



# Full wwPDB X-ray Structure Validation Report ⓘ

Jun 16, 2020 – 12:09 am BST

PDB ID : 5EXK  
Title : Crystal structure of M. tuberculosis lipoyl synthase with 6-thiooctanoyl peptide intermediate  
Authors : McLaughlin, M.I.; Lanz, N.D.; Goldman, P.J.; Lee, K.-H.; Booker, S.J.; Drennan, C.L.  
Deposited on : 2015-11-23  
Resolution : 1.86 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](mailto: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.

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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 : 2.11  
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) : 2.11

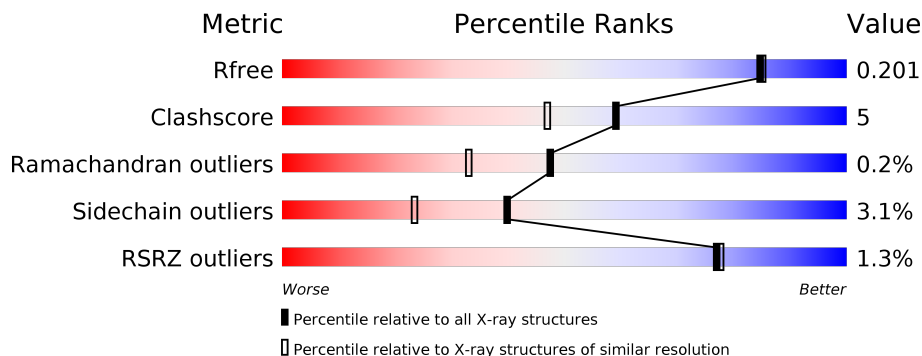
# 1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.86 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	2469 (1.86-1.86)
Clashscore	141614	2625 (1.86-1.86)
Ramachandran outliers	138981	2592 (1.86-1.86)
Sidechain outliers	138945	2592 (1.86-1.86)
RSRZ outliers	127900	2436 (1.86-1.86)

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  $\geq 3$ , 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	331	
1	C	331	
1	E	331	
1	G	331	
1	I	331	
1	K	331	

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Mol	Chain	Length	Quality of chain
2	B	8	
2	D	8	
2	F	8	
2	H	8	
2	J	8	
2	L	8	

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
9	IPA	C	405	-	-	-	X

## 2 Entry composition [i](#)

There are 10 unique types of molecules in this entry. The entry contains 16721 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 Lipoyl synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	306	Total 2381	C 1493	N 435	O 445	S 8	0	4	0
1	C	300	Total 2346	C 1476	N 422	O 440	S 8	0	7	0
1	E	302	Total 2385	C 1497	N 432	O 448	S 8	0	7	0
1	G	296	Total 2282	C 1438	N 402	O 434	S 8	0	6	0
1	I	296	Total 2290	C 1440	N 408	O 434	S 8	0	4	0
1	K	296	Total 2300	C 1442	N 416	O 434	S 8	0	4	0

There are 120 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	initiating methionine	UNP P9WK91
A	-18	GLY	-	expression tag	UNP P9WK91
A	-17	SER	-	expression tag	UNP P9WK91
A	-16	SER	-	expression tag	UNP P9WK91
A	-15	HIS	-	expression tag	UNP P9WK91
A	-14	HIS	-	expression tag	UNP P9WK91
A	-13	HIS	-	expression tag	UNP P9WK91
A	-12	HIS	-	expression tag	UNP P9WK91
A	-11	HIS	-	expression tag	UNP P9WK91
A	-10	HIS	-	expression tag	UNP P9WK91
A	-9	SER	-	expression tag	UNP P9WK91
A	-8	SER	-	expression tag	UNP P9WK91
A	-7	GLY	-	expression tag	UNP P9WK91
A	-6	LEU	-	expression tag	UNP P9WK91
A	-5	VAL	-	expression tag	UNP P9WK91
A	-4	PRO	-	expression tag	UNP P9WK91
A	-3	ARG	-	expression tag	UNP P9WK91

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	expression tag	UNP P9WK91
A	-1	SER	-	expression tag	UNP P9WK91
A	0	HIS	-	expression tag	UNP P9WK91
C	-19	MET	-	initiating methionine	UNP P9WK91
C	-18	GLY	-	expression tag	UNP P9WK91
C	-17	SER	-	expression tag	UNP P9WK91
C	-16	SER	-	expression tag	UNP P9WK91
C	-15	HIS	-	expression tag	UNP P9WK91
C	-14	HIS	-	expression tag	UNP P9WK91
C	-13	HIS	-	expression tag	UNP P9WK91
C	-12	HIS	-	expression tag	UNP P9WK91
C	-11	HIS	-	expression tag	UNP P9WK91
C	-10	HIS	-	expression tag	UNP P9WK91
C	-9	SER	-	expression tag	UNP P9WK91
C	-8	SER	-	expression tag	UNP P9WK91
C	-7	GLY	-	expression tag	UNP P9WK91
C	-6	LEU	-	expression tag	UNP P9WK91
C	-5	VAL	-	expression tag	UNP P9WK91
C	-4	PRO	-	expression tag	UNP P9WK91
C	-3	ARG	-	expression tag	UNP P9WK91
C	-2	GLY	-	expression tag	UNP P9WK91
C	-1	SER	-	expression tag	UNP P9WK91
C	0	HIS	-	expression tag	UNP P9WK91
E	-19	MET	-	initiating methionine	UNP P9WK91
E	-18	GLY	-	expression tag	UNP P9WK91
E	-17	SER	-	expression tag	UNP P9WK91
E	-16	SER	-	expression tag	UNP P9WK91
E	-15	HIS	-	expression tag	UNP P9WK91
E	-14	HIS	-	expression tag	UNP P9WK91
E	-13	HIS	-	expression tag	UNP P9WK91
E	-12	HIS	-	expression tag	UNP P9WK91
E	-11	HIS	-	expression tag	UNP P9WK91
E	-10	HIS	-	expression tag	UNP P9WK91
E	-9	SER	-	expression tag	UNP P9WK91
E	-8	SER	-	expression tag	UNP P9WK91
E	-7	GLY	-	expression tag	UNP P9WK91
E	-6	LEU	-	expression tag	UNP P9WK91
E	-5	VAL	-	expression tag	UNP P9WK91
E	-4	PRO	-	expression tag	UNP P9WK91
E	-3	ARG	-	expression tag	UNP P9WK91
E	-2	GLY	-	expression tag	UNP P9WK91
E	-1	SER	-	expression tag	UNP P9WK91

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Chain	Residue	Modelled	Actual	Comment	Reference
E	0	HIS	-	expression tag	UNP P9WK91
G	-19	MET	-	initiating methionine	UNP P9WK91
G	-18	GLY	-	expression tag	UNP P9WK91
G	-17	SER	-	expression tag	UNP P9WK91
G	-16	SER	-	expression tag	UNP P9WK91
G	-15	HIS	-	expression tag	UNP P9WK91
G	-14	HIS	-	expression tag	UNP P9WK91
G	-13	HIS	-	expression tag	UNP P9WK91
G	-12	HIS	-	expression tag	UNP P9WK91
G	-11	HIS	-	expression tag	UNP P9WK91
G	-10	HIS	-	expression tag	UNP P9WK91
G	-9	SER	-	expression tag	UNP P9WK91
G	-8	SER	-	expression tag	UNP P9WK91
G	-7	GLY	-	expression tag	UNP P9WK91
G	-6	LEU	-	expression tag	UNP P9WK91
G	-5	VAL	-	expression tag	UNP P9WK91
G	-4	PRO	-	expression tag	UNP P9WK91
G	-3	ARG	-	expression tag	UNP P9WK91
G	-2	GLY	-	expression tag	UNP P9WK91
G	-1	SER	-	expression tag	UNP P9WK91
G	0	HIS	-	expression tag	UNP P9WK91
I	-19	MET	-	initiating methionine	UNP P9WK91
I	-18	GLY	-	expression tag	UNP P9WK91
I	-17	SER	-	expression tag	UNP P9WK91
I	-16	SER	-	expression tag	UNP P9WK91
I	-15	HIS	-	expression tag	UNP P9WK91
I	-14	HIS	-	expression tag	UNP P9WK91
I	-13	HIS	-	expression tag	UNP P9WK91
I	-12	HIS	-	expression tag	UNP P9WK91
I	-11	HIS	-	expression tag	UNP P9WK91
I	-10	HIS	-	expression tag	UNP P9WK91
I	-9	SER	-	expression tag	UNP P9WK91
I	-8	SER	-	expression tag	UNP P9WK91
I	-7	GLY	-	expression tag	UNP P9WK91
I	-6	LEU	-	expression tag	UNP P9WK91
I	-5	VAL	-	expression tag	UNP P9WK91
I	-4	PRO	-	expression tag	UNP P9WK91
I	-3	ARG	-	expression tag	UNP P9WK91
I	-2	GLY	-	expression tag	UNP P9WK91
I	-1	SER	-	expression tag	UNP P9WK91
I	0	HIS	-	expression tag	UNP P9WK91
K	-19	MET	-	initiating methionine	UNP P9WK91

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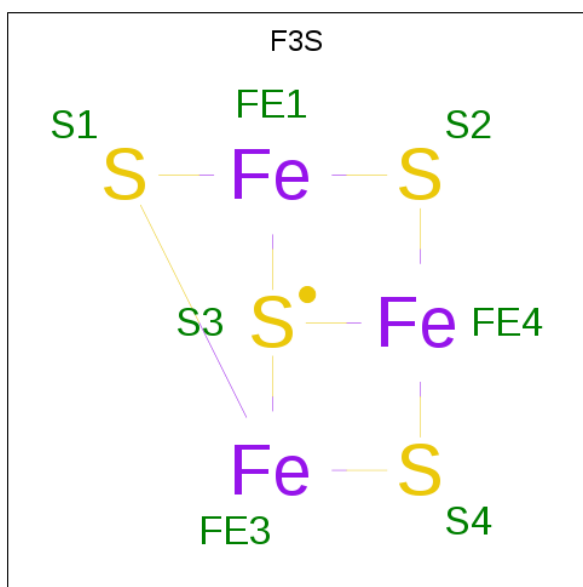
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Chain	Residue	Modelled	Actual	Comment	Reference
K	-18	GLY	-	expression tag	UNP P9WK91
K	-17	SER	-	expression tag	UNP P9WK91
K	-16	SER	-	expression tag	UNP P9WK91
K	-15	HIS	-	expression tag	UNP P9WK91
K	-14	HIS	-	expression tag	UNP P9WK91
K	-13	HIS	-	expression tag	UNP P9WK91
K	-12	HIS	-	expression tag	UNP P9WK91
K	-11	HIS	-	expression tag	UNP P9WK91
K	-10	HIS	-	expression tag	UNP P9WK91
K	-9	SER	-	expression tag	UNP P9WK91
K	-8	SER	-	expression tag	UNP P9WK91
K	-7	GLY	-	expression tag	UNP P9WK91
K	-6	LEU	-	expression tag	UNP P9WK91
K	-5	VAL	-	expression tag	UNP P9WK91
K	-4	PRO	-	expression tag	UNP P9WK91
K	-3	ARG	-	expression tag	UNP P9WK91
K	-2	GLY	-	expression tag	UNP P9WK91
K	-1	SER	-	expression tag	UNP P9WK91
K	0	HIS	-	expression tag	UNP P9WK91

