



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 22, 2025 – 01:12 AM EDT

PDB ID : 8DSJ / pdb_00008dsj
Title : Peptidylglycine alpha hydroxylating monooxygenase anaerobic
Authors : Arias, R.J.; Blackburn, N.J.
Deposited on : 2022-07-22
Resolution : 2.80 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 : 4.02b-467
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0rc1
EDS : 3.0
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4 : 9.0.006 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.42

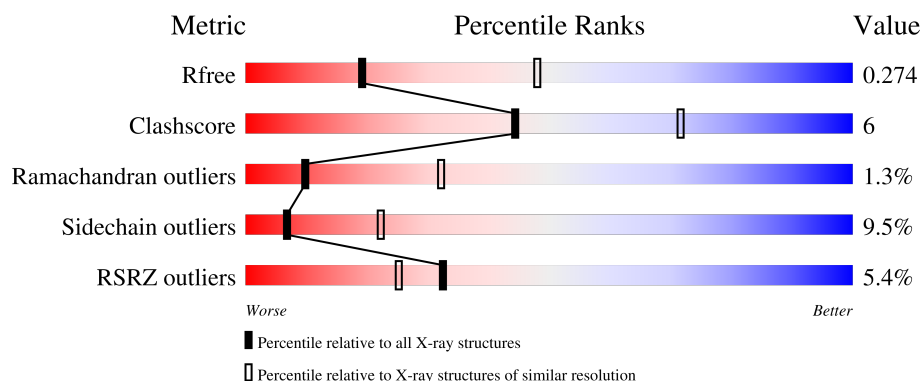
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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}	164625	3657 (2.80-2.80)
Clashscore	180529	4123 (2.80-2.80)
Ramachandran outliers	177936	4071 (2.80-2.80)
Sidechain outliers	177891	4073 (2.80-2.80)
RSRZ outliers	164620	3659 (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	310	 5% 80% 18% ..
1	K	310	 6% 77% 19% ..

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 4832 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.

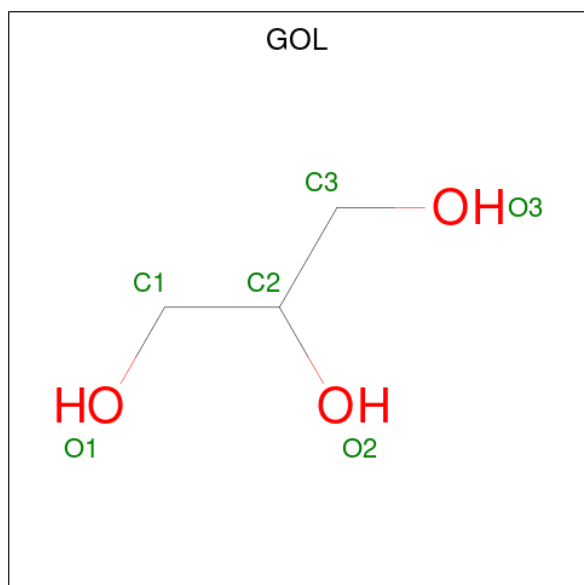
- Molecule 1 is a protein called Peptidylglycine alpha-amidating monooxygenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	309	Total	C	N	O	S	0	0	0
			2409	1534	408	441	26			
1	K	306	Total	C	N	O	S	0	0	0
			2387	1518	404	439	26			

- 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	3	Total	Cu	0	0
			3	3		
2	K	3	Total	Cu	0	0
			3	3		

- Molecule 3 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).

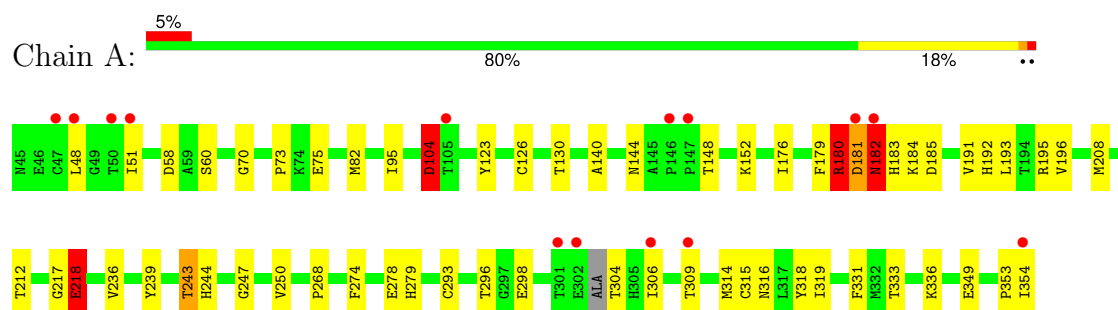


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total 6	C 3	O 3	0	0
3	A	1	Total 6	C 3	O 3	0	0
3	K	1	Total 6	C 3	O 3	0	0
3	K	1	Total 6	C 3	O 3	0	0
3	K	1	Total 6	C 3	O 3	0	0

