



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 10:17 AM UTC

PDB ID : 6Z8O / pdb_00006z8o
Title : Structure of [NiFeSe] hydrogenase G491A variant from *Desulfovibrio vulgaris* Hildenborough pressurized with Krypton gas - structure G491A-Kr
Authors : Zacarias, S.; Temporao, A.; Carpentier, P.; van der Linden, P.; Pereira, I.A.C.; Matias, P.M.
Deposited on : 2020-06-02
Resolution : 2.20 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

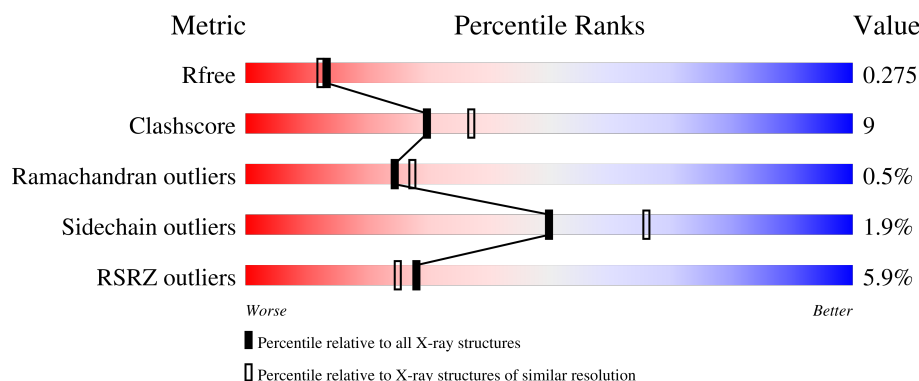
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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	283	 6% 78% 19% ..
1	C	283	 3% 76% 22% .
2	B	484	 7% 79% 21% .
2	D	484	 6% 80% 19% ..

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard

residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
5	KR	A	306	-	-	X	-
5	KR	C	310	-	-	X	-
6	FCO	D	501	-	-	X	-

2 Entry composition

There are 11 unique types of molecules in this entry. The entry contains 23485 atoms, of which 11622 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 Periplasmic [NiFeSe] hydrogenase, small subunit.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	277	Total	C	H	N	O	S	0	3	0
			4130	1335	2037	349	389	20			
1	C	278	Total	C	H	N	O	S	0	3	0
			4145	1340	2043	350	392	20			

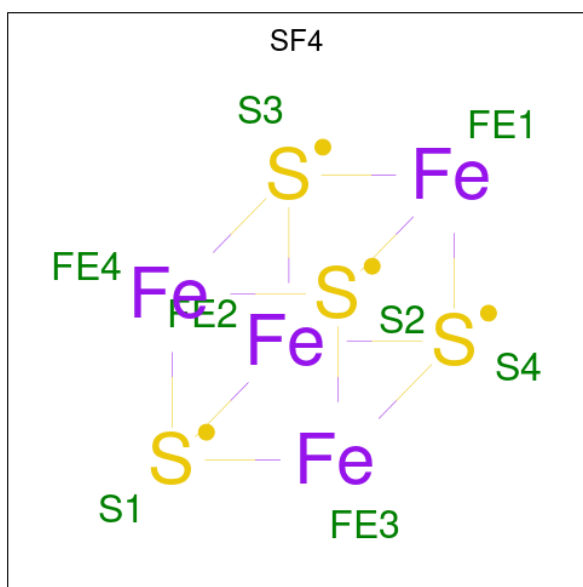
- Molecule 2 is a protein called Periplasmic [NiFeSe] hydrogenase, large subunit, selenocysteine-containing.

Mol	Chain	Residues	Atoms							ZeroOcc	AltConf	Trace
2	B	481	Total	C	H	N	O	S	Se	0	5	0
			7532	2401	3772	654	683	19	3			
2	D	481	Total	C	H	N	O	S	Se	0	5	0
			7530	2401	3770	654	683	19	3			

There are 2 discrepancies between the modelled and reference sequences:

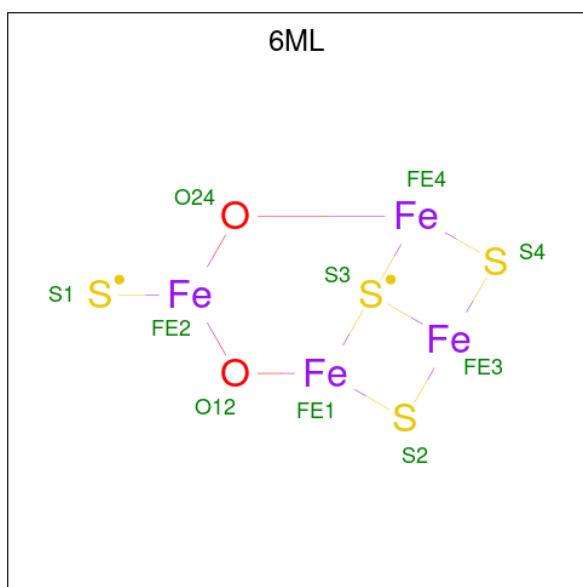
Chain	Residue	Modelled	Actual	Comment	Reference
B	491	ALA	GLY	engineered mutation	UNP Q72AS3
D	491	ALA	GLY	engineered mutation	UNP Q72AS3

- Molecule 3 is IRON/SULFUR CLUSTER (CCD ID: SF4) (formula: Fe₄S₄).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	Fe	S	0	0
			8	4	4		
3	A	1	Total	Fe	S	0	0
			8	4	4		
3	A	1	Total	Fe	S	0	1
			8	4	4		
3	C	1	Total	Fe	S	0	0
			8	4	4		
3	C	1	Total	Fe	S	0	0
			8	4	4		
3	C	1	Total	Fe	S	0	1
			8	4	4		

- Molecule 4 is oxygen-damaged SF4 (CCD ID: 6ML) (formula: $\text{Fe}_4\text{O}_2\text{S}_4$).

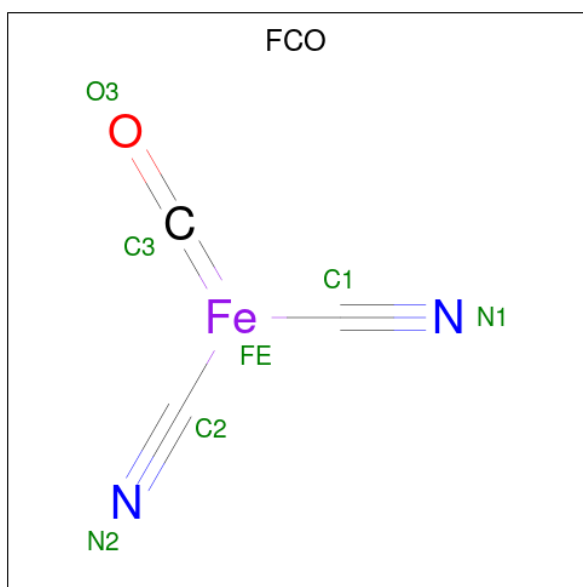


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	Fe	O	S	0	1
			10	4	2	4		
4	C	1	Total	Fe	O	S	0	1
			10	4	2	4		

- Molecule 5 is KRYPTON (CCD ID: KR) (formula: Kr) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	6	Total	Kr	0	0
			6	6		
5	B	5	Total	Kr	0	2
			5	5		
5	C	6	Total	Kr	0	0
			6	6		
5	D	3	Total	Kr	0	0
			3	3		

- Molecule 6 is CARBONMONOXIDE-(DICYANO) IRON (CCD ID: FCO) (formula: C₃FeN₂O).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
6	B	1	Total	C	Fe	N	O	0	0
			7	3	1	2	1		
6	D	1	Total	C	Fe	N	O	0	0
			7	3	1	2	1		

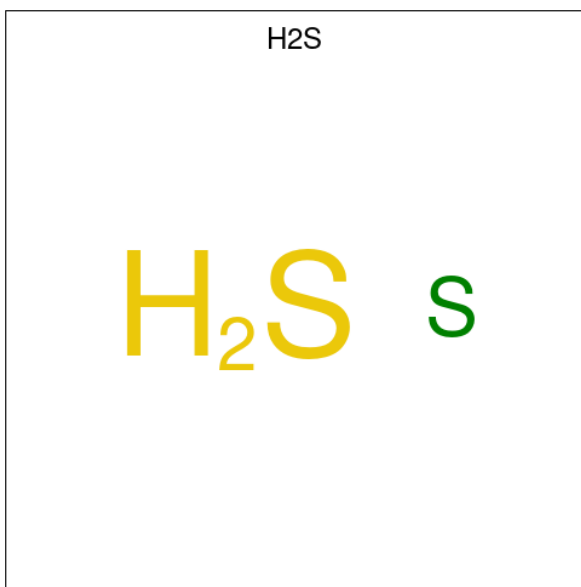
- Molecule 7 is NICKEL (II) ION (CCD ID: NI) (formula: Ni).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	B	1	Total	Ni	0	0
			1	1		
7	D	1	Total	Ni	0	0
			1	1		

- Molecule 8 is FE (II) ION (CCD ID: FE2) (formula: Fe).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	B	1	Total	Fe	0	0
			1	1		
8	D	1	Total	Fe	0	0
			1	1		

- Molecule 9 is HYDROSULFURIC ACID (CCD ID: H2S) (formula: H₂S).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
9	B	1	Total S 1 1	0	0
9	D	1	Total S 1 1	0	0

- Molecule 10 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
10	B	1	Total Cl 1 1	0	0
10	D	1	Total Cl 1 1	0	0

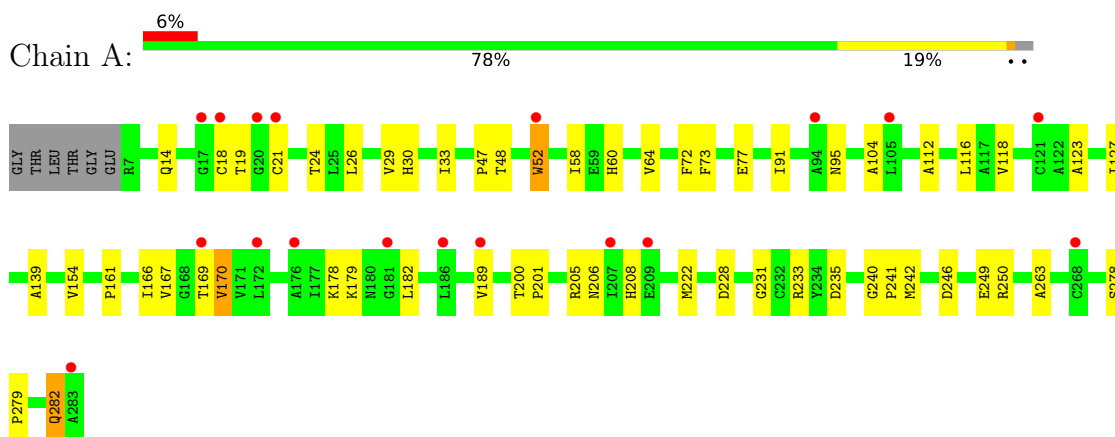
- Molecule 11 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
11	A	6	Total O 6 6	0	0
11	B	9	Total O 9 9	0	0
11	C	11	Total O 11 11	0	0
11	D	12	Total O 12 12	0	0

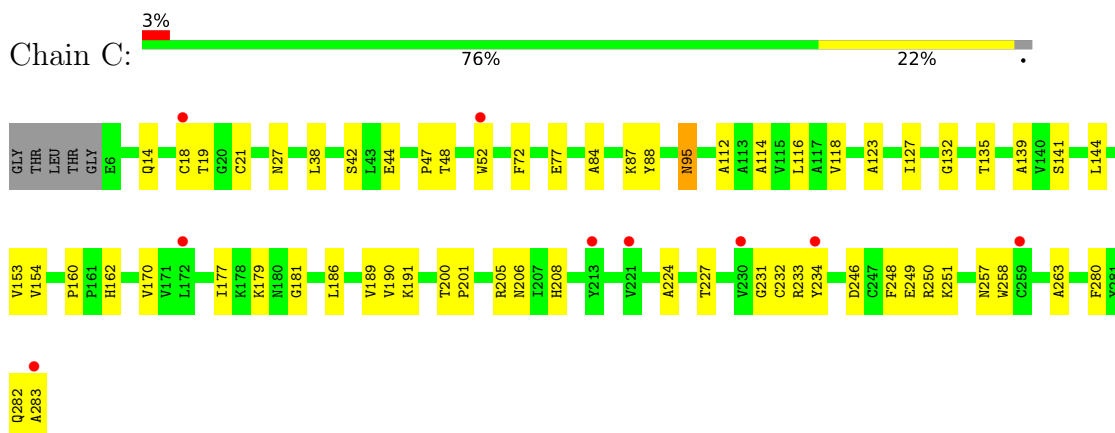
3 Residue-property plots [i](#)

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.

