



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 8, 2020 – 01:44 AM BST

PDB ID : 3CYQ
Title : The crystal structure of the complex of the C-terminal domain of Helicobacter pylori MotB (residues 125-256) with N-acetylmuramic acid
Authors : Roujeinikova, A.
Deposited on : 2008-04-26
Resolution : 2.30 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.13.1
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.13.1

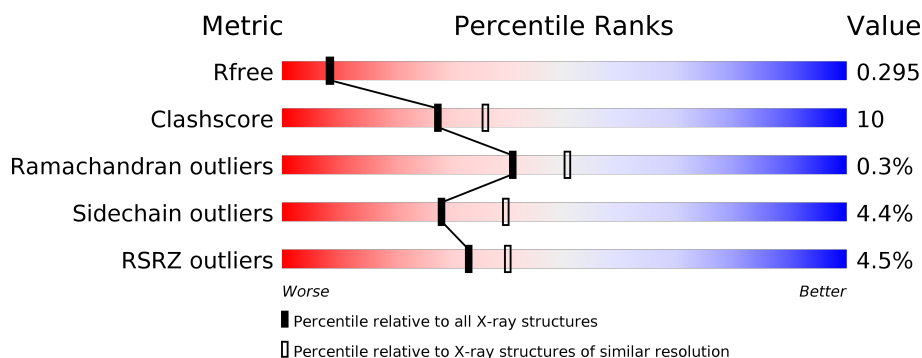
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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}	130704	5042 (2.30-2.30)
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (2.30-2.30)

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 on 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	138	<div> <div>5%</div> <div> <div></div> <div>74%</div> <div>21%</div> <div>• •</div> </div> </div>
1	B	138	<div> <div>4%</div> <div> <div></div> <div>75%</div> <div>21%</div> <div>•</div> </div> </div>
1	C	138	<div> <div>4%</div> <div> <div></div> <div>74%</div> <div>20%</div> <div>• •</div> </div> </div>
1	D	138	<div> <div>%</div> <div> <div></div> <div>75%</div> <div>18%</div> <div>• •</div> </div> </div>
1	E	138	<div> <div>4%</div> <div> <div></div> <div>65%</div> <div>30%</div> <div>• •</div> </div> </div>
1	F	138	<div> <div>3%</div> <div> <div></div> <div>83%</div> <div>12%</div> <div>• •</div> </div> </div>

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Mol	Chain	Length	Quality of chain
1	G	138	
1	H	138	
1	I	138	
1	J	138	
1	K	138	
1	L	138	
1	M	138	
1	N	138	
1	O	138	
1	P	138	

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 18460 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 Chemotaxis protein motB.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	133	Total	C	N	O	S	0	2	0
			1090	688	193	205	4			
1	C	133	Total	C	N	O	S	0	2	0
			1088	687	191	206	4			
1	D	132	Total	C	N	O	S	0	0	0
			1064	671	187	202	4			
1	E	132	Total	C	N	O	S	0	0	0
			1064	671	187	202	4			
1	F	133	Total	C	N	O	S	0	0	0
			1071	676	188	203	4			
1	G	134	Total	C	N	O	S	0	0	0
			1081	682	191	204	4			
1	H	132	Total	C	N	O	S	0	0	0
			1064	671	187	202	4			
1	I	132	Total	C	N	O	S	0	0	0
			1064	671	187	202	4			
1	A	133	Total	C	N	O	S	0	0	0
			1071	676	188	203	4			
1	J	134	Total	C	N	O	S	0	0	0
			1081	682	191	204	4			
1	K	132	Total	C	N	O	S	0	0	0
			1064	671	187	202	4			
1	L	132	Total	C	N	O	S	0	0	0
			1064	671	187	202	4			
1	M	133	Total	C	N	O	S	0	0	0
			1071	676	188	203	4			
1	N	134	Total	C	N	O	S	0	0	0
			1081	682	191	204	4			
1	O	132	Total	C	N	O	S	0	0	0
			1064	671	187	202	4			
1	P	132	Total	C	N	O	S	0	0	0
			1064	671	187	202	4			

There are 96 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	119	GLY	-	expression tag	UNP P56427
B	120	ILE	-	expression tag	UNP P56427
B	121	ASP	-	expression tag	UNP P56427
B	122	PRO	-	expression tag	UNP P56427
B	123	PHE	-	expression tag	UNP P56427
B	124	THR	-	expression tag	UNP P56427
C	119	GLY	-	expression tag	UNP P56427
C	120	ILE	-	expression tag	UNP P56427
C	121	ASP	-	expression tag	UNP P56427
C	122	PRO	-	expression tag	UNP P56427
C	123	PHE	-	expression tag	UNP P56427
C	124	THR	-	expression tag	UNP P56427
D	119	GLY	-	expression tag	UNP P56427
D	120	ILE	-	expression tag	UNP P56427
D	121	ASP	-	expression tag	UNP P56427
D	122	PRO	-	expression tag	UNP P56427
D	123	PHE	-	expression tag	UNP P56427
D	124	THR	-	expression tag	UNP P56427
E	119	GLY	-	expression tag	UNP P56427
E	120	ILE	-	expression tag	UNP P56427
E	121	ASP	-	expression tag	UNP P56427
E	122	PRO	-	expression tag	UNP P56427
E	123	PHE	-	expression tag	UNP P56427
E	124	THR	-	expression tag	UNP P56427
F	119	GLY	-	expression tag	UNP P56427
F	120	ILE	-	expression tag	UNP P56427
F	121	ASP	-	expression tag	UNP P56427
F	122	PRO	-	expression tag	UNP P56427
F	123	PHE	-	expression tag	UNP P56427
F	124	THR	-	expression tag	UNP P56427
G	119	GLY	-	expression tag	UNP P56427
G	120	ILE	-	expression tag	UNP P56427
G	121	ASP	-	expression tag	UNP P56427
G	122	PRO	-	expression tag	UNP P56427
G	123	PHE	-	expression tag	UNP P56427
G	124	THR	-	expression tag	UNP P56427
H	119	GLY	-	expression tag	UNP P56427
H	120	ILE	-	expression tag	UNP P56427
H	121	ASP	-	expression tag	UNP P56427
H	122	PRO	-	expression tag	UNP P56427
H	123	PHE	-	expression tag	UNP P56427
H	124	THR	-	expression tag	UNP P56427

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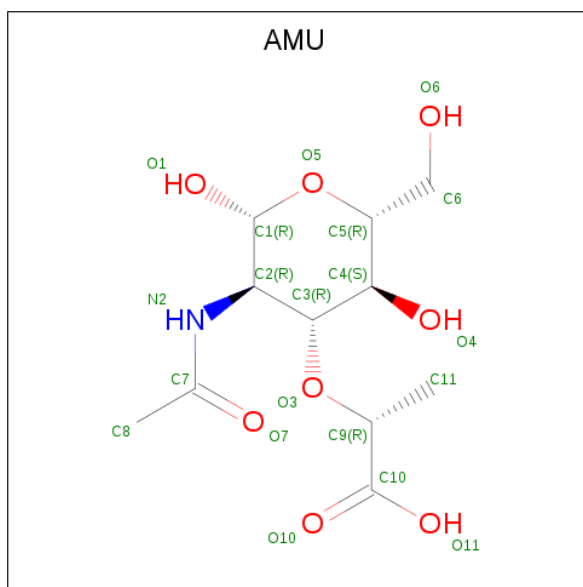
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I	119	GLY	-	expression tag	UNP P56427
I	120	ILE	-	expression tag	UNP P56427
I	121	ASP	-	expression tag	UNP P56427
I	122	PRO	-	expression tag	UNP P56427
I	123	PHE	-	expression tag	UNP P56427
I	124	THR	-	expression tag	UNP P56427
A	119	GLY	-	expression tag	UNP P56427
A	120	ILE	-	expression tag	UNP P56427
A	121	ASP	-	expression tag	UNP P56427
A	122	PRO	-	expression tag	UNP P56427
A	123	PHE	-	expression tag	UNP P56427
A	124	THR	-	expression tag	UNP P56427
J	119	GLY	-	expression tag	UNP P56427
J	120	ILE	-	expression tag	UNP P56427
J	121	ASP	-	expression tag	UNP P56427
J	122	PRO	-	expression tag	UNP P56427
J	123	PHE	-	expression tag	UNP P56427
J	124	THR	-	expression tag	UNP P56427
K	119	GLY	-	expression tag	UNP P56427
K	120	ILE	-	expression tag	UNP P56427
K	121	ASP	-	expression tag	UNP P56427
K	122	PRO	-	expression tag	UNP P56427
K	123	PHE	-	expression tag	UNP P56427
K	124	THR	-	expression tag	UNP P56427
L	119	GLY	-	expression tag	UNP P56427
L	120	ILE	-	expression tag	UNP P56427
L	121	ASP	-	expression tag	UNP P56427
L	122	PRO	-	expression tag	UNP P56427
L	123	PHE	-	expression tag	UNP P56427
L	124	THR	-	expression tag	UNP P56427
M	119	GLY	-	expression tag	UNP P56427
M	120	ILE	-	expression tag	UNP P56427
M	121	ASP	-	expression tag	UNP P56427
M	122	PRO	-	expression tag	UNP P56427
M	123	PHE	-	expression tag	UNP P56427
M	124	THR	-	expression tag	UNP P56427
N	119	GLY	-	expression tag	UNP P56427
N	120	ILE	-	expression tag	UNP P56427
N	121	ASP	-	expression tag	UNP P56427
N	122	PRO	-	expression tag	UNP P56427
N	123	PHE	-	expression tag	UNP P56427
N	124	THR	-	expression tag	UNP P56427

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Chain	Residue	Modelled	Actual	Comment	Reference
O	119	GLY	-	expression tag	UNP P56427
O	120	ILE	-	expression tag	UNP P56427
O	121	ASP	-	expression tag	UNP P56427
O	122	PRO	-	expression tag	UNP P56427
O	123	PHE	-	expression tag	UNP P56427
O	124	THR	-	expression tag	UNP P56427
P	119	GLY	-	expression tag	UNP P56427
P	120	ILE	-	expression tag	UNP P56427
P	121	ASP	-	expression tag	UNP P56427
P	122	PRO	-	expression tag	UNP P56427
P	123	PHE	-	expression tag	UNP P56427
P	124	THR	-	expression tag	UNP P56427

- Molecule 2 is N-acetyl-beta-muramic acid (three-letter code: AMU) (formula: $C_{11}H_{19}NO_8$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	D	1	Total	C	N	O	0	0
			20	11	1	8		
2	K	1	Total	C	N	O	0	0
			20	11	1	8		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	89	Total	O	0	0
			89	89		

