



Full wwPDB X-ray Structure Validation Report ⓘ

May 15, 2020 – 09:06 pm BST

PDB ID : 4QM1
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 D67
Authors : Kim, Y.; Makowska-Grzyska, M.; Gu, M.; Mandapati, K.; Gollapalli, D.; Gorla, S.K.; Zhang, M.; Hedstrom, L.; Anderson, W.F.; Joachimiak, A.; CSGID; Center for Structural Genomics of Infectious Diseases (CSGID)
Deposited on : 2014-06-14
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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.11
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

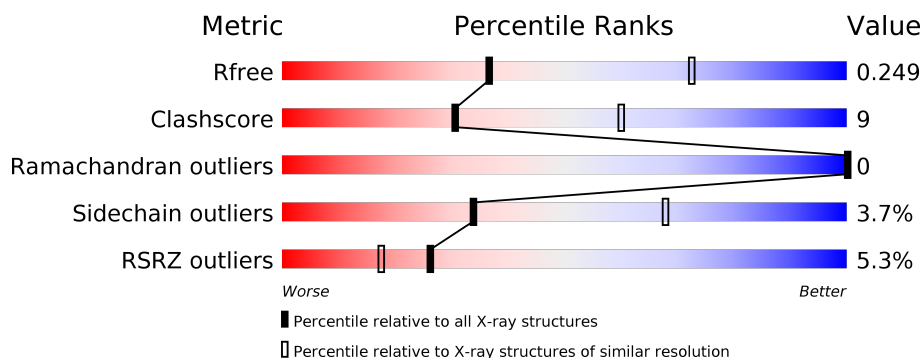
1 Overall quality at a glance

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	384	<div> <div>4%</div> <div> <div></div> <div>68%</div> <div>18%</div> <div>•</div> <div>12%</div> </div> </div>
1	B	384	<div> <div>5%</div> <div> <div></div> <div>73%</div> <div>16%</div> <div>•</div> <div>10%</div> </div> </div>
1	C	384	<div> <div>3%</div> <div> <div></div> <div>68%</div> <div>17%</div> <div>•</div> <div>14%</div> </div> </div>
1	D	384	<div> <div>6%</div> <div> <div></div> <div>65%</div> <div>21%</div> <div>•</div> <div>13%</div> </div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 10079 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	337	Total	C	N	O	S	0	0	0
			2462	1548	429	469	16			
1	B	345	Total	C	N	O	S	0	0	0
			2525	1587	441	481	16			
1	C	332	Total	C	N	O	S	0	0	0
			2417	1519	422	460	16			
1	D	336	Total	C	N	O	S	0	0	0
			2450	1539	426	469	16			

There are 104 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
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

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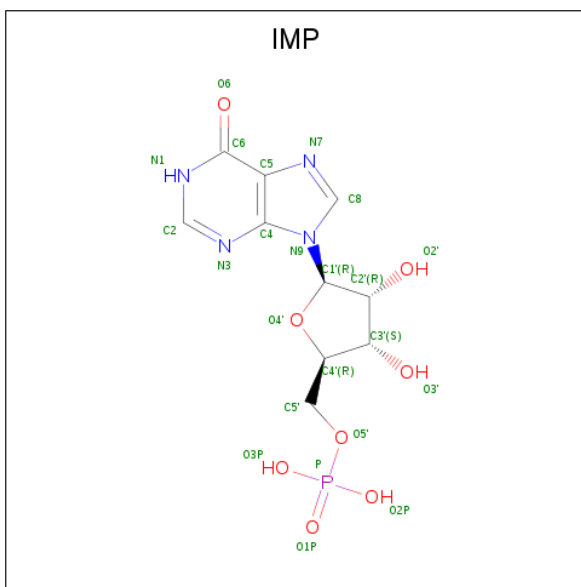
Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	SER	-	EXPRESSION TAG	UNP Q81W29
A	-1	ASN	-	EXPRESSION TAG	UNP Q81W29
A	0	ALA	-	EXPRESSION TAG	UNP Q81W29
A	92	GLY	-	EXPRESSION TAG	UNP Q81W29
A	220	GLY	-	EXPRESSION TAG	UNP Q81W29
B	-23	MET	-	EXPRESSION TAG	UNP Q81W29
B	-22	HIS	-	EXPRESSION TAG	UNP 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	-	EXPRESSION TAG	UNP Q81W29
B	220	GLY	-	EXPRESSION TAG	UNP Q81W29
C	-23	MET	-	EXPRESSION TAG	UNP Q81W29
C	-22	HIS	-	EXPRESSION TAG	UNP Q81W29
C	-21	HIS	-	EXPRESSION TAG	UNP Q81W29
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

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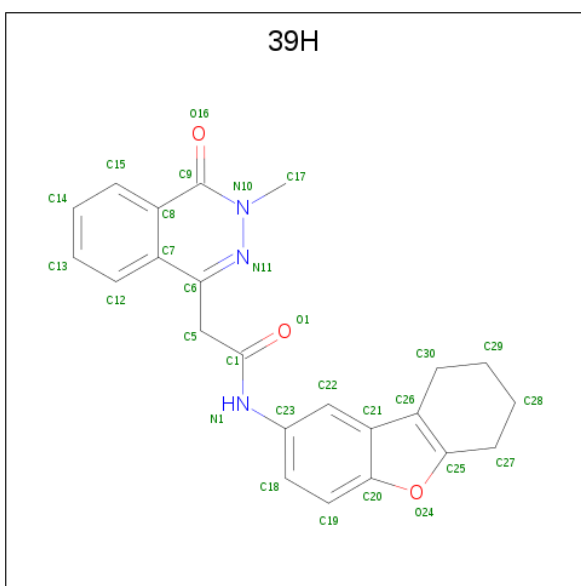
Chain	Residue	Modelled	Actual	Comment	Reference
C	-12	ASP	-	EXPRESSION TAG	UNP Q81W29
C	-11	LEU	-	EXPRESSION TAG	UNP Q81W29
C	-10	GLY	-	EXPRESSION TAG	UNP Q81W29
C	-9	THR	-	EXPRESSION TAG	UNP Q81W29
C	-8	GLU	-	EXPRESSION TAG	UNP Q81W29
C	-7	ASN	-	EXPRESSION TAG	UNP Q81W29
C	-6	LEU	-	EXPRESSION TAG	UNP 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	-	EXPRESSION TAG	UNP Q81W29
C	220	GLY	-	EXPRESSION TAG	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
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	-	EXPRESSION TAG	UNP Q81W29
D	220	GLY	-	EXPRESSION TAG	UNP Q81W29

- Molecule 2 is INOSINIC ACID (three-letter code: IMP) (formula: C₁₀H₁₃N₄O₈P).



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		

- Molecule 3 is 2-(3-methyl-4-oxo-3,4-dihydrophthalazin-1-yl)-N-(6,7,8,9-tetrahydrodibenzo[b,d]furan-2-yl)acetamide (three-letter code: 39H) (formula: $C_{23}H_{21}N_3O_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			29	23	3	3		
3	B	1	Total	C	N	O	0	0
			29	23	3	3		
3	C	1	Total	C	N	O	0	0
			29	23	3	3		
3	C	1	Total	C	N	O	0	0
			29	23	3	3		

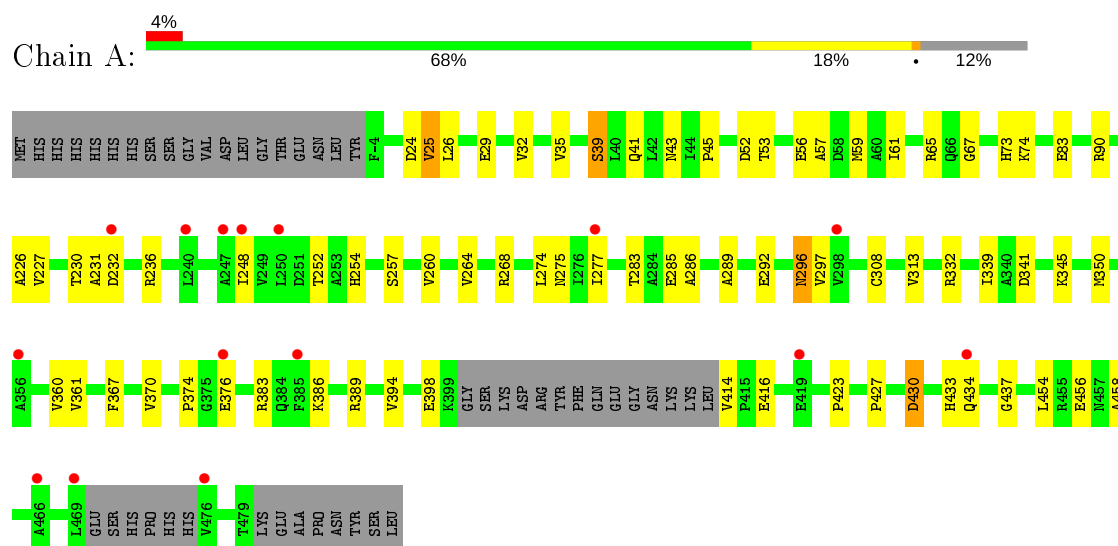
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	5	Total	O	0	0
			5	5		
4	B	4	Total	O	0	0
			4	4		
4	C	4	Total	O	0	0
			4	4		
4	D	4	Total	O	0	0
			4	4		

