Full wwPDB X-ray Structure Validation Report (i)

Sep 12, 2023 – 10:52 AM EDT

PDB ID	:	4MY9
Title	:	Crystal Structure of the Inosine 5'-monophosphate Dehydrogenase with an Internal Deletion of the CBS Domain from Bacillus anthracis str. Ames complexed with inhibitor C91
Authors	:	Kim, Y.; Makowska-Grzyska, M.; Gu, M.; Gorla, S.K.; Hedstrom, L.; Anderson, W.F.; Joachimiak, A.; CSGID; Center for Structural Genomics of Infectious Diseases (CSGID)
Deposited on	:	2013-09-27
Resolution	:	2.59 Å(reported)

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

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

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

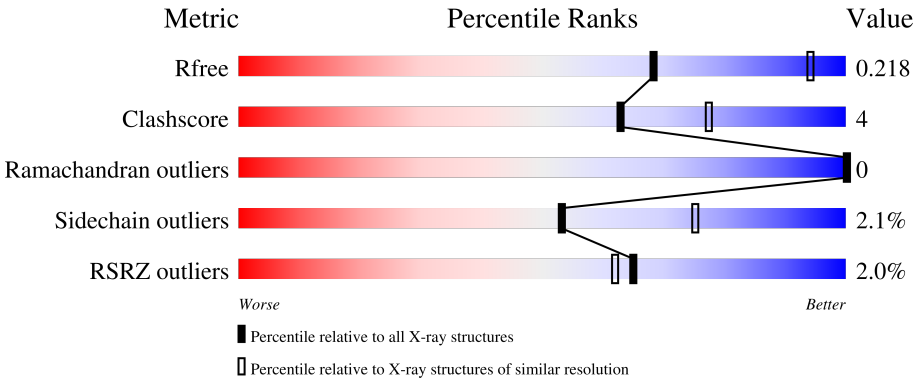
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.35.1
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)

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 2.59 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	3676 (2.60-2.56)
Clashscore	141614	4049 (2.60-2.56)
Ramachandran outliers	138981	3979 (2.60-2.56)
Sidechain outliers	138945	3979 (2.60-2.56)
RSRZ outliers	127900	3614 (2.60-2.56)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=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	384	83% 9% 9%
1	B	384	77% 14% 9%
1	C	384	80% 10% 10%

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Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.35.1

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Mol	Chain	Length	Quality of chain
1	D	384	3% 75% 15% 10%
1	E	384	% 80% 10% 9%
1	F	384	4% 80% 10% 9%
1	G	384	2% 81% 10% 8%
1	H	384	2% 80% 10% 9%

2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 21262 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 Inosine-5'-monophosphate dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	350	Total	C	N	O	S	0	0	0
			2567	1612	450	489	16			
1	B	351	Total	C	N	O	S	0	0	0
			2571	1614	451	490	16			
1	C	347	Total	C	N	O	S	0	0	0
			2544	1599	445	484	16			
1	D	347	Total	C	N	O	S	0	0	0
			2543	1597	445	485	16			
1	E	349	Total	C	N	O	S	0	0	0
			2558	1606	448	488	16			
1	F	348	Total	C	N	O	S	0	0	0
			2552	1603	447	486	16			
1	G	352	Total	C	N	O	S	0	0	0
			2577	1617	452	492	16			
1	H	348	Total	C	N	O	S	0	1	0
			2561	1608	448	489	16			

There are 208 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-23	MET	-	expression tag	UNP Q81W29
A	-22	HIS	-	expression tag	UNP Q81W29
A	-21	HIS	-	expression tag	UNP Q81W29
A	-20	HIS	-	expression tag	UNP Q81W29
A	-19	HIS	-	expression tag	UNP Q81W29
A	-18	HIS	-	expression tag	UNP Q81W29
A	-17	HIS	-	expression tag	UNP Q81W29
A	-16	SER	-	expression tag	UNP Q81W29
A	-15	SER	-	expression tag	UNP Q81W29
A	-14	GLY	-	expression tag	UNP Q81W29
A	-13	VAL	-	expression tag	UNP Q81W29
A	-12	ASP	-	expression tag	UNP Q81W29
A	-11	LEU	-	expression tag	UNP Q81W29

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-10	GLY	-	expression tag	UNP Q81W29
A	-9	THR	-	expression tag	UNP Q81W29
A	-8	GLU	-	expression tag	UNP Q81W29
A	-7	ASN	-	expression tag	UNP Q81W29
A	-6	LEU	-	expression tag	UNP Q81W29
A	-5	TYR	-	expression tag	UNP Q81W29
A	-4	PHE	-	expression tag	UNP Q81W29
A	-3	GLN	-	expression tag	UNP Q81W29
A	-2	SER	-	expression tag	UNP Q81W29
A	-1	ASN	-	expression tag	UNP Q81W29
A	0	ALA	-	expression tag	UNP Q81W29
A	92	GLY	-	linker	UNP Q81W29
A	220	GLY	-	linker	UNP Q81W29
B	-23	MET	-	expression tag	UNP Q81W29
B	-22	HIS	-	expression tag	UNP Q81W29
B	-21	HIS	-	expression tag	UNP Q81W29
B	-20	HIS	-	expression tag	UNP Q81W29
B	-19	HIS	-	expression tag	UNP Q81W29
B	-18	HIS	-	expression tag	UNP Q81W29
B	-17	HIS	-	expression tag	UNP Q81W29
B	-16	SER	-	expression tag	UNP Q81W29
B	-15	SER	-	expression tag	UNP Q81W29
B	-14	GLY	-	expression tag	UNP Q81W29
B	-13	VAL	-	expression tag	UNP Q81W29
B	-12	ASP	-	expression tag	UNP Q81W29
B	-11	LEU	-	expression tag	UNP Q81W29
B	-10	GLY	-	expression tag	UNP Q81W29
B	-9	THR	-	expression tag	UNP Q81W29
B	-8	GLU	-	expression tag	UNP Q81W29
B	-7	ASN	-	expression tag	UNP Q81W29
B	-6	LEU	-	expression tag	UNP Q81W29
B	-5	TYR	-	expression tag	UNP Q81W29
B	-4	PHE	-	expression tag	UNP Q81W29
B	-3	GLN	-	expression tag	UNP Q81W29
B	-2	SER	-	expression tag	UNP Q81W29
B	-1	ASN	-	expression tag	UNP Q81W29
B	0	ALA	-	expression tag	UNP Q81W29
B	92	GLY	-	linker	UNP Q81W29
B	220	GLY	-	linker	UNP Q81W29
C	-23	MET	-	expression tag	UNP Q81W29
C	-22	HIS	-	expression tag	UNP Q81W29
C	-21	HIS	-	expression tag	UNP Q81W29

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-20	HIS	-	expression tag	UNP Q81W29
C	-19	HIS	-	expression tag	UNP Q81W29
C	-18	HIS	-	expression tag	UNP Q81W29
C	-17	HIS	-	expression tag	UNP Q81W29
C	-16	SER	-	expression tag	UNP Q81W29
C	-15	SER	-	expression tag	UNP Q81W29
C	-14	GLY	-	expression tag	UNP Q81W29
C	-13	VAL	-	expression tag	UNP Q81W29
C	-12	ASP	-	expression tag	UNP Q81W29
C	-11	LEU	-	expression tag	UNP Q81W29
C	-10	GLY	-	expression tag	UNP Q81W29
C	-9	THR	-	expression tag	UNP Q81W29
C	-8	GLU	-	expression tag	UNP Q81W29
C	-7	ASN	-	expression tag	UNP Q81W29
C	-6	LEU	-	expression tag	UNP Q81W29
C	-5	TYR	-	expression tag	UNP Q81W29
C	-4	PHE	-	expression tag	UNP Q81W29
C	-3	GLN	-	expression tag	UNP Q81W29
C	-2	SER	-	expression tag	UNP Q81W29
C	-1	ASN	-	expression tag	UNP Q81W29
C	0	ALA	-	expression tag	UNP Q81W29
C	92	GLY	-	linker	UNP Q81W29
C	220	GLY	-	linker	UNP Q81W29
D	-23	MET	-	expression tag	UNP Q81W29
D	-22	HIS	-	expression tag	UNP Q81W29
D	-21	HIS	-	expression tag	UNP Q81W29
D	-20	HIS	-	expression tag	UNP Q81W29
D	-19	HIS	-	expression tag	UNP Q81W29
D	-18	HIS	-	expression tag	UNP Q81W29
D	-17	HIS	-	expression tag	UNP Q81W29
D	-16	SER	-	expression tag	UNP Q81W29
D	-15	SER	-	expression tag	UNP Q81W29
D	-14	GLY	-	expression tag	UNP Q81W29
D	-13	VAL	-	expression tag	UNP Q81W29
D	-12	ASP	-	expression tag	UNP Q81W29
D	-11	LEU	-	expression tag	UNP Q81W29
D	-10	GLY	-	expression tag	UNP Q81W29
D	-9	THR	-	expression tag	UNP Q81W29
D	-8	GLU	-	expression tag	UNP Q81W29
D	-7	ASN	-	expression tag	UNP Q81W29
D	-6	LEU	-	expression tag	UNP Q81W29
D	-5	TYR	-	expression tag	UNP Q81W29

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-4	PHE	-	expression tag	UNP Q81W29
D	-3	GLN	-	expression tag	UNP Q81W29
D	-2	SER	-	expression tag	UNP Q81W29
D	-1	ASN	-	expression tag	UNP Q81W29
D	0	ALA	-	expression tag	UNP Q81W29
D	92	GLY	-	linker	UNP Q81W29
D	220	GLY	-	linker	UNP Q81W29
E	-23	MET	-	expression tag	UNP Q81W29
E	-22	HIS	-	expression tag	UNP Q81W29
E	-21	HIS	-	expression tag	UNP Q81W29
E	-20	HIS	-	expression tag	UNP Q81W29
E	-19	HIS	-	expression tag	UNP Q81W29
E	-18	HIS	-	expression tag	UNP Q81W29
E	-17	HIS	-	expression tag	UNP Q81W29
E	-16	SER	-	expression tag	UNP Q81W29
E	-15	SER	-	expression tag	UNP Q81W29
E	-14	GLY	-	expression tag	UNP Q81W29
E	-13	VAL	-	expression tag	UNP Q81W29
E	-12	ASP	-	expression tag	UNP Q81W29
E	-11	LEU	-	expression tag	UNP Q81W29
E	-10	GLY	-	expression tag	UNP Q81W29
E	-9	THR	-	expression tag	UNP Q81W29
E	-8	GLU	-	expression tag	UNP Q81W29
E	-7	ASN	-	expression tag	UNP Q81W29
E	-6	LEU	-	expression tag	UNP Q81W29
E	-5	TYR	-	expression tag	UNP Q81W29
E	-4	PHE	-	expression tag	UNP Q81W29
E	-3	GLN	-	expression tag	UNP Q81W29
E	-2	SER	-	expression tag	UNP Q81W29
E	-1	ASN	-	expression tag	UNP Q81W29
E	0	ALA	-	expression tag	UNP Q81W29
E	92	GLY	-	linker	UNP Q81W29
E	220	GLY	-	linker	UNP Q81W29
F	-23	MET	-	expression tag	UNP Q81W29
F	-22	HIS	-	expression tag	UNP Q81W29
F	-21	HIS	-	expression tag	UNP Q81W29
F	-20	HIS	-	expression tag	UNP Q81W29
F	-19	HIS	-	expression tag	UNP Q81W29
F	-18	HIS	-	expression tag	UNP Q81W29
F	-17	HIS	-	expression tag	UNP Q81W29
F	-16	SER	-	expression tag	UNP Q81W29
F	-15	SER	-	expression tag	UNP Q81W29

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Chain	Residue	Modelled	Actual	Comment	Reference
F	-14	GLY	-	expression tag	UNP Q81W29
F	-13	VAL	-	expression tag	UNP Q81W29
F	-12	ASP	-	expression tag	UNP Q81W29
F	-11	LEU	-	expression tag	UNP Q81W29
F	-10	GLY	-	expression tag	UNP Q81W29
F	-9	THR	-	expression tag	UNP Q81W29
F	-8	GLU	-	expression tag	UNP Q81W29
F	-7	ASN	-	expression tag	UNP Q81W29
F	-6	LEU	-	expression tag	UNP Q81W29
F	-5	TYR	-	expression tag	UNP Q81W29
F	-4	PHE	-	expression tag	UNP Q81W29
F	-3	GLN	-	expression tag	UNP Q81W29
F	-2	SER	-	expression tag	UNP Q81W29
F	-1	ASN	-	expression tag	UNP Q81W29
F	0	ALA	-	expression tag	UNP Q81W29
F	92	GLY	-	linker	UNP Q81W29
F	220	GLY	-	linker	UNP Q81W29
G	-23	MET	-	expression tag	UNP Q81W29
G	-22	HIS	-	expression tag	UNP Q81W29
G	-21	HIS	-	expression tag	UNP Q81W29
G	-20	HIS	-	expression tag	UNP Q81W29
G	-19	HIS	-	expression tag	UNP Q81W29
G	-18	HIS	-	expression tag	UNP Q81W29
G	-17	HIS	-	expression tag	UNP Q81W29
G	-16	SER	-	expression tag	UNP Q81W29
G	-15	SER	-	expression tag	UNP Q81W29
G	-14	GLY	-	expression tag	UNP Q81W29
G	-13	VAL	-	expression tag	UNP Q81W29
G	-12	ASP	-	expression tag	UNP Q81W29
G	-11	LEU	-	expression tag	UNP Q81W29
G	-10	GLY	-	expression tag	UNP Q81W29
G	-9	THR	-	expression tag	UNP Q81W29
G	-8	GLU	-	expression tag	UNP Q81W29
G	-7	ASN	-	expression tag	UNP Q81W29
G	-6	LEU	-	expression tag	UNP Q81W29
G	-5	TYR	-	expression tag	UNP Q81W29
G	-4	PHE	-	expression tag	UNP Q81W29
G	-3	GLN	-	expression tag	UNP Q81W29
G	-2	SER	-	expression tag	UNP Q81W29
G	-1	ASN	-	expression tag	UNP Q81W29
G	0	ALA	-	expression tag	UNP Q81W29
G	92	GLY	-	linker	UNP Q81W29

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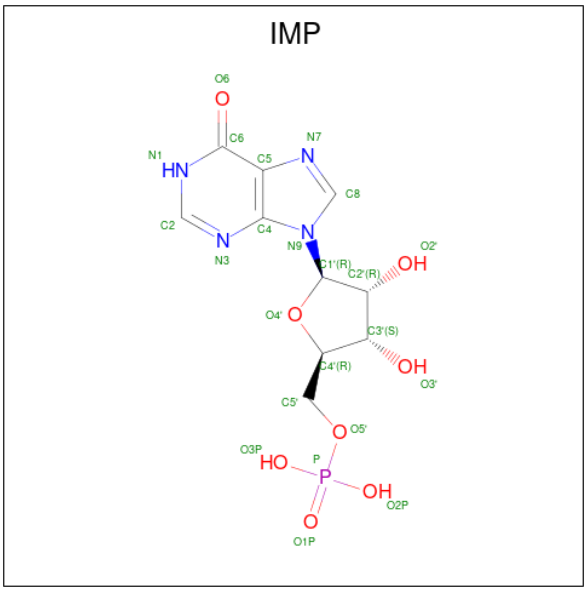
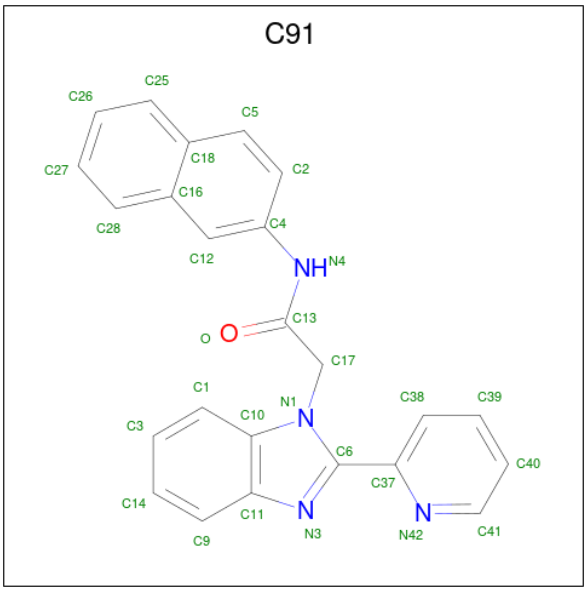
Chain	Residue	Modelled	Actual	Comment	Reference
G	220	GLY	-	linker	UNP Q81W29
H	-23	MET	-	expression tag	UNP Q81W29
H	-22	HIS	-	expression tag	UNP Q81W29
H	-21	HIS	-	expression tag	UNP Q81W29
H	-20	HIS	-	expression tag	UNP Q81W29
H	-19	HIS	-	expression tag	UNP Q81W29
H	-18	HIS	-	expression tag	UNP Q81W29
H	-17	HIS	-	expression tag	UNP Q81W29
H	-16	SER	-	expression tag	UNP Q81W29
H	-15	SER	-	expression tag	UNP Q81W29
H	-14	GLY	-	expression tag	UNP Q81W29
H	-13	VAL	-	expression tag	UNP Q81W29
H	-12	ASP	-	expression tag	UNP Q81W29
H	-11	LEU	-	expression tag	UNP Q81W29
H	-10	GLY	-	expression tag	UNP Q81W29
H	-9	THR	-	expression tag	UNP Q81W29
H	-8	GLU	-	expression tag	UNP Q81W29
H	-7	ASN	-	expression tag	UNP Q81W29
H	-6	LEU	-	expression tag	UNP Q81W29
H	-5	TYR	-	expression tag	UNP Q81W29
H	-4	PHE	-	expression tag	UNP Q81W29
H	-3	GLN	-	expression tag	UNP Q81W29
H	-2	SER	-	expression tag	UNP Q81W29
H	-1	ASN	-	expression tag	UNP Q81W29
H	0	ALA	-	expression tag	UNP Q81W29
H	92	GLY	-	linker	UNP Q81W29
H	220	GLY	-	linker	UNP Q81W29

- Molecule 2 is INOSINIC ACID (three-letter code: IMP) (formula: $C_{10}H_{13}N_4O_8P$).