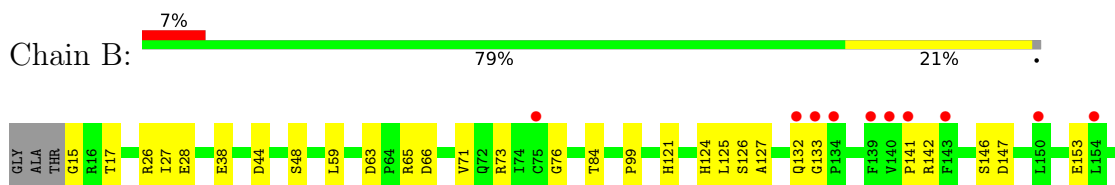
- Molecule 1: Periplasmic [NiFeSe] hydrogenase, small subunit

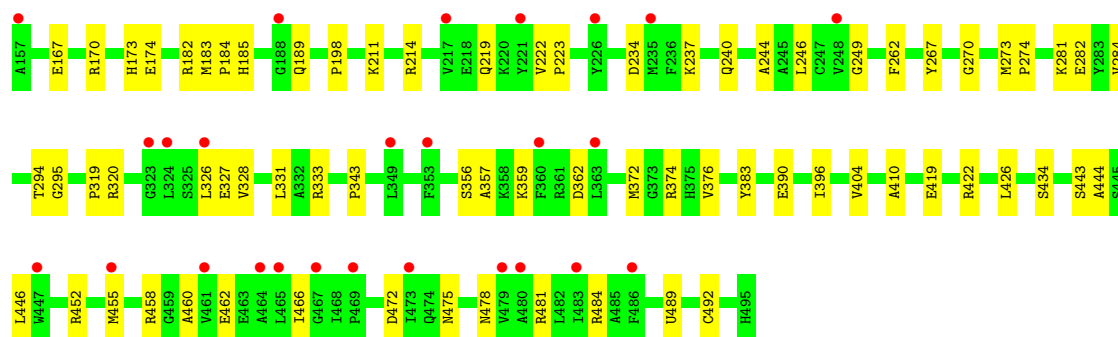


- Molecule 1: Periplasmic [NiFeSe] hydrogenase, small subunit

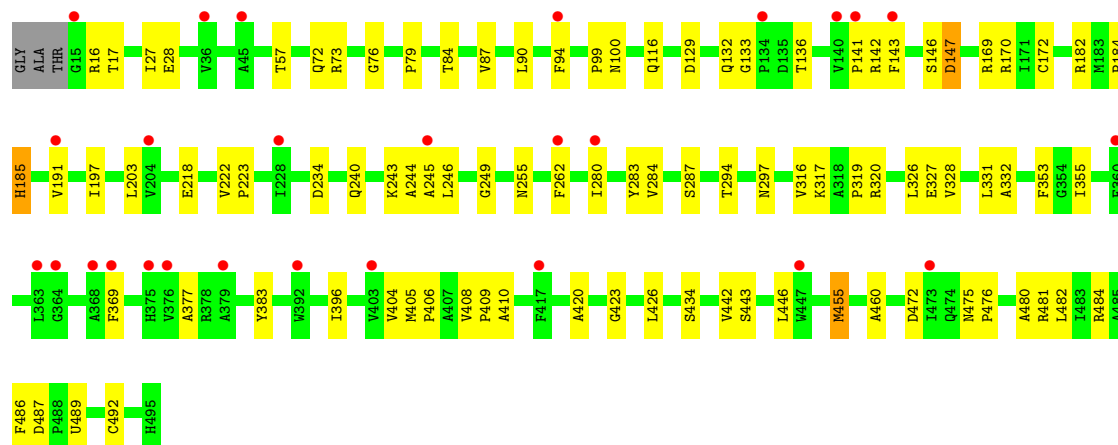
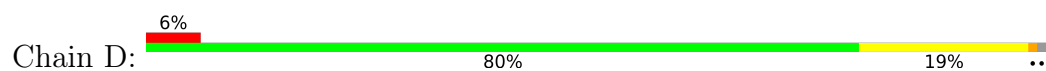


- Molecule 2: Periplasmic [NiFeSe] hydrogenase, large subunit, selenocysteine-containing





- Molecule 2: Periplasmic [NiFeSe] hydrogenase, large subunit, selenocysteine-containing



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	63.72Å 97.02Å 121.32Å 90.00° 104.66° 90.00°	Depositor
Resolution (Å)	39.12 – 2.20 39.12 – 2.20	Depositor EDS
% Data completeness (in resolution range)	46.5 (39.12-2.20) 46.7 (39.12-2.20)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.01 (at 2.20Å)	Xtriage
Refinement program	PHENIX 1.18.2_3874	Depositor
R, R_{free}	0.226 , 0.275 0.227 , 0.275	Depositor DCC
R_{free} test set	1686 reflections (2.31%)	wwPDB-VP
Wilson B-factor (Å ²)	46.9	Xtriage
Anisotropy	0.094	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.42 , 33.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.026 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	23485	wwPDB-VP
Average B, all atoms (Å ²)	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 33.69 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 7.6786e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: FE2, SEC, KR, H2S, FCO, CL, NI, 6ML, SF4

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	0.19	0/2157	0.47	0/2931
1	C	0.17	0/2166	0.46	0/2943
2	B	0.19	0/3852	0.48	0/5206
2	D	0.21	0/3852	0.50	0/5206
All	All	0.19	0/12027	0.48	0/16286

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2093	2037	2030	37	2
1	C	2102	2043	2036	41	1
2	B	3760	3772	3761	64	1
2	D	3760	3770	3761	73	2
3	A	24	0	0	1	0
3	C	24	0	0	1	0
4	A	10	0	0	2	0
4	C	10	0	0	2	0
5	A	6	0	0	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	B	5	0	0	1	0
5	C	6	0	0	4	0
5	D	3	0	0	0	0
6	B	7	0	0	1	0
6	D	7	0	0	3	0
7	B	1	0	0	0	0
7	D	1	0	0	0	0
8	B	1	0	0	0	0
8	D	1	0	0	0	0
9	B	1	0	0	1	0
9	D	1	0	0	0	0
10	B	1	0	0	0	0
10	D	1	0	0	0	0
11	A	6	0	0	0	0
11	B	9	0	0	3	0
11	C	11	0	0	1	0
11	D	12	0	0	4	0
All	All	11863	11622	11588	200	3

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 9.