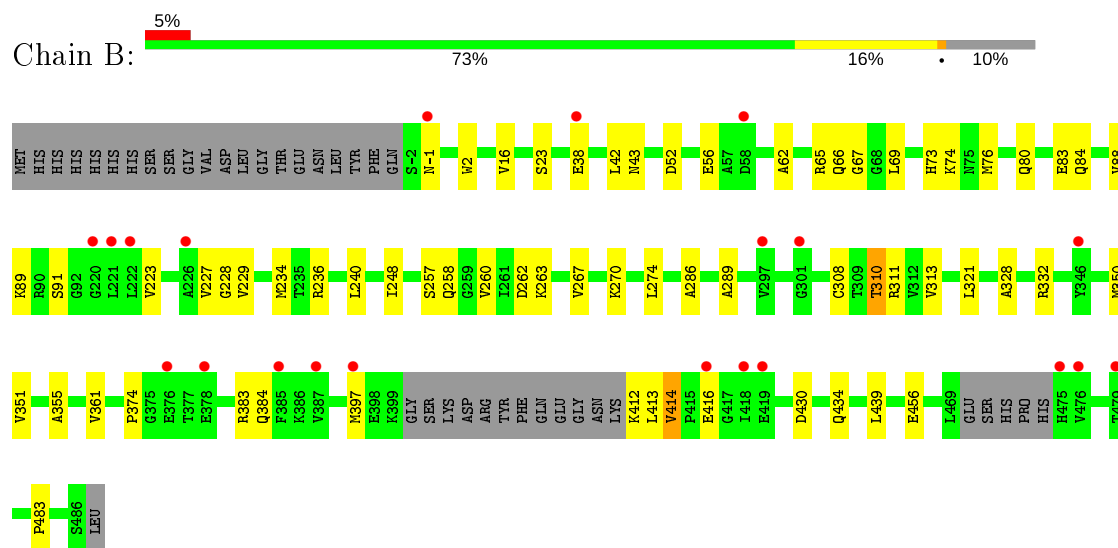
3 Residue-property plots [i](#)

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

- Molecule 1: Inosine-5'-monophosphate dehydrogenase

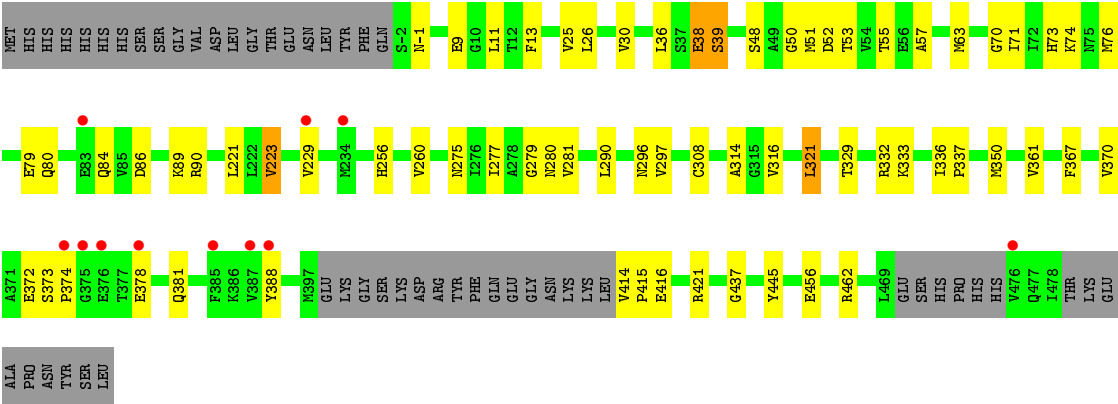


- Molecule 1: Inosine-5'-monophosphate dehydrogenase

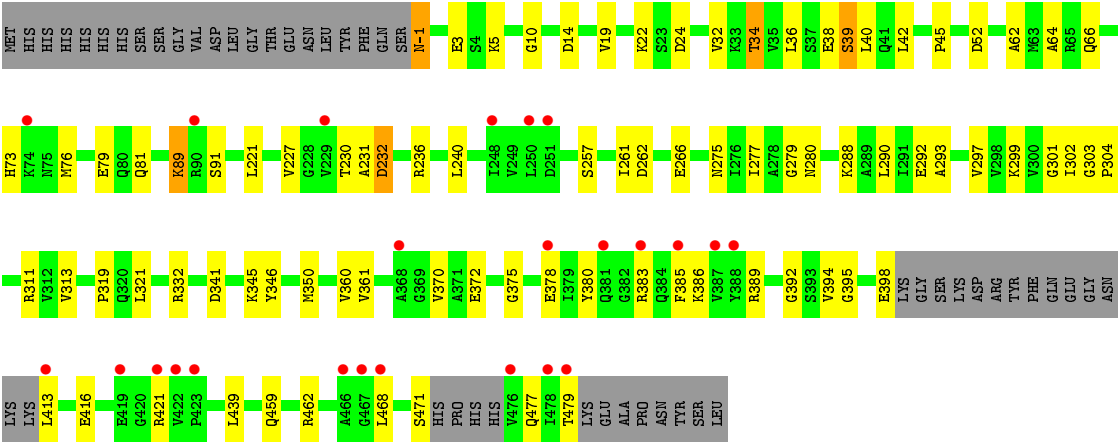


- Molecule 1: Inosine-5'-monophosphate dehydrogenase





• Molecule 1: Inosine-5'-monophosphate dehydrogenase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	83.13Å 101.33Å 87.27Å 90.00° 109.57° 90.00°	Depositor
Resolution (Å)	40.95 – 2.80 44.18 – 2.80	Depositor EDS
% Data completeness (in resolution range)	97.5 (40.95-2.80) 97.6 (44.18-2.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.15	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.35 (at 2.81Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.9_1692)	Depositor
R, R_{free}	0.192 , 0.248 0.193 , 0.249	Depositor DCC
R_{free} test set	1682 reflections (5.09%)	wwPDB-VP
Wilson B-factor (Å ²)	63.9	Xtriage
Anisotropy	0.324	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 33.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.014 for l,-k,h	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	10079	wwPDB-VP
Average B, all atoms (Å ²)	73.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.48% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: IMP, 39H

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	0.46	0/2492	0.59	0/3365
1	B	0.49	0/2557	0.61	1/3453 (0.0%)
1	C	0.45	0/2446	0.59	0/3304
1	D	0.50	0/2479	0.61	0/3349
All	All	0.47	0/9974	0.60	1/13471 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	414	VAL	C-N-CD	5.81	140.60	128.40

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	2462	0	2528	49	0
1	B	2525	0	2593	42	0
1	C	2417	0	2485	45	0
1	D	2450	0	2515	57	0
2	A	23	0	11	4	0
2	B	23	0	11	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	C	23	0	11	2	0
2	D	23	0	11	4	0
3	A	29	0	21	3	0
3	B	29	0	21	1	0
3	C	58	0	42	10	0
4	A	5	0	0	0	0
4	B	4	0	0	0	0
4	C	4	0	0	0	0
4	D	4	0	0	0	0
All	All	10079	0	10249	187	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 9.