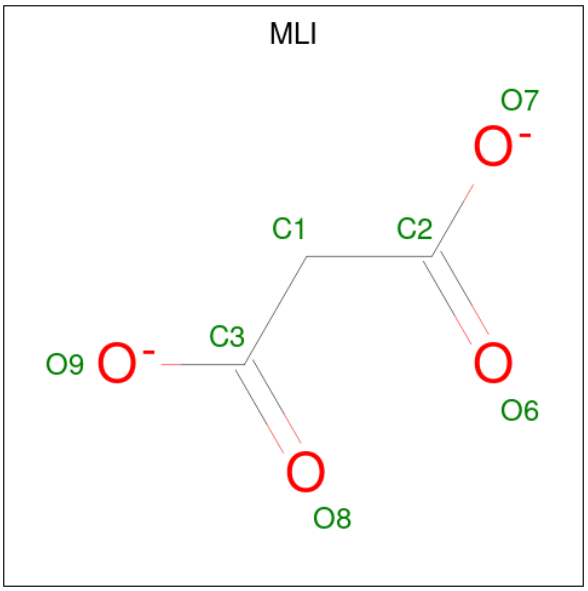
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			23	10	4	8	1		
2	B	1	Total	C	N	O	P	0	0
			23	10	4	8	1		
2	C	1	Total	C	N	O	P	0	0
			23	10	4	8	1		
2	D	1	Total	C	N	O	P	0	0
			23	10	4	8	1		
2	E	1	Total	C	N	O	P	0	0
			23	10	4	8	1		
2	F	1	Total	C	N	O	P	0	0
			23	10	4	8	1		
2	G	1	Total	C	N	O	P	0	0
			23	10	4	8	1		
2	H	1	Total	C	N	O	P	0	0
			23	10	4	8	1		

- Molecule 3 is N-(naphthalen-2-yl)-2-[2-(pyridin-2-yl)-1H-benzimidazol-1-yl]acetamide (three-letter code: C91) (formula: $C_{24}H_{18}N_4O$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C N O 29 24 4 1	0	0
3	B	1	Total C N O 29 24 4 1	0	0
3	C	1	Total C N O 29 24 4 1	0	0
3	D	1	Total C N O 29 24 4 1	0	0
3	E	1	Total C N O 29 24 4 1	0	0
3	F	1	Total C N O 29 24 4 1	0	0
3	G	1	Total C N O 29 24 4 1	0	0
3	H	1	Total C N O 29 24 4 1	0	0

- Molecule 4 is MALONATE ION (three-letter code: MLI) (formula: $C_3H_2O_4$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C O 7 3 4	0	0
4	B	1	Total C O 7 3 4	0	0
4	B	1	Total C O 7 3 4	0	0
4	C	1	Total C O 7 3 4	0	0
4	C	1	Total C O 7 3 4	0	0
4	D	1	Total C O 7 3 4	0	0
4	E	1	Total C O 7 3 4	0	0
4	G	1	Total C O 7 3 4	0	0
4	H	1	Total C O 7 3 4	0	0

- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	39	Total O 39 39	0	0
5	B	52	Total O 52 52	0	0
5	C	42	Total O 42 42	0	0

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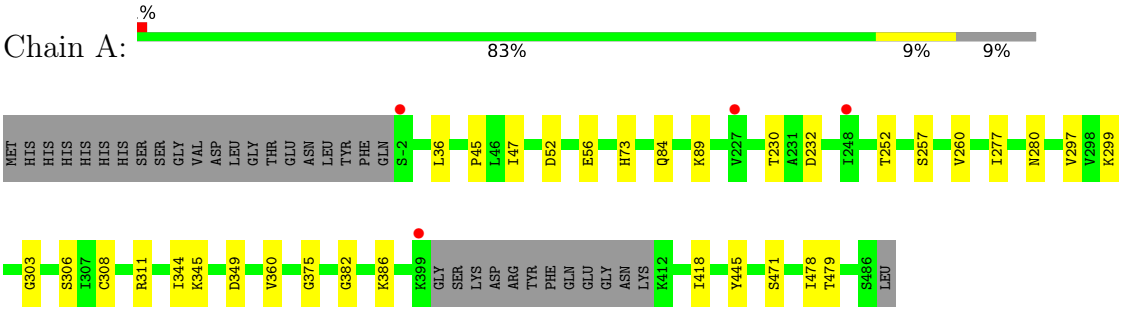
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	D	35	Total O 35 35	0	0
5	E	41	Total O 41 41	0	0
5	F	34	Total O 34 34	0	0
5	G	47	Total O 47 47	0	0
5	H	20	Total O 20 20	0	0

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.

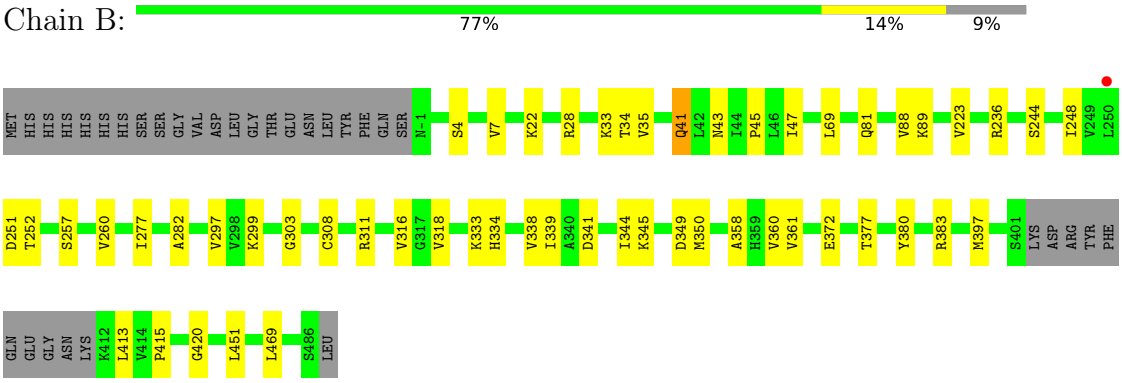
• Molecule 1: Inosine-5'-monophosphate dehydrogenase

Chain A:



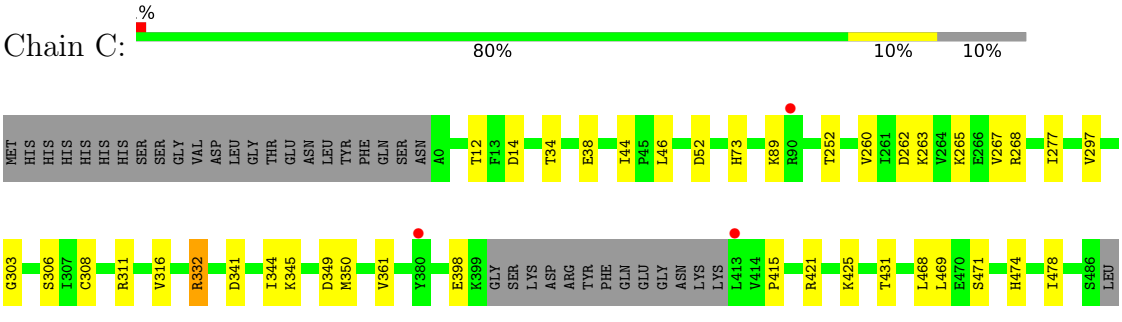
• Molecule 1: Inosine-5'-monophosphate dehydrogenase

Chain B:



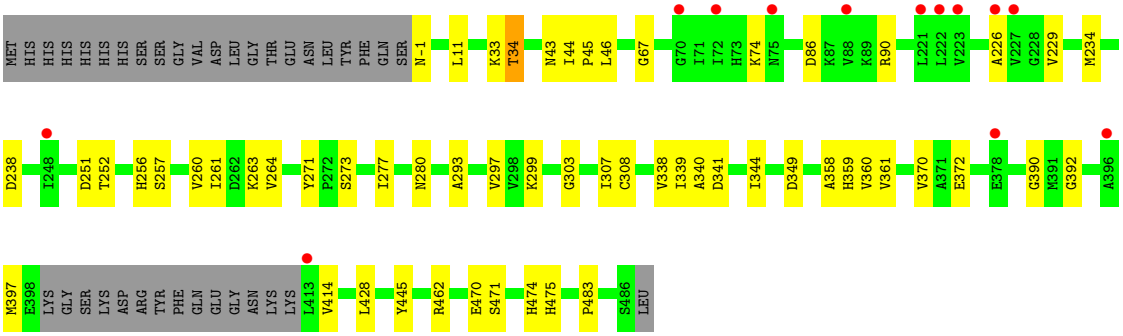
• Molecule 1: Inosine-5'-monophosphate dehydrogenase

Chain C:



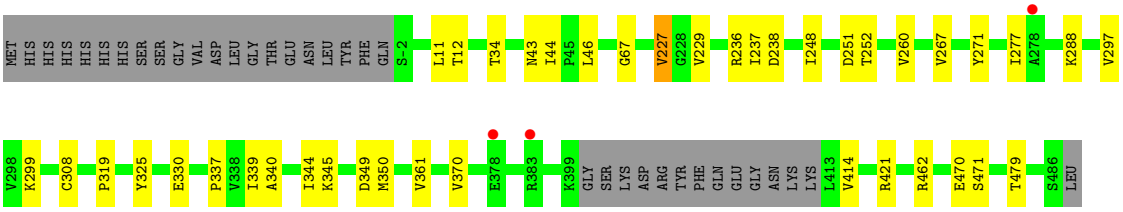
• Molecule 1: Inosine-5'-monophosphate dehydrogenase

Chain D:



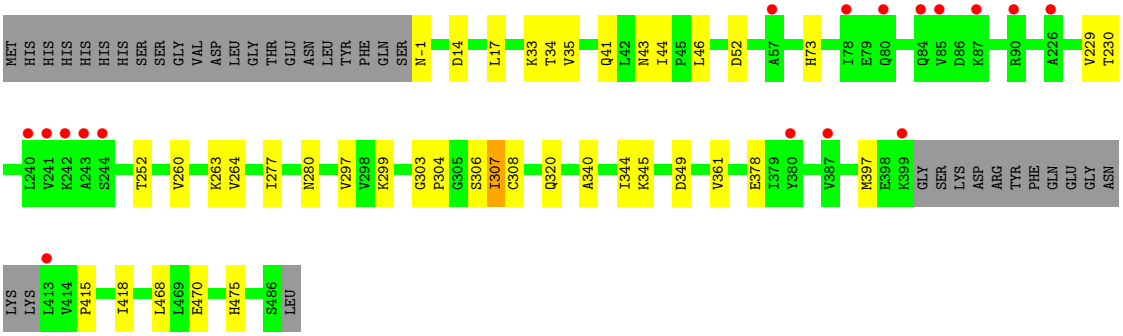
• Molecule 1: Inosine-5'-monophosphate dehydrogenase

Chain E:



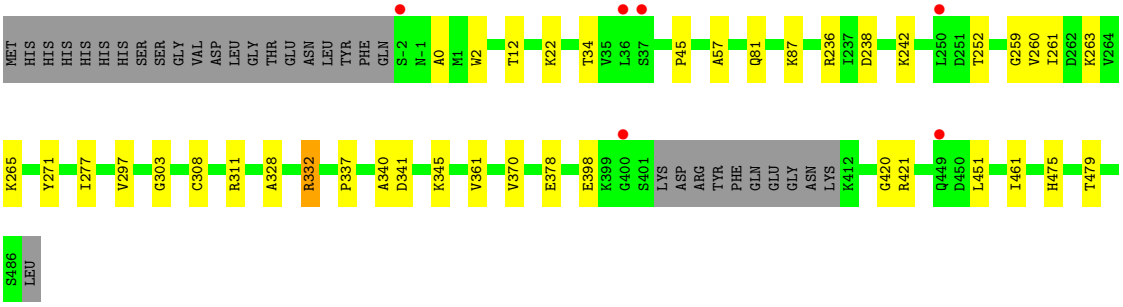
• Molecule 1: Inosine-5'-monophosphate dehydrogenase

Chain F:



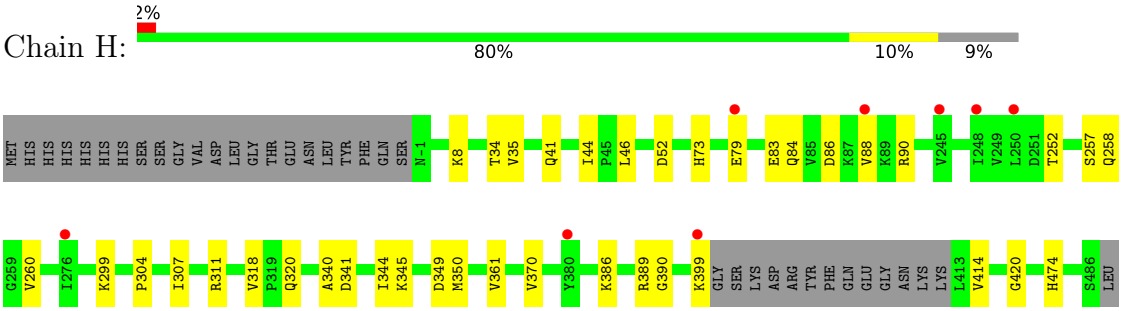
• Molecule 1: Inosine-5'-monophosphate dehydrogenase

Chain G:



• Molecule 1: Inosine-5'-monophosphate dehydrogenase

Chain H:


2%
80%
10%
9%
MET HIS HIS HIS HIS HIS HIS SER SER GLY VAL ASP LEU GLY THR GLU ASN LEU TYR PHE GLN SER N-1 K8 T34 V35 Q41 I44 P45 L46 D52 H73 E79 E83 Q84 V85 D86 K87 V88 K89 R90 V245 I248 V249 L250 D251 T252 S257 Q258
G259 V260 I276 K299 P304 I307 R311 V318 P319 Q320 A340 D341 I344 K345 D349 M350 V361 V370 Y380 K386 R389 G390 K399 GLY SER LYS ASP ARG TYR PHE GLN GLU GLY ASN LYS LYS L413 V414 G420 H474 S486 LEU

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α, β, γ	84.93Å 89.88Å 104.62Å 98.70° 90.32° 96.46°	Depositor
Resolution (Å)	38.75 – 2.59 48.11 – 2.59	Depositor EDS
% Data completeness (in resolution range)	96.6 (38.75-2.59) 96.6 (48.11-2.59)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
$< I/\sigma(I) >$ [1]	1.13 (at 2.58Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.1_1161)	Depositor
R, R_{free}	0.170 , 0.216 0.172 , 0.218	Depositor DCC
R_{free} test set	4584 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å^2)	46.8	Xtriage
Anisotropy	0.533	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.35 , 41.5	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	21262	wwPDB-VP
Average B, all atoms (Å^2)	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.69% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: C91, IMP, MLI

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	0.22	0/2603	0.42	0/3518
1	B	0.22	0/2607	0.42	0/3523
1	C	0.23	0/2580	0.43	0/3488
1	D	0.21	0/2579	0.41	0/3488
1	E	0.22	0/2594	0.42	0/3507
1	F	0.22	0/2588	0.41	0/3499
1	G	0.22	0/2613	0.42	0/3531
1	H	0.22	0/2597	0.40	0/3511
All	All	0.22	0/20761	0.42	0/28065

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

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	2567	0	2626	23	0
1	B	2571	0	2629	27	0
1	C	2544	0	2602	24	0
1	D	2543	0	2595	33	0
1	E	2558	0	2613	25	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	F	2552	0	2608	22	0
1	G	2577	0	2634	26	0
1	H	2561	0	2613	24	0
2	A	23	0	11	1	0
2	B	23	0	11	1	0
2	C	23	0	11	1	0
2	D	23	0	11	1	0
2	E	23	0	11	1	0
2	F	23	0	11	0	0
2	G	23	0	11	1	0
2	H	23	0	11	1	0
3	A	29	0	18	1	0
3	B	29	0	18	1	0
3	C	29	0	18	0	0
3	D	29	0	18	0	0
3	E	29	0	18	1	0
3	F	29	0	18	1	0
3	G	29	0	18	0	0
3	H	29	0	18	0	0
4	A	7	0	2	0	0
4	B	14	0	4	0	0
4	C	14	0	4	1	0
4	D	7	0	2	1	0
4	E	7	0	2	0	0
4	G	7	0	2	0	0
4	H	7	0	2	0	0
5	A	39	0	0	0	0
5	B	52	0	0	0	0
5	C	42	0	0	1	0
5	D	35	0	0	1	0
5	E	41	0	0	0	0
5	F	34	0	0	0	0
5	G	47	0	0	1	0
5	H	20	0	0	0	0
All	All	21262	0	21170	181	0