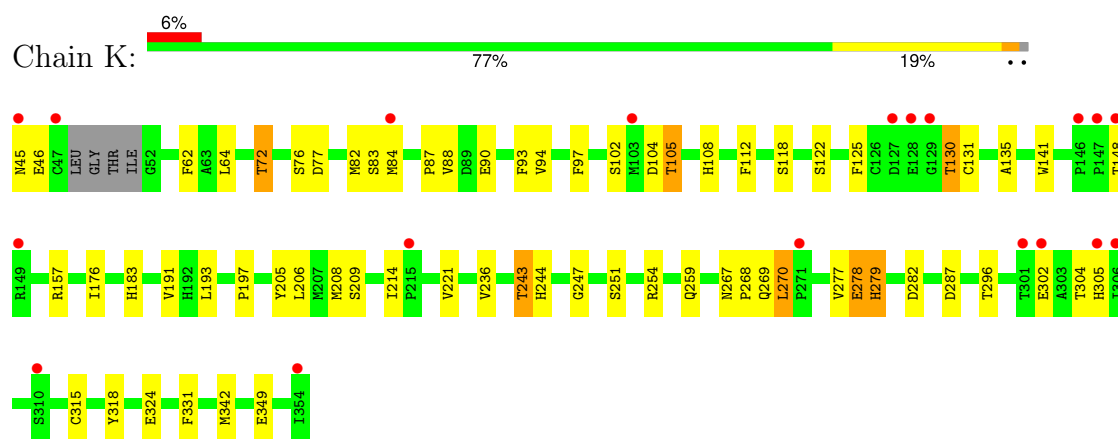
3 Residue-property plots [i](#)

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: Peptidylglycine alpha-amidating monooxygenase



- Molecule 1: Peptidylglycine alpha-amidating monooxygenase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	38.14Å 53.26Å 86.42Å 84.85° 89.96° 78.50°	Depositor
Resolution (Å)	37.40 – 2.80 37.40 – 2.80	Depositor EDS
% Data completeness (in resolution range)	100.0 (37.40-2.80) 100.0 (37.40-2.80)	Depositor EDS
R_{merge}	0.17	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.77 (at 2.81Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.197 , 0.275 0.201 , 0.274	Depositor DCC
R_{free} test set	812 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å ²)	36.1	Xtriage
Anisotropy	0.072	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 37.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	4832	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.25% 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: CU, GOL

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	0.63	0/2476	0.82	0/3366
1	K	0.64	0/2454	0.82	0/3337
All	All	0.63	0/4930	0.82	0/6703

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	K	0	1
All	All	0	2

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	104	ASP	Peptide
1	K	104	ASP	Peptide

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	2409	0	2327	27	1
1	K	2387	0	2293	29	1
2	A	3	0	0	0	0
2	K	3	0	0	0	0
3	A	12	0	16	0	0
3	K	18	0	24	1	0
All	All	4832	0	4660	56	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 6.

All (56) 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:243:THR:HG23	1:A:244:HIS:O	1.84	0.76
1:K:243:THR:HG21	1:K:247:GLY:HA3	1.70	0.73
1:K:278:GLU:HG2	1:K:279:HIS:H	1.54	0.73
1:A:180:ARG:O	1:A:182:ASN:N	2.16	0.72
1:A:180:ARG:C	1:A:182:ASN:H	1.95	0.69
1:K:102:SER:HB2	1:K:105:THR:HG22	1.76	0.67
1:K:72:THR:HG22	1:K:72:THR:O	1.95	0.66
1:K:214:ILE:HG23	1:K:221:VAL:HG21	1.76	0.66
1:A:179:PHE:HZ	1:A:185:ASP:HB2	1.62	0.63
1:A:243:THR:HG22	1:A:268:PRO:HB2	1.80	0.62
1:A:181:ASP:O	1:A:183:HIS:N	2.35	0.59
1:K:208:MET:HA	1:K:315:CYS:O	2.03	0.59
1:A:48:LEU:HD22	1:A:51:ILE:HD11	1.86	0.58
1:K:267:ASN:HB3	1:K:270:LEU:HD22	1.88	0.56
1:K:205:TYR:OH	3:K:406:GOL:O2	2.25	0.54
1:A:179:PHE:CZ	1:A:185:ASP:HB2	2.42	0.53
1:K:93:PHE:CE1	1:K:197:PRO:HA	2.44	0.52
1:K:112:PHE:CD1	1:K:135:ALA:HB1	2.44	0.52
1:A:180:ARG:C	1:A:182:ASN:N	2.58	0.51
1:K:349:GLU:OE2	1:K:349:GLU:N	2.41	0.51
1:K:72:THR:O	1:K:72:THR:CG2	2.60	0.50
1:A:104:ASP:OD1	1:A:104:ASP:N	2.43	0.49
1:A:243:THR:HA	1:A:314:MET:O	2.12	0.49
1:A:208:MET:HA	1:A:315:CYS:O	2.12	0.49
1:A:243:THR:HG21	1:A:247:GLY:HA3	1.95	0.49
1:K:72:THR:OG1	1:K:183:HIS:O	2.29	0.48
1:K:206:LEU:HD13	1:K:318:TYR:CE2	2.48	0.48
1:A:191:VAL:HG12	1:A:193:LEU:CD2	2.44	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:254:ARG:NH1	1:K:282:ASP:O	2.37	0.48
1:K:76:SER:HA	1:K:176:ILE:HG23	1.94	0.48
1:K:102:SER:CB	1:K:105:THR:HG22	2.44	0.48
1:A:250:VAL:HG22	1:A:293:CYS:SG	2.54	0.48
1:A:239:TYR:HA	1:A:318:TYR:O	2.14	0.48
1:K:130:THR:HG23	1:K:131:CYS:SG	2.54	0.47
1:A:236:VAL:HG13	1:A:319:ILE:HG23	1.97	0.47
1:A:123:TYR:C	1:A:123:TYR:CD1	2.89	0.45
1:K:236:VAL:HG12	1:K:277:VAL:HG21	1.99	0.45
1:K:87:PRO:O	1:K:88:VAL:HG13	2.17	0.44
1:K:83:SER:O	1:K:84:MET:SD	2.75	0.44
1:K:243:THR:HG22	1:K:268:PRO:HB2	1.99	0.44
1:A:73:PRO:HD2	1:A:179:PHE:CG	2.53	0.44
1:A:140:ALA:HB3	1:A:274:PHE:HD1	1.83	0.43
1:A:75:GLU:O	1:A:176:ILE:HG21	2.18	0.43
1:A:95:ILE:HD11	1:A:192:HIS:CD2	2.53	0.43
1:K:254:ARG:HD2	1:K:287:ASP:OD2	2.19	0.42
1:A:212:THR:HB	1:A:306:ILE:HD11	2.02	0.42
1:K:342:MET:HE2	1:K:342:MET:HB3	1.90	0.42
1:K:94:VAL:HG11	1:K:97:PHE:CZ	2.53	0.42
1:K:268:PRO:HG2	1:K:269:GLN:HE21	1.85	0.42
1:K:108:HIS:HB2	1:K:141:TRP:O	2.20	0.42
1:K:243:THR:HG23	1:K:244:HIS:N	2.35	0.42
1:A:60:SER:HB2	1:A:196:VAL:HG23	2.02	0.41
1:K:191:VAL:HG12	1:K:193:LEU:CD2	2.51	0.41
1:A:236:VAL:HG22	1:A:319:ILE:CG2	2.51	0.41
1:A:208:MET:HB3	1:A:316:ASN:ND2	2.36	0.41
1:A:217:GLY:O	1:A:218:GLU:C	2.58	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:A:182:ASN:O	1:K:125:PHE:CE1[1_556]	2.15	0.05