All (187) 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:C:53:THR:HG21	1:C:388:TYR:HA	1.49	0.93
1:A:433:HIS:CE1	1:B:412:LYS:HB3	2.07	0.90
1:A:433:HIS:NE2	1:B:412:LYS:HB3	1.90	0.85
1:D:89:LYS:HE2	1:D:221:LEU:O	1.82	0.78
1:A:433:HIS:CE1	1:B:412:LYS:CB	2.71	0.74
1:C:26:LEU:HD12	3:C:501:39H:H16	1.69	0.74
1:D:350:MET:HG3	1:D:361:VAL:HG21	1.70	0.73
1:A:433:HIS:NE2	1:B:412:LYS:CB	2.54	0.69
1:B:332:ARG:NH2	1:B:456:GLU:OE1	2.25	0.69
2:A:500:IMP:C2	3:C:501:39H:H2	2.23	0.69
1:D:230:THR:HG21	1:D:232:ASP:OD2	1.93	0.69
1:D:36:LEU:HD12	1:D:40:LEU:HD23	1.76	0.68
1:A:275:ASN:HA	1:A:296:ASN:ND2	2.09	0.68
1:A:24:ASP:OD1	1:B:257:SER:HB2	1.94	0.67
1:C:39:SER:HB2	1:C:275:ASN:HD21	1.59	0.67
1:D:89:LYS:CE	1:D:221:LEU:O	2.47	0.63
1:A:257:SER:HB2	1:A:260:VAL:HG23	1.80	0.63
1:B:397:MET:SD	1:B:413:LEU:HD21	2.38	0.63
1:A:427:PRO:HG2	1:A:430:ASP:OD1	1.99	0.63
1:D:39:SER:HB2	1:D:275:ASN:HD21	1.65	0.62
1:A:430:ASP:O	1:A:434:GLN:HG2	2.01	0.61
1:A:56:GLU:HG3	1:A:374:PRO:HG3	1.83	0.60
1:D:299:LYS:HE3	1:D:341:ASP:OD2	2.00	0.60
1:A:394:VAL:O	1:A:398:GLU:HG3	2.02	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:229:VAL:HG21	1:B:260:VAL:HG22	1.84	0.59
1:C:89:LYS:HE2	1:C:223:VAL:HG23	1.84	0.59
1:D:383:ARG:HH21	1:D:385:PHE:HE2	1.50	0.59
1:A:230:THR:HG22	1:A:231:ALA:H	1.68	0.59
1:D:24:ASP:OD1	1:D:24:ASP:N	2.36	0.58
1:C:86:ASP:OD1	1:C:90:ARG:HD3	2.03	0.58
1:D:230:THR:HG22	1:D:231:ALA:H	1.68	0.58
1:C:372:GLU:N	1:C:372:GLU:OE1	2.37	0.57
2:A:500:IMP:N1	3:C:501:39H:H2	2.19	0.57
1:C:51:MET:HE1	2:C:502:IMP:H5'2	1.86	0.57
1:C:89:LYS:NZ	1:C:221:LEU:O	2.30	0.57
1:C:332:ARG:NH2	1:C:456:GLU:OE1	2.37	0.57
1:D:-1:ASN:O	1:D:3:GLU:OE1	2.23	0.57
1:A:277:ILE:HG13	1:A:297:VAL:HB	1.86	0.57
1:D:230:THR:HG22	1:D:231:ALA:N	2.20	0.56
1:D:341:ASP:OD1	2:D:500:IMP:O3'	2.20	0.56
1:B:414:VAL:HG12	1:B:414:VAL:O	2.06	0.56
1:D:313:VAL:HG21	1:D:416:GLU:HG2	1.88	0.56
1:A:383:ARG:NH1	1:A:423:PRO:HB3	2.21	0.55
1:A:308:CYS:SG	2:A:500:IMP:H2	2.46	0.55
1:C:297:VAL:HG13	1:C:337:PRO:HG2	1.89	0.55
1:D:468:LEU:HD12	1:D:471:SER:HB3	1.89	0.55
1:D:34:THR:HG22	1:D:42:LEU:HB2	1.88	0.55
1:B:313:VAL:HG21	1:B:416:GLU:HG2	1.88	0.55
1:D:230:THR:CG2	1:D:232:ASP:OD2	2.54	0.54
1:C:48:SER:HB3	1:C:70:GLY:HA2	1.88	0.54
1:A:313:VAL:HG21	1:A:416:GLU:HG2	1.90	0.54
1:A:283:THR:HG22	1:A:286:ALA:H	1.72	0.54
1:A:52:ASP:HA	1:A:73:HIS:CD2	2.43	0.54
1:B:258:GLN:HG3	1:B:262:ASP:OD1	2.07	0.54
1:A:437:GLY:HA3	1:B:414:VAL:HG21	1.89	0.53
1:B:-1:ASN:O	1:B:2:TRP:N	2.42	0.53
1:C:55:THR:HG21	1:C:71:ILE:O	2.08	0.53
1:A:332:ARG:NH2	1:A:456:GLU:OE1	2.42	0.53
1:B:56:GLU:HB2	1:B:374:PRO:HG3	1.90	0.53
1:C:367:PHE:O	1:C:370:VAL:HG22	2.10	0.52
1:A:43:ASN:HB2	1:A:67:GLY:HA3	1.90	0.52
1:B:350:MET:HE1	1:B:439:LEU:HB2	1.92	0.52
1:B:355:ALA:O	1:D:5:LYS:HE3	2.10	0.52
1:D:277:ILE:HG12	1:D:297:VAL:HB	1.91	0.52
1:D:79:GLU:H	1:D:79:GLU:CD	2.13	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:350:MET:HG3	1:C:361:VAL:HG21	1.92	0.52
1:C:26:LEU:CD1	3:C:501:39H:H16	2.39	0.52
1:C:25:VAL:HG21	1:C:30:VAL:HG12	1.92	0.51
1:A:268:ARG:NH2	1:A:296:ASN:OD1	2.43	0.51
1:A:25:VAL:HG23	1:A:29:GLU:HG3	1.93	0.51
1:A:248:ILE:HG12	1:A:274:LEU:HD21	1.92	0.51
1:C:416:GLU:OE1	3:C:503:39H:H11	2.11	0.51
1:A:252:THR:HG21	1:A:260:VAL:HG21	1.93	0.51
3:A:501:39H:H12	3:A:501:39H:O1	2.10	0.51
1:B:62:ALA:O	1:B:66:GLN:HG2	2.11	0.51
1:B:227:VAL:HG22	1:B:236:ARG:HD3	1.93	0.50
1:B:43:ASN:HB2	1:B:67:GLY:HA3	1.94	0.50
1:A:275:ASN:HA	1:A:296:ASN:HD21	1.76	0.50
3:C:503:39H:H12	3:C:503:39H:O1	2.13	0.49
1:C:57:ALA:N	1:C:84:GLN:OE1	2.44	0.49
1:D:-1:ASN:N	1:D:-1:ASN:HD22	2.10	0.49
1:D:280:ASN:OD1	1:D:299:LYS:HE2	2.11	0.49
1:C:308:CYS:SG	2:C:502:IMP:H2	2.52	0.49
1:B:308:CYS:SG	1:B:310:THR:HB	2.53	0.49
1:C:48:SER:HB2	1:C:63:MET:HG3	1.94	0.49
1:D:378:GLU:OE1	1:D:421:ARG:HD2	2.13	0.49
1:B:42:LEU:HD21	1:B:69:LEU:HB2	1.93	0.49
1:D:303:GLY:HA3	1:D:311:ARG:HE	1.78	0.48
1:D:64:ALA:HB3	1:D:221:LEU:HD13	1.95	0.48
1:C:381:GLN:O	1:C:381:GLN:HG2	2.13	0.48
1:B:88:VAL:HG11	1:B:223:VAL:HB	1.96	0.48
1:A:383:ARG:HH12	1:A:423:PRO:HB3	1.79	0.48
1:A:350:MET:HG3	1:A:361:VAL:HG21	1.94	0.48
1:A:35:VAL:HG22	1:A:41:GLN:HG2	1.96	0.48
1:B:65:ARG:NH2	1:B:91:SER:O	2.47	0.48
1:A:74:LYS:HB3	1:A:226:ALA:O	2.13	0.47
1:D:389:ARG:NH1	1:D:395:GLY:HA3	2.29	0.47
1:D:394:VAL:O	1:D:398:GLU:HB3	2.14	0.47
1:D:227:VAL:HG13	1:D:236:ARG:HD2	1.97	0.47
1:A:26:LEU:O	1:A:29:GLU:HG2	2.14	0.47
1:C:321:LEU:HA	1:C:321:LEU:HD12	1.76	0.47
1:C:229:VAL:HG21	1:C:260:VAL:HG22	1.97	0.47
1:D:372:GLU:OE1	1:D:372:GLU:N	2.48	0.47
1:D:380:TYR:O	1:D:383:ARG:HG3	2.15	0.46
1:D:10:GLY:HA3	1:D:319:PRO:HG2	1.98	0.46
1:D:52:ASP:HA	1:D:73:HIS:CD2	2.50	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:39:SER:CB	1:C:275:ASN:HD21	2.27	0.46
1:C:9:GLU:OE2	1:D:462:ARG:NH2	2.48	0.46
1:A:414:VAL:HG21	1:C:437:GLY:HA3	1.98	0.46
3:B:501:39H:H2	2:D:500:IMP:C2	2.45	0.46
1:A:289:ALA:O	1:A:292:GLU:HG2	2.16	0.46
1:A:433:HIS:CE1	1:B:412:LYS:HB2	2.51	0.46
1:C:256:HIS:CD2	1:D:22:LYS:HD3	2.51	0.45
1:A:227:VAL:HG22	1:A:236:ARG:HD2	1.97	0.45
1:A:257:SER:OG	3:C:501:39H:H17	2.15	0.45
1:D:261:ILE:HG23	1:D:293:ALA:HB2	1.98	0.45
1:C:-1:ASN:OD1	1:D:332:ARG:NH1	2.50	0.45
1:C:277:ILE:HG12	1:C:297:VAL:HB	1.98	0.45
1:C:11:LEU:HD21	1:C:462:ARG:HD3	1.98	0.45
1:D:303:GLY:N	1:D:304:PRO:HD3	2.32	0.45
1:D:341:ASP:OD2	2:D:500:IMP:O2'	2.29	0.45
1:D:375:GLY:O	1:D:386:LYS:NZ	2.42	0.45
1:C:52:ASP:HA	1:C:73:HIS:CD2	2.52	0.45
1:C:314:ALA:O	1:C:316:VAL:HG23	2.16	0.44
1:C:281:VAL:HG11	1:C:290:LEU:HD22	1.99	0.44
1:C:36:LEU:N	1:C:36:LEU:HD23	2.32	0.44
3:A:501:39H:H2	2:B:500:IMP:C2	2.47	0.44
1:A:283:THR:CG2	1:A:285:GLU:HB2	2.47	0.44
1:A:345:LYS:HB2	1:A:345:LYS:HE2	1.77	0.44
1:B:16:VAL:HG21	1:B:321:LEU:HD21	1.99	0.44
1:B:350:MET:HG3	1:B:361:VAL:HG11	1.99	0.44
1:C:296:ASN:O	1:C:336:ILE:HG23	2.18	0.44
1:D:262:ASP:O	1:D:266:GLU:HG3	2.18	0.44
1:C:421:ARG:NH1	1:D:479:THR:OG1	2.50	0.44
1:B:328:ALA:O	1:B:332:ARG:HG3	2.18	0.44
1:A:341:ASP:OD1	2:A:500:IMP:O3'	2.36	0.43
1:B:248:ILE:HG12	1:B:274:LEU:HD21	2.00	0.43
1:B:23:SER:O	1:D:257:SER:HA	2.18	0.43
1:D:261:ILE:HD13	1:D:290:LEU:HD23	2.01	0.43
1:D:350:MET:HE1	1:D:439:LEU:HB2	2.00	0.43
1:A:59:MET:CE	1:A:367:PHE:HB3	2.49	0.43
1:B:74:LYS:NZ	1:B:228:GLY:H	2.16	0.43
1:A:39:SER:HB2	1:A:275:ASN:OD1	2.19	0.43
1:D:302:ILE:C	1:D:304:PRO:HD3	2.39	0.43
3:A:501:39H:H8	3:A:501:39H:H7	1.83	0.43
1:A:339:ILE:HG12	1:A:360:VAL:CG1	2.49	0.42
1:D:280:ASN:ND2	1:D:301:GLY:O	2.40	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:81:GLN:HG2	1:D:240:LEU:HD21	2.00	0.42
1:D:19:VAL:HG22	1:D:459:GLN:HB2	2.01	0.42
1:C:329:THR:O	1:C:333:LYS:HE3	2.19	0.42
1:D:345:LYS:HE2	1:D:346:TYR:CE2	2.55	0.42
1:D:62:ALA:HB2	1:D:372:GLU:HG2	2.01	0.42
1:A:43:ASN:HB2	1:A:67:GLY:CA	2.50	0.42
1:D:321:LEU:HD12	1:D:321:LEU:HA	1.76	0.42
1:B:52:ASP:HA	1:B:73:HIS:CD2	2.54	0.42
1:D:288:LYS:HE3	1:D:292:GLU:OE2	2.20	0.42
1:D:62:ALA:O	1:D:66:GLN:HG2	2.20	0.42
1:B:236:ARG:O	1:B:240:LEU:HG	2.20	0.42
1:B:397:MET:SD	1:B:413:LEU:CD2	3.07	0.41
1:B:84:GLN:O	1:B:88:VAL:HG23	2.20	0.41
1:C:373:SER:HA	1:C:374:PRO:HD2	1.82	0.41
1:C:25:VAL:O	3:C:501:39H:H20	2.20	0.41
1:D:14:ASP:HB3	1:D:468:LEU:HD23	2.00	0.41
1:B:350:MET:CE	1:B:439:LEU:HB2	2.49	0.41
1:D:32:VAL:O	1:D:45:PRO:HD3	2.19	0.41
1:B:263:LYS:O	1:B:267:VAL:HG23	2.20	0.41
1:A:32:VAL:O	1:A:45:PRO:HD3	2.21	0.41
1:C:13:PHE:HA	1:C:321:LEU:HD22	2.02	0.41
3:C:501:39H:H12	3:C:501:39H:O1	2.21	0.41
1:A:260:VAL:O	1:A:264:VAL:HG23	2.20	0.41
1:B:234:MET:HE1	1:B:270:LYS:HD3	2.02	0.41
1:D:279:GLY:HA3	1:D:280:ASN:HA	1.92	0.41
1:A:454:LEU:HD12	1:A:458:ALA:HB2	2.02	0.41
1:B:76:MET:HB2	1:B:80:GLN:HG3	2.02	0.41
1:C:414:VAL:HA	1:C:415:PRO:HD3	1.92	0.41
1:A:254:HIS:CD2	1:C:445:TYR:HA	2.56	0.40
1:B:286:ALA:O	1:B:289:ALA:HB3	2.22	0.40
1:B:351:VAL:HG22	1:B:439:LEU:HA	2.03	0.40
1:B:383:ARG:HH12	1:B:483:PRO:HG3	1.86	0.40
1:C:50:GLY:HA2	1:C:71:ILE:O	2.21	0.40
1:A:53:THR:HG21	1:A:389:ARG:HG2	2.03	0.40
1:C:38:GLU:HG3	1:C:38:GLU:H	1.46	0.40
1:B:430:ASP:O	1:B:434:GLN:HG2	2.22	0.40
3:C:501:39H:H8	3:C:501:39H:H7	1.79	0.40
1:A:57:ALA:O	1:A:61:ILE:HG13	2.21	0.40
1:C:279:GLY:HA3	1:C:280:ASN:HA	1.90	0.40
1:D:392:GLY:N	2:D:500:IMP:O6	2.51	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	331/384 (86%)	322 (97%)	9 (3%)	0	100	100
1	B	339/384 (88%)	331 (98%)	8 (2%)	0	100	100
1	C	326/384 (85%)	314 (96%)	12 (4%)	0	100	100
1	D	330/384 (86%)	322 (98%)	8 (2%)	0	100	100
All	All	1326/1536 (86%)	1289 (97%)	37 (3%)	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	256/298 (86%)	245 (96%)	11 (4%)	29	62
1	B	263/298 (88%)	257 (98%)	6 (2%)	50	82
1	C	251/298 (84%)	242 (96%)	9 (4%)	35	69
1	D	255/298 (86%)	243 (95%)	12 (5%)	26	59
All	All	1025/1192 (86%)	987 (96%)	38 (4%)	34	68

