



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 10:24 AM UTC

PDB ID : 6K1F / pdb_00006k1f
Title : Crystal structure of the L-fucose isomerase from Raoultella sp.
Authors : Kim, I.J.; Kim, D.H.; Nam, K.H.; Kim, K.H.
Deposited on : 2019-05-10
Resolution : 2.50 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

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

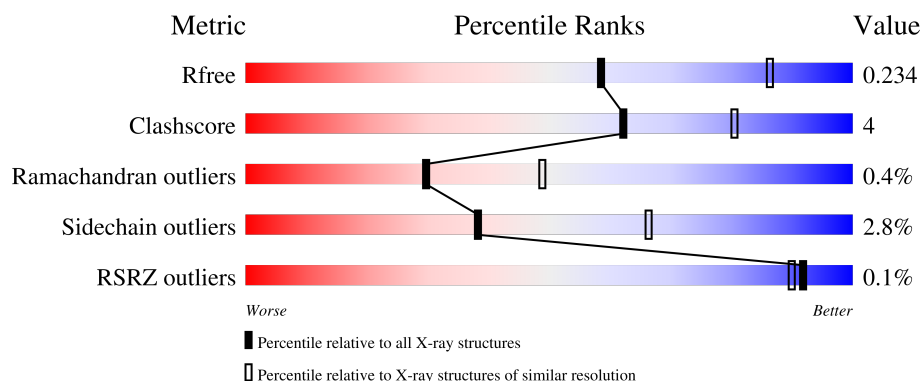
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	612	 85% 9% . .
1	B	612	 83% 12% . .
1	C	612	 87% 8% .
1	D	612	 87% 8% . .
1	E	612	 82% 13% . .

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Mol	Chain	Length	Quality of chain
1	F	612	87% 8% . .

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 28145 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 L-fucose isomerase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	587	Total	C	N	O	S	0	0	0
			4520	2832	794	859	35			
1	B	587	Total	C	N	O	S	0	0	0
			4520	2832	794	859	35			
1	C	587	Total	C	N	O	S	0	0	0
			4520	2832	794	859	35			
1	D	587	Total	C	N	O	S	0	0	0
			4520	2832	794	859	35			
1	E	587	Total	C	N	O	S	0	0	0
			4520	2832	794	859	35			
1	F	587	Total	C	N	O	S	0	0	0
			4520	2832	794	859	35			

There are 126 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-20	MET	-	initiating methionine	UNP A0A377T0E7
A	-19	GLY	-	expression tag	UNP A0A377T0E7
A	-18	SER	-	expression tag	UNP A0A377T0E7
A	-17	SER	-	expression tag	UNP A0A377T0E7
A	-16	HIS	-	expression tag	UNP A0A377T0E7
A	-15	HIS	-	expression tag	UNP A0A377T0E7
A	-14	HIS	-	expression tag	UNP A0A377T0E7
A	-13	HIS	-	expression tag	UNP A0A377T0E7
A	-12	HIS	-	expression tag	UNP A0A377T0E7
A	-11	HIS	-	expression tag	UNP A0A377T0E7
A	-10	SER	-	expression tag	UNP A0A377T0E7
A	-9	SER	-	expression tag	UNP A0A377T0E7
A	-8	GLY	-	expression tag	UNP A0A377T0E7
A	-7	LEU	-	expression tag	UNP A0A377T0E7
A	-6	VAL	-	expression tag	UNP A0A377T0E7
A	-5	PRO	-	expression tag	UNP A0A377T0E7
A	-4	ARG	-	expression tag	UNP A0A377T0E7

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-3	GLY	-	expression tag	UNP A0A377T0E7
A	-2	SER	-	expression tag	UNP A0A377T0E7
A	-1	HIS	-	expression tag	UNP A0A377T0E7
A	0	MET	-	expression tag	UNP A0A377T0E7
B	-20	MET	-	initiating methionine	UNP A0A377T0E7
B	-19	GLY	-	expression tag	UNP A0A377T0E7
B	-18	SER	-	expression tag	UNP A0A377T0E7
B	-17	SER	-	expression tag	UNP A0A377T0E7
B	-16	HIS	-	expression tag	UNP A0A377T0E7
B	-15	HIS	-	expression tag	UNP A0A377T0E7
B	-14	HIS	-	expression tag	UNP A0A377T0E7
B	-13	HIS	-	expression tag	UNP A0A377T0E7
B	-12	HIS	-	expression tag	UNP A0A377T0E7
B	-11	HIS	-	expression tag	UNP A0A377T0E7
B	-10	SER	-	expression tag	UNP A0A377T0E7
B	-9	SER	-	expression tag	UNP A0A377T0E7
B	-8	GLY	-	expression tag	UNP A0A377T0E7
B	-7	LEU	-	expression tag	UNP A0A377T0E7
B	-6	VAL	-	expression tag	UNP A0A377T0E7
B	-5	PRO	-	expression tag	UNP A0A377T0E7
B	-4	ARG	-	expression tag	UNP A0A377T0E7
B	-3	GLY	-	expression tag	UNP A0A377T0E7
B	-2	SER	-	expression tag	UNP A0A377T0E7
B	-1	HIS	-	expression tag	UNP A0A377T0E7
B	0	MET	-	expression tag	UNP A0A377T0E7
C	-20	MET	-	initiating methionine	UNP A0A377T0E7
C	-19	GLY	-	expression tag	UNP A0A377T0E7
C	-18	SER	-	expression tag	UNP A0A377T0E7
C	-17	SER	-	expression tag	UNP A0A377T0E7
C	-16	HIS	-	expression tag	UNP A0A377T0E7
C	-15	HIS	-	expression tag	UNP A0A377T0E7
C	-14	HIS	-	expression tag	UNP A0A377T0E7
C	-13	HIS	-	expression tag	UNP A0A377T0E7
C	-12	HIS	-	expression tag	UNP A0A377T0E7
C	-11	HIS	-	expression tag	UNP A0A377T0E7
C	-10	SER	-	expression tag	UNP A0A377T0E7
C	-9	SER	-	expression tag	UNP A0A377T0E7
C	-8	GLY	-	expression tag	UNP A0A377T0E7
C	-7	LEU	-	expression tag	UNP A0A377T0E7
C	-6	VAL	-	expression tag	UNP A0A377T0E7
C	-5	PRO	-	expression tag	UNP A0A377T0E7
C	-4	ARG	-	expression tag	UNP A0A377T0E7

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-3	GLY	-	expression tag	UNP A0A377T0E7
C	-2	SER	-	expression tag	UNP A0A377T0E7
C	-1	HIS	-	expression tag	UNP A0A377T0E7
C	0	MET	-	expression tag	UNP A0A377T0E7
D	-20	MET	-	initiating methionine	UNP A0A377T0E7
D	-19	GLY	-	expression tag	UNP A0A377T0E7
D	-18	SER	-	expression tag	UNP A0A377T0E7
D	-17	SER	-	expression tag	UNP A0A377T0E7
D	-16	HIS	-	expression tag	UNP A0A377T0E7
D	-15	HIS	-	expression tag	UNP A0A377T0E7
D	-14	HIS	-	expression tag	UNP A0A377T0E7
D	-13	HIS	-	expression tag	UNP A0A377T0E7
D	-12	HIS	-	expression tag	UNP A0A377T0E7
D	-11	HIS	-	expression tag	UNP A0A377T0E7
D	-10	SER	-	expression tag	UNP A0A377T0E7
D	-9	SER	-	expression tag	UNP A0A377T0E7
D	-8	GLY	-	expression tag	UNP A0A377T0E7
D	-7	LEU	-	expression tag	UNP A0A377T0E7
D	-6	VAL	-	expression tag	UNP A0A377T0E7
D	-5	PRO	-	expression tag	UNP A0A377T0E7
D	-4	ARG	-	expression tag	UNP A0A377T0E7
D	-3	GLY	-	expression tag	UNP A0A377T0E7
D	-2	SER	-	expression tag	UNP A0A377T0E7
D	-1	HIS	-	expression tag	UNP A0A377T0E7
D	0	MET	-	expression tag	UNP A0A377T0E7
E	-20	MET	-	initiating methionine	UNP A0A377T0E7
E	-19	GLY	-	expression tag	UNP A0A377T0E7
E	-18	SER	-	expression tag	UNP A0A377T0E7
E	-17	SER	-	expression tag	UNP A0A377T0E7
E	-16	HIS	-	expression tag	UNP A0A377T0E7
E	-15	HIS	-	expression tag	UNP A0A377T0E7
E	-14	HIS	-	expression tag	UNP A0A377T0E7
E	-13	HIS	-	expression tag	UNP A0A377T0E7
E	-12	HIS	-	expression tag	UNP A0A377T0E7
E	-11	HIS	-	expression tag	UNP A0A377T0E7
E	-10	SER	-	expression tag	UNP A0A377T0E7
E	-9	SER	-	expression tag	UNP A0A377T0E7
E	-8	GLY	-	expression tag	UNP A0A377T0E7
E	-7	LEU	-	expression tag	UNP A0A377T0E7
E	-6	VAL	-	expression tag	UNP A0A377T0E7
E	-5	PRO	-	expression tag	UNP A0A377T0E7
E	-4	ARG	-	expression tag	UNP A0A377T0E7