- Molecule 2 is a protein called Octanoylated peptide from M. tuberculosis H protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	8	66	40	9	16	1	0	0	0
2	D	8	69	41	9	18	1	0	0	0
2	F	6	51	32	7	11	1	0	0	0
2	H	6	51	32	7	11	1	0	0	0
2	J	6	48	31	7	9	1	0	0	0
2	L	6	51	32	7	11	1	0	0	0

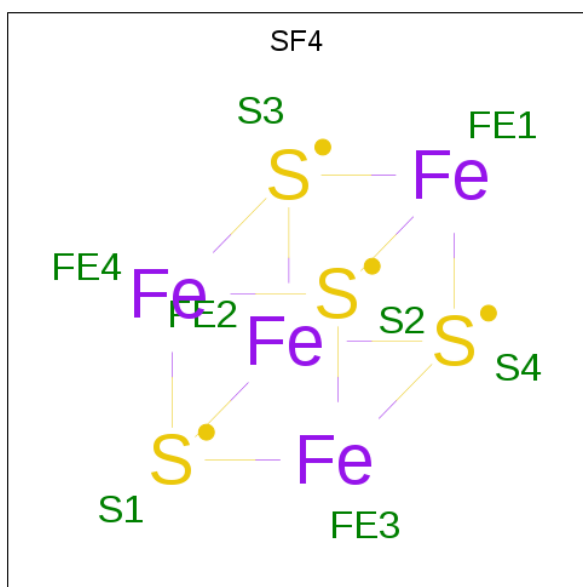
- Molecule 3 is FE3-S4 CLUSTER (three-letter code: F3S) (formula: Fe<sub>3</sub>S<sub>4</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	Fe	S	0	0
			6	3	3		
3	C	1	Total	Fe	S	0	0
			6	3	3		
3	E	1	Total	Fe	S	0	0
			6	3	3		
3	G	1	Total	Fe	S	0	0
			6	3	3		
3	I	1	Total	Fe	S	0	0
			6	3	3		
3	K	1	Total	Fe	S	0	0
			6	3	3		

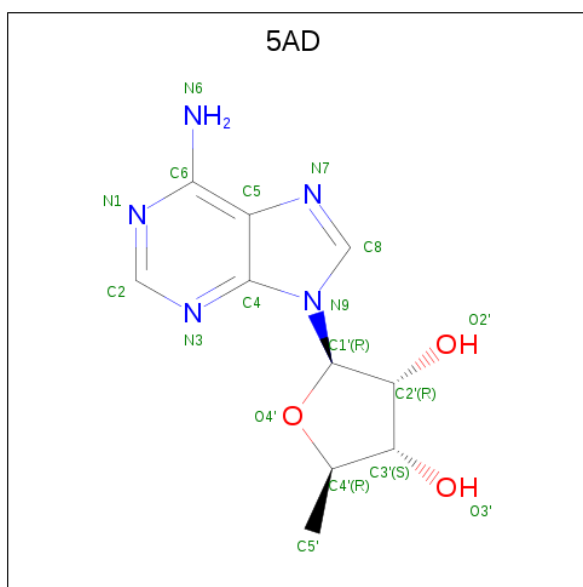
- Molecule 4 is IRON/SULFUR CLUSTER (three-letter code: SF4) (formula: Fe<sub>4</sub>S<sub>4</sub>).





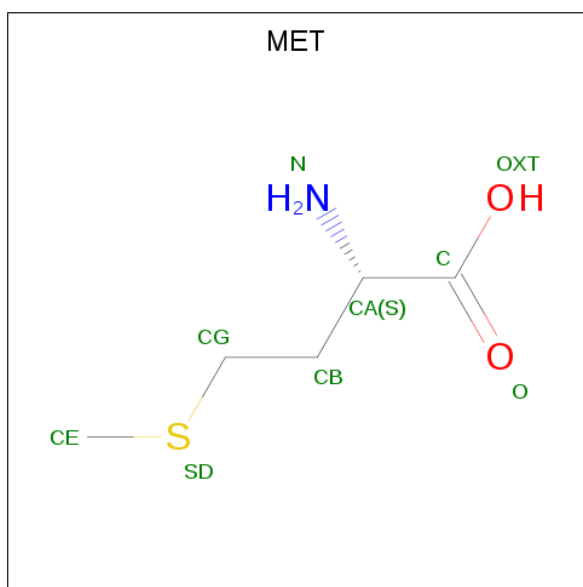
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf	
4	A	1	Total	Fe	S	0	0
			8	4	4		
4	C	1	Total	Fe	S	0	0
			8	4	4		
4	E	1	Total	Fe	S	0	0
			8	4	4		
4	G	1	Total	Fe	S	0	0
			8	4	4		
4	I	1	Total	Fe	S	0	0
			8	4	4		
4	K	1	Total	Fe	S	0	0
			8	4	4		

- Molecule 5 is 5'-DEOXYADENOSINE (three-letter code: 5AD) (formula:  $C_{10}H_{13}N_5O_3$ ).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
5	A	1	18	10	5	3	0	0
5	C	1	18	10	5	3	0	0
5	E	1	18	10	5	3	0	0
5	G	1	18	10	5	3	0	0
5	I	1	18	10	5	3	0	0
5	K	1	18	10	5	3	0	0

- Molecule 6 is METHIONINE (three-letter code: MET) (formula: C<sub>5</sub>H<sub>11</sub>NO<sub>2</sub>S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	S		
6	A	1	Total 9	5	1	2	1	0	0
6	C	1	Total 9	5	1	2	1	0	0
6	E	1	Total 9	5	1	2	1	0	0
6	G	1	Total 9	5	1	2	1	0	0
6	I	1	Total 9	5	1	2	1	0	0
6	K	1	Total 9	5	1	2	1	0	0

- Molecule 7 is IMIDAZOLE (three-letter code: IMD) (formula: C<sub>3</sub>H<sub>5</sub>N<sub>2</sub>).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	1	Total C N 5 3 2	0	0
7	E	1	Total C N 5 3 2	0	0

- Molecule 8 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	A	1	Total Cl 1 1	0	0
8	K	1	Total Cl 1 1	0	0

- Molecule 9 is ISOPROPYL ALCOHOL (three-letter code: IPA) (formula: C<sub>3</sub>H<sub>8</sub>O).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
9	C	1	Total C O 4 3 1	0	0

- Molecule 10 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
10	A	401	Total O 401 401	0	0
10	B	6	Total O 6 6	0	0
10	C	357	Total O 357 357	0	0
10	D	17	Total O 17 17	0	0
10	E	348	Total O 348 348	0	0
10	F	10	Total O 10 10	0	0
10	G	265	Total O 265 265	0	0
10	H	9	Total O 9 9	0	0
10	I	364	Total O 364 364	0	0
10	J	6	Total O 6 6	0	0
10	K	348	Total O 348 348	0	0

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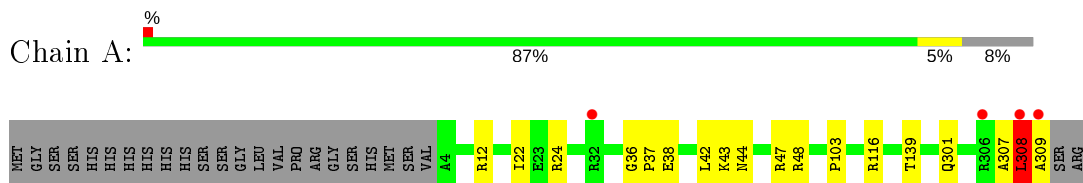
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<b>Mol</b>	<b>Chain</b>	<b>Residues</b>	<b>Atoms</b>		<b>ZeroOcc</b>	<b>AltConf</b>
10	L	8	Total	O	0	0
			8	8		

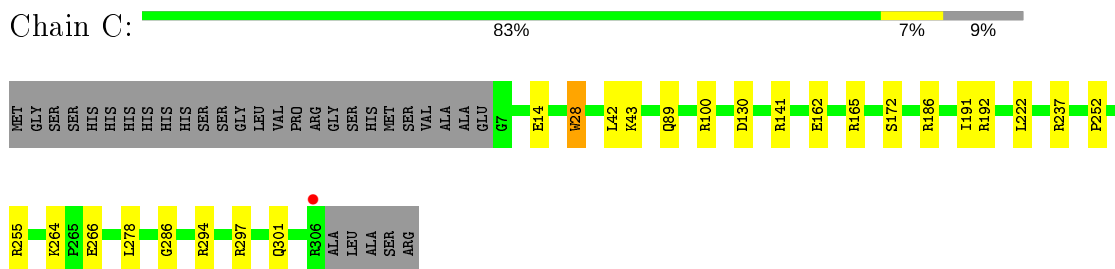
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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.

