



# Full wwPDB X-ray Structure Validation Report ⓘ

Apr 7, 2022 – 06:02 PM EDT

PDB ID : 3MLE  
Title : Crystal structure of dethiobiotin synthetase (BioD) from *Helicobacter pylori* cocrystallized with ATP  
Authors : Nicholls, R.; Porebski, P.J.; Klimecka, M.M.; Chruszcz, M.; Murzyn, K.; Joachimiak, A.; Murshudov, G.; Minor, W.; Midwest Center for Structural Genomics (MCSG)  
Deposited on : 2010-04-16  
Resolution : 2.80 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 : 2.27  
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) : 2.27

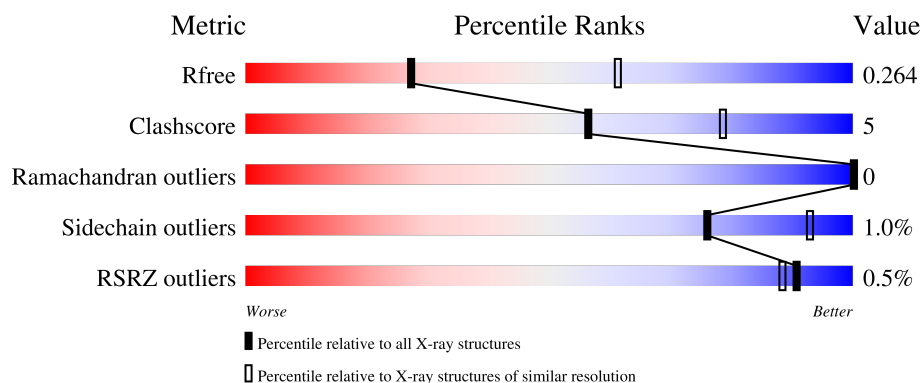
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*






The reported resolution of this entry is 2.80 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	3140 (2.80-2.80)
Clashscore	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)
RSRZ outliers	127900	3078 (2.80-2.80)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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	242	 82% 9% 9%
1	B	242	 83% 9% 7%
1	C	242	 86% 7% 8%
1	D	242	 83% 8% 8%
1	E	242	 81% 10% 9%

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Mol	Chain	Length	Quality of chain
1	F	242	 84% 8% 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
2	PO4	E	221	-	X	-	-

## 2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 10803 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 Dethiobiotin synthetase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	221	Total	C	N	O	S	0	3	0
			1728	1103	291	325	9			
1	B	224	Total	C	N	O	S	0	3	0
			1761	1122	300	330	9			
1	C	223	Total	C	N	O	S	0	2	0
			1731	1103	292	327	9			
1	D	223	Total	C	N	O	S	0	1	0
			1740	1112	294	325	9			
1	E	221	Total	C	N	O	S	0	1	0
			1723	1099	292	323	9			
1	F	223	Total	C	N	O	S	0	1	0
			1735	1107	294	325	9			

There are 144 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-21	MET	-	expression tag	UNP O24872
A	-20	GLY	-	expression tag	UNP O24872
A	-19	SER	-	expression tag	UNP O24872
A	-18	SER	-	expression tag	UNP O24872
A	-17	HIS	-	expression tag	UNP O24872
A	-16	HIS	-	expression tag	UNP O24872
A	-15	HIS	-	expression tag	UNP O24872
A	-14	HIS	-	expression tag	UNP O24872
A	-13	HIS	-	expression tag	UNP O24872
A	-12	HIS	-	expression tag	UNP O24872
A	-11	SER	-	expression tag	UNP O24872
A	-10	SER	-	expression tag	UNP O24872
A	-9	GLY	-	expression tag	UNP O24872
A	-8	ARG	-	expression tag	UNP O24872
A	-7	GLU	-	expression tag	UNP O24872
A	-6	ASN	-	expression tag	UNP O24872
A	-5	LEU	-	expression tag	UNP O24872

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-4	TYR	-	expression tag	UNP O24872
A	-3	PHE	-	expression tag	UNP O24872
A	-2	GLN	-	expression tag	UNP O24872
A	-1	GLY	-	expression tag	UNP O24872
A	0	HIS	-	expression tag	UNP O24872
A	219	GLY	-	expression tag	UNP O24872
A	220	SER	-	expression tag	UNP O24872
B	-21	MET	-	expression tag	UNP O24872
B	-20	GLY	-	expression tag	UNP O24872
B	-19	SER	-	expression tag	UNP O24872
B	-18	SER	-	expression tag	UNP O24872
B	-17	HIS	-	expression tag	UNP O24872
B	-16	HIS	-	expression tag	UNP O24872
B	-15	HIS	-	expression tag	UNP O24872
B	-14	HIS	-	expression tag	UNP O24872
B	-13	HIS	-	expression tag	UNP O24872
B	-12	HIS	-	expression tag	UNP O24872
B	-11	SER	-	expression tag	UNP O24872
B	-10	SER	-	expression tag	UNP O24872
B	-9	GLY	-	expression tag	UNP O24872
B	-8	ARG	-	expression tag	UNP O24872
B	-7	GLU	-	expression tag	UNP O24872
B	-6	ASN	-	expression tag	UNP O24872
B	-5	LEU	-	expression tag	UNP O24872
B	-4	TYR	-	expression tag	UNP O24872
B	-3	PHE	-	expression tag	UNP O24872
B	-2	GLN	-	expression tag	UNP O24872
B	-1	GLY	-	expression tag	UNP O24872
B	0	HIS	-	expression tag	UNP O24872
B	219	GLY	-	expression tag	UNP O24872
B	220	SER	-	expression tag	UNP O24872
C	-21	MET	-	expression tag	UNP O24872
C	-20	GLY	-	expression tag	UNP O24872
C	-19	SER	-	expression tag	UNP O24872
C	-18	SER	-	expression tag	UNP O24872
C	-17	HIS	-	expression tag	UNP O24872
C	-16	HIS	-	expression tag	UNP O24872
C	-15	HIS	-	expression tag	UNP O24872
C	-14	HIS	-	expression tag	UNP O24872
C	-13	HIS	-	expression tag	UNP O24872
C	-12	HIS	-	expression tag	UNP O24872
C	-11	SER	-	expression tag	UNP O24872

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-10	SER	-	expression tag	UNP O24872
C	-9	GLY	-	expression tag	UNP O24872
C	-8	ARG	-	expression tag	UNP O24872
C	-7	GLU	-	expression tag	UNP O24872
C	-6	ASN	-	expression tag	UNP O24872
C	-5	LEU	-	expression tag	UNP O24872
C	-4	TYR	-	expression tag	UNP O24872
C	-3	PHE	-	expression tag	UNP O24872
C	-2	GLN	-	expression tag	UNP O24872
C	-1	GLY	-	expression tag	UNP O24872
C	0	HIS	-	expression tag	UNP O24872
C	219	GLY	-	expression tag	UNP O24872
C	220	SER	-	expression tag	UNP O24872
D	-21	MET	-	expression tag	UNP O24872
D	-20	GLY	-	expression tag	UNP O24872
D	-19	SER	-	expression tag	UNP O24872
D	-18	SER	-	expression tag	UNP O24872
D	-17	HIS	-	expression tag	UNP O24872
D	-16	HIS	-	expression tag	UNP O24872
D	-15	HIS	-	expression tag	UNP O24872
D	-14	HIS	-	expression tag	UNP O24872
D	-13	HIS	-	expression tag	UNP O24872
D	-12	HIS	-	expression tag	UNP O24872
D	-11	SER	-	expression tag	UNP O24872
D	-10	SER	-	expression tag	UNP O24872
D	-9	GLY	-	expression tag	UNP O24872
D	-8	ARG	-	expression tag	UNP O24872
D	-7	GLU	-	expression tag	UNP O24872
D	-6	ASN	-	expression tag	UNP O24872
D	-5	LEU	-	expression tag	UNP O24872
D	-4	TYR	-	expression tag	UNP O24872
D	-3	PHE	-	expression tag	UNP O24872
D	-2	GLN	-	expression tag	UNP O24872
D	-1	GLY	-	expression tag	UNP O24872
D	0	HIS	-	expression tag	UNP O24872
D	219	GLY	-	expression tag	UNP O24872
D	220	SER	-	expression tag	UNP O24872
E	-21	MET	-	expression tag	UNP O24872
E	-20	GLY	-	expression tag	UNP O24872
E	-19	SER	-	expression tag	UNP O24872
E	-18	SER	-	expression tag	UNP O24872
E	-17	HIS	-	expression tag	UNP O24872