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Chain	Residue	Modelled	Actual	Comment	Reference
E	-3	GLY	-	expression tag	UNP A0A377T0E7
E	-2	SER	-	expression tag	UNP A0A377T0E7
E	-1	HIS	-	expression tag	UNP A0A377T0E7
E	0	MET	-	expression tag	UNP A0A377T0E7
F	-20	MET	-	initiating methionine	UNP A0A377T0E7
F	-19	GLY	-	expression tag	UNP A0A377T0E7
F	-18	SER	-	expression tag	UNP A0A377T0E7
F	-17	SER	-	expression tag	UNP A0A377T0E7
F	-16	HIS	-	expression tag	UNP A0A377T0E7
F	-15	HIS	-	expression tag	UNP A0A377T0E7
F	-14	HIS	-	expression tag	UNP A0A377T0E7
F	-13	HIS	-	expression tag	UNP A0A377T0E7
F	-12	HIS	-	expression tag	UNP A0A377T0E7
F	-11	HIS	-	expression tag	UNP A0A377T0E7
F	-10	SER	-	expression tag	UNP A0A377T0E7
F	-9	SER	-	expression tag	UNP A0A377T0E7
F	-8	GLY	-	expression tag	UNP A0A377T0E7
F	-7	LEU	-	expression tag	UNP A0A377T0E7
F	-6	VAL	-	expression tag	UNP A0A377T0E7
F	-5	PRO	-	expression tag	UNP A0A377T0E7
F	-4	ARG	-	expression tag	UNP A0A377T0E7
F	-3	GLY	-	expression tag	UNP A0A377T0E7
F	-2	SER	-	expression tag	UNP A0A377T0E7
F	-1	HIS	-	expression tag	UNP A0A377T0E7
F	0	MET	-	expression tag	UNP A0A377T0E7

- Molecule 2 is MANGANESE (II) ION (CCD ID: MN) (formula: Mn) (labeled as "Ligand of Interest" by depositor).

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

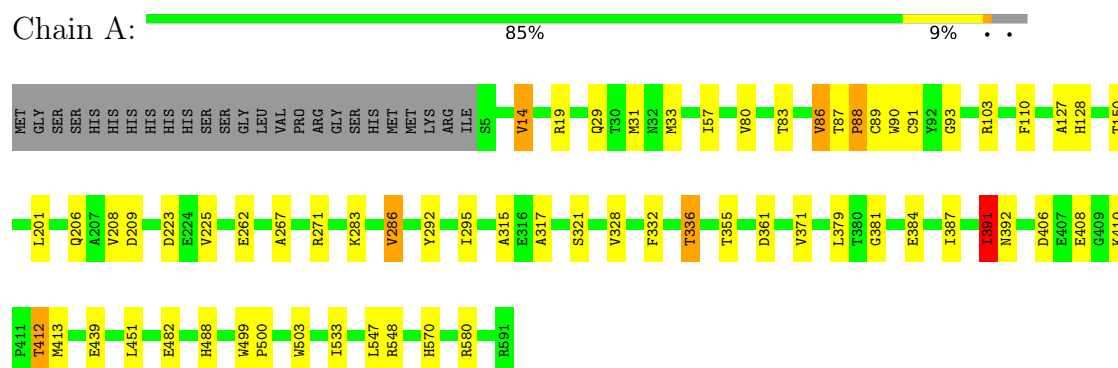
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	148	Total 148	O 148	0	0
3	B	167	Total 167	O 167	0	0
3	C	163	Total 163	O 163	0	0
3	D	216	Total 216	O 216	0	0
3	E	139	Total 139	O 139	0	0
3	F	186	Total 186	O 186	0	0

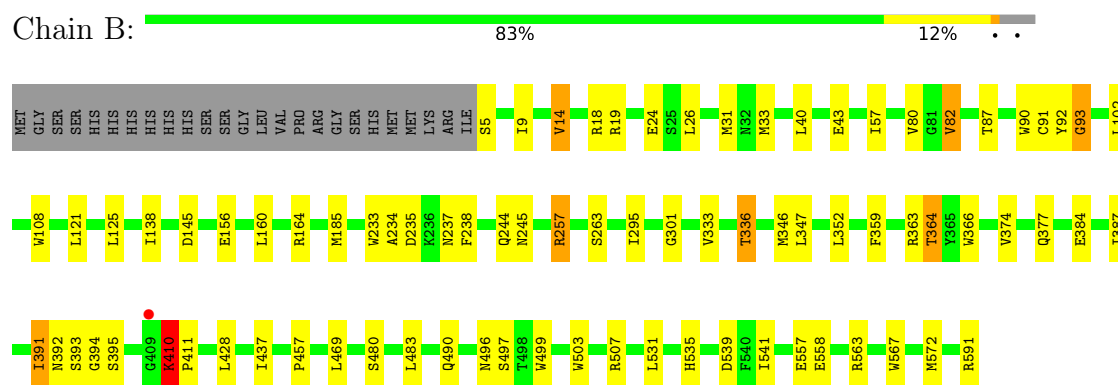
3 Residue-property plots

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.

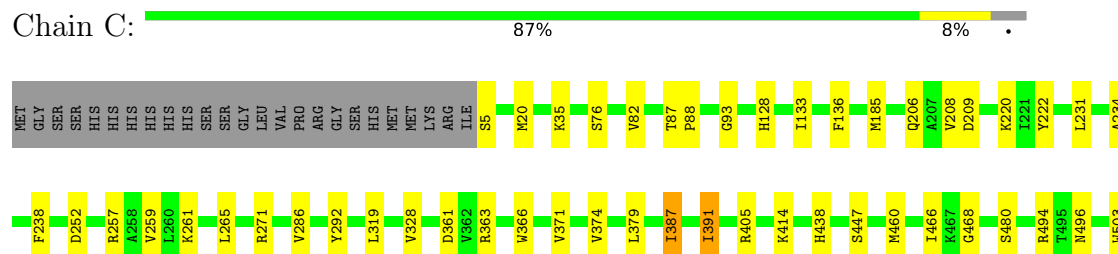
• Molecule 1: L-fucose isomerase



• Molecule 1: L-fucose isomerase



• Molecule 1: L-fucose isomerase





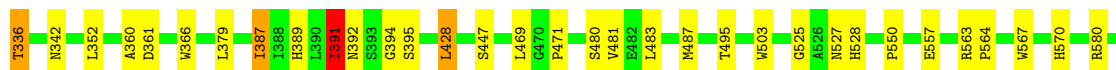
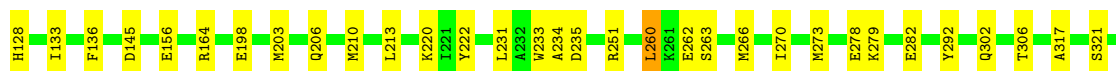
- Molecule 1: L-fucose isomerase

Chain D: 87% 8% . .



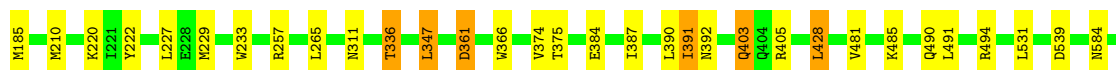
- Molecule 1: L-fucose isomerase

Chain E: 82% 13% . .