All (200) 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)
2:D:489[A]:SEC:SE	6:D:501:FCO:C1	2.72	0.88
2:B:489[A]:SEC:SE	6:B:501:FCO:C1	2.71	0.88
2:D:486:PHE:O	11:D:601:HOH:O	2.05	0.75
2:D:169:ARG:NH1	11:D:602:HOH:O	2.20	0.72
2:B:189:GLN:O	11:B:602:HOH:O	2.08	0.72
2:B:66:ASP:OD2	11:B:601:HOH:O	2.07	0.71
2:D:94:PHE:CD2	2:D:191:VAL:HG21	2.25	0.71
2:B:214:ARG:NH2	2:B:390:GLU:OE1	2.23	0.70
2:B:282:GLU:OE2	11:B:603:HOH:O	2.12	0.68
1:C:18:CYS:O	2:D:76:GLY:N	2.25	0.67
1:C:21[B]:CYS:N	4:C:304[B]:6ML:S1	2.69	0.66
1:C:132:GLY:O	2:D:57:THR:OG1	2.10	0.66
2:D:94:PHE:CD1	2:D:409:PRO:HD2	2.32	0.65
2:D:240:GLN:HA	2:D:369:PHE:O	1.97	0.65
1:A:19:THR:HG22	1:A:19:THR:O	1.98	0.63
2:B:357:ALA:HB1	2:B:362:ASP:HB2	1.81	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:280:PHE:O	11:C:401:HOH:O	2.16	0.62
2:D:94:PHE:CE1	2:D:408:VAL:HA	2.35	0.61
2:D:94:PHE:HE1	2:D:408:VAL:HA	1.66	0.61
3:A:302:SF4:S2	2:B:182:ARG:NH1	2.74	0.60
1:C:19:THR:HG22	1:C:19:THR:O	2.02	0.58
2:D:203:LEU:HD21	2:D:396:ILE:HG21	1.86	0.58
2:B:333:ARG:NH1	2:B:419:GLU:OE1	2.31	0.56
1:C:123:ALA:O	1:C:139:ALA:HB1	2.05	0.56
2:B:237:LYS:C	2:B:240:GLN:HE22	2.14	0.56
1:C:21[B]:CYS:SG	1:C:118:VAL:HG12	2.46	0.56
2:D:203:LEU:CD2	2:D:396:ILE:HG21	2.36	0.56
1:C:177:ILE:O	1:C:181:GLY:N	2.39	0.56
1:C:234:TYR:CE2	1:C:283:ALA:HB1	2.41	0.55
2:D:94:PHE:CD1	2:D:409:PRO:CD	2.89	0.55
2:B:234:ASP:N	2:B:234:ASP:OD1	2.39	0.55
2:B:472:ASP:OD2	2:B:475:ASN:ND2	2.37	0.54
2:B:410:ALA:HA	2:B:434:SER:OG	2.08	0.54
2:B:28:GLU:O	2:B:489[A]:SEC:HA	2.09	0.53
2:D:94:PHE:CE1	2:D:409:PRO:HD2	2.43	0.53
2:D:443:SER:OG	2:D:492:CYS:O	2.19	0.53
2:B:170:ARG:NH1	2:B:174:GLU:OE2	2.42	0.53
2:D:280:ILE:O	2:D:297:ASN:ND2	2.38	0.53
2:B:489[C]:SEC:CB	9:B:504:H2S:S	2.98	0.52
1:C:42:SER:OG	1:C:44:GLU:OE2	2.28	0.52
2:D:28:GLU:HB3	2:D:489[A]:SEC:CB	2.40	0.52
2:D:99:PRO:HD3	2:D:404:VAL:HG13	1.90	0.52
1:A:21[B]:CYS:SG	1:A:118:VAL:HG12	2.50	0.51
1:A:170:VAL:HG21	5:A:307:KR:KR	2.70	0.51
2:D:443:SER:O	2:D:446:LEU:HB3	2.10	0.51
2:B:249:GLY:HA2	2:B:262:PHE:O	2.11	0.50
2:B:267:TYR:OH	2:B:270:GLY:HA2	2.10	0.50
1:A:19:THR:HA	1:A:77:GLU:OE2	2.11	0.50
2:B:15:GLY:O	2:B:38:GLU:HA	2.11	0.50
2:B:284:VAL:HG22	2:B:294:THR:HB	1.94	0.50
2:D:249:GLY:HA2	2:D:262:PHE:O	2.11	0.50
1:C:48:THR:HG23	2:D:27:ILE:O	2.10	0.50
1:A:231:GLY:O	1:A:233:ARG:NH1	2.40	0.50
2:D:129:ASP:HB3	2:D:481:ARG:HG2	1.94	0.50
2:B:99:PRO:HD3	2:B:404:VAL:HG13	1.94	0.49
1:C:170:VAL:HG21	5:C:308:KR:KR	2.74	0.49
2:B:63:ASP:OD2	2:B:65:ARG:NH2	2.39	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:458:ARG:HD2	2:B:462:GLU:HB3	1.94	0.49
1:A:18:CYS:O	2:B:76:GLY:N	2.45	0.49
1:A:205:ARG:HH22	1:A:249:GLU:CD	2.21	0.49
2:D:482:LEU:C	2:D:482:LEU:HD13	2.38	0.49
1:A:123:ALA:O	1:A:139:ALA:HB1	2.12	0.49
1:A:21[A]:CYS:HB3	1:A:118:VAL:HG12	1.94	0.48
1:C:72:PHE:CE1	1:C:112:ALA:HB2	2.48	0.48
2:B:331:LEU:HD22	2:B:374:ARG:HG2	1.96	0.48
2:B:443:SER:OG	2:B:492:CYS:O	2.25	0.48
1:A:29:VAL:O	1:A:30:HIS:HD2	1.98	0.47
2:D:410:ALA:HA	2:D:434:SER:OG	2.14	0.47
1:C:52:TRP:HB2	2:D:141:PRO:O	2.15	0.47
1:A:14:GLN:OE1	1:A:47:PRO:HG2	2.15	0.47
2:D:353:PHE:HB3	2:D:355:ILE:HD12	1.96	0.47
1:A:200:THR:N	1:A:201:PRO:CD	2.78	0.47
2:B:121:HIS:NE2	2:B:484:ARG:O	2.45	0.47
1:C:206:ASN:OD1	1:C:208:HIS:N	2.47	0.47
1:C:114:ALA:HB3	5:C:307:KR:KR	2.75	0.47
1:A:73:PHE:CE1	5:A:306:KR:KR	3.29	0.47
3:C:302:SF4:S2	2:D:182:ARG:NH1	2.88	0.47
2:D:245:ALA:HB1	2:D:332:ALA:HB1	1.97	0.46
1:C:27:ASN:HA	2:D:170:ARG:HG3	1.98	0.46
1:A:73:PHE:CD1	5:A:306:KR:KR	3.30	0.46
2:D:246:LEU:HB2	2:D:326:LEU:HD13	1.96	0.46
2:B:167:GLU:OE1	2:B:170:ARG:NE	2.39	0.46
4:C:304[B]:6ML:S1	4:C:304[B]:6ML:S2	3.14	0.46
2:D:284:VAL:HG22	2:D:294:THR:HB	1.97	0.46
2:D:218:GLU:HG2	2:D:383:TYR:CE1	2.50	0.46
1:A:21[B]:CYS:N	4:A:304[B]:6ML:S1	2.89	0.46
1:C:19:THR:HA	1:C:77:GLU:OE2	2.16	0.45
1:C:14:GLN:OE1	1:C:47:PRO:HG2	2.16	0.45
2:D:331:LEU:N	2:D:377:ALA:HB1	2.31	0.45
2:D:116:GLN:HB2	2:D:172:CYS:CB	2.46	0.45
2:D:184:PRO:O	2:D:185:HIS:HB2	2.17	0.45
2:D:331:LEU:HA	2:D:377:ALA:HB1	1.98	0.45
2:B:219:GLN:O	2:B:223:PRO:HG2	2.16	0.45
2:B:443:SER:O	2:B:446:LEU:HB3	2.17	0.45
2:D:317:LYS:NZ	11:D:604:HOH:O	2.31	0.45
1:A:48:THR:O	2:B:26:ARG:HA	2.17	0.45
2:B:273[A]:MET:HB2	2:B:274:PRO:HD2	1.99	0.45
2:D:319:PRO:O	2:D:320:ARG:NH1	2.36	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:72:PHE:CE1	1:A:112:ALA:HB2	2.52	0.45
2:B:59:LEU:HD11	2:B:71:VAL:HG12	1.99	0.45
2:D:234:ASP:N	2:D:234:ASP:OD1	2.50	0.45
2:B:28:GLU:OE1	2:B:489[C]:SEC:SE	2.85	0.45
2:B:319:PRO:O	2:B:320:ARG:NH1	2.41	0.44
2:D:146:SER:O	2:D:147:ASP:HB3	2.17	0.44
2:D:244:ALA:O	2:D:327:GLU:HG2	2.18	0.44
1:A:166:ILE:O	1:A:169:THR:N	2.51	0.44
1:C:162:HIS:NE2	1:C:251:LYS:O	2.51	0.44
2:D:353:PHE:HB3	2:D:355:ILE:CD1	2.47	0.44
1:C:248:PHE:CD1	1:C:249:GLU:N	2.85	0.44
1:A:91:ILE:HD11	1:A:104:ALA:HB2	2.00	0.44
2:B:28:GLU:CD	2:B:28:GLU:C	2.85	0.44
2:B:121:HIS:O	2:B:125:LEU:HB2	2.17	0.44
1:C:38:LEU:CD2	2:D:136:THR:HG21	2.47	0.44
1:C:144:LEU:HD12	1:C:153:VAL:HG21	1.99	0.43
1:A:52:TRP:CZ3	2:B:141:PRO:HB2	2.54	0.43
1:A:228:ASP:OD1	1:A:228:ASP:C	2.62	0.43
1:C:246:ASP:OD1	1:C:250:ARG:NE	2.48	0.43
1:A:206:ASN:OD1	1:A:208:HIS:N	2.51	0.43
2:B:343:PRO:HD2	2:B:383:TYR:CZ	2.53	0.43
2:B:472:ASP:CG	2:B:475:ASN:HD22	2.25	0.43
1:A:48:THR:HG23	2:B:27:ILE:O	2.17	0.43
2:B:184:PRO:O	2:B:185:HIS:HB2	2.18	0.43
1:A:246:ASP:OD1	1:A:250:ARG:NE	2.45	0.43
4:A:304[B]:6ML:S1	4:A:304[B]:6ML:S2	3.16	0.43
1:C:88:TYR:CG	1:C:135:THR:HG22	2.54	0.43
2:B:222:VAL:HB	2:B:223:PRO:HD3	2.01	0.43
1:C:116:LEU:HD23	1:C:154:VAL:HB	2.00	0.43
2:D:222:VAL:HB	2:D:223:PRO:HD3	2.01	0.43
2:D:331:LEU:CA	2:D:377:ALA:HB1	2.48	0.43
1:C:200:THR:N	1:C:201:PRO:CD	2.82	0.43
2:D:487:ASP:OD2	11:D:603:HOH:O	2.21	0.43
2:B:146:SER:O	2:B:147:ASP:HB3	2.18	0.43
2:D:100:ASN:OD1	2:D:255:ASN:ND2	2.49	0.43
2:D:132:GLN:OE1	2:D:142:ARG:HD3	2.19	0.43
2:B:44:ASP:OD1	2:B:452:ARG:NH2	2.52	0.42
2:B:478:ASN:OD1	2:B:481:ARG:NH1	2.49	0.42
1:A:24:THR:HG21	1:A:161:PRO:O	2.19	0.42
2:B:244:ALA:O	2:B:327:GLU:HG2	2.19	0.42
2:B:198:PRO:HG2	2:B:396:ILE:HD11	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:84:ALA:O	1:C:87:LYS:HG3	2.19	0.42
1:C:227:THR:CG2	2:D:197:ILE:HG13	2.50	0.42
2:D:489[A]:SEC:SE	6:D:501:FCO:N1	3.01	0.42
2:D:404:VAL:HG23	2:D:406:PRO:HD3	2.00	0.42
2:B:132:GLN:OE1	2:B:142:ARG:HD3	2.19	0.42
1:C:127:ILE:HD12	2:D:73:ARG:HG2	2.00	0.42
1:C:154:VAL:HG23	5:C:310:KR:KR	2.80	0.42
1:A:116:LEU:HD23	1:A:154:VAL:HB	2.02	0.42
2:B:173:HIS:CD2	2:B:183:MET:HG2	2.55	0.42
2:B:458:ARG:NH2	2:B:466:ILE:HG13	2.34	0.42
1:C:154:VAL:CG2	5:C:310:KR:KR	3.29	0.42
1:C:257:ASN:OD1	1:C:258:TRP:N	2.50	0.42
2:D:283:TYR:CZ	2:D:455:MET:HE1	2.55	0.42
2:D:28:GLU:CD	2:D:28:GLU:C	2.87	0.42
2:D:316:VAL:CG1	2:D:446:LEU:HD12	2.50	0.42
1:C:224:ALA:O	2:D:405:MET:HE2	2.19	0.42
2:D:129:ASP:OD1	2:D:484:ARG:NH1	2.52	0.42
2:D:326:LEU:O	2:D:460:ALA:HB2	2.20	0.42
2:D:480:ALA:O	2:D:484:ARG:HG3	2.20	0.42
1:A:166:ILE:O	1:A:167:VAL:C	2.62	0.42
2:B:28:GLU:HB3	2:B:489[B]:SEC:SE	2.70	0.42
2:B:246:LEU:HB3	2:B:328:VAL:HG12	2.02	0.42
1:C:52:TRP:CZ3	2:D:141:PRO:HB2	2.55	0.42
2:D:420:ALA:HB1	6:D:501:FCO:N2	2.35	0.42
2:B:126:SER:O	2:B:127:ALA:C	2.63	0.41
2:B:326:LEU:O	2:B:460:ALA:HB2	2.20	0.41
1:C:95:ASN:OD1	1:C:95:ASN:N	2.54	0.41
2:B:124:HIS:HB3	5:B:508[B]:KR:KR	2.81	0.41
1:C:160:PRO:HG3	2:D:184:PRO:HD2	2.02	0.41
1:C:231:GLY:O	1:C:233:ARG:NH1	2.50	0.41
2:D:472:ASP:OD2	2:D:475:ASN:ND2	2.45	0.41
1:C:186:LEU:O	1:C:190:VAL:HG23	2.20	0.41
1:A:116:LEU:CD2	1:A:154:VAL:HB	2.51	0.41
2:D:243:LYS:HA	2:D:369:PHE:CD2	2.56	0.41
1:A:240:GLY:N	1:A:241:PRO:CD	2.83	0.41
2:B:244:ALA:HB1	2:B:326:LEU:CD2	2.50	0.41
2:B:422:ARG:C	2:B:444:ALA:HB1	2.46	0.41
2:D:328:VAL:HB	2:D:423:GLY:HA3	2.02	0.41
1:C:232:CYS:SG	1:C:233:ARG:N	2.93	0.41
2:D:143:PHE:O	2:D:146:SER:OG	2.39	0.41
2:B:281:LYS:HD3	2:B:295:GLY:O	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:426:LEU:HD12	2:B:426:LEU:C	2.46	0.41
2:D:72:GLN:HB3	2:D:79:PRO:HB2	2.02	0.41
2:D:87:VAL:HA	2:D:90:LEU:HD12	2.03	0.41
1:A:26:LEU:HD23	1:A:33:ILE:HG12	2.03	0.41
1:A:279:PRO:HB2	1:A:282:GLN:HG3	2.02	0.41
1:A:60:HIS:O	1:A:64:VAL:HG23	2.20	0.40
2:B:28:GLU:HB3	2:B:489[A]:SEC:CB	2.52	0.40
2:D:284:VAL:O	2:D:284:VAL:HG23	2.21	0.40
2:D:409:PRO:O	2:D:434:SER:HA	2.21	0.40
1:A:127:ILE:HD12	2:B:73:ARG:HG2	2.02	0.40
1:A:222:MET:HG3	1:A:235:ASP:OD2	2.21	0.40
1:A:278:SER:HA	1:A:279:PRO:C	2.45	0.40
2:B:147:ASP:O	2:B:481:ARG:NH2	2.54	0.40
2:B:372:MET:O	2:B:376:VAL:HG23	2.20	0.40
1:C:205:ARG:HH22	1:C:249:GLU:CD	2.30	0.40
2:D:143:PHE:CZ	2:D:476:PRO:HD2	2.56	0.40
2:D:28:GLU:O	2:D:489[A]:SEC:HA	2.22	0.40
2:D:426:LEU:HG	2:D:442:VAL:HB	2.04	0.40
1:A:60:HIS:ND1	2:B:141:PRO:HG3	2.37	0.40

All (3) 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:178:LYS:O	2:D:16:ARG:NH1[2_545]	1.72	0.48
1:A:178:LYS:O	2:D:16:ARG:HH12[2_545]	1.37	0.23
2:B:153:GLU:OE1	1:C:191:LYS:NZ[1_655]	2.09	0.11

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	278/283 (98%)	257 (92%)	19 (7%)	2 (1%)	18	19
1	C	279/283 (99%)	256 (92%)	22 (8%)	1 (0%)	30	34
2	B	483/484 (100%)	450 (93%)	32 (7%)	1 (0%)	43	51
2	D	483/484 (100%)	452 (94%)	28 (6%)	3 (1%)	21	23
All	All	1523/1534 (99%)	1415 (93%)	101 (7%)	7 (0%)	24	27

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	182	LEU
1	C	263	ALA
2	D	185	HIS
1	A	263	ALA
2	B	133	GLY
2	D	133	GLY
2	D	147	ASP

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	225/226 (100%)	218 (97%)	7 (3%)	35	48
1	C	226/226 (100%)	221 (98%)	5 (2%)	45	61
2	B	396/392 (101%)	388 (98%)	8 (2%)	48	64
2	D	396/392 (101%)	392 (99%)	4 (1%)	68	81
All	All	1243/1236 (101%)	1219 (98%)	24 (2%)	50	66

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	52	TRP
1	A	58	ILE
1	A	95	ASN
1	A	170	VAL

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Mol	Chain	Res	Type
1	A	179	LYS
1	A	189	VAL
1	A	282	GLN
2	B	17	THR
2	B	48[A]	SER
2	B	48[B]	SER
2	B	84	THR
2	B	211	LYS
2	B	356	SER
2	B	359	LYS
2	B	455	MET
1	C	95	ASN
1	C	141	SER
1	C	179	LYS
1	C	189	VAL
1	C	282	GLN
2	D	17	THR
2	D	84	THR
2	D	287	SER
2	D	455	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	30	HIS
1	A	97	HIS
1	A	253	ASN
2	B	30	HIS
2	B	128	GLN
2	B	161	GLN
2	B	189	GLN
1	C	97	HIS
1	C	111	ASN
1	C	147	ASN
2	D	161	GLN
2	D	189	GLN

5.3.3 RNA ⓘ

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 38 ligands modelled in this entry, 26 are monoatomic and 2 are modelled with single atom - leaving 10 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	$\# Z > 2$
6	FCO	B	501	2	0,6,6	-	-	-		
3	SF4	C	303[A]	1	0,12,12	-	-	-		
3	SF4	A	301	1	0,12,12	-	-	-		
3	SF4	C	301	1	0,12,12	-	-	-		
3	SF4	C	302	1	0,12,12	-	-	-		
3	SF4	A	303[A]	1	0,12,12	-	-	-		
4	6ML	A	304[B]	1	0,12,12	-	-	-		
4	6ML	C	304[B]	1	0,12,12	-	-	-		
6	FCO	D	501	2	0,6,6	-	-	-		
3	SF4	A	302	1	0,12,12	-	-	-		

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	SF4	C	303[A]	1	-	-	0/6/5/5
3	SF4	A	301	1	-	-	0/6/5/5
3	SF4	C	301	1	-	-	0/6/5/5
3	SF4	C	302	1	-	-	0/6/5/5

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	SF4	A	303[A]	1	-	-	0/6/5/5
4	6ML	A	304[B]	1	-	-	0/2/3/3
4	6ML	C	304[B]	1	-	-	0/2/3/3
3	SF4	A	302	1	-	-	0/6/5/5

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.