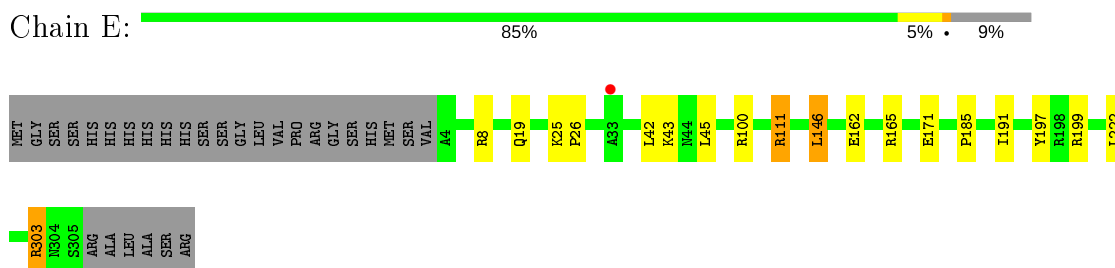
- Molecule 1: Lipoyl synthase



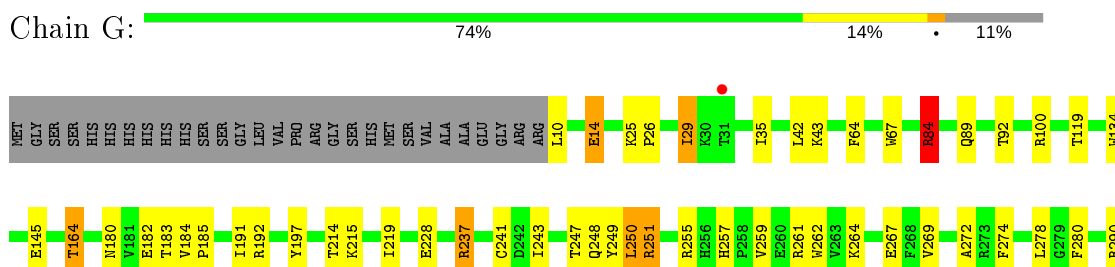
- Molecule 1: Lipoyl synthase

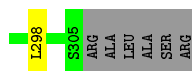


- Molecule 1: Lipoyl synthase

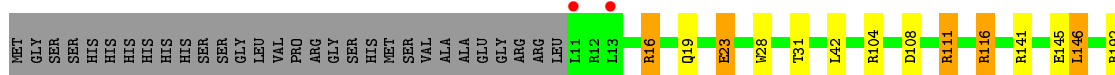
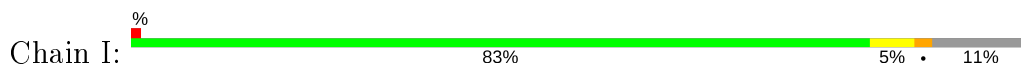


- Molecule 1: Lipoyl synthase

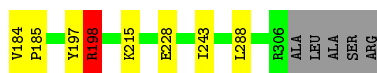
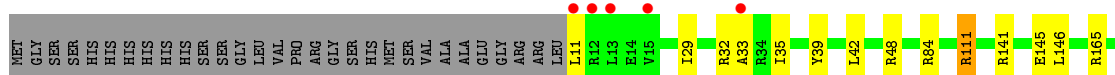
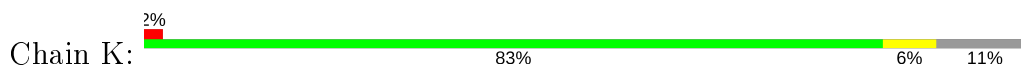




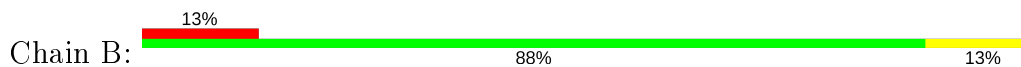
• Molecule 1: Lipoyl synthase



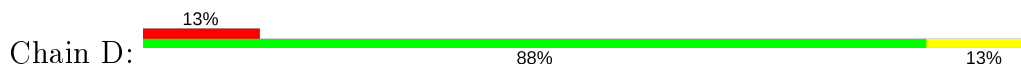
• Molecule 1: Lipoyl synthase



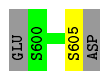
• Molecule 2: Octanoylated peptide from *M. tuberculosis* H protein



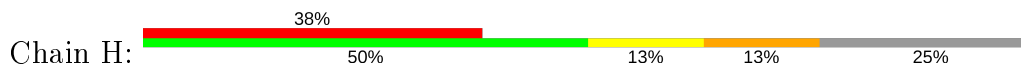
• Molecule 2: Octanoylated peptide from *M. tuberculosis* H protein



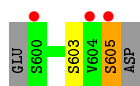
• Molecule 2: Octanoylated peptide from *M. tuberculosis* H protein



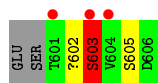
• Molecule 2: Octanoylated peptide from *M. tuberculosis* H protein



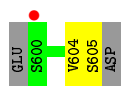




- Molecule 2: Octanoylated peptide from *M. tuberculosis* H protein



- Molecule 2: Octanoylated peptide from *M. tuberculosis* H protein



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	81.77Å 95.97Å 114.33Å 90.00° 90.46° 90.00°	Depositor
Resolution (Å)	47.03 – 1.86 47.03 – 1.86	Depositor EDS
% Data completeness (in resolution range)	96.8 (47.03-1.86) 90.3 (47.03-1.86)	Depositor EDS
$R_{merge}$	0.10	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	5.66 (at 1.86Å)	Xtrriage
Refinement program	PHENIX	Depositor
R, $R_{free}$	0.160 , 0.201 0.160 , 0.201	Depositor DCC
$R_{free}$ test set	7130 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	14.1	Xtrriage
Anisotropy	0.366	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 43.8	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	0.026 for h,-k,-l	Xtrriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	16721	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	19.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.71% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup> Intensities estimated from amplitudes.

<sup>2</sup> Theoretical 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: IPA, IMD, CL, SF4, 5AD, F3S, XOK

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	0.50	0/2437	0.61	2/3312 (0.1%)
1	C	0.43	0/2405	0.56	1/3270 (0.0%)
1	E	0.47	0/2438	0.59	0/3315
1	G	0.44	0/2338	0.60	1/3188 (0.0%)
1	I	0.44	0/2349	0.56	0/3197
1	K	0.44	0/2350	0.59	1/3198 (0.0%)
2	B	0.39	0/45	0.50	0/57
2	D	0.38	0/48	0.51	0/61
2	F	0.36	0/30	0.61	0/38
2	H	0.34	0/30	0.34	0/38
2	J	0.43	0/27	0.65	0/34
2	L	0.42	0/30	0.87	0/38
All	All	0.45	0/14527	0.59	5/19746 (0.0%)

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	C	0	1
1	I	0	1
All	All	0	2

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	308	LEU	CA-CB-CG	6.81	130.96	115.30
1	C	28	TRP	O-C-N	-6.40	112.46	122.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	36	GLY	C-N-CD	6.08	141.18	128.40
1	K	198	ARG	NE-CZ-NH1	6.01	123.31	120.30
1	G	84	ARG	NE-CZ-NH1	5.32	122.96	120.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	28	TRP	Mainchain
1	I	28	TRP	Mainchain

## 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	2381	0	2328	15	0
1	C	2346	0	2282	19	0
1	E	2385	0	2324	27	0
1	G	2282	0	2203	44	0
1	I	2290	0	2226	14	0
1	K	2300	0	2217	17	0
2	B	66	0	38	1	0
2	D	69	0	40	1	0
2	F	51	0	30	1	0
2	H	51	0	30	2	0
2	J	48	0	22	2	0
2	L	51	0	30	2	0
3	A	6	0	0	0	0
3	C	6	0	0	0	0
3	E	6	0	0	0	0
3	G	6	0	0	0	0
3	I	6	0	0	0	0
3	K	6	0	0	0	0
4	A	8	0	0	0	0
4	C	8	0	0	0	0
4	E	8	0	0	0	0
4	G	8	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	I	8	0	0	0	0
4	K	8	0	0	0	0
5	A	18	0	13	0	0
5	C	18	0	13	0	0
5	E	18	0	13	0	0
5	G	18	0	13	0	0
5	I	18	0	13	0	0
5	K	18	0	13	0	0
6	A	9	0	8	0	0
6	C	9	0	8	0	0
6	E	9	0	8	0	0
6	G	9	0	8	0	0
6	I	9	0	8	0	0
6	K	9	0	8	0	0
7	A	5	0	5	1	0
7	E	5	0	5	0	0
8	A	1	0	0	0	0
8	K	1	0	0	0	0
9	C	4	0	8	2	0
10	A	401	0	0	2	0
10	B	6	0	0	0	0
10	C	357	0	0	4	0
10	D	17	0	0	0	0
10	E	348	0	0	7	0
10	F	10	0	0	0	0
10	G	265	0	0	8	0
10	H	9	0	0	0	0
10	I	364	0	0	8	0
10	J	6	0	0	1	0
10	K	348	0	0	6	0
10	L	8	0	0	0	0
All	All	16721	0	13914	133	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 5.

All (133) 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:162[B]:GLU:OE2	1:E:165:ARG:CZ	1.83	1.24
1:E:162[B]:GLU:OE2	1:E:165:ARG:NH1	1.88	1.06