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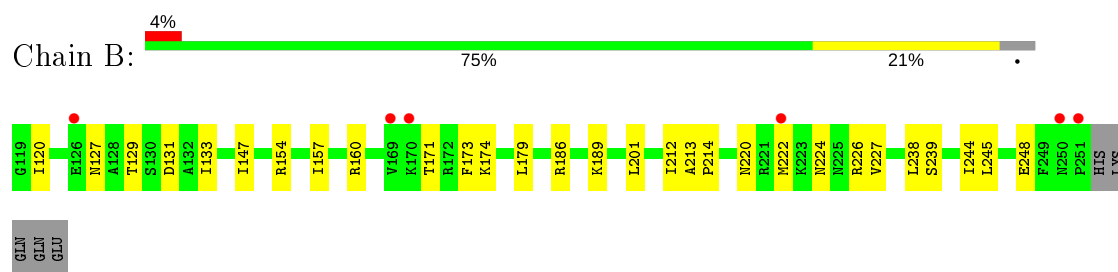
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	C	70	Total	O	0	0
			70	70		
3	D	99	Total	O	0	0
			99	99		
3	E	78	Total	O	0	0
			78	78		
3	F	89	Total	O	0	0
			89	89		
3	G	85	Total	O	0	0
			85	85		
3	H	70	Total	O	0	0
			70	70		
3	I	59	Total	O	0	0
			59	59		
3	A	68	Total	O	0	0
			68	68		
3	J	76	Total	O	0	0
			76	76		
3	K	96	Total	O	0	0
			96	96		
3	L	83	Total	O	0	0
			83	83		
3	M	92	Total	O	0	0
			92	92		
3	N	77	Total	O	0	0
			77	77		
3	O	74	Total	O	0	0
			74	74		
3	P	69	Total	O	0	0
			69	69		

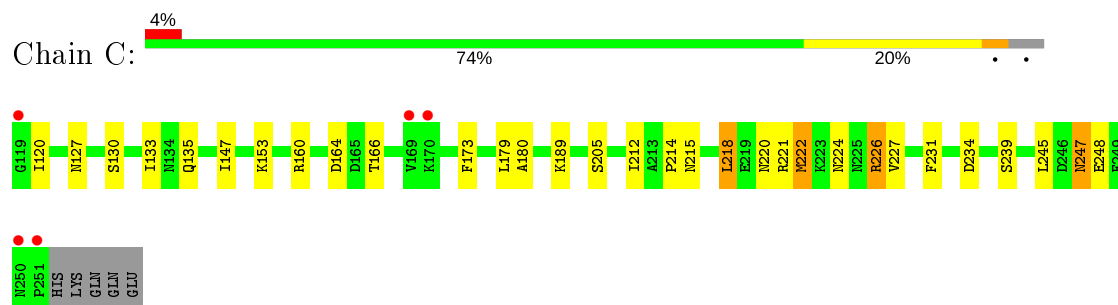
3 Residue-property plots [i](#)

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.

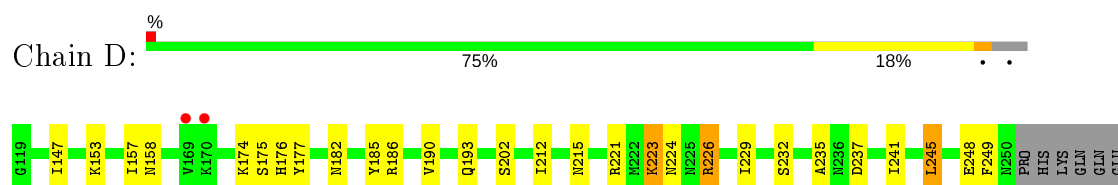
- Molecule 1: Chemotaxis protein motB



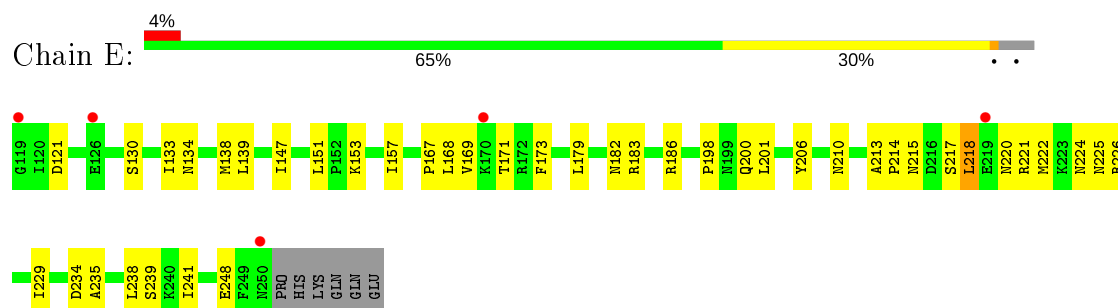
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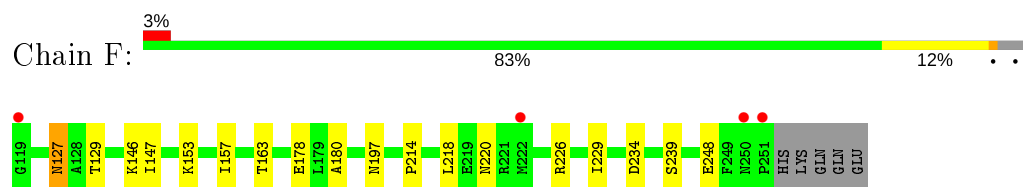
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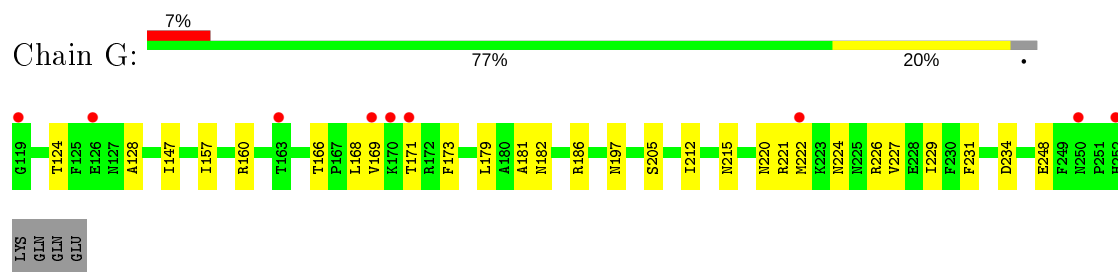
- Molecule 1: Chemotaxis protein motB



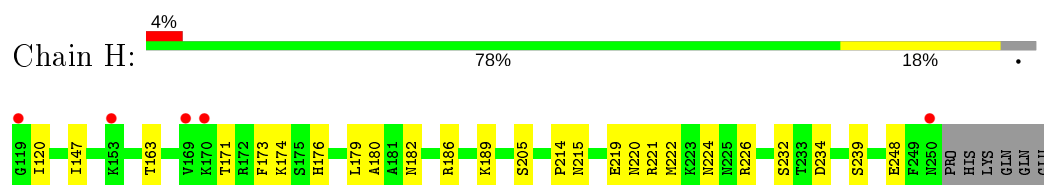
- Molecule 1: Chemotaxis protein motB



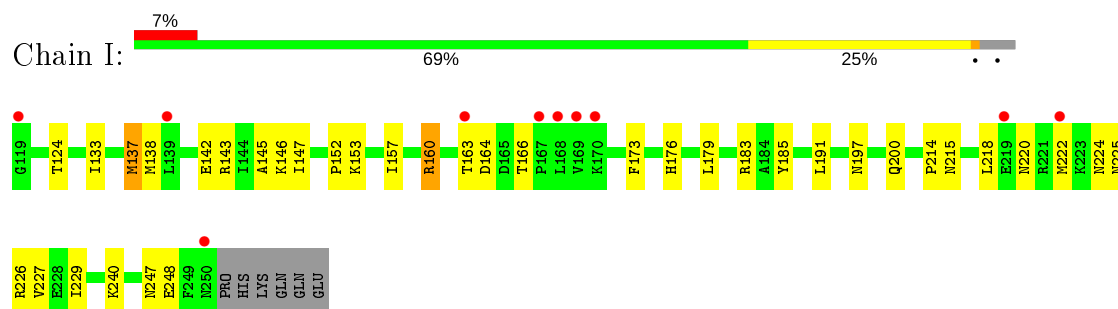
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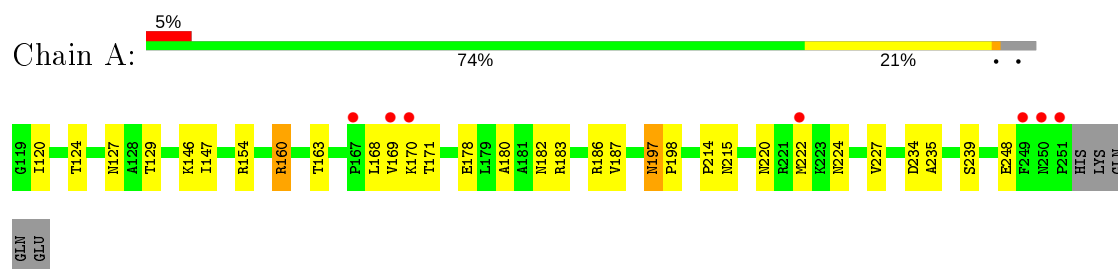
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- Molecule 1: Chemotaxis protein motB

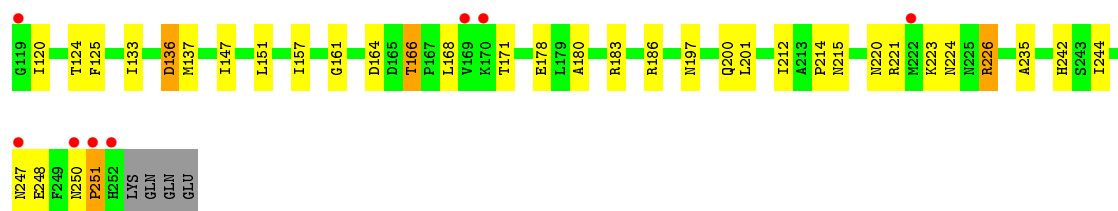


- Molecule 1: Chemotaxis protein motB

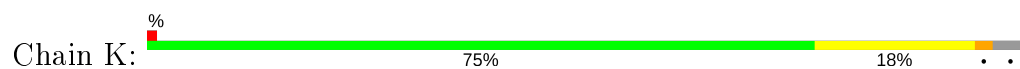


- Molecule 1: Chemotaxis protein motB

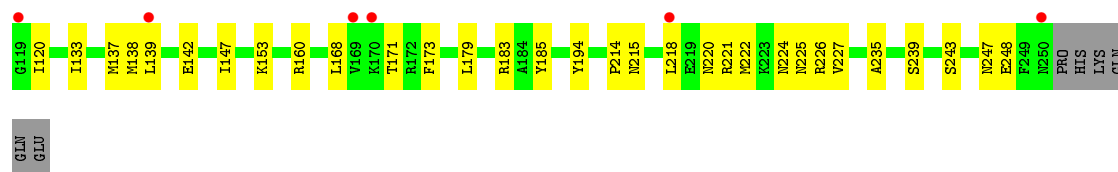
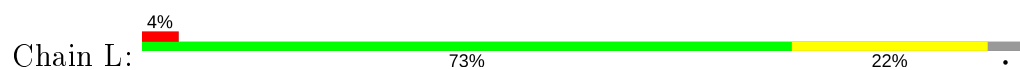