6 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	B	501	FCO	1	0
3	C	302	SF4	1	0
4	A	304[B]	6ML	2	0
4	C	304[B]	6ML	2	0
6	D	501	FCO	3	0
3	A	302	SF4	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	277/283 (97%)	0.59	18 (6%)	25 22	25, 43, 63, 84	2 (0%)
1	C	278/283 (98%)	0.50	9 (3%)	50 47	16, 44, 63, 77	3 (1%)
2	B	480/484 (99%)	0.67	36 (7%)	20 17	24, 50, 74, 89	6 (1%)
2	D	480/484 (99%)	0.73	27 (5%)	30 27	23, 53, 72, 91	5 (1%)
All	All	1515/1534 (98%)	0.64	90 (5%)	28 25	16, 48, 72, 91	16 (1%)

All (90) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	363	LEU	6.8
1	A	21[A]	CYS	5.9
2	D	368	ALA	4.9
2	B	140	VAL	4.4
2	D	140	VAL	4.4
1	A	172	LEU	4.0
2	D	141	PRO	3.9
1	A	207	ILE	3.7
1	C	283	ALA	3.6
2	B	483	ILE	3.5
2	B	217	VAL	3.5
2	B	360	PHE	3.5
1	A	17	GLY	3.3
1	C	221	VAL	3.1
2	B	141	PRO	3.1
2	D	364	GLY	3.0
2	B	465	LEU	3.0
2	D	403	VAL	3.0
1	A	186	LEU	2.9
2	B	479	VAL	2.9
2	D	94	PHE	2.8

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Mol	Chain	Res	Type	RSRZ
2	B	248	VAL	2.8
2	B	133	GLY	2.8
1	A	105	LEU	2.8
2	D	447	TRP	2.7
2	D	280	ILE	2.7
2	B	139	PHE	2.7
1	A	181	GLY	2.6
2	D	379	ALA	2.6
2	B	235	MET	2.6
2	D	376	VAL	2.6
2	B	323	GLY	2.6
2	B	473	ILE	2.6
1	A	52	TRP	2.5
2	B	353	PHE	2.5
2	D	392	TRP	2.5
2	B	226	TYR	2.5
2	D	369	PHE	2.5
1	A	268	CYS	2.4
2	B	486	PHE	2.4
1	A	283	ALA	2.4
1	C	230	VAL	2.4
2	D	143	PHE	2.4
2	B	157	ALA	2.4
2	B	143	PHE	2.4
2	B	461	VAL	2.4
1	A	94	ALA	2.4
2	D	36	VAL	2.4
2	B	221	TYR	2.3
2	B	464	ALA	2.3
2	B	480	ALA	2.3
2	D	473	ILE	2.3
2	B	134	PRO	2.3
1	A	20	GLY	2.3
2	B	150	LEU	2.3
1	A	169	THR	2.3
1	A	189	VAL	2.3
2	D	134	PRO	2.3
2	B	132	GLN	2.3
2	B	349	LEU	2.2
2	D	360	PHE	2.2
1	A	18	CYS	2.2
2	D	228	ILE	2.2

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Mol	Chain	Res	Type	RSRZ
2	D	375	HIS	2.2
1	C	172	LEU	2.2
2	B	154	LEU	2.2
2	B	467	GLY	2.2
1	A	209	GLU	2.2
1	A	176	ALA	2.2
2	D	262	PHE	2.1
1	C	213	TYR	2.1
2	B	326	LEU	2.1
2	D	204	VAL	2.1
2	D	45	ALA	2.1
2	B	363	LEU	2.1
2	B	188	GLY	2.1
1	C	18	CYS	2.1
2	D	15	GLY	2.1
1	C	52	TRP	2.1
2	B	447	TRP	2.1
1	C	259	CYS	2.0
2	D	245	ALA	2.0
1	C	234	TYR	2.0
2	D	191	VAL	2.0
2	D	417	PHE	2.0
2	B	455	MET	2.0
1	A	121	CYS	2.0
2	B	75	CYS	2.0
2	B	324	LEU	2.0
2	B	469	PRO	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

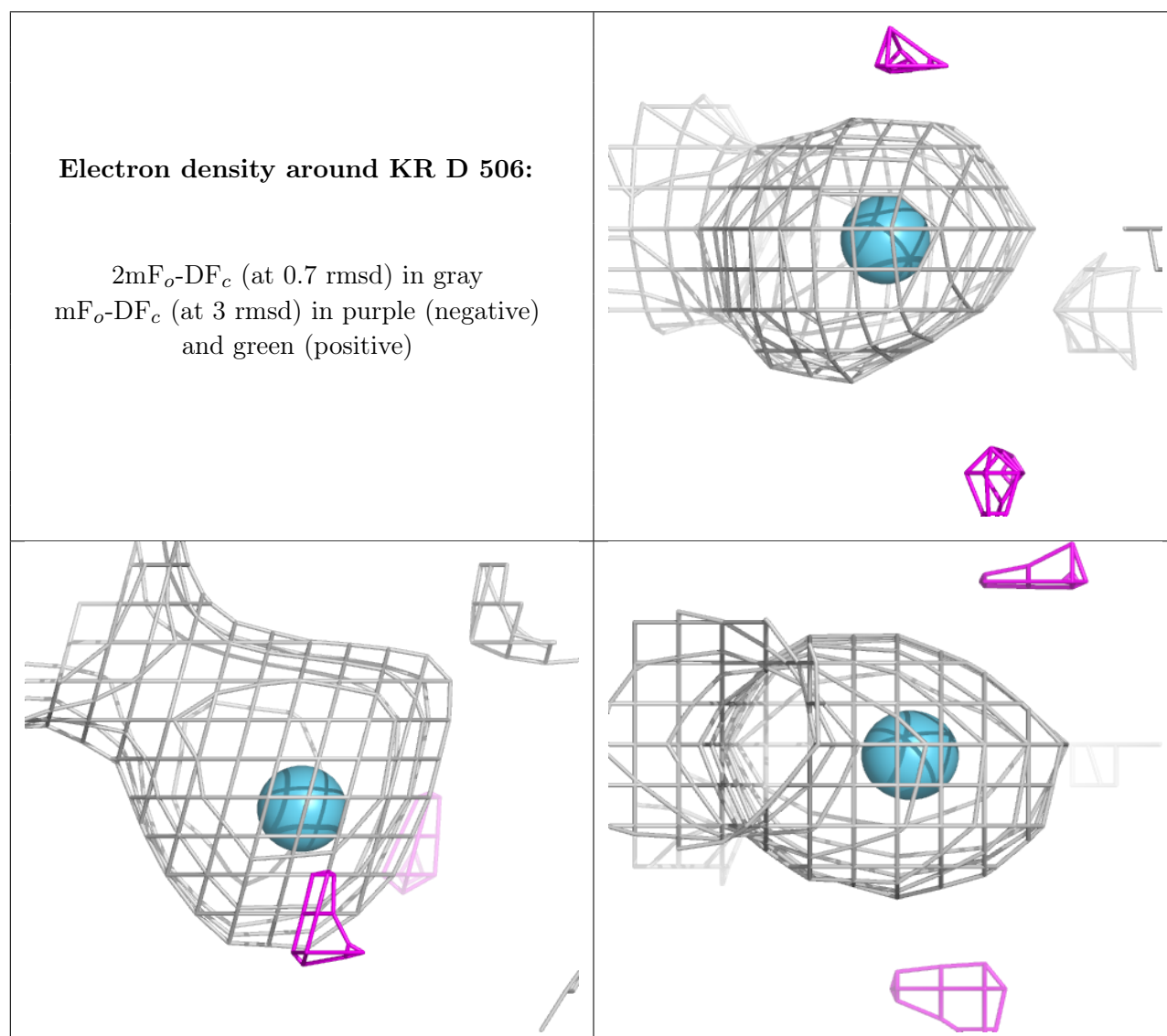
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum,

median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
5	KR	D	506	1/1	0.85	0.11	58,58,58,58	1
5	KR	C	307	1/1	0.89	0.10	42,42,42,42	1
5	KR	B	510	1/1	0.89	0.08	93,93,93,93	0
5	KR	B	506	1/1	0.91	0.07	43,43,43,43	1
5	KR	D	507	1/1	0.91	0.07	43,43,43,43	1
7	NI	D	502	1/1	0.92	0.07	22,22,22,22	1
10	CL	D	505	1/1	0.92	0.14	37,37,37,37	0
5	KR	B	508[B]	1/1	0.93	0.19	36,36,36,36	1
5	KR	B	509	1/1	0.93	0.08	45,45,45,45	1
5	KR	D	508	1/1	0.93	0.10	53,53,53,53	1
5	KR	A	305	1/1	0.93	0.09	57,57,57,57	1
8	FE2	B	503	1/1	0.93	0.04	22,22,22,22	0
5	KR	B	507[A]	1/1	0.93	0.08	40,40,40,40	1
5	KR	A	306	1/1	0.94	0.08	58,58,58,58	1
5	KR	C	306	1/1	0.95	0.05	48,48,48,48	1
9	H2S	D	504	1/1	0.95	0.21	38,38,38,38	1
5	KR	A	310	1/1	0.95	0.07	36,36,36,36	1
4	6ML	A	304[B]	10/10	0.96	0.09	25,29,37,39	10
5	KR	C	308	1/1	0.96	0.06	51,51,51,51	1
3	SF4	C	303[A]	8/8	0.96	0.08	21,24,27,29	8
5	KR	C	309	1/1	0.97	0.06	52,52,52,52	1
3	SF4	A	301	8/8	0.97	0.06	25,41,51,55	0
5	KR	A	307	1/1	0.97	0.07	45,45,45,45	1
5	KR	A	309	1/1	0.97	0.09	46,46,46,46	1
5	KR	C	305	1/1	0.97	0.04	59,59,59,59	1
3	SF4	A	303[A]	8/8	0.97	0.08	25,29,34,38	8
8	FE2	D	503	1/1	0.97	0.02	23,23,23,23	0
4	6ML	C	304[B]	10/10	0.97	0.09	21,23,28,29	10
10	CL	B	505	1/1	0.97	0.04	29,29,29,29	0
3	SF4	C	301	8/8	0.97	0.05	40,47,56,63	0
5	KR	C	310	1/1	0.98	0.04	39,39,39,39	1
3	SF4	A	302	8/8	0.98	0.09	24,30,34,41	0
3	SF4	C	302	8/8	0.98	0.08	25,31,37,50	0
9	H2S	B	504	1/1	0.98	0.12	31,31,31,31	0
5	KR	A	308	1/1	0.98	0.05	52,52,52,52	1
6	FCO	B	501	7/7	0.98	0.13	21,30,34,41	0
6	FCO	D	501	7/7	0.98	0.13	20,33,37,39	0
7	NI	B	502	1/1	0.99	0.07	23,23,23,23	1

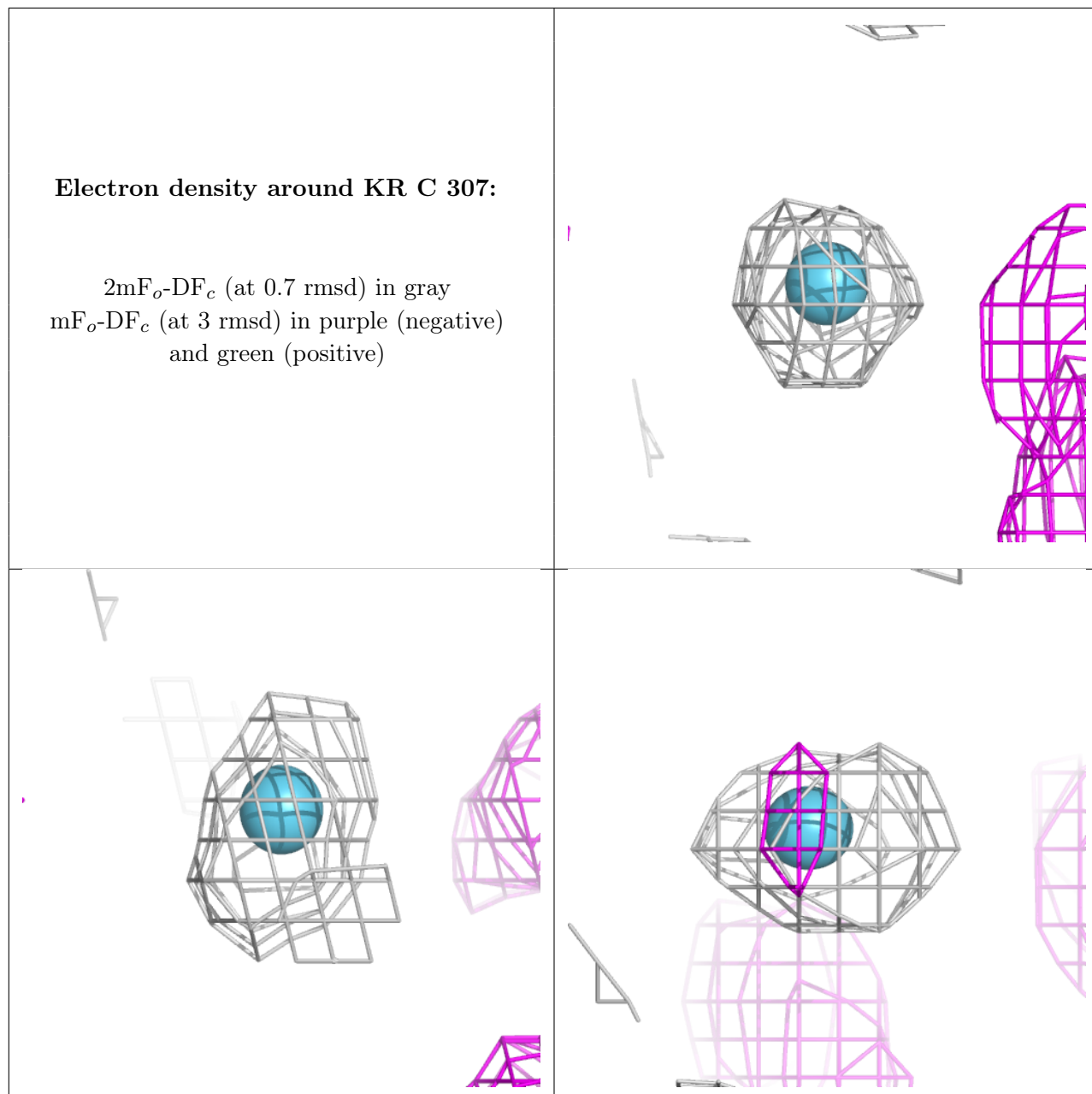
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers

as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.