All (38) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	25	VAL
1	A	39	SER

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Mol	Chain	Res	Type
1	A	65	ARG
1	A	83	GLU
1	A	90	ARG
1	A	232	ASP
1	A	296	ASN
1	A	370	VAL
1	A	376	GLU
1	A	386	LYS
1	A	430	ASP
1	B	38	GLU
1	B	83	GLU
1	B	89	LYS
1	B	310	THR
1	B	311	ARG
1	B	384	GLN
1	C	38	GLU
1	C	39	SER
1	C	74	LYS
1	C	76	MET
1	C	79	GLU
1	C	80	GLN
1	C	223	VAL
1	C	321	LEU
1	C	378	GLU
1	D	-1	ASN
1	D	34	THR
1	D	38	GLU
1	D	39	SER
1	D	76	MET
1	D	89	LYS
1	D	91	SER
1	D	232	ASP
1	D	360	VAL
1	D	370	VAL
1	D	413	LEU
1	D	477	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	256	HIS
1	B	258	GLN

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Mol	Chain	Res	Type
1	D	-1	ASN

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

8 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	IMP	C	502	-	21,25,25	1.40	3 (14%)	23,38,38	1.35	2 (8%)
2	IMP	A	500	-	21,25,25	1.36	3 (14%)	23,38,38	1.43	3 (13%)
3	39H	A	501	-	25,33,33	1.73	6 (24%)	25,48,48	1.58	4 (16%)
3	39H	C	501	-	25,33,33	1.63	5 (20%)	25,48,48	1.87	6 (24%)
3	39H	B	501	-	25,33,33	1.71	6 (24%)	25,48,48	1.75	4 (16%)
3	39H	C	503	-	25,33,33	1.70	5 (20%)	25,48,48	1.20	2 (8%)
2	IMP	B	500	-	21,25,25	1.28	3 (14%)	23,38,38	1.58	3 (13%)
2	IMP	D	500	-	21,25,25	1.34	3 (14%)	23,38,38	1.45	2 (8%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the

Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns.
'-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	IMP	C	502	-	-	0/6/26/26	0/3/3/3
2	IMP	A	500	-	-	2/6/26/26	0/3/3/3
3	39H	A	501	-	-	0/8/15/15	0/5/5/5
3	39H	C	501	-	-	0/8/15/15	0/5/5/5
3	39H	B	501	-	-	0/8/15/15	0/5/5/5
3	39H	C	503	-	-	0/8/15/15	0/5/5/5
2	IMP	B	500	-	-	0/6/26/26	0/3/3/3
2	IMP	D	500	-	-	0/6/26/26	0/3/3/3