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:247[A]:THR:HG21	10:G:583:HOH:O	1.57	1.02
1:A:308:LEU:HG	1:A:309:ALA:H	1.32	0.94
1:G:248:GLN:NE2	10:G:501:HOH:O	2.06	0.87
1:E:162[B]:GLU:OE2	1:E:165:ARG:NE	2.09	0.84
1:E:162[B]:GLU:CD	1:E:165:ARG:NE	2.32	0.82
1:I:104:ARG:NH1	10:I:501:HOH:O	2.11	0.82
1:E:162[B]:GLU:HB2	1:E:165:ARG:HD2	1.62	0.82
1:G:261:ARG:NH2	1:G:267:GLU:OE1	2.17	0.78
1:K:84[B]:ARG:NH2	10:K:502:HOH:O	2.22	0.73
1:C:43:LYS:NZ	2:D:599:GLU:OE1	2.23	0.72
1:G:237:ARG:HG3	1:G:278:LEU:HB3	1.72	0.71
1:E:162[B]:GLU:CD	1:E:165:ARG:CZ	2.58	0.71
1:G:14:GLU:HG2	1:G:255:ARG:HB3	1.72	0.71
1:G:84:ARG:HG3	1:G:84:ARG:HH11	1.55	0.70
1:G:261:ARG:O	10:G:501:HOH:O	2.10	0.68
1:A:37:PRO:HG2	1:A:38:GLU:CD	2.13	0.68
1:G:84:ARG:NH1	1:G:92:THR:OG1	2.28	0.67
1:A:308:LEU:HG	1:A:309:ALA:N	2.09	0.66
1:C:14:GLU:OE1	1:C:255:ARG:NH1	2.30	0.64
1:E:162[B]:GLU:CD	1:E:165:ARG:HD2	2.18	0.62
1:E:199[C]:ARG:HD2	10:E:660:HOH:O	1.99	0.62
1:G:25:LYS:HG2	1:G:29:ILE:HD11	1.81	0.62
1:E:162[B]:GLU:CD	1:E:165:ARG:CD	2.69	0.60
1:E:162[B]:GLU:OE2	1:E:165:ARG:CD	2.49	0.60
1:E:162[B]:GLU:OE1	1:E:165:ARG:NE	2.34	0.60
1:G:259:VAL:N	10:G:509:HOH:O	2.34	0.60
2:J:602:XOK:O	2:J:603:SER:HB2	2.00	0.59
1:K:198:ARG:HH11	1:K:198:ARG:HG2	1.67	0.59
1:G:264:LYS:HB2	1:G:267:GLU:HG3	1.84	0.59
1:E:19:GLN:NE2	10:E:506:HOH:O	2.36	0.58
1:K:48:ARG:NH1	10:K:512:HOH:O	2.37	0.56
1:I:208[B]:ARG:NH2	10:I:507:HOH:O	2.38	0.56
1:E:199[C]:ARG:CD	10:E:660:HOH:O	2.54	0.56
1:G:100:ARG:HG3	1:G:134:TRP:CH2	2.40	0.55
1:C:172:SER:HA	9:C:405:IPA:H13	1.88	0.55
1:A:43:LYS:NZ	2:B:599:GLU:OE1	2.29	0.55
1:G:219:ILE:HG22	1:G:249:TYR:HB2	1.89	0.55
1:G:84:ARG:HH11	1:G:84:ARG:CG	2.18	0.55
1:C:100:ARG:NH1	10:C:504:HOH:O	2.39	0.54
1:K:111:ARG:NH2	10:K:513:HOH:O	2.39	0.54
1:G:274:PHE:CE1	1:G:278:LEU:HD11	2.42	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:162[A]:GLU:HB3	1:E:165:ARG:HD2	1.88	0.54
1:I:16:ARG:NH2	10:I:508:HOH:O	2.42	0.53
1:E:162[B]:GLU:CB	1:E:165:ARG:HD2	2.36	0.53
1:C:165[B]:ARG:NH1	10:C:507:HOH:O	2.41	0.53
1:C:162:GLU:HB3	1:C:165[B]:ARG:HG2	1.90	0.53
1:K:165[B]:ARG:NH2	10:K:511:HOH:O	2.36	0.52
1:G:250:LEU:HD21	2:H:603:SER:O	2.09	0.52
1:G:237:ARG:HG2	1:G:280:PHE:CE1	2.45	0.52
1:E:303:ARG:NH1	10:E:511:HOH:O	2.41	0.52
1:K:11:LEU:N	10:K:518:HOH:O	2.42	0.51
1:C:294:ARG:NH2	10:C:510:HOH:O	2.43	0.51
1:G:261:ARG:NH1	10:G:515:HOH:O	2.42	0.51
1:I:16:ARG:NH2	10:I:511:HOH:O	2.43	0.51
1:G:259:VAL:HG23	10:G:509:HOH:O	2.11	0.50
1:G:185:PRO:HD3	1:G:197:TYR:CE1	2.47	0.50
1:C:191[B]:ILE:HD13	1:C:222:LEU:HD13	1.94	0.49
1:K:198:ARG:CG	1:K:198:ARG:HH11	2.25	0.49
1:C:237[B]:ARG:HD3	1:C:278:LEU:O	2.11	0.49
1:I:116:ARG:NH1	10:I:516:HOH:O	2.46	0.49
1:K:29:ILE:O	2:L:605:SER:HB2	2.12	0.49
1:E:185:PRO:HG3	1:E:197:TYR:CD2	2.48	0.49
1:K:32:ARG:O	10:K:501:HOH:O	2.20	0.49
1:A:37:PRO:HG2	1:A:38:GLU:H	1.78	0.48
1:G:247[A]:THR:HG23	1:G:290:ARG:NH2	2.28	0.48
1:I:111:ARG:HD2	1:I:146:LEU:HD12	1.93	0.48
1:E:199[B]:ARG:NH2	10:E:510:HOH:O	2.41	0.48
1:C:264:LYS:HB3	1:C:266[B]:GLU:OE2	2.14	0.48
1:G:259:VAL:HG11	1:G:262:TRP:CZ2	2.48	0.48
1:I:141:ARG:NH1	10:I:502:HOH:O	2.26	0.48
1:A:22:ILE:HG13	1:G:145:GLU:HG2	1.96	0.48
1:A:47:ARG:HH12	1:E:146:LEU:HD23	1.79	0.48
1:K:35:ILE:HG23	1:K:39:TYR:CD2	2.49	0.48
1:G:183:THR:OG1	1:G:184:VAL:N	2.47	0.47
1:A:307:ALA:O	1:A:308:LEU:HB3	2.14	0.47
1:G:261:ARG:HH21	1:G:267:GLU:CD	2.15	0.47
1:C:266[B]:GLU:CD	1:C:266[B]:GLU:H	2.17	0.47
1:A:37:PRO:HG2	1:A:38:GLU:N	2.30	0.47
1:C:89:GLN:HB3	1:C:252:PRO:HB3	1.95	0.47
1:E:191[B]:ILE:HD13	1:E:222:LEU:HD13	1.96	0.46
1:G:237:ARG:NH1	1:G:241:CYS:O	2.43	0.46
1:G:191:ILE:HG12	10:G:504:HOH:O	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:22:ILE:CG1	1:G:145:GLU:HG2	2.45	0.46
1:A:37:PRO:HD2	1:A:38:GLU:OE2	2.15	0.46
1:G:25:LYS:HG3	1:G:262:TRP:CH2	2.51	0.46
1:C:286:GLY:CA	1:G:164[A]:THR:HG21	2.46	0.46
1:I:266:GLU:CD	1:I:266:GLU:H	2.19	0.46
1:G:29:ILE:O	1:G:29:ILE:HG13	2.16	0.45
1:K:184:VAL:HG11	1:K:228:GLU:HB3	1.99	0.45
1:G:191:ILE:HD11	1:G:257:HIS:CD2	2.52	0.45
1:G:25:LYS:HB3	2:H:605:SER:HB3	1.98	0.45
1:G:214:THR:OG1	1:G:241:CYS:HA	2.17	0.45
1:E:111:ARG:O	1:E:111:ARG:HD3	2.17	0.45
1:G:35:ILE:HG12	1:G:64:PHE:CE1	2.52	0.45
7:A:405:IMD:H5	10:A:798:HOH:O	2.17	0.44
1:G:184:VAL:HG11	1:G:228:GLU:HB3	1.99	0.44
1:I:141:ARG:O	1:I:145:GLU:HG3	2.18	0.44
1:C:237[B]:ARG:NH1	10:C:513:HOH:O	2.49	0.44
1:E:26:PRO:O	2:F:605:SER:OG	2.35	0.44
1:A:44:ASN:O	1:A:48:ARG:HB2	2.18	0.43
1:I:219:ILE:HG22	1:I:249:TYR:HB2	2.00	0.43
1:A:301:GLN:NE2	10:A:509:HOH:O	2.40	0.43
1:E:8:ARG:NH1	10:E:512:HOH:O	2.42	0.43
1:I:23:GLU:H	1:I:23:GLU:CD	2.22	0.43
1:K:141:ARG:O	1:K:145:GLU:HG2	2.18	0.43
1:E:199[C]:ARG:HD3	1:E:199[C]:ARG:HA	1.59	0.43
1:K:185:PRO:HD3	1:K:197:TYR:CE1	2.54	0.43
1:G:269:VAL:O	1:G:272:ALA:HB3	2.19	0.42
1:C:186:ARG:HD2	1:I:277:GLY:O	2.18	0.42
1:I:297:ARG:NH1	10:I:526:HOH:O	2.51	0.42
1:K:215:LYS:HA	1:K:243:ILE:O	2.18	0.42
1:C:286:GLY:HA2	1:G:164[A]:THR:HG21	2.00	0.42
1:E:100:ARG:NH1	10:E:524:HOH:O	2.52	0.42
1:G:26:PRO:HD3	1:G:262:TRP:CD2	2.55	0.42
1:C:141:ARG:HD3	9:C:405:IPA:H2	2.01	0.42
1:G:251:ARG:HD2	10:G:509:HOH:O	2.20	0.42
1:C:294:ARG:HH21	1:C:297:ARG:CZ	2.33	0.42
1:G:215:LYS:HA	1:G:243:ILE:O	2.20	0.42
1:G:84:ARG:HG3	1:G:84:ARG:NH1	2.25	0.41
1:K:146:LEU:HA	1:K:146:LEU:HD12	1.91	0.41
1:K:29:ILE:O	2:L:604:VAL:HG13	2.20	0.41
1:E:162[B]:GLU:OE2	1:E:165:ARG:HD2	2.18	0.41
1:G:43:LYS:HG2	1:G:67:TRP:HZ3	1.84	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:29:ILE:HG23	1:K:288:LEU:HD11	2.02	0.41
1:A:103:PRO:HB3	1:A:139:THR:HA	2.03	0.41
1:A:24:ARG:HE	1:A:24:ARG:HB2	1.57	0.41
1:C:130:ASP:HA	1:C:165[B]:ARG:CZ	2.51	0.41
1:I:208[B]:ARG:NE	10:I:534:HOH:O	2.53	0.41
2:J:603:SER:N	10:J:702:HOH:O	2.53	0.41
1:E:162[A]:GLU:OE1	1:E:165:ARG:HG3	2.21	0.40
1:G:180:ASN:HB3	1:G:182:GLU:CD	2.42	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	308/331 (93%)	303 (98%)	4 (1%)	1 (0%)	41	26
1	C	305/331 (92%)	301 (99%)	4 (1%)	0	100	100
1	E	308/331 (93%)	302 (98%)	6 (2%)	0	100	100
1	G	300/331 (91%)	293 (98%)	7 (2%)	0	100	100
1	I	298/331 (90%)	294 (99%)	4 (1%)	0	100	100
1	K	298/331 (90%)	291 (98%)	6 (2%)	1 (0%)	41	26
2	B	5/8 (62%)	4 (80%)	1 (20%)	0	100	100
2	D	5/8 (62%)	4 (80%)	1 (20%)	0	100	100
2	F	3/8 (38%)	2 (67%)	1 (33%)	0	100	100
2	H	3/8 (38%)	2 (67%)	1 (33%)	0	100	100
2	J	3/8 (38%)	2 (67%)	0	1 (33%)	0	0
2	L	3/8 (38%)	3 (100%)	0	0	100	100
All	All	1839/2034 (90%)	1801 (98%)	35 (2%)	3 (0%)	47	33

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	308	LEU
1	K	33	ALA
2	J	603	SER