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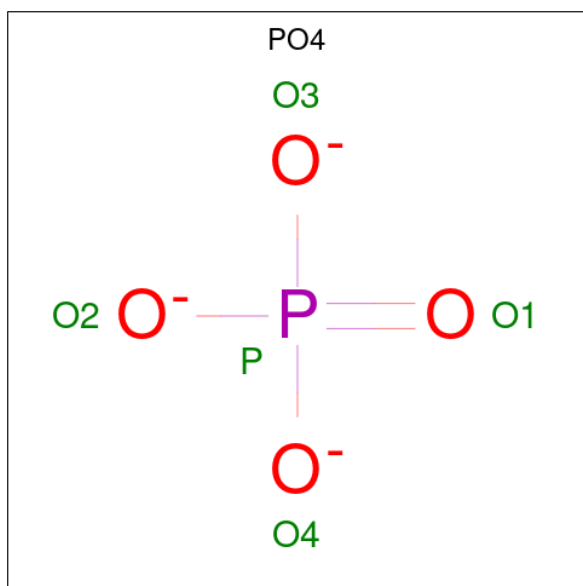
Chain	Residue	Modelled	Actual	Comment	Reference
E	-16	HIS	-	expression tag	UNP O24872
E	-15	HIS	-	expression tag	UNP O24872
E	-14	HIS	-	expression tag	UNP O24872
E	-13	HIS	-	expression tag	UNP O24872
E	-12	HIS	-	expression tag	UNP O24872
E	-11	SER	-	expression tag	UNP O24872
E	-10	SER	-	expression tag	UNP O24872
E	-9	GLY	-	expression tag	UNP O24872
E	-8	ARG	-	expression tag	UNP O24872
E	-7	GLU	-	expression tag	UNP O24872
E	-6	ASN	-	expression tag	UNP O24872
E	-5	LEU	-	expression tag	UNP O24872
E	-4	TYR	-	expression tag	UNP O24872
E	-3	PHE	-	expression tag	UNP O24872
E	-2	GLN	-	expression tag	UNP O24872
E	-1	GLY	-	expression tag	UNP O24872
E	0	HIS	-	expression tag	UNP O24872
E	219	GLY	-	expression tag	UNP O24872
E	220	SER	-	expression tag	UNP O24872
F	-21	MET	-	expression tag	UNP O24872
F	-20	GLY	-	expression tag	UNP O24872
F	-19	SER	-	expression tag	UNP O24872
F	-18	SER	-	expression tag	UNP O24872
F	-17	HIS	-	expression tag	UNP O24872
F	-16	HIS	-	expression tag	UNP O24872
F	-15	HIS	-	expression tag	UNP O24872
F	-14	HIS	-	expression tag	UNP O24872
F	-13	HIS	-	expression tag	UNP O24872
F	-12	HIS	-	expression tag	UNP O24872
F	-11	SER	-	expression tag	UNP O24872
F	-10	SER	-	expression tag	UNP O24872
F	-9	GLY	-	expression tag	UNP O24872
F	-8	ARG	-	expression tag	UNP O24872
F	-7	GLU	-	expression tag	UNP O24872
F	-6	ASN	-	expression tag	UNP O24872
F	-5	LEU	-	expression tag	UNP O24872
F	-4	TYR	-	expression tag	UNP O24872
F	-3	PHE	-	expression tag	UNP O24872
F	-2	GLN	-	expression tag	UNP O24872
F	-1	GLY	-	expression tag	UNP O24872
F	0	HIS	-	expression tag	UNP O24872
F	219	GLY	-	expression tag	UNP O24872

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Chain	Residue	Modelled	Actual	Comment	Reference
F	220	SER	-	expression tag	UNP O24872

- Molecule 2 is PHOSPHATE ION (three-letter code: PO4) (formula: O<sub>4</sub>P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	P	0	0
			5	4	1		
2	B	1	Total	O	P	0	0
			5	4	1		
2	B	1	Total	O	P	0	1
			5	4	1		
2	C	1	Total	O	P	0	0
			5	4	1		
2	D	1	Total	O	P	0	0
			5	4	1		
2	D	1	Total	O	P	0	1
			5	4	1		
2	E	1	Total	O	P	0	0
			5	4	1		
2	F	1	Total	O	P	0	0
			5	4	1		

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Mg	0	0
			1	1		

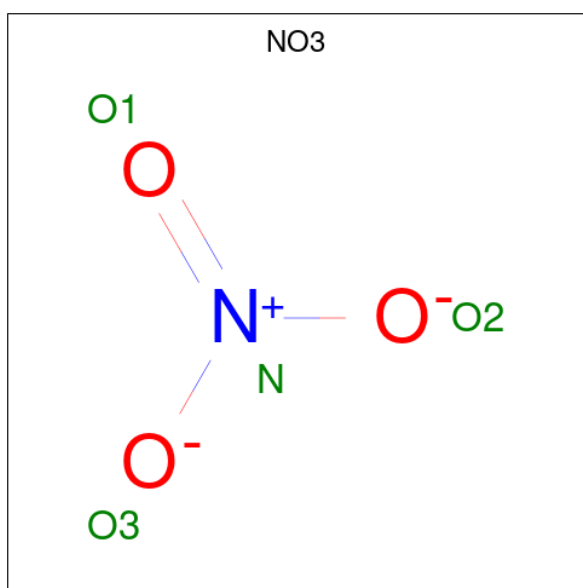
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Mg	0	0
			1	1		
3	C	1	Total	Mg	0	0
			1	1		
3	D	1	Total	Mg	0	0
			1	1		
3	E	1	Total	Mg	0	0
			1	1		
3	F	1	Total	Mg	0	0
			1	1		

- Molecule 4 is NITRATE ION (three-letter code: NO3) (formula: NO<sub>3</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	N	O	0	0
			4	1	3		
4	B	1	Total	N	O	0	0
			4	1	3		
4	F	1	Total	N	O	0	0
			4	1	3		

- Molecule 5 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

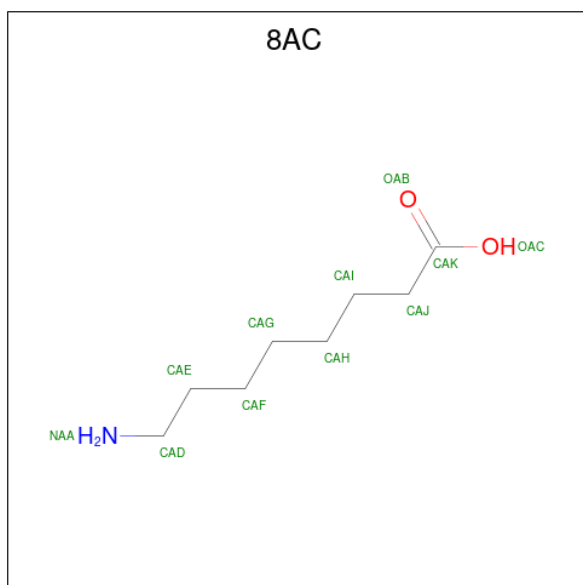
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total	Cl	0	0
			1	1		

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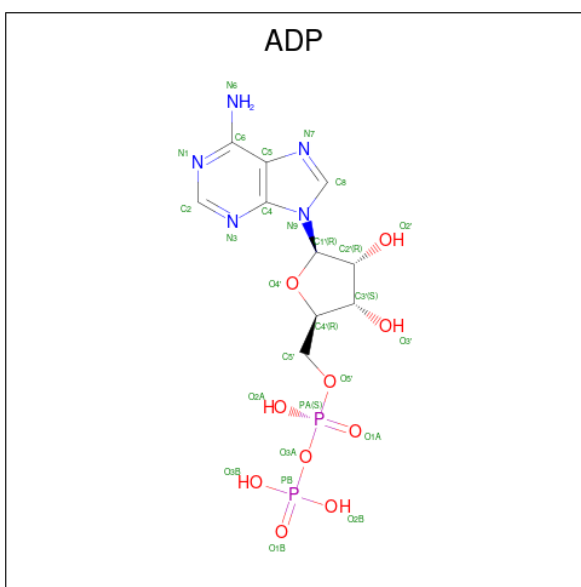
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	C	1	Total Cl 1 1	0	0
5	F	1	Total Cl 1 1	0	0

- Molecule 6 is 8-aminooctanoic acid (three-letter code: 8AC) (formula:  $C_8H_{17}NO_2$ ).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	1	Total C N O 11 8 1 2	0	0
6	B	1	Total C N O 11 8 1 2	0	0
6	C	1	Total C N O 11 8 1 2	0	0
6	D	1	Total C N O 11 8 1 2	0	0
6	E	1	Total C N O 11 8 1 2	0	0
6	F	1	Total C N O 11 8 1 2	0	0

- Molecule 7 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula:  $C_{10}H_{15}N_5O_{10}P_2$ ).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
7	A	1	Total 27	C 10	N 5	O 10	P 2	0	0
7	B	1	Total 27	C 10	N 5	O 10	P 2	0	1
7	C	1	Total 27	C 10	N 5	O 10	P 2	0	0
7	D	1	Total 27	C 10	N 5	O 10	P 2	0	1
7	E	1	Total 27	C 10	N 5	O 10	P 2	0	0
7	F	1	Total 27	C 10	N 5	O 10	P 2	0	0

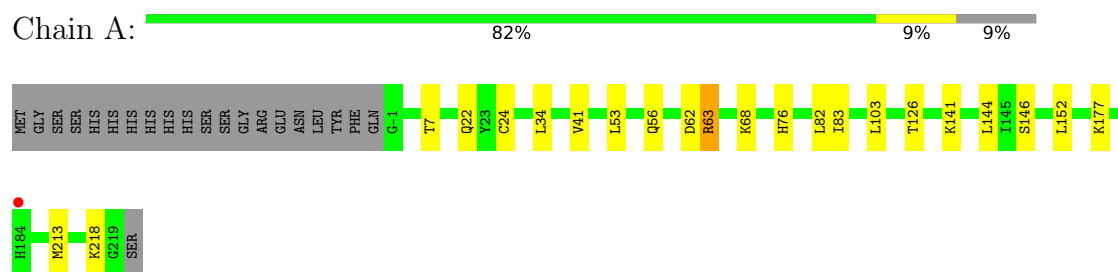
- Molecule 8 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	A	24	Total O 24 24	0	0
8	B	13	Total O 13 13	0	0
8	C	16	Total O 16 16	0	0
8	D	9	Total O 9 9	0	0
8	E	16	Total O 16 16	0	0
8	F	18	Total O 18 18	0	0

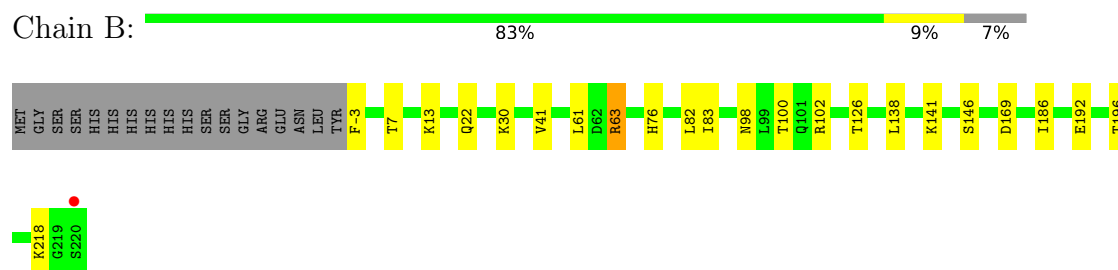
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