- Molecule 1: L-fucose isomerase

Chain F: 87% 8% . .



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	113.99Å 127.61Å 257.30Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.01 – 2.50 50.01 – 2.50	Depositor EDS
% Data completeness (in resolution range)	96.3 (50.01-2.50) 96.3 (50.01-2.50)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.04 (at 2.51Å)	Xtriage
Refinement program	REFMAC 5.8.0218	Depositor
R, R_{free}	0.177 , 0.233 0.183 , 0.234	Depositor DCC
R_{free} test set	6337 reflections (4.87%)	wwPDB-VP
Wilson B-factor (Å ²)	32.5	Xtriage
Anisotropy	0.021	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 24.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	28145	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.90% 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: MN

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	0.78	1/4621 (0.0%)	0.90	4/6261 (0.1%)
1	B	0.81	1/4621 (0.0%)	0.96	5/6261 (0.1%)
1	C	0.77	1/4621 (0.0%)	0.91	2/6261 (0.0%)
1	D	0.81	0/4621	0.91	2/6261 (0.0%)
1	E	0.78	2/4621 (0.0%)	0.92	6/6261 (0.1%)
1	F	0.77	0/4621	0.92	3/6261 (0.0%)
All	All	0.79	5/27726 (0.0%)	0.92	22/37566 (0.1%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	86	VAL	C-O	-7.80	1.16	1.24
1	E	495	THR	CA-C	6.23	1.56	1.52
1	E	395	SER	C-O	-5.97	1.16	1.23
1	C	466	ILE	C-O	-5.47	1.18	1.24
1	B	395	SER	C-O	-5.35	1.17	1.23

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	410	LYS	N-CA-C	11.43	122.75	109.60
1	B	410	LYS	CA-C-N	9.77	129.53	119.76
1	B	410	LYS	C-N-CA	9.77	129.53	119.76
1	B	394	GLY	N-CA-C	8.02	124.47	114.37
1	E	87	THR	N-CA-C	7.90	115.33	108.22
1	A	87	THR	N-CA-C	6.93	114.46	108.22
1	E	87	THR	CA-C-N	6.58	128.07	119.84
1	E	87	THR	C-N-CA	6.58	128.07	119.84
1	A	87	THR	CA-C-N	5.82	127.11	119.84
1	A	87	THR	C-N-CA	5.82	127.11	119.84

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	391	ILE	CB-CA-C	-5.63	104.03	110.73
1	F	87	THR	N-CA-C	5.58	115.30	108.11
1	E	342	ASN	N-CA-C	-5.56	105.22	111.28
1	B	384	GLU	N-CA-C	5.41	118.68	111.75
1	F	53	ILE	CB-CA-C	-5.36	103.35	111.33
1	F	590	LYS	N-CA-C	5.22	117.01	108.76
1	C	87	THR	N-CA-C	5.18	114.79	108.11
1	A	89	CYS	N-CA-C	5.17	115.87	108.74
1	E	394	GLY	N-CA-C	5.16	121.63	114.92
1	D	361	ASP	N-CA-C	-5.10	101.65	109.76
1	D	80	VAL	N-CA-CB	5.03	116.01	110.53
1	C	438	HIS	N-CA-C	5.00	117.38	111.33

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	4520	0	4385	38	0
1	B	4520	0	4385	47	0
1	C	4520	0	4385	33	0
1	D	4520	0	4385	34	0
1	E	4520	0	4385	51	0
1	F	4520	0	4385	43	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
3	A	148	0	0	3	0
3	B	167	0	0	5	0
3	C	163	0	0	2	0
3	D	216	0	0	1	0
3	E	139	0	0	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	F	186	0	0	2	0
All	All	28145	0	26310	217	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 4.

