



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 10:33 AM UTC

PDB ID : 6DKC / pdb_00006dkc
Title : Yeast Ddi2 Cyanamide Hydratase, T157V mutant, apo structure
Authors : Moore, S.A.; Xiao, W.; Li, J.
Deposited on : 2018-05-29
Resolution : 2.90 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

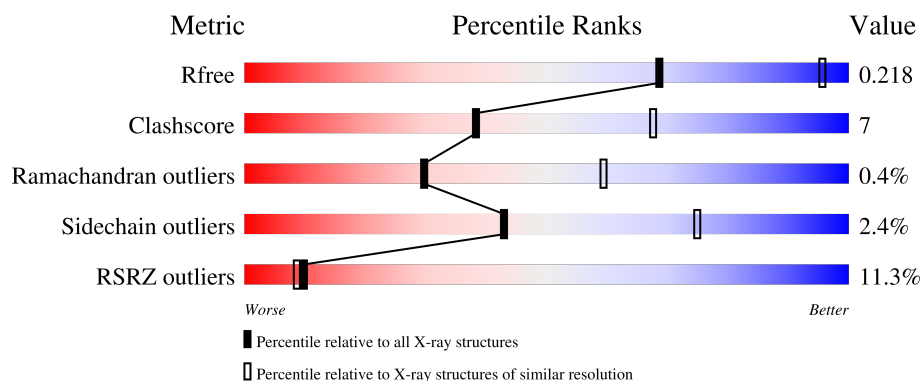
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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	2481 (2.90-2.90)
Clashscore	190562	2690 (2.90-2.90)
Ramachandran outliers	187476	2623 (2.90-2.90)
Sidechain outliers	187428	2625 (2.90-2.90)
RSRZ outliers	180081	2481 (2.90-2.90)

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	234	3% 86% 12% .
1	B	234	2% 89% 8% ..
1	C	234	2% 86% 11% .
1	D	234	% 84% 12% ..
1	E	234	2% 88% 9% ..

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Mol	Chain	Length	Quality of chain
1	F	234	
1	G	234	
1	H	234	
1	I	234	

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
3	SO4	G	302	-	-	X	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 32430 atoms, of which 15869 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 DNA damage-inducible protein.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	230	Total	C	H	N	O	S	0	0	0
			3548	1140	1757	306	339	6			
1	B	230	Total	C	H	N	O	S	0	0	0
			3548	1140	1757	306	339	6			
1	C	229	Total	C	H	N	O	S	0	0	0
			3578	1147	1774	309	342	6			
1	D	229	Total	C	H	N	O	S	0	0	0
			3574	1147	1770	309	342	6			
1	E	229	Total	C	H	N	O	S	0	0	0
			3574	1147	1770	309	342	6			
1	F	232	Total	C	H	N	O	S	0	0	0
			3595	1153	1783	311	342	6			
1	G	229	Total	C	H	N	O	S	0	0	0
			3554	1140	1762	308	338	6			
1	H	227	Total	C	H	N	O	S	0	0	0
			3512	1128	1741	303	335	5			
1	I	226	Total	C	H	N	O	S	0	0	0
			3534	1132	1755	306	335	6			

There are 81 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-7	GLY	-	expression tag	UNP A7A1Y4
A	-6	PRO	-	expression tag	UNP A7A1Y4
A	-5	LEU	-	expression tag	UNP A7A1Y4
A	-4	GLY	-	expression tag	UNP A7A1Y4
A	-3	SER	-	expression tag	UNP A7A1Y4
A	-2	PRO	-	expression tag	UNP A7A1Y4
A	-1	GLU	-	expression tag	UNP A7A1Y4
A	0	PHE	-	expression tag	UNP A7A1Y4
A	157	VAL	THR	engineered mutation	UNP A7A1Y4
B	-7	GLY	-	expression tag	UNP A7A1Y4
B	-6	PRO	-	expression tag	UNP A7A1Y4

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-5	LEU	-	expression tag	UNP A7A1Y4
B	-4	GLY	-	expression tag	UNP A7A1Y4
B	-3	SER	-	expression tag	UNP A7A1Y4
B	-2	PRO	-	expression tag	UNP A7A1Y4
B	-1	GLU	-	expression tag	UNP A7A1Y4
B	0	PHE	-	expression tag	UNP A7A1Y4
B	157	VAL	THR	engineered mutation	UNP A7A1Y4
C	-7	GLY	-	expression tag	UNP A7A1Y4
C	-6	PRO	-	expression tag	UNP A7A1Y4
C	-5	LEU	-	expression tag	UNP A7A1Y4
C	-4	GLY	-	expression tag	UNP A7A1Y4
C	-3	SER	-	expression tag	UNP A7A1Y4
C	-2	PRO	-	expression tag	UNP A7A1Y4
C	-1	GLU	-	expression tag	UNP A7A1Y4
C	0	PHE	-	expression tag	UNP A7A1Y4
C	157	VAL	THR	engineered mutation	UNP A7A1Y4
D	-7	GLY	-	expression tag	UNP A7A1Y4
D	-6	PRO	-	expression tag	UNP A7A1Y4
D	-5	LEU	-	expression tag	UNP A7A1Y4
D	-4	GLY	-	expression tag	UNP A7A1Y4
D	-3	SER	-	expression tag	UNP A7A1Y4
D	-2	PRO	-	expression tag	UNP A7A1Y4
D	-1	GLU	-	expression tag	UNP A7A1Y4
D	0	PHE	-	expression tag	UNP A7A1Y4
D	157	VAL	THR	engineered mutation	UNP A7A1Y4
E	-7	GLY	-	expression tag	UNP A7A1Y4
E	-6	PRO	-	expression tag	UNP A7A1Y4
E	-5	LEU	-	expression tag	UNP A7A1Y4
E	-4	GLY	-	expression tag	UNP A7A1Y4
E	-3	SER	-	expression tag	UNP A7A1Y4
E	-2	PRO	-	expression tag	UNP A7A1Y4
E	-1	GLU	-	expression tag	UNP A7A1Y4
E	0	PHE	-	expression tag	UNP A7A1Y4
E	157	VAL	THR	engineered mutation	UNP A7A1Y4
F	-7	GLY	-	expression tag	UNP A7A1Y4
F	-6	PRO	-	expression tag	UNP A7A1Y4
F	-5	LEU	-	expression tag	UNP A7A1Y4
F	-4	GLY	-	expression tag	UNP A7A1Y4
F	-3	SER	-	expression tag	UNP A7A1Y4
F	-2	PRO	-	expression tag	UNP A7A1Y4
F	-1	GLU	-	expression tag	UNP A7A1Y4
F	0	PHE	-	expression tag	UNP A7A1Y4

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Chain	Residue	Modelled	Actual	Comment	Reference
F	157	VAL	THR	engineered mutation	UNP A7A1Y4
G	-7	GLY	-	expression tag	UNP A7A1Y4
G	-6	PRO	-	expression tag	UNP A7A1Y4
G	-5	LEU	-	expression tag	UNP A7A1Y4
G	-4	GLY	-	expression tag	UNP A7A1Y4
G	-3	SER	-	expression tag	UNP A7A1Y4
G	-2	PRO	-	expression tag	UNP A7A1Y4
G	-1	GLU	-	expression tag	UNP A7A1Y4
G	0	PHE	-	expression tag	UNP A7A1Y4
G	157	VAL	THR	engineered mutation	UNP A7A1Y4
H	-7	GLY	-	expression tag	UNP A7A1Y4
H	-6	PRO	-	expression tag	UNP A7A1Y4
H	-5	LEU	-	expression tag	UNP A7A1Y4
H	-4	GLY	-	expression tag	UNP A7A1Y4
H	-3	SER	-	expression tag	UNP A7A1Y4
H	-2	PRO	-	expression tag	UNP A7A1Y4
H	-1	GLU	-	expression tag	UNP A7A1Y4
H	0	PHE	-	expression tag	UNP A7A1Y4
H	157	VAL	THR	engineered mutation	UNP A7A1Y4
I	-7	GLY	-	expression tag	UNP A7A1Y4
I	-6	PRO	-	expression tag	UNP A7A1Y4
I	-5	LEU	-	expression tag	UNP A7A1Y4
I	-4	GLY	-	expression tag	UNP A7A1Y4
I	-3	SER	-	expression tag	UNP A7A1Y4
I	-2	PRO	-	expression tag	UNP A7A1Y4
I	-1	GLU	-	expression tag	UNP A7A1Y4
I	0	PHE	-	expression tag	UNP A7A1Y4
I	157	VAL	THR	engineered mutation	UNP A7A1Y4

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn) (labeled as "Ligand of Interest" by depositor).

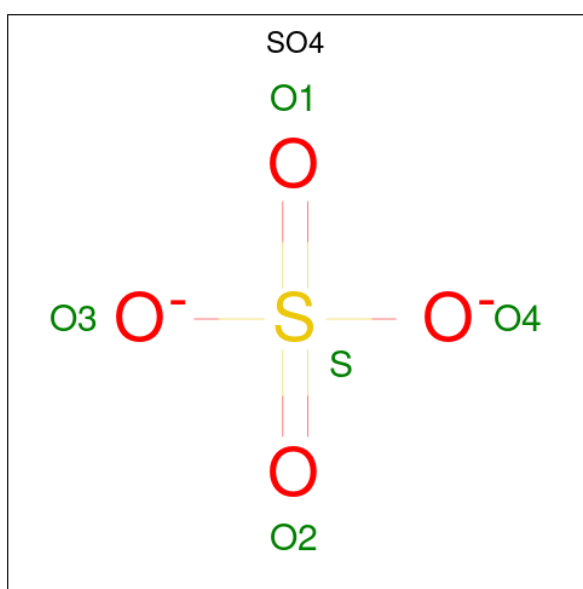
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Zn 1 1	0	0
2	B	1	Total Zn 1 1	0	0
2	C	1	Total Zn 1 1	0	0
2	D	1	Total Zn 1 1	0	0
2	E	1	Total Zn 1 1	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	F	1	Total	Zn	0	0
			1	1		
2	G	1	Total	Zn	0	0
			1	1		
2	H	1	Total	Zn	0	0
			1	1		
2	I	1	Total	Zn	0	0
			1	1		

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		
3	E	1	Total	O	S	0	0
			5	4	1		
3	E	1	Total	O	S	0	0
			5	4	1		
3	E	1	Total	O	S	0	0
			5	4	1		
3	E	1	Total	O	S	0	0
			5	4	1		
3	E	1	Total	O	S	0	0
			5	4	1		
3	E	1	Total	O	S	0	0
			5	4	1		
3	F	1	Total	O	S	0	0
			5	4	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	F	1	Total	O	S	0	0
			5	4	1		
3	F	1	Total	O	S	0	0
			5	4	1		
3	G	1	Total	O	S	0	0
			5	4	1		
3	G	1	Total	O	S	0	0
			5	4	1		

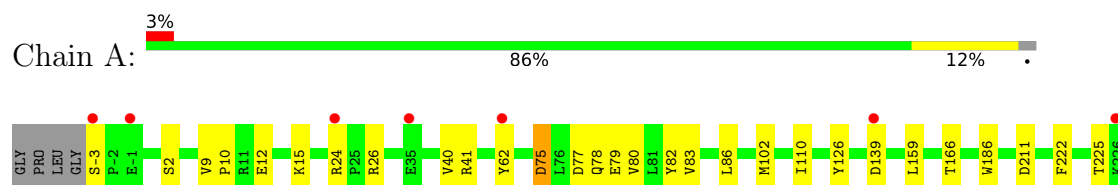
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	36	Total	O	0	0
			36	36		
4	B	34	Total	O	0	0
			34	34		
4	C	48	Total	O	0	0
			48	48		
4	D	36	Total	O	0	0
			36	36		
4	E	45	Total	O	0	0
			45	45		
4	F	28	Total	O	0	0
			28	28		
4	G	17	Total	O	0	0
			17	17		

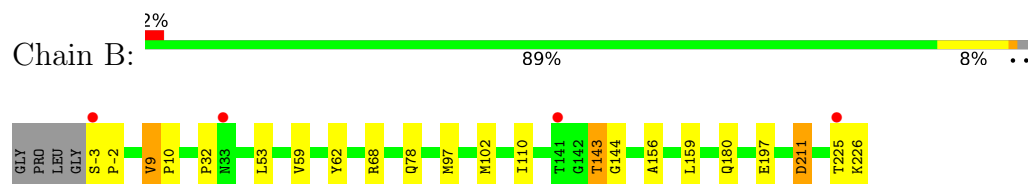
3 Residue-property plots [i](#)

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.