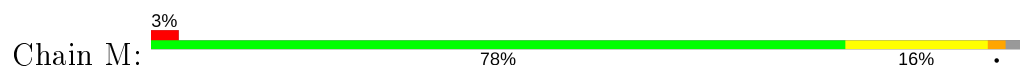
• Molecule 1: Chemotaxis protein motB



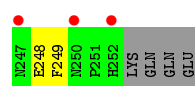
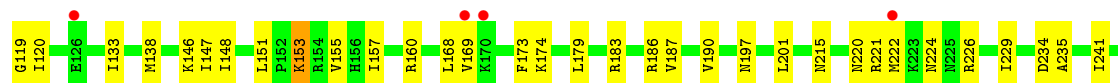
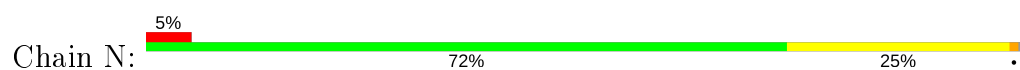
• Molecule 1: Chemotaxis protein motB



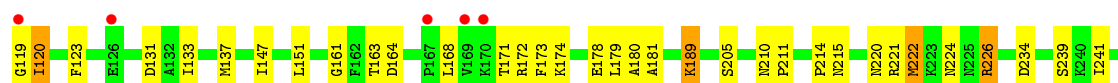
• Molecule 1: Chemotaxis protein motB

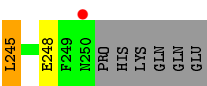


• Molecule 1: Chemotaxis protein motB

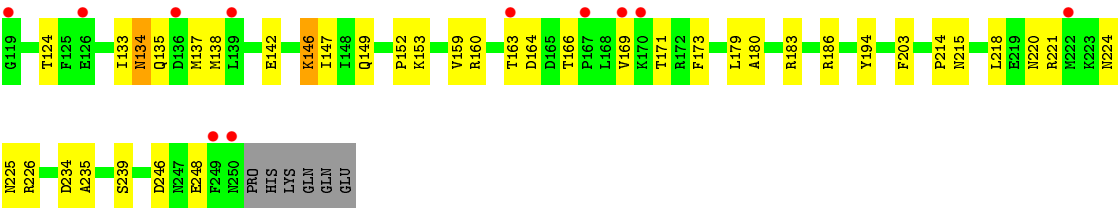


• Molecule 1: Chemotaxis protein motB





● Molecule 1: Chemotaxis protein motB



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	101.78 Å 110.06 Å 113.66 Å 90.00° 104.58° 90.00°	Depositor
Resolution (Å)	15.00 – 2.30 15.00 – 2.30	Depositor EDS
% Data completeness (in resolution range)	96.1 (15.00-2.30) 96.4 (15.00-2.30)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.48 (at 2.29 Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.217 , 0.289 0.224 , 0.295	Depositor DCC
R_{free} test set	5187 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	27.7	Xtriage
Anisotropy	0.638	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 61.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	18460	wwPDB-VP
Average B, all atoms (Å ²)	22.0	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 90.57 % 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 2.1772e-08. 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: AMU

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	0.62	0/1093	0.87	2/1480 (0.1%)
1	B	0.59	0/1112	0.84	1/1505 (0.1%)
1	C	0.59	0/1110	0.86	1/1503 (0.1%)
1	D	0.63	0/1085	0.92	1/1468 (0.1%)
1	E	0.56	0/1085	0.86	2/1468 (0.1%)
1	F	0.57	0/1093	0.83	1/1480 (0.1%)
1	G	0.59	0/1104	0.86	1/1495 (0.1%)
1	H	0.58	0/1085	0.84	2/1468 (0.1%)
1	I	0.56	0/1085	0.83	1/1468 (0.1%)
1	J	0.57	0/1104	0.84	1/1495 (0.1%)
1	K	0.60	0/1085	0.91	2/1468 (0.1%)
1	L	0.56	0/1085	0.84	0/1468
1	M	0.61	0/1093	0.87	2/1480 (0.1%)
1	N	0.60	0/1104	0.87	2/1495 (0.1%)
1	O	0.58	0/1085	0.84	1/1468 (0.1%)
1	P	0.55	0/1085	0.83	1/1468 (0.1%)
All	All	0.59	0/17493	0.86	21/23677 (0.1%)

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	K	0	1

There are no bond length outliers.

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	234	ASP	CB-CG-OD1	8.77	126.19	118.30
1	C	234	ASP	CB-CG-OD1	6.67	124.30	118.30
1	E	234	ASP	CB-CG-OD2	-6.66	112.31	118.30
1	K	160	ARG	NE-CZ-NH2	-6.62	116.99	120.30
1	M	188	MET	CG-SD-CE	6.53	110.65	100.20
1	H	226	ARG	NE-CZ-NH1	5.87	123.23	120.30
1	B	154	ARG	NE-CZ-NH2	-5.76	117.42	120.30
1	A	160	ARG	NE-CZ-NH2	-5.70	117.45	120.30
1	O	234	ASP	CB-CG-OD1	5.56	123.30	118.30
1	D	186	ARG	NE-CZ-NH1	5.49	123.05	120.30
1	N	234	ASP	CB-CG-OD1	5.48	123.23	118.30
1	I	160	ARG	NE-CZ-NH1	5.44	123.02	120.30
1	F	234	ASP	CB-CG-OD1	5.40	123.16	118.30
1	M	234	ASP	CB-CG-OD1	5.35	123.11	118.30
1	P	234	ASP	CB-CG-OD1	5.28	123.06	118.30
1	H	234	ASP	CB-CG-OD1	5.20	122.97	118.30
1	N	160	ARG	NE-CZ-NH2	-5.18	117.71	120.30
1	A	234	ASP	CB-CG-OD1	5.15	122.94	118.30
1	K	226	ARG	NE-CZ-NH1	5.05	122.83	120.30
1	G	234	ASP	CB-CG-OD1	5.04	122.83	118.30
1	J	186	ARG	NE-CZ-NH2	-5.02	117.79	120.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	K	249	PHE	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	1071	0	1052	23	0
1	B	1090	0	1074	22	0
1	C	1088	0	1069	24	0
1	D	1064	0	1045	24	0
1	E	1064	0	1045	27	0
1	F	1071	0	1052	10	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	G	1081	0	1059	20	0
1	H	1064	0	1045	16	0
1	I	1064	0	1045	24	0
1	J	1081	0	1059	35	0
1	K	1064	0	1045	24	0
1	L	1064	0	1045	22	0
1	M	1071	0	1052	14	0
1	N	1081	0	1059	24	0
1	O	1064	0	1045	25	0
1	P	1064	0	1045	25	0
2	D	20	0	18	2	0
2	K	20	0	18	4	0
3	A	68	0	0	5	0
3	B	89	0	0	1	0
3	C	70	0	0	3	0
3	D	99	0	0	6	0
3	E	78	0	0	2	0
3	F	89	0	0	1	0
3	G	85	0	0	3	0
3	H	70	0	0	4	0
3	I	59	0	0	2	0
3	J	76	0	0	2	0
3	K	96	0	0	3	0
3	L	83	0	0	5	0
3	M	92	0	0	3	0
3	N	77	0	0	3	0
3	O	74	0	0	4	0
3	P	69	0	0	5	0
All	All	18460	0	16872	339	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 10.

All (339) 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:H:163:THR:HG23	3:H:321:HOH:O	1.12	1.28
1:H:163:THR:CG2	3:H:321:HOH:O	1.80	1.01
1:F:147:ILE:HD11	1:F:248:GLU:HG2	1.45	0.95
1:C:215:ASN:HD22	1:C:221:ARG:HG2	1.33	0.94
1:L:147:ILE:HD11	1:L:248:GLU:HG2	1.54	0.90

