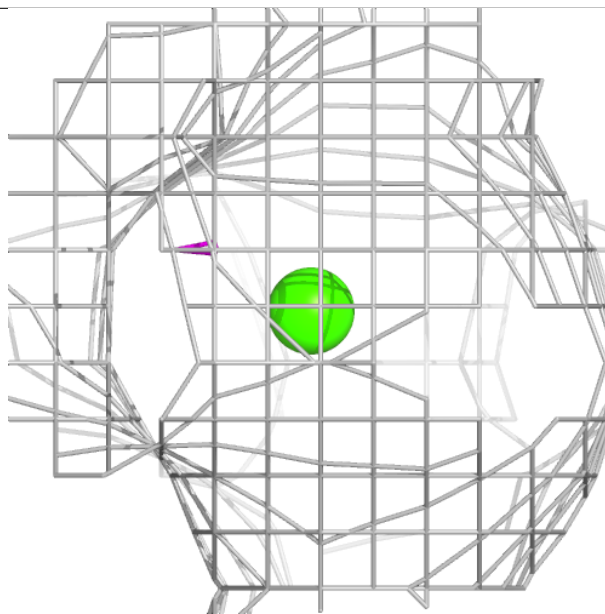
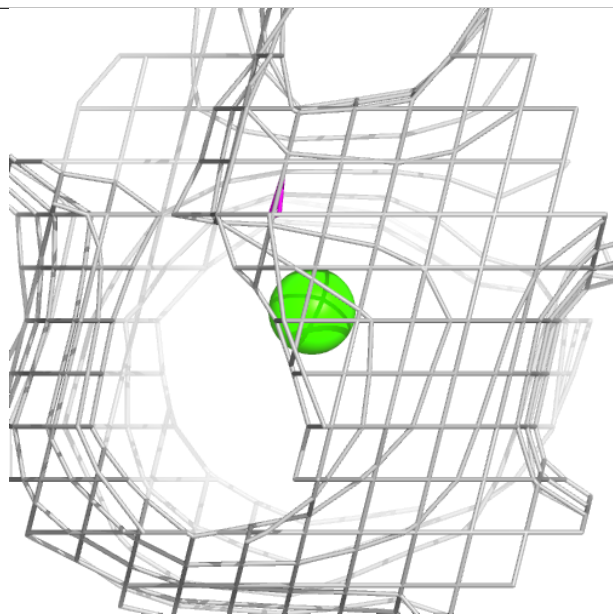
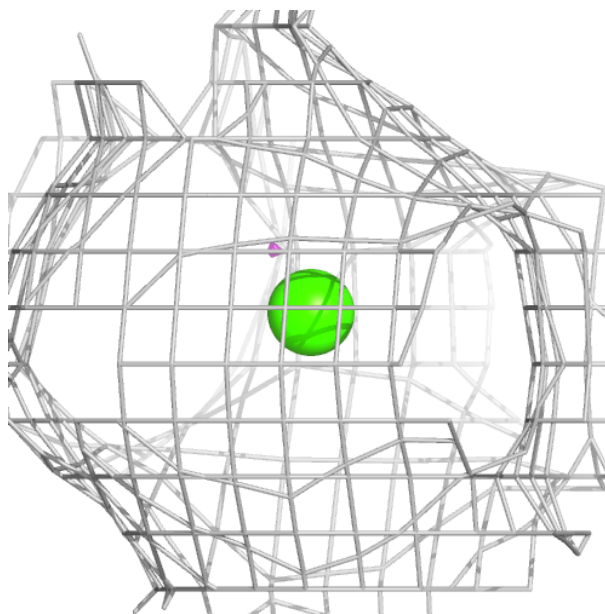
Electron density around CA C 602:

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



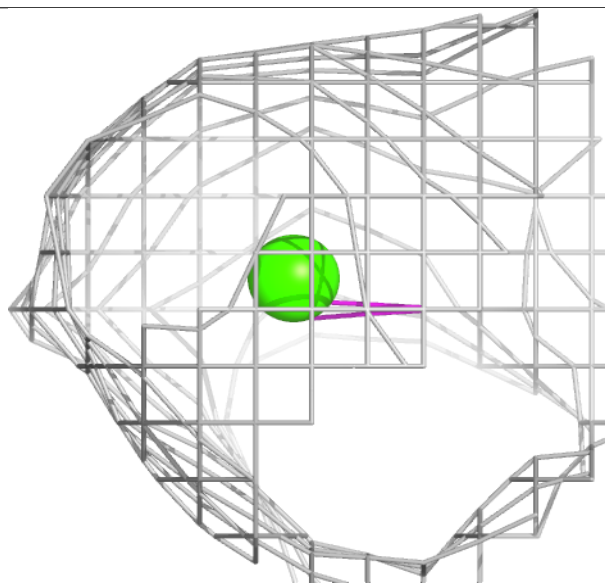
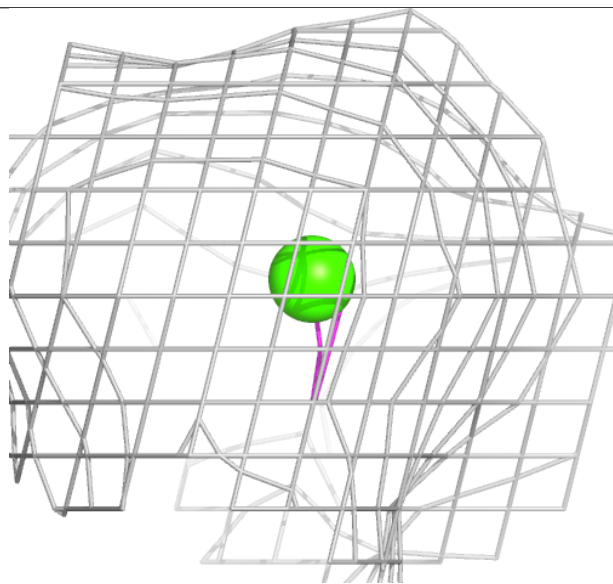
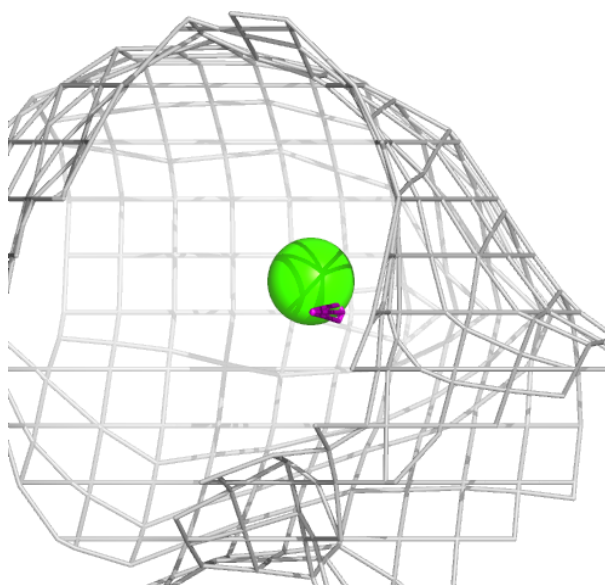
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and green (positive)



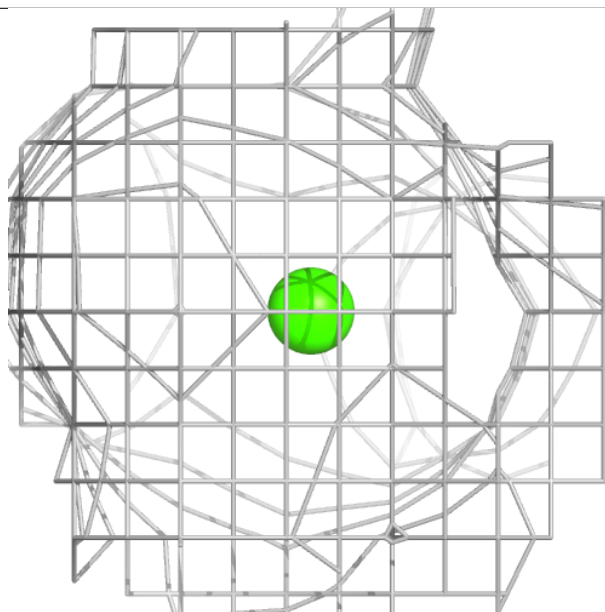
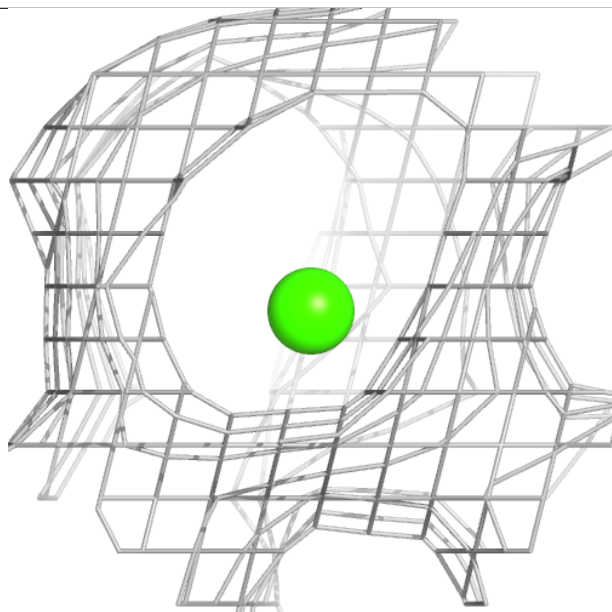
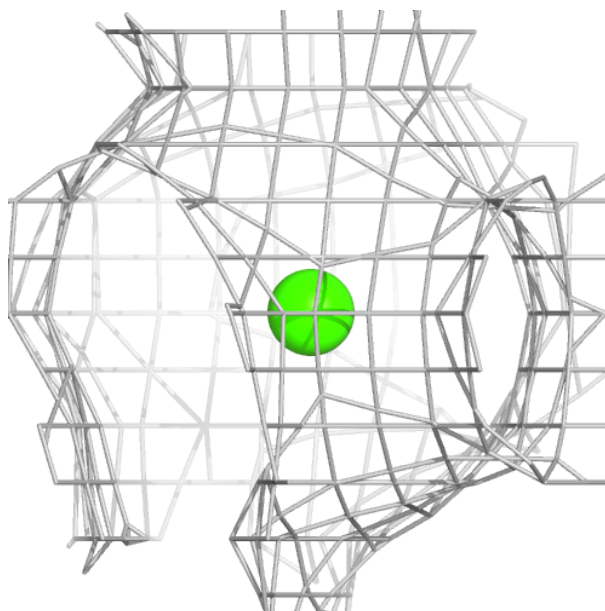
Electron density around CA D 602:

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and green (positive)



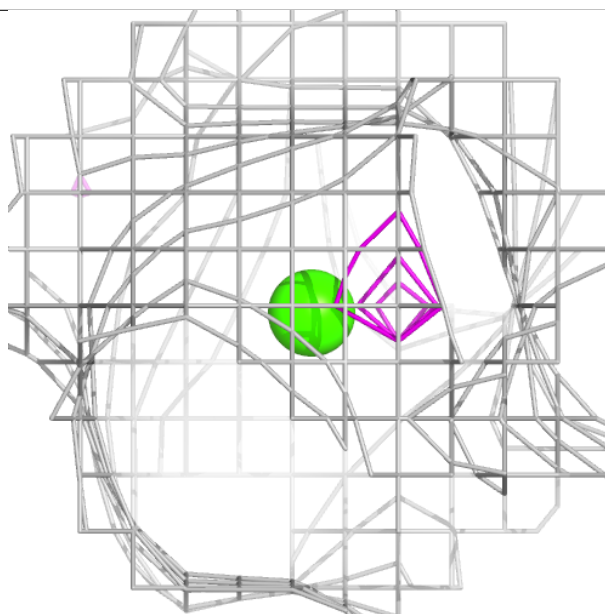
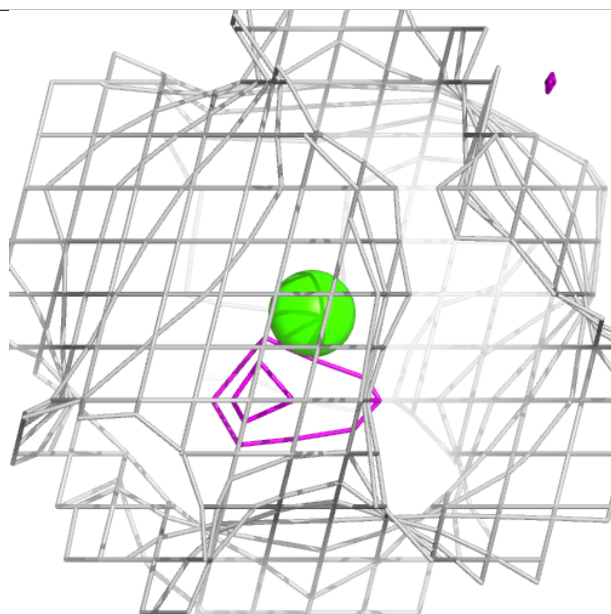
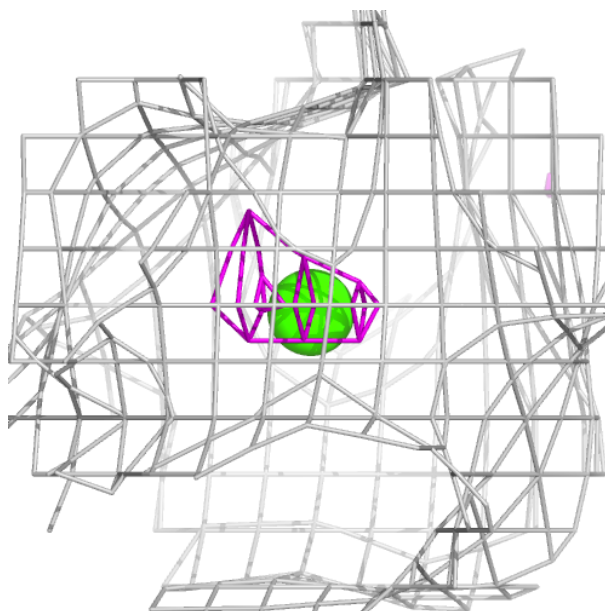
Electron density around CA D 601:

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and green (positive)



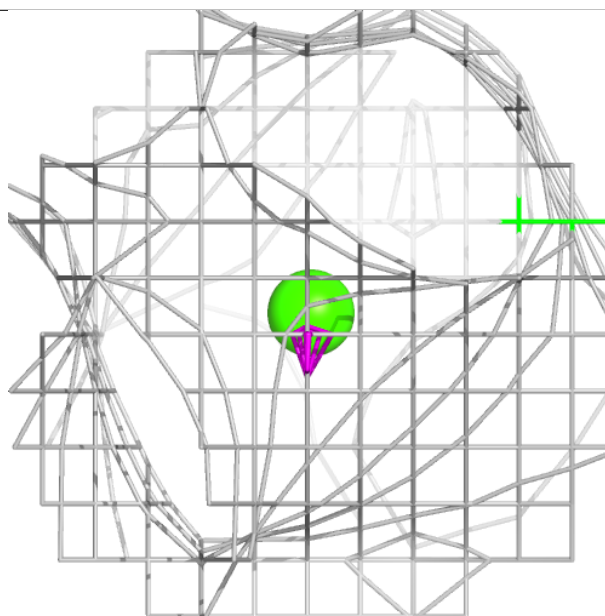
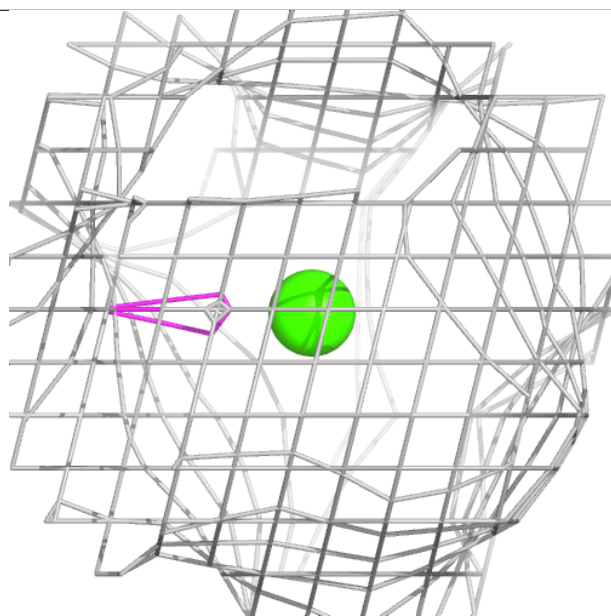
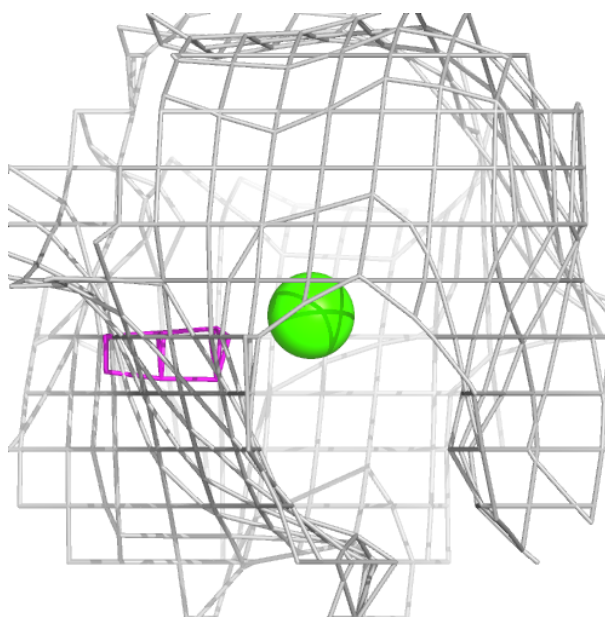
Electron density around CA C 601:

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



Electron density around CA 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)



6.5 Other polymers ⓘ

There are no such residues in this entry.