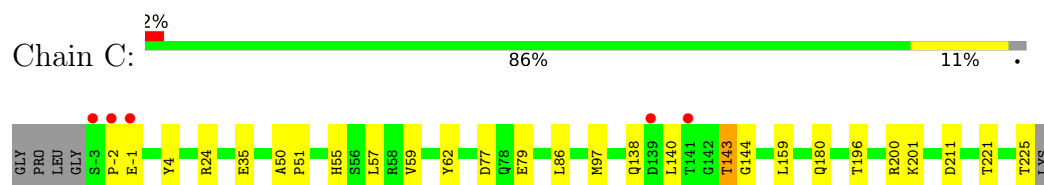
- Molecule 1: DNA damage-inducible protein



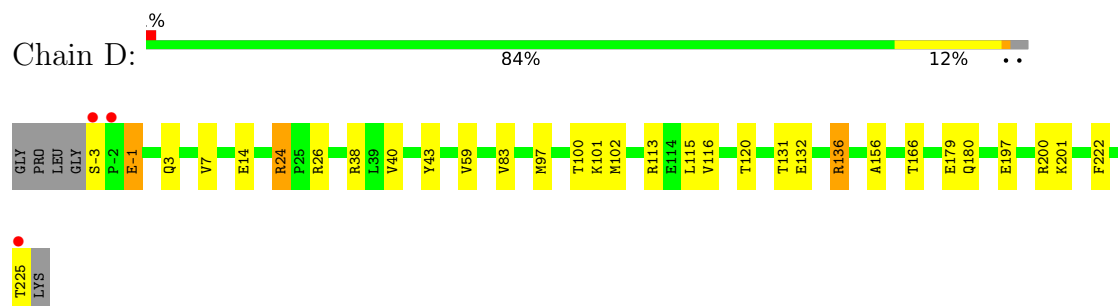
- Molecule 1: DNA damage-inducible protein



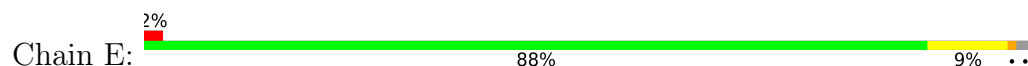
- Molecule 1: DNA damage-inducible protein



- Molecule 1: DNA damage-inducible protein



- Molecule 1: DNA damage-inducible protein





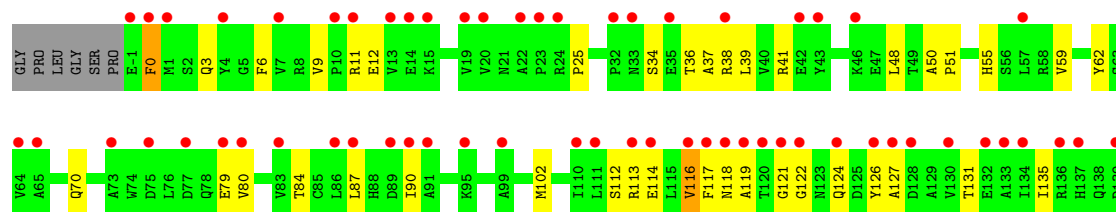
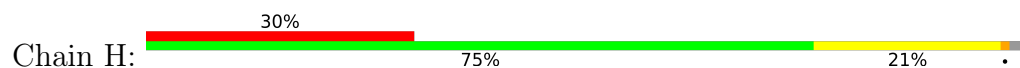
- Molecule 1: DNA damage-inducible protein



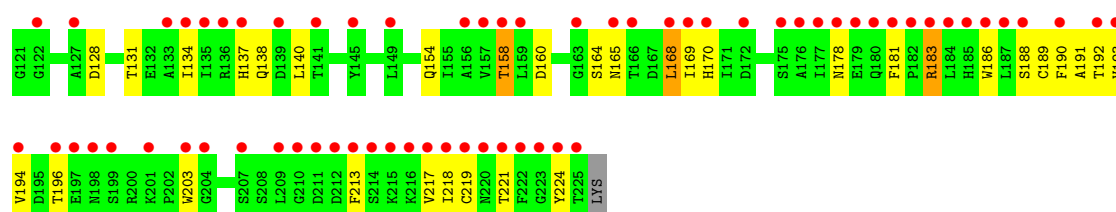
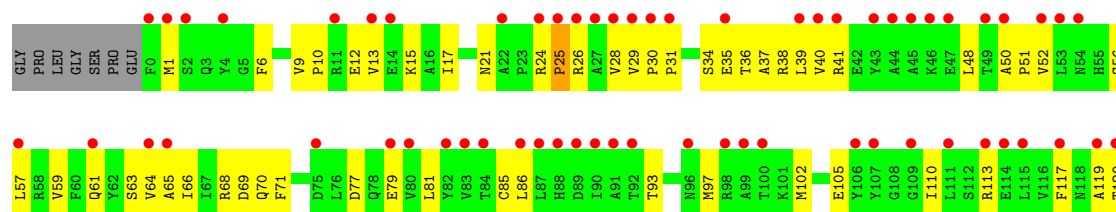
- Molecule 1: DNA damage-inducible protein



- Molecule 1: DNA damage-inducible protein



- Molecule 1: DNA damage-inducible protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 3 2 1	Depositor
Cell constants a, b, c, α , β , γ	263.61Å 263.61Å 119.03Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	39.30 – 2.90 39.30 – 2.90	Depositor EDS
% Data completeness (in resolution range)	97.0 (39.30-2.90) 97.0 (39.30-2.90)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.07 (at 2.90Å)	Xtriage
Refinement program	PHENIX (1.13_2998)	Depositor
R, R_{free}	0.193 , 0.215 0.196 , 0.218	Depositor DCC
R_{free} test set	2578 reflections (2.53%)	wwPDB-VP
Wilson B-factor (Å ²)	53.4	Xtriage
Anisotropy	0.513	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 56.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.011 for -h,-k,l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	32430	wwPDB-VP
Average B, all atoms (Å ²)	76.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.40% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, SO4

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	0.33	0/1833	0.49	0/2506
1	B	0.30	0/1833	0.48	0/2506
1	C	0.33	0/1846	0.51	0/2520
1	D	0.31	0/1846	0.47	0/2520
1	E	0.35	0/1846	0.53	0/2520
1	F	0.29	0/1855	0.47	0/2534
1	G	0.26	0/1834	0.46	0/2506
1	H	0.28	0/1812	0.52	0/2476
1	I	0.30	0/1820	0.57	0/2484
All	All	0.31	0/16525	0.50	0/22572

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	D	0	1
1	H	0	1
All	All	0	3

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	-3	SER	Peptide

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Mol	Chain	Res	Type	Group
1	D	-3	SER	Peptide
1	H	38	ARG	Peptide

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	1791	1757	1746	16	1
1	B	1791	1757	1746	14	0
1	C	1804	1774	1774	15	0
1	D	1804	1770	1774	17	0
1	E	1804	1770	1774	18	0
1	F	1812	1783	1779	11	0
1	G	1792	1762	1755	21	0
1	H	1771	1741	1733	37	1
1	I	1779	1755	1755	85	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
2	E	1	0	0	0	0
2	F	1	0	0	0	0
2	G	1	0	0	0	0
2	H	1	0	0	0	0
2	I	1	0	0	0	0
3	A	15	0	0	1	0
3	B	25	0	0	1	0
3	C	30	0	0	1	0
3	D	35	0	0	1	0
3	E	30	0	0	3	0
3	F	15	0	0	1	0
3	G	10	0	0	4	0
4	A	36	0	0	1	0
4	B	34	0	0	0	0
4	C	48	0	0	1	0
4	D	36	0	0	1	0
4	E	45	0	0	0	0
4	F	28	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	G	17	0	0	1	0
All	All	16561	15869	15836	222	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 7.