### 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	242/272 (89%)	238 (98%)	4 (2%)	60	47
1	C	239/272 (88%)	236 (99%)	3 (1%)	69	58
1	E	243/272 (89%)	235 (97%)	8 (3%)	38	21
1	G	233/272 (86%)	218 (94%)	15 (6%)	17	4
1	I	236/272 (87%)	225 (95%)	11 (5%)	26	10
1	K	233/272 (86%)	230 (99%)	3 (1%)	69	58
2	B	6/7 (86%)	6 (100%)	0	100	100
2	D	7/7 (100%)	7 (100%)	0	100	100
2	F	5/7 (71%)	5 (100%)	0	100	100
2	H	5/7 (71%)	4 (80%)	1 (20%)	1	0
2	J	3/7 (43%)	1 (33%)	2 (67%)	0	0
2	L	5/7 (71%)	5 (100%)	0	100	100
All	All	1457/1674 (87%)	1410 (97%)	47 (3%)	40	22

All (47) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	12	ARG
1	A	42	LEU
1	A	116	ARG
1	A	308	LEU
1	C	42	LEU
1	C	192	ARG

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Mol	Chain	Res	Type
1	C	301	GLN
1	E	25	LYS
1	E	42	LEU
1	E	43	LYS
1	E	45	LEU
1	E	111	ARG
1	E	146	LEU
1	E	171	GLU
1	E	303	ARG
1	G	10	LEU
1	G	14	GLU
1	G	29	ILE
1	G	42	LEU
1	G	84	ARG
1	G	89	GLN
1	G	119[A]	THR
1	G	119[B]	THR
1	G	164[A]	THR
1	G	164[B]	THR
1	G	192	ARG
1	G	237	ARG
1	G	250	LEU
1	G	251	ARG
1	G	298	LEU
2	H	605	SER
1	I	16	ARG
1	I	19	GLN
1	I	23	GLU
1	I	31	THR
1	I	42	LEU
1	I	108	ASP
1	I	111	ARG
1	I	116	ARG
1	I	146	LEU
1	I	192	ARG
1	I	288	LEU
2	J	603	SER
2	J	605	SER
1	K	42	LEU
1	K	111	ARG
1	K	198	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such

sidechains are listed below:

Mol	Chain	Res	Type
1	E	19	GLN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

6 non-standard protein/DNA/RNA residues 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	XOK	B	602	3,2	16,18,19	1.09	1 (6%)	12,20,22	1.40	1 (8%)
2	XOK	F	602	3,2	16,18,19	1.23	1 (6%)	12,20,22	1.72	1 (8%)
2	XOK	D	602	3,2	16,18,19	1.18	1 (6%)	12,20,22	1.73	1 (8%)
2	XOK	J	602	3,2	16,18,19	1.19	1 (6%)	12,20,22	1.60	2 (16%)
2	XOK	H	602	3,2	16,18,19	1.27	1 (6%)	12,20,22	1.66	1 (8%)
2	XOK	L	602	3,2	16,18,19	1.16	1 (6%)	12,20,22	1.76	1 (8%)

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	XOK	B	602	3,2	-	1/17/19/21	-
2	XOK	F	602	3,2	-	0/17/19/21	-
2	XOK	D	602	3,2	-	1/17/19/21	-
2	XOK	J	602	3,2	-	3/17/19/21	-
2	XOK	H	602	3,2	-	1/17/19/21	-
2	XOK	L	602	3,2	-	0/17/19/21	-

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	H	602	XOK	C1-NZ	3.89	1.42	1.33
2	L	602	XOK	C1-NZ	3.77	1.42	1.33
2	D	602	XOK	C1-NZ	3.66	1.41	1.33
2	F	602	XOK	C1-NZ	3.50	1.41	1.33
2	J	602	XOK	C1-NZ	3.44	1.41	1.33
2	B	602	XOK	C1-NZ	3.10	1.40	1.33

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	602	XOK	CE-NZ-C1	-5.53	112.57	122.84
2	L	602	XOK	CE-NZ-C1	-5.41	112.79	122.84
2	F	602	XOK	CE-NZ-C1	-5.34	112.92	122.84
2	H	602	XOK	CE-NZ-C1	-5.22	113.14	122.84
2	B	602	XOK	CE-NZ-C1	-4.44	114.58	122.84
2	J	602	XOK	CE-NZ-C1	-4.10	115.23	122.84
2	J	602	XOK	CD-CE-NZ	-3.01	103.60	112.21

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	J	602	XOK	N-CA-CB-CG
2	J	602	XOK	C-CA-CB-CG
2	J	602	XOK	CA-CB-CG-CD
2	H	602	XOK	C4-C5-C6-C7
2	B	602	XOK	CG-CD-CE-NZ
2	D	602	XOK	N-CA-CB-CG

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	J	602	XOK	1	0

## 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

Of 29 ligands modelled in this entry, 2 are monoatomic - leaving 27 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	F3S	G	401	1,2	0,7,9	0.00	-	-		
4	SF4	C	402	1,6	0,12,12	0.00	-	-		
3	F3S	A	401	1,2	0,7,9	0.00	-	-		
3	F3S	C	401	1,2	0,7,9	0.00	-	-		
5	5AD	A	403	-	17,20,20	0.62	0	15,30,30	0.89	1 (6%)
4	SF4	E	402	1,6	0,12,12	0.00	-	-		
4	SF4	K	402	1,6	0,12,12	0.00	-	-		
4	SF4	G	402	1,6	0,12,12	0.00	-	-		
5	5AD	E	403	-	17,20,20	0.65	0	15,30,30	0.82	1 (6%)
5	5AD	G	403	-	17,20,20	0.64	0	15,30,30	0.92	1 (6%)
5	5AD	I	403	-	17,20,20	0.66	0	15,30,30	0.86	1 (6%)
5	5AD	K	403	-	17,20,20	0.66	0	15,30,30	0.93	1 (6%)
5	5AD	C	403	-	17,20,20	0.58	0	15,30,30	0.85	1 (6%)
4	SF4	I	402	1,6	0,12,12	0.00	-	-		
9	IPA	C	405	-	3,3,3	0.53	0	3,3,3	0.40	0
7	IMD	E	405	-	3,5,5	0.32	0	4,5,5	0.61	0
7	IMD	A	405	-	3,5,5	0.37	0	4,5,5	0.59	0
3	F3S	I	401	1,2	0,7,9	0.00	-	-		
3	F3S	K	401	1,2	0,7,9	0.00	-	-		
4	SF4	A	402	1,6	0,12,12	0.00	-	-		
3	F3S	E	401	1,2	0,7,9	0.00	-	-		

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 Torsions and Rings columns. '2' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	F3S	G	401	1,2	-	-	0/2/2/3
4	SF4	C	402	1,6	-	-	0/6/5/5

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	F3S	A	401	1,2	-	-	0/2/2/3
3	F3S	C	401	1,2	-	-	0/2/2/3
4	SF4	E	402	1,6	-	-	0/6/5/5
5	5AD	A	403	-	-	0/0/20/20	0/3/3/3
5	5AD	C	403	-	-	0/0/20/20	0/3/3/3
4	SF4	G	402	1,6	-	-	0/6/5/5
5	5AD	E	403	-	-	0/0/20/20	0/3/3/3
5	5AD	G	403	-	-	0/0/20/20	0/3/3/3
5	5AD	I	403	-	-	0/0/20/20	0/3/3/3
5	5AD	K	403	-	-	0/0/20/20	0/3/3/3
4	SF4	K	402	1,6	-	-	0/6/5/5
4	SF4	I	402	1,6	-	-	0/6/5/5
7	IMD	E	405	-	-	-	0/1/1/1
7	IMD	A	405	-	-	-	0/1/1/1
3	F3S	I	401	1,2	-	-	0/2/2/3
3	F3S	K	401	1,2	-	-	0/2/2/3
4	SF4	A	402	1,6	-	-	0/6/5/5
3	F3S	E	401	1,2	-	-	0/2/2/3

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	K	403	5AD	C5-C6-N6	2.33	123.89	120.35
5	A	403	5AD	C5-C6-N6	2.23	123.75	120.35
5	G	403	5AD	C5-C6-N6	2.23	123.74	120.35
5	C	403	5AD	C5-C6-N6	2.15	123.62	120.35
5	I	403	5AD	C5-C6-N6	2.10	123.54	120.35
5	E	403	5AD	C5-C6-N6	2.01	123.41	120.35

There are no chirality outliers.

There are no torsion outliers.

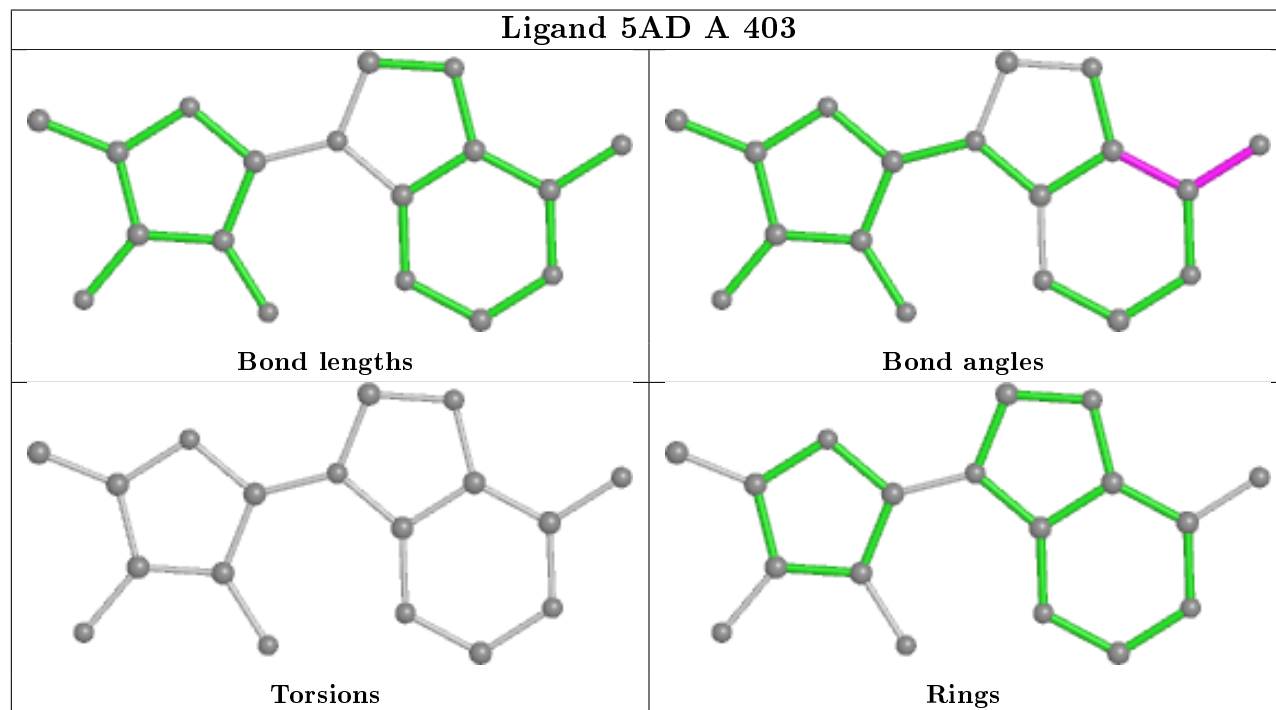
There are no ring outliers.