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	305/310 (98%)	282 (92%)	16 (5%)	7 (2%)	5	19
1	K	302/310 (97%)	280 (93%)	21 (7%)	1 (0%)	37	67
All	All	607/620 (98%)	562 (93%)	37 (6%)	8 (1%)	10	32

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	182	ASN
1	K	278	GLU
1	A	70	GLY
1	A	181	ASP
1	A	218	GLU
1	A	353	PRO
1	A	58	ASP
1	A	180	ARG

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	264/265 (100%)	240 (91%)	24 (9%)	7	24
1	K	261/265 (98%)	235 (90%)	26 (10%)	6	20
All	All	525/530 (99%)	475 (90%)	50 (10%)	7	22

All (50) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	82	MET
1	A	104	ASP
1	A	126	CYS
1	A	130	THR
1	A	144	ASN
1	A	148	THR
1	A	152	LYS
1	A	180	ARG
1	A	182	ASN
1	A	184	LYS
1	A	195	ARG
1	A	218	GLU
1	A	243	THR
1	A	278	GLU
1	A	279	HIS
1	A	296	THR
1	A	298	GLU
1	A	304	THR
1	A	309	THR
1	A	331	PHE
1	A	333	THR
1	A	336	LYS
1	A	349	GLU
1	A	354	ILE
1	K	45	ASN
1	K	46	GLU
1	K	62	PHE
1	K	64	LEU
1	K	72	THR
1	K	77	ASP
1	K	82	MET
1	K	90	GLU
1	K	105	THR
1	K	118	SER
1	K	122	SER
1	K	130	THR
1	K	148	THR
1	K	157	ARG
1	K	209	SER
1	K	243	THR
1	K	251	SER
1	K	259	GLN
1	K	270	LEU

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Mol	Chain	Res	Type
1	K	279	HIS
1	K	296	THR
1	K	302	GLU
1	K	304	THR
1	K	305	HIS
1	K	324	GLU
1	K	331	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	316	ASN
1	K	45	ASN
1	K	222	ASN
1	K	259	GLN
1	K	316	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 11 ligands modelled in this entry, 6 are monoatomic - leaving 5 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
3	GOL	A	405	-	5,5,5	0.15	0	5,5,5	0.35	0
3	GOL	K	404	-	5,5,5	0.11	0	5,5,5	0.30	0
3	GOL	K	405	-	5,5,5	0.13	0	5,5,5	0.38	0
3	GOL	A	404	-	5,5,5	0.16	0	5,5,5	0.45	0
3	GOL	K	406	-	5,5,5	0.14	0	5,5,5	0.33	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
3	GOL	A	405	-	-	2/4/4/4	-
3	GOL	K	404	-	-	1/4/4/4	-
3	GOL	K	405	-	-	3/4/4/4	-
3	GOL	A	404	-	-	2/4/4/4	-
3	GOL	K	406	-	-	4/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (12) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	405	GOL	O1-C1-C2-C3
3	K	406	GOL	O1-C1-C2-C3
3	A	404	GOL	C1-C2-C3-O3
3	K	405	GOL	O1-C1-C2-C3
3	K	405	GOL	C1-C2-C3-O3
3	K	406	GOL	C1-C2-C3-O3
3	A	405	GOL	O1-C1-C2-O2
3	K	405	GOL	O1-C1-C2-O2
3	K	406	GOL	O2-C2-C3-O3
3	A	404	GOL	O2-C2-C3-O3
3	K	404	GOL	O1-C1-C2-C3
3	K	406	GOL	O1-C1-C2-O2