All (222) 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:I:178:ASN:OD1	1:I:183:ARG:NH1	1.99	0.95
1:C:24:ARG:NH2	3:C:307:SO4:O1	2.01	0.93
1:I:190:PHE:HB3	1:I:217:VAL:HG11	1.52	0.89
1:F:2:SER:OG	1:F:75:ASP:OD2	1.94	0.85
1:E:2:SER:OG	1:E:75:ASP:OD2	1.98	0.80
1:I:36:THR:HG21	1:I:120:THR:HG22	1.62	0.80
1:I:66:ILE:HD13	1:I:168:LEU:HD21	1.63	0.80
1:A:2:SER:OG	1:A:75:ASP:OD2	1.98	0.80
1:I:165:ASN:O	1:I:168:LEU:HD22	1.82	0.79
1:I:29:VAL:H	1:I:61:GLN:HE22	1.33	0.77
1:C:35:GLU:OE1	4:C:401:HOH:O	2.03	0.76
1:B:68:ARG:NH2	3:B:306:SO4:O2	2.22	0.72
1:F:136:ARG:NH1	4:F:401:HOH:O	2.23	0.72
1:A:12:GLU:OE1	1:A:15:LYS:NZ	2.21	0.71
1:H:117:PHE:HB2	1:H:127:ALA:HB2	1.72	0.71
1:F:1:MET:HE3	1:F:6:PHE:O	1.90	0.71
1:I:154:GLN:O	1:I:158:THR:OG1	2.07	0.71
1:H:114:GLU:O	1:H:118:ASN:HB2	1.92	0.69
1:A:62:TYR:CD2	1:A:159:LEU:HD23	2.28	0.69
1:I:34:SER:O	1:I:38:ARG:N	2.25	0.69
1:I:189:CYS:O	1:I:193:VAL:HG23	1.93	0.69
1:D:101:LYS:NZ	1:E:1:MET:HB3	2.08	0.68
1:D:131:THR:HB	1:E:102:MET:HE2	1.73	0.68
1:E:24:ARG:NH1	3:E:307:SO4:O1	2.27	0.68
1:G:186:TRP:N	3:G:302:SO4:O4	2.27	0.68
1:I:93:THR:O	1:I:97:MET:HG2	1.95	0.67
3:G:302:SO4:O1	4:G:401:HOH:O	2.10	0.67
1:F:32:PRO:O	1:F:41:ARG:NH2	2.28	0.66
1:H:102:MET:HE2	1:I:131:THR:HB	1.77	0.65
1:B:102:MET:HE1	1:B:110:ILE:HD12	1.78	0.65
1:D:43:TYR:CE2	1:D:115:LEU:HD11	2.32	0.65
1:I:41:ARG:NH1	1:I:57:LEU:HD21	2.10	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:139:ASP:OD1	4:A:401:HOH:O	2.15	0.65
1:I:12:GLU:OE1	1:I:15:LYS:NZ	2.29	0.65
1:I:66:ILE:CD1	1:I:168:LEU:HD21	2.27	0.64
1:I:25:PRO:HG3	1:I:170:HIS:ND1	2.14	0.62
1:I:28:VAL:HG13	1:I:61:GLN:NE2	2.14	0.62
1:I:63:SER:CB	1:I:81:LEU:HD21	2.30	0.62
1:E:75:ASP:OD1	1:E:75:ASP:N	2.29	0.61
1:I:36:THR:CG2	1:I:120:THR:HG22	2.29	0.61
1:B:143:THR:HG23	1:B:144:GLY:N	2.16	0.60
1:H:114:GLU:OE2	1:H:118:ASN:ND2	2.34	0.60
1:G:97:MET:HE3	1:G:196:THR:HG22	1.81	0.60
1:I:134:ILE:O	1:I:137:HIS:ND1	2.34	0.60
1:F:1:MET:HE1	1:H:200:ARG:HG2	1.83	0.60
1:I:30:PRO:HB2	1:I:31:PRO:HD2	1.84	0.60
1:I:41:ARG:HH12	1:I:57:LEU:HD11	1.68	0.59
1:H:183:ARG:NH1	1:H:224:TYR:CG	2.71	0.59
1:I:12:GLU:OE1	1:I:15:LYS:CE	2.50	0.59
1:H:37:ALA:HB2	1:H:79:GLU:HG3	1.85	0.58
1:B:9:VAL:HG23	1:B:10:PRO:HD2	1.86	0.58
1:I:26:ARG:HH21	1:I:69:ASP:N	2.03	0.57
1:I:168:LEU:C	1:I:168:LEU:HD23	2.30	0.57
1:A:24:ARG:O	1:A:26:ARG:HG3	2.04	0.56
1:B:211:ASP:OD1	1:B:211:ASP:N	2.30	0.56
1:I:9:VAL:HG22	1:I:10:PRO:HD2	1.88	0.56
1:A:40:VAL:HG11	1:A:83:VAL:HG13	1.88	0.55
1:H:80:VAL:HG22	1:H:126:TYR:CE1	2.41	0.55
1:B:62:TYR:CG	1:B:159:LEU:HD23	2.40	0.55
1:I:102:MET:HE1	1:I:110:ILE:HD12	1.87	0.55
1:I:26:ARG:NH2	1:I:69:ASP:CG	2.65	0.55
1:I:160:ASP:O	1:I:224:TYR:OH	2.23	0.55
1:C:143:THR:CG2	1:C:144:GLY:N	2.70	0.55
1:I:1:MET:HE3	1:I:6:PHE:O	2.07	0.54
1:D:132:GLU:O	1:D:136:ARG:HG3	2.08	0.54
1:E:43:TYR:CE1	1:E:115:LEU:HD11	2.42	0.54
1:H:112:SER:O	1:H:116:VAL:HG23	2.06	0.54
1:I:1:MET:HG2	1:I:6:PHE:HB2	1.90	0.54
1:I:63:SER:HB2	1:I:81:LEU:HD21	1.90	0.54
1:I:63:SER:HB3	1:I:81:LEU:HD21	1.89	0.54
1:I:213:PHE:O	1:I:217:VAL:HG23	2.07	0.54
1:A:12:GLU:OE1	1:A:15:LYS:HD2	2.08	0.54
1:C:143:THR:HG23	1:C:144:GLY:N	2.22	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:117:PHE:HB2	1:H:127:ALA:CB	2.38	0.53
1:I:105:GLU:OE1	1:I:138:GLN:NE2	2.37	0.53
1:D:-1:GLU:N	1:D:3:GLN:OE1	2.42	0.53
1:G:12:GLU:O	1:G:13:VAL:HB	2.07	0.52
1:G:59:VAL:HG13	1:G:156:ALA:HB3	1.91	0.52
1:H:84:THR:OG1	1:H:149:LEU:HD21	2.09	0.52
1:I:36:THR:HG23	1:I:119:ALA:O	2.09	0.52
1:B:143:THR:CG2	1:B:144:GLY:N	2.73	0.51
1:H:117:PHE:CE1	1:H:122:GLY:HA2	2.45	0.51
1:G:97:MET:HE3	1:G:196:THR:CG2	2.41	0.51
1:F:1:MET:HG2	1:F:6:PHE:HB2	1.93	0.51
1:H:50:ALA:HB3	1:H:51:PRO:HD3	1.93	0.51
1:C:97:MET:HE3	1:C:196:THR:HG22	1.92	0.51
1:F:1:MET:CE	1:H:200:ARG:HG2	2.41	0.51
1:H:41:ARG:HH11	1:H:41:ARG:HG2	1.76	0.50
1:G:62:TYR:CD2	1:G:159:LEU:HD23	2.47	0.50
1:I:25:PRO:HG3	1:I:170:HIS:CE1	2.46	0.50
1:I:29:VAL:HG11	1:I:64:VAL:HG11	1.94	0.50
1:D:43:TYR:CZ	1:D:115:LEU:HD11	2.47	0.50
1:D:100:THR:OG1	1:D:102:MET:HG2	2.11	0.50
1:H:25:PRO:HB3	1:H:172:ASP:HB2	1.94	0.50
1:I:28:VAL:HA	1:I:61:GLN:NE2	2.26	0.50
1:C:62:TYR:CD2	1:C:159:LEU:HD23	2.46	0.50
1:B:225:THR:O	1:B:226:LYS:CB	2.60	0.49
1:D:101:LYS:HZ2	1:E:1:MET:HB3	1.73	0.49
1:G:98:ARG:HA	1:G:200:ARG:NH2	2.26	0.49
1:I:29:VAL:N	1:I:61:GLN:HE22	2.06	0.49
1:G:12:GLU:O	1:G:13:VAL:CB	2.59	0.49
1:I:24:ARG:HG3	1:I:69:ASP:OD1	2.12	0.49
1:F:211:ASP:N	1:F:211:ASP:OD1	2.46	0.49
1:E:97:MET:HE2	1:E:196:THR:HG22	1.95	0.48
1:H:102:MET:HE1	1:I:113:ARG:HD2	1.95	0.48
1:H:48:LEU:HD21	1:H:90:ILE:HA	1.96	0.48
1:H:55:HIS:O	1:H:59:VAL:HG23	2.12	0.48
1:I:34:SER:O	1:I:38:ARG:HB2	2.13	0.48
1:I:57:LEU:HD13	1:I:181:PHE:CD1	2.48	0.48
1:H:12:GLU:OE2	1:H:143:THR:HG22	2.13	0.48
1:A:9:VAL:HG23	1:A:10:PRO:HD2	1.95	0.47
1:I:21:ASN:H	1:I:70:GLN:HE22	1.62	0.47
1:I:39:LEU:C	1:I:39:LEU:HD23	2.38	0.47
1:I:56:SER:HB3	1:I:86:LEU:HA	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:77:ASP:OD1	1:A:79:GLU:N	2.45	0.47
1:A:80:VAL:HG22	1:A:126:TYR:CE1	2.49	0.47
1:A:186:TRP:N	3:A:303:SO4:O1	2.47	0.47
1:A:102:MET:HE1	1:A:110:ILE:HD12	1.96	0.47
1:I:35:GLU:HA	1:I:38:ARG:HB3	1.96	0.47
3:E:305:SO4:O1	1:G:1:MET:HB3	2.15	0.47
1:C:200:ARG:HG2	1:E:1:MET:HE2	1.97	0.46
1:H:116:VAL:HG12	1:H:117:PHE:N	2.31	0.46
1:E:97:MET:HE1	1:E:197:GLU:HA	1.96	0.46
1:I:34:SER:O	1:I:38:ARG:CB	2.64	0.46
1:C:62:TYR:CG	1:C:159:LEU:HD23	2.51	0.46
1:C:143:THR:HG23	1:C:144:GLY:H	1.81	0.46
1:I:26:ARG:NH2	1:I:68:ARG:HB2	2.31	0.46
1:I:57:LEU:HD13	1:I:181:PHE:CE1	2.50	0.46
1:I:77:ASP:OD1	1:I:79:GLU:N	2.47	0.46
1:G:9:VAL:CG2	1:G:10:PRO:HD2	2.46	0.46
1:H:87:LEU:O	1:H:90:ILE:HG22	2.15	0.46
1:I:12:GLU:OE1	1:I:15:LYS:HE3	2.15	0.46
1:I:37:ALA:HB2	1:I:79:GLU:HG3	1.96	0.46
1:G:9:VAL:HG22	1:G:10:PRO:HD2	1.97	0.46
1:H:183:ARG:NH1	1:H:224:TYR:CD1	2.83	0.46
1:H:183:ARG:HH21	1:H:186:TRP:HD1	1.62	0.46
1:B:225:THR:O	1:B:225:THR:HG22	2.15	0.46
1:D:113:ARG:NH1	4:D:405:HOH:O	2.49	0.46
1:H:11:ARG:NH1	1:H:154:GLN:OE1	2.49	0.46
1:C:2:PRO:HA	1:C:4:TYR:OH	2.16	0.45
1:I:50:ALA:HB3	1:I:51:PRO:HD3	1.98	0.45
1:I:186:TRP:CZ2	1:I:190:PHE:CE1	3.04	0.45
1:F:186:TRP:N	3:F:304:SO4:O2	2.46	0.45
1:G:2:SER:HB3	1:G:7:VAL:CG2	2.46	0.45
1:H:0:PHE:HB2	1:H:3:GLN:HG3	1.98	0.45
1:H:6:PHE:HB3	1:I:203:TRP:CH2	2.52	0.45
1:I:26:ARG:CD	1:I:65:ALA:HA	2.45	0.45
1:C:55:HIS:O	1:C:59:VAL:HG23	2.17	0.45
1:I:26:ARG:HD2	1:I:65:ALA:HA	1.98	0.45
1:E:50:ALA:HB3	1:E:51:PRO:HD3	1.99	0.45
1:C:138:GLN:HA	1:C:140:LEU:HG	1.97	0.45
1:E:43:TYR:CZ	1:E:115:LEU:HD11	2.51	0.45
1:D:97:MET:HE1	1:D:197:GLU:OE1	2.17	0.44
1:I:17:ILE:HD11	1:I:71:PHE:CZ	2.52	0.44
1:I:191:ALA:CB	1:I:218:ILE:HD11	2.47	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:117:PHE:CE1	1:H:124:GLN:NE2	2.86	0.44
1:G:59:VAL:HG13	1:G:156:ALA:CB	2.47	0.44
1:G:187:LEU:N	3:G:302:SO4:O4	2.38	0.44
1:G:224:TYR:O	1:G:225:THR:C	2.60	0.44
1:I:188:SER:HA	1:I:218:ILE:CD1	2.48	0.44
1:H:183:ARG:HH21	1:H:186:TRP:CD1	2.36	0.44
1:B:32:PRO:HG3	1:B:78:GLN:HG2	1.99	0.43
1:H:41:ARG:HG2	1:H:41:ARG:NH1	2.34	0.43
1:A:41:ARG:HA	1:A:86:LEU:HD13	2.00	0.43
1:G:35:GLU:O	1:G:39:LEU:HB2	2.18	0.43
1:I:66:ILE:HG23	1:I:168:LEU:HG	1.99	0.43
1:C:50:ALA:HB3	1:C:51:PRO:HD3	2.01	0.43
1:H:131:THR:HG22	1:H:135:ILE:HD12	1.99	0.43
1:I:158:THR:HG22	1:I:164:SER:HB2	2.00	0.43
1:E:77:ASP:OD1	1:E:79:GLU:N	2.51	0.43
1:H:113:ARG:CD	1:I:110:ILE:HD13	2.49	0.43
1:I:190:PHE:O	1:I:194:VAL:HG23	2.19	0.43
1:D:24:ARG:O	1:D:26:ARG:HG3	2.19	0.43
1:D:116:VAL:O	1:D:120:THR:HG23	2.19	0.43
1:E:215:LYS:HG3	1:G:18:PRO:HG3	2.00	0.43
1:G:12:GLU:C	1:G:13:VAL:HG23	2.44	0.43
1:I:40:VAL:HG12	1:I:86:LEU:CD1	2.49	0.43
1:C:57:LEU:HD21	1:C:86:LEU:HD21	2.01	0.43
1:A:166:THR:HG21	1:A:222:PHE:CE1	2.53	0.42
1:E:26:ARG:HE	1:E:69:ASP:CG	2.27	0.42
1:H:117:PHE:CB	1:H:127:ALA:HB2	2.46	0.42
1:I:97:MET:HE3	1:I:196:THR:HG22	2.00	0.42
1:I:160:ASP:C	1:I:224:TYR:OH	2.63	0.42
1:I:117:PHE:CD1	1:I:117:PHE:C	2.98	0.42
1:I:158:THR:HG22	1:I:164:SER:CB	2.50	0.42
1:H:59:VAL:HG13	1:H:156:ALA:HB3	2.01	0.42
1:D:200:ARG:NH2	3:D:301:SO4:O1	2.53	0.42
1:B:97:MET:HE1	1:B:197:GLU:OE1	2.19	0.42
1:D:166:THR:HG21	1:D:222:PHE:CE1	2.54	0.42
1:F:50:ALA:HB3	1:F:51:PRO:HD3	2.02	0.42
1:I:13:VAL:O	1:I:13:VAL:HG12	2.19	0.42
1:I:97:MET:CE	1:I:196:THR:HG22	2.50	0.42
1:E:9:VAL:CG2	1:E:10:PRO:HD2	2.50	0.42
1:I:59:VAL:HG12	1:I:85:CYS:SG	2.59	0.42
1:I:190:PHE:HA	1:I:193:VAL:HB	2.02	0.42
1:I:219:CYS:O	1:I:221:THR:HG23	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:68:ARG:NH2	3:G:303:SO4:O4	2.53	0.41
1:I:28:VAL:HA	1:I:61:GLN:HE22	1.85	0.41
1:C:77:ASP:OD1	1:C:79:GLU:N	2.53	0.41
1:H:62:TYR:CD2	1:H:159:LEU:HD23	2.55	0.41
1:I:26:ARG:NH2	1:I:68:ARG:CB	2.84	0.41
1:F:34:SER:OG	1:F:79:GLU:OE1	2.34	0.41
1:A:211:ASP:OD1	1:A:211:ASP:N	2.53	0.41
1:H:102:MET:CE	1:I:128:ASP:HA	2.50	0.41
1:H:183:ARG:HE	1:H:183:ARG:HB3	1.70	0.41
1:E:186:TRP:N	3:E:303:SO4:O1	2.50	0.41
1:A:75:ASP:OD1	1:A:75:ASP:N	2.53	0.41
1:B:62:TYR:CD2	1:B:159:LEU:HD23	2.56	0.41
1:I:35:GLU:O	1:I:39:LEU:N	2.31	0.41
1:I:186:TRP:CZ2	1:I:190:PHE:CD1	3.09	0.41
1:D:40:VAL:HG11	1:D:83:VAL:HG13	2.03	0.41
1:D:59:VAL:HG13	1:D:156:ALA:CB	2.51	0.41
1:D:101:LYS:HZ3	1:E:1:MET:HB3	1.81	0.41
1:B:-3:SER:CB	1:B:-2:PRO:HD3	2.51	0.40
1:B:59:VAL:HG13	1:B:156:ALA:HB3	2.03	0.40
1:G:166:THR:HG21	1:G:222:PHE:CE1	2.56	0.40
1:I:97:MET:CE	1:I:196:THR:CG2	2.98	0.40
1:I:48:LEU:HD22	1:I:52:VAL:HG12	2.04	0.40
1:I:138:GLN:HA	1:I:140:LEU:HG	2.03	0.40
1:G:62:TYR:CG	1:G:159:LEU:HD23	2.57	0.40
1:I:117:PHE:C	1:I:117:PHE:HD1	2.29	0.40
1:I:168:LEU:HD23	1:I:169:ILE:N	2.36	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:82:TYR:OH	1:H:121:GLY:O[3_566]	1.83	0.37