2 monomers are involved in 3 short contacts:

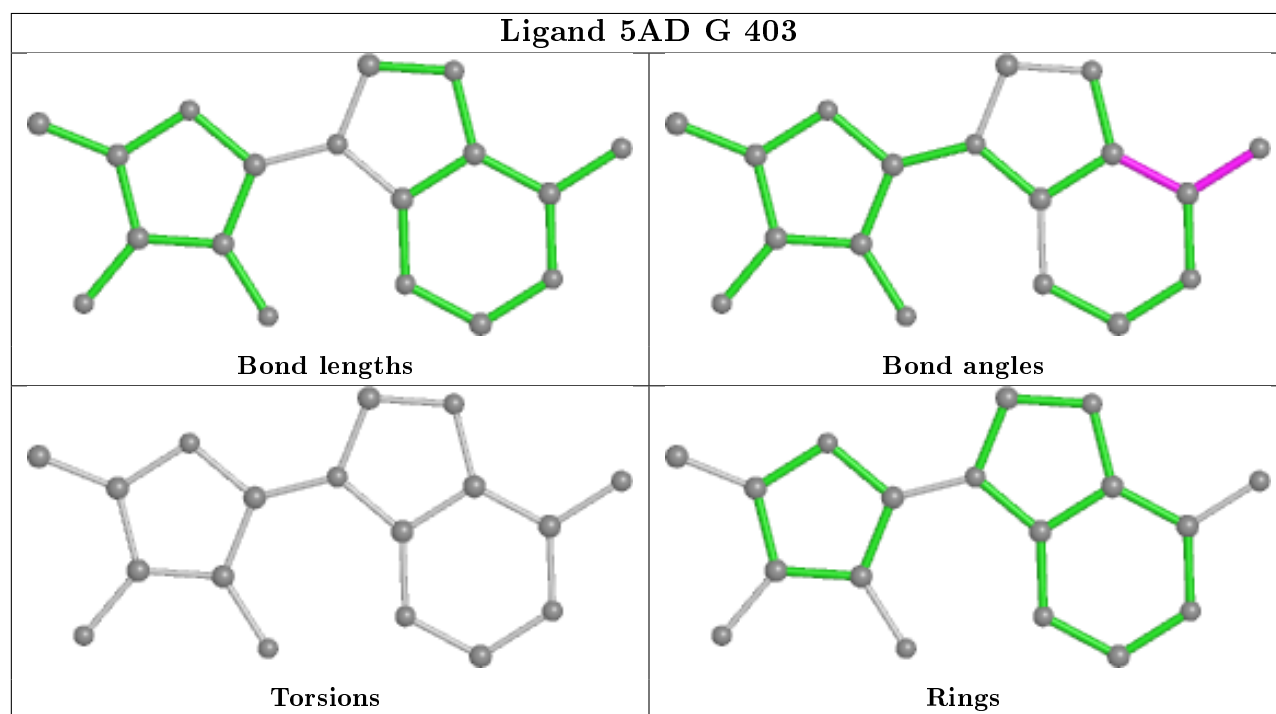
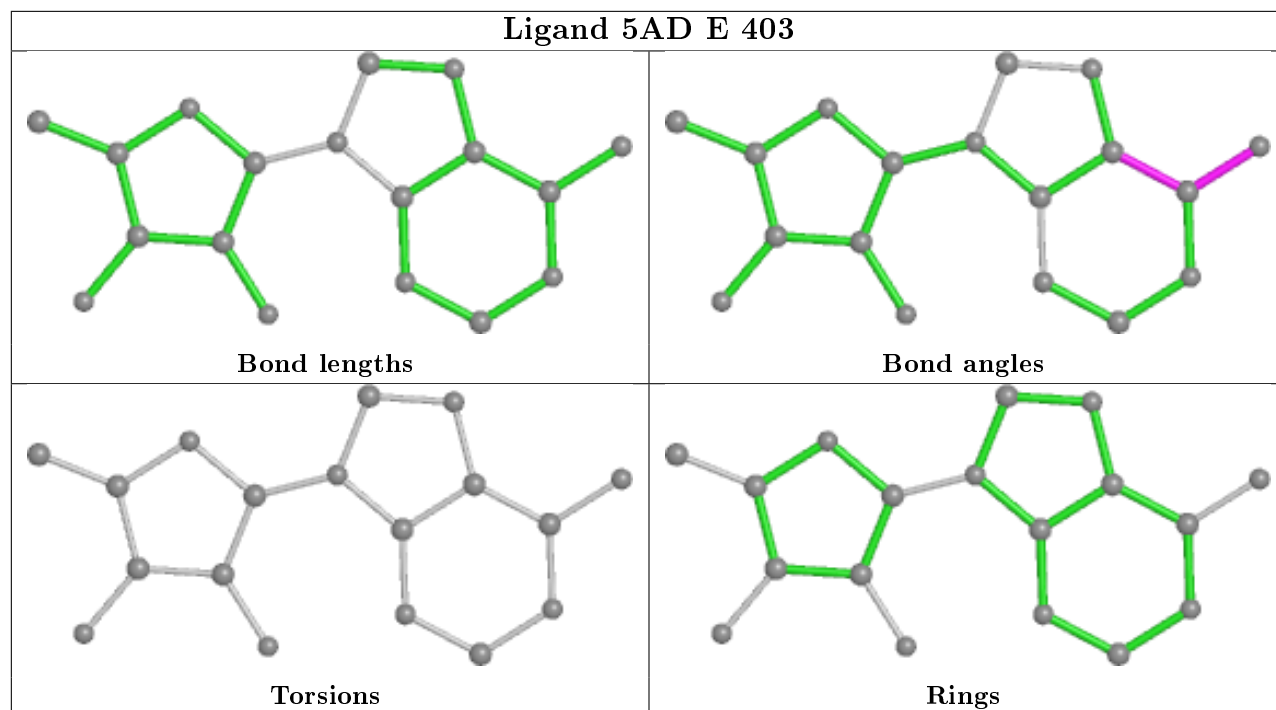
Mol	Chain	Res	Type	Clashes	Symm-Clashes
9	C	405	IPA	2	0
7	A	405	IMD	1	0

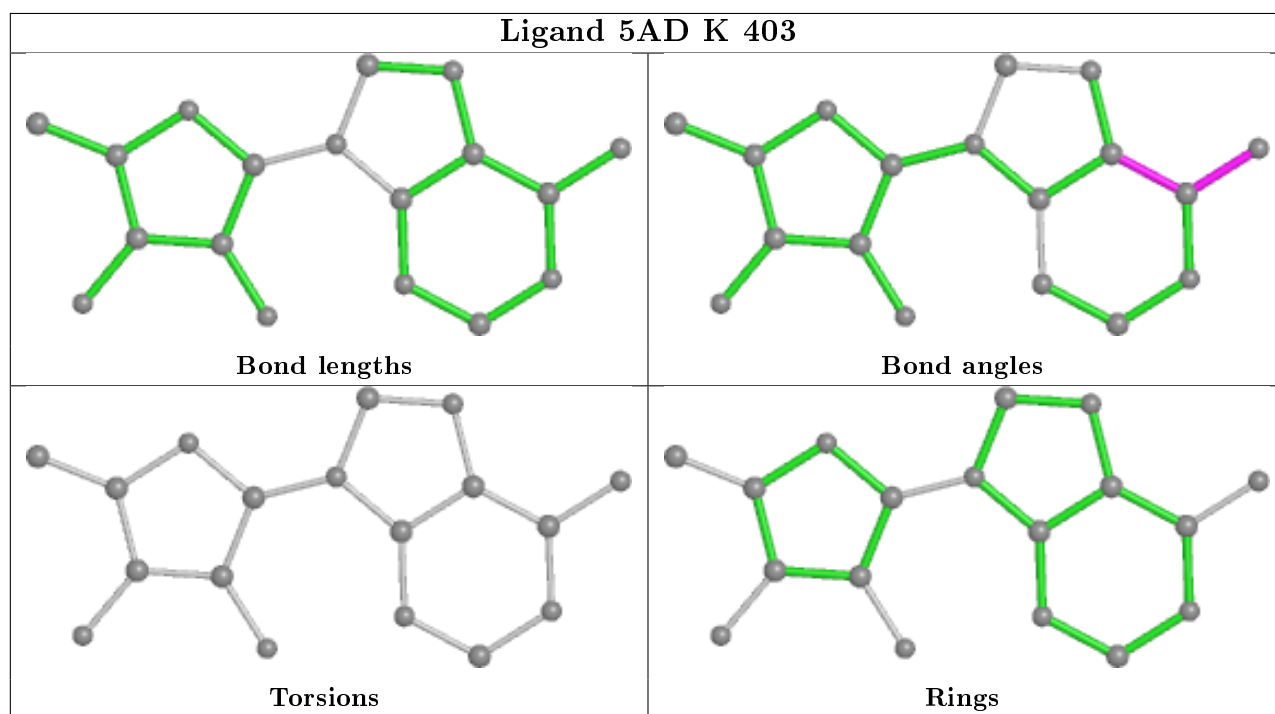
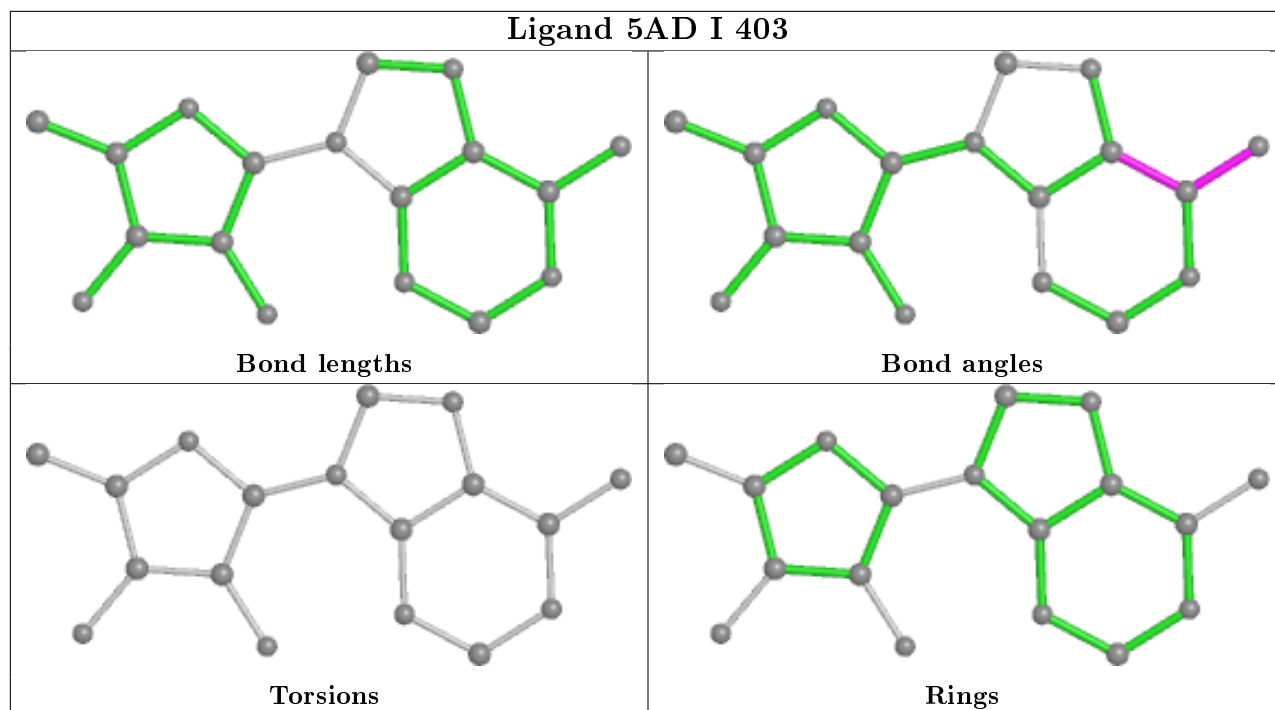
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In