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:147:ILE:HD11	1:I:248:GLU:HG2	1.54	0.87
1:O:179:LEU:HD12	3:O:390:HOH:O	1.74	0.87
1:B:129:THR:OG1	1:B:171:THR:HG21	1.75	0.86
1:C:120[A]:ILE:HD11	1:C:245:LEU:HD22	1.57	0.85
1:M:147:ILE:HD11	1:M:248:GLU:HG2	1.55	0.85
1:J:124:THR:O	1:J:137:MET:CE	2.26	0.82
1:E:168:LEU:O	1:E:171:THR:HG22	1.79	0.82
1:J:215:ASN:HD22	1:J:221:ARG:HG2	1.44	0.82
1:B:120[B]:ILE:HG21	1:E:235:ALA:HB3	1.60	0.81
1:I:183:ARG:HH12	1:I:225:ASN:HD22	1.29	0.81
1:E:147:ILE:HD11	1:E:248:GLU:HG2	1.61	0.80
1:E:215:ASN:HD22	1:E:221:ARG:HG2	1.47	0.79
1:P:215:ASN:HD21	1:P:224:ASN:HD22	1.31	0.79
1:P:215:ASN:ND2	1:P:224:ASN:HD22	1.81	0.77
1:D:176:HIS:ND1	2:D:5:AMU:H3	2.00	0.77
1:H:215:ASN:ND2	1:H:224:ASN:HD22	1.84	0.75
1:B:222:MET:HE3	3:B:341:HOH:O	1.86	0.74
1:P:169:VAL:O	1:P:171:THR:HG22	1.88	0.74
1:L:183:ARG:HH12	1:L:225:ASN:HD22	1.34	0.74
1:O:179:LEU:O	3:O:390:HOH:O	2.04	0.73
1:H:147:ILE:HD11	1:H:248:GLU:HG2	1.69	0.73
1:L:215:ASN:HD22	1:L:221:ARG:HG2	1.52	0.72
1:L:142:GLU:OE1	3:L:334:HOH:O	2.08	0.72
1:O:147:ILE:HD11	1:O:248:GLU:HG2	1.72	0.71
1:H:180:ALA:HB2	3:H:321:HOH:O	1.91	0.68
1:H:182:ASN:O	1:H:186:ARG:HG3	1.93	0.68
1:J:120:ILE:HD13	1:K:235:ALA:HB1	1.75	0.68
1:J:164:ASP:OD2	1:J:166:THR:HG23	1.94	0.67
1:N:235:ALA:HB2	1:O:120:ILE:HD12	1.77	0.67
1:G:186:ARG:HD2	3:G:277:HOH:O	1.95	0.67
1:O:215:ASN:HD22	1:O:221:ARG:HG2	1.58	0.66
1:B:147:ILE:HD12	1:B:244:ILE:HG22	1.76	0.66
1:D:147:ILE:HD11	1:D:248:GLU:HG2	1.77	0.66
1:K:238:LEU:HA	1:K:241:ILE:HD12	1.78	0.66
1:F:178:GLU:HG2	1:G:197:ASN:HD21	1.61	0.66
1:K:215:ASN:ND2	1:K:224:ASN:HD22	1.94	0.66
1:L:218:LEU:HD22	1:L:222:MET:CE	2.26	0.66
1:N:173:PHE:CE1	1:N:179:LEU:HD13	2.31	0.66
1:B:120[B]:ILE:CG2	1:E:235:ALA:HB3	2.26	0.65
1:J:168:LEU:HB3	1:J:171:THR:HG21	1.79	0.65
1:O:222:MET:HA	1:O:222:MET:HE3	1.78	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:147:ILE:HD11	1:F:248:GLU:CG	2.23	0.65
1:B:147:ILE:HD11	1:B:248:GLU:HB2	1.79	0.65
1:K:147:ILE:HD11	1:K:248:GLU:HG2	1.78	0.64
1:J:124:THR:O	1:J:137:MET:HE3	1.97	0.64
1:I:197:ASN:OD1	3:I:309:HOH:O	2.15	0.64
1:B:133:ILE:HD11	1:B:186[A]:ARG:HB3	1.79	0.63
1:C:133:ILE:HG23	3:C:260:HOH:O	1.98	0.63
1:J:164:ASP:CG	1:J:166:THR:HG23	2.20	0.62
1:K:192:ILE:HD12	3:L:333:HOH:O	2.00	0.62
1:I:215:ASN:ND2	1:I:224:ASN:HD22	1.98	0.62
1:D:158:ASN:ND2	1:D:202:SER:OG	2.24	0.62
1:E:183:ARG:HH12	1:E:225:ASN:HD22	1.47	0.62
1:K:223:LYS:HD3	3:K:292:HOH:O	1.99	0.62
1:J:151:LEU:HD23	1:J:244:ILE:HD12	1.82	0.62
1:D:215:ASN:ND2	1:D:224:ASN:HD22	1.98	0.61
1:A:168:LEU:HB3	1:A:171:THR:HG21	1.81	0.61
1:N:169:VAL:HA	3:N:338:HOH:O	2.00	0.61
1:K:193:GLN:NE2	3:K:341:HOH:O	2.32	0.61
1:D:241:ILE:HG22	1:D:245:LEU:HD22	1.83	0.61
1:K:215:ASN:HD21	1:K:224:ASN:HD22	1.46	0.61
1:J:124:THR:O	1:J:137:MET:HE2	2.00	0.60
1:J:133:ILE:HG23	1:J:137:MET:SD	2.41	0.60
1:A:178:GLU:HG2	1:J:197:ASN:HD21	1.65	0.60
1:I:133:ILE:HG22	1:I:138:MET:HG2	1.83	0.60
1:L:218:LEU:HD22	1:L:222:MET:HE1	1.82	0.60
1:P:124:THR:O	1:P:137:MET:CE	2.48	0.60
1:G:215:ASN:HD22	1:G:221:ARG:HG2	1.66	0.60
1:P:159:VAL:O	1:P:203:PHE:HA	2.02	0.60
1:G:169:VAL:HA	3:G:323:HOH:O	2.02	0.60
1:H:180:ALA:HB3	1:H:205:SER:OG	2.02	0.60
1:P:142:GLU:HG3	1:P:194:TYR:CD2	2.37	0.60
1:P:169:VAL:C	1:P:171:THR:HG22	2.22	0.59
1:M:142:GLU:HB2	1:M:194:TYR:CE1	2.36	0.59
1:K:147:ILE:CD1	1:K:249:PHE:CE2	2.86	0.59
1:N:215:ASN:HD22	1:N:221:ARG:HG2	1.68	0.58
1:D:185:TYR:HB3	3:D:354:HOH:O	2.02	0.58
1:G:215:ASN:ND2	1:G:224:ASN:HD22	2.01	0.58
1:G:124:THR:HG21	1:G:222:MET:CE	2.34	0.58
1:E:157:ILE:HG23	1:E:229:ILE:HG23	1.85	0.57
1:H:215:ASN:HD21	1:H:224:ASN:HD22	1.50	0.57
1:I:215:ASN:HD21	1:I:224:ASN:HD22	1.51	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:215:ASN:ND2	1:J:224:ASN:HD22	2.03	0.57
1:C:120[A]:ILE:CD1	1:C:245:LEU:HD22	2.31	0.57
1:P:183:ARG:HH12	1:P:225:ASN:HD22	1.51	0.57
1:I:157:ILE:HG23	1:I:229:ILE:HG23	1.85	0.56
1:K:175:SER:OG	2:K:2:AMU:C10	2.53	0.56
1:D:182:ASN:HA	3:D:354:HOH:O	2.04	0.56
1:J:147:ILE:HD11	1:J:248:GLU:CB	2.36	0.56
1:O:241:ILE:HG22	1:O:245:LEU:HD22	1.87	0.56
1:K:240:LYS:NZ	3:K:337:HOH:O	2.39	0.56
1:A:235:ALA:HB1	1:L:120:ILE:HD13	1.88	0.56
1:K:215:ASN:HD22	1:K:221:ARG:HG2	1.70	0.56
1:A:178:GLU:HG2	1:J:197:ASN:ND2	2.21	0.56
1:P:124:THR:O	1:P:137:MET:HE1	2.06	0.56
1:B:120[B]:ILE:HG21	1:E:235:ALA:CB	2.34	0.55
1:D:190:VAL:HA	1:D:193:GLN:HE21	1.71	0.55
1:A:169:VAL:HG23	1:A:170:LYS:H	1.71	0.55
1:C:215:ASN:ND2	1:C:224:ASN:HD22	2.05	0.55
3:D:354:HOH:O	1:E:198:PRO:CB	2.55	0.55
1:C:147:ILE:HD11	1:C:248:GLU:HG2	1.88	0.55
1:H:215:ASN:ND2	1:H:224:ASN:ND2	2.53	0.55
1:G:157:ILE:HG23	1:G:229:ILE:HG23	1.89	0.55
1:J:147:ILE:HD11	1:J:248:GLU:HB2	1.88	0.55
1:B:212:ILE:HG22	1:B:224:ASN:OD1	2.06	0.55
1:N:147:ILE:HD11	1:N:248:GLU:CB	2.37	0.54
1:B:214:PRO:O	1:B:220:ASN:HB3	2.08	0.54
1:C:247:ASN:N	1:C:247:ASN:HD22	2.04	0.54
1:D:147:ILE:HD11	1:D:248:GLU:CB	2.37	0.54
1:H:176:HIS:O	3:H:321:HOH:O	2.18	0.54
1:B:127:ASN:ND2	1:B:129:THR:HG22	2.23	0.54
1:E:147:ILE:HD11	1:E:248:GLU:CG	2.35	0.54
1:K:175:SER:HG	2:K:2:AMU:C10	2.22	0.53
1:C:164:ASP:CG	1:C:166:THR:HG23	2.28	0.53
1:C:218:LEU:O	1:C:222:MET:HG2	2.09	0.53
1:L:138:MET:CE	1:L:194:TYR:OH	2.57	0.53
1:D:147:ILE:HD11	1:D:248:GLU:CG	2.37	0.53
1:L:147:ILE:CD1	1:L:248:GLU:HG2	2.34	0.53
1:H:215:ASN:HD22	1:H:221:ARG:HG2	1.74	0.53
1:I:142:GLU:OE2	1:I:146:LYS:NZ	2.42	0.53
1:E:215:ASN:ND2	1:E:224:ASN:HD22	2.07	0.53
1:F:214:PRO:O	1:F:220:ASN:HB3	2.09	0.53
1:J:212:ILE:HG23	1:J:226:ARG:NH2	2.24	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:M:127:ASN:HD22	1:M:129:THR:H	1.57	0.52
1:P:214:PRO:O	1:P:220:ASN:HB3	2.09	0.52
1:I:152:PRO:HA	3:I:286:HOH:O	2.08	0.52
1:J:136:ASP:N	1:J:136:ASP:OD2	2.43	0.52
1:H:214:PRO:O	1:H:220:ASN:HB3	2.10	0.52
1:K:157:ILE:HG23	1:K:229:ILE:HG23	1.92	0.52
1:E:167:PRO:HB2	1:E:169:VAL:HG23	1.90	0.52
1:A:197:ASN:HD22	1:A:198:PRO:HD2	1.75	0.52
1:G:147:ILE:HD11	1:G:248:GLU:HG2	1.91	0.52
1:N:235:ALA:CB	1:O:120:ILE:HD12	2.39	0.51
1:N:168:LEU:HD22	1:N:173:PHE:HB2	1.92	0.51
1:P:173:PHE:CE1	1:P:179:LEU:HD13	2.46	0.51
1:G:147:ILE:HD11	1:G:248:GLU:CB	2.40	0.51
1:D:212:ILE:HG13	1:D:226:ARG:CZ	2.40	0.51
1:L:139:LEU:HD23	3:L:334:HOH:O	2.09	0.51
1:C:214:PRO:O	1:C:220:ASN:HB3	2.10	0.51
1:F:127:ASN:HD22	1:F:129:THR:H	1.58	0.51
1:I:218:LEU:O	1:I:222:MET:HG2	2.09	0.51
1:J:220:ASN:HD22	1:J:223:LYS:CE	2.24	0.51
1:B:189:LYS:HD2	1:C:189:LYS:HE3	1.93	0.51
1:D:223:LYS:HG3	3:D:302:HOH:O	2.10	0.51
1:A:160:ARG:O	1:A:227:VAL:HA	2.11	0.51
1:D:147:ILE:HD11	1:D:248:GLU:HB3	1.92	0.51
1:K:147:ILE:HD13	1:K:249:PHE:HE2	1.76	0.51
1:O:163:THR:HG23	1:O:180:ALA:HB2	1.92	0.51
1:B:157:ILE:HB	1:B:201:LEU:HD23	1.93	0.51
1:D:157:ILE:HG23	1:D:229:ILE:HG23	1.93	0.51
1:A:214:PRO:O	1:A:220:ASN:HB3	2.10	0.51
1:C:173:PHE:CE1	1:C:179:LEU:HD13	2.46	0.51
1:M:185:TYR:CZ	1:M:189:LYS:HD3	2.46	0.51
1:O:151:LEU:HD22	1:O:241:ILE:HG12	1.92	0.51
1:C:130:SER:OG	3:C:317:HOH:O	2.19	0.50
1:A:168:LEU:HB3	1:A:171:THR:CG2	2.42	0.50
1:K:176:HIS:ND1	2:K:2:AMU:H3	2.26	0.50
1:D:177:TYR:OH	3:D:343:HOH:O	2.16	0.50
1:D:215:ASN:HD21	1:D:224:ASN:HD22	1.58	0.50
1:H:173:PHE:CE1	1:H:179:LEU:HD13	2.47	0.50
1:O:168:LEU:HD22	1:O:171:THR:HG21	1.94	0.50
1:A:215:ASN:OD1	1:A:224:ASN:ND2	2.44	0.50
1:K:188:MET:HE3	3:L:333:HOH:O	2.11	0.50
1:O:164:ASP:HB2	1:O:215:ASN:ND2	2.27	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:131:ASP:O	1:B:186[A]:ARG:NE	2.45	0.50
1:I:164:ASP:HB2	1:I:215:ASN:ND2	2.27	0.50
1:I:183:ARG:HH12	1:I:225:ASN:ND2	2.05	0.50
1:J:235:ALA:HB1	1:K:120:ILE:HD13	1.93	0.50
1:L:173:PHE:CE1	1:L:179:LEU:HD13	2.47	0.50
1:I:145:ALA:HB2	1:I:191:LEU:HD23	1.93	0.49
1:N:157:ILE:HB	1:N:201:LEU:HD23	1.94	0.49
1:A:147:ILE:HD11	1:A:248:GLU:HG2	1.94	0.49
1:I:133:ILE:HG23	1:I:137:MET:HG3	1.93	0.49
1:O:222:MET:HA	1:O:222:MET:CE	2.41	0.49