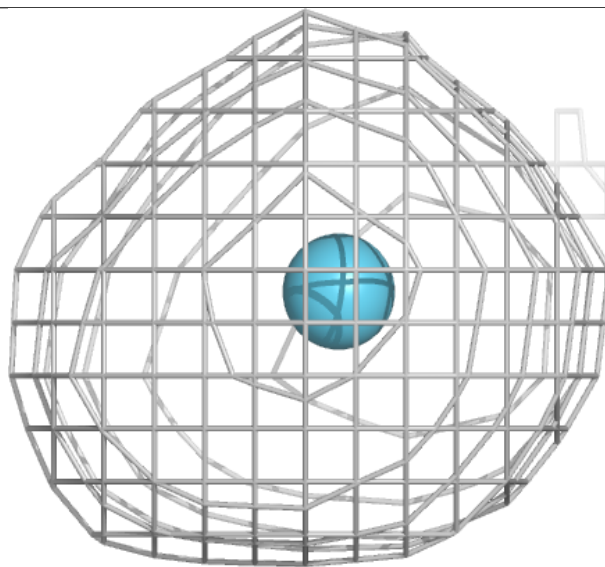
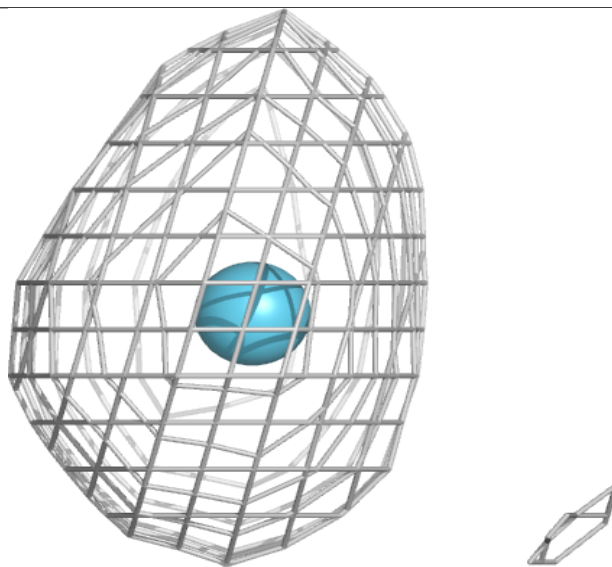
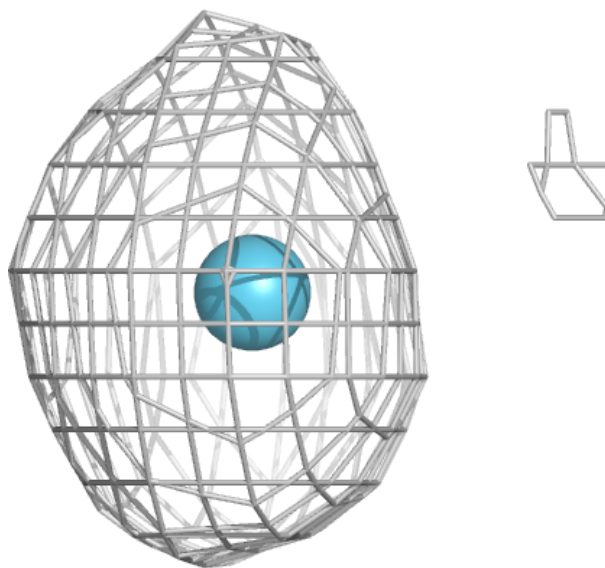
Electron density around KR C 307:

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



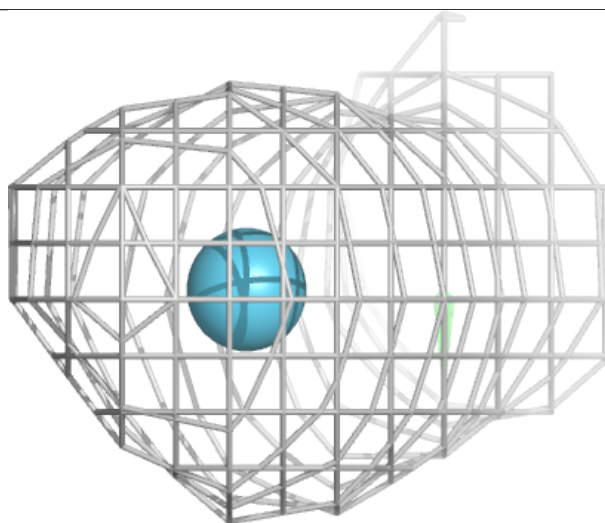
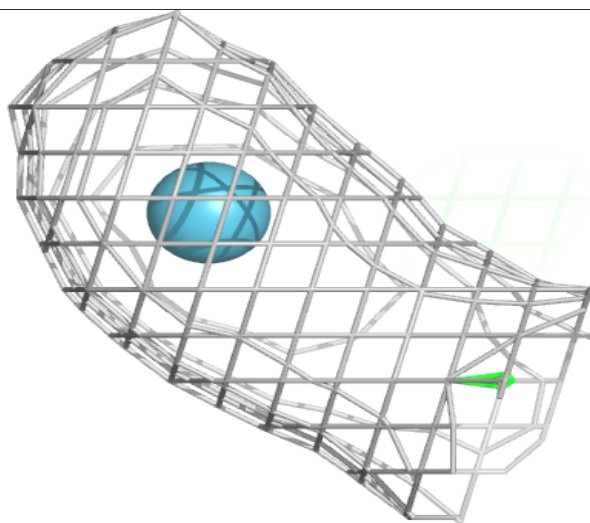
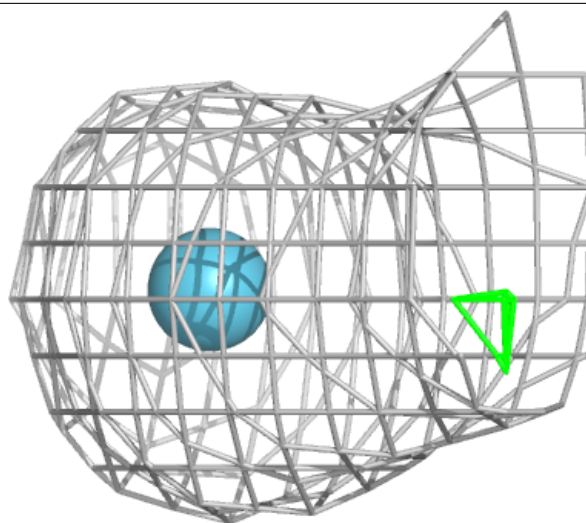
Electron density around KR B 510:

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



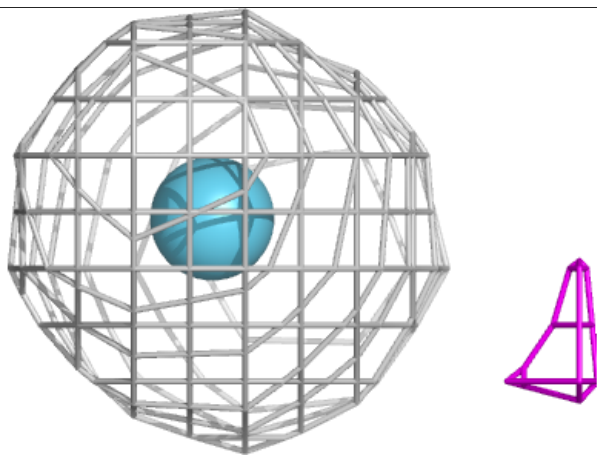
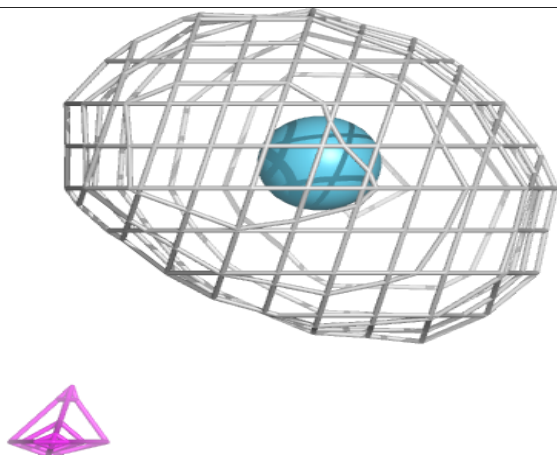
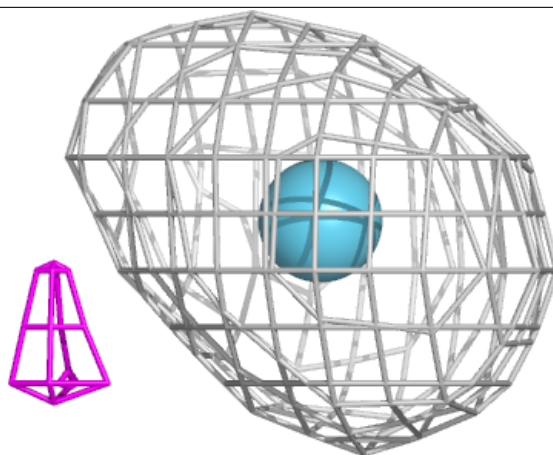
Electron density around KR B 506:

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



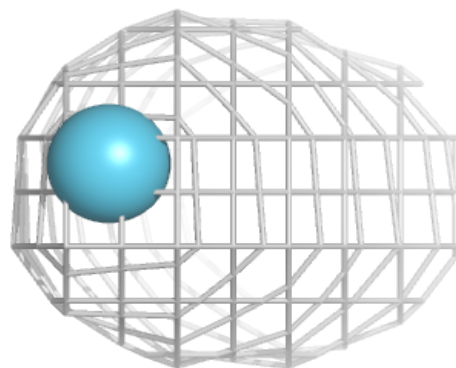
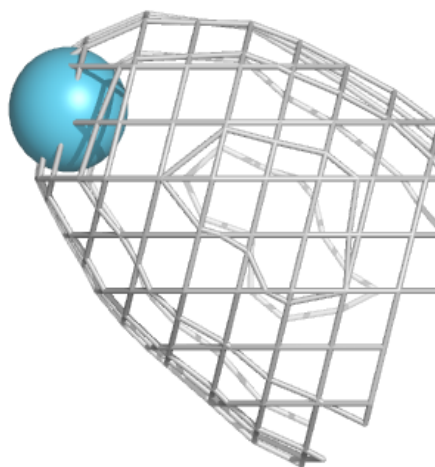
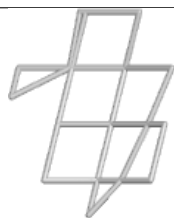
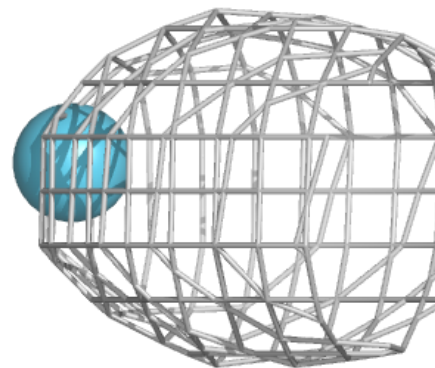
Electron density around KR D 507:

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



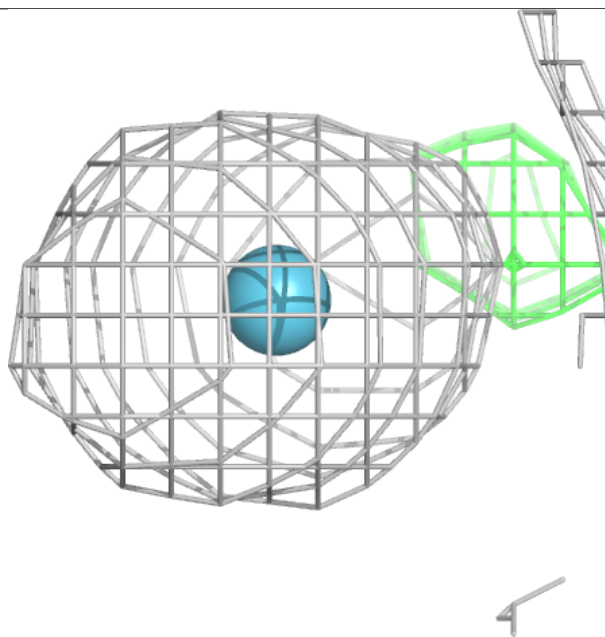
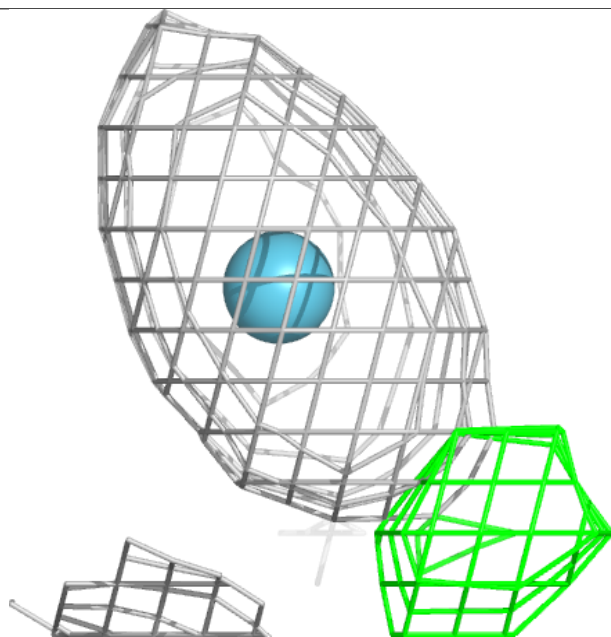
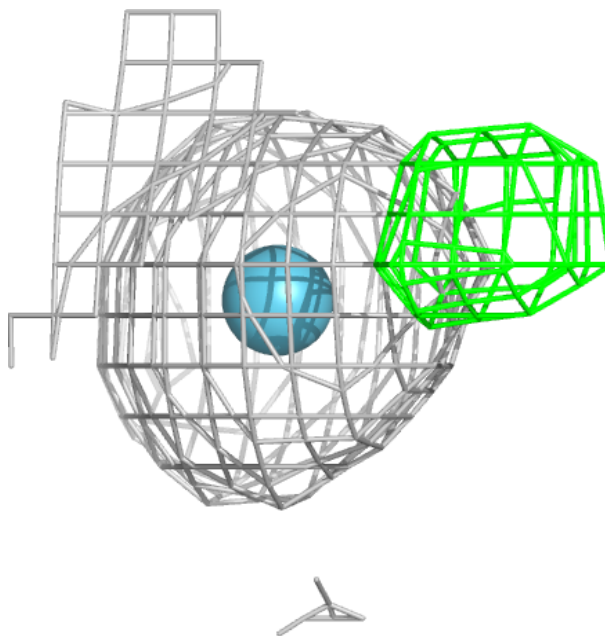
Electron density around KR B 508 (B):