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	228/234 (97%)	222 (97%)	6 (3%)	0	100	100
1	B	228/234 (97%)	224 (98%)	4 (2%)	0	100	100
1	C	227/234 (97%)	223 (98%)	3 (1%)	1 (0%)	30	59
1	D	227/234 (97%)	224 (99%)	3 (1%)	0	100	100
1	E	227/234 (97%)	223 (98%)	4 (2%)	0	100	100
1	F	230/234 (98%)	224 (97%)	4 (2%)	2 (1%)	14	41
1	G	227/234 (97%)	223 (98%)	3 (1%)	1 (0%)	30	59
1	H	225/234 (96%)	214 (95%)	7 (3%)	4 (2%)	6	25
1	I	224/234 (96%)	216 (96%)	7 (3%)	1 (0%)	30	59
All	All	2043/2106 (97%)	1993 (98%)	41 (2%)	9 (0%)	30	59

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	-1	GLU
1	G	13	VAL
1	H	116	VAL
1	H	119	ALA
1	F	-3	SER
1	H	0	PHE
1	H	34	SER
1	F	139	ASP
1	I	25	PRO

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	194/202 (96%)	191 (98%)	3 (2%)	57	84
1	B	194/202 (96%)	189 (97%)	5 (3%)	40	73

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	198/202 (98%)	192 (97%)	6 (3%)	36	70
1	D	198/202 (98%)	188 (95%)	10 (5%)	21	53
1	E	198/202 (98%)	194 (98%)	4 (2%)	48	78
1	F	198/202 (98%)	196 (99%)	2 (1%)	68	89
1	G	195/202 (96%)	192 (98%)	3 (2%)	57	84
1	H	192/202 (95%)	187 (97%)	5 (3%)	40	73
1	I	195/202 (96%)	191 (98%)	4 (2%)	47	77
All	All	1762/1818 (97%)	1720 (98%)	42 (2%)	43	75

All (42) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	75	ASP
1	A	78	GLN
1	A	225	THR
1	B	9	VAL
1	B	53	LEU
1	B	143	THR
1	B	180	GLN
1	B	211	ASP
1	C	143	THR
1	C	180	GLN
1	C	201	LYS
1	C	211	ASP
1	C	221	THR
1	C	225	THR
1	D	-1	GLU
1	D	7	VAL
1	D	14	GLU
1	D	24	ARG
1	D	38	ARG
1	D	136	ARG
1	D	179	GLU
1	D	180	GLN
1	D	201	LYS
1	D	225	THR
1	E	1	MET
1	E	75	ASP
1	E	117	PHE
1	E	201	LYS

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Mol	Chain	Res	Type
1	F	201	LYS
1	F	211	ASP
1	G	2	SER
1	G	12	GLU
1	G	39	LEU
1	H	9	VAL
1	H	36	THR
1	H	39	LEU
1	H	70	GLN
1	H	225	THR
1	I	158	THR
1	I	168	LEU
1	I	183	ARG
1	I	192	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	78	GLN
1	B	124	GLN
1	B	165	ASN
1	B	180	GLN
1	E	3	GLN
1	H	124	GLN
1	H	165	ASN
1	I	61	GLN
1	I	70	GLN
1	I	165	ASN
1	I	198	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 41 ligands modelled in this entry, 9 are monoatomic - leaving 32 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	SO4	E	302	-	4,4,4	0.28	0	6,6,6	0.09	0
3	SO4	B	301	-	4,4,4	0.26	0	6,6,6	0.15	0
3	SO4	C	306	-	4,4,4	0.28	0	6,6,6	0.13	0
3	SO4	F	301	-	4,4,4	0.20	0	6,6,6	0.21	0
3	SO4	C	307	-	4,4,4	0.27	0	6,6,6	0.17	0
3	SO4	E	305	-	4,4,4	0.21	0	6,6,6	0.36	0
3	SO4	F	303	-	4,4,4	0.24	0	6,6,6	0.18	0
3	SO4	D	304	-	4,4,4	0.17	0	6,6,6	0.24	0
3	SO4	E	304	-	4,4,4	0.29	0	6,6,6	0.22	0
3	SO4	G	303	-	4,4,4	0.28	0	6,6,6	0.21	0
3	SO4	D	305	-	4,4,4	0.26	0	6,6,6	0.16	0
3	SO4	D	308	-	4,4,4	0.30	0	6,6,6	0.19	0
3	SO4	E	306	-	4,4,4	0.30	0	6,6,6	0.18	0
3	SO4	D	306	-	4,4,4	0.26	0	6,6,6	0.16	0
3	SO4	E	307	-	4,4,4	0.13	0	6,6,6	0.40	0
3	SO4	B	305	-	4,4,4	0.27	0	6,6,6	0.17	0
3	SO4	C	303	-	4,4,4	0.29	0	6,6,6	0.16	0
3	SO4	B	306	-	4,4,4	0.16	0	6,6,6	0.23	0
3	SO4	G	302	-	4,4,4	0.29	0	6,6,6	0.20	0
3	SO4	F	304	-	4,4,4	0.18	0	6,6,6	0.12	0
3	SO4	A	303	-	4,4,4	0.23	0	6,6,6	0.32	0
3	SO4	D	307	-	4,4,4	0.26	0	6,6,6	0.14	0
3	SO4	B	303	-	4,4,4	0.21	0	6,6,6	0.16	0
3	SO4	D	301	-	4,4,4	0.31	0	6,6,6	0.19	0
3	SO4	C	302	-	4,4,4	0.31	0	6,6,6	0.23	0
3	SO4	E	303	-	4,4,4	0.28	0	6,6,6	0.20	0
3	SO4	A	304	-	4,4,4	0.18	0	6,6,6	0.19	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	SO4	B	304	-	4,4,4	0.21	0	6,6,6	0.18	0
3	SO4	C	304	-	4,4,4	0.23	0	6,6,6	0.12	0
3	SO4	D	303	-	4,4,4	0.27	0	6,6,6	0.08	0
3	SO4	A	302	-	4,4,4	0.25	0	6,6,6	0.10	0
3	SO4	C	305	-	4,4,4	0.22	0	6,6,6	0.30	0

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.