All (217) 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:E:366:TRP:HB2	1:E:387:ILE:HG22	1.51	0.92
1:C:468:GLY:HA2	1:F:591:ARG:NH2	1.90	0.86
3:B:724:HOH:O	1:F:79:ASN:HB2	1.77	0.83
1:B:145:ASP:OD1	1:C:494:ARG:NH2	2.14	0.79
1:C:468:GLY:CA	1:F:591:ARG:NH2	2.47	0.78
1:C:468:GLY:HA2	1:F:591:ARG:HH22	1.49	0.77
1:F:374:VAL:O	1:F:490:GLN:NE2	2.17	0.77
1:B:387:ILE:HD11	1:B:503:TRP:HB3	1.66	0.77
1:F:366:TRP:HB2	1:F:387:ILE:HG22	1.69	0.75
1:C:468:GLY:CA	1:F:591:ARG:HH22	2.03	0.69
1:A:14:VAL:HG12	1:A:91:CYS:SG	2.32	0.69
1:C:405:ARG:O	1:C:414:LYS:NZ	2.25	0.69
1:F:336:THR:HG21	1:F:392:ASN:HD21	1.58	0.69
1:A:570:HIS:O	1:A:580:ARG:NH1	2.26	0.69
1:B:90:TRP:HE1	1:C:185:MET:HE2	1.61	0.66
1:E:90:TRP:HE1	1:F:185:MET:HE2	1.60	0.65
1:C:468:GLY:N	1:F:591:ARG:HH22	1.94	0.64
1:E:145:ASP:OD1	1:F:494:ARG:NH2	2.31	0.64
1:C:387:ILE:HD11	1:C:503:TRP:HB3	1.80	0.64
1:A:317:ALA:O	1:A:321:SER:OG	2.15	0.63
1:E:14:VAL:HG12	1:E:87:THR:HB	1.78	0.63
1:F:387:ILE:HD11	1:F:481:VAL:HG21	1.80	0.63
1:E:317:ALA:O	1:E:321:SER:OG	2.15	0.63
1:C:468:GLY:N	1:F:591:ARG:NH2	2.47	0.62
1:D:381:GLY:O	1:D:384:GLU:HB2	1.99	0.62
1:E:282:GLU:HB3	3:E:813:HOH:O	1.99	0.62
1:A:371:VAL:HG11	1:A:379:LEU:HD13	1.82	0.62
1:D:14:VAL:HG12	1:D:91:CYS:SG	2.39	0.62
1:A:90:TRP:HE1	1:B:185:MET:HE2	1.65	0.61
1:E:483:LEU:HD22	1:E:487:MET:HE2	1.81	0.61
1:A:31:MET:HE1	1:A:57:ILE:HG21	1.82	0.61
1:E:389:HIS:CE1	1:E:391:ILE:HD11	2.36	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:31:MET:HE1	1:B:57:ILE:CG2	2.30	0.61
1:E:570:HIS:O	1:E:580:ARG:NH1	2.32	0.60
1:A:271:ARG:HG3	1:A:332:PHE:CE2	2.36	0.60
1:A:223:ASP:OD1	1:A:225:VAL:HG12	2.02	0.60
1:C:222:TYR:CE2	1:C:265:LEU:HD21	2.36	0.60
1:E:379:LEU:HD23	1:E:487:MET:HE1	1.83	0.59
1:E:251:ARG:NH1	3:E:702:HOH:O	2.34	0.59
1:F:233:TRP:CZ2	1:F:428:LEU:HD11	2.38	0.58
1:E:336:THR:HG21	1:E:392:ASN:HD21	1.69	0.58
1:C:222:TYR:CD2	1:C:265:LEU:HD21	2.39	0.57
1:A:336:THR:HG21	1:A:392:ASN:HD21	1.70	0.57
1:B:31:MET:HE1	1:B:57:ILE:HG21	1.84	0.57
1:C:231:LEU:HD11	1:C:261:LYS:HG3	1.85	0.57
1:B:336:THR:HG21	1:B:392:ASN:HD21	1.71	0.56
1:E:113:THR:OG1	1:F:494:ARG:NH1	2.37	0.56
1:A:90:TRP:HE1	1:B:185:MET:CE	2.19	0.56
1:B:391:ILE:HD11	1:B:499:TRP:HB2	1.89	0.55
1:D:200:TRP:CE2	1:D:587:PRO:HG3	2.42	0.55
1:E:164:ARG:HD3	1:E:352:LEU:O	2.07	0.55
1:F:387:ILE:HD11	1:F:481:VAL:CG2	2.37	0.55
1:D:200:TRP:CZ2	1:D:587:PRO:HG3	2.42	0.54
1:F:374:VAL:HG23	1:F:375:THR:HG23	1.89	0.54
1:E:40:LEU:HD21	1:E:156:GLU:HG3	1.88	0.54
1:E:387:ILE:HD11	1:E:481:VAL:HG21	1.90	0.54
1:D:563:ARG:HD3	1:D:567:TRP:CD1	2.42	0.54
1:F:14:VAL:HG21	1:F:96:THR:HB	1.89	0.54
1:F:403:GLN:HE21	1:F:405:ARG:HH12	1.57	0.53
1:F:227:LEU:HD23	3:F:842:HOH:O	2.09	0.53
1:F:366:TRP:CZ2	1:F:491:LEU:HD22	2.44	0.53
1:D:407:GLU:HB2	3:D:803:HOH:O	2.09	0.53
1:D:494:ARG:NH1	1:F:113:THR:OG1	2.42	0.52
1:E:14:VAL:CG1	1:E:87:THR:HB	2.39	0.52
1:D:270:ILE:HA	1:D:273:MET:HE3	1.90	0.52
1:F:46:ARG:NH1	3:F:707:HOH:O	2.42	0.52
1:F:361:ASP:HB2	1:F:391:ILE:O	2.10	0.51
1:E:115:ARG:NH1	3:E:707:HOH:O	2.44	0.51
1:D:366:TRP:HB2	1:D:387:ILE:HG22	1.93	0.51
1:E:361:ASP:HB2	1:E:391:ILE:O	2.10	0.51
1:E:206:GLN:HG2	1:E:292:TYR:CD1	2.46	0.51
1:D:231:LEU:HD23	1:D:260:LEU:HD12	1.93	0.51
1:E:220:LYS:HA	1:E:222:TYR:CE2	2.46	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:361:ASP:OD1	1:E:528:HIS:CD2	2.63	0.51
1:C:460:MET:HA	1:C:530:VAL:O	2.10	0.51
1:D:103:ARG:O	1:D:105:LYS:HG3	2.11	0.51
1:B:121:LEU:CD2	1:B:125:LEU:HD11	2.41	0.50
1:A:355:THR:HG21	1:A:533:ILE:HG21	1.93	0.50
1:E:90:TRP:HE1	1:F:185:MET:CE	2.24	0.50
1:F:53:ILE:HD12	1:F:167:LEU:HD13	1.92	0.50
1:B:352:LEU:HD13	1:B:539:ASP:HB3	1.93	0.50
1:A:83:THR:HG23	1:A:103:ARG:CD	2.42	0.50
1:A:499:TRP:HB3	1:A:500:PRO:HD2	1.94	0.49
1:C:271:ARG:HB2	1:C:319:LEU:HD21	1.93	0.49
1:E:483:LEU:HD22	1:E:487:MET:CE	2.42	0.49
1:D:486:ALA:O	1:D:490:GLN:HG3	2.13	0.49
1:A:387:ILE:HD12	1:A:503:TRP:CE3	2.47	0.49
1:C:286:VAL:HG23	1:E:19:ARG:NE	2.26	0.49
1:A:83:THR:HG23	1:A:103:ARG:HD2	1.93	0.49
1:B:19:ARG:NH1	3:B:704:HOH:O	2.43	0.49
1:B:572:MET:HE2	3:C:761:HOH:O	2.11	0.49
1:B:233:TRP:CZ2	1:B:428:LEU:HD11	2.48	0.48
1:B:14:VAL:HG12	1:B:91:CYS:SG	2.52	0.48
1:C:128:HIS:HD1	1:C:133:ILE:HG13	1.78	0.48
1:E:80:VAL:HG23	1:E:80:VAL:O	2.14	0.48
1:C:234:ALA:O	1:C:238:PHE:HB2	2.13	0.48
1:B:160:LEU:O	1:B:164:ARG:HB2	2.13	0.48
1:F:14:VAL:CG2	1:F:96:THR:HB	2.44	0.48
1:E:527:ASN:HD21	1:E:528:HIS:CE1	2.31	0.47
1:B:374:VAL:O	1:B:490:GLN:NE2	2.43	0.47
1:B:235:ASP:OD1	1:B:257:ARG:NH1	2.39	0.47
1:A:381:GLY:O	1:A:384:GLU:HB2	2.13	0.47
1:B:33:MET:HE1	1:B:87:THR:O	2.14	0.47
1:B:563:ARG:HD3	1:B:567:TRP:CD1	2.49	0.47
1:A:451:LEU:HD22	1:A:488:HIS:CE1	2.50	0.47
1:B:92:TYR:O	1:B:93:GLY:C	2.58	0.47
1:B:591:ARG:HG3	1:B:591:ARG:OXT	2.14	0.47
1:B:393:SER:HA	1:B:499:TRP:CD2	2.50	0.47
1:A:286:VAL:HG23	1:D:19:ARG:CD	2.45	0.46
1:B:366:TRP:HB2	1:B:387:ILE:CG2	2.45	0.46
1:D:235:ASP:CG	1:D:257:ARG:HH12	2.23	0.46
1:E:233:TRP:CZ2	1:E:428:LEU:HD11	2.50	0.46
1:E:360:ALA:HA	1:E:392:ASN:HB2	1.95	0.46
1:A:127:ALA:O	1:A:128:HIS:C	2.55	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:206:GLN:HG2	1:C:292:TYR:CD1	2.50	0.46
1:A:33:MET:HE2	1:A:86:VAL:HB	1.96	0.46
1:E:563:ARG:HD3	1:E:567:TRP:CD1	2.51	0.46
1:B:234:ALA:O	1:B:238:PHE:HB2	2.16	0.46
1:B:295:ILE:HD12	1:B:347:LEU:HD11	1.98	0.46
1:E:14:VAL:HG21	1:E:96:THR:HB	1.97	0.45
1:B:333:VAL:HG13	1:B:346:MET:HE2	1.98	0.45
1:D:311:ASN:C	1:D:311:ASN:OD1	2.59	0.45
1:E:480:SER:HA	1:E:503:TRP:O	2.15	0.45
1:F:83:THR:HG23	1:F:103:ARG:CD	2.47	0.45
1:A:286:VAL:HG23	1:D:19:ARG:NE	2.31	0.45
1:F:220:LYS:HA	1:F:222:TYR:CE1	2.50	0.45
1:A:31:MET:HE1	1:A:57:ILE:CG2	2.46	0.45
1:A:283:LYS:O	1:A:283:LYS:HG2	2.17	0.45
1:B:393:SER:HA	1:B:499:TRP:CE3	2.50	0.45
1:C:371:VAL:HG11	1:C:379:LEU:HG	1.98	0.45
1:F:8:LYS:HD2	1:F:78:GLN:O	2.17	0.45
1:E:270:ILE:HA	1:E:273:MET:HE3	1.98	0.45
1:D:165:ALA:O	1:D:168:ALA:HB3	2.17	0.45
1:E:80:VAL:O	1:E:80:VAL:CG2	2.65	0.45
1:D:14:VAL:CG1	1:D:91:CYS:SG	3.05	0.44
1:F:164:ARG:NH1	1:F:539:ASP:OD1	2.50	0.44
1:B:14:VAL:CG1	1:B:91:CYS:SG	3.05	0.44
1:E:302:GLN:HG2	1:E:306:THR:OG1	2.18	0.44
1:B:245:ASN:OD1	1:B:437:ILE:HA	2.17	0.44
1:E:231:LEU:HD23	1:E:260:LEU:CD1	2.48	0.44
1:B:90:TRP:HE1	1:C:185:MET:CE	2.29	0.44
1:B:507:ARG:HH21	1:B:535:HIS:CE1	2.35	0.44
1:F:311:ASN:OD1	1:F:311:ASN:C	2.61	0.44
1:D:75:PHE:HB3	1:D:80:VAL:CG2	2.47	0.44
1:B:363:ARG:HG3	1:B:364:THR:OG1	2.17	0.44
1:E:87:THR:HA	1:E:88:PRO:HD3	1.74	0.44
1:B:457:PRO:HB2	1:B:535:HIS:CD2	2.53	0.43
1:C:133:ILE:HG13	1:C:133:ILE:O	2.18	0.43
1:C:286:VAL:CG2	1:E:19:ARG:NE	2.81	0.43
1:F:53:ILE:HD12	1:F:167:LEU:CD1	2.48	0.43
1:B:480:SER:HA	1:B:503:TRP:O	2.18	0.43
1:C:366:TRP:HB2	1:C:387:ILE:HG23	2.00	0.43
1:D:233:TRP:CZ2	1:D:428:LEU:HD11	2.53	0.43
1:E:91:CYS:HB2	1:E:120:TYR:CD1	2.53	0.43
1:E:198:GLU:HA	1:E:203:MET:O	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:361:ASP:HB2	1:A:391:ILE:O	2.19	0.43
1:F:336:THR:HG21	1:F:392:ASN:ND2	2.31	0.43
1:A:547:LEU:C	1:A:548:ARG:HD2	2.43	0.43
1:E:83:THR:HG23	1:E:103:ARG:CD	2.49	0.43
1:E:234:ALA:O	1:E:235:ASP:C	2.62	0.43
1:B:237:ASN:HB3	1:B:428:LEU:HD13	2.00	0.43
1:C:468:GLY:H	1:F:591:ARG:HH22	1.64	0.43
1:E:213:LEU:HD21	1:E:266:MET:HG2	2.01	0.43
1:F:366:TRP:HB2	1:F:387:ILE:CG2	2.46	0.43
1:B:18:ARG:NH1	3:B:714:HOH:O	2.52	0.43
1:B:244:GLN:NE2	1:B:497:SER:OG	2.43	0.43
1:D:126:ALA:HB2	1:E:525:GLY:O	2.19	0.43
1:E:128:HIS:ND1	1:E:133:ILE:HG13	2.33	0.43
1:A:206:GLN:HG2	1:A:292:TYR:CD1	2.54	0.43
1:A:209:ASP:OD2	1:E:210:MET:HB2	2.19	0.43
1:B:108:TRP:CD1	1:B:138:ILE:HG22	2.54	0.43
1:A:271:ARG:HG3	1:A:332:PHE:CZ	2.54	0.42
1:D:460:MET:HA	1:D:530:VAL:O	2.19	0.42
1:F:390:LEU:CD1	1:F:531:LEU:HD13	2.49	0.42
1:D:185:MET:HE2	1:F:92:TYR:CE1	2.54	0.42
1:C:209:ASP:OD2	1:F:210:MET:HB2	2.19	0.42
1:E:136:PHE:CE1	1:E:564:PRO:HG3	2.55	0.42
1:F:265:LEU:HD23	1:F:265:LEU:HA	1.90	0.42
1:E:389:HIS:NE2	1:E:391:ILE:HD11	2.35	0.42
1:B:410:LYS:HA	1:B:411:PRO:HD3	1.77	0.42
1:A:19:ARG:NE	1:D:286:VAL:HG23	2.34	0.42
1:C:136:PHE:CE1	1:C:546:MET:HE2	2.54	0.42
1:F:347:LEU:HD23	1:F:347:LEU:HA	1.91	0.42
1:B:387:ILE:HD12	1:B:483:LEU:HD21	2.02	0.42
1:C:366:TRP:HB3	1:C:371:VAL:HG23	2.02	0.41
1:A:29:GLN:HG3	3:A:749:HOH:O	2.19	0.41
1:C:252:ASP:OD1	1:C:252:ASP:C	2.62	0.41
1:D:361:ASP:HB2	1:D:391:ILE:O	2.20	0.41
1:B:359:PHE:O	1:B:392:ASN:HB2	2.20	0.41
1:C:363:ARG:HD3	3:C:802:HOH:O	2.19	0.41
1:A:19:ARG:NH2	3:A:724:HOH:O	2.54	0.41
1:A:412:THR:HG23	1:A:413:MET:N	2.35	0.41
1:D:100:ASP:O	1:D:105:LYS:CE	2.69	0.41
1:D:92:TYR:O	1:D:93:GLY:C	2.63	0.41
1:D:185:MET:CE	1:F:18:ARG:HH22	2.33	0.41
1:D:371:VAL:HG21	1:D:379:LEU:HD11	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:406:ASP:OD1	1:A:410:LYS:N	2.54	0.41
1:E:278:GLU:O	1:E:279:LYS:C	2.64	0.41
1:B:24:GLU:HG3	3:B:704:HOH:O	2.20	0.41
1:A:439:GLU:HA	1:C:20:MET:HE2	2.03	0.41
1:A:548:ARG:HD2	1:A:548:ARG:N	2.36	0.41
1:B:557:GLU:O	1:B:558:GLU:C	2.62	0.41
1:D:206:GLN:HG2	1:D:292:TYR:CD1	2.56	0.41
1:A:482:GLU:N	3:A:712:HOH:O	2.47	0.41
1:B:9:ILE:HG12	1:B:82:VAL:HG13	2.03	0.41
1:E:471:PRO:O	1:E:550:PRO:HD2	2.21	0.41
1:B:40:LEU:HD21	1:B:156:GLU:HG3	2.03	0.40
1:C:563:ARG:NH2	1:C:574:THR:O	2.54	0.40
1:D:358:VAL:HB	1:D:531:LEU:HB2	2.03	0.40
1:A:267:ALA:HB2	1:A:315:ALA:HA	2.03	0.40
1:D:336:THR:HG21	1:D:392:ASN:HD21	1.85	0.40
1:D:468:GLY:HA3	1:F:584:ASN:OD1	2.21	0.40
1:E:103:ARG:O	1:E:105:LYS:NZ	2.52	0.40
1:A:88:PRO:HA	1:A:110:PHE:HB3	2.02	0.40
1:C:480:SER:HA	1:C:503:TRP:O	2.21	0.40
1:D:193:ASP:C	1:D:193:ASP:OD1	2.64	0.40
1:B:377:GLN:NE2	3:B:718:HOH:O	2.54	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	585/612 (96%)	562 (96%)	20 (3%)	3 (0%)	24	43
1	B	585/612 (96%)	548 (94%)	35 (6%)	2 (0%)	36	55
1	C	585/612 (96%)	563 (96%)	19 (3%)	3 (0%)	24	43
1	D	585/612 (96%)	565 (97%)	17 (3%)	3 (0%)	24	43