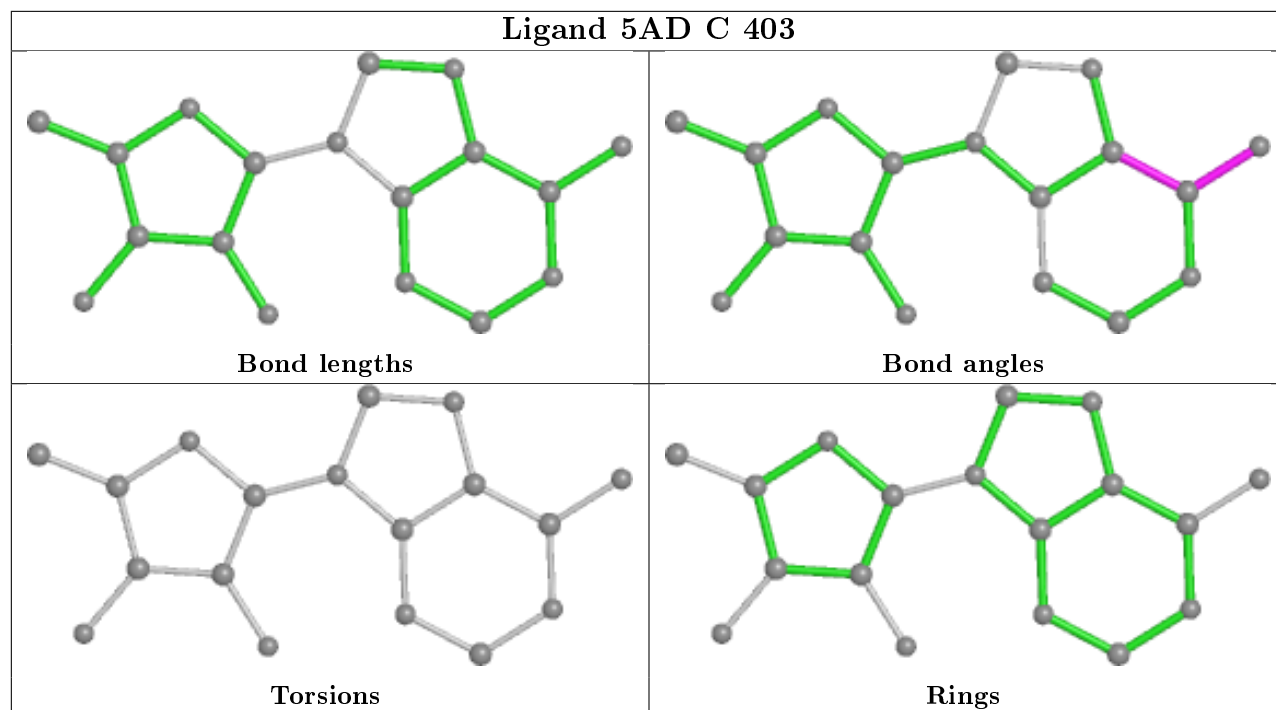
addition, ligands with molecular weight  $> 250$  and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. 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 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	306/331 (92%)	-0.48	4 (1%) 77 78	7, 14, 33, 53	0
1	C	300/331 (90%)	-0.56	1 (0%) 94 93	8, 13, 32, 56	0
1	E	302/331 (91%)	-0.46	1 (0%) 94 93	8, 15, 34, 47	0
1	G	296/331 (89%)	-0.05	1 (0%) 94 93	10, 22, 39, 62	0
1	I	296/331 (89%)	-0.46	2 (0%) 87 88	8, 15, 32, 44	0
1	K	296/331 (89%)	-0.43	5 (1%) 70 70	9, 16, 37, 50	0
2	B	7/8 (87%)	0.78	1 (14%) 2 2	26, 27, 43, 51	0
2	D	7/8 (87%)	0.66	1 (14%) 2 2	20, 23, 37, 43	0
2	F	5/8 (62%)	0.39	0 100 100	24, 27, 34, 42	0
2	H	5/8 (62%)	2.13	3 (60%) 0 0	33, 40, 46, 51	0
2	J	5/8 (62%)	2.34	3 (60%) 0 0	44, 50, 51, 53	0
2	L	5/8 (62%)	0.67	1 (20%) 1 1	27, 29, 39, 44	0
All	All	1830/2034 (89%)	-0.38	23 (1%) 77 78	7, 16, 37, 62	0

All (23) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	33	ALA	4.8
1	K	11	LEU	4.0
1	G	31	THR	3.8
1	A	309	ALA	3.7
2	H	605	SER	3.5
1	A	32	ARG	3.3
2	J	601	THR	3.3
2	J	604	VAL	3.0
2	B	599	GLU	2.9
1	K	15	VAL	2.9
1	A	306	ARG	2.9

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Mol	Chain	Res	Type	RSRZ
1	I	11	LEU	2.8
2	J	603	SER	2.7
1	K	33	ALA	2.5
2	H	604	VAL	2.4
2	D	599	GLU	2.4
1	C	306	ARG	2.4
1	A	308	LEU	2.1
2	H	600	SER	2.1
2	L	600	SER	2.1
1	K	13	LEU	2.1
1	I	13	LEU	2.0
1	K	12	ARG	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains [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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
2	XOK	J	602	19/20	0.96	0.11	7,10,45,48	0
2	XOK	H	602	19/20	0.96	0.10	13,20,28,34	0
2	XOK	D	602	19/20	0.97	0.08	6,10,16,19	0
2	XOK	B	602	19/20	0.97	0.09	5,11,24,25	0
2	XOK	F	602	19/20	0.97	0.08	6,10,20,24	0
2	XOK	L	602	19/20	0.98	0.07	10,12,22,24	0

## 6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 6.4 Ligands [i](#)

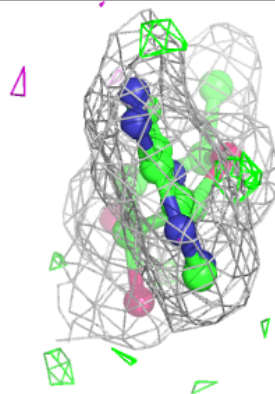
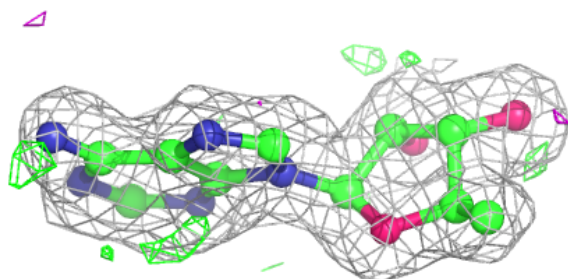
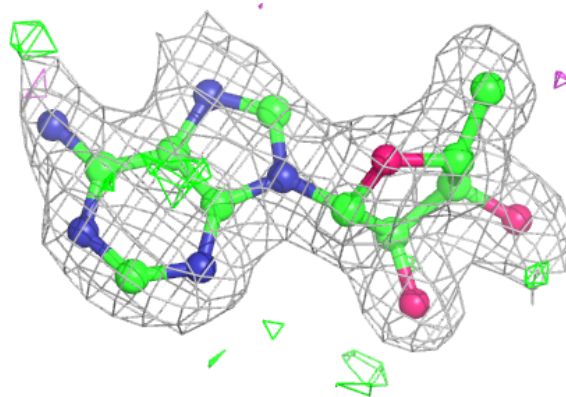
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
9	IPA	C	405	4/4	0.77	0.44	30,37,40,44	0
7	IMD	E	405	5/5	0.92	0.10	13,14,17,17	5
5	5AD	G	403	18/18	0.95	0.08	12,22,32,34	0
6	MET	G	404	9/9	0.96	0.11	10,11,13,14	0
5	5AD	A	403	18/18	0.97	0.07	7,9,12,13	0
4	SF4	G	402	8/8	0.97	0.07	12,14,16,16	0
5	5AD	E	403	18/18	0.98	0.06	6,9,11,12	0
5	5AD	K	403	18/18	0.98	0.06	8,11,12,13	0
5	5AD	I	403	18/18	0.98	0.06	5,9,11,12	0
6	MET	I	404	9/9	0.98	0.07	8,9,10,10	0
7	IMD	A	405	5/5	0.98	0.06	10,12,15,16	0
6	MET	K	404	9/9	0.98	0.07	9,11,13,14	0
3	F3S	G	401	6/7	0.98	0.05	17,17,17,18	0
5	5AD	C	403	18/18	0.98	0.07	6,9,10,11	0
8	CL	A	406	1/1	0.99	0.05	27,27,27,27	0
6	MET	C	404	9/9	0.99	0.07	5,7,10,10	0
6	MET	E	404	9/9	0.99	0.07	6,8,10,12	0
8	CL	K	405	1/1	0.99	0.04	22,22,22,22	0
4	SF4	K	402	8/8	0.99	0.06	12,13,14,14	0
3	F3S	K	401	6/7	0.99	0.06	11,12,13,13	0
4	SF4	I	402	8/8	0.99	0.06	10,11,13,14	0
6	MET	A	404	9/9	0.99	0.07	5,8,10,11	0
3	F3S	A	401	6/7	1.00	0.06	10,11,11,12	0
4	SF4	A	402	8/8	1.00	0.06	9,10,11,11	0
4	SF4	C	402	8/8	1.00	0.06	8,9,10,10	0
3	F3S	C	401	6/7	1.00	0.05	9,10,11,11	0
4	SF4	E	402	8/8	1.00	0.06	9,10,11,11	0
3	F3S	I	401	6/7	1.00	0.06	9,10,11,12	0
3	F3S	E	401	6/7	1.00	0.05	10,12,13,14	0