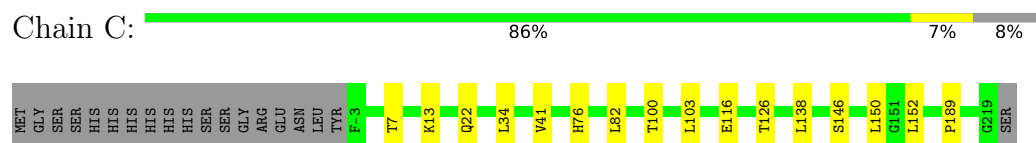
- Molecule 1: Dethiobiotin synthetase



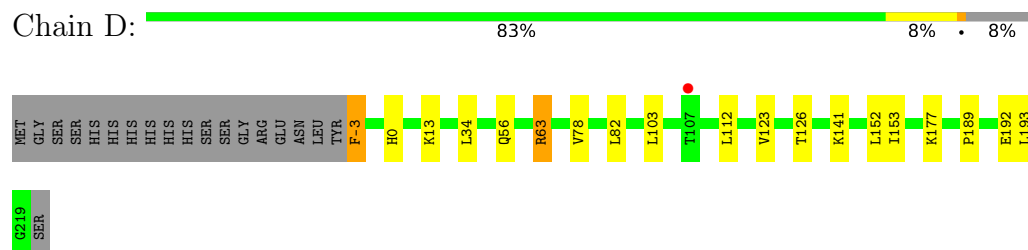
- Molecule 1: Dethiobiotin synthetase



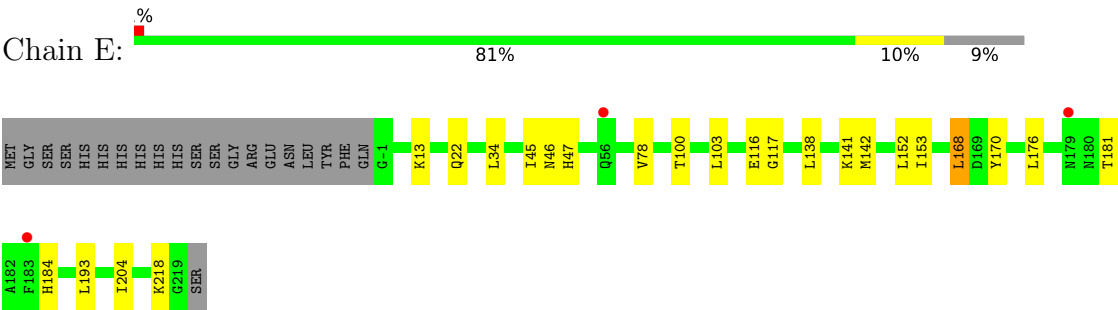
- Molecule 1: Dethiobiotin synthetase



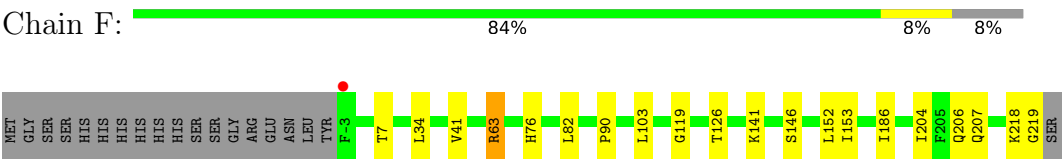
- Molecule 1: Dethiobiotin synthetase



• Molecule 1: Dethiobiotin synthetase



• Molecule 1: Dethiobiotin synthetase



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	79.77Å 131.91Å 133.10Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.80 37.20 – 2.79	Depositor EDS
% Data completeness (in resolution range)	99.9 (50.00-2.80) 99.7 (37.20-2.79)	Depositor EDS
$R_{merge}$	0.10	Depositor
$R_{sym}$	0.10	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.23 (at 2.81Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, $R_{free}$	0.201 , 0.246 0.219 , 0.264	Depositor DCC
$R_{free}$ test set	1777 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	61.0	Xtriage
Anisotropy	0.217	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.30 , 41.9	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.51$ , $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	0.000 for -h,l,k	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	10803	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	59.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.02% 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: CL, NO3, ADP, 8AC, PO4, MG

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	0.73	0/1767	0.70	0/2400
1	B	0.69	0/1798	0.69	0/2441
1	C	0.65	1/1768 (0.1%)	0.69	0/2404
1	D	0.63	0/1774	0.69	0/2408
1	E	0.61	0/1756	0.71	2/2384 (0.1%)
1	F	0.58	0/1767	0.66	0/2401
All	All	0.65	1/10630 (0.0%)	0.69	2/14438 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	116	GLU	CB-CG	5.05	1.61	1.52

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	47	HIS	N-CA-CB	-6.00	99.80	110.60
1	E	46	ASN	CB-CA-C	-5.06	100.28	110.40