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	K	406	GOL	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	309/310 (99%)	0.06	14 (4%) 39 31	10, 33, 84, 119	0
1	K	306/310 (98%)	0.16	19 (6%) 28 21	12, 34, 90, 126	0
All	All	615/620 (99%)	0.11	33 (5%) 32 25	10, 34, 89, 126	0

All (33) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	354	ILE	3.9
1	K	147	PRO	3.8
1	K	354	ILE	3.4
1	K	305	HIS	3.4
1	K	146	PRO	3.3
1	A	48	LEU	3.1
1	A	182	ASN	3.1
1	A	146	PRO	3.1
1	K	306	ILE	3.0
1	A	306	ILE	3.0
1	A	51	ILE	2.8
1	K	128	GLU	2.8
1	K	45	ASN	2.8
1	A	147	PRO	2.5
1	A	302	GLU	2.5
1	K	148	THR	2.5
1	K	271	PRO	2.4
1	A	181	ASP	2.4
1	K	301	THR	2.3
1	K	310	SER	2.3
1	K	47	CYS	2.2
1	K	84	MET	2.2
1	A	105	THR	2.2
1	A	47	CYS	2.1

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Mol	Chain	Res	Type	RSRZ
1	K	215	PRO	2.1
1	A	301	THR	2.1
1	K	127	ASP	2.1
1	K	103	MET	2.1
1	A	309	THR	2.0
1	K	302	GLU	2.0
1	A	50	THR	2.0
1	K	149	ARG	2.0
1	K	129	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 monosaccharides 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.

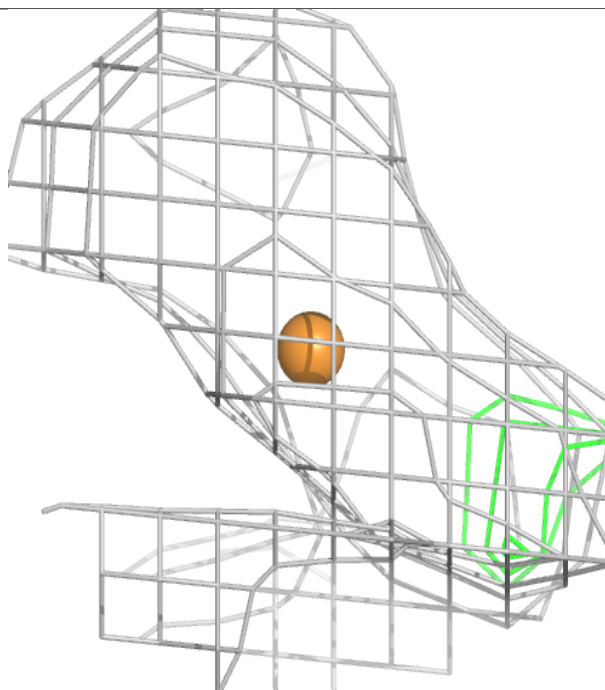
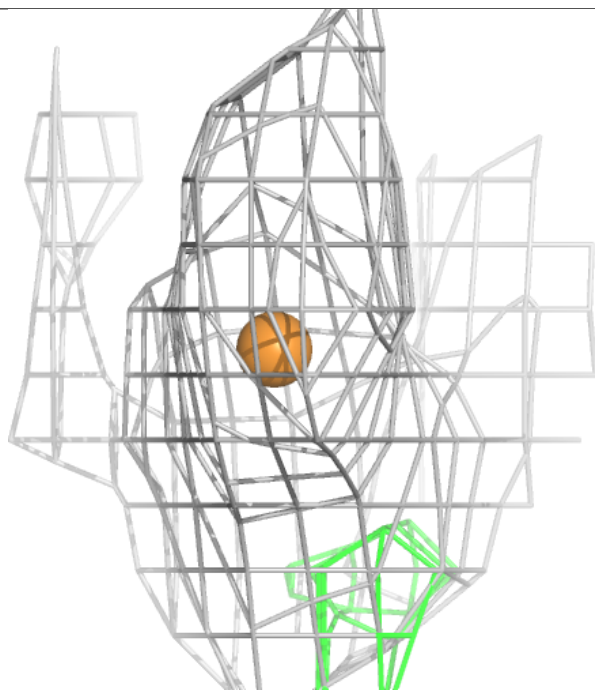
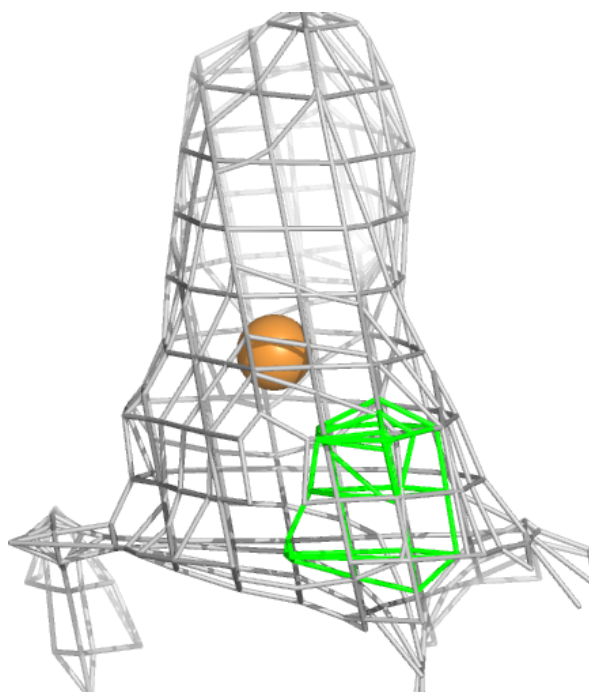
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	CU	K	403	1/1	0.74	0.16	124,124,124,124	0
2	CU	A	403	1/1	0.85	0.12	102,102,102,102	0
3	GOL	A	404	6/6	0.89	0.12	26,29,30,31	0
3	GOL	A	405	6/6	0.89	0.12	35,39,40,41	0
3	GOL	K	404	6/6	0.89	0.19	66,68,81,84	0
3	GOL	K	406	6/6	0.89	0.14	51,54,59,60	0
3	GOL	K	405	6/6	0.91	0.11	38,40,41,43	0
2	CU	K	401	1/1	0.93	0.12	79,79,79,79	0
2	CU	K	402	1/1	0.95	0.05	62,62,62,62	0
2	CU	A	402	1/1	0.97	0.04	61,61,61,61	0
2	CU	A	401	1/1	0.97	0.09	65,65,65,65	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.