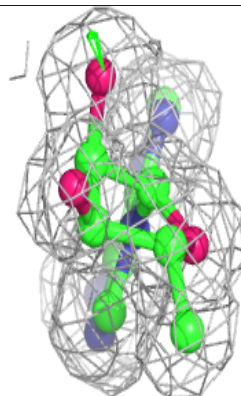
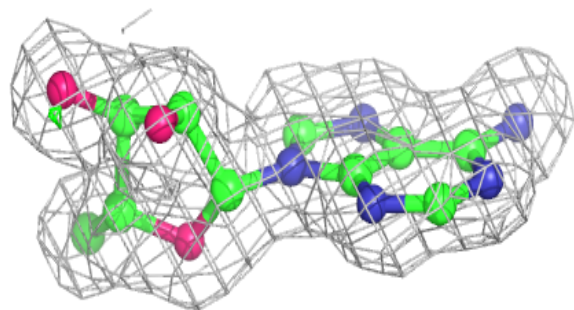
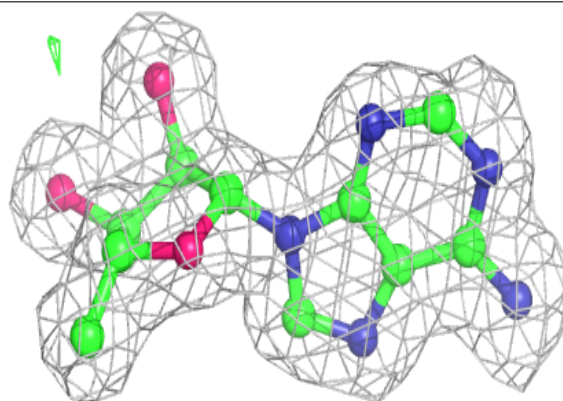
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

**Electron density around 5AD G 403:**

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

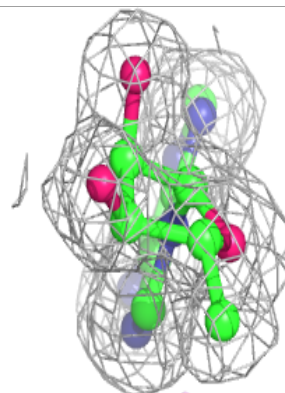
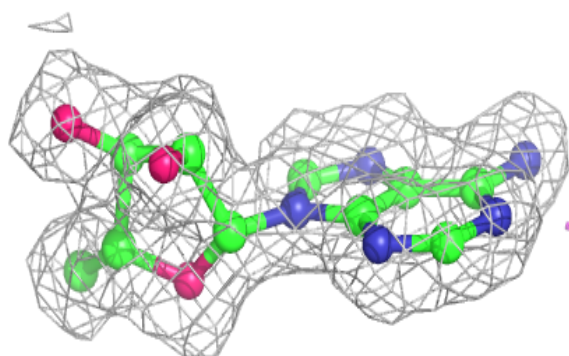
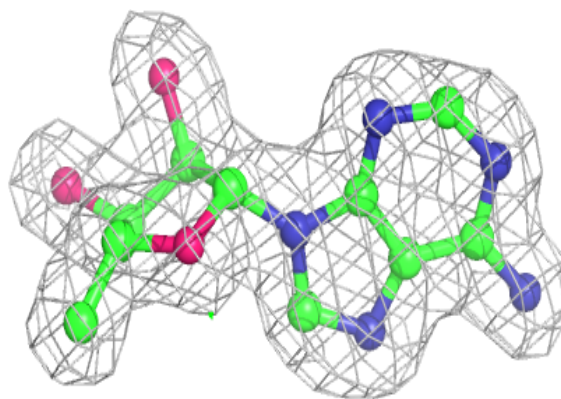
**Electron density around 5AD A 403:**

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

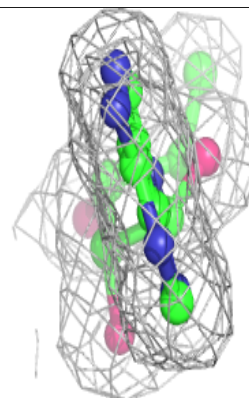
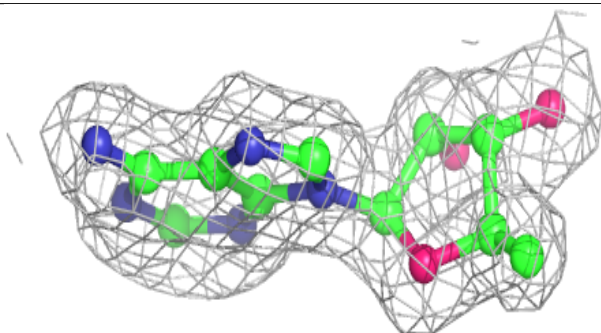
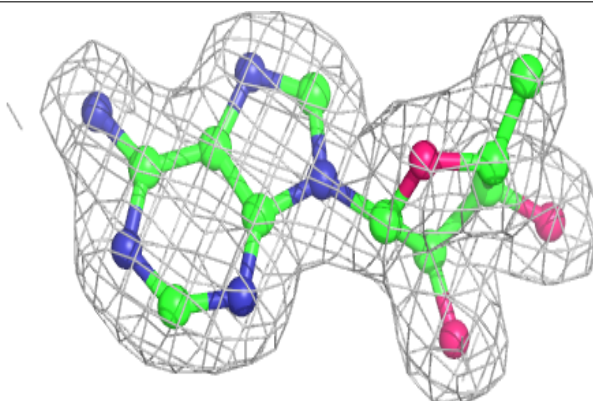


**Electron density around 5AD E 403:**

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

**Electron density around 5AD K 403:**

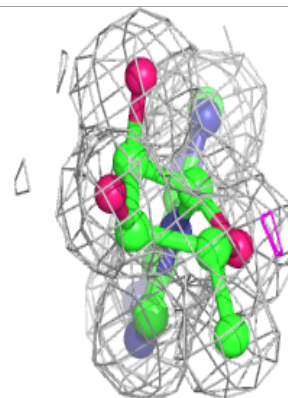
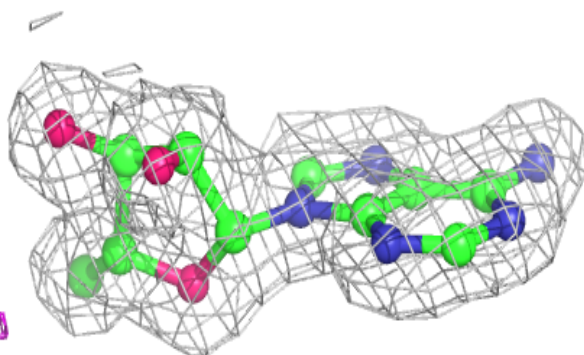
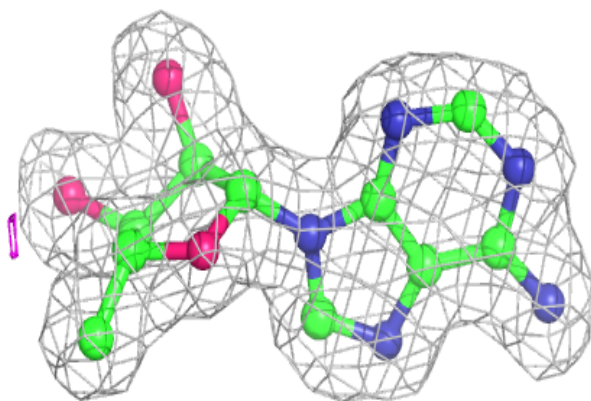
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



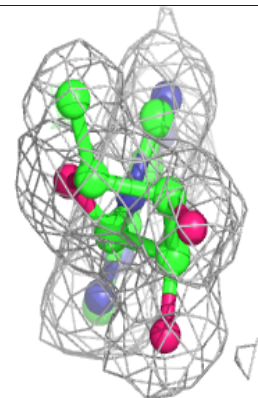
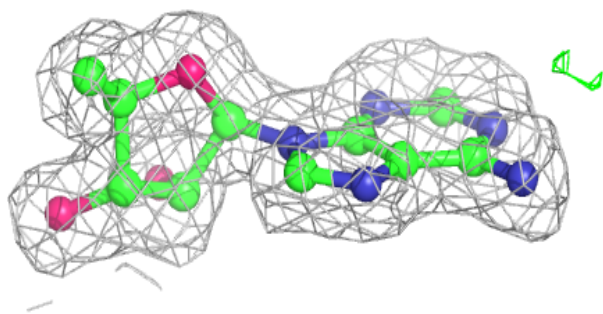
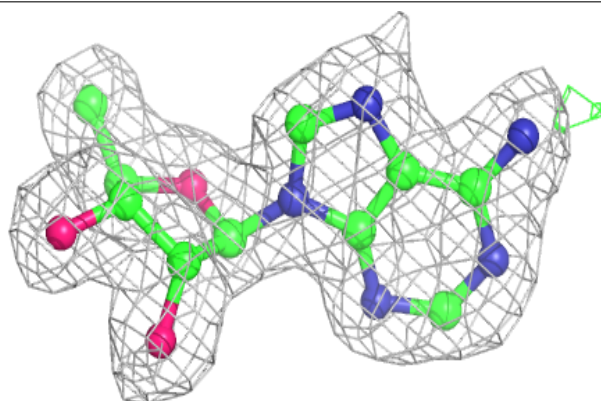


**Electron density around 5AD I 403:**

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

**Electron density around 5AD C 403:**

$2mF_o-DF_c$  (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.