1:D:237:ASP:O	1:D:241:ILE:HG13	2.11	0.49
1:N:120:ILE:HD11	1:N:249:PHE:HB3	1.95	0.49
1:D:185:TYR:CB	3:D:354:HOH:O	2.59	0.49
1:J:214:PRO:O	1:J:220:ASN:HB3	2.12	0.49
1:P:215:ASN:ND2	1:P:224:ASN:ND2	2.57	0.49
1:N:190:VAL:HG23	3:N:394:HOH:O	2.12	0.49
1:A:124:THR:CG2	3:A:337:HOH:O	2.61	0.49
1:J:164:ASP:OD1	1:J:166:THR:HG23	2.13	0.49
1:O:173:PHE:CE1	1:O:179:LEU:HD13	2.48	0.48
1:M:197:ASN:HD22	1:M:198:PRO:HD2	1.78	0.48
1:P:142:GLU:O	1:P:146:LYS:HD3	2.13	0.48
1:I:163:THR:OG1	1:I:176:HIS:HB3	2.13	0.48
1:I:214:PRO:O	1:I:220:ASN:HB3	2.13	0.48
1:N:155:VAL:HG21	1:N:241:ILE:HD11	1.95	0.48
1:B:129:THR:OG1	1:B:171:THR:CG2	2.53	0.48
1:G:124:THR:HG21	1:G:222:MET:HE3	1.94	0.48
1:H:189:LYS:HG2	1:I:185:TYR:OH	2.14	0.48
1:J:147:ILE:HD11	1:J:248:GLU:HG2	1.94	0.48
1:C:231:PHE:O	1:D:232:SER:HA	2.14	0.48
1:I:160:ARG:O	1:I:227:VAL:HA	2.14	0.48
1:D:175:SER:HB2	2:D:5:AMU:H113	1.96	0.48
1:O:214:PRO:O	1:O:220:ASN:HB3	2.14	0.48
1:O:189:LYS:NZ	3:O:368:HOH:O	2.14	0.48
1:G:124:THR:HG21	1:G:222:MET:HE1	1.96	0.48
1:N:120:ILE:HD11	1:N:249:PHE:CB	2.44	0.48
1:D:215:ASN:HD22	1:D:221:ARG:HG2	1.78	0.47
1:F:146:LYS:HD3	1:F:146:LYS:N	2.28	0.47
1:L:147:ILE:HD11	1:L:248:GLU:CG	2.36	0.47
1:N:147:ILE:HD11	1:N:248:GLU:HB2	1.97	0.47
1:A:124:THR:HG22	3:A:337:HOH:O	2.14	0.47
1:A:182:ASN:O	1:A:186:ARG:HG3	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:133:ILE:HG23	1:I:137:MET:CG	2.43	0.47
1:H:182:ASN:O	1:H:186:ARG:CG	2.62	0.47
1:B:173:PHE:CE2	1:B:179:LEU:HD13	2.50	0.47
1:E:173:PHE:CE1	1:E:179:LEU:HD13	2.50	0.47
1:L:133:ILE:HG23	1:L:137:MET:HB3	1.96	0.47
1:D:147:ILE:CD1	1:D:248:GLU:HG2	2.44	0.47
1:O:161:GLY:C	1:O:180:ALA:HB1	2.35	0.47
1:N:155:VAL:HG11	1:N:241:ILE:HD11	1.97	0.47
1:A:146:LYS:CD	3:A:365:HOH:O	2.63	0.46
1:C:120[A]:ILE:HD12	1:D:235:ALA:CB	2.46	0.46
1:B:147:ILE:HD12	1:B:244:ILE:CG2	2.45	0.46
1:E:133:ILE:HG22	1:E:138:MET:HG2	1.96	0.46
1:O:133:ILE:HG23	1:O:137:MET:HB3	1.98	0.46
1:E:218:LEU:O	1:E:222:MET:HG2	2.16	0.46
1:F:153:LYS:HD3	3:F:298:HOH:O	2.15	0.46
1:O:178:GLU:HG3	3:P:315:HOH:O	2.16	0.46
1:C:218:LEU:O	1:C:222:MET:HE3	2.16	0.46
1:G:147:ILE:HD11	1:G:248:GLU:HB2	1.97	0.46
1:G:173:PHE:CE1	1:G:179:LEU:HD13	2.51	0.46
1:I:197:ASN:HB3	1:I:200:GLN:HE21	1.81	0.46
1:K:147:ILE:HD13	1:K:249:PHE:CE2	2.50	0.45
1:L:168:LEU:O	1:L:171:THR:HG22	2.16	0.45
1:J:147:ILE:HD11	1:J:248:GLU:CG	2.47	0.45
1:E:139:LEU:HD11	3:E:325:HOH:O	2.16	0.45
1:G:147:ILE:HD11	1:G:248:GLU:CG	2.47	0.45
1:P:137:MET:HE3	3:P:318:HOH:O	2.17	0.45
1:I:143:ARG:O	1:I:147:ILE:HD12	2.16	0.45
1:L:218:LEU:O	1:L:222:MET:HG2	2.17	0.45
1:J:164:ASP:OD2	1:J:166:THR:CG2	2.65	0.45
1:N:151:LEU:HD22	1:N:241:ILE:HG12	1.99	0.45
1:J:151:LEU:HD23	1:J:244:ILE:CD1	2.47	0.45
1:K:173:PHE:CE1	1:K:179:LEU:HD13	2.52	0.45
1:K:147:ILE:CD1	1:K:248:GLU:HG2	2.46	0.45
1:M:169:VAL:HG12	3:M:284:HOH:O	2.16	0.45
1:N:215:ASN:ND2	1:N:224:ASN:HD22	2.15	0.45
1:C:164:ASP:OD2	1:C:166:THR:HG23	2.17	0.44
1:P:138:MET:CE	3:P:301:HOH:O	2.64	0.44
1:C:222:MET:HE3	1:C:222:MET:N	2.31	0.44
1:C:147:ILE:HD11	1:C:248:GLU:CG	2.47	0.44
1:N:157:ILE:HG23	1:N:229:ILE:HG23	1.99	0.44
1:P:218:LEU:HD12	1:P:221:ARG:HD2	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:185:TYR:HE1	3:L:333:HOH:O	1.98	0.44
1:E:157:ILE:HB	1:E:201:LEU:HD23	1.99	0.44
1:E:206:TYR:HD1	1:E:210:ASN:HD22	1.64	0.44
1:E:206:TYR:HD1	1:E:210:ASN:ND2	2.15	0.44
1:E:151:LEU:HD22	1:E:241:ILE:HG12	1.99	0.44
1:L:160:ARG:O	1:L:227:VAL:HA	2.17	0.44
1:P:164:ASP:OD2	1:P:166:THR:HG23	2.17	0.44
1:E:182:ASN:O	1:E:186:ARG:HG2	2.18	0.44
1:A:197:ASN:HD21	1:J:178:GLU:HG3	1.82	0.44
1:N:183:ARG:O	1:N:187:VAL:HG23	2.17	0.44
1:P:147:ILE:HD11	1:P:248:GLU:HG2	2.00	0.44
1:P:152:PRO:HA	3:P:319:HOH:O	2.17	0.44
1:B:213:ALA:HB1	1:B:214:PRO:HD2	2.00	0.44
1:K:147:ILE:HD11	1:K:249:PHE:CE2	2.53	0.44
1:P:164:ASP:CG	1:P:166:THR:HG23	2.38	0.44
1:L:218:LEU:HD22	1:L:222:MET:HE2	1.98	0.43
1:M:147:ILE:HG23	1:M:244:ILE:HG22	1.99	0.43
1:P:135:GLN:NE2	1:P:135:GLN:O	2.50	0.43
1:A:146:LYS:HD2	3:A:365:HOH:O	2.17	0.43
1:K:241:ILE:HG22	1:K:245:LEU:HD22	1.99	0.43
1:A:120:ILE:HD13	1:L:235:ALA:HB1	1.99	0.43
1:O:131:ASP:HB3	3:O:390:HOH:O	2.17	0.43
1:P:133:ILE:HG22	1:P:138:MET:HG2	1.99	0.43
1:A:154:ARG:HD3	3:A:326:HOH:O	2.17	0.43
1:F:163:THR:HG23	1:F:180:ALA:HB2	2.00	0.43
1:J:242:HIS:CD2	3:J:332:HOH:O	2.70	0.43
1:C:160:ARG:O	1:C:227:VAL:HA	2.19	0.43
1:I:147:ILE:CD1	1:I:248:GLU:HG2	2.36	0.43
1:J:161:GLY:O	1:J:180:ALA:HB1	2.19	0.43
1:G:182:ASN:ND2	3:G:340:HOH:O	2.49	0.43
1:J:157:ILE:HB	1:J:201:LEU:HD23	2.01	0.43
1:N:173:PHE:CD1	1:N:179:LEU:HD13	2.53	0.43
1:A:197:ASN:HD22	1:A:198:PRO:CD	2.31	0.43
1:E:200:GLN:NE2	3:E:304:HOH:O	2.52	0.43
1:N:119:GLY:O	1:N:120:ILE:HD13	2.18	0.43
1:C:212:ILE:HG23	1:C:226:ARG:NH2	2.34	0.43
1:E:217:SER:HB3	1:E:220:ASN:HD22	1.83	0.43
1:L:214:PRO:O	1:L:220:ASN:HB3	2.19	0.43
1:P:163:THR:HG23	1:P:180:ALA:HB2	2.00	0.43
1:C:180:ALA:HB3	1:C:205:SER:OG	2.19	0.42
1:A:183:ARG:O	1:A:187:VAL:HG23	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:147:ILE:HD11	1:B:248:GLU:CB	2.48	0.42
1:F:157:ILE:HG23	1:F:229:ILE:HG23	2.00	0.42
1:B:238:LEU:O	1:E:238:LEU:HD21	2.19	0.42
1:J:166:THR:HG22	1:J:221:ARG:HE	1.84	0.42
1:J:215:ASN:HD21	1:J:224:ASN:HD22	1.68	0.42
1:M:175:SER:HB2	3:M:313:HOH:O	2.17	0.42
1:L:215:ASN:ND2	1:L:221:ARG:HG2	2.29	0.42
1:M:176:HIS:ND1	3:M:286:HOH:O	2.35	0.42
1:B:120[A]:ILE:CD1	1:B:245:LEU:HD22	2.49	0.42
1:A:163:THR:HG23	1:A:180:ALA:HB2	2.01	0.42
1:E:215:ASN:HD21	1:E:224:ASN:HD22	1.66	0.42
1:J:125:PHE:CZ	1:J:183:ARG:HB3	2.54	0.42
1:M:214:PRO:O	1:M:220:ASN:HB3	2.19	0.42
1:C:135[B]:GLN:HE21	1:C:135[B]:GLN:H	1.67	0.41
1:F:147:ILE:HD11	1:F:248:GLU:CB	2.50	0.41
1:N:133:ILE:HG22	1:N:138:MET:HG2	2.00	0.41
1:D:147:ILE:CD1	1:D:249:PHE:HE2	2.34	0.41
1:N:148:ILE:O	3:N:379:HOH:O	2.22	0.41
1:G:168:LEU:HB3	1:G:171:THR:HG21	2.03	0.41
1:M:147:ILE:HD11	1:M:248:GLU:CG	2.38	0.41
1:G:181:ALA:HB2	1:G:205:SER:HB2	2.02	0.41
1:J:250:ASN:N	1:J:251:PRO:HD3	2.35	0.41
1:L:215:ASN:ND2	1:L:224:ASN:HD22	2.18	0.41
1:K:190:VAL:HA	1:K:193:GLN:HE21	1.86	0.41
1:C:212:ILE:HD11	3:C:267:HOH:O	2.20	0.41
1:A:127:ASN:HD22	1:A:129:THR:H	1.68	0.41
1:M:178:GLU:HG2	1:N:197:ASN:HD21	1.85	0.41
1:J:200:GLN:OE1	3:J:294:HOH:O	2.22	0.41
1:P:149:GLN:HB3	3:P:325:HOH:O	2.20	0.41
1:I:173:PHE:CE1	1:I:179:LEU:HD13	2.56	0.41
1:O:123:PHE:O	1:O:226:ARG:HA	2.20	0.41
1:O:181:ALA:HB2	1:O:205:SER:HB2	2.03	0.41
1:G:231:PHE:O	1:H:232:SER:HA	2.20	0.40
1:O:210:ASN:N	1:O:211:PRO:CD	2.84	0.40
1:I:164:ASP:OD2	1:I:166:THR:HG23	2.21	0.40
2:K:2:AMU:H81	2:K:2:AMU:H2	1.81	0.40
1:M:142:GLU:HB2	1:M:194:TYR:CZ	2.56	0.40
1:B:160:ARG:O	1:B:227:VAL:HA	2.22	0.40
1:E:213:ALA:HB1	1:E:214:PRO:HD2	2.03	0.40
1:J:247:ASN:HD22	1:J:247:ASN:N	2.19	0.40
1:M:120:ILE:HD13	1:P:235:ALA:HB1	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:N:186:ARG:O	1:N:190:VAL:HG23	2.22	0.40
1:G:128:ALA:HB1	1:G:166:THR:OG1	2.21	0.40
1:O:215:ASN:ND2	1:O:224:ASN:HD22	2.20	0.40
1:E:147:ILE:CD1	1:E:248:GLU:HG2	2.43	0.40
1:G:160:ARG:O	1:G:227:VAL:HA	2.22	0.40
1:O:119:GLY:O	1:O:120:ILE:C	2.60	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	131/138 (95%)	123 (94%)	8 (6%)	0	100	100
1	B	133/138 (96%)	128 (96%)	5 (4%)	0	100	100
1	C	133/138 (96%)	127 (96%)	6 (4%)	0	100	100
1	D	130/138 (94%)	126 (97%)	4 (3%)	0	100	100
1	E	130/138 (94%)	125 (96%)	4 (3%)	1 (1%)	19	23
1	F	131/138 (95%)	128 (98%)	3 (2%)	0	100	100
1	G	132/138 (96%)	128 (97%)	4 (3%)	0	100	100
1	H	130/138 (94%)	125 (96%)	4 (3%)	1 (1%)	19	23
1	I	130/138 (94%)	123 (95%)	7 (5%)	0	100	100
1	J	132/138 (96%)	127 (96%)	4 (3%)	1 (1%)	19	23
1	K	130/138 (94%)	126 (97%)	4 (3%)	0	100	100
1	L	130/138 (94%)	123 (95%)	7 (5%)	0	100	100
1	M	131/138 (95%)	129 (98%)	2 (2%)	0	100	100
1	N	132/138 (96%)	128 (97%)	3 (2%)	1 (1%)	19	23
1	O	130/138 (94%)	126 (97%)	3 (2%)	1 (1%)	19	23