All (34) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	501	39H	C6-C7	-5.03	1.37	1.43
3	C	501	39H	C6-C7	-4.69	1.38	1.43
3	A	501	39H	C6-C7	-4.66	1.38	1.43
3	C	503	39H	C6-C7	-4.38	1.38	1.43
2	A	500	IMP	C2-N3	4.12	1.38	1.32
2	C	502	IMP	C2-N3	4.04	1.38	1.32
2	B	500	IMP	C2-N3	3.87	1.38	1.32
2	D	500	IMP	C2-N3	3.86	1.38	1.32
2	C	502	IMP	C6-N1	3.59	1.39	1.33
3	C	503	39H	C21-C20	-3.58	1.35	1.43
3	B	501	39H	C21-C20	-3.35	1.36	1.43
2	D	500	IMP	C6-N1	3.33	1.38	1.33
3	C	503	39H	C23-N1	-3.22	1.35	1.41
3	C	501	39H	C21-C20	-3.08	1.36	1.43
3	A	501	39H	C21-C20	-3.08	1.36	1.43
3	A	501	39H	C23-N1	-3.04	1.35	1.41
2	A	500	IMP	C6-N1	3.01	1.38	1.33
2	B	500	IMP	C6-N1	2.91	1.38	1.33
2	B	500	IMP	C2-N1	2.75	1.39	1.33
2	C	502	IMP	C2-N1	2.71	1.39	1.33
3	C	501	39H	C23-N1	-2.70	1.36	1.41
3	A	501	39H	C7-C8	-2.68	1.37	1.42
2	A	500	IMP	C2-N1	2.54	1.38	1.33
3	C	503	39H	C7-C8	-2.51	1.38	1.42
3	B	501	39H	C23-N1	-2.49	1.36	1.41
2	D	500	IMP	C2-N1	2.45	1.38	1.33
3	A	501	39H	C9-C8	-2.40	1.37	1.41
3	B	501	39H	C7-C8	-2.35	1.38	1.42
3	B	501	39H	C9-C8	-2.29	1.37	1.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	501	39H	C12-C7	-2.27	1.37	1.42
3	B	501	39H	C27-C25	-2.25	1.49	1.50
3	C	501	39H	C7-C8	-2.24	1.38	1.42
3	C	503	39H	C9-C8	-2.07	1.38	1.41
3	C	501	39H	C9-C8	-2.06	1.38	1.41

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	500	IMP	N3-C2-N1	-5.49	120.10	128.68
2	D	500	IMP	N3-C2-N1	-5.25	120.47	128.68
2	C	502	IMP	N3-C2-N1	-5.03	120.82	128.68
2	A	500	IMP	N3-C2-N1	-4.99	120.89	128.68
3	C	501	39H	C12-C7-C6	-4.48	119.35	123.35
3	C	501	39H	C29-C30-C26	-4.38	105.85	112.49
3	B	501	39H	C28-C27-C25	-4.32	106.43	113.53
3	B	501	39H	C12-C7-C6	-4.31	119.51	123.35
3	A	501	39H	C12-C7-C6	-4.15	119.65	123.35
3	C	501	39H	C6-C7-C8	3.89	120.31	118.19
3	A	501	39H	C6-C7-C8	3.82	120.27	118.19
3	B	501	39H	C6-C7-C8	3.80	120.26	118.19
3	C	501	39H	C28-C27-C25	-3.34	108.03	113.53
3	C	503	39H	C12-C7-C6	-2.98	120.70	123.35
2	B	500	IMP	C2-N1-C6	2.90	120.74	115.88
3	C	503	39H	C23-C22-C21	-2.78	117.05	120.78
3	A	501	39H	C23-C22-C21	-2.56	117.34	120.78
2	D	500	IMP	C2-N1-C6	2.55	120.15	115.88
3	B	501	39H	C23-C22-C21	-2.48	117.44	120.78
3	A	501	39H	C28-C27-C25	-2.46	109.48	113.53
2	C	502	IMP	C2-N1-C6	2.44	119.97	115.88
2	A	500	IMP	C3'-C2'-C1'	2.36	104.53	100.98
2	A	500	IMP	O2P-P-O1P	2.26	119.53	110.68
2	B	500	IMP	C4-C5-N7	-2.14	107.17	109.40
3	C	501	39H	C23-C22-C21	-2.10	117.95	120.78
3	C	501	39H	C5-C6-N11	2.02	119.86	116.12

There are no chirality outliers.

All (2) torsion outliers are listed below:

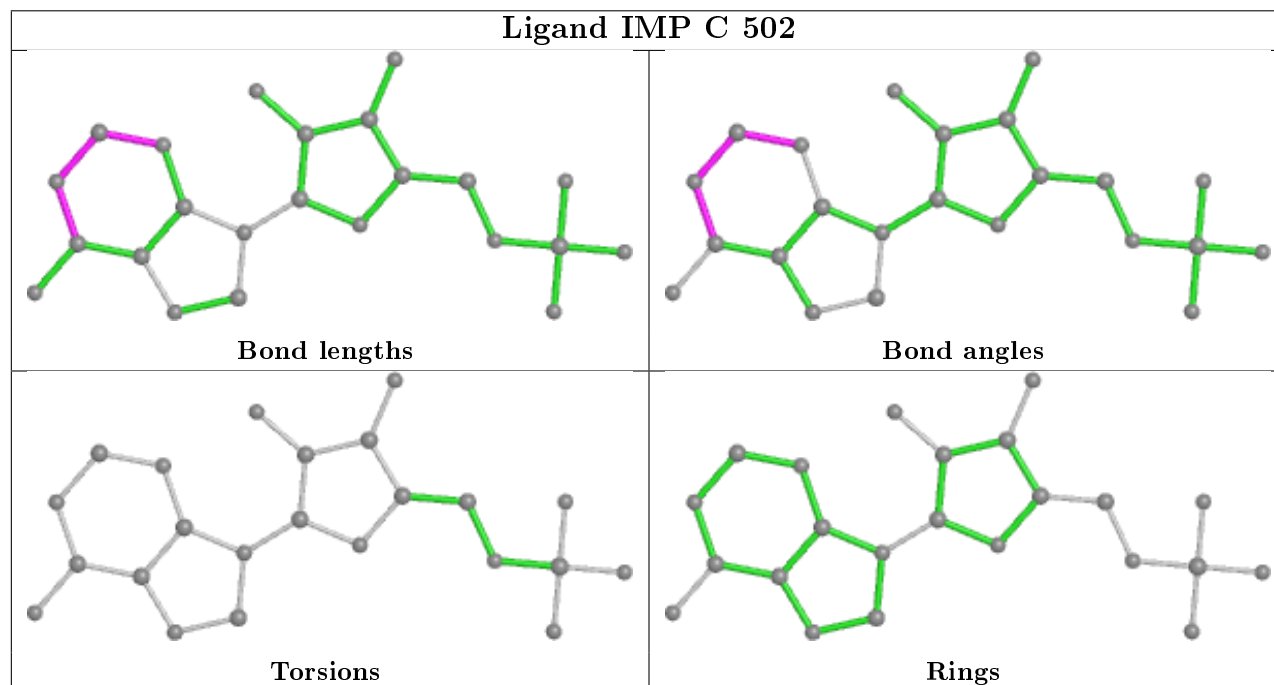
Mol	Chain	Res	Type	Atoms
2	A	500	IMP	C3'-C4'-C5'-O5'
2	A	500	IMP	O4'-C4'-C5'-O5'

There are no ring outliers.

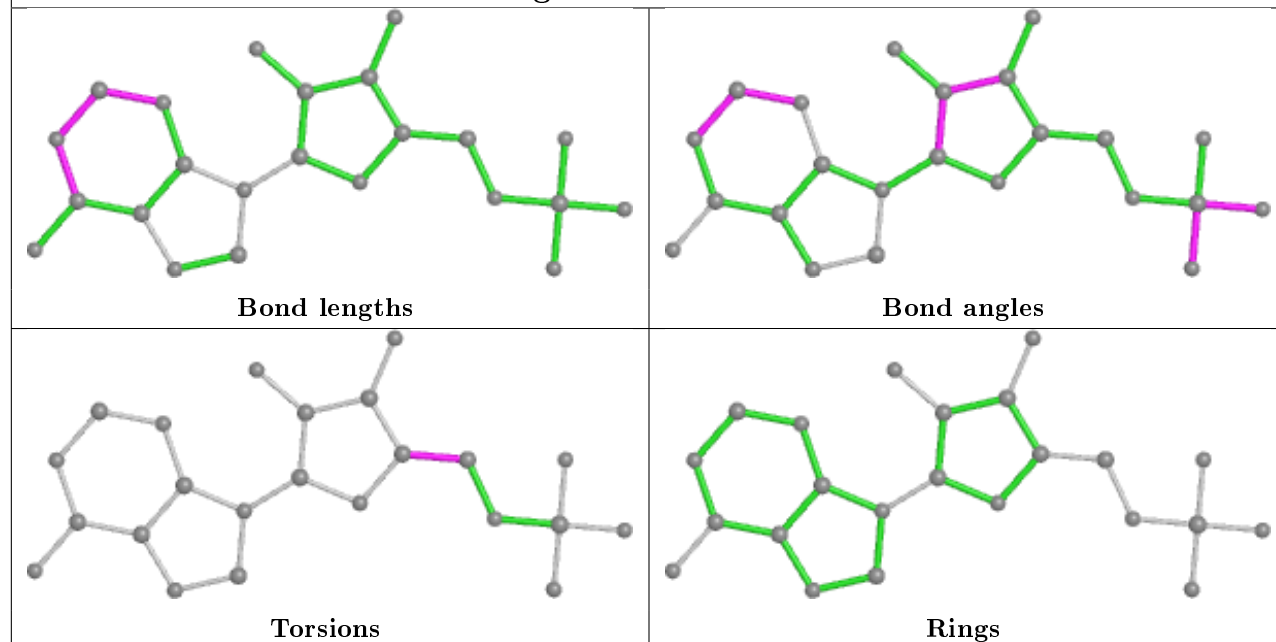
8 monomers are involved in 21 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	502	IMP	2	0
2	A	500	IMP	4	0
3	A	501	39H	3	0
3	C	501	39H	8	0
3	B	501	39H	1	0
3	C	503	39H	2	0
2	B	500	IMP	1	0
2	D	500	IMP	4	0