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

All (181) 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:252:THR:HG21	1:E:260:VAL:HG21	1.69	0.75
1:F:252:THR:HG21	1:F:260:VAL:HG21	1.77	0.67
1:A:375:GLY:O	1:A:386:LYS:NZ	2.25	0.66
1:E:277:ILE:HG13	1:E:297:VAL:HB	1.78	0.64
1:G:461:ILE:HD12	1:H:8:LYS:HE3	1.81	0.63
1:C:44:ILE:HD12	1:C:46:LEU:HD12	1.82	0.62
4:D:502:MLI:O6	5:D:618:HOH:O	2.16	0.61
1:F:307:ILE:H	1:F:307:ILE:HD13	1.65	0.61
1:E:344:ILE:HG23	1:E:349:ASP:HB2	1.82	0.60
1:E:12:THR:OG1	1:F:470:GLU:OE1	2.21	0.59
1:B:35:VAL:HG22	1:B:41:GLN:HG3	1.86	0.58
1:D:44:ILE:HD12	1:D:46:LEU:HD12	1.85	0.58
1:E:229:VAL:HG21	1:E:260:VAL:HG22	1.85	0.58
1:D:341:ASP:OD2	2:D:500:IMP:O2'	2.21	0.58
1:B:344:ILE:HG23	1:B:349:ASP:HB2	1.86	0.58
1:G:238:ASP:OD1	1:G:271:TYR:OH	2.20	0.57
1:G:378:GLU:OE1	1:G:421:ARG:NH1	2.38	0.57
1:G:252:THR:HG21	1:G:260:VAL:HG21	1.87	0.57
1:D:252:THR:HG21	1:D:260:VAL:HG21	1.86	0.57
1:G:475:HIS:CE1	1:H:345:LYS:HD2	2.40	0.57
1:A:471:SER:HA	1:B:311:ARG:HD2	1.86	0.56
1:B:277:ILE:HG12	1:B:297:VAL:HB	1.88	0.56
1:B:413:LEU:HD12	3:B:501:C91:H15	1.85	0.56
1:E:340:ALA:HB3	1:E:361:VAL:HG12	1.88	0.56
1:D:344:ILE:HG23	1:D:349:ASP:HB2	1.88	0.56
1:H:341:ASP:OD2	2:H:500:IMP:O2'	2.24	0.56
1:G:297:VAL:HG22	1:G:337:PRO:HG2	1.89	0.55
1:G:341:ASP:OD2	2:G:500:IMP:O2'	2.23	0.55
1:E:470:GLU:OE1	1:G:12:THR:OG1	2.25	0.54
1:D:338:VAL:HG23	1:D:358:ALA:HA	1.88	0.54
1:G:261:ILE:HG22	1:G:265:LYS:HE3	1.90	0.54
1:C:341:ASP:OD2	2:C:500:IMP:O2'	2.26	0.54
1:D:86:ASP:O	1:D:90:ARG:HG3	2.08	0.53
1:E:44:ILE:HD12	1:E:46:LEU:HD12	1.90	0.53
1:H:344:ILE:HG23	1:H:349:ASP:HB2	1.90	0.53
1:A:418:ILE:HD13	1:C:478:ILE:HG12	1.90	0.53
1:C:332:ARG:NH2	5:C:614:HOH:O	2.41	0.53
1:C:303:GLY:HA2	1:C:308:CYS:SG	2.49	0.53
1:D:303:GLY:HA2	1:D:308:CYS:SG	2.49	0.53
1:F:340:ALA:HB3	1:F:361:VAL:HG12	1.90	0.52
1:A:252:THR:HG21	1:A:260:VAL:HG21	1.91	0.52
1:G:22:LYS:NZ	1:H:258:GLN:HG2	2.25	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:252:THR:HG21	1:H:260:VAL:HG21	1.90	0.52
1:D:238:ASP:OD1	1:D:271:TYR:OH	2.24	0.52
1:C:52:ASP:HA	1:C:73:HIS:CD2	2.45	0.52
1:H:86:ASP:O	1:H:90:ARG:HG3	2.10	0.51
1:H:44:ILE:HD12	1:H:46:LEU:HD12	1.92	0.51
1:H:389:ARG:HH22	1:H:399:LYS:HD3	1.74	0.51
1:H:307:ILE:HD13	1:H:390:GLY:HA2	1.91	0.51
1:E:308:CYS:SG	2:E:500:IMP:H2	2.51	0.51
1:D:229:VAL:HG13	1:D:263:LYS:HD2	1.91	0.51
1:D:11:LEU:HD11	1:D:462:ARG:HD3	1.91	0.51
1:H:370:VAL:O	1:H:386:LYS:NZ	2.43	0.50
1:A:280:ASN:OD1	1:A:299:LYS:HE3	2.12	0.50
1:A:382:GLY:HA3	1:G:0:ALA:HB3	1.94	0.50
1:E:350:MET:HG3	1:E:361:VAL:HG21	1.92	0.50
1:A:257:SER:HB3	1:A:260:VAL:HG23	1.94	0.50
1:A:303:GLY:HA3	1:A:311:ARG:HE	1.77	0.50
1:B:45:PRO:HG3	1:B:451:LEU:HD11	1.94	0.49
1:F:280:ASN:OD1	1:F:299:LYS:HE3	2.12	0.49
1:H:257:SER:HB3	1:H:260:VAL:HG23	1.94	0.49
1:B:88:VAL:HG11	1:B:223:VAL:HB	1.94	0.49
1:C:12:THR:OG1	1:D:470:GLU:OE1	2.29	0.49
1:C:311:ARG:HD2	1:D:471:SER:HA	1.95	0.49
1:C:345:LYS:HD2	1:D:475:HIS:CE1	2.47	0.49
1:G:479:THR:HG23	1:H:420:GLY:HA2	1.95	0.49
1:F:44:ILE:HD12	1:F:46:LEU:HD12	1.95	0.49
1:C:268:ARG:NH1	4:C:502:MLI:O7	2.46	0.48
1:D:33:LYS:NZ	1:D:43:ASN:OD1	2.46	0.48
1:A:277:ILE:HG12	1:A:297:VAL:HB	1.93	0.48
1:C:277:ILE:HG13	1:C:297:VAL:HB	1.94	0.48
1:G:242:LYS:NZ	5:G:639:HOH:O	2.45	0.48
1:A:47:ILE:HG13	1:A:360:VAL:HG11	1.94	0.48
1:E:297:VAL:HG22	1:E:337:PRO:HG2	1.96	0.48
1:G:303:GLY:HA2	1:G:308:CYS:SG	2.54	0.48
1:F:33:LYS:HG2	1:F:43:ASN:HA	1.96	0.48
1:G:259:GLY:O	1:G:263:LYS:HG2	2.14	0.48
1:G:22:LYS:HZ1	1:H:258:GLN:HG2	1.79	0.48
1:H:340:ALA:HB3	1:H:361:VAL:HG12	1.96	0.47
1:E:471:SER:HA	1:G:311:ARG:HD2	1.95	0.47
1:A:52:ASP:HA	1:A:73:HIS:CD2	2.50	0.47
1:F:344:ILE:HG23	1:F:349:ASP:HB2	1.97	0.47
1:G:340:ALA:HB3	1:G:361:VAL:HG12	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:479:THR:HG23	1:B:420:GLY:HA2	1.96	0.47
1:B:45:PRO:HA	1:B:360:VAL:HG12	1.97	0.47
1:C:350:MET:HG3	1:C:361:VAL:HG21	1.96	0.47
1:D:340:ALA:HB3	1:D:361:VAL:HG12	1.95	0.47
1:A:56:GLU:HB3	1:A:84:GLN:HE22	1.80	0.47
1:B:33:LYS:HG2	1:B:43:ASN:HA	1.96	0.47
1:F:397:MET:HE3	1:F:415:PRO:HA	1.96	0.46
1:B:341:ASP:OD2	2:B:500:IMP:O2'	2.33	0.46
1:A:303:GLY:HA2	1:A:308:CYS:SG	2.56	0.46
1:E:479:THR:HG23	1:G:420:GLY:HA2	1.97	0.46
1:F:14:ASP:HB3	1:F:468:LEU:HD22	1.98	0.46
1:D:260:VAL:O	1:D:264:VAL:HG23	2.16	0.45
1:B:22:LYS:HG3	1:D:256:HIS:CE1	2.52	0.45
1:D:45:PRO:C	1:D:360:VAL:HG23	2.37	0.45
1:B:81:GLN:OE1	1:B:236:ARG:NH1	2.40	0.45
1:D:370:VAL:HG12	1:D:372:GLU:H	1.82	0.45
1:E:345:LYS:HD2	1:F:475:HIS:CE1	2.52	0.45
1:B:338:VAL:HG23	1:B:358:ALA:HA	1.99	0.45
1:H:311:ARG:NH2	1:H:318:VAL:O	2.49	0.45
1:B:257:SER:HB3	1:B:260:VAL:HG23	1.98	0.45
1:B:303:GLY:HA2	1:B:308:CYS:SG	2.57	0.45
1:B:380:TYR:O	1:B:383:ARG:HG2	2.16	0.44
1:H:83:GLU:OE2	1:H:90:ARG:NH1	2.49	0.44
1:H:52:ASP:HA	1:H:73:HIS:CD2	2.52	0.44
1:A:306:SER:HB2	1:C:474:HIS:O	2.17	0.44
1:D:257:SER:HB3	1:D:260:VAL:HG23	1.99	0.44
1:D:307:ILE:HD13	1:D:390:GLY:HA2	1.99	0.44
1:E:43:ASN:HB2	1:E:67:GLY:HA3	1.99	0.44
1:E:345:LYS:HE3	1:E:345:LYS:HB2	1.87	0.44
1:E:325:TYR:OH	1:G:2:TRP:O	2.23	0.44
1:D:280:ASN:OD1	1:D:299:LYS:HE2	2.18	0.43
1:E:248:ILE:HD11	1:E:267:VAL:HG11	1.99	0.43
1:B:4:SER:O	1:B:7:VAL:HG22	2.18	0.43
1:C:262:ASP:HA	1:C:265:LYS:HE3	1.98	0.43
1:E:227:VAL:HG11	1:E:237:ILE:HG13	2.00	0.43
1:F:52:ASP:HA	1:F:73:HIS:CD2	2.54	0.43
1:G:328:ALA:O	1:G:332:ARG:HB2	2.18	0.43
1:A:344:ILE:HG23	1:A:349:ASP:HB2	1.99	0.43
1:C:316:VAL:HG11	1:D:445:TYR:HB3	2.00	0.43
1:F:303:GLY:HA2	1:F:308:CYS:SG	2.58	0.43
1:B:333:LYS:HB2	1:B:334:HIS:HD2	1.82	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:261:ILE:HG23	1:D:293:ALA:HB2	2.01	0.43
1:B:397:MET:HE3	1:B:415:PRO:HA	2.00	0.43
1:D:34:THR:HG21	1:D:359:HIS:O	2.19	0.43
1:F:277:ILE:HG12	1:F:297:VAL:HB	2.00	0.43
1:D:277:ILE:HG12	1:D:297:VAL:HB	2.00	0.43
1:H:389:ARG:NH2	1:H:399:LYS:HD3	2.33	0.43
1:A:36:LEU:HD21	1:A:360:VAL:HG23	2.01	0.42
1:B:282:ALA:HB1	1:B:318:VAL:HB	2.01	0.42
1:H:350:MET:HG3	1:H:361:VAL:HG21	2.01	0.42
1:F:306:SER:HB2	1:H:474:HIS:O	2.18	0.42
1:A:45:PRO:C	1:A:360:VAL:HG13	2.40	0.42
1:A:308:CYS:SG	2:A:500:IMP:H2	2.59	0.42
1:E:227:VAL:HG13	1:E:236:ARG:HD2	1.99	0.42
1:F:230:THR:O	1:F:263:LYS:NZ	2.52	0.42
3:F:501:C91:H8	3:F:501:C91:O	2.19	0.42
1:C:14:ASP:HB3	1:C:468:LEU:HD22	2.01	0.42
1:D:299:LYS:HG3	1:D:339:ILE:HB	2.01	0.42
1:C:89:LYS:HD3	1:C:89:LYS:HA	1.78	0.42
1:B:372:GLU:OE1	1:B:372:GLU:N	2.52	0.42
1:F:229:VAL:HG21	1:F:260:VAL:HG22	2.02	0.42
1:G:252:THR:HG21	1:G:260:VAL:CG2	2.50	0.42
1:A:311:ARG:HD2	1:C:471:SER:HA	2.02	0.42
1:D:392:GLY:O	1:D:397:MET:HE3	2.20	0.42
1:A:230:THR:OG1	1:A:232:ASP:OD1	2.32	0.41
1:B:350:MET:HG3	1:B:361:VAL:HG21	2.02	0.41
1:E:299:LYS:HG3	1:E:339:ILE:HB	2.01	0.41
1:A:89:LYS:HD3	1:A:89:LYS:HA	1.84	0.41
1:C:306:SER:HB2	1:D:474:HIS:O	2.20	0.41
1:E:11:LEU:HD11	1:E:462:ARG:HD3	2.02	0.41
1:F:260:VAL:O	1:F:264:VAL:HG23	2.20	0.41
1:G:57:ALA:HB1	1:G:87:LYS:HE2	2.01	0.41
1:H:84:GLN:O	1:H:88:VAL:HG23	2.20	0.41
1:H:304:PRO:HB3	1:H:320:GLN:HB2	2.02	0.41
1:D:74:LYS:HB3	1:D:226:ALA:O	2.20	0.41
1:F:35:VAL:HG22	1:F:41:GLN:HG2	2.01	0.41
1:H:35:VAL:HG22	1:H:41:GLN:HG2	2.02	0.41
1:A:445:TYR:HB3	1:B:316:VAL:HG11	2.03	0.41
1:C:425:LYS:HE2	1:C:431:THR:OG1	2.20	0.41
1:E:288:LYS:HD3	1:E:330:GLU:OE2	2.20	0.41
1:F:304:PRO:HB3	1:F:320:GLN:HB2	2.01	0.41
1:G:277:ILE:HG23	1:G:297:VAL:HG12	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:47:ILE:HG12	1:B:69:LEU:HB3	2.03	0.41
1:G:45:PRO:HG3	1:G:451:LEU:HD11	2.02	0.41
1:D:370:VAL:HG11	1:D:428:LEU:HB2	2.03	0.41
1:B:89:LYS:HD2	1:B:244:SER:O	2.21	0.41
1:C:252:THR:HG21	1:C:260:VAL:HG22	2.02	0.41
1:C:344:ILE:HG23	1:C:349:ASP:HB2	2.03	0.41
1:E:319:PRO:HD3	1:F:17:LEU:HD12	2.03	0.41
1:F:263:LYS:HA	1:F:263:LYS:HD2	1.78	0.41
1:B:299:LYS:HG3	1:B:339:ILE:HB	2.04	0.41
3:A:501:C91:H8	3:A:501:C91:O	2.22	0.40
1:C:415:PRO:HG3	1:D:483:PRO:HD2	2.03	0.40
1:D:43:ASN:HB2	1:D:67:GLY:HA3	2.02	0.40
1:G:81:GLN:OE1	1:G:236:ARG:NH1	2.38	0.40
1:E:238:ASP:OD1	1:E:271:TYR:OH	2.28	0.40
1:C:263:LYS:O	1:C:267:VAL:HG23	2.21	0.40
3:E:501:C91:O	3:E:501:C91:H8	2.22	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	346/384 (90%)	338 (98%)	8 (2%)	0	100	100
1	B	347/384 (90%)	337 (97%)	10 (3%)	0	100	100
1	C	343/384 (89%)	337 (98%)	6 (2%)	0	100	100
1	D	343/384 (89%)	336 (98%)	7 (2%)	0	100	100
1	E	345/384 (90%)	336 (97%)	9 (3%)	0	100	100
1	F	344/384 (90%)	338 (98%)	6 (2%)	0	100	100
1	G	348/384 (91%)	340 (98%)	8 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	H	345/384 (90%)	336 (97%)	9 (3%)	0	100	100
All	All	2761/3072 (90%)	2698 (98%)	63 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	268/298 (90%)	266 (99%)	2 (1%)	84	93
1	B	268/298 (90%)	259 (97%)	9 (3%)	37	60
1	C	265/298 (89%)	259 (98%)	6 (2%)	50	73
1	D	265/298 (89%)	259 (98%)	6 (2%)	50	73
1	E	267/298 (90%)	261 (98%)	6 (2%)	52	74
1	F	266/298 (89%)	260 (98%)	6 (2%)	50	73
1	G	269/298 (90%)	264 (98%)	5 (2%)	57	77
1	H	267/298 (90%)	262 (98%)	5 (2%)	57	77
All	All	2135/2384 (90%)	2090 (98%)	45 (2%)	53	75

All (45) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	345	LYS
1	A	478	ILE
1	B	28	ARG
1	B	34	THR
1	B	41	GLN
1	B	248	ILE
1	B	251	ASP
1	B	252	THR
1	B	345	LYS
1	B	377	THR

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Mol	Chain	Res	Type
1	B	469	LEU
1	C	34	THR
1	C	38	GLU
1	C	332	ARG
1	C	398	GLU
1	C	421	ARG
1	C	469	LEU
1	D	-1	ASN
1	D	34	THR
1	D	234	MET
1	D	251	ASP
1	D	273	SER
1	D	414	VAL
1	E	34	THR
1	E	227	VAL
1	E	251	ASP
1	E	370	VAL
1	E	414	VAL
1	E	421	ARG
1	F	-1	ASN
1	F	34	THR
1	F	307	ILE
1	F	345	LYS
1	F	378	GLU
1	F	418	ILE
1	G	34	THR
1	G	332	ARG
1	G	345	LYS
1	G	370	VAL
1	G	398	GLU
1	H	34	THR
1	H	79[A]	GLU
1	H	79[B]	GLU
1	H	299	LYS
1	H	414	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	84	GLN
1	B	334	HIS
1	D	381	GLN

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Mol	Chain	Res	Type
1	D	475	HIS

5.3.3 RNA

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates

There are no monosaccharides in this entry.

5.6 Ligand geometry

25 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
3	C91	G	501	-	30,33,33	1.71	8 (26%)	38,46,46	1.10	4 (10%)
3	C91	H	501	-	30,33,33	1.71	8 (26%)	38,46,46	1.15	4 (10%)
4	MLI	C	502	-	6,6,6	1.24	0	7,7,7	1.29	0
4	MLI	C	503	-	6,6,6	1.18	0	7,7,7	1.61	2 (28%)
2	IMP	E	500	-	21,25,25	1.46	2 (9%)	24,38,38	1.24	2 (8%)
4	MLI	A	502	-	6,6,6	1.21	0	7,7,7	1.33	0
3	C91	C	501	-	30,33,33	1.68	8 (26%)	38,46,46	1.14	4 (10%)
2	IMP	H	500	-	21,25,25	1.47	2 (9%)	24,38,38	1.36	4 (16%)
4	MLI	B	503	-	6,6,6	1.20	0	7,7,7	1.32	0
4	MLI	E	502	-	6,6,6	1.23	0	7,7,7	1.31	0
3	C91	E	501	-	30,33,33	1.71	8 (26%)	38,46,46	1.12	4 (10%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\|>2$	Counts	RMSZ	$\#\|Z\|>2$
2	IMP	B	500	-	21,25,25	1.47	2 (9%)	24,38,38	1.29	5 (20%)
2	IMP	D	500	-	21,25,25	1.47	2 (9%)	24,38,38	1.30	5 (20%)
3	C91	A	501	-	30,33,33	1.70	9 (30%)	38,46,46	1.15	4 (10%)
4	MLI	B	502	-	6,6,6	1.18	0	7,7,7	1.34	0
2	IMP	G	500	-	21,25,25	1.46	2 (9%)	24,38,38	1.20	3 (12%)
4	MLI	D	502	-	6,6,6	1.13	0	7,7,7	1.64	2 (28%)
3	C91	B	501	-	30,33,33	1.71	8 (26%)	38,46,46	1.14	4 (10%)
4	MLI	G	502	-	6,6,6	1.21	0	7,7,7	1.32	0
3	C91	F	501	-	30,33,33	1.69	8 (26%)	38,46,46	1.12	4 (10%)
4	MLI	H	502	-	6,6,6	1.22	0	7,7,7	1.29	0
2	IMP	A	500	-	21,25,25	1.45	2 (9%)	24,38,38	1.26	3 (12%)
2	IMP	C	500	-	21,25,25	1.48	2 (9%)	24,38,38	1.24	2 (8%)
3	C91	D	501	-	30,33,33	1.69	8 (26%)	38,46,46	1.06	4 (10%)
2	IMP	F	500	-	21,25,25	1.46	2 (9%)	24,38,38	1.29	4 (16%)

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
3	C91	G	501	-	-	0/12/12/12	0/5/5/5
3	C91	H	501	-	-	0/12/12/12	0/5/5/5
4	MLI	C	502	-	-	4/4/4/4	-
4	MLI	C	503	-	-	3/4/4/4	-
2	IMP	E	500	-	-	3/6/26/26	0/3/3/3
4	MLI	A	502	-	-	2/4/4/4	-
3	C91	C	501	-	-	0/12/12/12	0/5/5/5
2	IMP	H	500	-	-	5/6/26/26	0/3/3/3
4	MLI	B	503	-	-	2/4/4/4	-
4	MLI	E	502	-	-	0/4/4/4	-
3	C91	E	501	-	-	0/12/12/12	0/5/5/5
2	IMP	B	500	-	-	5/6/26/26	0/3/3/3
2	IMP	D	500	-	-	4/6/26/26	0/3/3/3
3	C91	A	501	-	-	0/12/12/12	0/5/5/5
4	MLI	B	502	-	-	2/4/4/4	-
2	IMP	G	500	-	-	4/6/26/26	0/3/3/3
4	MLI	D	502	-	-	2/4/4/4	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	C91	B	501	-	-	0/12/12/12	0/5/5/5
4	MLI	G	502	-	-	2/4/4/4	-
3	C91	F	501	-	-	0/12/12/12	0/5/5/5
4	MLI	H	502	-	-	0/4/4/4	-
2	IMP	A	500	-	-	5/6/26/26	0/3/3/3
2	IMP	C	500	-	-	4/6/26/26	0/3/3/3
3	C91	D	501	-	-	3/12/12/12	0/5/5/5
2	IMP	F	500	-	-	5/6/26/26	0/3/3/3

All (81) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	500	IMP	C2-N3	4.88	1.38	1.29
2	H	500	IMP	C2-N3	4.81	1.38	1.29
2	C	500	IMP	C2-N3	4.81	1.38	1.29
2	G	500	IMP	C2-N3	4.81	1.38	1.29
2	B	500	IMP	C2-N3	4.78	1.38	1.29
2	F	500	IMP	C2-N3	4.74	1.38	1.29
2	A	500	IMP	C2-N3	4.73	1.38	1.29
2	D	500	IMP	C2-N3	4.61	1.38	1.29
2	D	500	IMP	C5-C6	-3.98	1.39	1.47
2	C	500	IMP	C5-C6	-3.97	1.39	1.47
2	B	500	IMP	C5-C6	-3.94	1.39	1.47
2	H	500	IMP	C5-C6	-3.85	1.39	1.47
2	E	500	IMP	C5-C6	-3.84	1.39	1.47
2	F	500	IMP	C5-C6	-3.83	1.39	1.47
2	G	500	IMP	C5-C6	-3.81	1.39	1.47
2	A	500	IMP	C5-C6	-3.77	1.39	1.47
3	G	501	C91	C13-N4	3.64	1.43	1.35
3	E	501	C91	C13-N4	3.63	1.43	1.35
3	D	501	C91	C13-N4	3.62	1.43	1.35
3	A	501	C91	C13-N4	3.60	1.43	1.35
3	H	501	C91	C13-N4	3.60	1.43	1.35
3	B	501	C91	C13-N4	3.59	1.43	1.35
3	F	501	C91	C13-N4	3.52	1.43	1.35
3	C	501	C91	C13-N4	3.50	1.43	1.35
3	H	501	C91	C37-C6	3.45	1.54	1.47
3	G	501	C91	C37-C6	3.44	1.54	1.47
3	A	501	C91	C37-C6	3.42	1.54	1.47
3	C	501	C91	C37-C6	3.36	1.54	1.47
3	F	501	C91	C37-C6	3.34	1.54	1.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	501	C91	C37-C6	3.33	1.54	1.47
3	B	501	C91	C37-C6	3.31	1.54	1.47
3	E	501	C91	C37-C6	3.30	1.54	1.47
3	B	501	C91	C26-C25	2.89	1.43	1.36
3	G	501	C91	C26-C25	2.85	1.43	1.36
3	A	501	C91	C26-C25	2.81	1.43	1.36
3	D	501	C91	C5-C2	2.80	1.42	1.36
3	E	501	C91	C5-C2	2.79	1.42	1.36
3	F	501	C91	C26-C25	2.78	1.43	1.36
3	C	501	C91	C26-C25	2.78	1.43	1.36
3	D	501	C91	C26-C25	2.77	1.43	1.36
3	H	501	C91	C26-C25	2.76	1.43	1.36
3	H	501	C91	C5-C2	2.76	1.42	1.36
3	D	501	C91	C16-C18	-2.76	1.36	1.42
3	E	501	C91	C26-C25	2.74	1.43	1.36
3	C	501	C91	C5-C2	2.74	1.42	1.36
3	F	501	C91	C5-C2	2.73	1.42	1.36
3	C	501	C91	C16-C18	-2.73	1.36	1.42
3	G	501	C91	C5-C2	2.72	1.42	1.36
3	F	501	C91	C16-C18	-2.72	1.36	1.42
3	H	501	C91	C16-C18	-2.70	1.36	1.42
3	A	501	C91	C16-C18	-2.68	1.36	1.42
3	E	501	C91	C16-C18	-2.68	1.36	1.42
3	B	501	C91	C16-C18	-2.67	1.36	1.42
3	A	501	C91	C5-C2	2.66	1.42	1.36
3	B	501	C91	C5-C2	2.66	1.42	1.36
3	G	501	C91	C16-C18	-2.65	1.36	1.42
3	A	501	C91	C3-C1	2.43	1.42	1.36
3	F	501	C91	C3-C1	2.39	1.42	1.36
3	C	501	C91	C3-C1	2.39	1.42	1.36
3	E	501	C91	C3-C1	2.38	1.42	1.36
3	G	501	C91	C3-C1	2.35	1.42	1.36
3	H	501	C91	C3-C1	2.34	1.42	1.36
3	D	501	C91	C3-C1	2.33	1.42	1.36
3	B	501	C91	C3-C1	2.32	1.42	1.36
3	B	501	C91	C28-C16	2.31	1.47	1.41
3	G	501	C91	C28-C16	2.26	1.47	1.41
3	E	501	C91	C9-C11	2.23	1.45	1.41
3	A	501	C91	C9-C11	2.22	1.45	1.41
3	D	501	C91	C9-C11	2.22	1.45	1.41
3	E	501	C91	C28-C16	2.20	1.47	1.41
3	C	501	C91	C28-C16	2.20	1.47	1.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	H	501	C91	C28-C16	2.20	1.47	1.41
3	D	501	C91	C28-C16	2.19	1.47	1.41
3	B	501	C91	C9-C11	2.18	1.45	1.41
3	F	501	C91	C28-C16	2.18	1.47	1.41
3	H	501	C91	C9-C11	2.18	1.45	1.41
3	G	501	C91	C9-C11	2.18	1.45	1.41
3	A	501	C91	C28-C16	2.17	1.47	1.41
3	C	501	C91	C9-C11	2.14	1.45	1.41
3	F	501	C91	C9-C11	2.12	1.45	1.41
3	A	501	C91	C27-C28	2.01	1.41	1.36