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1728	0	1735	16	0
1	B	1761	0	1761	22	0
1	C	1731	0	1719	12	0
1	D	1740	0	1745	22	0
1	E	1723	0	1732	16	0
1	F	1735	0	1716	16	0
2	A	5	0	0	0	0
2	B	10	0	0	0	0
2	C	5	0	0	0	0
2	D	10	0	0	2	0
2	E	5	0	0	1	0
2	F	5	0	0	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
3	E	1	0	0	0	0
3	F	1	0	0	0	0
4	A	4	0	0	0	0
4	B	4	0	0	0	0
4	F	4	0	0	0	0
5	A	1	0	0	0	0
5	C	1	0	0	0	0
5	F	1	0	0	0	0
6	A	11	0	16	0	0
6	B	11	0	16	2	0
6	C	11	0	16	1	0
6	D	11	0	16	0	0
6	E	11	0	16	1	0
6	F	11	0	16	2	0
7	A	27	0	12	1	0
7	B	27	0	12	1	0
7	C	27	0	12	1	0
7	D	27	0	12	6	0
7	E	27	0	12	1	0
7	F	27	0	12	0	0
8	A	24	0	0	0	0
8	B	13	0	0	0	0
8	C	16	0	0	0	0
8	D	9	0	0	1	0
8	E	16	0	0	0	0
8	F	18	0	0	1	0
All	All	10803	0	10576	106	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 (106) 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:B:63[A]:ARG:HH11	1:B:63[A]:ARG:CG	1.66	1.05
1:F:63:ARG:HD2	1:F:63:ARG:O	1.58	1.03
1:B:63[A]:ARG:HG3	1:B:63[A]:ARG:NH1	1.56	1.03
1:B:63[A]:ARG:HH11	1:B:63[A]:ARG:HG3	0.75	0.90
1:B:63[A]:ARG:H	1:B:63[A]:ARG:CD	1.91	0.84
1:D:63:ARG:O	1:D:63:ARG:HD2	1.76	0.84
7:D:225[A]:ADP:H8	7:D:225[A]:ADP:H5'1	1.42	0.82
1:F:63:ARG:HD2	1:F:63:ARG:C	1.98	0.77
7:D:225[A]:ADP:H5'1	7:D:225[A]:ADP:C8	2.20	0.76
1:A:63:ARG:O	1:A:63:ARG:HD2	1.85	0.76
1:B:63[A]:ARG:H	1:B:63[A]:ARG:HD2	1.49	0.74
7:D:225[A]:ADP:H8	7:D:225[A]:ADP:C5'	2.03	0.72
1:C:22:GLN:HG3	1:E:22:GLN:HG3	1.75	0.68
7:D:225[A]:ADP:C8	7:D:225[A]:ADP:C5'	2.76	0.67
1:E:168:LEU:HD21	1:E:170:TYR:HB3	1.74	0.67
1:A:68[B]:LYS:H	1:A:68[B]:LYS:HD2	1.62	0.62
1:B:141:LYS:NZ	1:B:218:LYS:O	2.33	0.61
1:D:-3:PHE:C	1:D:-3:PHE:CD1	2.77	0.57
1:A:141:LYS:NZ	1:A:218:LYS:O	2.37	0.57
1:D:63:ARG:HD2	1:D:63:ARG:C	2.25	0.57
1:D:152:LEU:C	1:D:152:LEU:HD12	2.27	0.55
1:E:152:LEU:C	1:E:152:LEU:HD12	2.29	0.53
1:F:141:LYS:NZ	1:F:218:LYS:O	2.41	0.53
1:A:63:ARG:O	1:A:63:ARG:CD	2.55	0.53
1:B:83:ILE:HD12	1:D:189:PRO:HB2	1.91	0.52
1:D:63:ARG:NH2	8:D:228:HOH:O	2.41	0.52
1:D:141:LYS:NZ	1:D:218:LYS:O	2.43	0.52
2:E:221:PO4:P	6:F:225:8AC:NAA	2.82	0.52
6:B:225:8AC:HAI	1:D:123:VAL:HG22	1.91	0.51
1:E:181:THR:O	1:E:184:HIS:ND1	2.42	0.51
1:B:98:ASN:O	1:B:102:ARG:HG3	2.11	0.51
1:D:63:ARG:O	1:D:63:ARG:CD	2.55	0.51
1:A:83:ILE:HD12	1:C:189:PRO:HB2	1.93	0.50
1:A:53:LEU:O	1:A:56[B]:GLN:HG2	2.12	0.49
1:B:63[A]:ARG:CG	1:B:63[A]:ARG:NH1	2.40	0.49
1:D:34:LEU:CD1	1:D:103:LEU:HD22	2.43	0.49
1:E:13:LYS:HB2	7:E:223:ADP:O2B	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:141:LYS:NZ	1:E:218:LYS:O	2.45	0.48
1:A:22:GLN:HG3	1:B:22:GLN:HG3	1.95	0.48
1:C:34:LEU:CD1	1:C:103:LEU:HD22	2.43	0.48
1:D:0:HIS:O	1:D:112:LEU:HD12	2.14	0.48
1:D:-3:PHE:CD1	1:D:-3:PHE:O	2.67	0.47
1:D:177:LYS:NZ	7:D:225[A]:ADP:C5	2.83	0.47
6:E:224:8AC:HAG	1:F:119:GLY:HA3	1.97	0.47
1:D:152:LEU:HD12	1:D:153:ILE:N	2.30	0.47
1:E:168:LEU:HD23	1:E:168:LEU:C	2.36	0.46
1:F:152:LEU:HD12	1:F:152:LEU:C	2.34	0.46
1:F:218:LYS:HG3	1:F:219:GLY:N	2.30	0.46
1:C:152:LEU:HD12	1:C:152:LEU:C	2.36	0.46
1:F:41:VAL:O	1:F:76:HIS:HA	2.16	0.46
7:A:226:ADP:C2	1:B:63[A]:ARG:NH1	2.84	0.45
1:F:204:ILE:HB	1:F:207:GLN:CG	2.47	0.45
1:B:13:LYS:HB2	7:B:226[A]:ADP:O2B	2.17	0.45
1:D:34:LEU:HD13	1:D:103:LEU:HD22	1.98	0.45
1:F:34:LEU:CD1	1:F:103:LEU:HD22	2.47	0.45
1:E:45:ILE:O	1:E:45:ILE:HG22	2.16	0.45
1:E:152:LEU:HD12	1:E:153:ILE:N	2.32	0.45
1:E:78:VAL:O	1:F:186:ILE:HD13	2.17	0.44
1:E:176:LEU:HD22	1:E:204:ILE:HD13	1.99	0.44
1:D:193:LEU:HD23	1:D:193:LEU:C	2.37	0.44
6:F:225:8AC:HNA	6:F:225:8AC:HAF	1.68	0.44
1:E:34:LEU:CD1	1:E:103:LEU:HD22	2.48	0.44
1:D:82:LEU:HD13	1:D:126:THR:HA	2.00	0.44
1:A:62:ASP:OD1	1:A:62:ASP:C	2.57	0.43
1:A:63:ARG:HD2	1:A:63:ARG:C	2.39	0.43
1:B:41:VAL:O	1:B:76:HIS:HA	2.18	0.43
1:C:150:LEU:HD11	6:C:224:8AC:HAH	1.99	0.43
1:A:68[B]:LYS:H	1:A:68[B]:LYS:CD	2.30	0.43
1:D:0:HIS:CD2	1:D:217:LEU:HA	2.53	0.43
6:B:225:8AC:HADA	2:D:221:PO4:O1	2.19	0.43
1:E:142:MET:HB2	1:E:168:LEU:HD11	2.00	0.43
1:B:186:ILE:HD12	1:D:78:VAL:O	2.18	0.43
1:F:7:THR:HG23	1:F:146:SER:HA	2.00	0.43
1:A:7:THR:HG23	1:A:146:SER:HA	2.01	0.43
1:E:100:THR:HG23	1:E:138:LEU:HD23	2.01	0.43
1:A:82:LEU:HD13	1:A:126:THR:HA	2.01	0.42
1:F:63:ARG:HD3	1:F:63:ARG:HA	1.88	0.42
1:B:82:LEU:HD13	1:B:126:THR:HA	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:13:LYS:HB2	7:C:225:ADP:O2B	2.19	0.42
1:C:100:THR:HG23	1:C:138:LEU:HD23	2.02	0.42
1:D:192:GLU:O	1:D:196:THR:HG23	2.19	0.42
1:A:41:VAL:O	1:A:76:HIS:HA	2.20	0.42
1:C:82:LEU:HD13	1:C:126:THR:HA	2.01	0.42
1:F:206:GLN:N	1:F:206:GLN:OE1	2.52	0.42
1:F:152:LEU:HD12	1:F:153:ILE:N	2.34	0.42
1:A:152:LEU:C	1:A:152:LEU:HD12	2.40	0.42
1:B:192:GLU:O	1:B:196:THR:HG23	2.20	0.42
1:C:7:THR:HG23	1:C:146:SER:HA	2.02	0.42
1:D:34:LEU:HD13	1:D:103:LEU:CD2	2.50	0.42
1:B:7:THR:HG23	1:B:146:SER:HA	2.02	0.41
1:C:34:LEU:HD12	1:C:103:LEU:HD22	2.01	0.41
1:C:41:VAL:O	1:C:76:HIS:HA	2.20	0.41
1:F:82:LEU:HD13	1:F:126:THR:HA	2.01	0.41
1:A:24:CYS:SG	1:A:213:MET:HG3	2.60	0.41
1:B:30:LYS:HG2	1:B:61:LEU:HD22	2.03	0.41
1:A:34:LEU:CD1	1:A:103:LEU:HD22	2.51	0.41
1:E:116:GLU:HG2	1:E:117:GLY:O	2.21	0.41
1:B:100:THR:HG23	1:B:138:LEU:HD23	2.03	0.41
7:D:225[A]:ADP:C8	7:D:225[A]:ADP:H5'2	2.54	0.41
1:F:90:PRO:HD2	8:F:231:HOH:O	2.20	0.40
1:E:193:LEU:C	1:E:193:LEU:HD23	2.41	0.40
1:C:34:LEU:HD13	1:C:103:LEU:CD2	2.51	0.40
1:D:13:LYS:NZ	2:D:223[B]:PO4:O3	2.38	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles

### 5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	222/242 (92%)	220 (99%)	2 (1%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	225/242 (93%)	223 (99%)	2 (1%)	0	100	100
1	C	223/242 (92%)	221 (99%)	2 (1%)	0	100	100
1	D	222/242 (92%)	220 (99%)	2 (1%)	0	100	100
1	E	220/242 (91%)	215 (98%)	5 (2%)	0	100	100
1	F	222/242 (92%)	221 (100%)	1 (0%)	0	100	100
All	All	1334/1452 (92%)	1320 (99%)	14 (1%)	0	100	100

There are no Ramachandran outliers to report.

### 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	192/214 (90%)	189 (98%)	3 (2%)	62	88
1	B	195/214 (91%)	191 (98%)	4 (2%)	53	84
1	C	191/214 (89%)	191 (100%)	0	100	100
1	D	192/214 (90%)	189 (98%)	3 (2%)	62	88
1	E	191/214 (89%)	190 (100%)	1 (0%)	88	96
1	F	189/214 (88%)	188 (100%)	1 (0%)	88	96
All	All	1150/1284 (90%)	1138 (99%)	12 (1%)	76	93

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	63	ARG
1	A	144	LEU
1	A	177	LYS
1	B	-3	PHE
1	B	63[A]	ARG
1	B	63[B]	ARG
1	B	169	ASP
1	D	-3	PHE

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Mol	Chain	Res	Type
1	D	56	GLN
1	D	63	ARG
1	E	168	LEU
1	F	63	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	D	200	ASN
1	F	0	HIS

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

Of 32 ligands modelled in this entry, 9 are monoatomic - leaving 23 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	$\# Z  > 2$
2	PO4	E	221	3	4,4,4	1.10	1 (25%)	6,6,6	1.85	3 (50%)
7	ADP	C	225	3	24,29,29	1.16	2 (8%)	29,45,45	1.51	5 (17%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
2	PO4	A	221	3	4,4,4	1.34	1 (25%)	6,6,6	1.62	1 (16%)
2	PO4	B	224[B]	3	4,4,4	1.11	0	6,6,6	0.19	0
6	8AC	E	224	-	7,10,10	0.69	0	6,10,10	0.28	0
7	ADP	E	223	3	24,29,29	1.42	5 (20%)	29,45,45	1.72	6 (20%)
6	8AC	D	224	-	7,10,10	0.68	0	6,10,10	0.42	0
7	ADP	F	226	3	24,29,29	1.18	3 (12%)	29,45,45	1.61	5 (17%)
6	8AC	A	225	-	7,10,10	0.37	0	6,10,10	0.84	0
7	ADP	D	225[A]	3	24,29,29	1.32	3 (12%)	29,45,45	1.42	4 (13%)
2	PO4	C	221	3	4,4,4	0.73	0	6,6,6	0.97	0
2	PO4	B	221	3	4,4,4	1.42	1 (25%)	6,6,6	0.82	0
2	PO4	D	221	3	4,4,4	0.68	0	6,6,6	0.89	0
2	PO4	F	221	3	4,4,4	0.73	0	6,6,6	0.65	0
6	8AC	B	225	-	7,10,10	0.63	0	6,10,10	0.56	0
6	8AC	F	225	-	7,10,10	1.06	0	6,10,10	0.37	0
4	NO3	B	223	-	1,3,3	3.81	1 (100%)	0,3,3	-	-
4	NO3	F	223	-	1,3,3	4.38	1 (100%)	0,3,3	-	-
4	NO3	A	223	-	1,3,3	3.59	1 (100%)	0,3,3	-	-
7	ADP	B	226[A]	3	24,29,29	1.22	3 (12%)	29,45,45	1.69	7 (24%)
7	ADP	A	226	3	24,29,29	1.26	3 (12%)	29,45,45	1.45	5 (17%)
2	PO4	D	223[B]	3	4,4,4	1.18	0	6,6,6	0.92	0
6	8AC	C	224	-	7,10,10	0.55	0	6,10,10	0.58	0

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
7	ADP	E	223	3	-	3/12/32/32	0/3/3/3
6	8AC	D	224	-	-	4/6/8/8	-
7	ADP	F	226	3	-	1/12/32/32	0/3/3/3
6	8AC	A	225	-	-	4/6/8/8	-
6	8AC	B	225	-	-	6/6/8/8	-
7	ADP	D	225[A]	3	-	7/12/32/32	0/3/3/3
7	ADP	C	225	3	-	2/12/32/32	0/3/3/3
7	ADP	B	226[A]	3	-	2/12/32/32	0/3/3/3
7	ADP	A	226	3	-	2/12/32/32	0/3/3/3
6	8AC	F	225	-	-	3/6/8/8	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	8AC	C	224	-	-	3/6/8/8	-
6	8AC	E	224	-	-	4/6/8/8	-