10 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	307	SO4	1	0
3	E	305	SO4	1	0
3	G	303	SO4	1	0
3	E	307	SO4	1	0
3	B	306	SO4	1	0
3	G	302	SO4	3	0
3	F	304	SO4	1	0
3	A	303	SO4	1	0
3	D	301	SO4	1	0
3	E	303	SO4	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	230/234 (98%)	-0.22	7 (3%)	52	43	43, 60, 89, 135	0
1	B	230/234 (98%)	-0.40	4 (1%)	69	61	36, 56, 79, 131	0
1	C	229/234 (97%)	-0.43	5 (2%)	62	53	33, 47, 72, 114	0
1	D	229/234 (97%)	-0.47	3 (1%)	75	67	36, 49, 72, 105	0
1	E	229/234 (97%)	-0.47	4 (1%)	69	61	34, 47, 70, 145	0
1	F	232/234 (99%)	-0.28	9 (3%)	43	35	40, 54, 88, 139	0
1	G	229/234 (97%)	-0.14	5 (2%)	62	53	45, 71, 100, 119	0
1	H	227/234 (97%)	1.54	70 (30%)	1	1	74, 124, 168, 200	0
1	I	226/234 (96%)	2.32	126 (55%)	0	0	97, 163, 205, 261	0
All	All	2061/2106 (97%)	0.16	233 (11%)	10	9	33, 60, 169, 261	0

All (233) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	-6	PRO	7.8
1	I	91	ALA	6.9
1	C	139	ASP	6.8
1	E	-3	SER	6.8
1	B	-3	SER	6.7
1	H	225	THR	6.4
1	I	90	ILE	6.3
1	H	111	LEU	6.2
1	I	225	THR	5.9
1	D	-3	SER	5.9
1	H	0	PHE	5.8
1	I	57	LEU	5.2
1	E	225	THR	5.0
1	I	39	LEU	5.0
1	H	119	ALA	4.8

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Mol	Chain	Res	Type	RSRZ
1	I	192	THR	4.7
1	I	0	PHE	4.7
1	H	127	ALA	4.7
1	I	27	ALA	4.7
1	H	91	ALA	4.5
1	H	38	ARG	4.5
1	I	82	TYR	4.5
1	I	115	LEU	4.4
1	I	84	THR	4.4
1	F	-1	GLU	4.4
1	I	86	LEU	4.4
1	C	-3	SER	4.3
1	I	196	THR	4.3
1	I	43	TYR	4.3
1	H	120	THR	4.2
1	H	180	GLN	4.2
1	I	190	PHE	4.2
1	I	96	ASN	4.1
1	I	83	VAL	4.1
1	I	204	GLY	4.1
1	H	7	VAL	4.1
1	I	50	ALA	4.0
1	H	43	TYR	4.0
1	I	181	PHE	4.0
1	C	-2	PRO	4.0
1	I	201	LYS	4.0
1	I	213	PHE	4.0
1	I	193	VAL	4.0
1	I	176	ALA	4.0
1	I	177	ILE	4.0
1	G	225	THR	4.0
1	I	186	TRP	4.0
1	I	216	LYS	4.0
1	H	35	GLU	3.9
1	I	170	HIS	3.9
1	I	87	LEU	3.9
1	H	126	TYR	3.9
1	I	169	ILE	3.9
1	I	53	LEU	3.8
1	I	211	ASP	3.8
1	H	117	PHE	3.8
1	H	-1	GLU	3.7

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Mol	Chain	Res	Type	RSRZ
1	H	14	GLU	3.7
1	I	183	ARG	3.7
1	G	35	GLU	3.7
1	I	64	VAL	3.7
1	I	106	TYR	3.7
1	I	88	HIS	3.6
1	G	-3	SER	3.6
1	F	141	THR	3.6
1	I	2	SER	3.6
1	H	121	GLY	3.6
1	C	-1	GLU	3.6
1	I	1	MET	3.6
1	H	116	VAL	3.5
1	I	219	CYS	3.5
1	H	23	PRO	3.5
1	H	145	TYR	3.5
1	H	75	ASP	3.4
1	I	185	HIS	3.4
1	I	139	ASP	3.4
1	I	117	PHE	3.4
1	H	124	GLN	3.3
1	I	49	THR	3.3
1	I	203	TRP	3.3
1	I	221	THR	3.3
1	H	24	ARG	3.3
1	D	225	THR	3.3
1	I	212	ASP	3.3
1	H	83	VAL	3.3
1	I	157	VAL	3.3
1	I	41	ARG	3.2
1	I	14	GLU	3.2
1	I	179	GLU	3.2
1	A	-3	SER	3.2
1	A	24	ARG	3.1
1	I	199	SER	3.1
1	I	210	GLY	3.1
1	I	30	PRO	3.1
1	I	184	LEU	3.1
1	H	118	ASN	3.1
1	I	180	GLN	3.1
1	F	-4	GLY	3.1
1	I	47	GLU	3.1

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Mol	Chain	Res	Type	RSRZ
1	I	65	ALA	3.1
1	I	197	GLU	3.1
1	I	217	VAL	3.0
1	I	107	TYR	3.0
1	E	33	ASN	3.0
1	I	92	THR	3.0
1	H	185	HIS	3.0
1	I	120	THR	2.9
1	H	136	ARG	2.9
1	H	134	ILE	2.9
1	I	222	PHE	2.9
1	I	182	PRO	2.9
1	I	218	ILE	2.9
1	A	62	TYR	2.8
1	B	225	THR	2.8
1	I	220	ASN	2.8
1	H	86	LEU	2.8
1	I	31	PRO	2.8
1	I	26	ARG	2.8
1	I	141	THR	2.8
1	I	29	VAL	2.8
1	I	99	ALA	2.8
1	I	109	GLY	2.7
1	F	-3	SER	2.7
1	H	19	VAL	2.7
1	H	80	VAL	2.7
1	H	149	LEU	2.7
1	A	226	LYS	2.7
1	H	33	ASN	2.7
1	H	143	THR	2.7
1	I	35	GLU	2.7
1	H	87	LEU	2.7
1	I	22	ALA	2.7
1	I	133	ALA	2.7
1	I	100	THR	2.7
1	I	44	ALA	2.6
1	I	4	TYR	2.6
1	I	11	ARG	2.6
1	H	64	VAL	2.6
1	H	73	ALA	2.6
1	H	133	ALA	2.6
1	I	98	ARG	2.6

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Mol	Chain	Res	Type	RSRZ
1	H	168	LEU	2.6
1	H	20	VAL	2.6
1	H	114	GLU	2.6
1	H	132	GLU	2.6
1	H	46	LYS	2.5
1	F	1	MET	2.5
1	H	110	ILE	2.5
1	I	13	VAL	2.5
1	I	135	ILE	2.5
1	I	114	GLU	2.5
1	I	136	ARG	2.5
1	I	215	LYS	2.5
1	H	32	PRO	2.5
1	H	57	LEU	2.5
1	I	187	LEU	2.5
1	I	209	LEU	2.5
1	H	42	GLU	2.5
1	E	98	ARG	2.5
1	H	1	MET	2.5
1	I	149	LEU	2.5
1	H	13	VAL	2.5
1	I	28	VAL	2.5
1	I	46	LYS	2.5
1	I	178	ASN	2.5
1	I	145	TYR	2.4
1	F	225	THR	2.4
1	A	139	ASP	2.4
1	H	15	LYS	2.4
1	I	113	ARG	2.4
1	I	175	SER	2.4
1	H	22	ALA	2.4
1	I	89	ASP	2.4
1	I	188	SER	2.4
1	C	141	THR	2.4
1	I	166	THR	2.4
1	I	168	LEU	2.4
1	H	10	PRO	2.4
1	H	99	ALA	2.3
1	I	194	VAL	2.3
1	I	224	TYR	2.3
1	D	-2	PRO	2.3
1	I	24	ARG	2.3

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Mol	Chain	Res	Type	RSRZ
1	H	128	ASP	2.3
1	I	134	ILE	2.3
1	I	40	VAL	2.3
1	I	198	ASN	2.3
1	I	172	ASP	2.3
1	H	90	ILE	2.3
1	H	4	TYR	2.3
1	H	122	GLY	2.3
1	G	139	ASP	2.2
1	I	207	SER	2.2
1	I	45	ALA	2.2
1	I	159	LEU	2.2
1	H	77	ASP	2.2
1	I	122	GLY	2.2
1	H	95	LYS	2.2
1	F	211	ASP	2.2
1	B	33	ASN	2.2
1	I	61	GLN	2.2
1	H	89	ASP	2.2
1	H	157	VAL	2.2
1	I	111	LEU	2.2
1	I	137	HIS	2.1
1	H	141	THR	2.1
1	I	156	ALA	2.1
1	H	113	ARG	2.1
1	I	79	GLU	2.1
1	I	54	ASN	2.1
1	H	65	ALA	2.1
1	I	127	ALA	2.1
1	I	214	SER	2.1
1	H	137	HIS	2.1
1	H	130	VAL	2.1
1	I	223	GLY	2.1
1	H	147	THR	2.1
1	H	11	ARG	2.1
1	I	165	ASN	2.1
1	G	211	ASP	2.1
1	H	139	ASP	2.1
1	A	35	GLU	2.1
1	H	79	GLU	2.1
1	I	52	VAL	2.0
1	I	80	VAL	2.0

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Mol	Chain	Res	Type	RSRZ
1	I	25	PRO	2.0
1	F	-5	LEU	2.0
1	A	-1	GLU	2.0
1	I	163	GLY	2.0
1	B	141	THR	2.0
1	I	119	ALA	2.0
1	I	158	THR	2.0
1	I	75	ASP	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
2	ZN	I	500	1/1	0.76	0.18	95,95,95,95	1
3	SO4	B	305	5/5	0.78	0.20	104,106,109,113	5
3	SO4	B	306	5/5	0.83	0.17	78,81,83,84	5
3	SO4	C	305	5/5	0.85	0.25	68,68,76,78	5
3	SO4	C	304	5/5	0.86	0.14	84,87,88,92	5
3	SO4	C	307	5/5	0.86	0.15	65,73,77,79	5
3	SO4	D	308	5/5	0.87	0.14	76,79,82,82	5
3	SO4	G	302	5/5	0.87	0.14	65,71,73,74	5
2	ZN	H	500	1/1	0.88	0.12	100,100,100,100	1
3	SO4	D	306	5/5	0.88	0.14	73,81,82,86	5
3	SO4	B	301	5/5	0.90	0.14	82,89,97,98	0
3	SO4	D	301	5/5	0.90	0.14	85,90,92,94	0
3	SO4	E	306	5/5	0.90	0.19	59,63,63,65	5
3	SO4	D	303	5/5	0.90	0.17	95,97,99,101	0
3	SO4	G	303	5/5	0.90	0.12	59,63,66,66	5