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



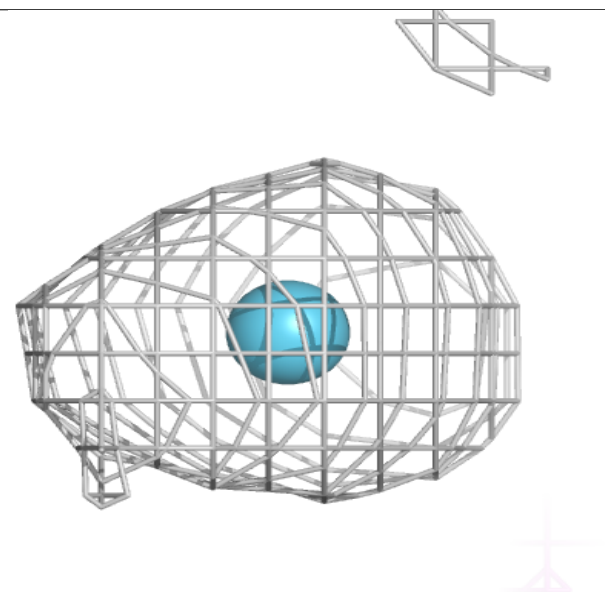
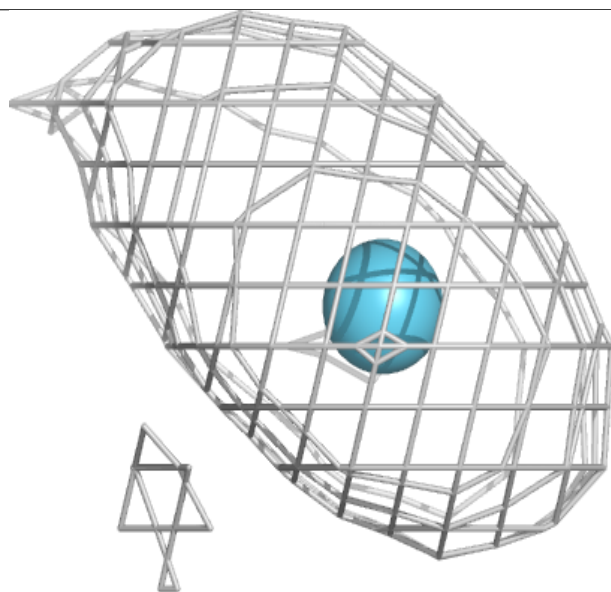
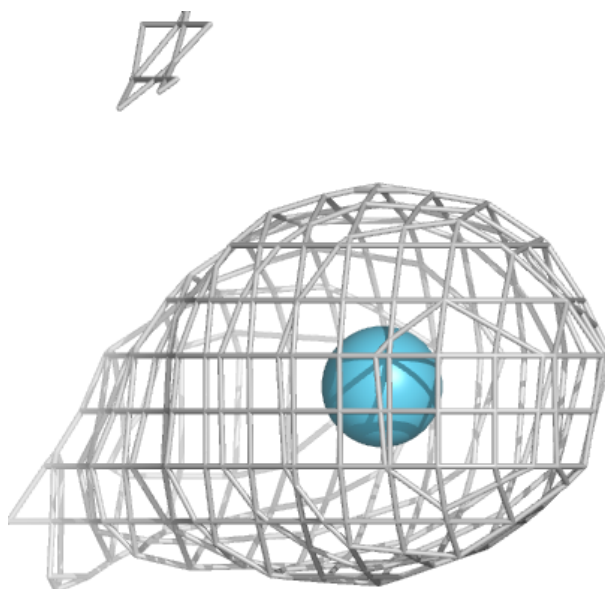
Electron density around KR B 509:

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



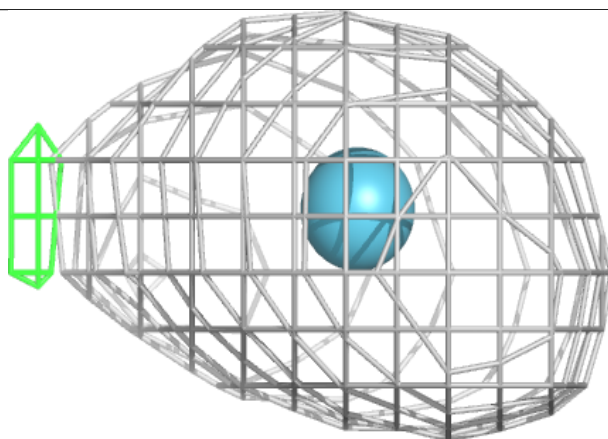
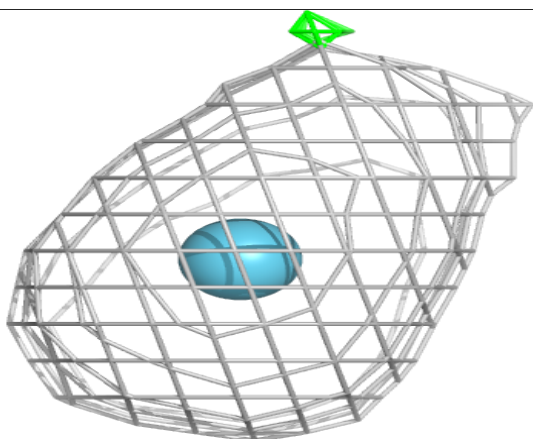
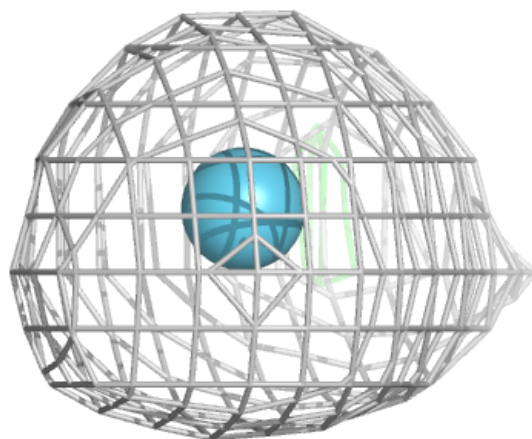
Electron density around KR D 508:

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



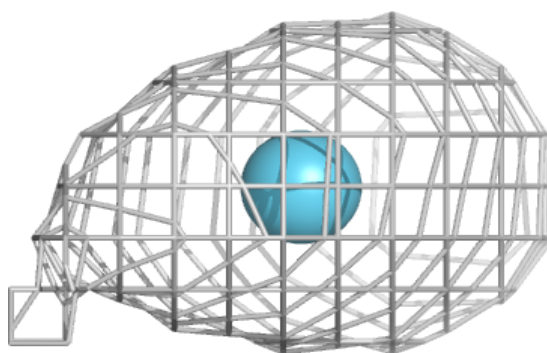
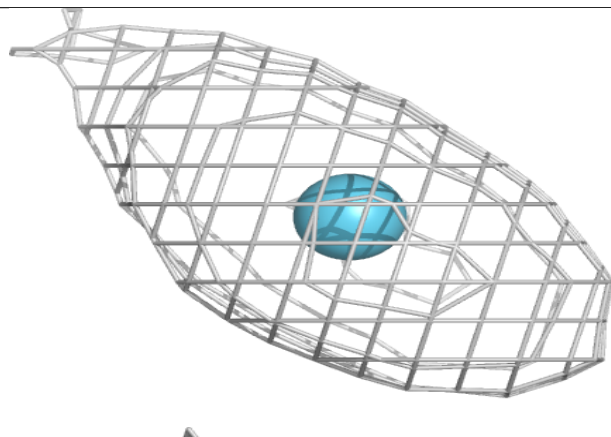
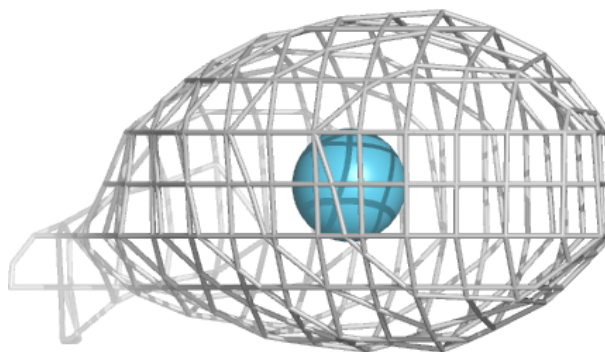
Electron density around KR A 305:

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



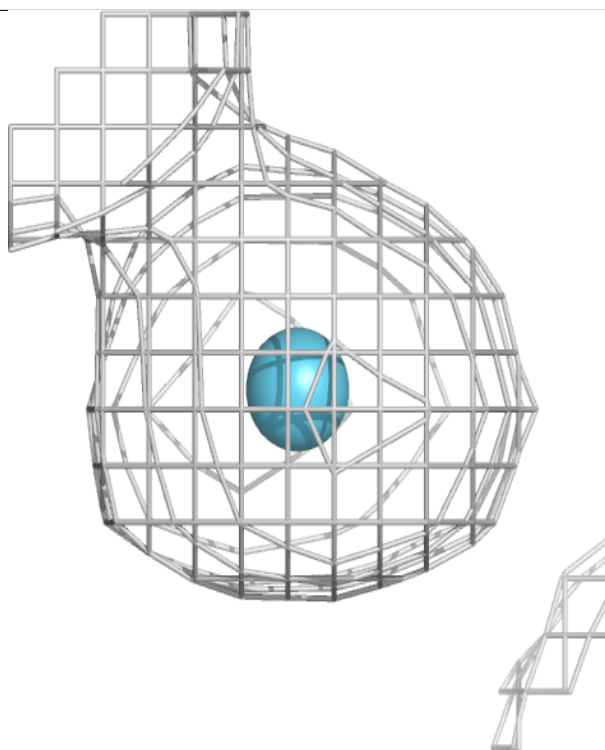
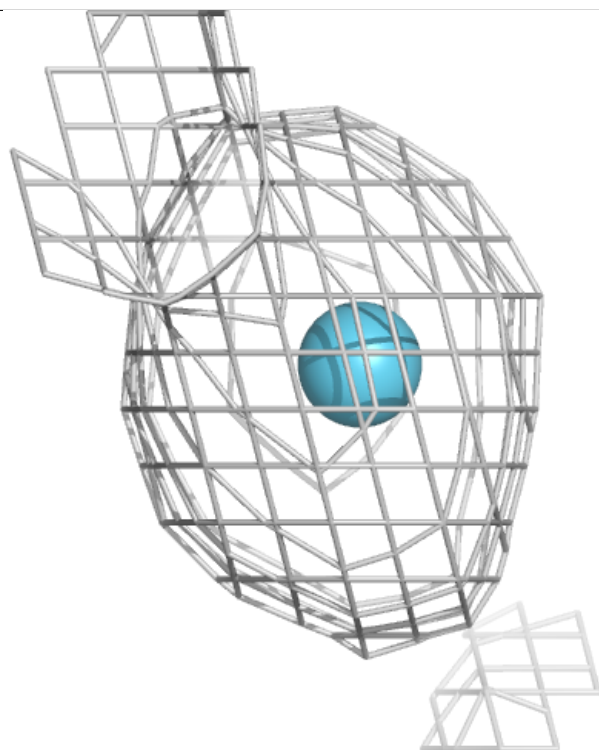
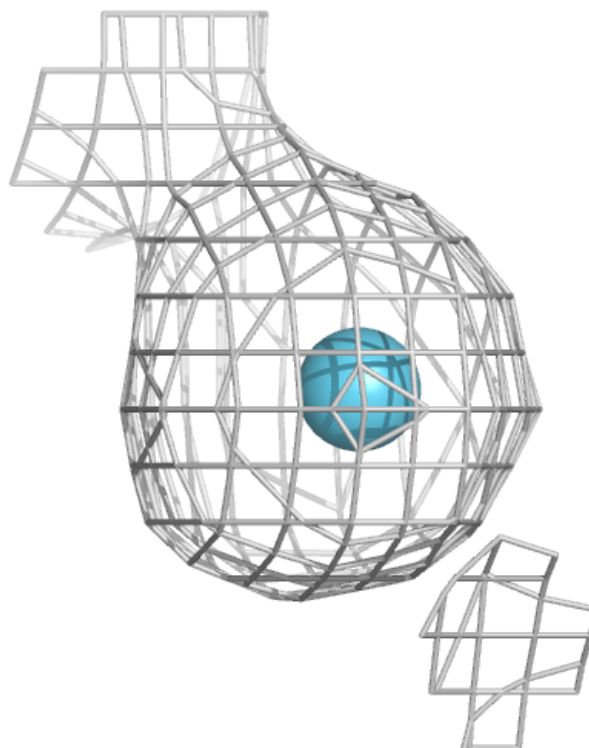
Electron density around KR B 507 (A):