All (25) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	F	223	NO3	O1-N	4.38	1.44	1.24
4	B	223	NO3	O1-N	3.81	1.41	1.24
4	A	223	NO3	O1-N	3.59	1.40	1.24
7	E	223	ADP	C5-C4	3.02	1.48	1.40
7	A	226	ADP	C5-C4	2.99	1.48	1.40
7	D	225[A]	ADP	C5-C4	2.97	1.48	1.40
7	B	226[A]	ADP	C5-C4	2.85	1.48	1.40
7	E	223	ADP	C2-N3	2.82	1.36	1.32
7	F	226	ADP	C5-C4	2.72	1.48	1.40
7	C	225	ADP	C5-C4	2.65	1.47	1.40
7	E	223	ADP	PB-O1B	2.63	1.59	1.50
7	D	225[A]	ADP	C2-N3	2.61	1.36	1.32
7	B	226[A]	ADP	C2-N3	2.57	1.36	1.32
7	D	225[A]	ADP	O4'-C1'	2.56	1.44	1.41
7	A	226	ADP	C2-N3	2.55	1.36	1.32
7	E	223	ADP	C4-N3	2.52	1.39	1.35
7	E	223	ADP	O4'-C1'	2.48	1.44	1.41
7	B	226[A]	ADP	O4'-C1'	2.43	1.44	1.41
7	F	226	ADP	C2-N3	2.43	1.36	1.32
2	A	221	PO4	P-O1	2.41	1.56	1.50
7	F	226	ADP	O4'-C1'	2.32	1.44	1.41
7	A	226	ADP	O4'-C1'	2.31	1.44	1.41
2	E	221	PO4	P-O1	2.18	1.56	1.50
7	C	225	ADP	C2-N3	2.17	1.35	1.32
2	B	221	PO4	P-O1	2.09	1.55	1.50

All (36) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	F	226	ADP	PA-O3A-PB	-4.79	116.39	132.83
7	C	225	ADP	N3-C2-N1	-4.57	121.53	128.68
7	D	225[A]	ADP	N3-C2-N1	-4.29	121.98	128.68
7	B	226[A]	ADP	N3-C2-N1	-4.24	122.05	128.68
7	A	226	ADP	N3-C2-N1	-4.05	122.34	128.68
7	E	223	ADP	N3-C2-N1	-4.03	122.37	128.68
7	E	223	ADP	PA-O3A-PB	-4.03	119.00	132.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	F	226	ADP	N3-C2-N1	-4.02	122.39	128.68
7	B	226[A]	ADP	PA-O3A-PB	-3.60	120.46	132.83
7	A	226	ADP	PA-O3A-PB	-3.43	121.06	132.83
7	C	225	ADP	C1'-N9-C4	-3.36	120.75	126.64
7	E	223	ADP	N6-C6-N1	2.99	124.77	118.57
7	C	225	ADP	PA-O3A-PB	-2.68	123.63	132.83
2	E	221	PO4	O2-P-O1	-2.59	101.41	110.89
7	B	226[A]	ADP	C2-N1-C6	2.50	123.04	118.75
7	B	226[A]	ADP	C3'-C2'-C1'	2.49	104.73	100.98
2	A	221	PO4	O2-P-O1	-2.47	101.86	110.89
7	C	225	ADP	C2-N1-C6	2.47	122.97	118.75
7	B	226[A]	ADP	C5'-C4'-C3'	-2.41	106.16	115.18
2	E	221	PO4	O4-P-O3	-2.35	100.44	107.97
7	B	226[A]	ADP	C4-C5-N7	-2.32	106.99	109.40
7	A	226	ADP	C4-C5-N7	-2.31	106.99	109.40
7	A	226	ADP	O3B-PB-O2B	2.29	116.39	107.64
7	A	226	ADP	C2-N1-C6	2.21	122.53	118.75
7	E	223	ADP	C5-C6-N6	-2.19	117.02	120.35
7	D	225[A]	ADP	C2-N1-C6	2.19	122.51	118.75
7	D	225[A]	ADP	C4-C5-N7	-2.14	107.17	109.40
7	C	225	ADP	C4-C5-N7	-2.10	107.21	109.40
7	E	223	ADP	C2-N1-C6	2.09	122.32	118.75
7	B	226[A]	ADP	O2A-PA-O1A	2.06	122.42	112.24
7	E	223	ADP	O2A-PA-O1A	2.06	122.42	112.24
7	F	226	ADP	N6-C6-N1	2.05	122.84	118.57
7	F	226	ADP	C2-N1-C6	2.05	122.26	118.75
7	D	225[A]	ADP	C3'-C2'-C1'	2.05	104.06	100.98
2	E	221	PO4	O4-P-O2	2.03	114.49	107.97
7	F	226	ADP	O3B-PB-O1B	2.00	118.51	110.68

There are no chirality outliers.

All (41) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	A	225	8AC	CAH-CAI-CAJ-CAK
6	C	224	8AC	CAH-CAI-CAJ-CAK
6	D	224	8AC	CAH-CAI-CAJ-CAK
6	E	224	8AC	CAH-CAI-CAJ-CAK
7	A	226	ADP	O4'-C4'-C5'-O5'
7	B	226[A]	ADP	O4'-C4'-C5'-O5'
7	C	225	ADP	O4'-C4'-C5'-O5'
7	D	225[A]	ADP	C5'-O5'-PA-O1A

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Mol	Chain	Res	Type	Atoms
7	D	225[A]	ADP	C5'-O5'-PA-O2A
7	F	226	ADP	O4'-C4'-C5'-O5'
6	C	224	8AC	CAG-CAH-CAI-CAJ
6	B	225	8AC	CAG-CAH-CAI-CAJ
6	C	224	8AC	CAF-CAG-CAH-CAI
6	F	225	8AC	CAD-CAE-CAF-CAG
7	D	225[A]	ADP	C3'-C4'-C5'-O5'
6	E	224	8AC	CAD-CAE-CAF-CAG
6	B	225	8AC	CAF-CAG-CAH-CAI
7	A	226	ADP	PA-O3A-PB-O1B
6	A	225	8AC	CAD-CAE-CAF-CAG
6	B	225	8AC	CAE-CAF-CAG-CAH
6	F	225	8AC	NAA-CAD-CAE-CAF
6	F	225	8AC	CAE-CAF-CAG-CAH
6	B	225	8AC	CAD-CAE-CAF-CAG
7	D	225[A]	ADP	O4'-C4'-C5'-O5'
6	B	225	8AC	NAA-CAD-CAE-CAF
7	E	223	ADP	O4'-C4'-C5'-O5'
7	D	225[A]	ADP	PB-O3A-PA-O1A
7	E	223	ADP	PB-O3A-PA-O1A
6	B	225	8AC	CAH-CAI-CAJ-CAK
6	E	224	8AC	CAE-CAF-CAG-CAH
6	A	225	8AC	NAA-CAD-CAE-CAF
7	C	225	ADP	C3'-C4'-C5'-O5'
6	A	225	8AC	CAG-CAH-CAI-CAJ
6	D	224	8AC	CAD-CAE-CAF-CAG
6	D	224	8AC	CAF-CAG-CAH-CAI
7	B	226[A]	ADP	C3'-C4'-C5'-O5'
7	D	225[A]	ADP	C5'-O5'-PA-O3A
7	E	223	ADP	C4'-C5'-O5'-PA
6	E	224	8AC	NAA-CAD-CAE-CAF
7	D	225[A]	ADP	PB-O3A-PA-O2A
6	D	224	8AC	NAA-CAD-CAE-CAF

There are no ring outliers.

12 monomers are involved in 17 short contacts:

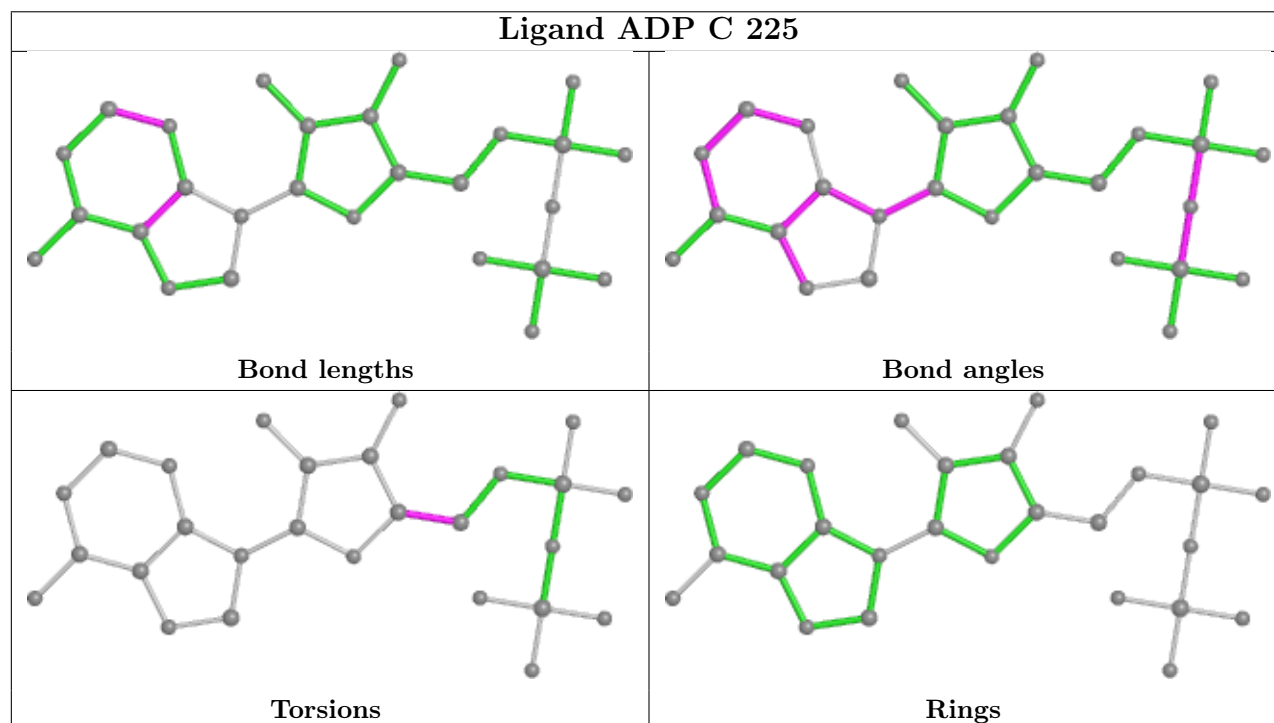
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	E	221	PO4	1	0
7	C	225	ADP	1	0
6	E	224	8AC	1	0
7	E	223	ADP	1	0