All (64) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	500	IMP	C8-N7-C5	3.20	109.09	102.99
2	E	500	IMP	C8-N7-C5	3.16	109.00	102.99
2	H	500	IMP	C8-N7-C5	3.15	109.00	102.99
2	G	500	IMP	C8-N7-C5	3.14	108.98	102.99
2	C	500	IMP	C8-N7-C5	3.14	108.96	102.99
2	F	500	IMP	C8-N7-C5	3.08	108.85	102.99
2	E	500	IMP	C5-C6-N1	3.07	119.37	113.95
2	A	500	IMP	C5-C6-N1	3.04	119.32	113.95
2	D	500	IMP	C8-N7-C5	3.03	108.75	102.99
3	C	501	C91	C6-N3-C11	3.01	109.74	103.78
2	B	500	IMP	C8-N7-C5	3.00	108.71	102.99
2	H	500	IMP	C5-C6-N1	3.00	119.25	113.95
3	H	501	C91	C6-N3-C11	2.99	109.70	103.78
2	C	500	IMP	C5-C6-N1	2.98	119.22	113.95
3	A	501	C91	C6-C37-N42	2.96	121.82	116.35
3	B	501	C91	C6-N3-C11	2.95	109.62	103.78
2	F	500	IMP	C5-C6-N1	2.94	119.14	113.95
3	F	501	C91	C6-N3-C11	2.94	109.61	103.78
3	A	501	C91	C6-N3-C11	2.93	109.60	103.78
2	B	500	IMP	C5-C6-N1	2.93	119.12	113.95
3	G	501	C91	C6-N3-C11	2.91	109.54	103.78
3	E	501	C91	C6-N3-C11	2.88	109.50	103.78
3	C	501	C91	C6-C37-N42	2.87	121.66	116.35
3	H	501	C91	C6-C37-N42	2.86	121.64	116.35
2	D	500	IMP	O3P-P-O2P	2.86	118.58	107.64
2	D	500	IMP	C5-C6-N1	2.86	119.00	113.95
3	B	501	C91	C6-C37-N42	2.86	121.64	116.35
2	H	500	IMP	O3P-P-O2P	2.84	118.49	107.64

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	G	500	IMP	C5-C6-N1	2.76	118.82	113.95
3	D	501	C91	C6-N3-C11	2.76	109.24	103.78
3	A	501	C91	C41-N42-C37	2.73	121.07	117.23
3	F	501	C91	C6-C37-N42	2.72	121.37	116.35
3	G	501	C91	C6-C37-N42	2.71	121.36	116.35
2	B	500	IMP	O3P-P-O2P	2.69	117.93	107.64
3	E	501	C91	C6-C37-N42	2.69	121.33	116.35
3	H	501	C91	C41-N42-C37	2.67	120.98	117.23
2	F	500	IMP	O3P-P-O2P	2.67	117.83	107.64
3	E	501	C91	C41-N42-C37	2.65	120.95	117.23
3	F	501	C91	C41-N42-C37	2.65	120.95	117.23
3	C	501	C91	N3-C6-N1	-2.63	108.51	115.11
3	F	501	C91	N3-C6-N1	-2.63	108.52	115.11
3	H	501	C91	N3-C6-N1	-2.62	108.53	115.11
3	C	501	C91	C41-N42-C37	2.59	120.87	117.23
3	B	501	C91	C41-N42-C37	2.58	120.85	117.23
3	A	501	C91	N3-C6-N1	-2.57	108.66	115.11
3	B	501	C91	N3-C6-N1	-2.55	108.72	115.11
3	E	501	C91	N3-C6-N1	-2.55	108.72	115.11
3	G	501	C91	N3-C6-N1	-2.54	108.73	115.11
3	D	501	C91	N3-C6-N1	-2.53	108.75	115.11
3	D	501	C91	C41-N42-C37	2.46	120.69	117.23
4	D	502	MLI	C3-C1-C2	2.35	121.08	112.87
3	G	501	C91	C41-N42-C37	2.32	120.49	117.23
2	H	500	IMP	O6-C6-C5	-2.32	119.84	124.37
3	D	501	C91	C6-C37-N42	2.25	120.51	116.35
2	B	500	IMP	O6-C6-C5	-2.16	120.14	124.37
2	D	500	IMP	O6-C6-C5	-2.15	120.17	124.37
2	F	500	IMP	O6-C6-C5	-2.12	120.23	124.37
4	C	503	MLI	O9-C3-C1	2.09	121.22	114.54
2	D	500	IMP	N1-C2-N3	-2.04	120.54	125.87
2	B	500	IMP	N1-C2-N3	-2.02	120.60	125.87
2	G	500	IMP	O2P-P-O1P	2.02	118.58	110.68
4	C	503	MLI	C3-C1-C2	2.02	119.93	112.87
2	A	500	IMP	N1-C2-N3	-2.02	120.61	125.87
4	D	502	MLI	O9-C3-C1	2.01	120.97	114.54

There are no chirality outliers.

All (55) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	500	IMP	C5'-O5'-P-O2P

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Mol	Chain	Res	Type	Atoms
2	A	500	IMP	C5'-O5'-P-O3P
2	B	500	IMP	C5'-O5'-P-O1P
2	B	500	IMP	C5'-O5'-P-O2P
2	B	500	IMP	C5'-O5'-P-O3P
2	B	500	IMP	C3'-C4'-C5'-O5'
2	C	500	IMP	C5'-O5'-P-O3P
2	D	500	IMP	C5'-O5'-P-O2P
2	D	500	IMP	C5'-O5'-P-O3P
2	E	500	IMP	C5'-O5'-P-O1P
2	F	500	IMP	C5'-O5'-P-O1P
2	F	500	IMP	C5'-O5'-P-O2P
2	F	500	IMP	C5'-O5'-P-O3P
2	G	500	IMP	C5'-O5'-P-O2P
2	G	500	IMP	C5'-O5'-P-O3P
2	H	500	IMP	C5'-O5'-P-O1P
2	H	500	IMP	C5'-O5'-P-O2P
2	H	500	IMP	C5'-O5'-P-O3P
2	H	500	IMP	C3'-C4'-C5'-O5'
3	D	501	C91	C38-C37-C6-N3
2	F	500	IMP	C3'-C4'-C5'-O5'
2	A	500	IMP	C3'-C4'-C5'-O5'
2	B	500	IMP	O4'-C4'-C5'-O5'
2	F	500	IMP	O4'-C4'-C5'-O5'
2	H	500	IMP	O4'-C4'-C5'-O5'
2	A	500	IMP	C5'-O5'-P-O1P
2	C	500	IMP	C5'-O5'-P-O1P
2	D	500	IMP	C5'-O5'-P-O1P
2	G	500	IMP	C5'-O5'-P-O1P
2	A	500	IMP	O4'-C4'-C5'-O5'
2	C	500	IMP	C3'-C4'-C5'-O5'
2	G	500	IMP	C3'-C4'-C5'-O5'
4	C	502	MLI	C3-C1-C2-O7
2	C	500	IMP	C5'-O5'-P-O2P
4	C	502	MLI	C3-C1-C2-O6
4	B	503	MLI	C2-C1-C3-O9
4	D	502	MLI	C3-C1-C2-O7
2	D	500	IMP	C3'-C4'-C5'-O5'
4	A	502	MLI	C3-C1-C2-O7
4	B	502	MLI	C2-C1-C3-O8
4	B	502	MLI	C2-C1-C3-O9
4	B	503	MLI	C2-C1-C3-O8
4	C	502	MLI	C2-C1-C3-O8

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Mol	Chain	Res	Type	Atoms
4	C	503	MLI	C2-C1-C3-O8
4	A	502	MLI	C3-C1-C2-O6
4	C	502	MLI	C2-C1-C3-O9
4	C	503	MLI	C2-C1-C3-O9
4	D	502	MLI	C3-C1-C2-O6
2	E	500	IMP	C5'-O5'-P-O3P
4	G	502	MLI	C3-C1-C2-O6
4	G	502	MLI	C3-C1-C2-O7
3	D	501	C91	N42-C37-C6-N1
3	D	501	C91	N42-C37-C6-N3
2	E	500	IMP	C3'-C4'-C5'-O5'
4	C	503	MLI	C3-C1-C2-O7

There are no ring outliers.

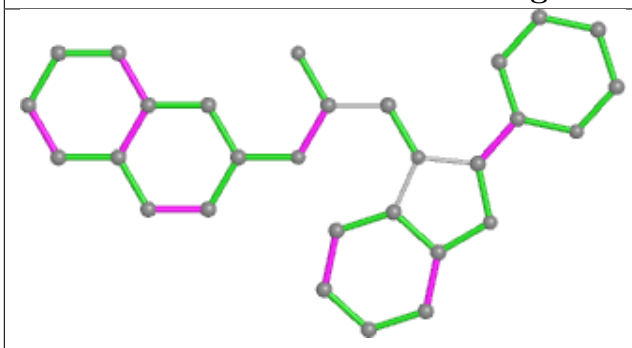
13 monomers are involved in 13 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	C	502	MLI	1	0
2	E	500	IMP	1	0
2	H	500	IMP	1	0
3	E	501	C91	1	0
2	B	500	IMP	1	0
2	D	500	IMP	1	0
3	A	501	C91	1	0
2	G	500	IMP	1	0
4	D	502	MLI	1	0
3	B	501	C91	1	0
3	F	501	C91	1	0
2	A	500	IMP	1	0
2	C	500	IMP	1	0

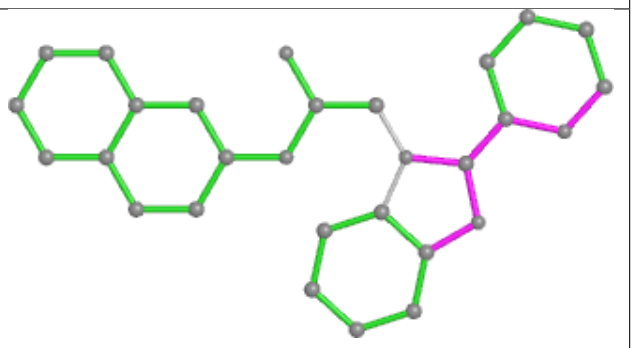
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

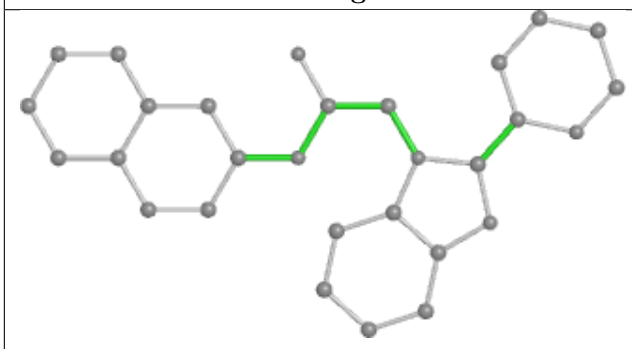
Ligand C91 G 501



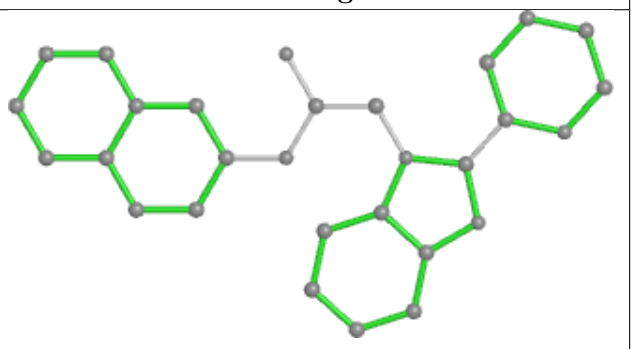
Bond lengths



Bond angles

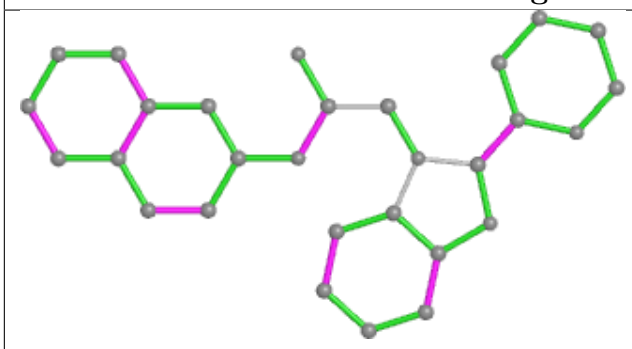


Torsions



Rings

Ligand C91 H 501



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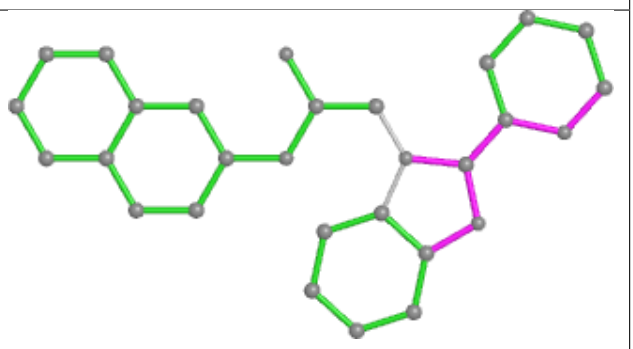
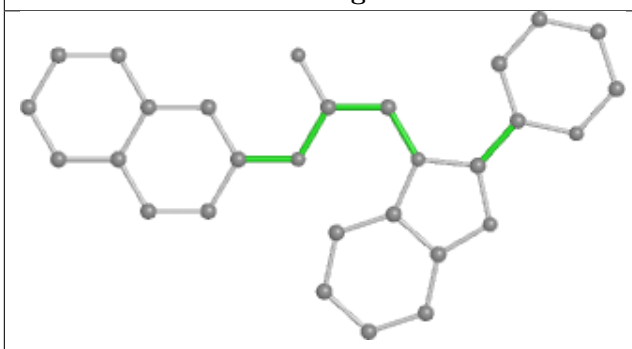
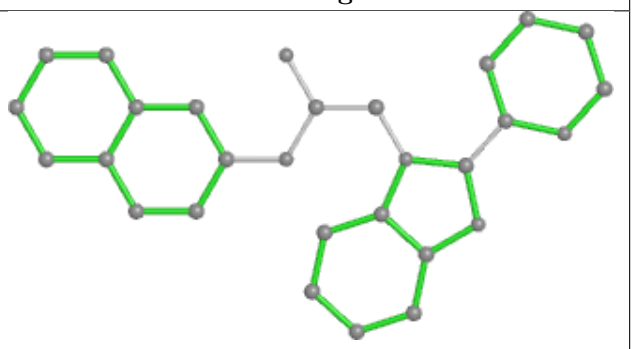


Bond angles

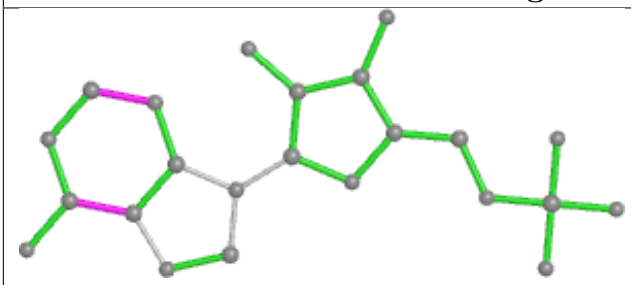


Torsions



Rings

Ligand IMP E 500



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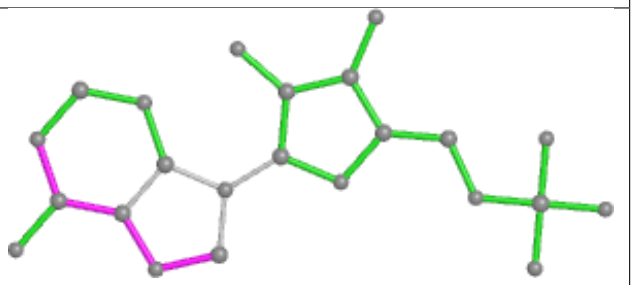
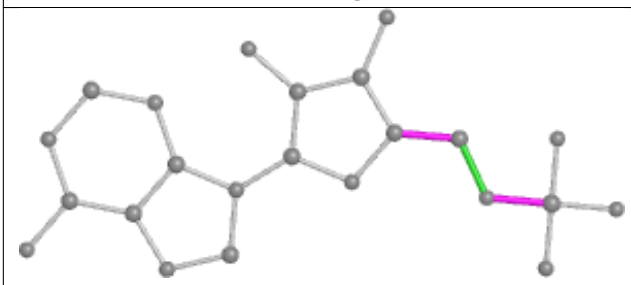


Bond angles



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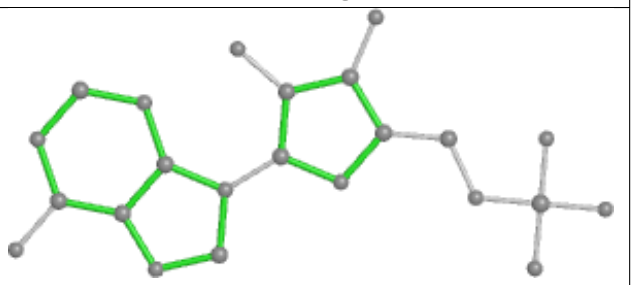
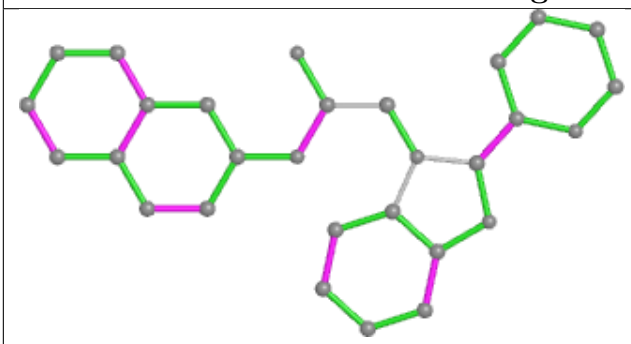