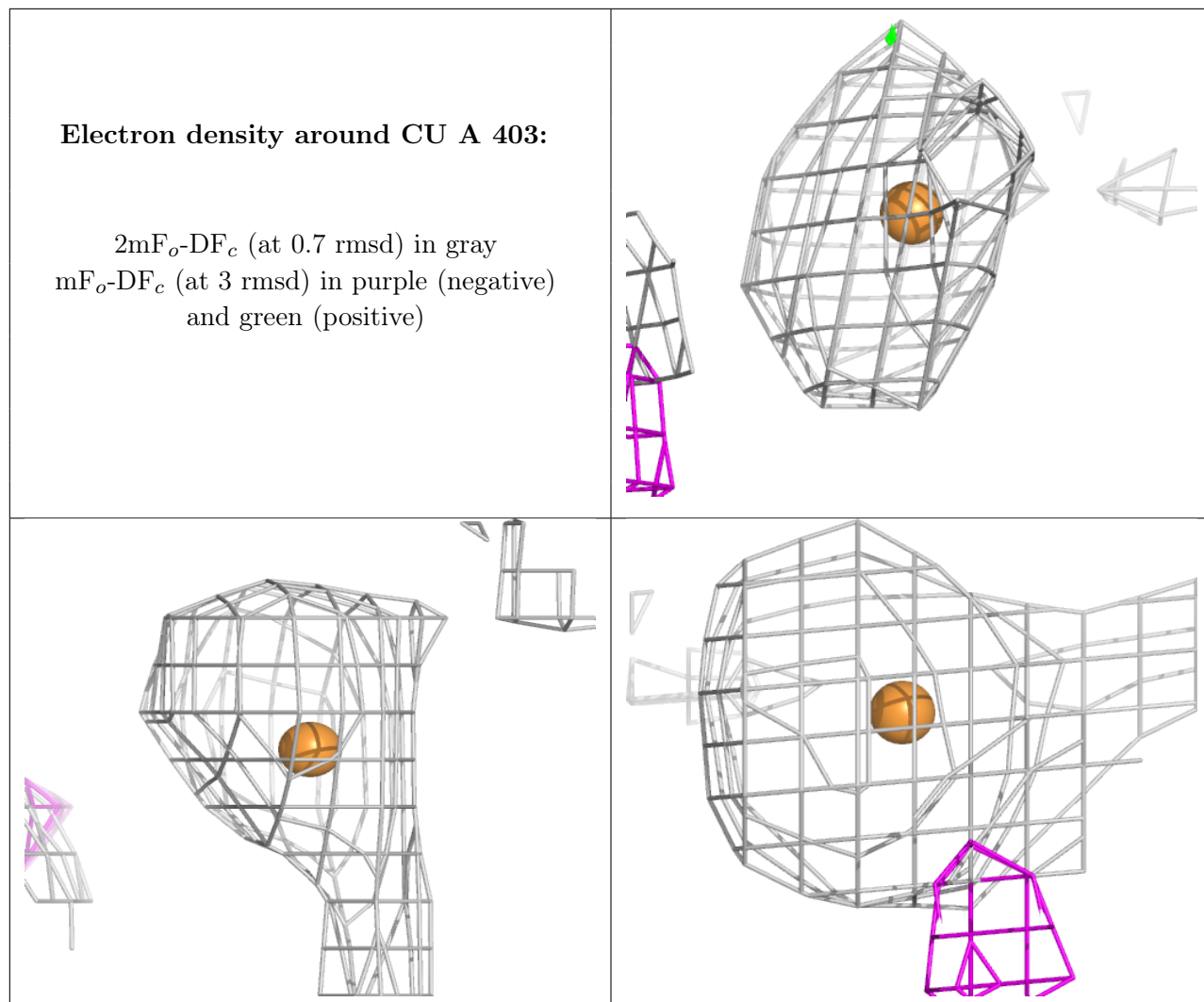
Electron density around CU K 403:

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



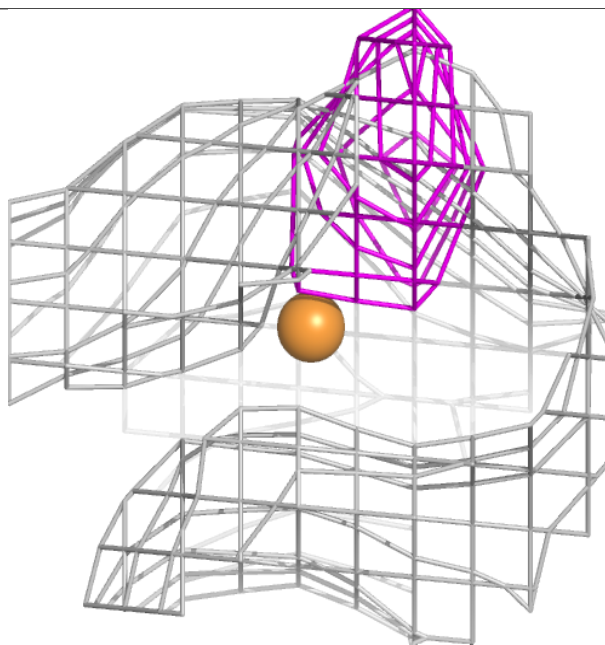
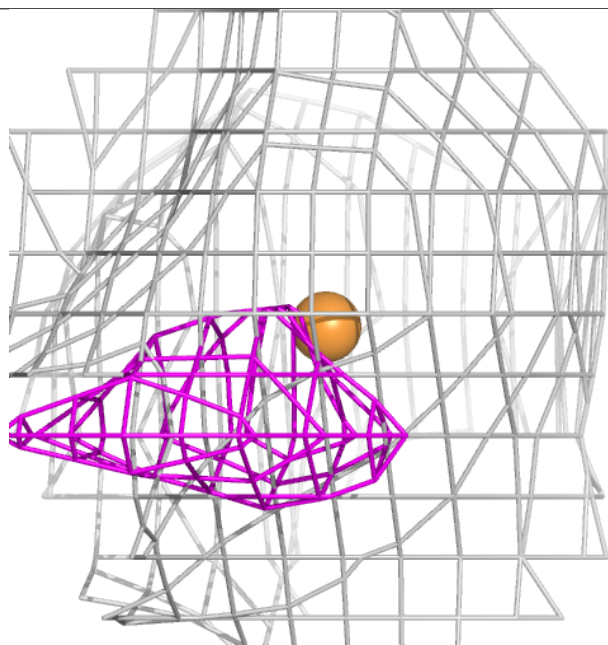
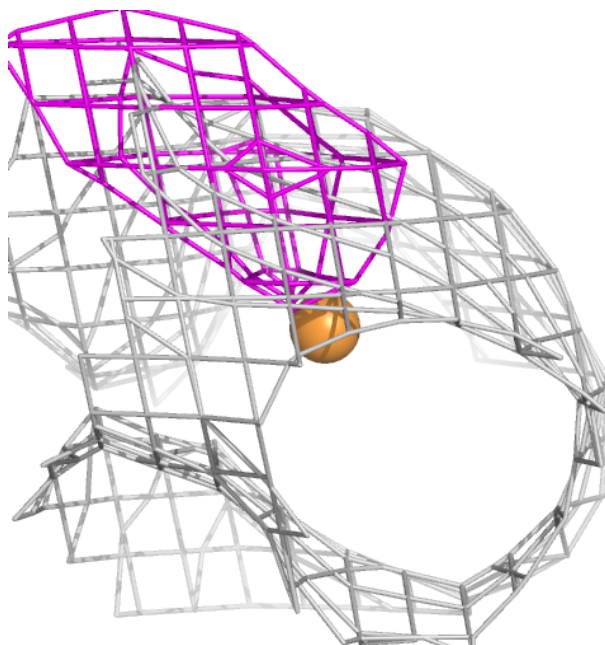
Electron density around CU A 403:

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



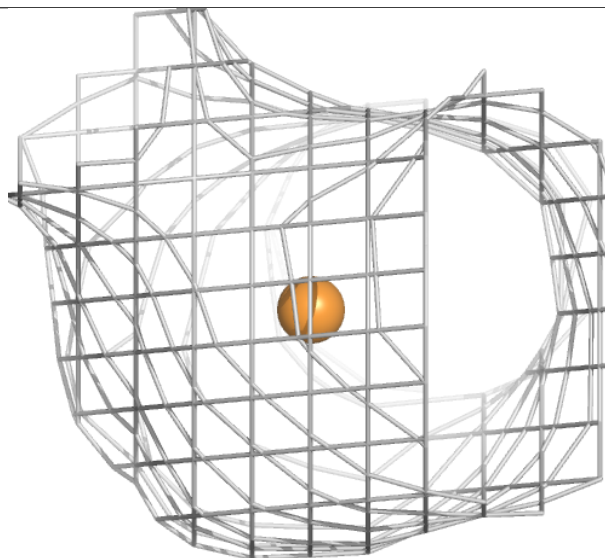
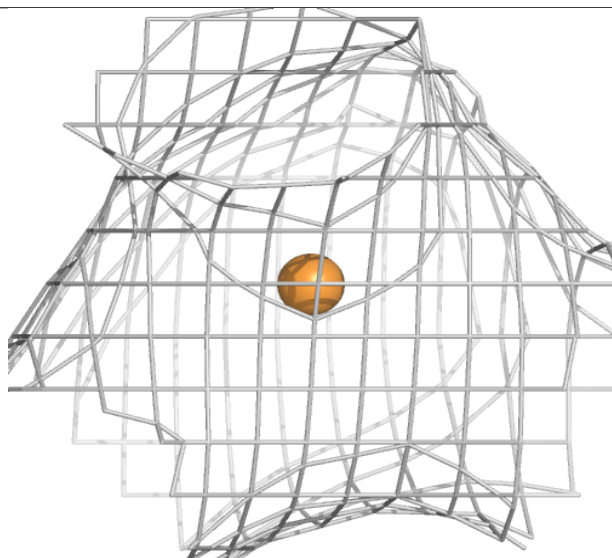
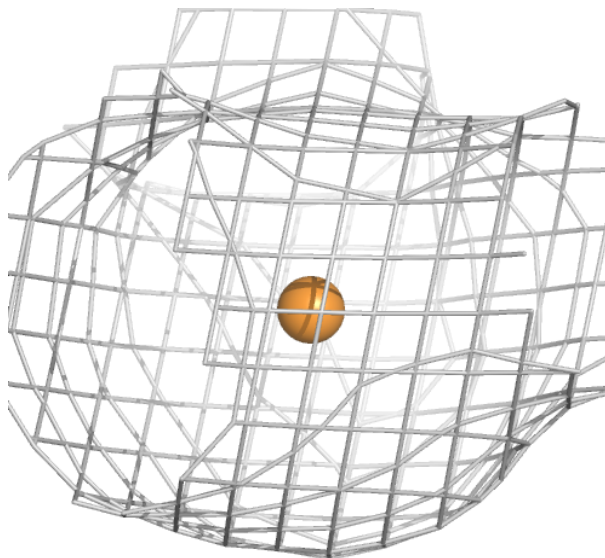
Electron density around CU K 401:

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



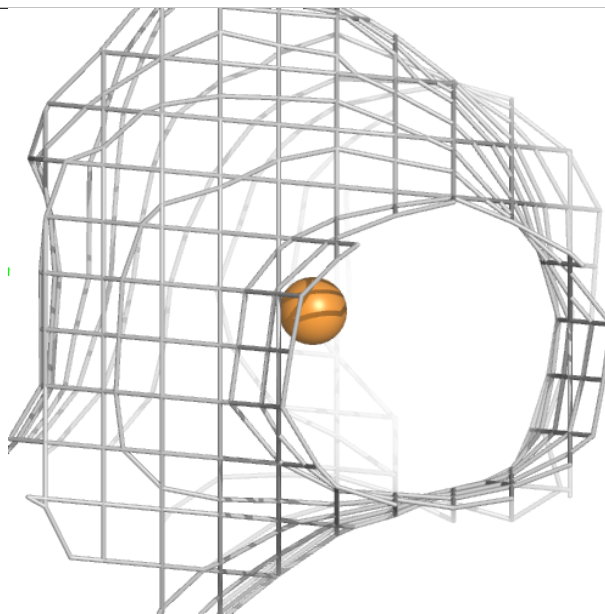
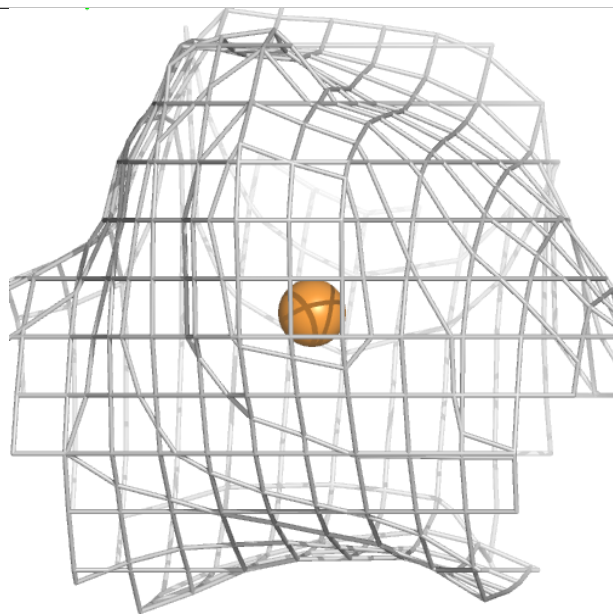
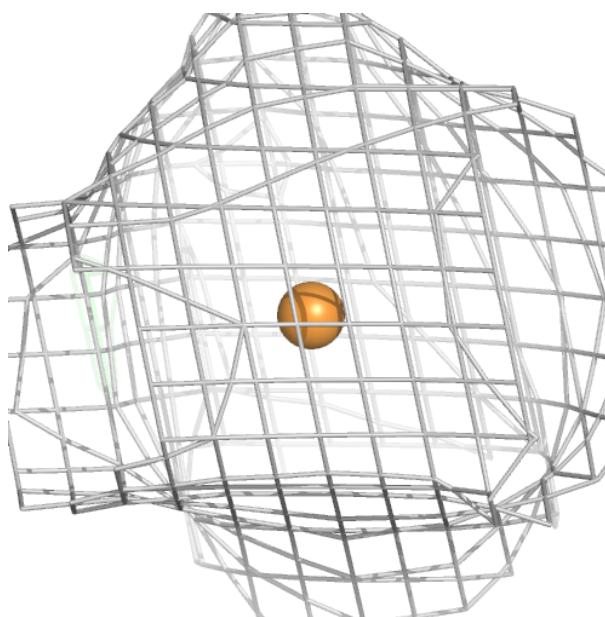
Electron density around CU K 402:

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



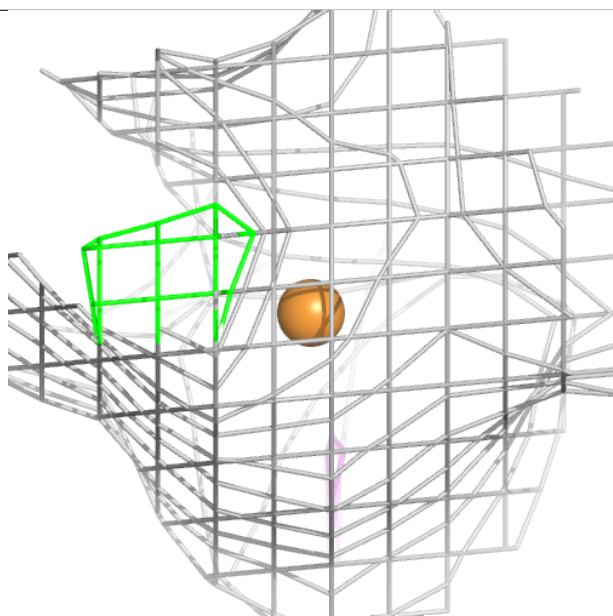
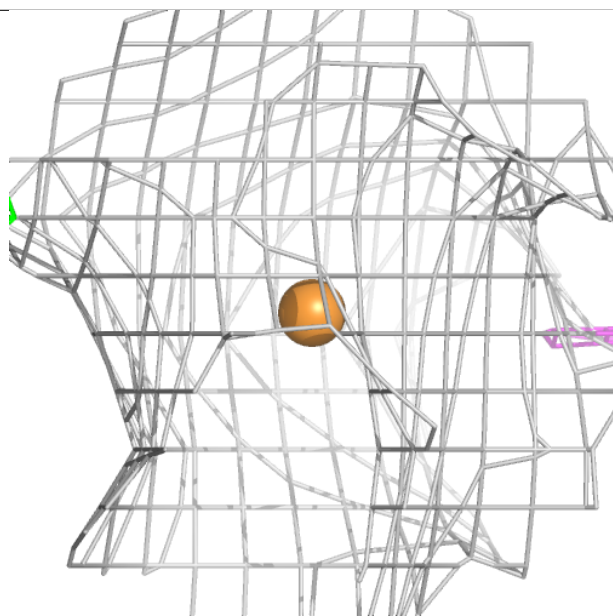
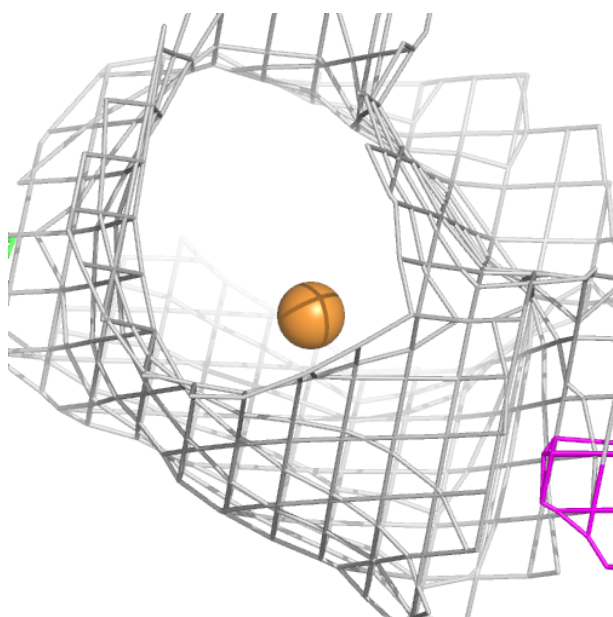
Electron density around CU A 402:

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



Electron density around CU A 401:

$2mF_o-DF_c$ (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.