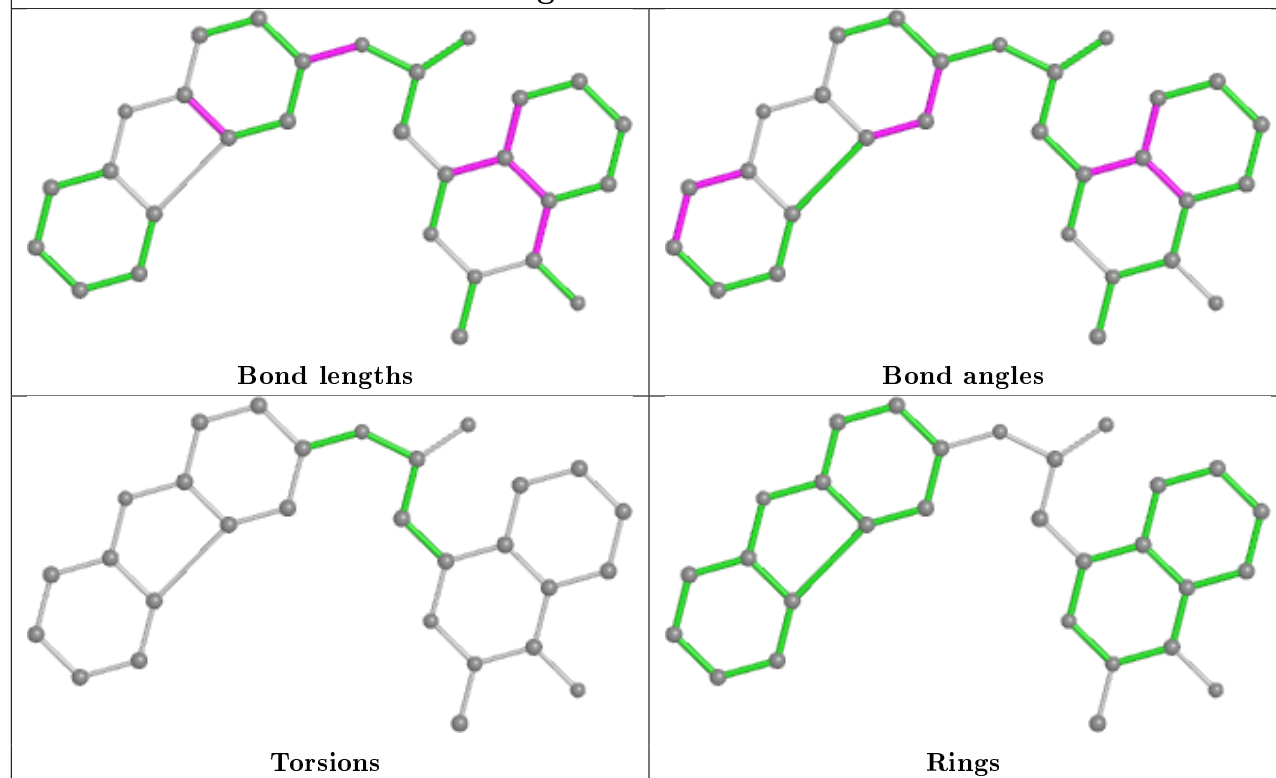
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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