Rings

Ligand C91 C 501



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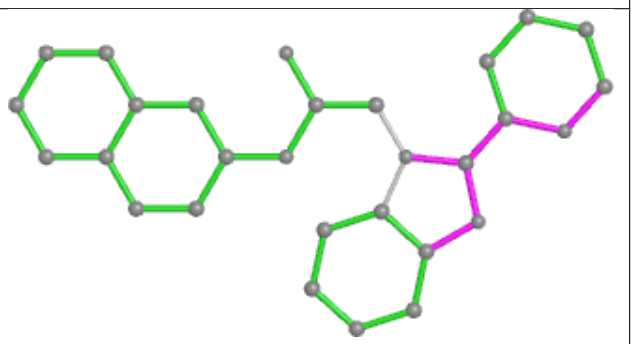
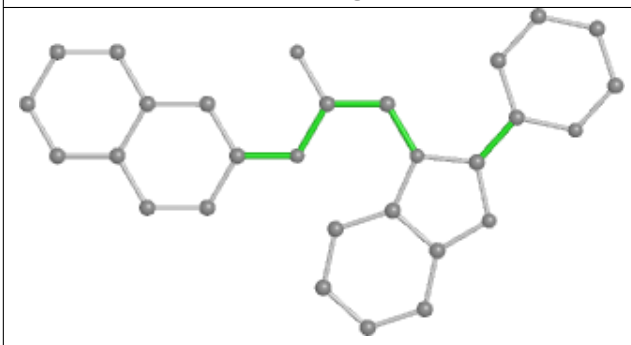
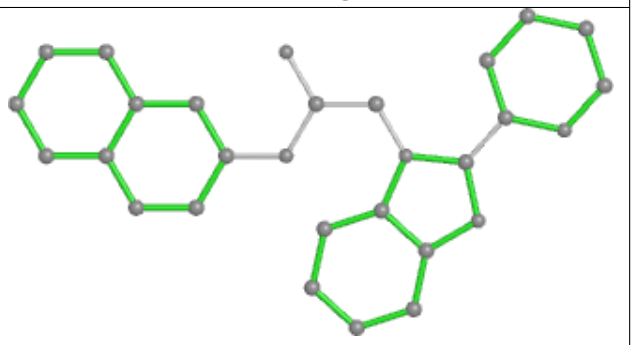


Bond angles

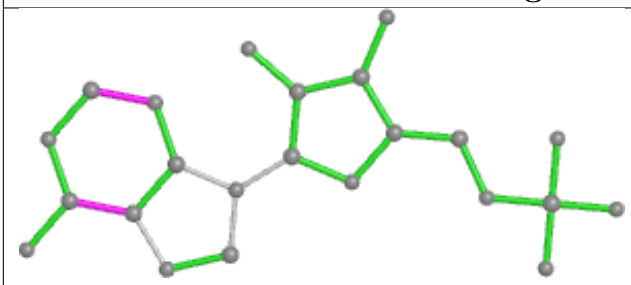


Torsions



Rings

Ligand IMP H 500



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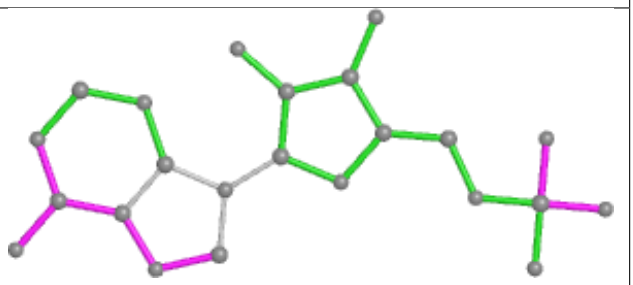

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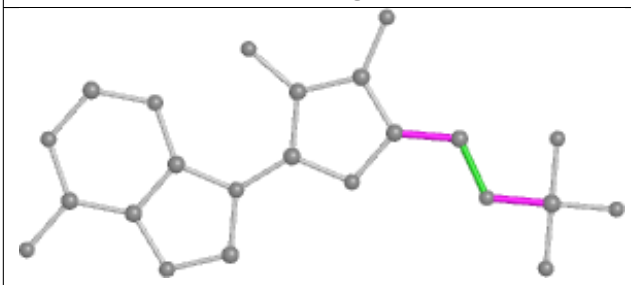

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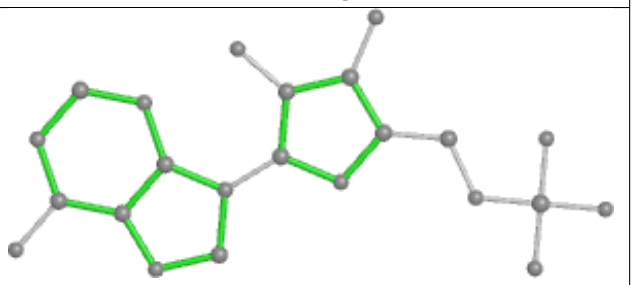

Rings

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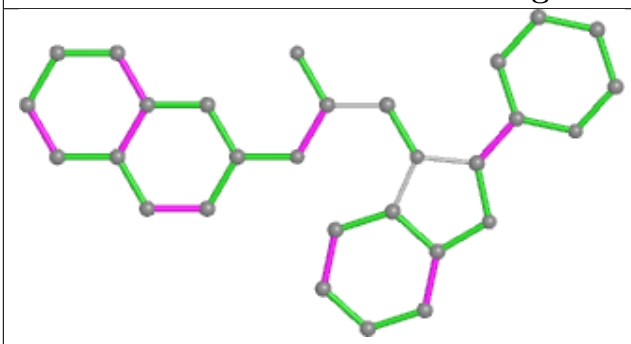

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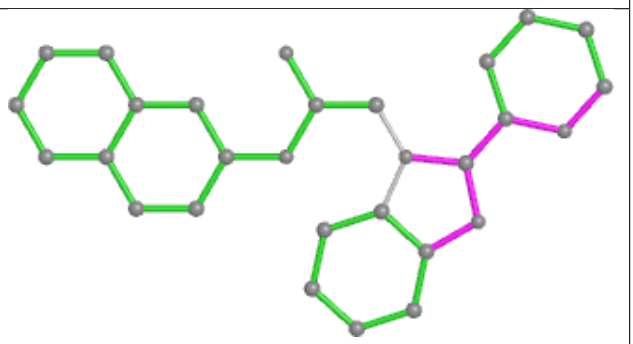
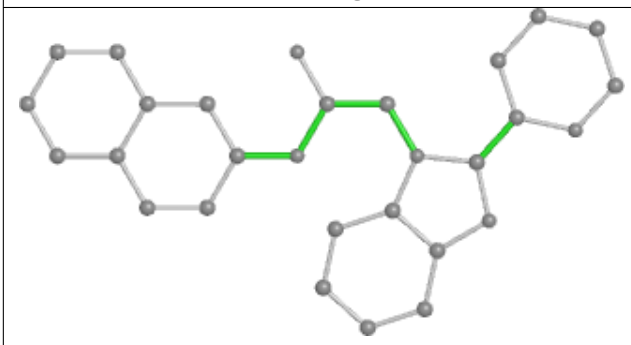


Bond angles



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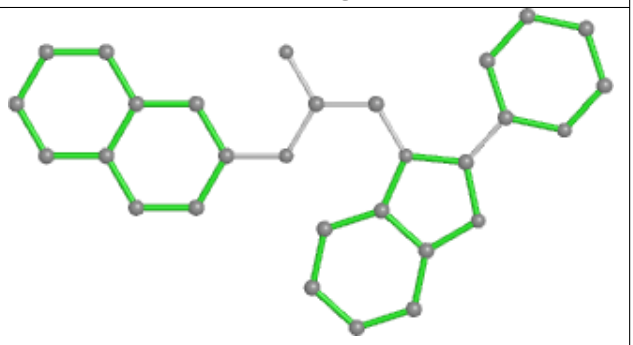
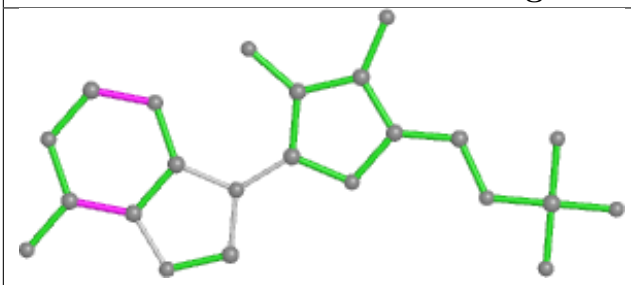


Rings

Ligand IMP B 500



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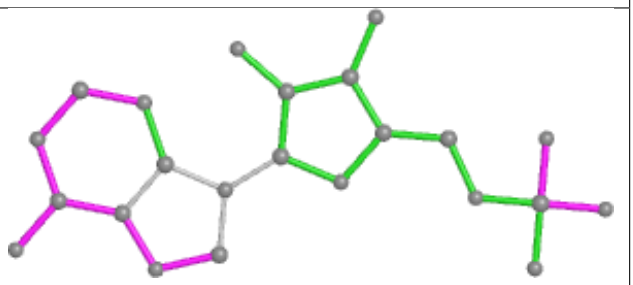
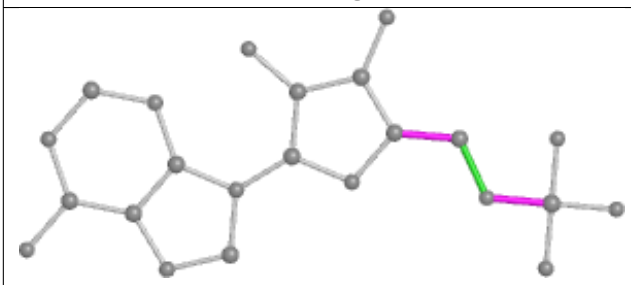


Bond angles



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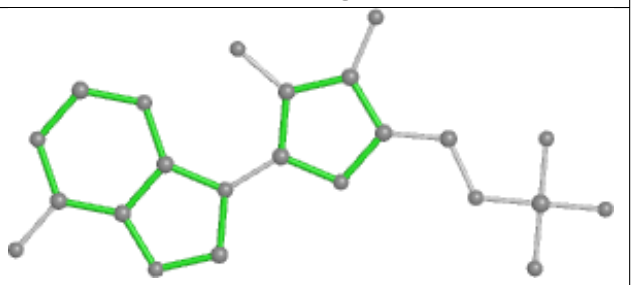
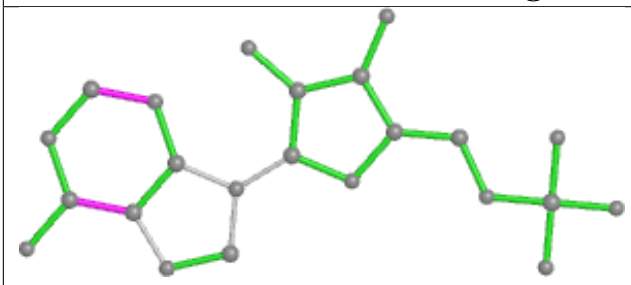
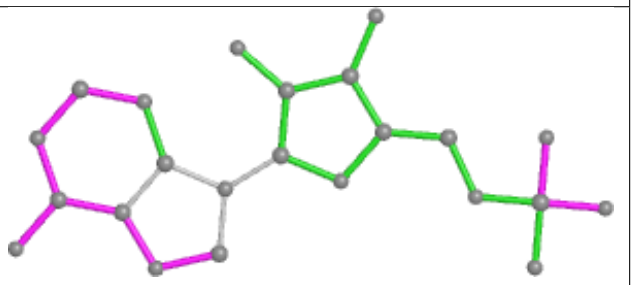


Rings

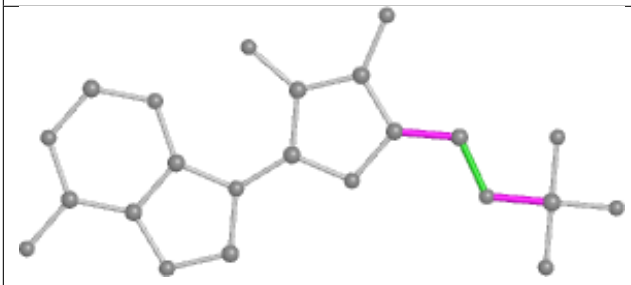
Ligand IMP D 500



Bond lengths



Bond angles



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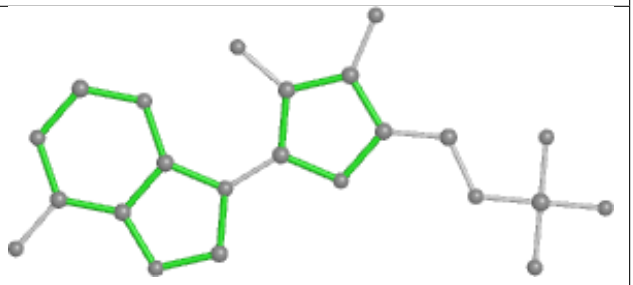
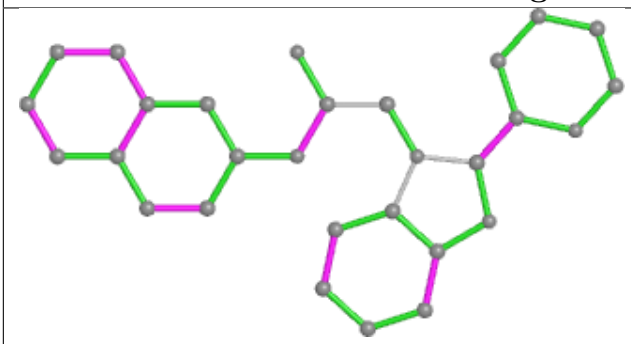


Rings

Ligand C91 A 501



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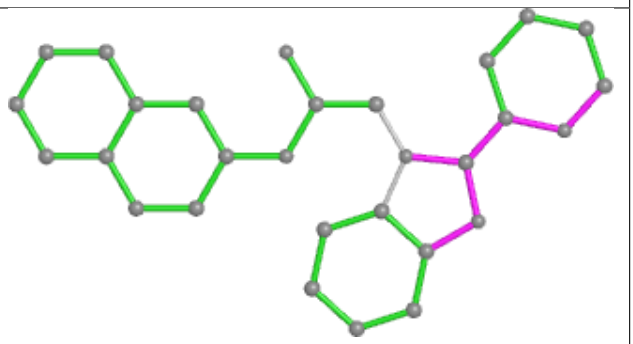
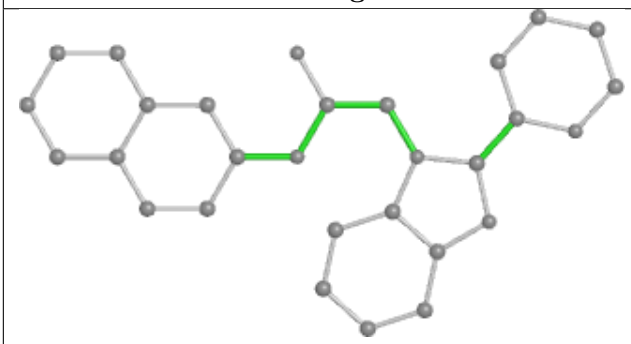


Bond angles



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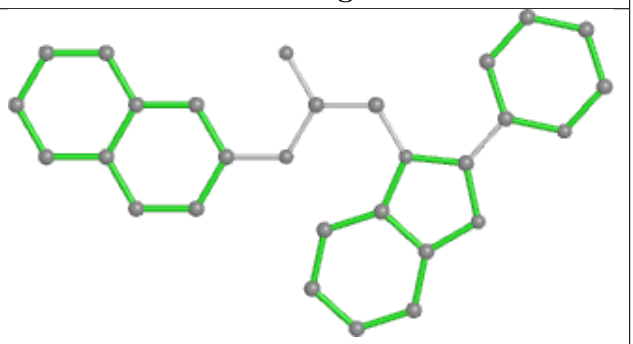
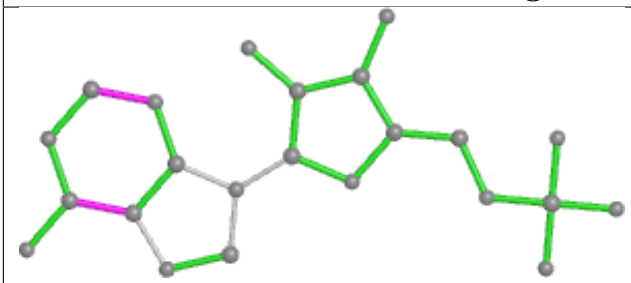


Rings

Ligand IMP G 500



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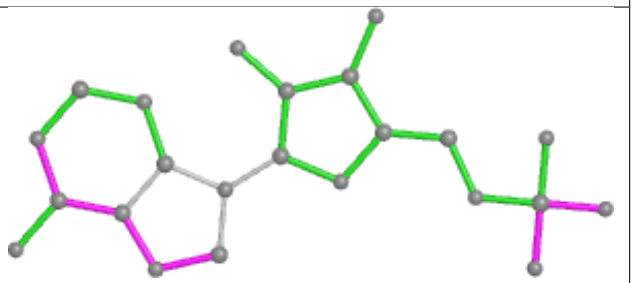
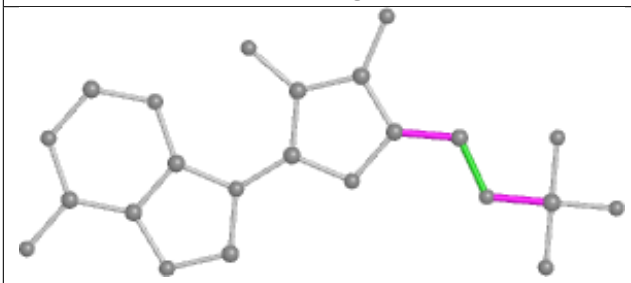
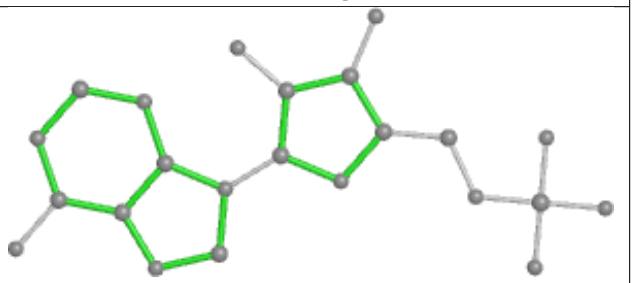