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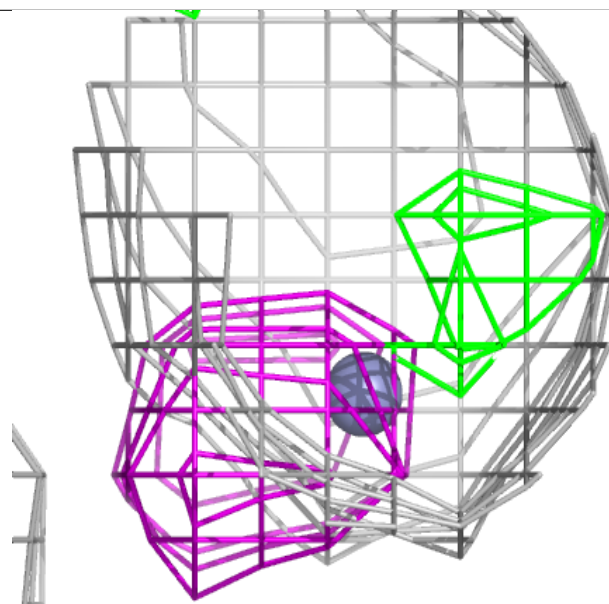
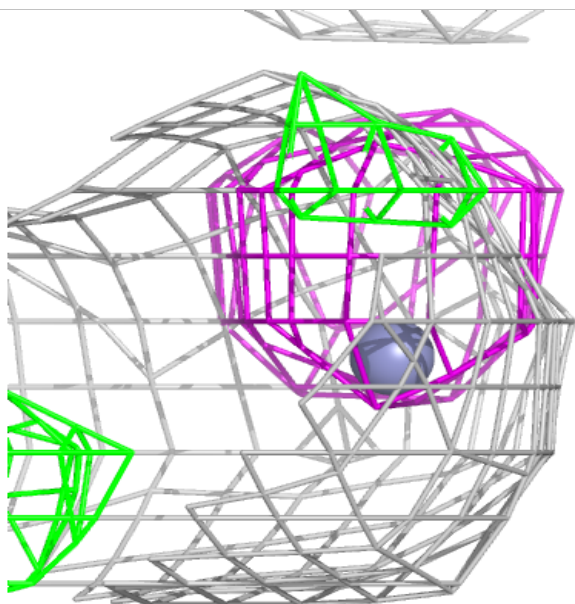
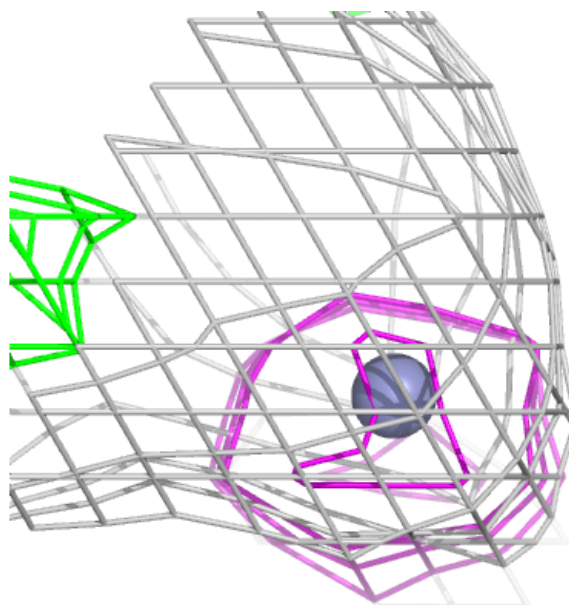
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	SO4	D	305	5/5	0.91	0.19	84,89,95,96	5
3	SO4	B	304	5/5	0.91	0.16	47,55,57,60	5
3	SO4	C	306	5/5	0.92	0.12	66,70,74,74	5
3	SO4	A	302	5/5	0.92	0.12	67,73,74,79	5
3	SO4	F	303	5/5	0.93	0.11	71,72,74,77	5
3	SO4	E	303	5/5	0.93	0.12	43,44,47,47	5
3	SO4	B	303	5/5	0.93	0.15	79,86,91,92	5
2	ZN	G	301	1/1	0.94	0.09	69,69,69,69	1
3	SO4	D	307	5/5	0.94	0.09	67,70,70,71	5
3	SO4	D	304	5/5	0.94	0.14	76,79,81,86	0
3	SO4	A	304	5/5	0.94	0.11	42,51,55,56	5
3	SO4	A	303	5/5	0.95	0.13	55,55,56,61	5
3	SO4	E	304	5/5	0.95	0.08	51,52,56,57	5
3	SO4	F	301	5/5	0.96	0.10	53,57,62,62	5
3	SO4	E	305	5/5	0.96	0.09	43,49,53,53	5
3	SO4	F	304	5/5	0.96	0.10	52,53,58,59	5
3	SO4	C	302	5/5	0.96	0.10	38,41,42,43	5
3	SO4	E	307	5/5	0.96	0.09	38,47,49,51	5
2	ZN	B	302	1/1	0.97	0.05	59,59,59,59	1
3	SO4	E	302	5/5	0.97	0.09	64,66,69,71	5
2	ZN	E	301	1/1	0.97	0.05	53,53,53,53	1
2	ZN	A	301	1/1	0.98	0.06	51,51,51,51	1
3	SO4	C	303	5/5	0.98	0.08	43,56,59,62	5
2	ZN	C	301	1/1	0.98	0.04	44,44,44,44	1
2	ZN	D	302	1/1	0.98	0.06	58,58,58,58	1
2	ZN	F	302	1/1	0.99	0.04	57,57,57,57	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.

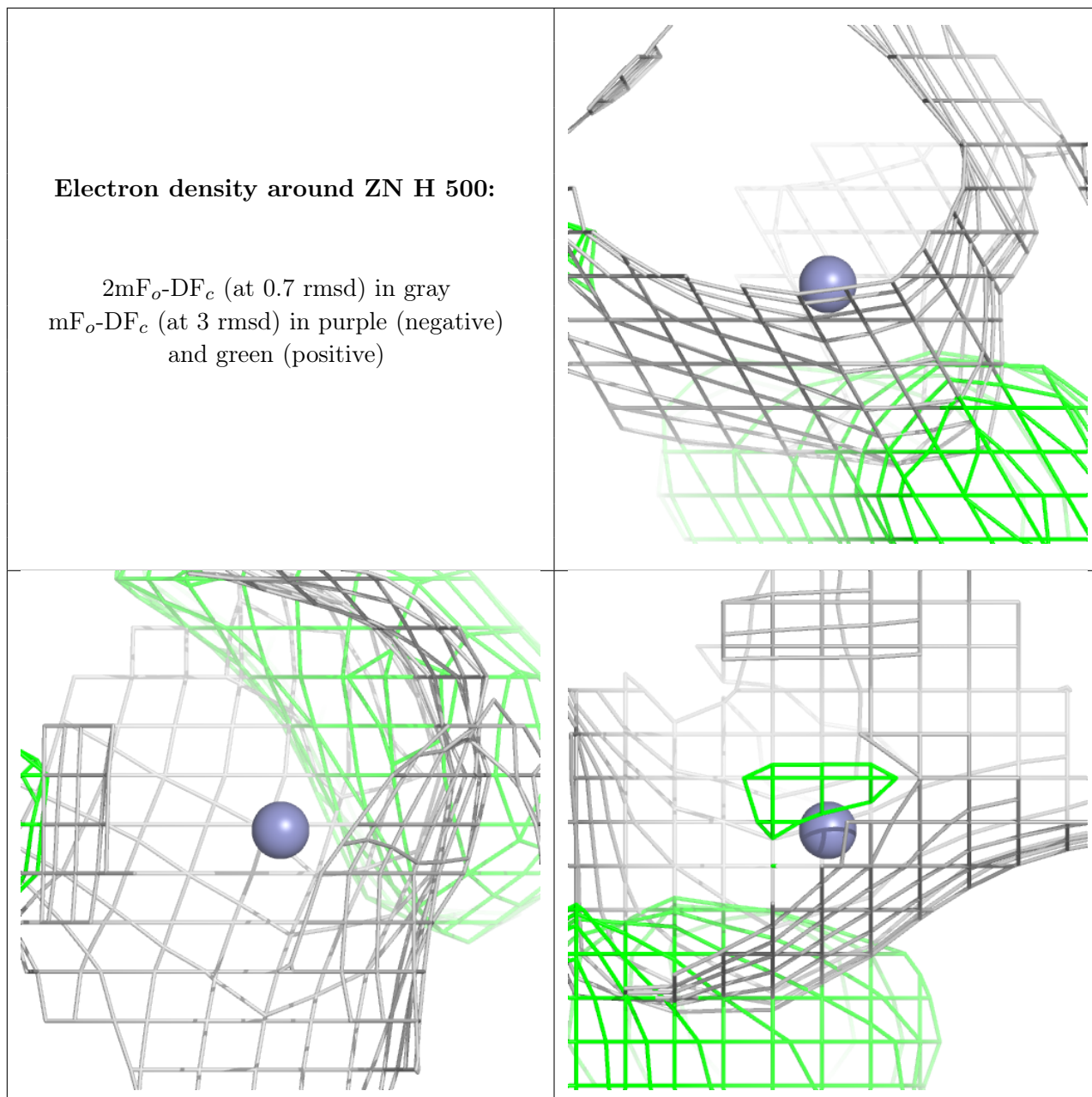
Electron density around ZN I 500:

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



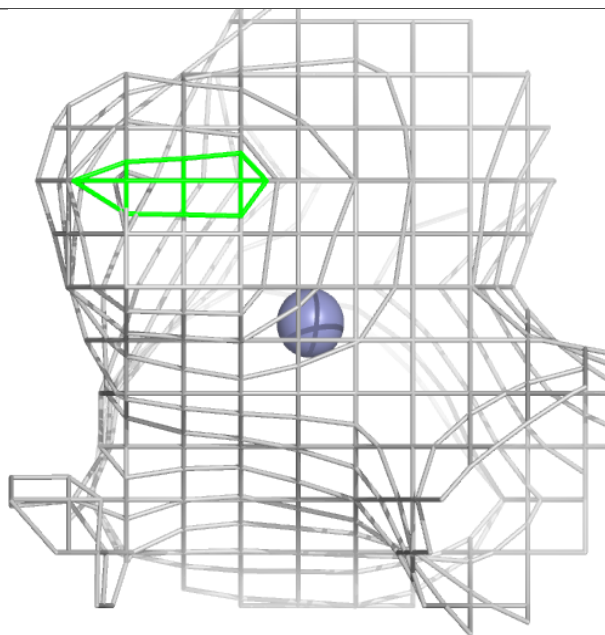
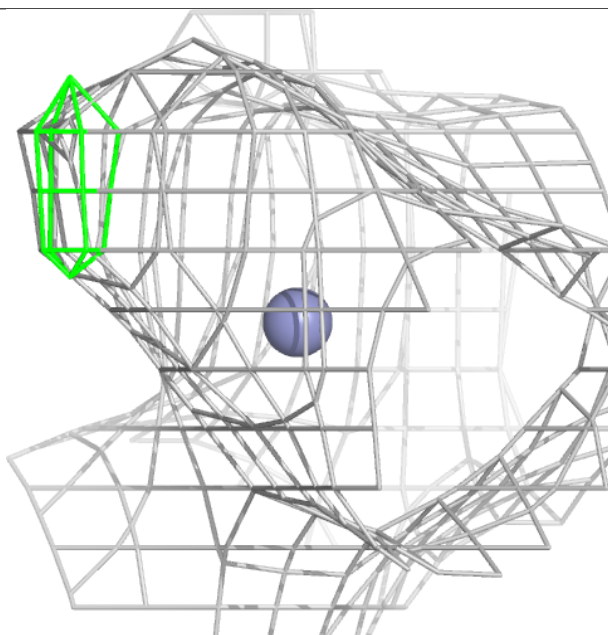
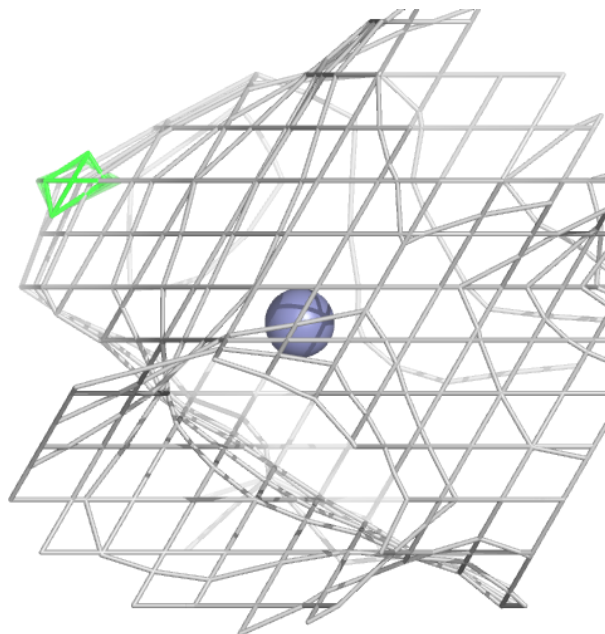
Electron density around ZN H 500:

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



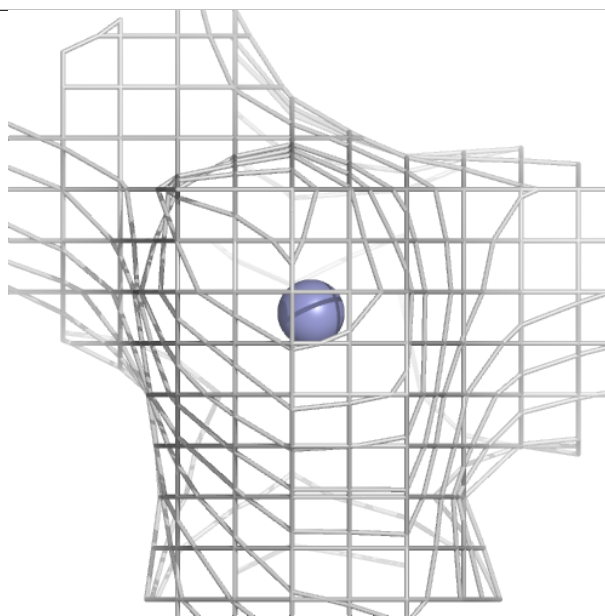
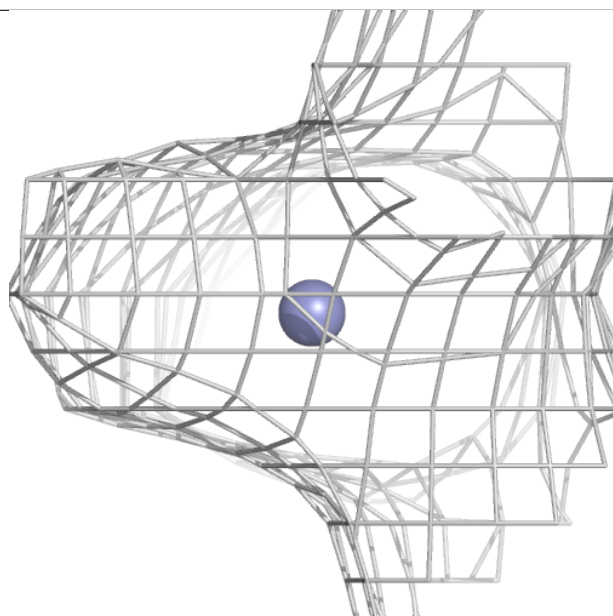
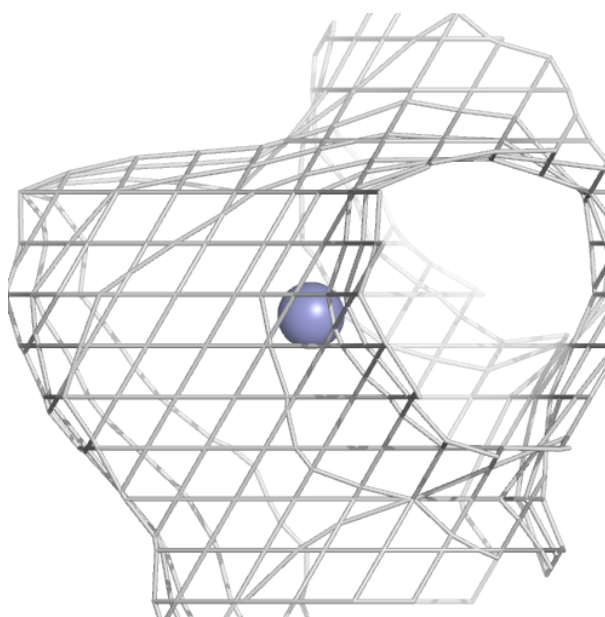
Electron density around ZN G 301:

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



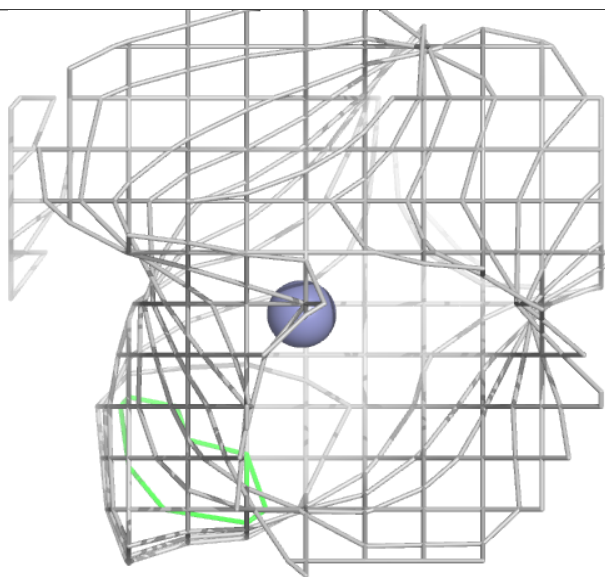
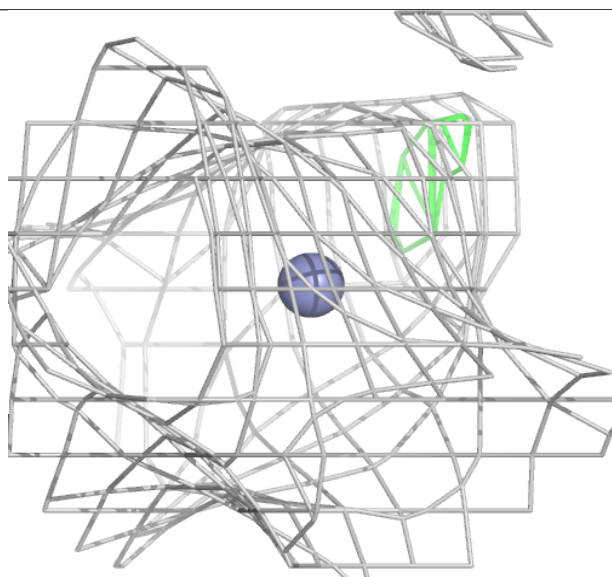
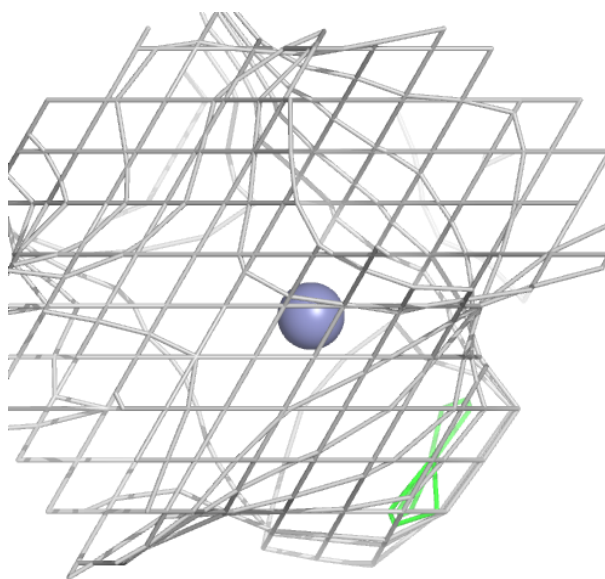
Electron density around ZN B 302:

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



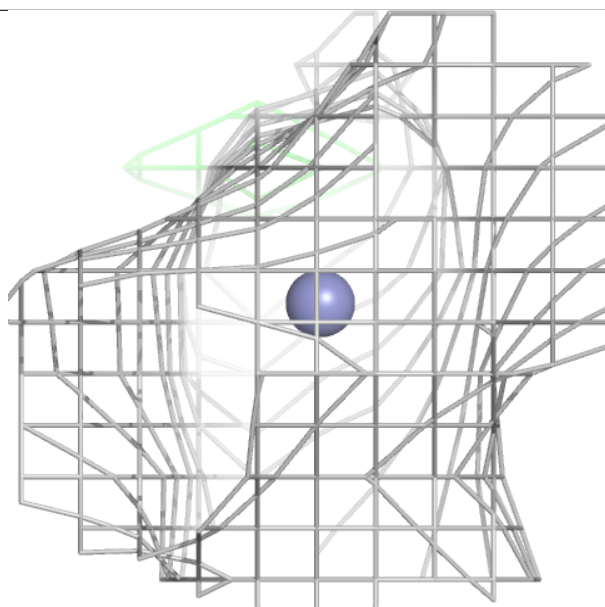
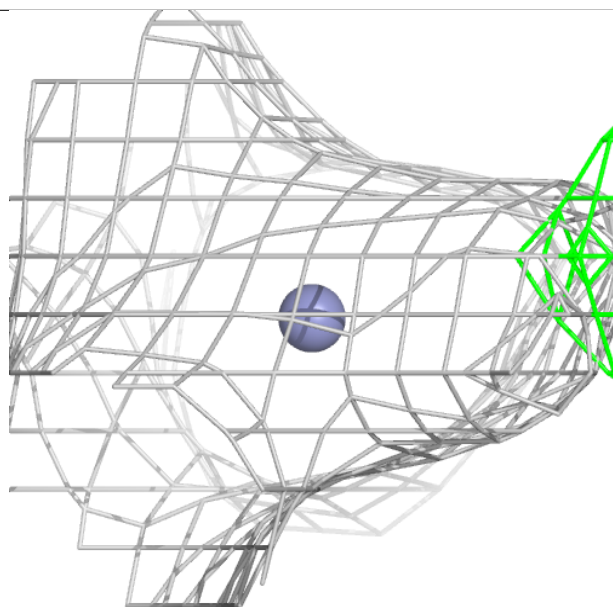
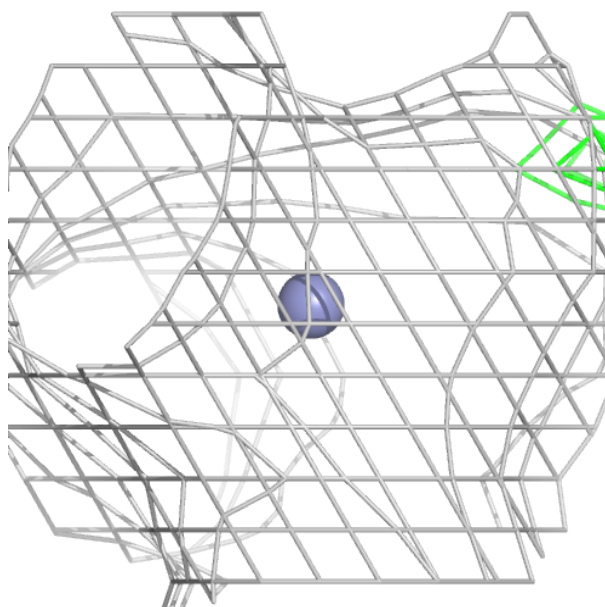
Electron density around ZN E 301:

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



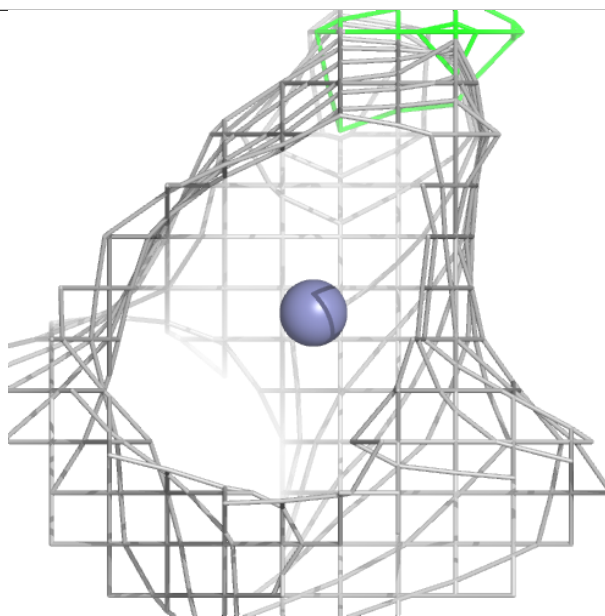
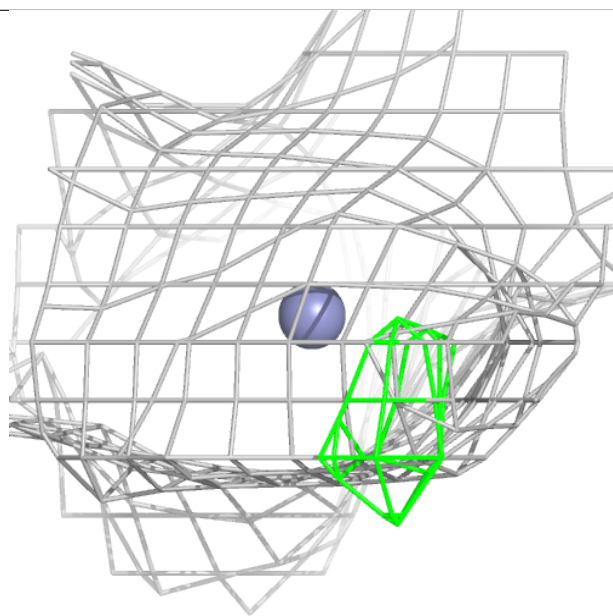
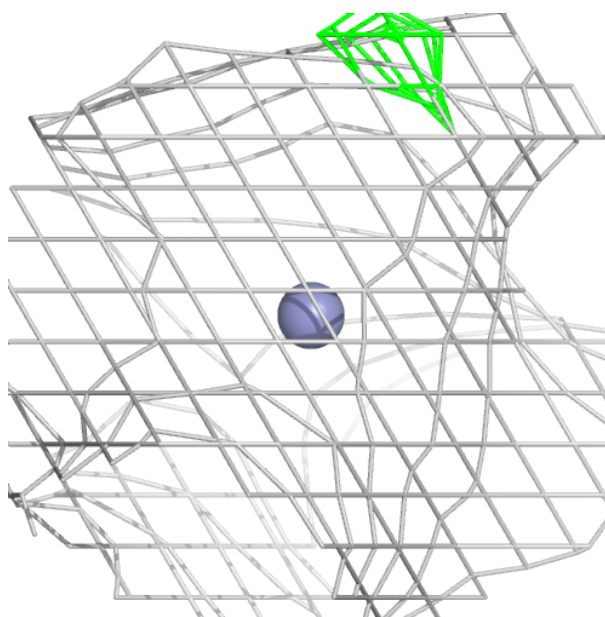
Electron density around ZN A 301:

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



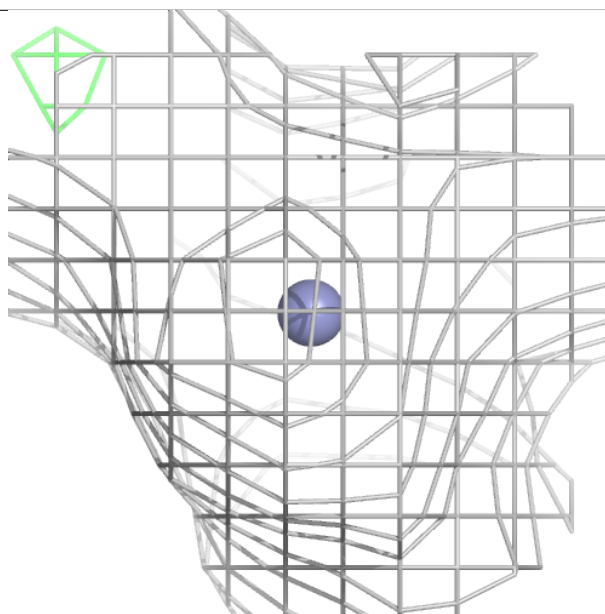
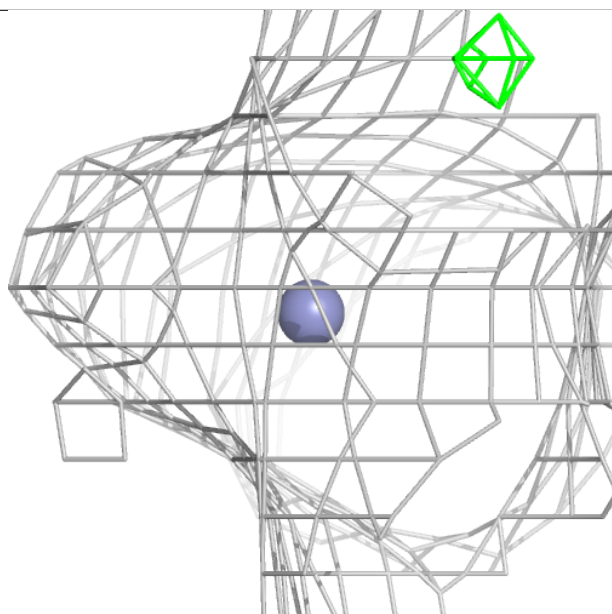
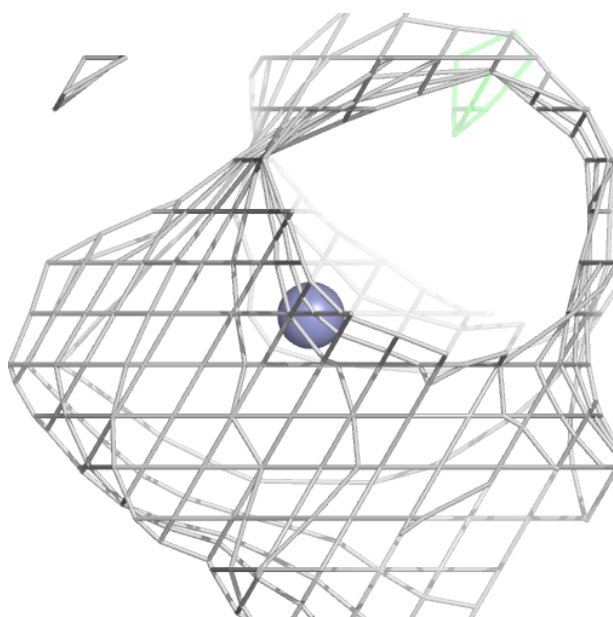
Electron density around ZN C 301:

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



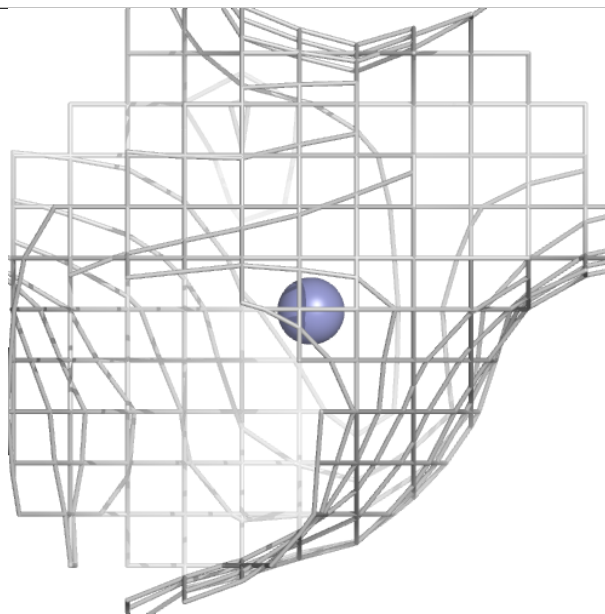
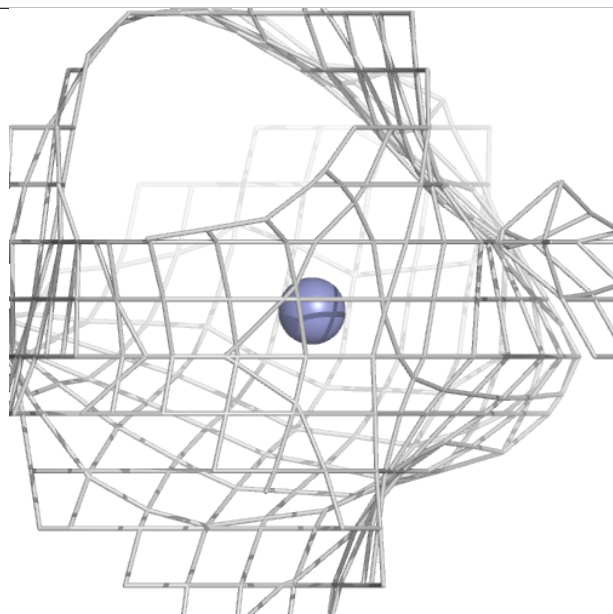
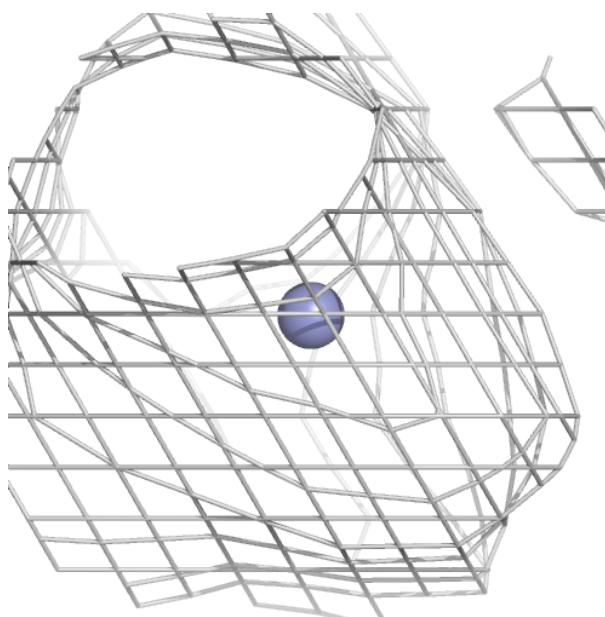
Electron density around ZN D 302:

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



Electron density around ZN F 302:

$2mF_o-DF_c$ (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.