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	P	130/138 (94%)	125 (96%)	4 (3%)	1 (1%)	19	23
All	All	2095/2208 (95%)	2017 (96%)	72 (3%)	6 (0%)	41	50

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	H	120	ILE
1	E	134	ASN
1	N	153	LYS
1	O	120	ILE
1	P	134	ASN
1	J	251	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	120/126 (95%)	117 (98%)	3 (2%)	47	65
1	B	122/126 (97%)	119 (98%)	3 (2%)	47	65
1	C	122/126 (97%)	115 (94%)	7 (6%)	20	28
1	D	119/126 (94%)	114 (96%)	5 (4%)	30	42
1	E	119/126 (94%)	113 (95%)	6 (5%)	24	34
1	F	120/126 (95%)	115 (96%)	5 (4%)	30	42
1	G	121/126 (96%)	118 (98%)	3 (2%)	47	65
1	H	119/126 (94%)	114 (96%)	5 (4%)	30	42
1	I	119/126 (94%)	113 (95%)	6 (5%)	24	34
1	J	121/126 (96%)	118 (98%)	3 (2%)	47	65
1	K	119/126 (94%)	114 (96%)	5 (4%)	30	42
1	L	119/126 (94%)	114 (96%)	5 (4%)	30	42
1	M	120/126 (95%)	112 (93%)	8 (7%)	16	21
1	N	121/126 (96%)	115 (95%)	6 (5%)	24	34

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	O	119/126 (94%)	112 (94%)	7 (6%)	19	27
1	P	119/126 (94%)	111 (93%)	8 (7%)	16	21
All	All	1919/2016 (95%)	1834 (96%)	85 (4%)	28	39

All (85) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	174	LYS
1	B	226	ARG
1	B	239	SER
1	C	127	ASN
1	C	153	LYS
1	C	218	LEU
1	C	222	MET
1	C	226	ARG
1	C	239	SER
1	C	247	ASN
1	D	153	LYS
1	D	174	LYS
1	D	223	LYS
1	D	226	ARG
1	D	245	LEU
1	E	121	ASP
1	E	130	SER
1	E	153	LYS
1	E	218	LEU
1	E	226	ARG
1	E	239	SER
1	F	127	ASN
1	F	197	ASN
1	F	218	LEU
1	F	226	ARG
1	F	239	SER
1	G	212	ILE
1	G	220	ASN
1	G	226	ARG
1	H	171	THR
1	H	174	LYS
1	H	219	GLU
1	H	222	MET
1	H	239	SER

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Mol	Chain	Res	Type
1	I	124	THR
1	I	137	MET
1	I	153	LYS
1	I	226	ARG
1	I	240	LYS
1	I	247	ASN
1	A	197	ASN
1	A	222	MET
1	A	239	SER
1	J	136	ASP
1	J	166	THR
1	J	226	ARG
1	K	135	GLN
1	K	153	LYS
1	K	218	LEU
1	K	226	ARG
1	K	245	LEU
1	L	153	LYS
1	L	226	ARG
1	L	239	SER
1	L	243	SER
1	L	247	ASN
1	M	127	ASN
1	M	134	ASN
1	M	197	ASN
1	M	214	PRO
1	M	218	LEU
1	M	226	ARG
1	M	239	SER
1	M	243	SER
1	N	146	LYS
1	N	153	LYS
1	N	174	LYS
1	N	220	ASN
1	N	222	MET
1	N	226	ARG
1	O	172	ARG
1	O	174	LYS
1	O	189	LYS
1	O	222	MET
1	O	226	ARG
1	O	239	SER

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Mol	Chain	Res	Type
1	O	245	LEU
1	P	134	ASN
1	P	146	LYS
1	P	153	LYS
1	P	160	ARG
1	P	186	ARG
1	P	226	ARG
1	P	239	SER
1	P	246	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (63) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	127	ASN
1	B	220	ASN
1	B	250	ASN
1	C	176	HIS
1	C	182	ASN
1	C	215	ASN
1	C	220	ASN
1	C	247	ASN
1	D	193	GLN
1	D	215	ASN
1	E	134	ASN
1	E	200	GLN
1	E	210	ASN
1	E	215	ASN
1	E	220	ASN
1	E	225	ASN
1	F	127	ASN
1	F	135	GLN
1	F	197	ASN
1	F	220	ASN
1	F	247	ASN
1	G	182	ASN
1	G	215	ASN
1	H	176	HIS
1	H	193	GLN
1	H	215	ASN
1	I	134	ASN
1	I	176	HIS
1	I	193	GLN