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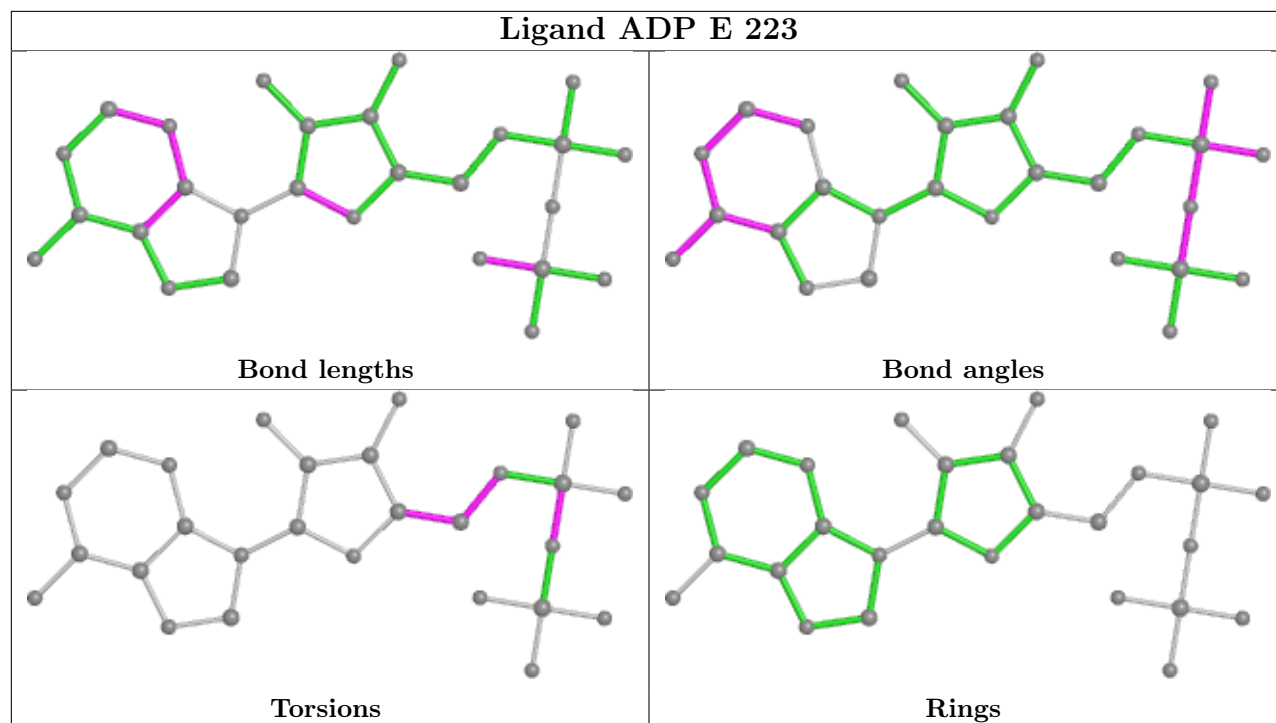
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Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	D	225[A]	ADP	6	0
2	D	221	PO4	1	0
6	B	225	8AC	2	0
6	F	225	8AC	2	0
7	B	226[A]	ADP	1	0
7	A	226	ADP	1	0
2	D	223[B]	PO4	1	0
6	C	224	8AC	1	0

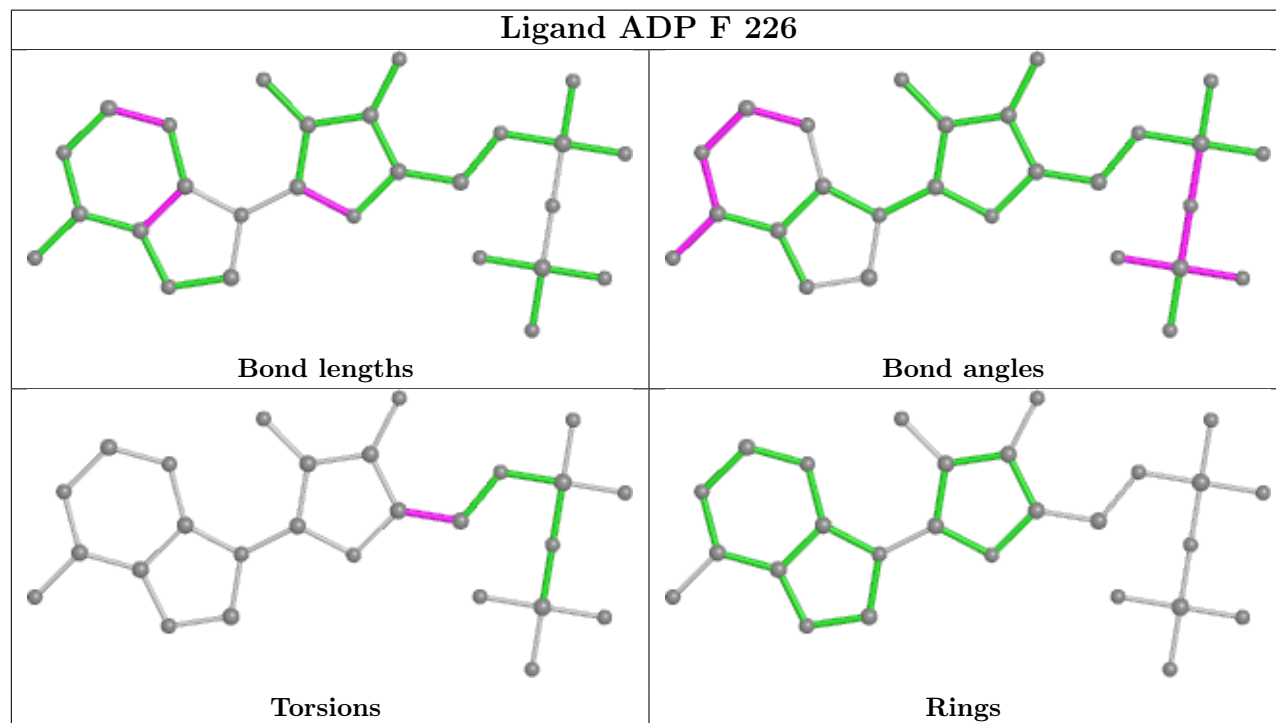
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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.