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



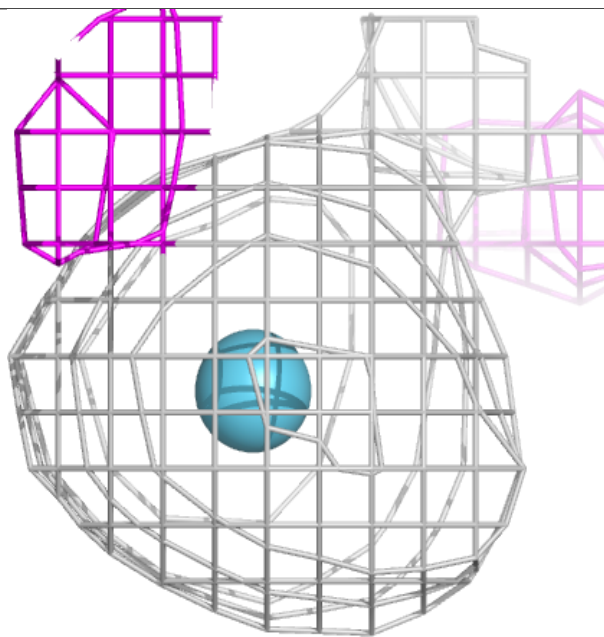
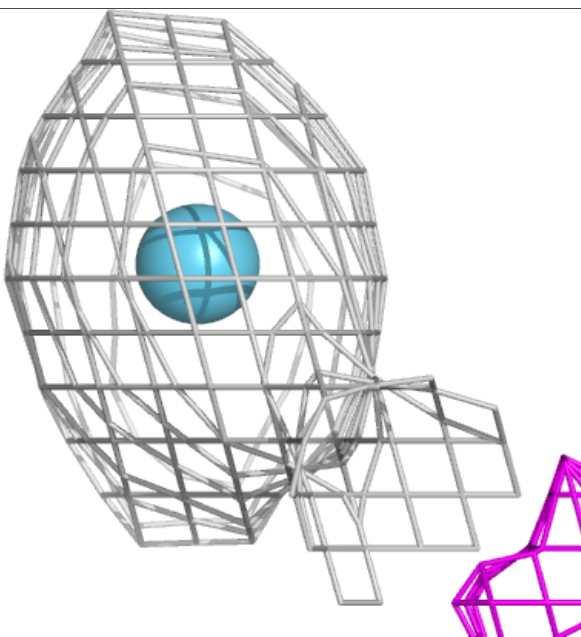
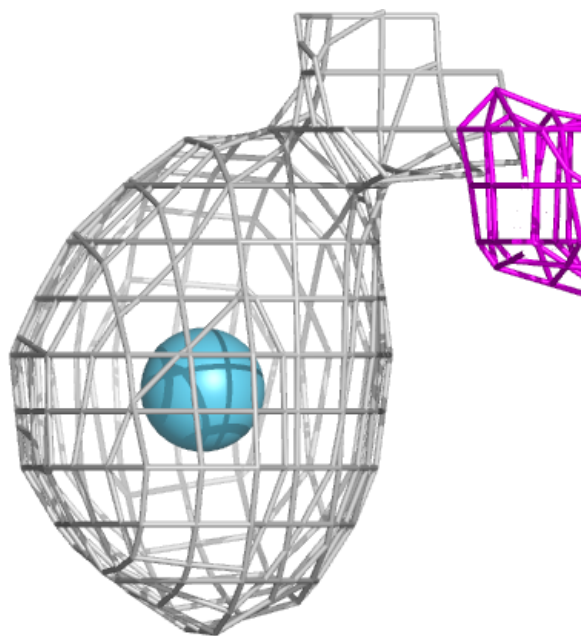
Electron density around KR A 306:

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



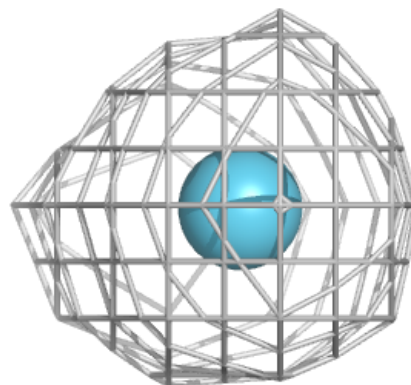
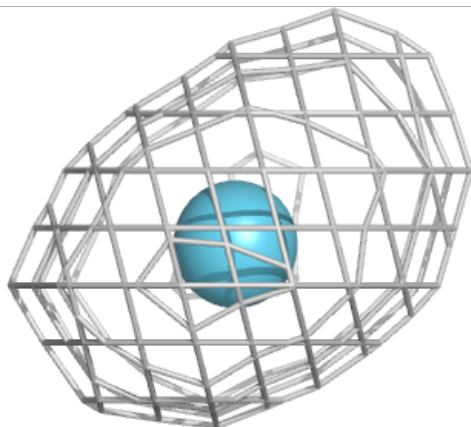
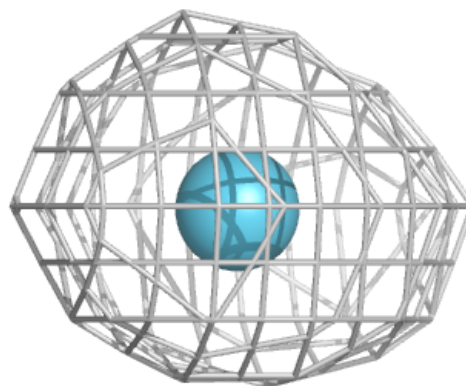
Electron density around KR C 306:

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



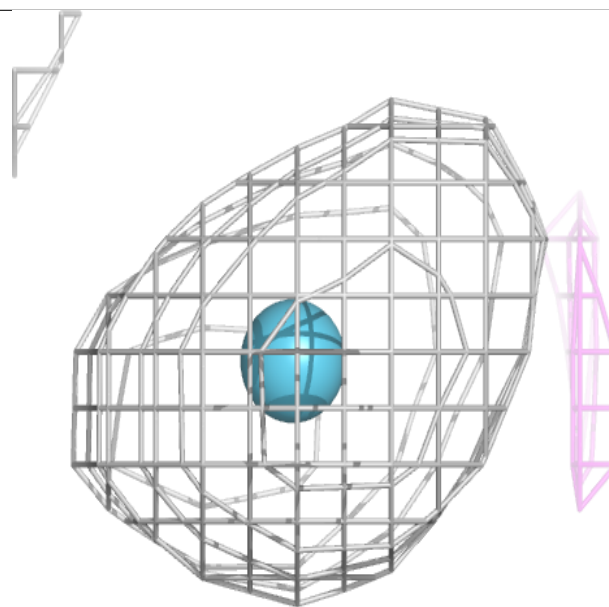
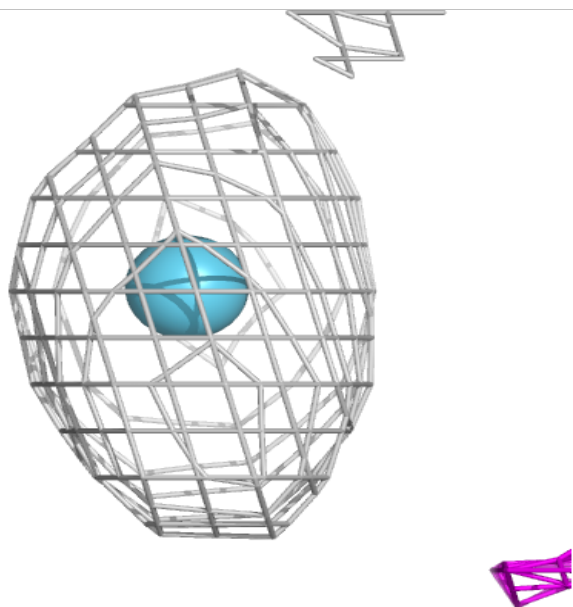
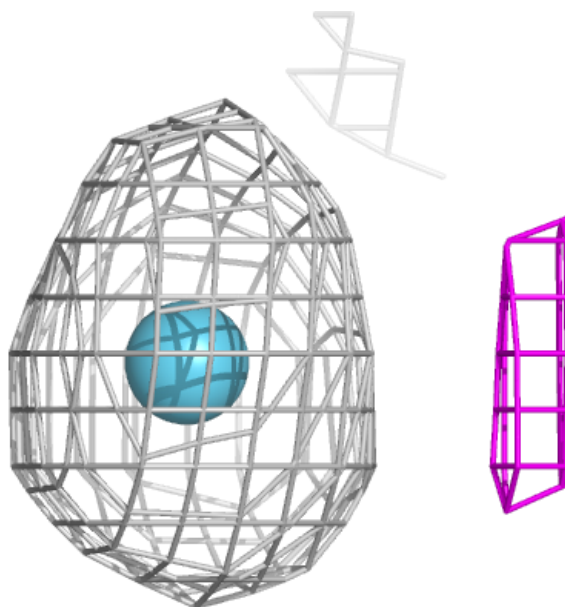
Electron density around KR A 310:

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



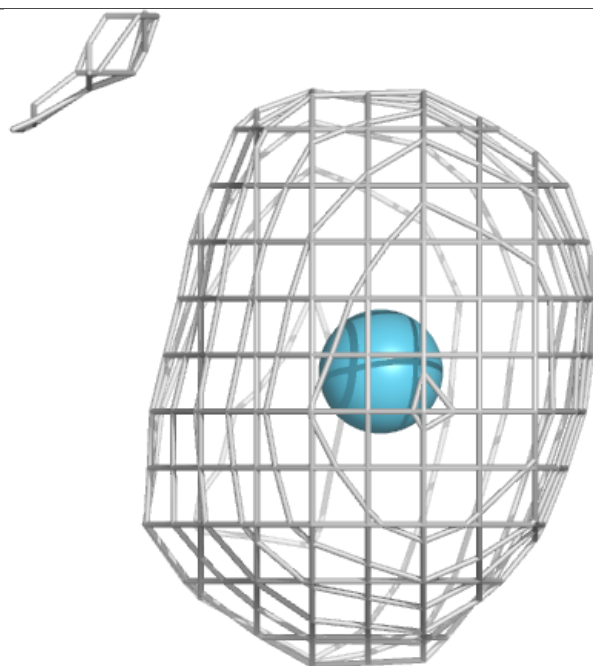
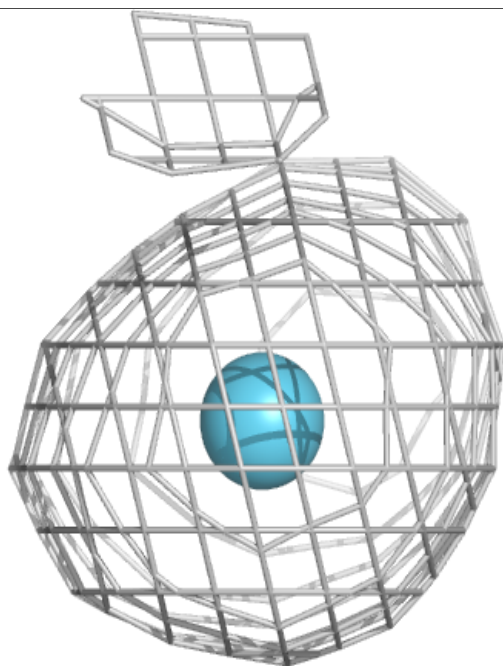
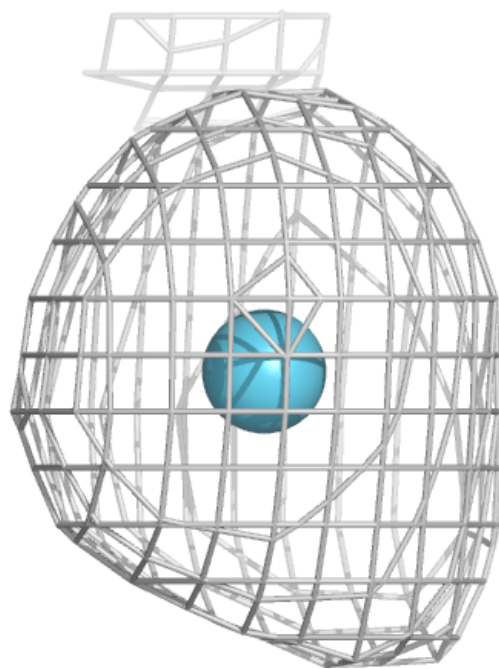
Electron density around KR C 308:

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



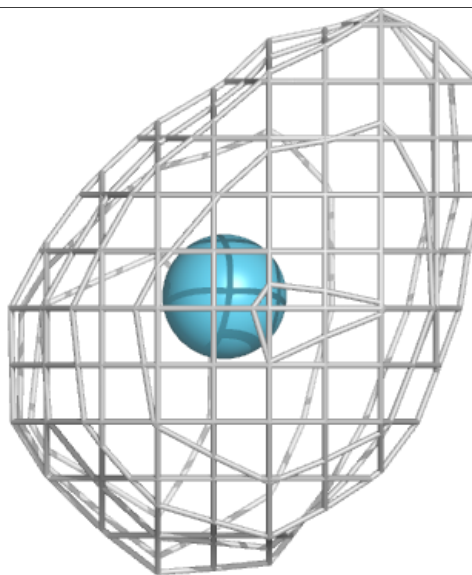
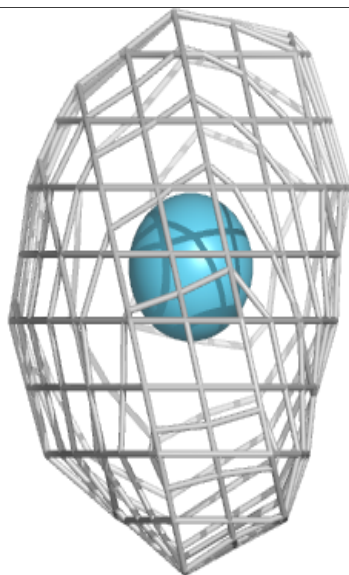
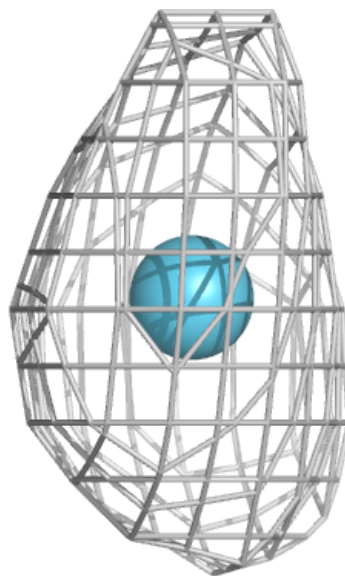
Electron density around KR C 309:

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



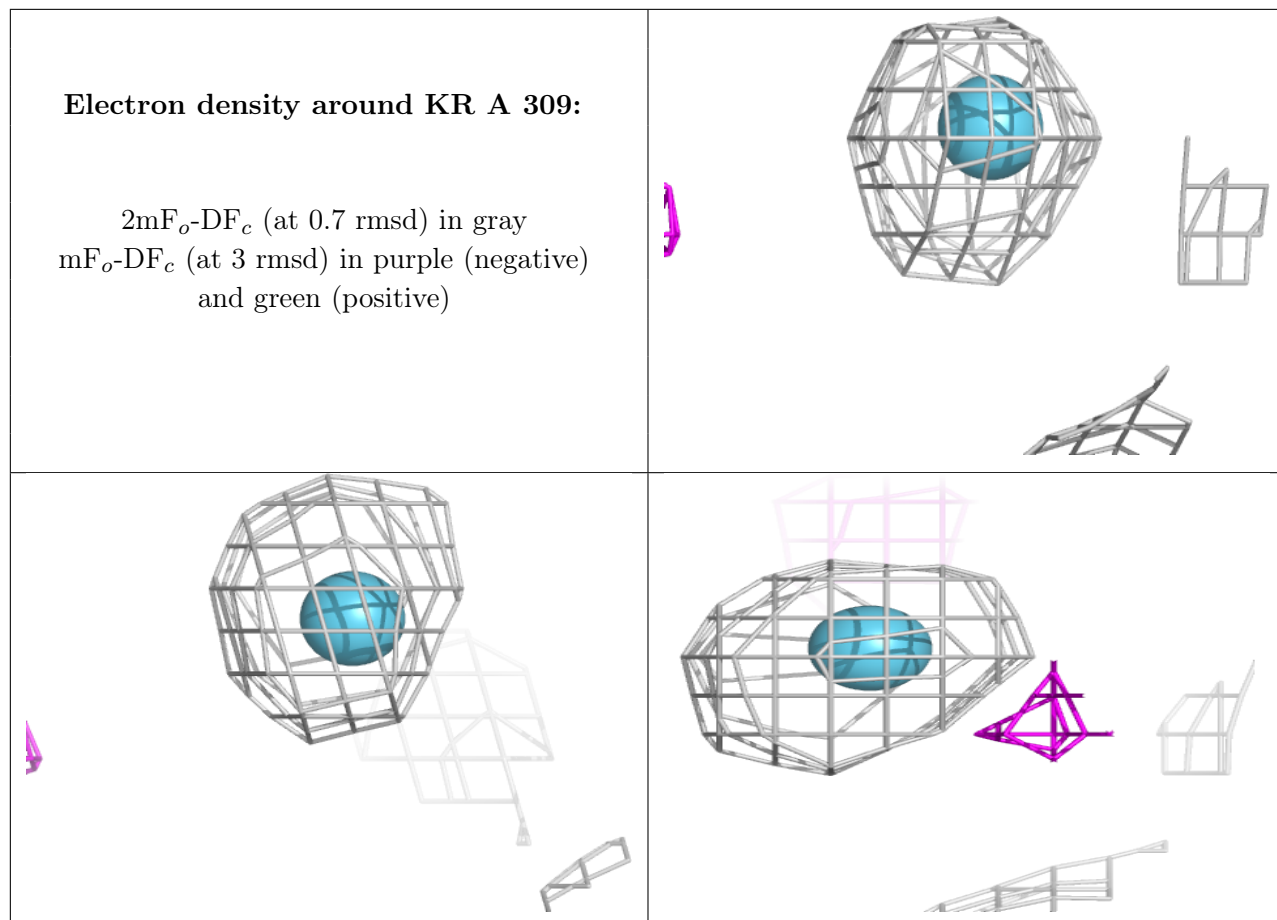
Electron density around KR A 307:

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



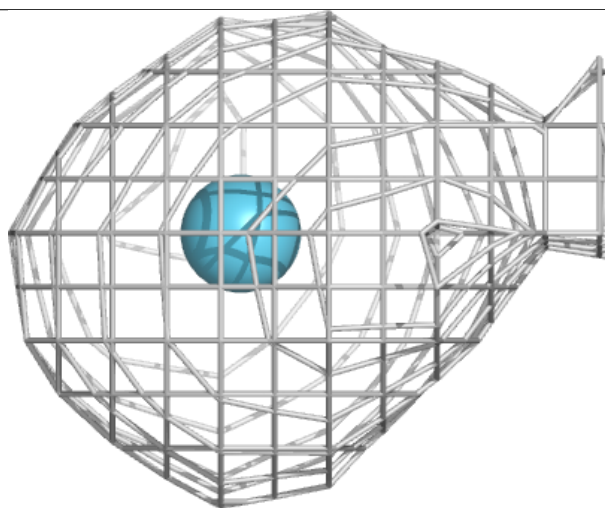
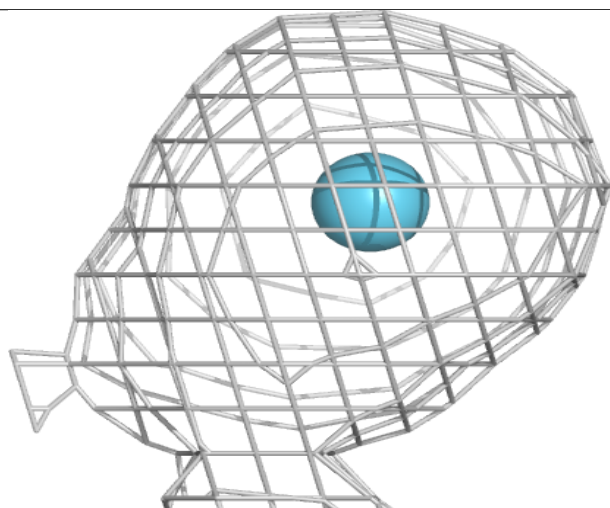
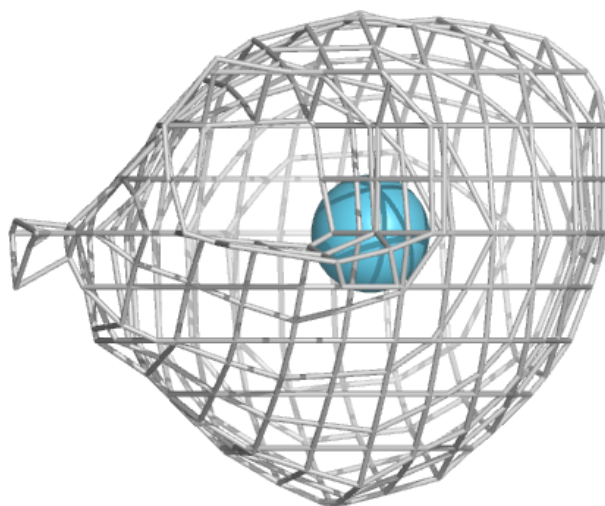
Electron density around KR A 309:

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



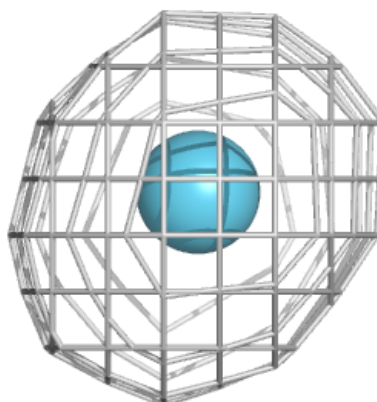
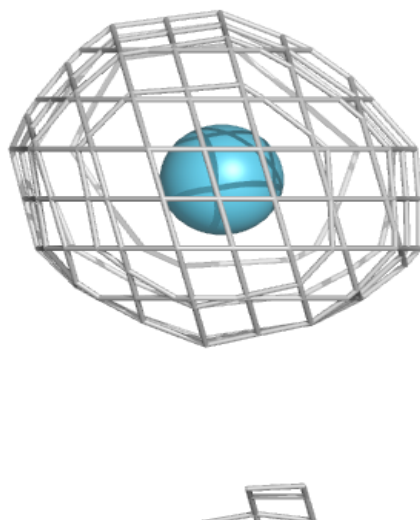
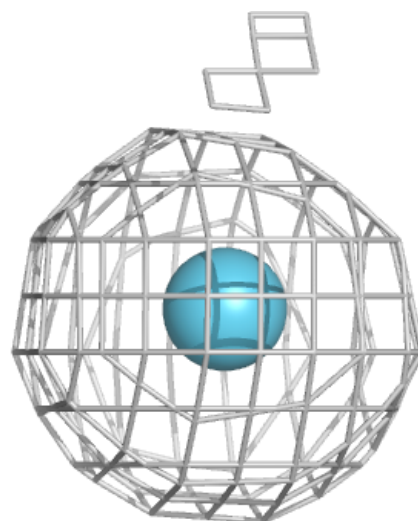
Electron density around KR C 305:

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



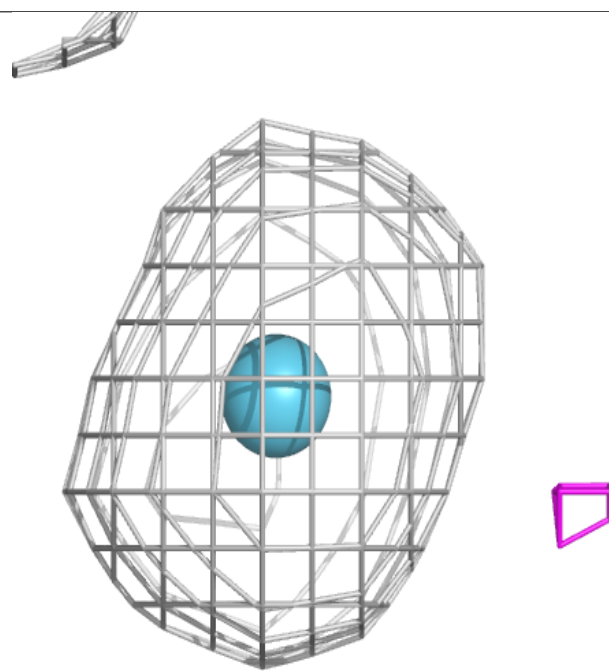
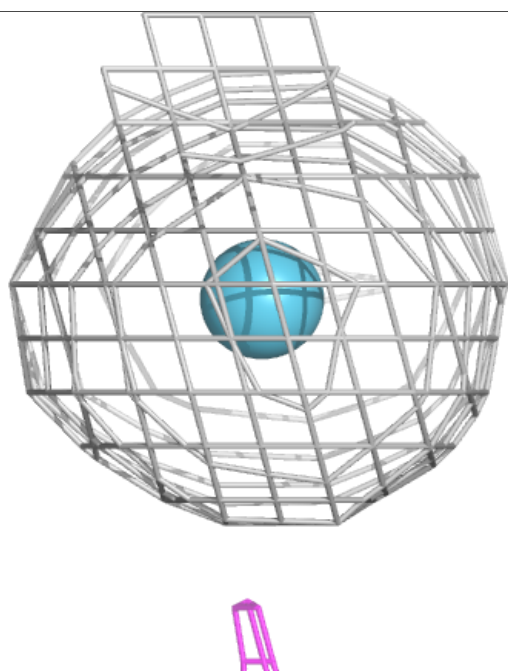
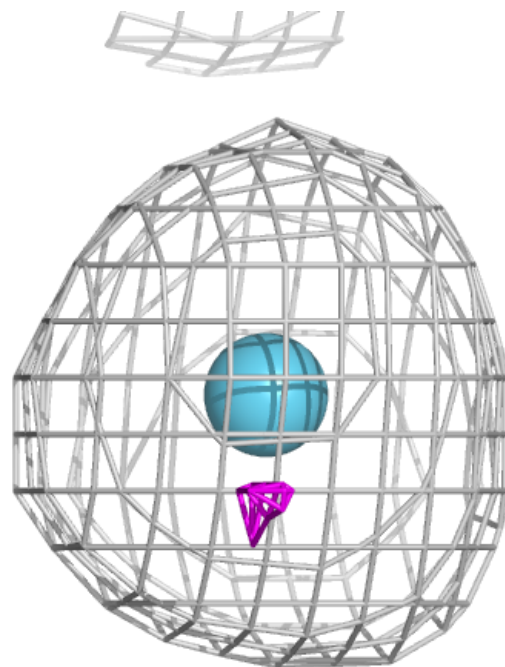
Electron density around KR C 310:

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



Electron density around KR A 308:

$2mF_o-DF_c$ (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.