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Mol	Chain	Res	Type
1	I	200	GLN
1	I	215	ASN
1	I	225	ASN
1	A	127	ASN
1	A	197	ASN
1	A	220	ASN
1	A	247	ASN
1	J	215	ASN
1	J	220	ASN
1	J	247	ASN
1	K	193	GLN
1	K	215	ASN
1	L	200	GLN
1	L	210	ASN
1	L	215	ASN
1	L	225	ASN
1	L	250	ASN
1	M	127	ASN
1	M	197	ASN
1	M	220	ASN
1	M	247	ASN
1	N	135	GLN
1	N	215	ASN
1	N	220	ASN
1	N	247	ASN
1	O	176	HIS
1	O	193	GLN
1	O	215	ASN
1	P	134	ASN
1	P	182	ASN
1	P	193	GLN
1	P	200	GLN
1	P	215	ASN
1	P	225	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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

2 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	AMU	D	5	-	17,20,20	0.49	0	22,28,28	1.64	4 (18%)
2	AMU	K	2	-	17,20,20	0.56	0	22,28,28	2.16	8 (36%)

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
2	AMU	D	5	-	-	5/10/34/34	0/1/1/1
2	AMU	K	2	-	-	5/10/34/34	0/1/1/1

There are no bond length outliers.

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	K	2	AMU	C11-C9-C10	-5.73	105.94	113.35
2	D	5	AMU	C11-C9-C10	-4.09	108.07	113.35
2	K	2	AMU	O5-C1-C2	-3.41	106.09	109.52
2	K	2	AMU	C1-C2-N2	3.37	114.63	110.73
2	K	2	AMU	C3-C2-N2	-3.34	105.36	110.91
2	K	2	AMU	C1-O5-C5	-2.70	108.57	113.66
2	K	2	AMU	O1-C1-C2	2.69	114.81	109.22
2	D	5	AMU	C1-C2-N2	2.68	113.83	110.73
2	D	5	AMU	C3-C2-N2	-2.46	106.83	110.91

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	K	2	AMU	C8-C7-N2	2.36	120.09	116.10
2	D	5	AMU	O1-C1-C2	2.04	113.46	109.22
2	K	2	AMU	C1-C2-C3	2.03	113.15	110.25

There are no chirality outliers.

All (10) torsion outliers are listed below:

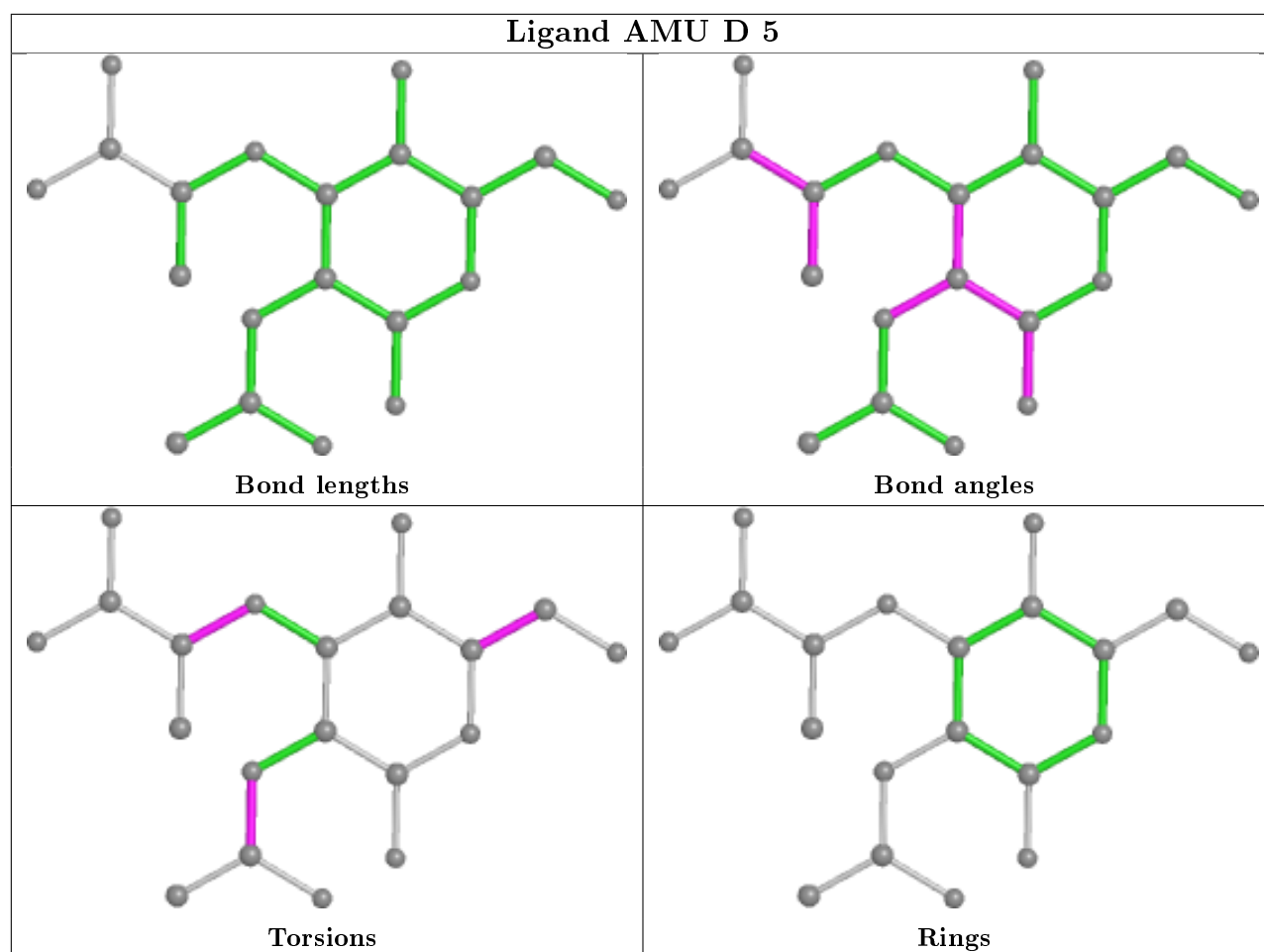
Mol	Chain	Res	Type	Atoms
2	D	5	AMU	C10-C9-O3-C3
2	D	5	AMU	C11-C9-O3-C3
2	K	2	AMU	C10-C9-O3-C3
2	K	2	AMU	C11-C9-O3-C3
2	D	5	AMU	C8-C7-N2-C2
2	D	5	AMU	O7-C7-N2-C2
2	K	2	AMU	C8-C7-N2-C2
2	K	2	AMU	O7-C7-N2-C2
2	K	2	AMU	O5-C5-C6-O6
2	D	5	AMU	O5-C5-C6-O6

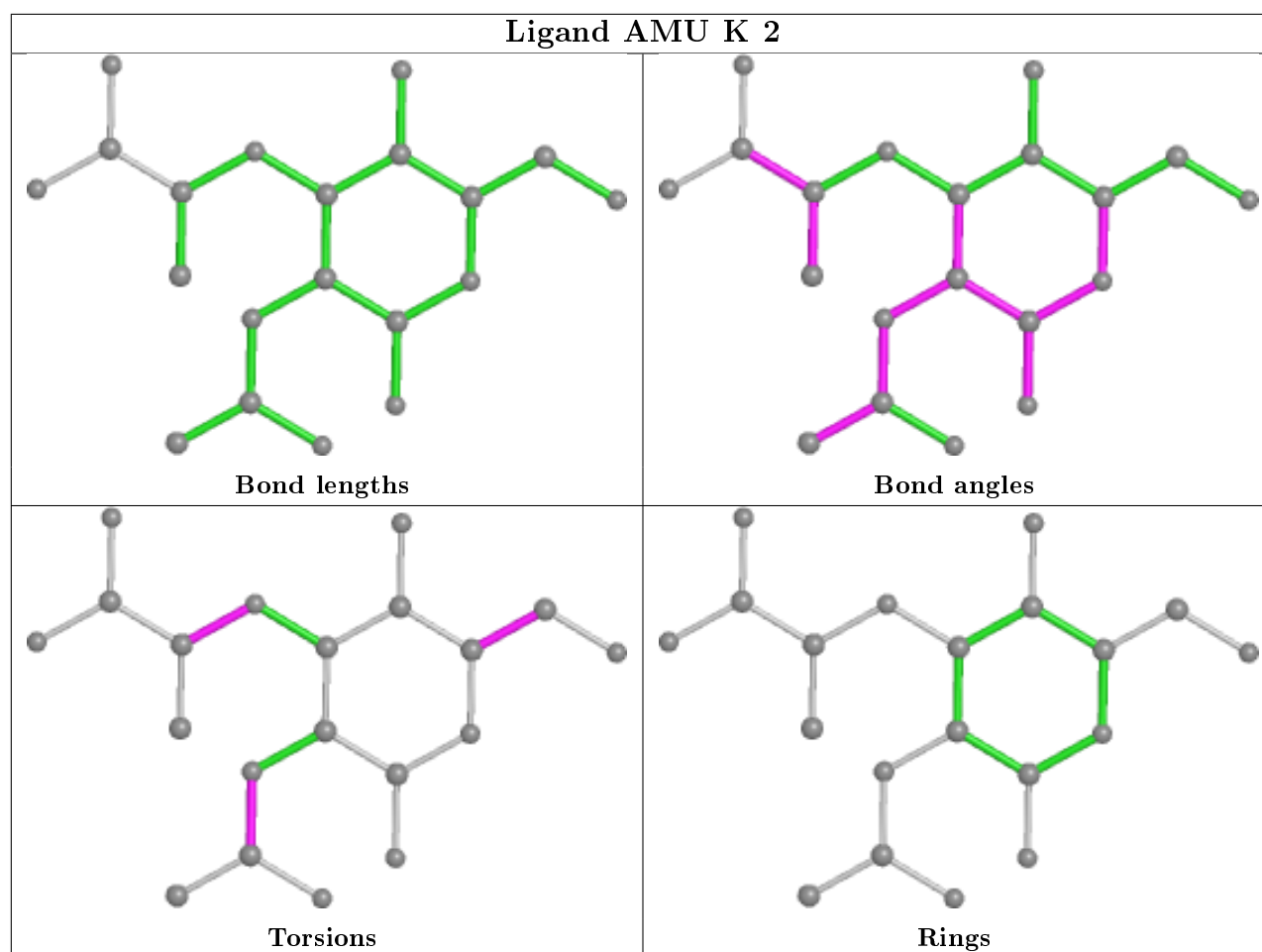
There are no ring outliers.