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	E	585/612 (96%)	564 (96%)	19 (3%)	2 (0%)	36	55
1	F	585/612 (96%)	563 (96%)	21 (4%)	1 (0%)	43	63
All	All	3510/3672 (96%)	3365 (96%)	131 (4%)	14 (0%)	30	49

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	93	GLY
1	B	93	GLY
1	C	93	GLY
1	C	391	ILE
1	F	391	ILE
1	E	88	PRO
1	A	88	PRO
1	C	88	PRO
1	A	391	ILE
1	D	93	GLY
1	D	391	ILE
1	B	301	GLY
1	D	88	PRO
1	E	116	PRO

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	467/489 (96%)	454 (97%)	13 (3%)	38	66
1	B	467/489 (96%)	450 (96%)	17 (4%)	31	58
1	C	467/489 (96%)	451 (97%)	16 (3%)	32	60
1	D	467/489 (96%)	461 (99%)	6 (1%)	61	82
1	E	467/489 (96%)	454 (97%)	13 (3%)	38	66
1	F	467/489 (96%)	454 (97%)	13 (3%)	38	66
All	All	2802/2934 (96%)	2724 (97%)	78 (3%)	38	66

All (78) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	14	VAL
1	A	80	VAL
1	A	150	THR
1	A	201	LEU
1	A	208	VAL
1	A	262	GLU
1	A	286	VAL
1	A	295	ILE
1	A	328	VAL
1	A	336	THR
1	A	391	ILE
1	A	408	GLU
1	A	412	THR
1	B	5	SER
1	B	14	VAL
1	B	26	LEU
1	B	43	GLU
1	B	80	VAL
1	B	82	VAL
1	B	102	LEU
1	B	257	ARG
1	B	263	SER
1	B	336	THR
1	B	364	THR
1	B	391	ILE
1	B	410	LYS
1	B	469	LEU
1	B	496	ASN
1	B	531	LEU
1	B	541	ILE
1	C	5	SER
1	C	35	LYS
1	C	76	SER
1	C	82	VAL
1	C	208	VAL
1	C	220	LYS
1	C	257	ARG
1	C	259	VAL
1	C	328	VAL
1	C	361	ASP
1	C	374	VAL
1	C	387	ILE