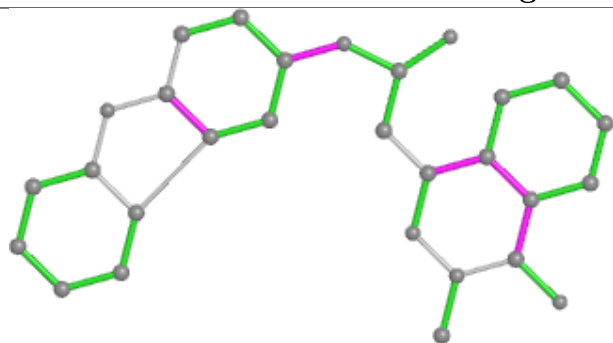
Ligand IMP A 500



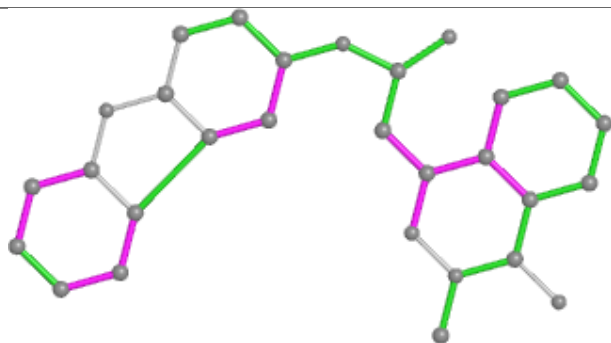
Ligand 39H A 501



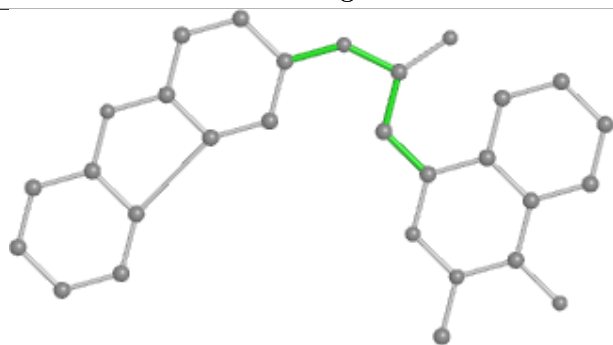
Ligand 39H C 501



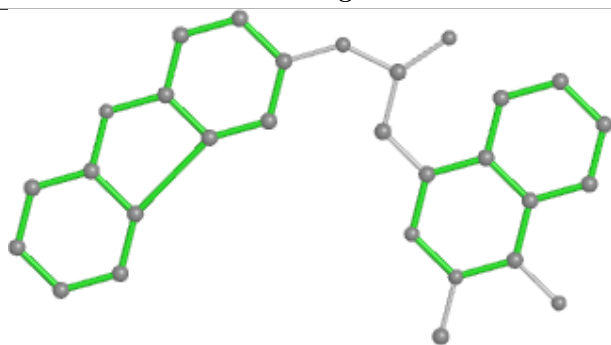
Bond lengths



Bond angles

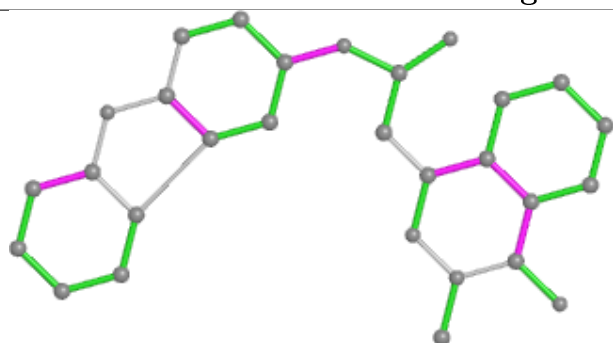


Torsions

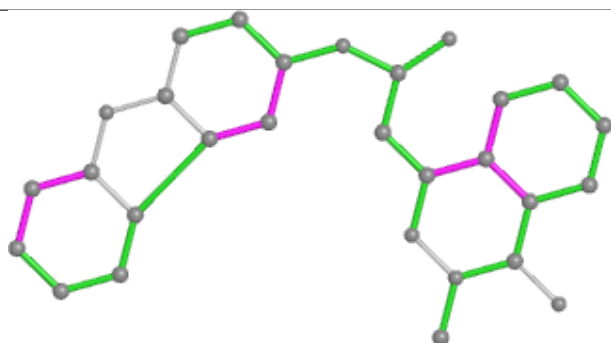


Rings

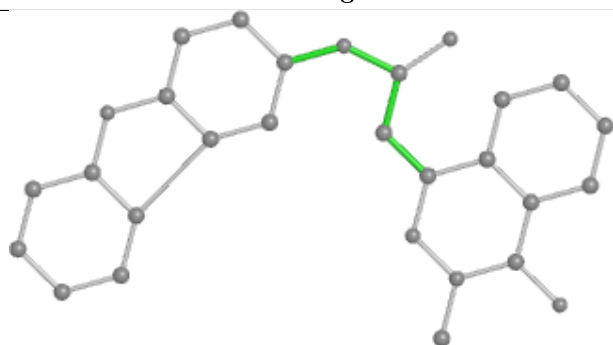
Ligand 39H B 501



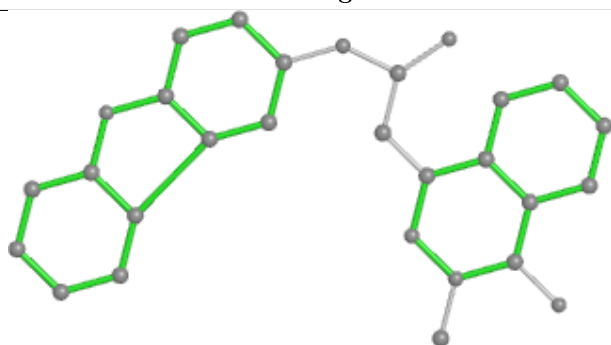
Bond lengths



Bond angles

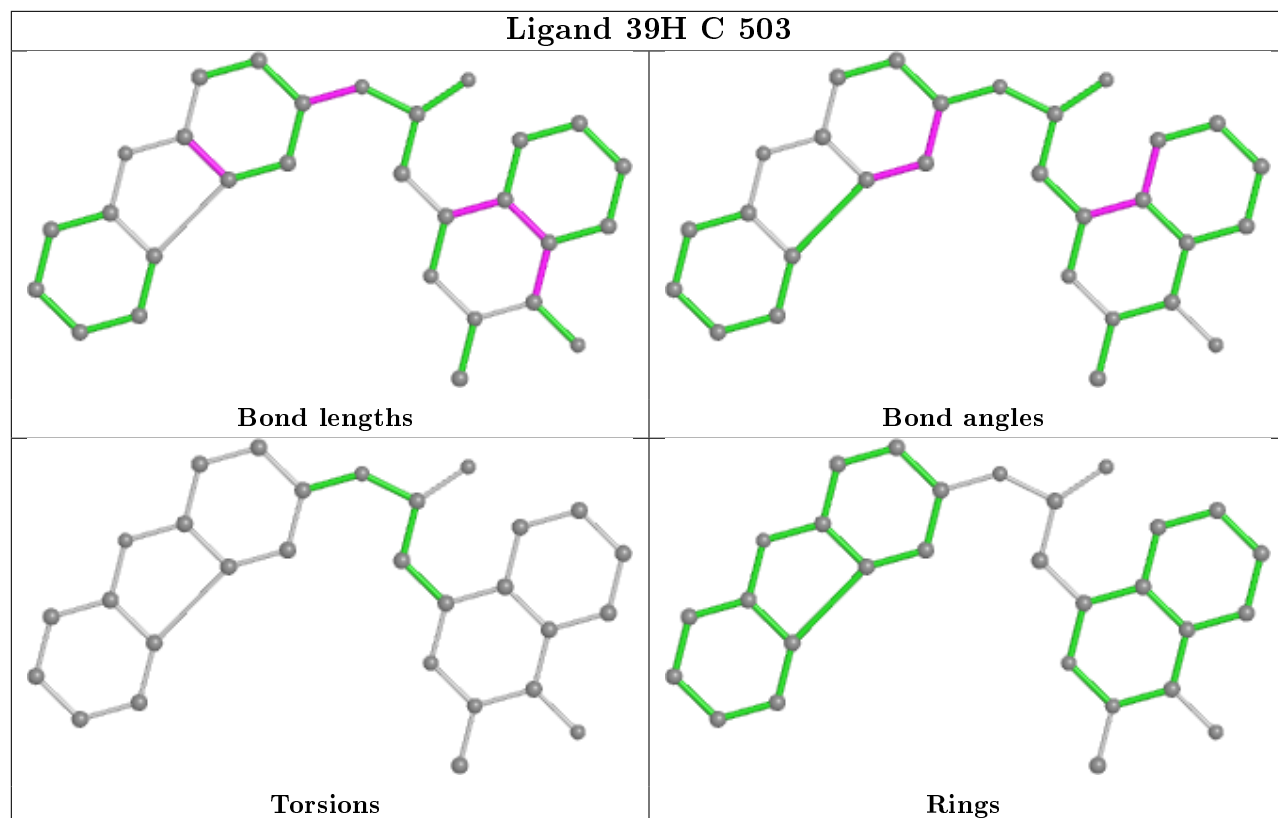


Torsions

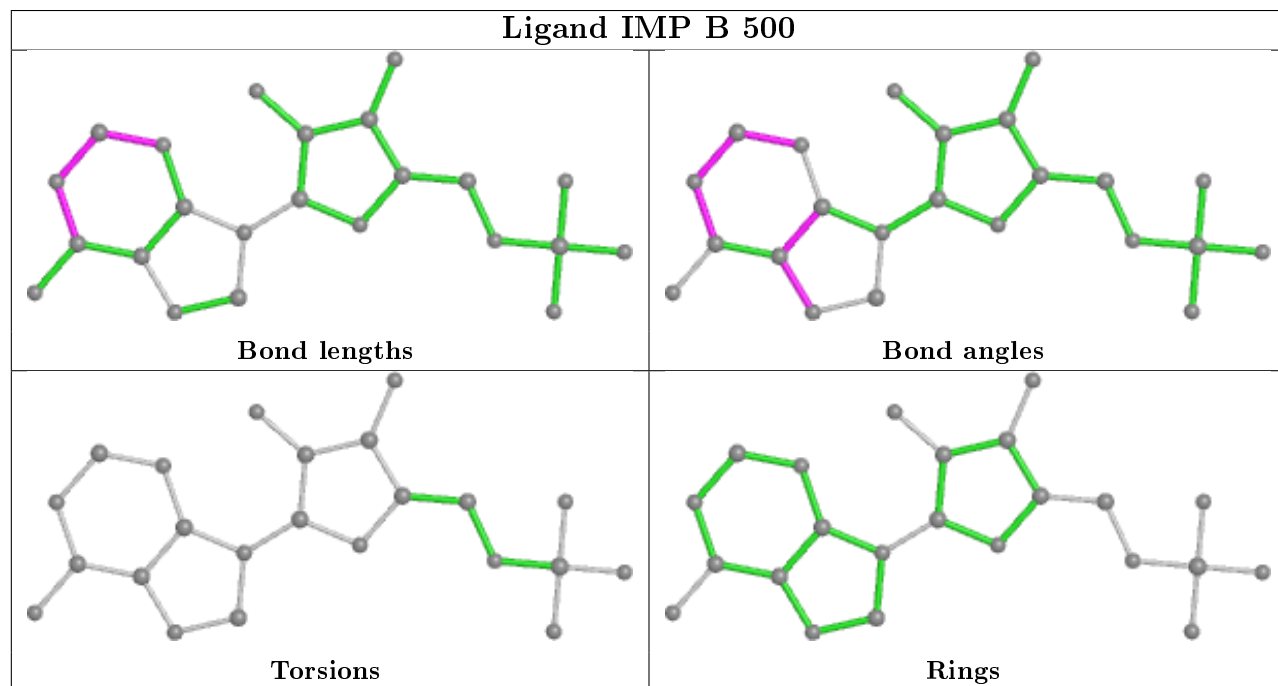


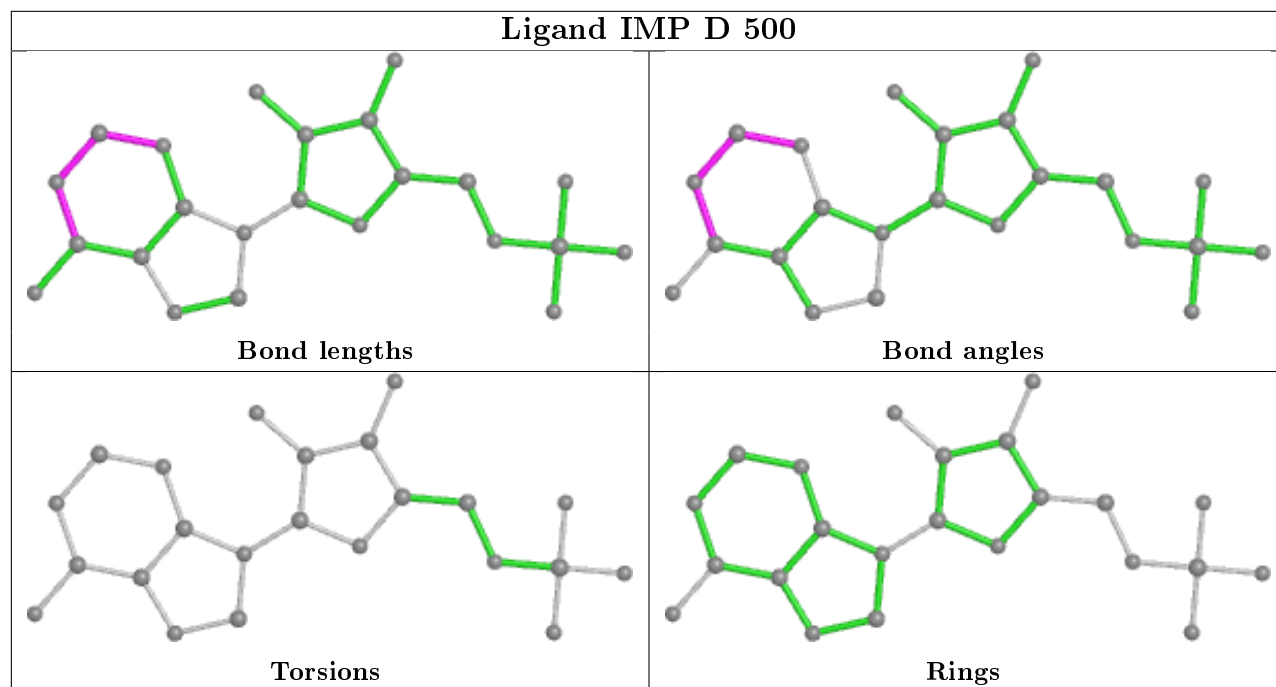
Rings

Ligand 39H C 503



Ligand IMP B 500





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2		OWAB(Å ²)	Q<0.9
1	A	337/384 (87%)	0.27	15 (4%)	33 23	48, 69, 113, 142	1 (0%)
1	B	345/384 (89%)	0.39	21 (6%)	21 13	46, 73, 120, 149	0
1	C	332/384 (86%)	0.38	11 (3%)	46 36	44, 70, 117, 144	2 (0%)
1	D	336/384 (87%)	0.33	24 (7%)	16 9	42, 63, 109, 131	1 (0%)
All	All	1350/1536 (87%)	0.34	71 (5%)	26 17	42, 69, 116, 149	4 (0%)

All (71) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	467	GLY	9.4
1	C	375	GLY	7.7
1	D	468	LEU	6.5
1	D	385	PHE	5.0
1	B	222	LEU	4.6
1	B	385	PHE	4.6
1	C	388	TYR	4.4
1	D	476	VAL	4.3
1	D	250	LEU	4.1
1	D	421	ARG	4.1
1	B	220	GLY	4.0
1	B	475	HIS	3.9
1	B	416	GLU	3.9
1	C	476	VAL	3.8
1	D	368	ALA	3.8
1	D	251	ASP	3.7
1	A	240	LEU	3.6
1	A	277	ILE	3.3
1	C	387	VAL	3.3
1	D	422	VAL	3.3
1	A	466	ALA	3.2

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Mol	Chain	Res	Type	RSRZ
1	A	247	ALA	3.2
1	C	378	GLU	3.1
1	D	413	LEU	3.1
1	C	229	VAL	3.1
1	C	376	GLU	3.1