2 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	5	AMU	2	0
2	K	2	AMU	4	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





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	133/138 (96%)	0.12	7 (5%)	26 33	7, 17, 44, 55	0
1	B	133/138 (96%)	0.12	6 (4%)	33 40	8, 16, 44, 59	0
1	C	133/138 (96%)	0.27	5 (3%)	40 47	7, 18, 45, 53	0
1	D	132/138 (95%)	-0.03	2 (1%)	73 79	7, 16, 32, 38	0
1	E	132/138 (95%)	0.16	5 (3%)	40 47	9, 21, 38, 50	0
1	F	133/138 (96%)	0.14	4 (3%)	50 57	9, 19, 37, 52	0
1	G	134/138 (97%)	0.21	9 (6%)	17 23	11, 22, 38, 45	0
1	H	132/138 (95%)	0.15	5 (3%)	40 47	8, 19, 41, 58	0
1	I	132/138 (95%)	0.45	10 (7%)	13 18	11, 23, 46, 55	0
1	J	134/138 (97%)	0.30	8 (5%)	21 28	7, 19, 49, 66	0
1	K	132/138 (95%)	-0.01	1 (0%)	86 89	7, 17, 31, 41	0
1	L	132/138 (95%)	0.16	6 (4%)	33 40	10, 21, 39, 54	0
1	M	133/138 (96%)	0.08	4 (3%)	50 57	9, 20, 37, 54	0
1	N	134/138 (97%)	0.19	7 (5%)	27 34	9, 23, 38, 48	0
1	O	132/138 (95%)	0.18	6 (4%)	33 40	8, 19, 44, 55	0
1	P	132/138 (95%)	0.44	11 (8%)	11 15	11, 24, 46, 56	0
All	All	2123/2208 (96%)	0.18	96 (4%)	33 40	7, 20, 42, 66	0

All (96) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	169	VAL	9.2
1	F	251	PRO	8.9
1	H	169	VAL	8.1
1	O	169	VAL	7.5
1	M	251	PRO	7.4

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Mol	Chain	Res	Type	RSRZ
1	J	252	HIS	7.0
1	P	119	GLY	6.1
1	J	169	VAL	6.0
1	J	251	PRO	5.3
1	O	250	ASN	5.3
1	I	119	GLY	5.2
1	N	169	VAL	5.2
1	B	251	PRO	5.1
1	B	170	LYS	4.8
1	K	169	VAL	4.6
1	B	169	VAL	4.6
1	H	170	LYS	4.5
1	D	169	VAL	4.5
1	M	250	ASN	4.4
1	G	169	VAL	4.3
1	E	250	ASN	4.2
1	P	169	VAL	4.1
1	G	252	HIS	4.0
1	A	251	PRO	4.0
1	A	169	VAL	4.0
1	L	250	ASN	3.8
1	A	170	LYS	3.7
1	L	218	LEU	3.7
1	I	250	ASN	3.6
1	B	250	ASN	3.6
1	P	170	LYS	3.6
1	I	169	VAL	3.5
1	N	252	HIS	3.4
1	E	119	GLY	3.4
1	M	169	VAL	3.3
1	L	169	VAL	3.3
1	P	250	ASN	3.2
1	C	250	ASN	3.2
1	B	126	GLU	3.2
1	C	251	PRO	3.1
1	H	250	ASN	3.1
1	I	163	THR	3.1
1	C	119	GLY	3.1
1	N	126	GLU	3.1
1	P	136	ASP	3.0
1	F	250	ASN	3.0
1	F	119	GLY	3.0

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Mol	Chain	Res	Type	RSRZ
1	L	119	GLY	3.0
1	J	247	ASN	2.9
1	L	170	LYS	2.9
1	P	163	THR	2.8
1	L	139	LEU	2.8
1	C	170	LYS	2.8
1	J	119	GLY	2.7
1	B	222	MET	2.7
1	P	222	MET	2.7
1	A	250	ASN	2.6
1	O	170	LYS	2.6
1	O	119	GLY	2.6
1	E	126	GLU	2.6
1	H	119	GLY	2.6
1	G	222	MET	2.6
1	I	222	MET	2.6
1	P	249	PHE	2.6
1	I	139	LEU	2.5
1	N	170	LYS	2.4
1	G	170	LYS	2.4
1	A	167	PRO	2.4
1	D	170	LYS	2.3
1	P	139	LEU	2.3
1	O	126	GLU	2.3
1	J	170	LYS	2.3
1	G	119	GLY	2.3
1	I	167	PRO	2.3
1	N	250	ASN	2.2
1	P	167	PRO	2.2
1	G	171	THR	2.2
1	F	222	MET	2.2
1	I	170	LYS	2.2
1	N	222	MET	2.2
1	M	119	GLY	2.2
1	E	219	GLU	2.2
1	G	250	ASN	2.1
1	A	249	PHE	2.1
1	P	126	GLU	2.1
1	H	153	LYS	2.1
1	J	250	ASN	2.1
1	G	126	GLU	2.1
1	E	170	LYS	2.1

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Mol	Chain	Res	Type	RSRZ
1	N	247	ASN	2.1
1	I	168	LEU	2.1
1	A	222	MET	2.1
1	I	219	GLU	2.1
1	G	163	THR	2.1
1	O	167	PRO	2.1
1	J	222	MET	2.0

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

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

6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

6.4 Ligands [i](#)

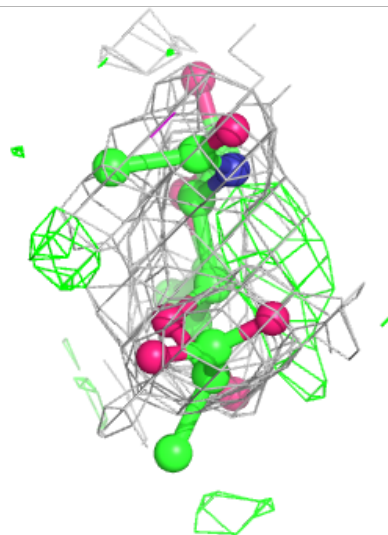
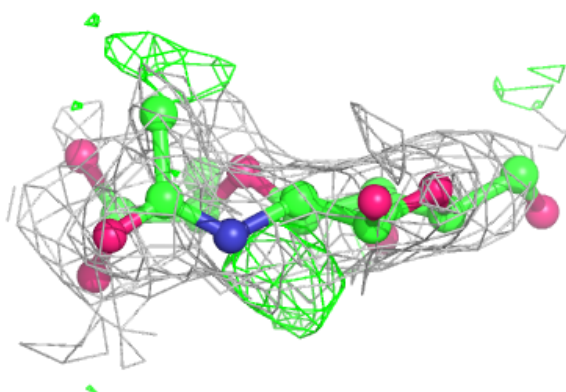
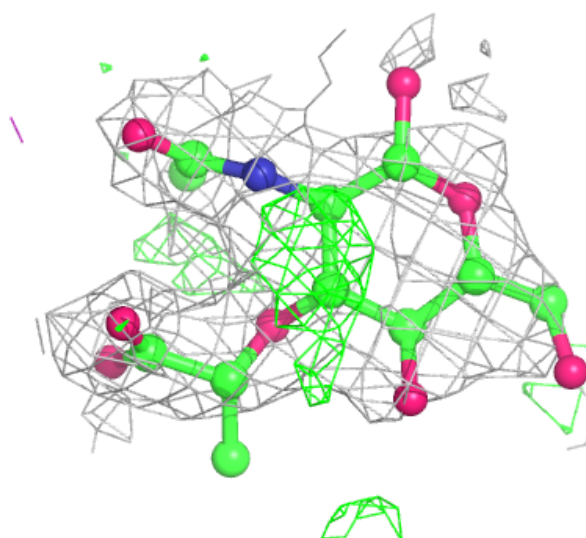
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	AMU	D	5	20/20	0.76	0.26	47,64,72,74	0
2	AMU	K	2	20/20	0.78	0.33	47,61,74,77	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.

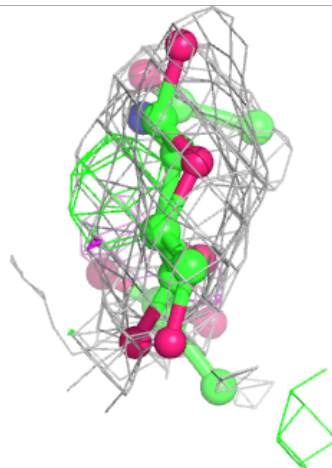
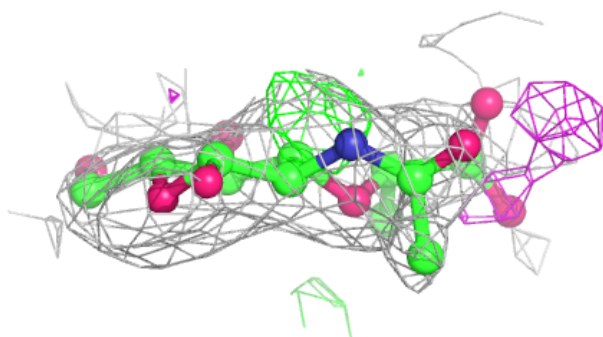
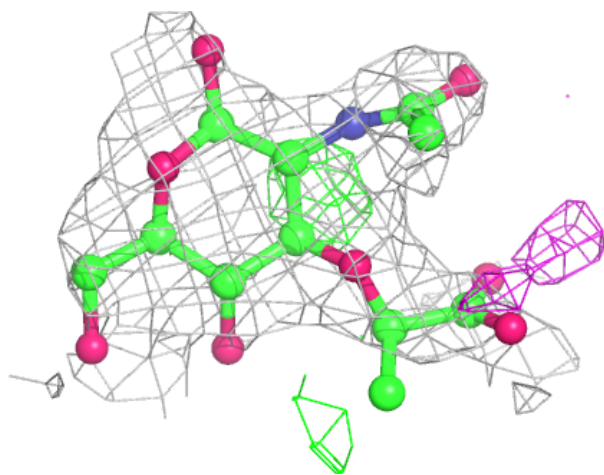
Electron density around AMU D 5:

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



Electron density around AMU K 2:

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



6.5 Other polymers [i](#)

There are no such residues in this entry.