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Mol	Chain	Res	Type
1	C	391	ILE
1	C	447	SER
1	C	496	ASN
1	C	563	ARG
1	D	14	VAL
1	D	286	VAL
1	D	336	THR
1	D	374	VAL
1	D	391	ILE
1	D	481	VAL
1	E	26	LEU
1	E	82	VAL
1	E	121	LEU
1	E	260	LEU
1	E	262	GLU
1	E	263	SER
1	E	336	THR
1	E	387	ILE
1	E	391	ILE
1	E	428	LEU
1	E	447	SER
1	E	469	LEU
1	E	557	GLU
1	F	5	SER
1	F	26	LEU
1	F	80	VAL
1	F	82	VAL
1	F	229	MET
1	F	257	ARG
1	F	336	THR
1	F	347	LEU
1	F	361	ASP
1	F	384	GLU
1	F	403	GLN
1	F	428	LEU
1	F	485	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (34) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	29	GLN
1	A	144	GLN

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Mol	Chain	Res	Type
1	A	237	ASN
1	A	351	GLN
1	B	144	GLN
1	B	255	GLN
1	B	351	GLN
1	B	392	ASN
1	B	523	ASN
1	B	577	GLN
1	C	29	GLN
1	C	206	GLN
1	C	237	ASN
1	C	308	GLN
1	C	351	GLN
1	C	385	HIS
1	D	302	GLN
1	D	351	GLN
1	D	377	GLN
1	D	385	HIS
1	D	421	GLN
1	E	78	GLN
1	E	194	HIS
1	E	219	GLN
1	E	351	GLN
1	E	438	HIS
1	E	490	GLN
1	E	527	ASN
1	F	144	GLN
1	F	206	GLN
1	F	392	ASN
1	F	403	GLN
1	F	523	ASN
1	F	577	GLN

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 6 ligands modelled in this entry, 6 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	587/612 (95%)	-0.42	0 100 100	21, 36, 57, 80	0
1	B	587/612 (95%)	-0.45	1 (0%) 91 89	19, 35, 55, 93	0
1	C	587/612 (95%)	-0.52	0 100 100	21, 35, 53, 73	0
1	D	587/612 (95%)	-0.67	0 100 100	18, 30, 44, 56	0
1	E	587/612 (95%)	-0.47	1 (0%) 91 89	22, 37, 55, 78	0
1	F	587/612 (95%)	-0.59	1 (0%) 91 89	17, 32, 52, 80	0
All	All	3522/3672 (95%)	-0.52	3 (0%) 92 90	17, 34, 53, 93	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	591	ARG	3.3
1	B	409	GLY	2.3
1	F	591	ARG	2.2

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,

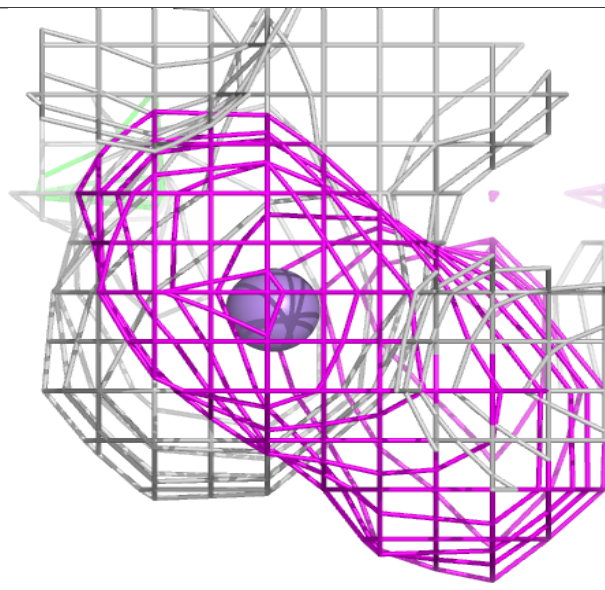
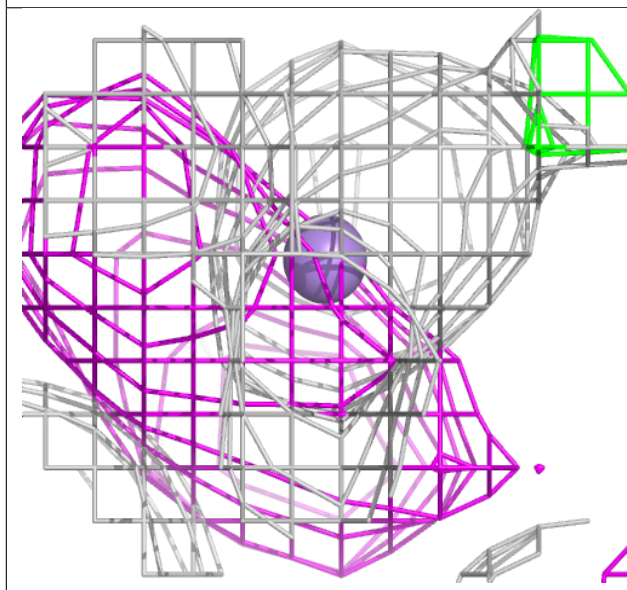
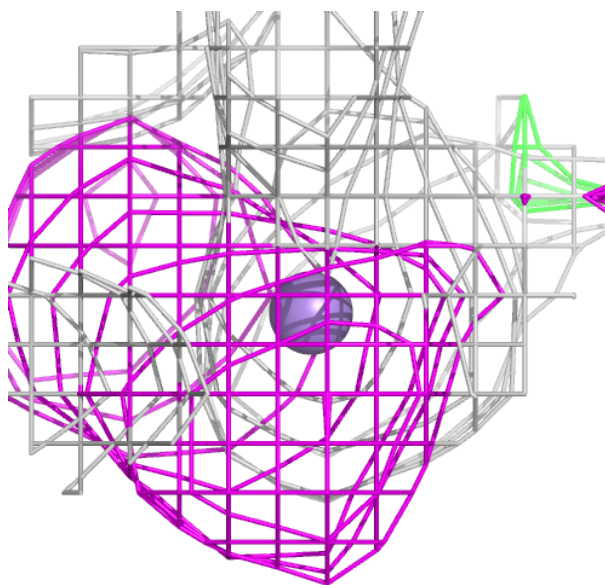
median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	MN	B	601	1/1	0.61	0.21	72,72,72,72	0
2	MN	F	601	1/1	0.79	0.14	74,74,74,74	0
2	MN	D	601	1/1	0.80	0.12	66,66,66,66	0
2	MN	E	601	1/1	0.81	0.14	75,75,75,75	0
2	MN	C	601	1/1	0.92	0.09	67,67,67,67	0
2	MN	A	601	1/1	0.93	0.08	68,68,68,68	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.