Bond angles

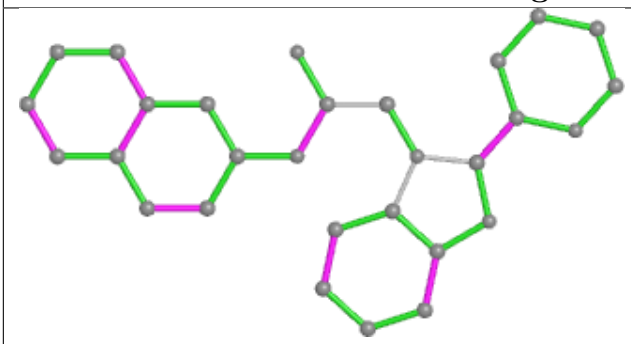


Torsions

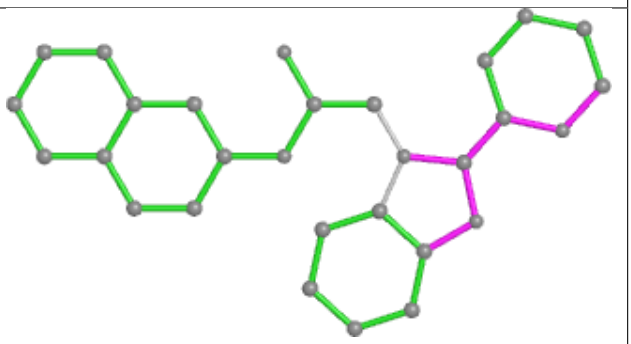


Rings

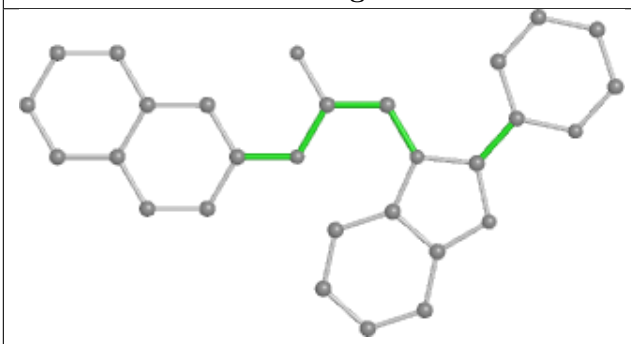
Ligand C91 B 501



Bond lengths



Bond angles



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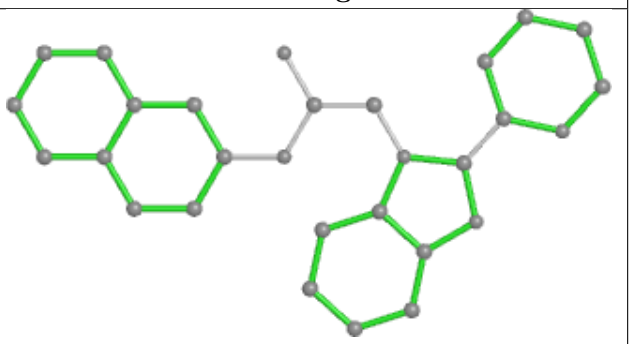
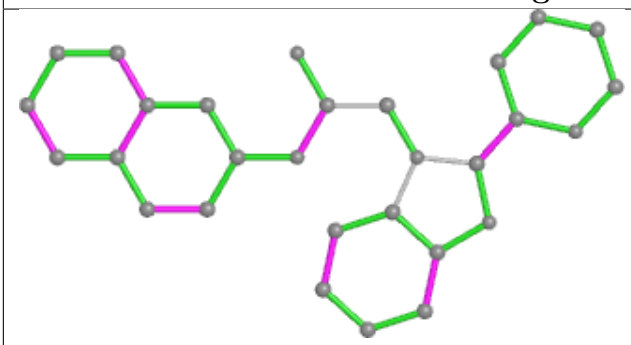
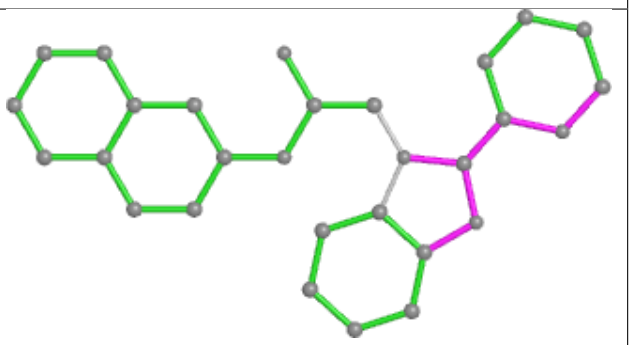


Rings

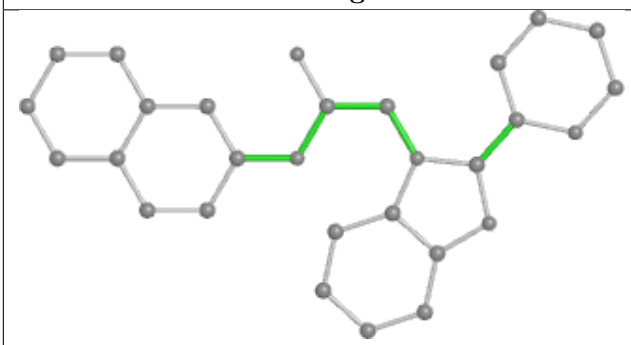
Ligand C91 F 501



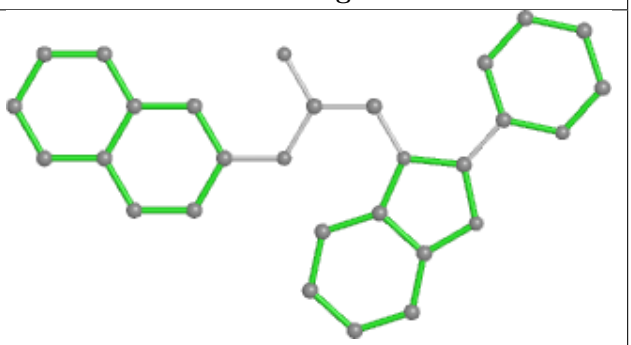
Bond lengths



Bond angles

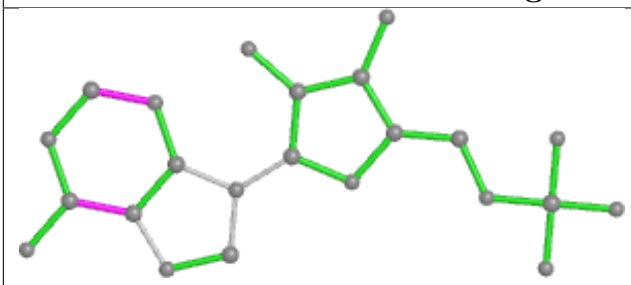


Torsions



Rings

Ligand IMP A 500



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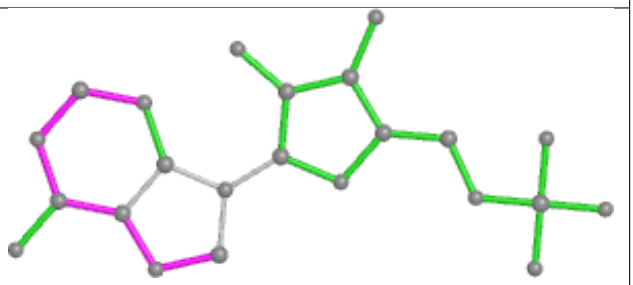
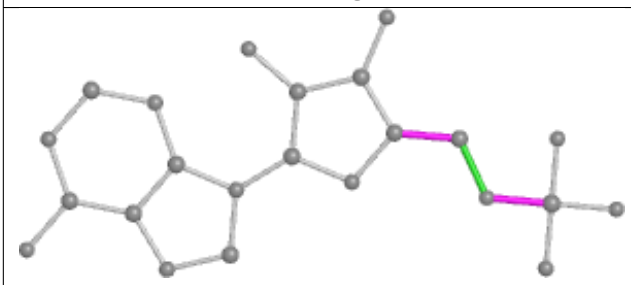


Bond angles



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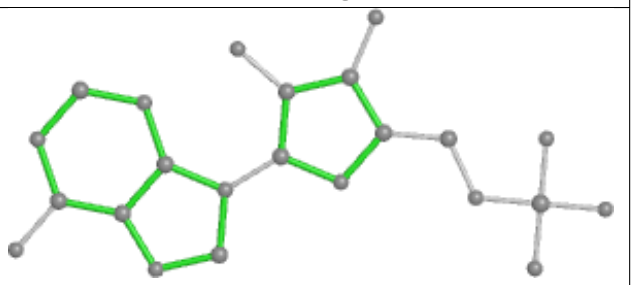
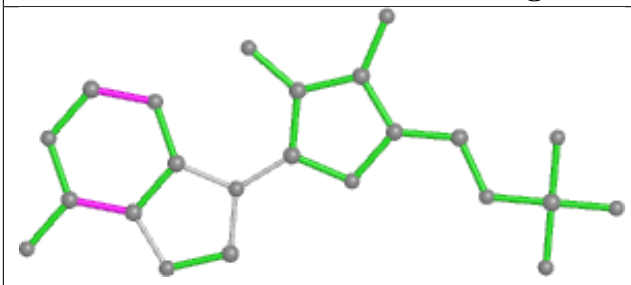


Rings

Ligand IMP C 500



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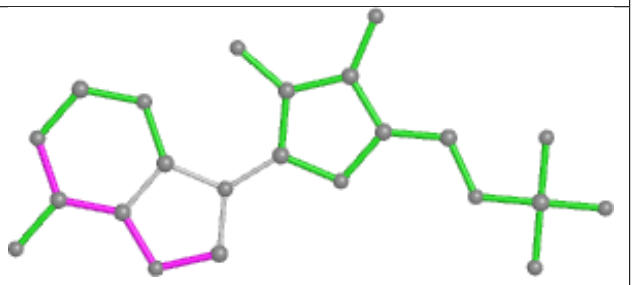
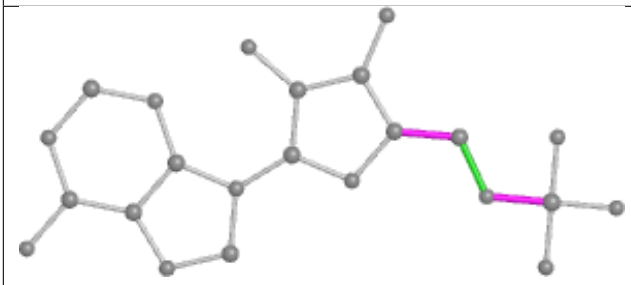
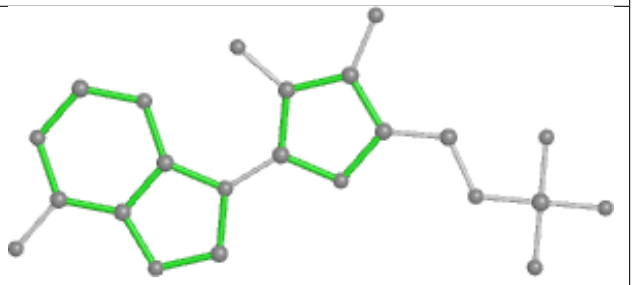


Bond angles

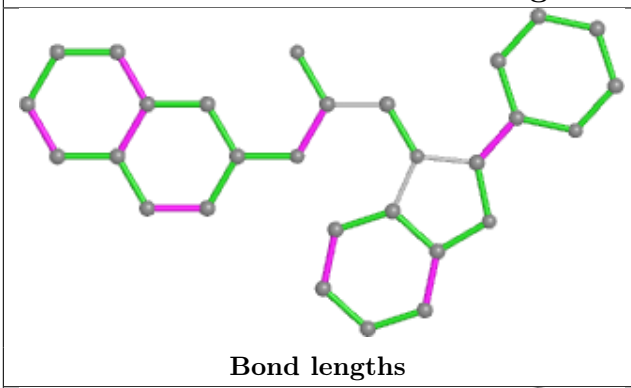


Torsions

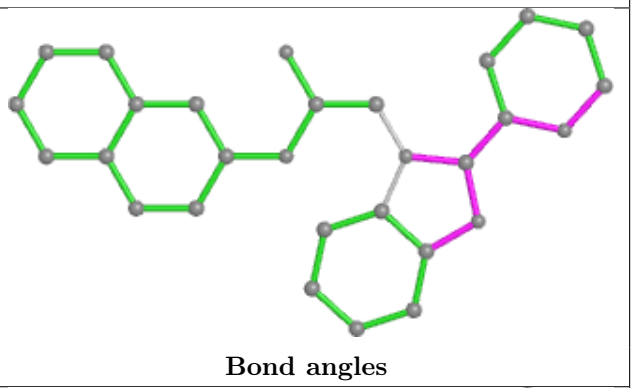


Rings

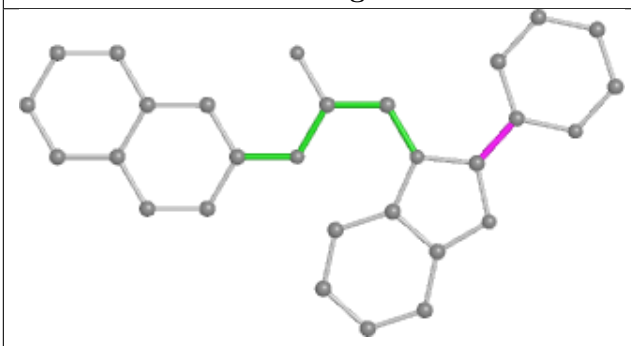
Ligand C91 D 501



Bond lengths



Bond angles



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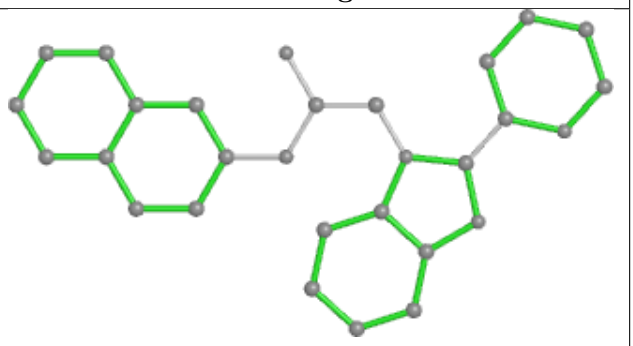
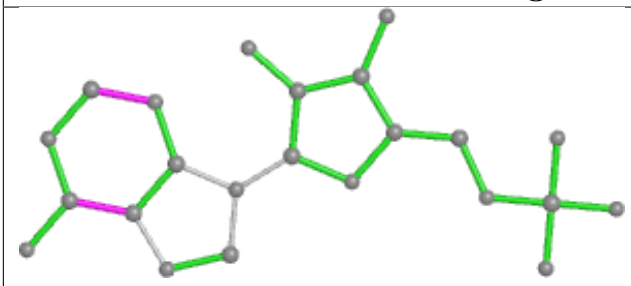
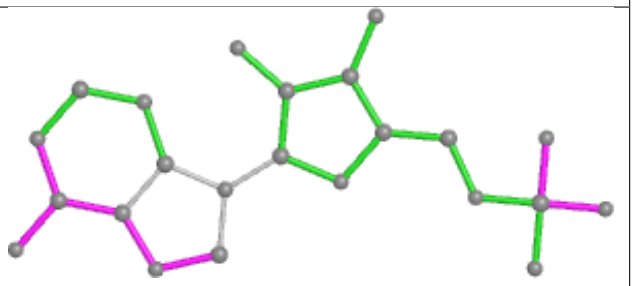


Rings

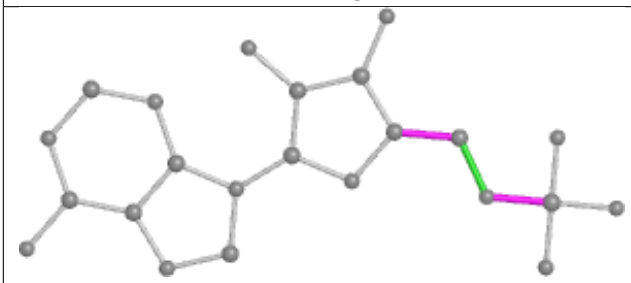
Ligand IMP F 500



Bond lengths



Bond angles



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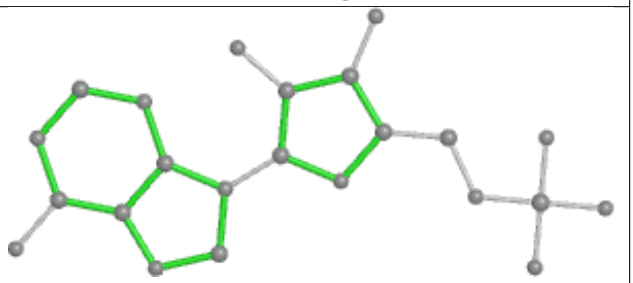

Rings

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled '#RSRZ> 2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95^{th} percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled 'Q< 0.9' lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å^2)	Q<0.9
1	A	350/384 (91%)	0.06	4 (1%)	80	79	27, 39, 64, 96	0
1	B	351/384 (91%)	0.07	1 (0%)	94	94	26, 40, 64, 78	1 (0%)
1	C	347/384 (90%)	0.12	3 (0%)	84	83	26, 39, 63, 82	0
1	D	347/384 (90%)	0.34	13 (3%)	41	37	25, 45, 67, 92	0
1	E	349/384 (90%)	0.19	3 (0%)	84	83	29, 43, 66, 92	0
1	F	348/384 (90%)	0.42	17 (4%)	29	26	28, 44, 66, 92	0
1	G	352/384 (91%)	0.22	6 (1%)	70	67	31, 44, 67, 94	1 (0%)
1	H	348/384 (90%)	0.34	8 (2%)	60	57	29, 48, 69, 94	1 (0%)
All	All	2792/3072 (90%)	0.22	55 (1%)	65	62	25, 43, 66, 96	3 (0%)

All (55) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	413	LEU	6.0
1	F	241	VAL	4.2
1	F	413	LEU	4.1
1	G	-2	SER	4.0
1	C	380	TYR	3.7
1	E	383	ARG	3.6
1	C	413	LEU	3.5
1	G	250	LEU	3.5
1	A	-2	SER	3.3
1	H	250	LEU	3.3
1	H	399	LYS	3.2
1	A	227	VAL	3.2
1	F	243	ALA	3.1
1	D	222	LEU	2.9
1	H	79[A]	GLU	2.9
1	D	221	LEU	2.8

Continued on next page...

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Mol	Chain	Res	Type	RSRZ
1	D	75	ASN	2.7
1	D	396	ALA	2.7
1	F	399	LYS	2.7
1	F	78	ILE	2.7
1	F	84	GLN	2.7
1	F	87	LYS	2.6
1	D	226	ALA	2.6
1	H	380	TYR	2.6
1	F	242	LYS	2.6
1	H	276	ILE	2.5
1	F	240	LEU	2.5
1	D	248	ILE	2.5
1	H	248	ILE	2.5
1	F	387	VAL	2.5
1	D	378	GLU	2.4
1	F	380	TYR	2.4
1	F	85	VAL	2.4
1	D	70	GLY	2.3
1	H	245	VAL	2.3
1	F	90	ARG	2.3
1	D	88	VAL	2.3
1	G	36	LEU	2.3
1	A	248	ILE	2.2
1	F	80	GLN	2.2
1	B	250	LEU	2.2
1	D	227	VAL	2.1
1	D	223	VAL	2.1
1	F	244	SER	2.1
1	C	90	ARG	2.1
1	G	400	GLY	2.1
1	E	278	ALA	2.1
1	F	226	ALA	2.1
1	D	72	ILE	2.1
1	F	57	ALA	2.1
1	H	88	VAL	2.0
1	G	449	GLN	2.0
1	E	378	GLU	2.0
1	A	399	LYS	2.0
1	G	37	SER	2.0

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

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

6.3 Carbohydrates (i)

There are no monosaccharides in this entry.