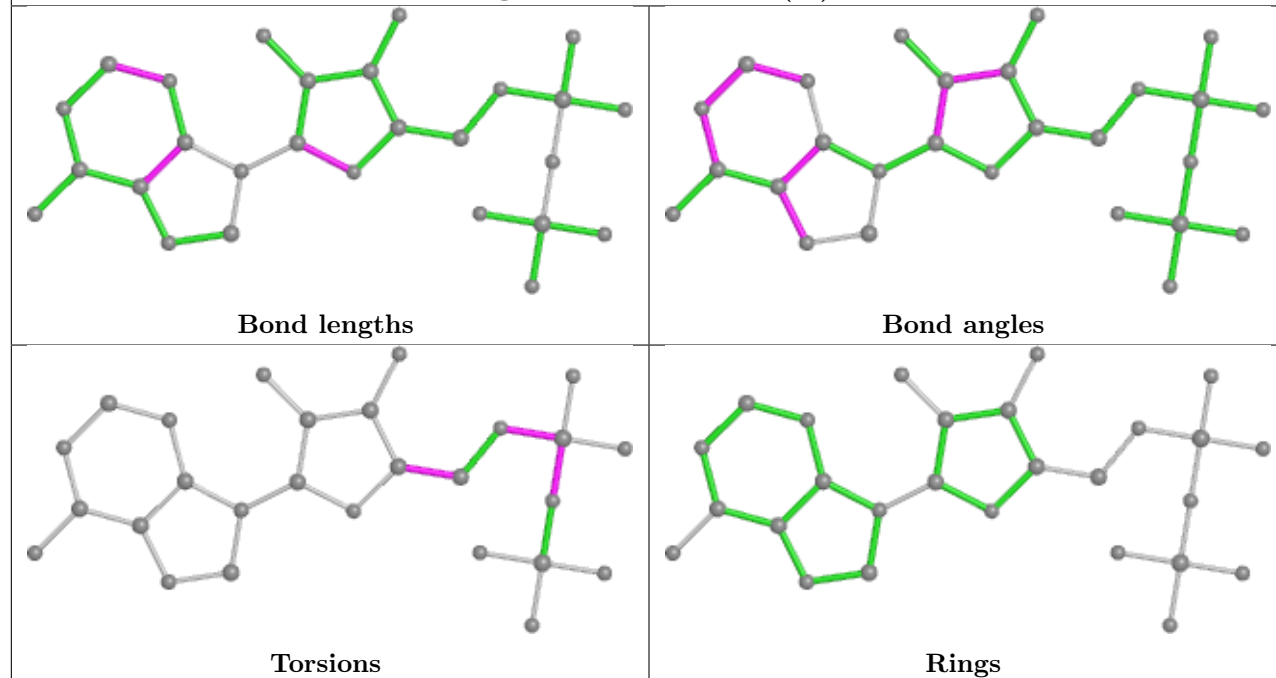
## Ligand ADP E 223



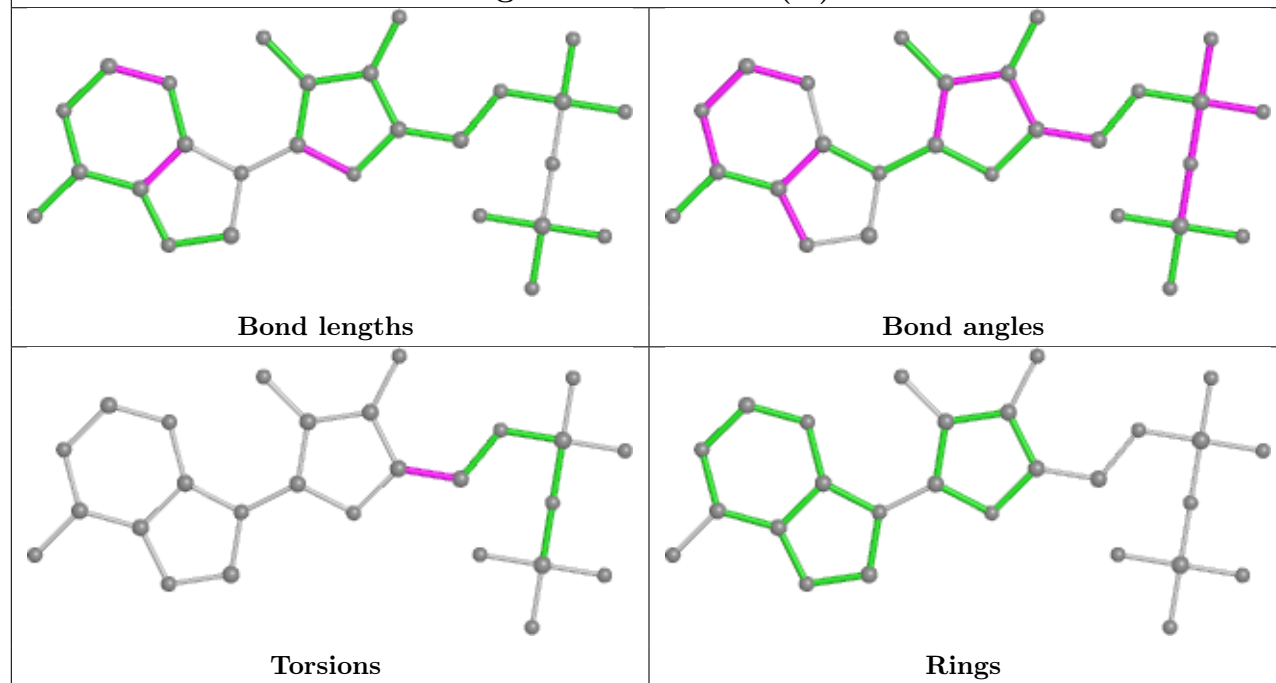
## Ligand ADP F 226

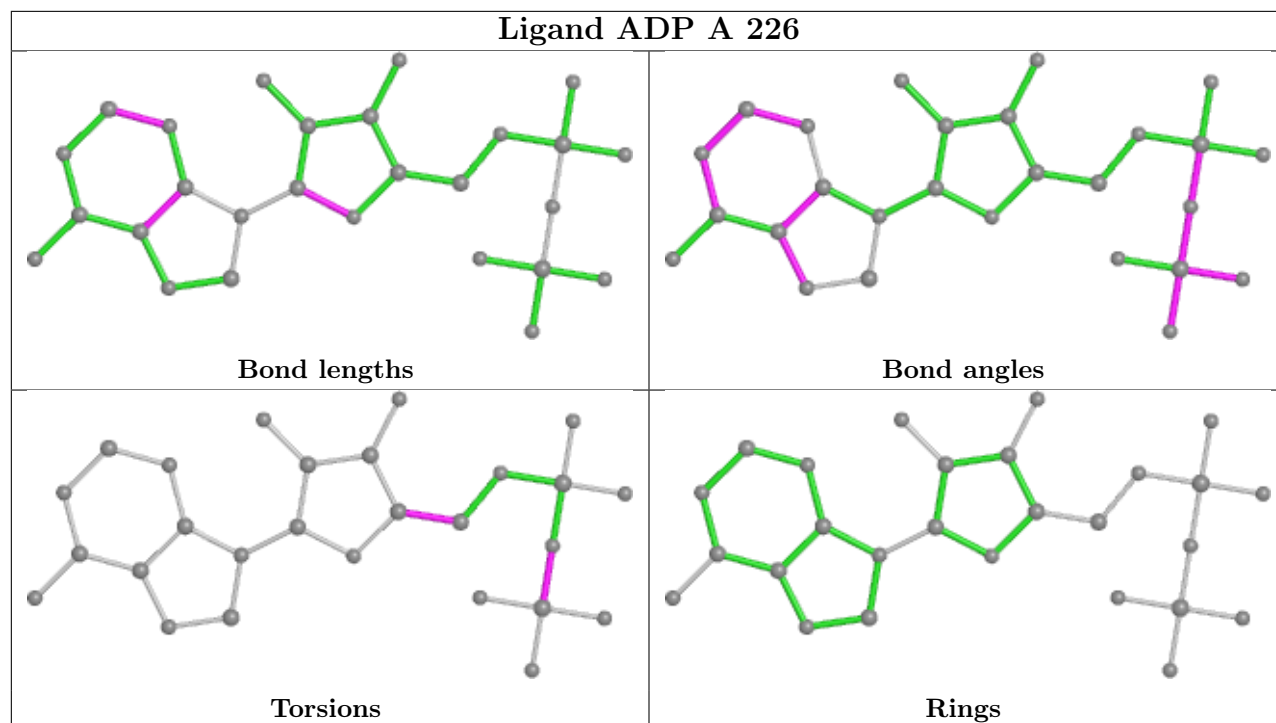


## Ligand ADP D 225 (A)



## Ligand ADP B 226 (A)





## 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	221/242 (91%)	-0.42	1 (0%) 91 88	36, 52, 77, 93	0
1	B	224/242 (92%)	-0.24	1 (0%) 92 91	37, 54, 77, 100	0
1	C	223/242 (92%)	-0.27	0 100 100	40, 57, 96, 131	0
1	D	223/242 (92%)	-0.17	1 (0%) 92 91	38, 59, 86, 117	0
1	E	221/242 (91%)	-0.17	3 (1%) 75 70	41, 62, 90, 112	0
1	F	223/242 (92%)	-0.22	1 (0%) 92 91	38, 59, 86, 107	0
All	All	1335/1452 (91%)	-0.25	7 (0%) 91 88	36, 57, 86, 131	0

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	220	SER	4.4
1	F	-3	PHE	2.9
1	D	107	THR	2.4
1	E	56[A]	GLN	2.4
1	E	183	PHE	2.4
1	A	184	HIS	2.2
1	E	179	ASN	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 ⓘ

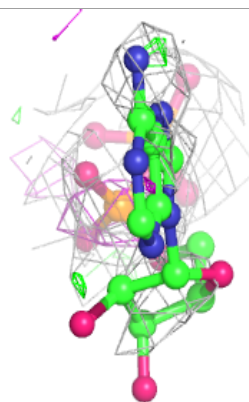
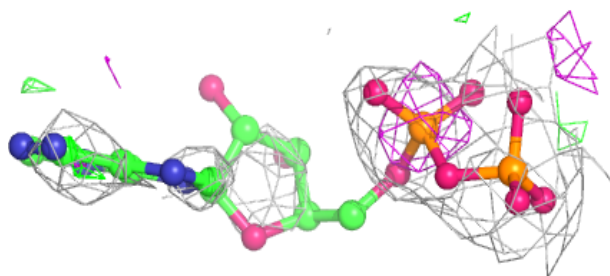
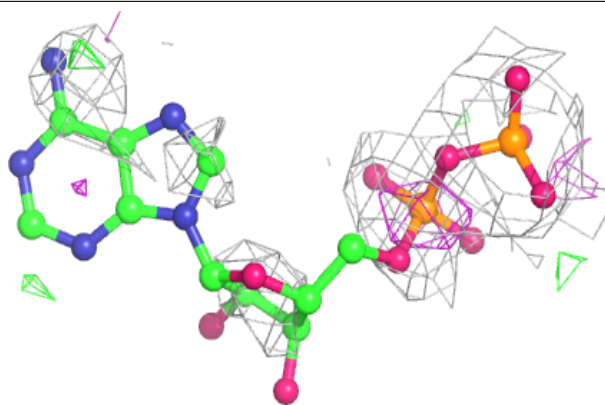
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 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
3	MG	E	222	1/1	0.53	0.09	30,30,30,30	0
6	8AC	F	225	11/11	0.80	0.32	35,44,52,55	0
3	MG	D	222	1/1	0.83	0.08	31,31,31,31	0
3	MG	B	222	1/1	0.85	0.13	24,24,24,24	0
5	CL	A	224	1/1	0.86	0.21	64,64,64,64	0
6	8AC	E	224	11/11	0.87	0.31	52,53,58,59	11
3	MG	C	222	1/1	0.88	0.14	32,32,32,32	0
7	ADP	F	226	27/27	0.88	0.29	24,85,100,101	27
4	NO3	A	223	4/4	0.89	0.24	55,59,61,61	0
4	NO3	F	223	4/4	0.89	0.62	55,58,58,58	0
7	ADP	A	226	27/27	0.89	0.26	20,81,88,91	0
7	ADP	C	225	27/27	0.89	0.27	27,95,115,124	0
3	MG	F	222	1/1	0.89	0.09	31,31,31,31	0
4	NO3	B	223	4/4	0.90	0.36	58,62,63,63	0
3	MG	A	222	1/1	0.90	0.16	29,29,29,29	0
6	8AC	B	225	11/11	0.91	0.21	42,51,55,55	0
5	CL	F	224	1/1	0.91	0.13	58,58,58,58	0
6	8AC	D	224	11/11	0.92	0.18	47,51,54,57	0
7	ADP	E	223	27/27	0.92	0.27	21,63,74,80	0
6	8AC	C	224	11/11	0.92	0.21	28,39,46,48	0
5	CL	C	223	1/1	0.93	0.08	67,67,67,67	0
6	8AC	A	225	11/11	0.94	0.19	27,34,47,48	0
7	ADP	D	225[A]	27/27	0.94	0.24	35,42,49,49	27
7	ADP	B	226[A]	27/27	0.95	0.25	27,46,52,52	27
2	PO4	E	221	5/5	0.96	0.14	29,35,42,45	0
2	PO4	C	221	5/5	0.97	0.16	31,32,39,41	0
2	PO4	D	221	5/5	0.97	0.10	35,35,41,43	0
2	PO4	A	221	5/5	0.97	0.09	22,22,28,31	0
2	PO4	F	221	5/5	0.97	0.11	41,44,46,50	0
2	PO4	B	221	5/5	0.98	0.09	27,28,30,36	0
2	PO4	D	223[B]	5/5	0.98	0.10	37,38,39,39	5
2	PO4	B	224[B]	5/5	0.99	0.10	28,33,34,35	5