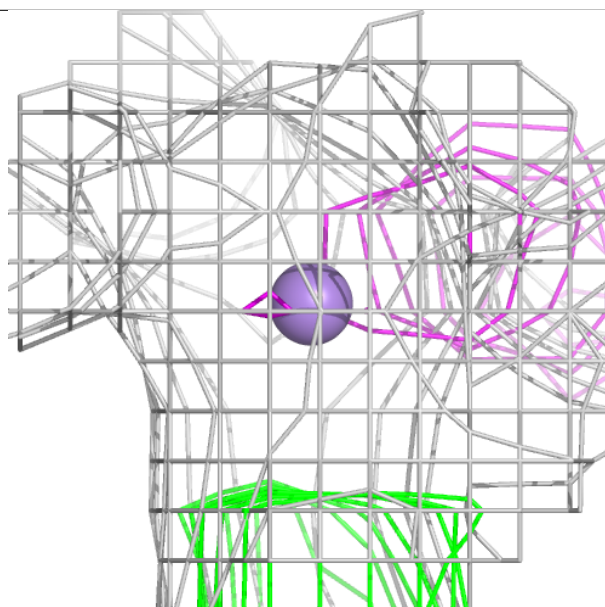
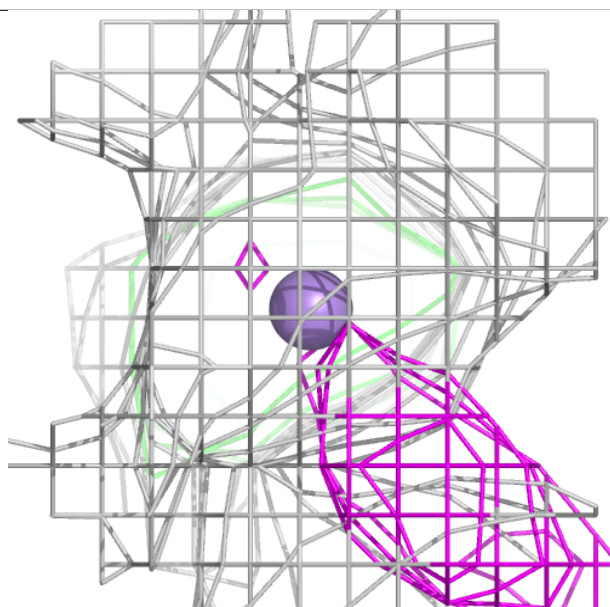
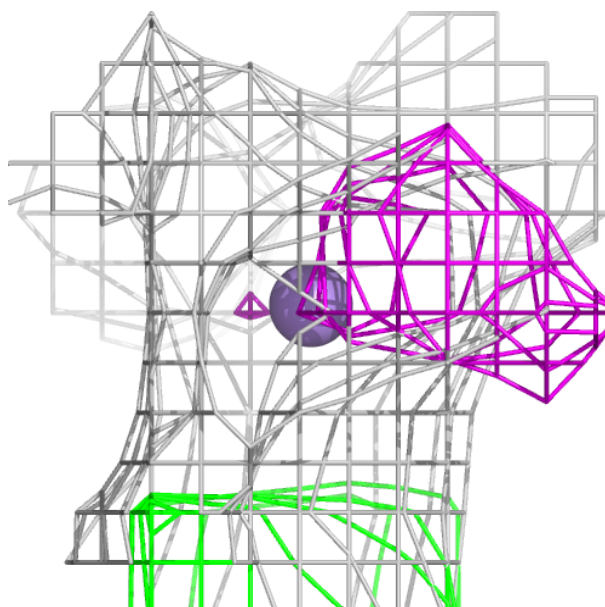
Electron density around MN B 601:

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



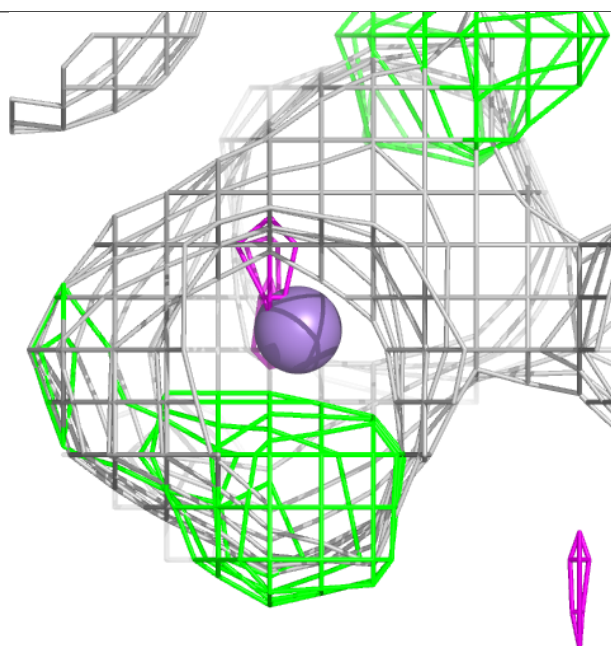
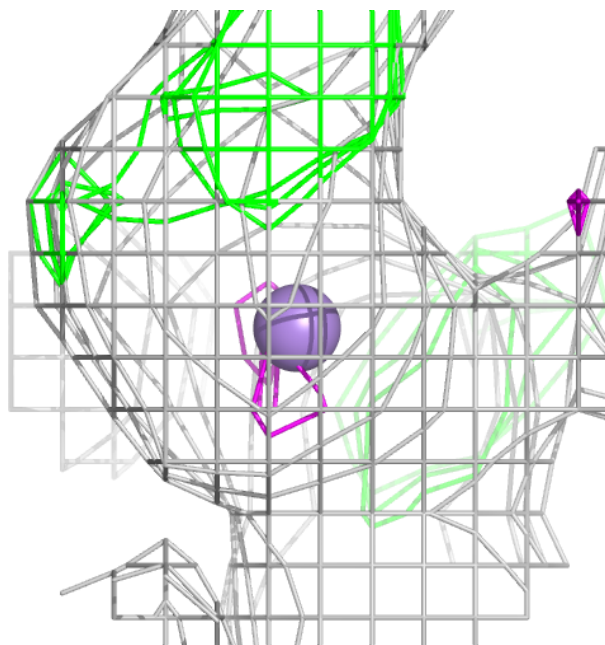
Electron density around MN F 601:

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



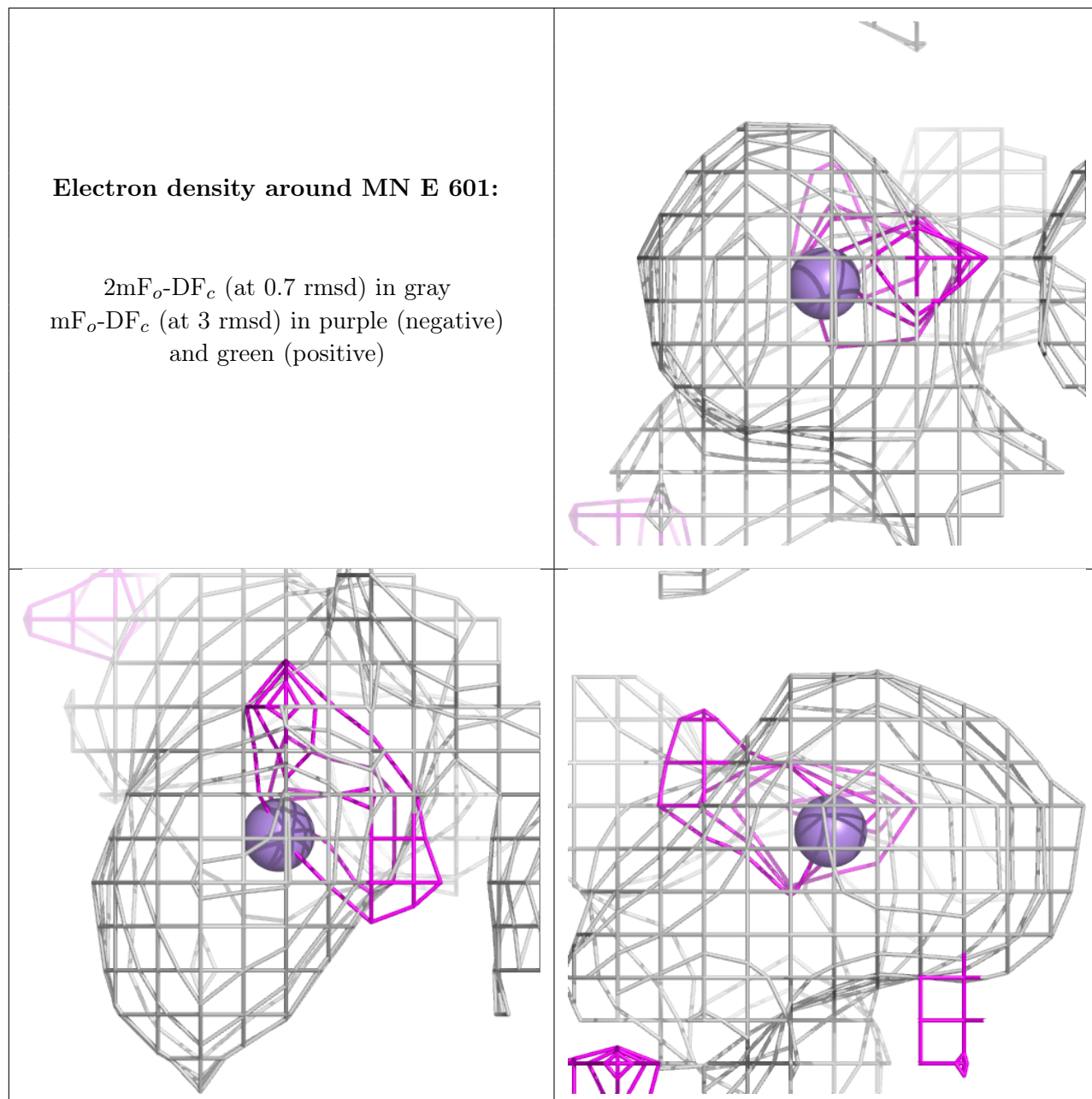
Electron density around MN D 601:

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



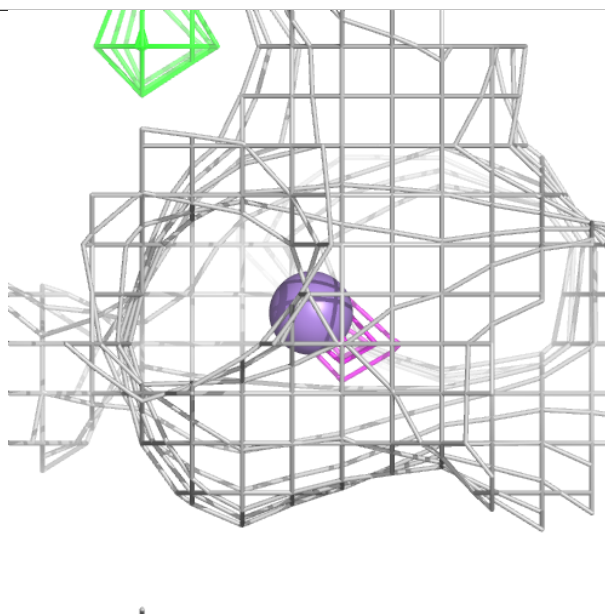
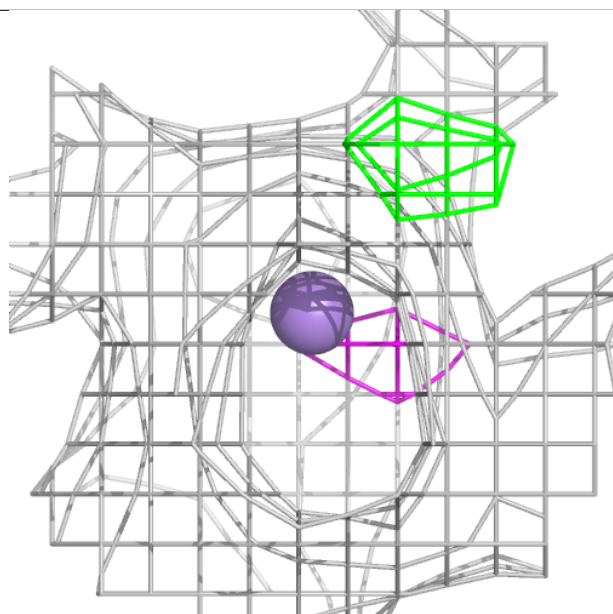
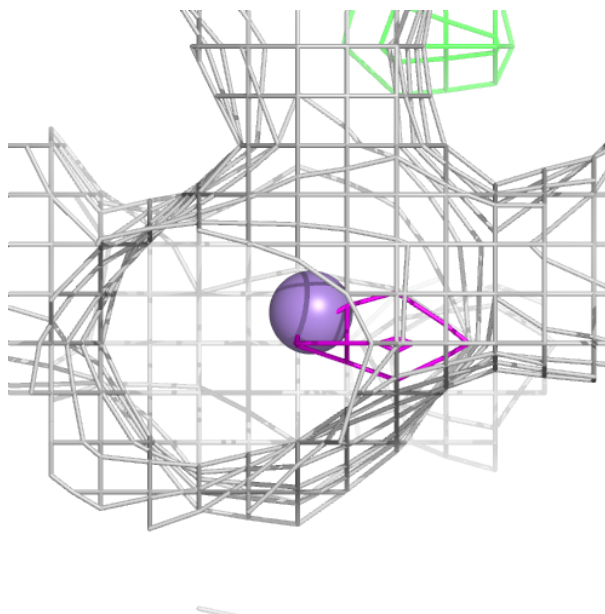
Electron density around MN E 601:

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



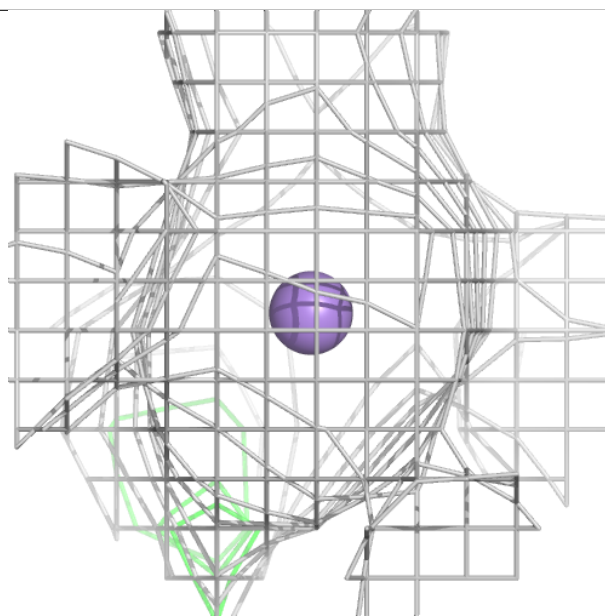
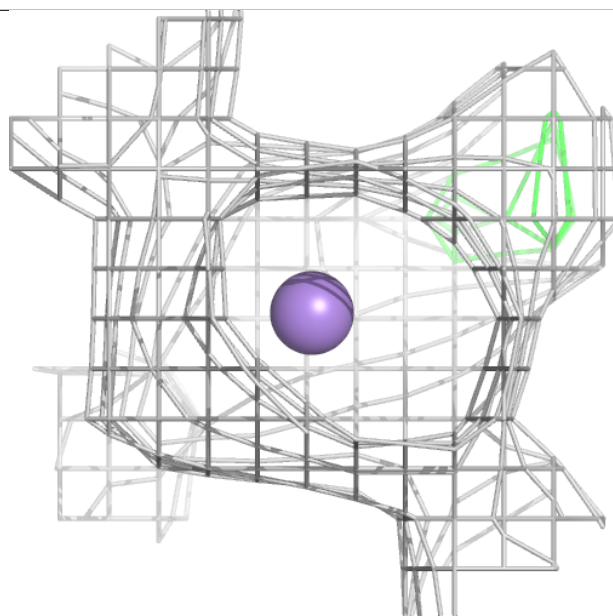
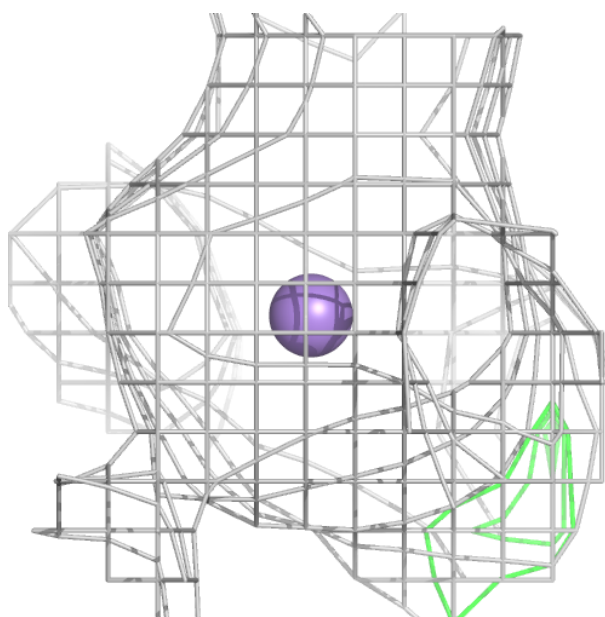
Electron density around MN C 601:

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



Electron density around MN A 601:

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



6.5 Other polymers ⓘ

There are no such residues in this entry.