6.4 Ligands (i)

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

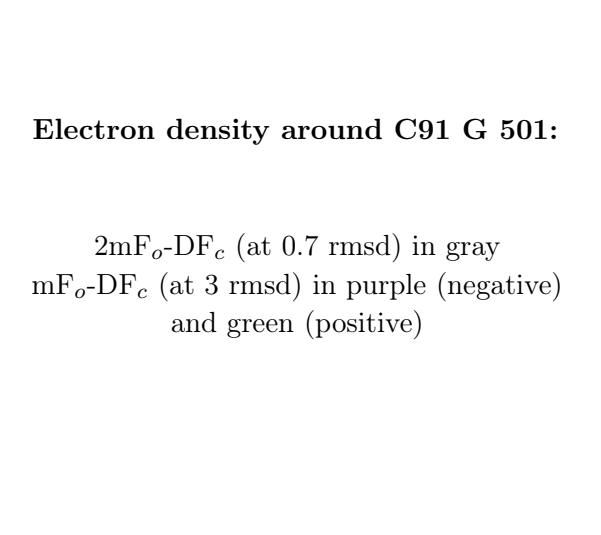
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
4	MLI	C	502	7/7	0.80	0.34	90,91,94,97	0
4	MLI	B	502	7/7	0.83	0.39	58,68,72,74	0
4	MLI	C	503	7/7	0.83	0.31	74,76,81,81	0
4	MLI	G	502	7/7	0.83	0.27	77,77,80,83	0
4	MLI	H	502	7/7	0.83	0.24	70,73,75,77	0
4	MLI	D	502	7/7	0.88	0.28	69,70,74,74	0
4	MLI	B	503	7/7	0.89	0.21	72,72,75,76	0
4	MLI	E	502	7/7	0.91	0.35	66,71,75,79	0
4	MLI	A	502	7/7	0.92	0.37	57,59,63,67	0
3	C91	G	501	29/29	0.93	0.26	38,45,52,53	0
3	C91	F	501	29/29	0.93	0.25	47,52,56,57	0
3	C91	E	501	29/29	0.94	0.18	34,43,47,50	0
3	C91	C	501	29/29	0.94	0.28	43,50,60,60	0
3	C91	D	501	29/29	0.94	0.27	47,55,66,67	0
3	C91	B	501	29/29	0.95	0.23	29,35,39,39	0
3	C91	A	501	29/29	0.95	0.20	34,44,48,49	0
3	C91	H	501	29/29	0.95	0.22	45,52,62,63	0
2	IMP	F	500	23/23	0.96	0.13	27,34,38,39	0
2	IMP	D	500	23/23	0.97	0.15	32,38,41,43	0
2	IMP	B	500	23/23	0.97	0.16	27,32,37,38	0
2	IMP	H	500	23/23	0.97	0.15	35,42,44,45	0
2	IMP	C	500	23/23	0.97	0.16	24,36,42,44	0
2	IMP	E	500	23/23	0.98	0.13	22,34,38,41	0
2	IMP	A	500	23/23	0.98	0.13	25,34,38,42	0
2	IMP	G	500	23/23	0.98	0.13	25,36,46,49	0

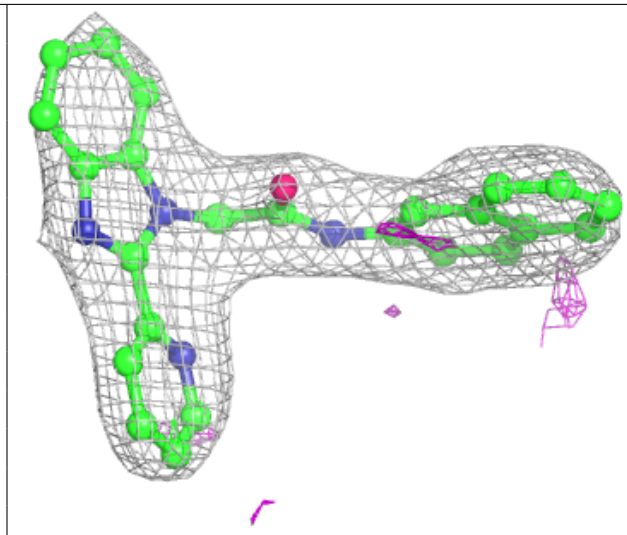
The following is a graphical depiction of the model fit to experimental electron density of all

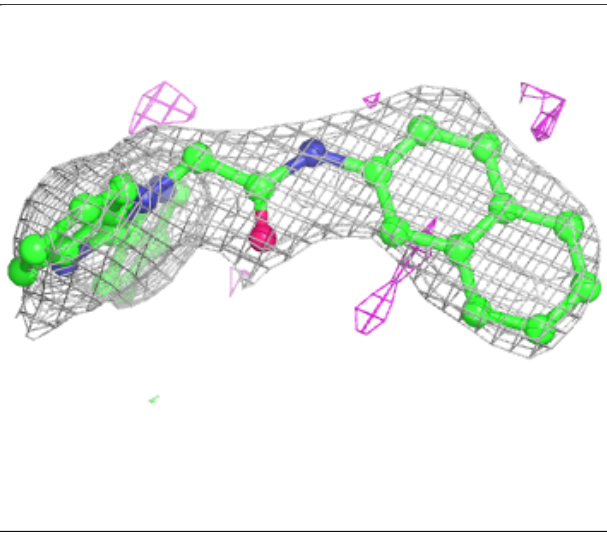
instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

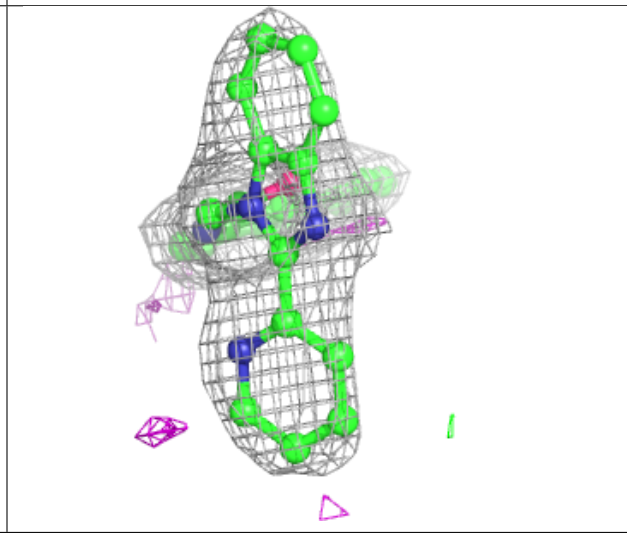
Electron density around C91 G 501:

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



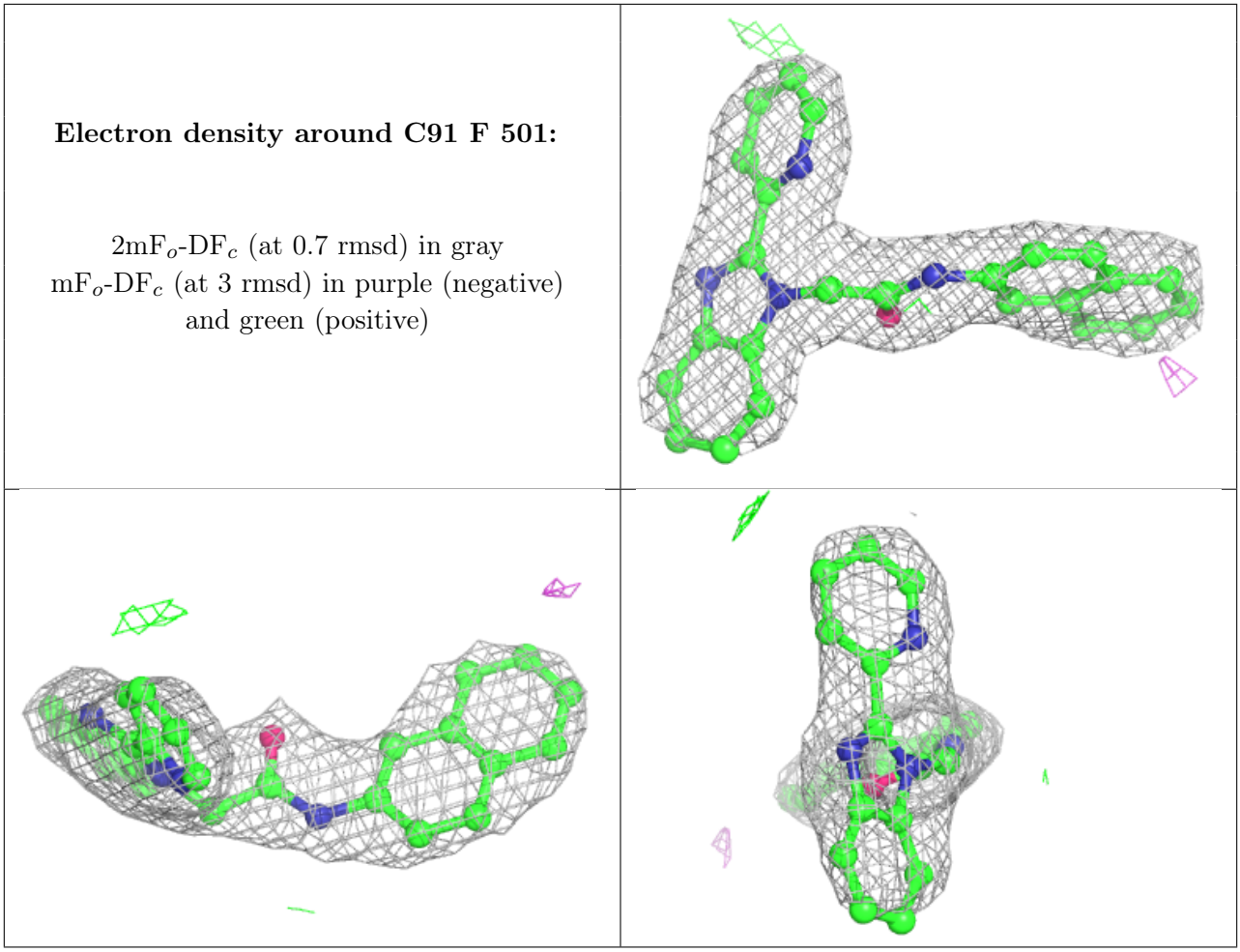






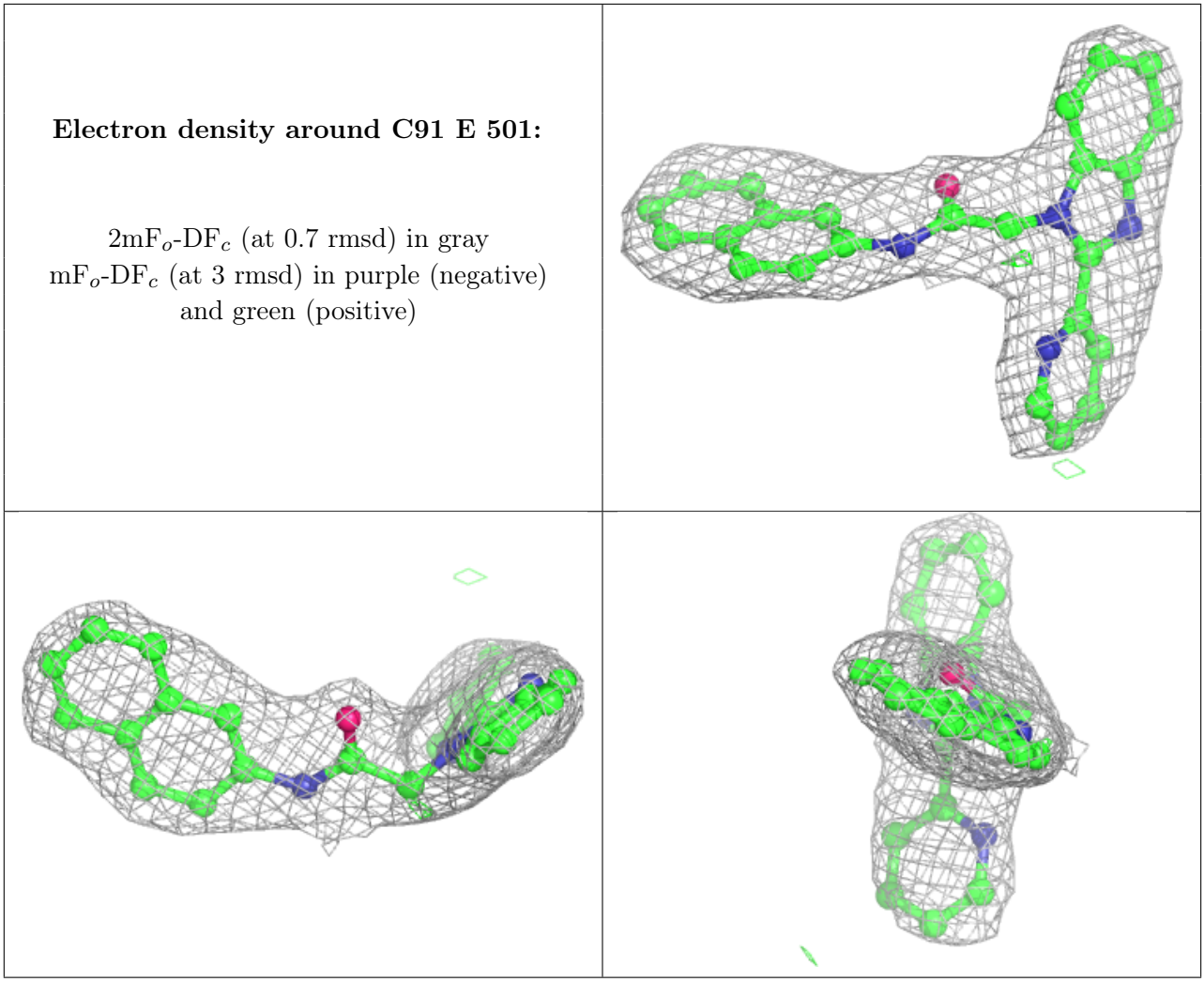
Electron density around C91 F 501:

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



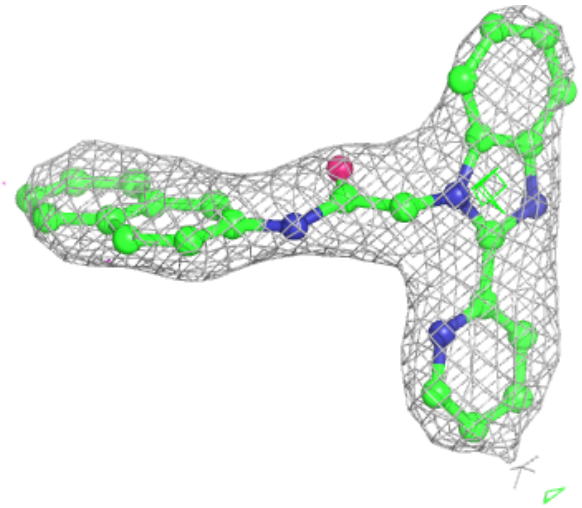
Electron density around C91 E 501:

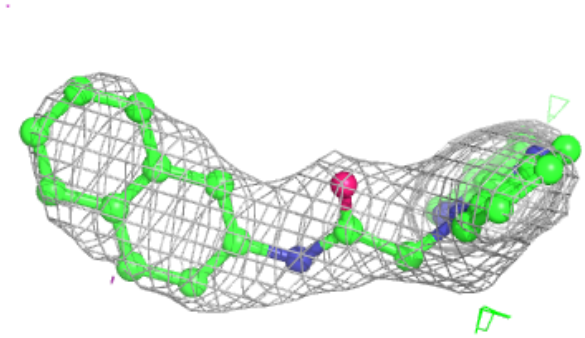
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

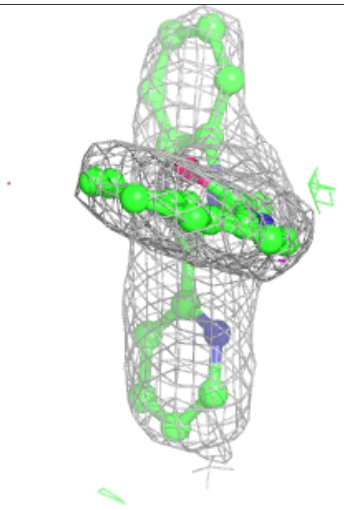


Electron density around C91 C 501:

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

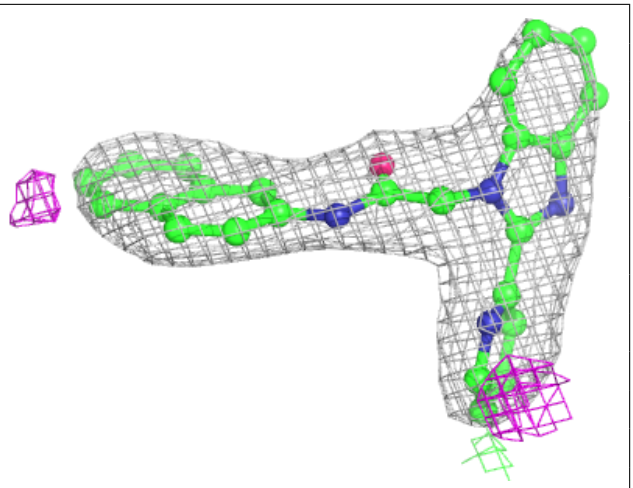


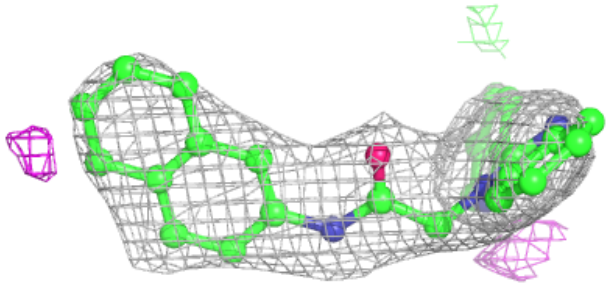


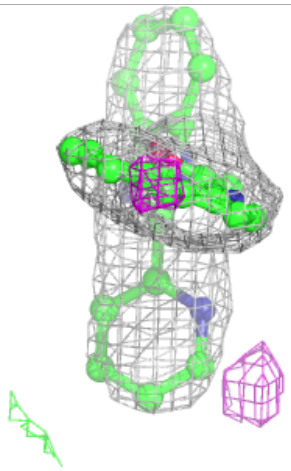


Electron density around C91 D 501:

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

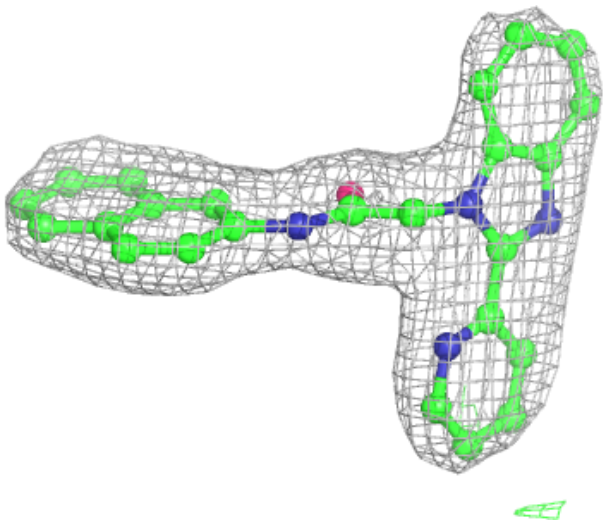


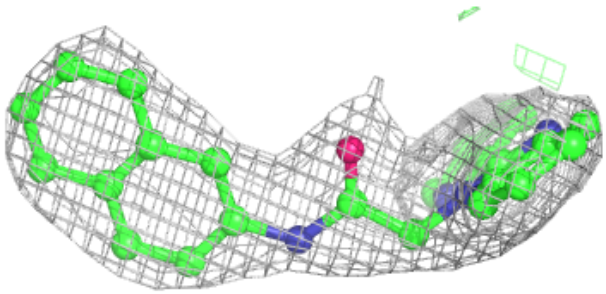


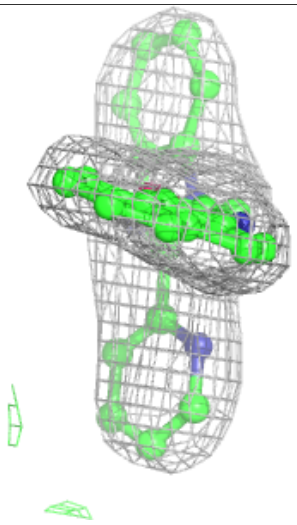


Electron density around C91 B 501:

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

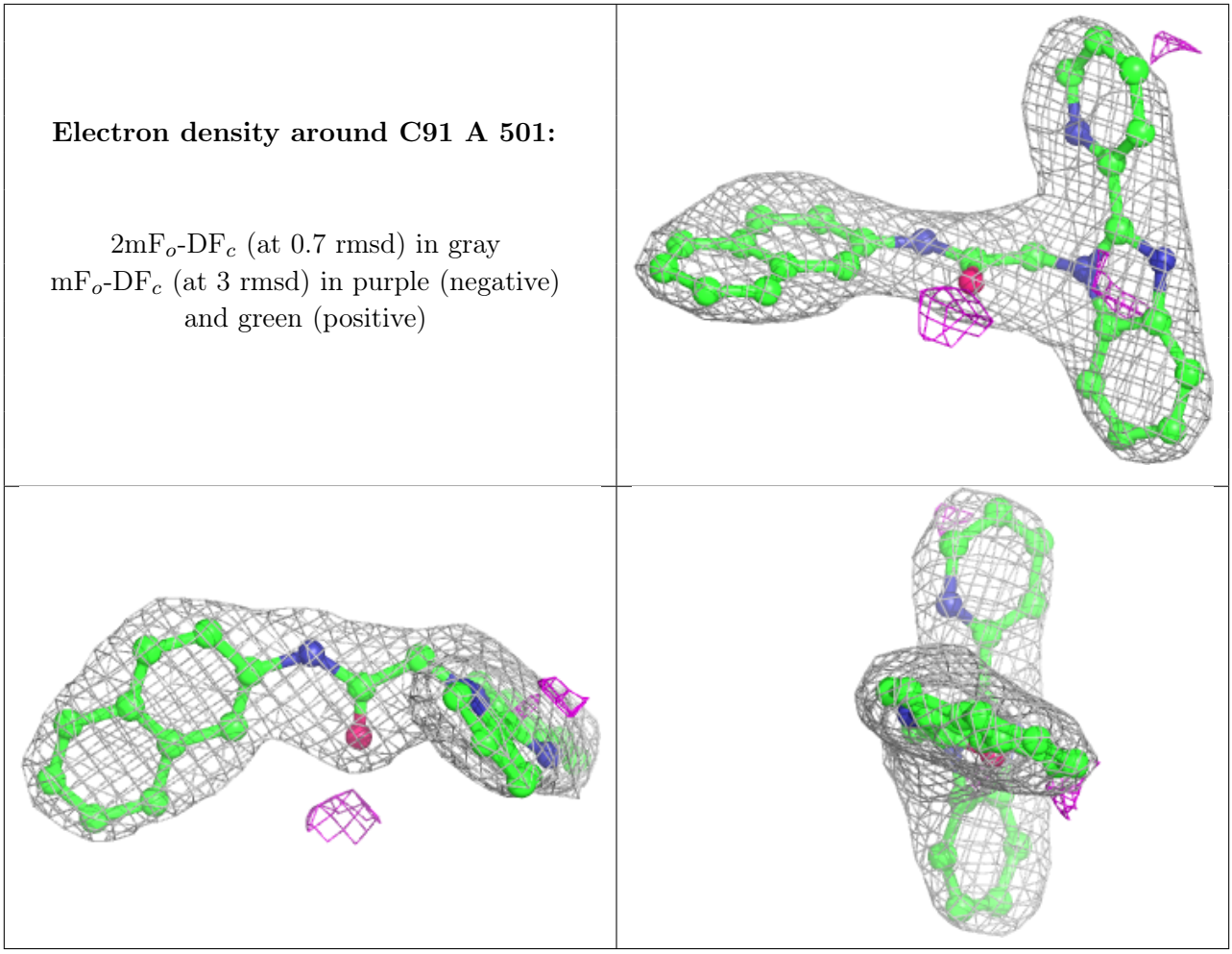






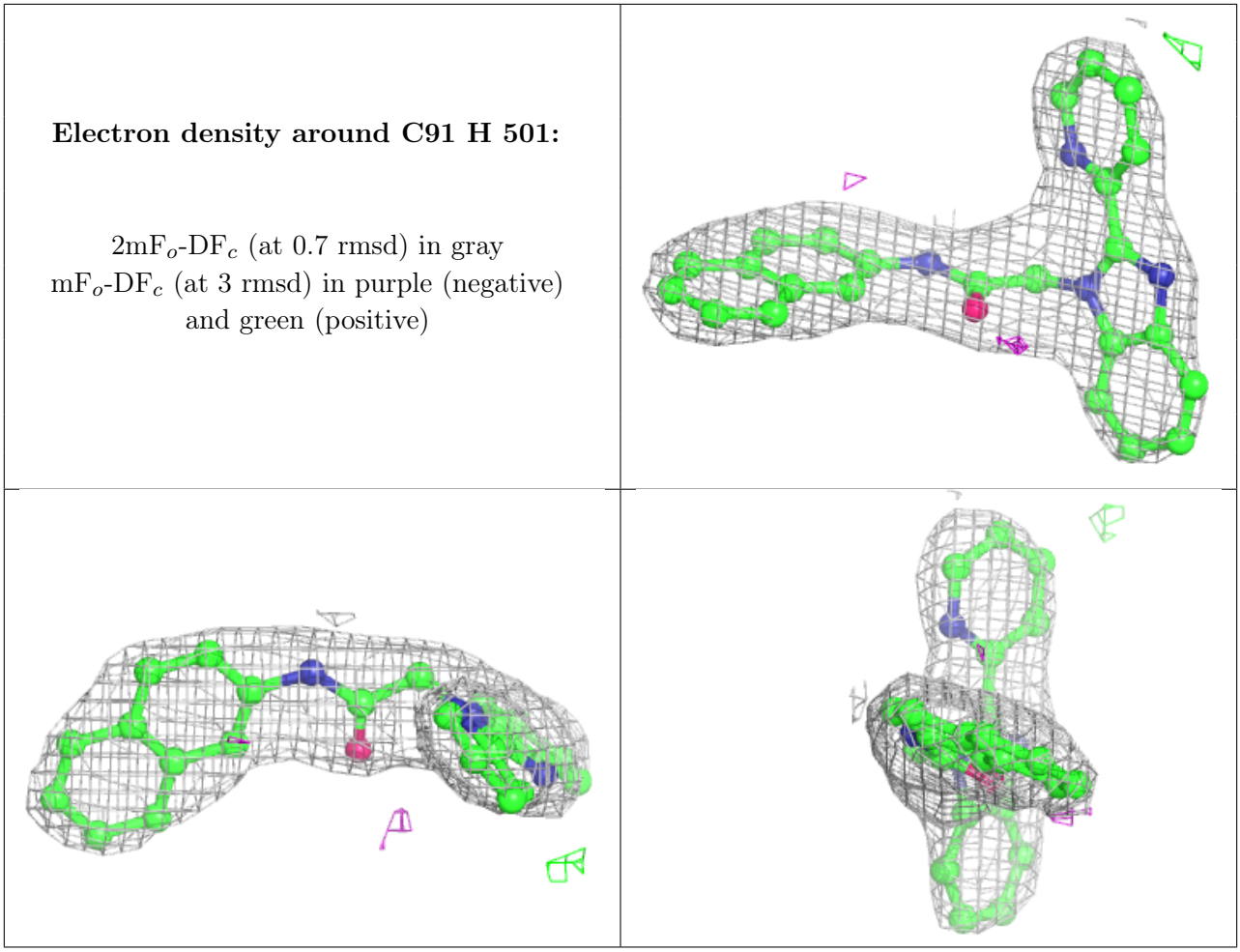
Electron density around C91 A 501:

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



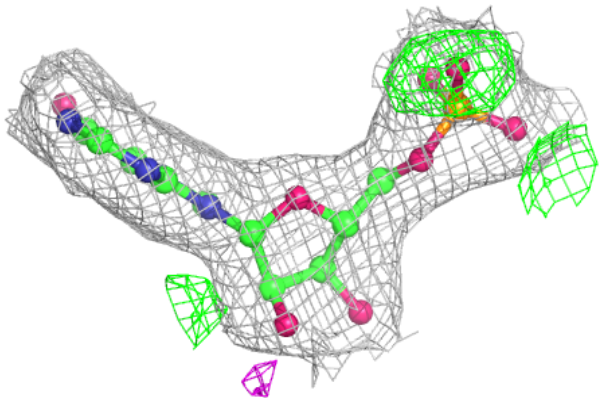
Electron density around C91 H 501:

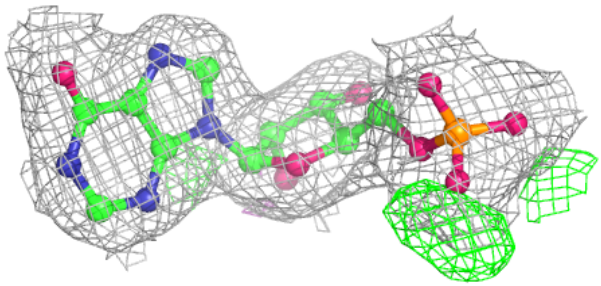
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

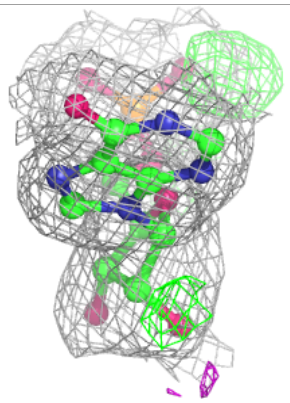


Electron density around IMP F 500:

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

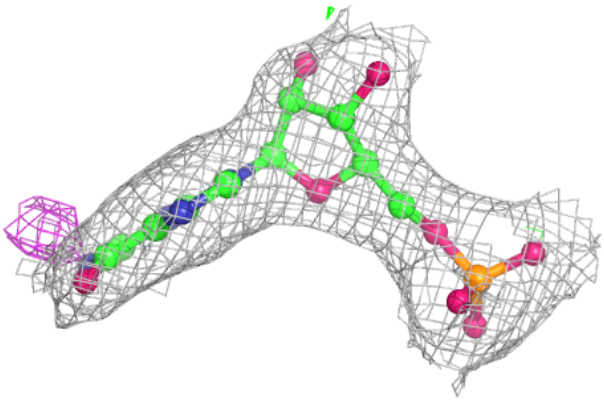


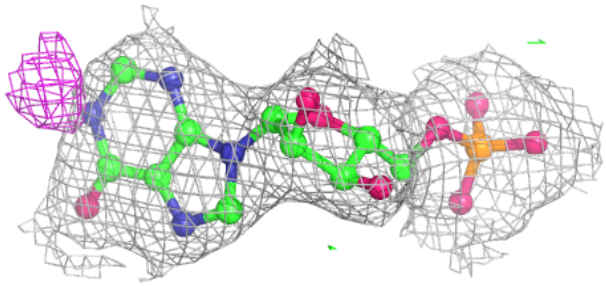


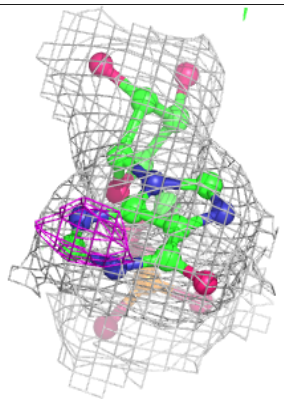


Electron density around IMP D 500:

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

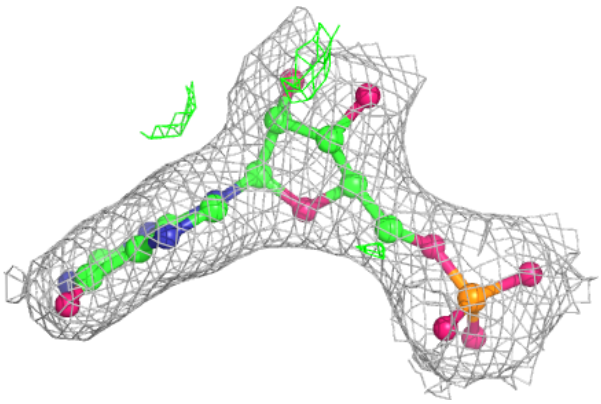


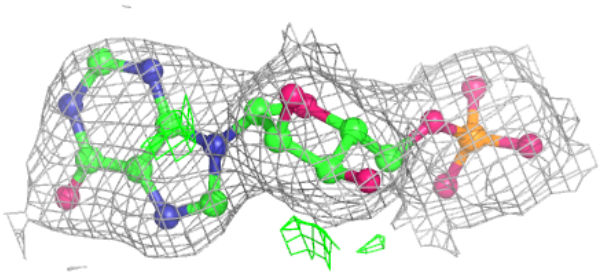


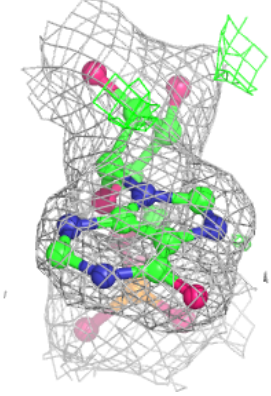


Electron density around IMP B 500:

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

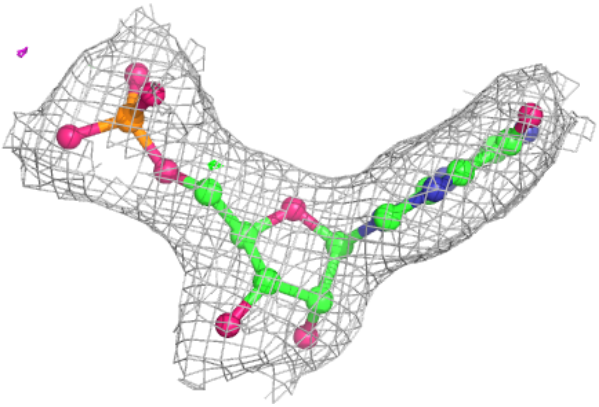


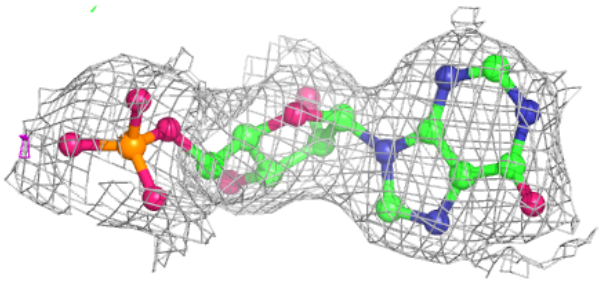


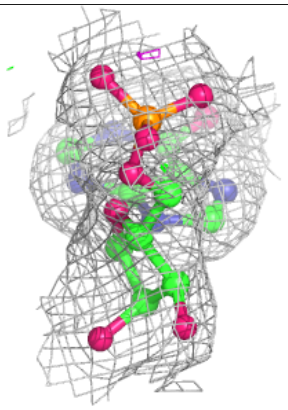


Electron density around IMP H 500:

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

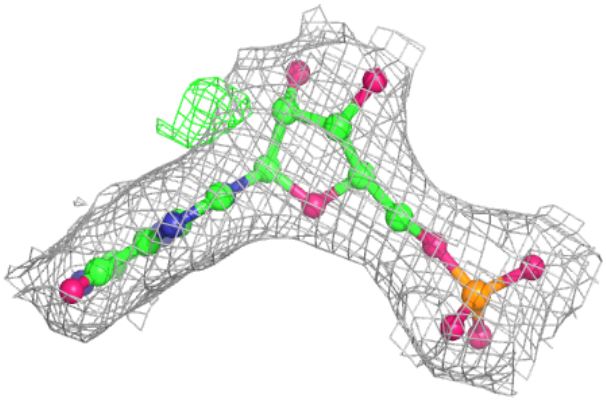


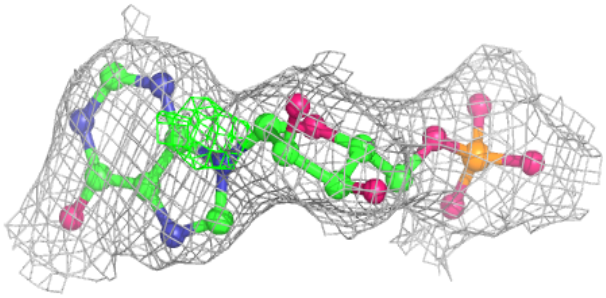


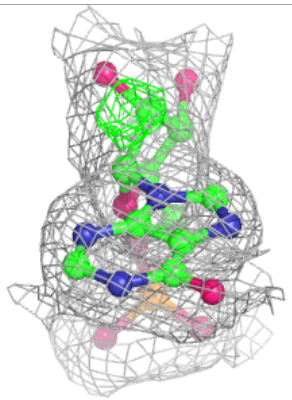


Electron density around IMP C 500:

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

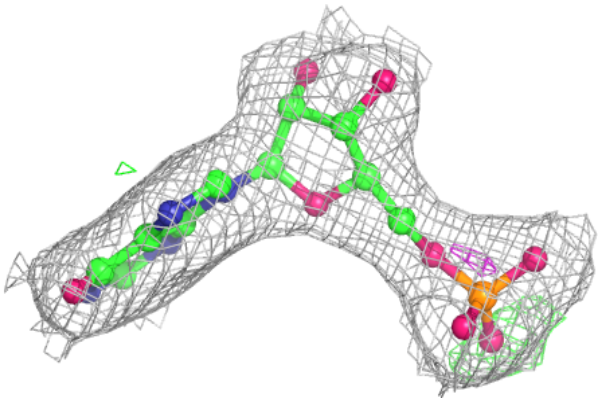


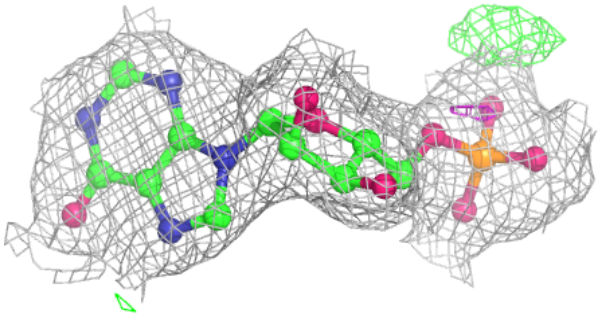


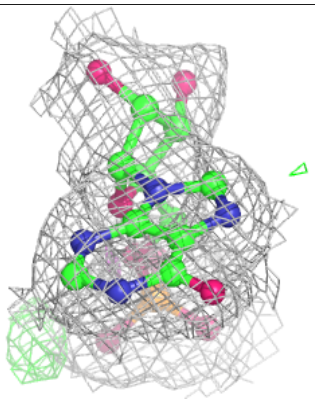


Electron density around IMP E 500:

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

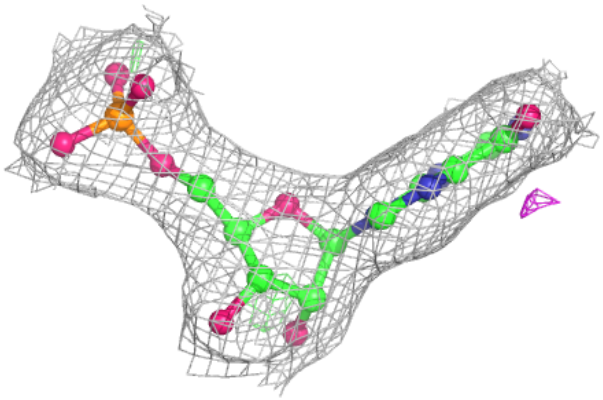


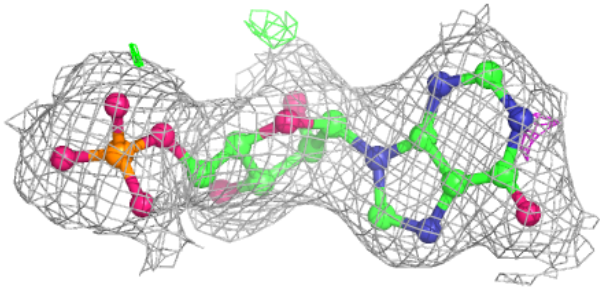


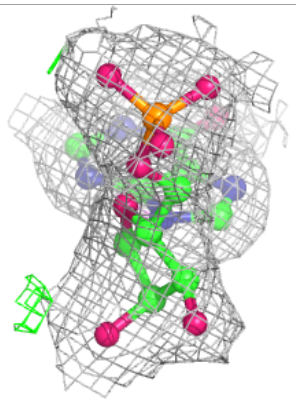


Electron density around IMP A 500:

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

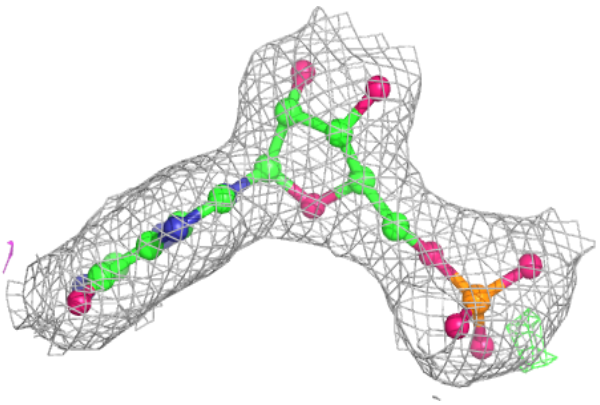


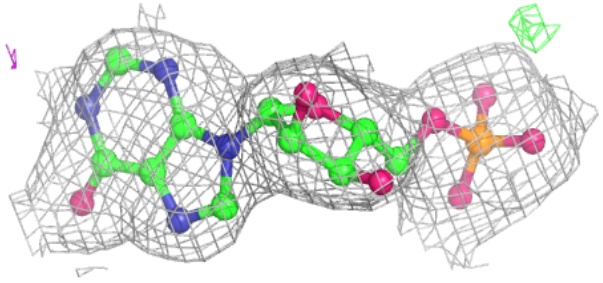


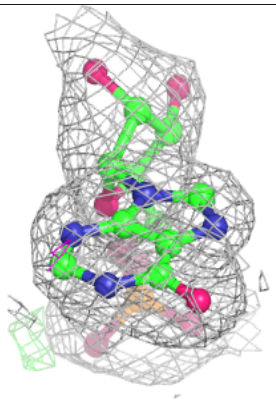


Electron density around IMP G 500:

$2mF_o$-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.