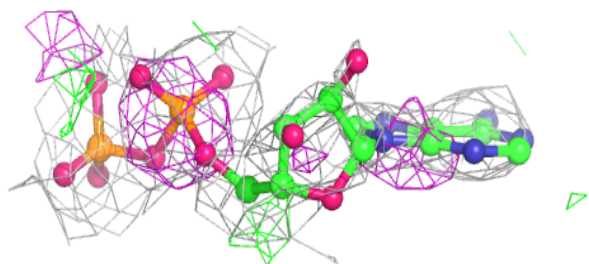
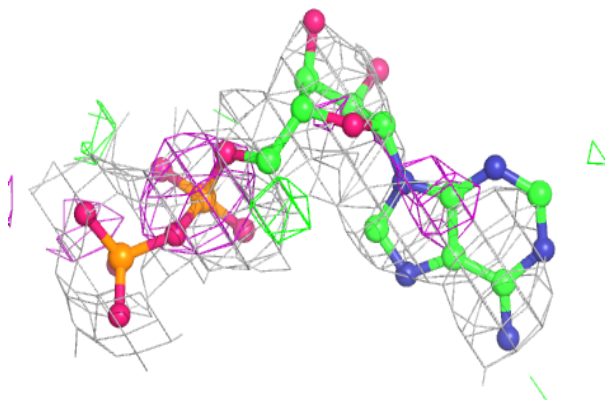
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

**Electron density around ADP F 226:**

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

**Electron density around ADP A 226:**

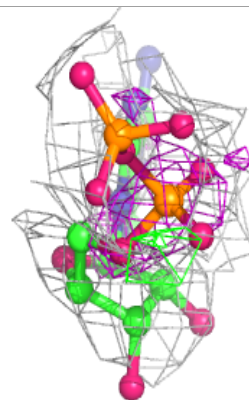
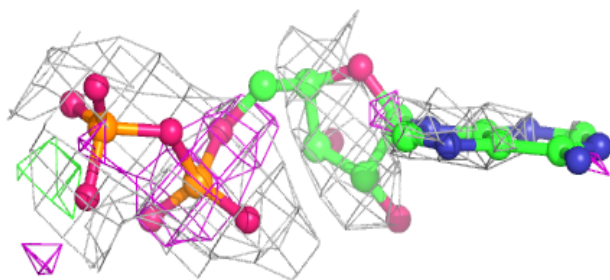
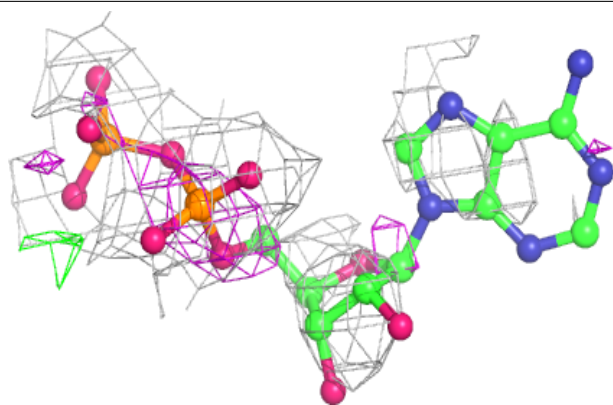
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



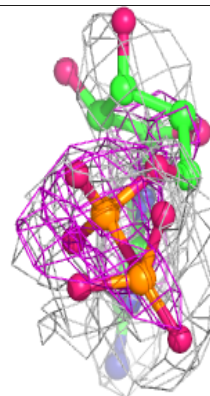
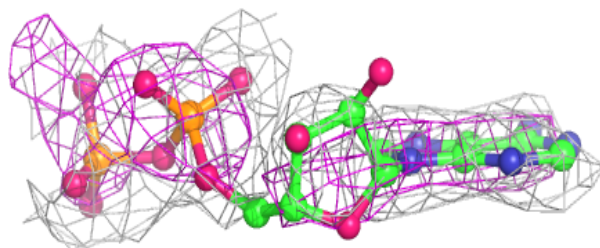
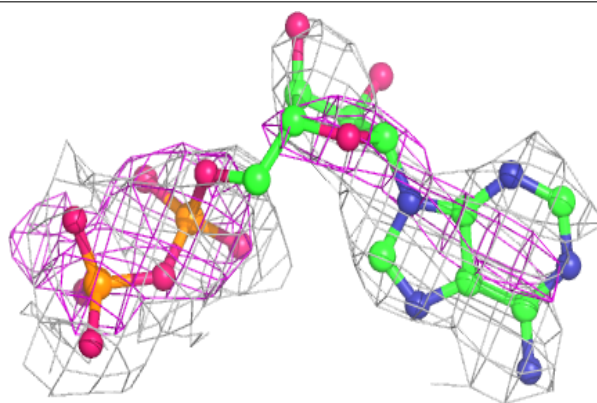


**Electron density around ADP C 225:**

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

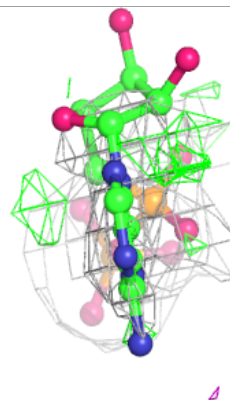
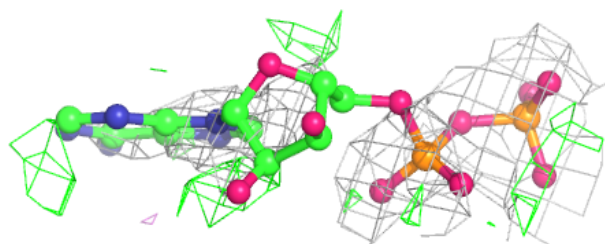
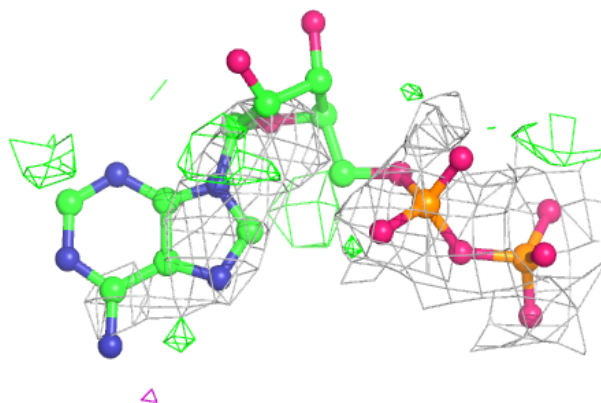
**Electron density around ADP E 223:**

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

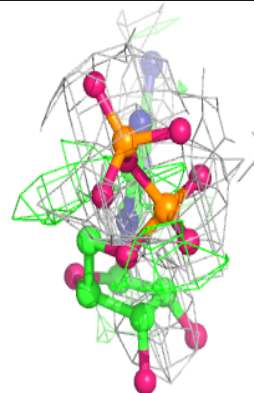
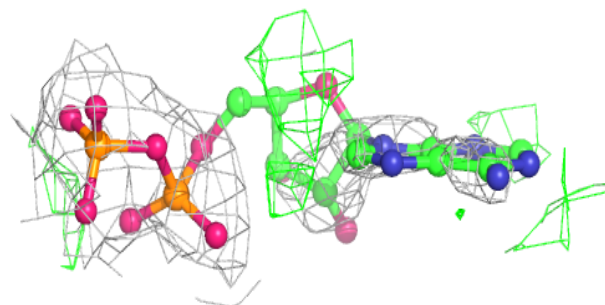
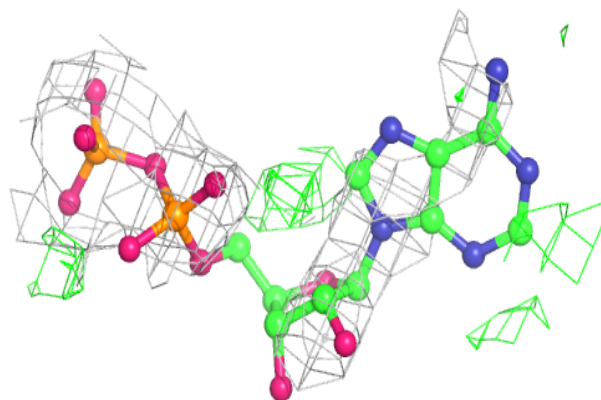


**Electron density around ADP D 225 (A):**

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

**Electron density around ADP B 226 (A):**

$2mF_o-DF_c$  (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.