1	B	418	ILE	3.1
1	D	378	GLU	3.1
1	D	387	VAL	3.0
1	B	376	GLU	3.0
1	A	248	ILE	2.9
1	A	250	LEU	2.9
1	B	378	GLU	2.9
1	A	434	GLN	2.8
1	A	419	GLU	2.8
1	B	476	VAL	2.8
1	D	479	THR	2.8
1	D	229	VAL	2.8
1	D	423	PRO	2.7
1	A	232	ASP	2.7
1	B	58	ASP	2.7
1	A	469	LEU	2.7
1	B	479	THR	2.7
1	B	301	GLY	2.6
1	B	38	GLU	2.6
1	D	478	ILE	2.5
1	D	388	TYR	2.5
1	D	419	GLU	2.5
1	A	298	VAL	2.5
1	C	385	PHE	2.4
1	D	74	LYS	2.4
1	D	383	ARG	2.4
1	D	90	ARG	2.4
1	B	221	LEU	2.4
1	D	248	ILE	2.4
1	D	381	GLN	2.3
1	C	374	PRO	2.3
1	A	385	PHE	2.3
1	B	226	ALA	2.3
1	A	356	ALA	2.2
1	D	466	ALA	2.2
1	B	397	MET	2.2
1	B	346	TYR	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	476	VAL	2.1
1	B	419	GLU	2.1
1	B	297	VAL	2.1
1	B	387	VAL	2.1
1	B	-1	ASN	2.1
1	C	83	GLU	2.1
1	C	234	MET	2.1
1	A	376	GLU	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 carbohydrates in this entry.

6.4 Ligands [i](#)

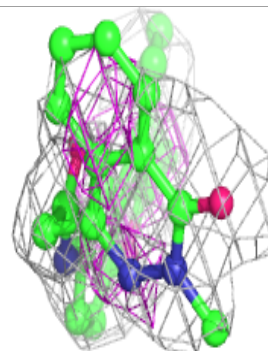
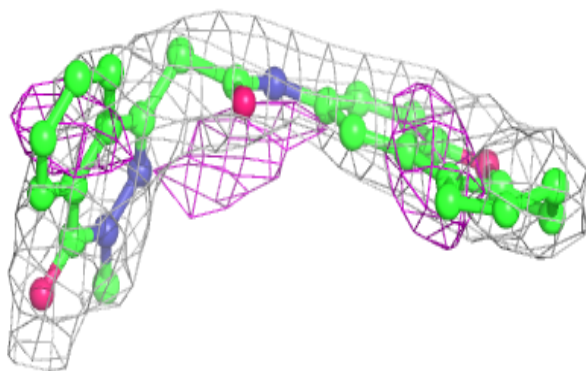
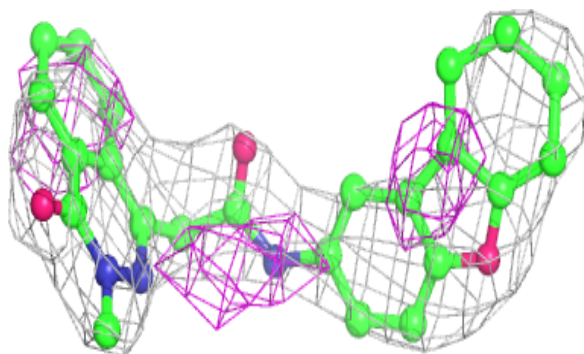
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	39H	B	501	29/29	0.84	0.39	71,74,75,77	0
3	39H	C	503	29/29	0.89	0.26	76,81,88,90	0
3	39H	A	501	29/29	0.90	0.45	68,72,74,76	0
3	39H	C	501	29/29	0.93	0.19	69,70,74,74	0
2	IMP	B	500	23/23	0.94	0.17	60,66,67,68	0
2	IMP	C	502	23/23	0.95	0.15	55,66,71,71	0
2	IMP	A	500	23/23	0.96	0.15	60,62,63,63	0
2	IMP	D	500	23/23	0.98	0.12	48,60,62,63	0

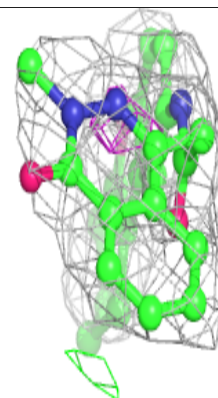
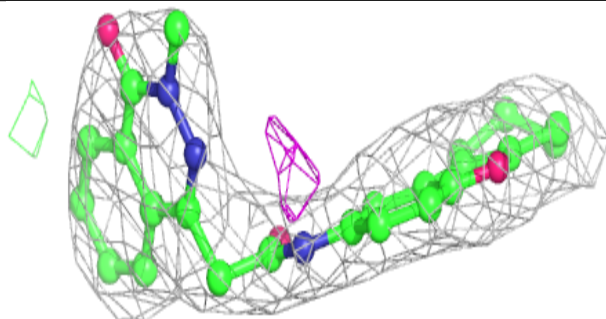
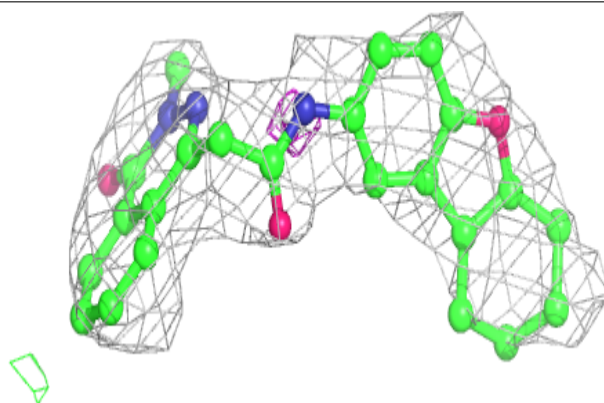
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around 39H 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)

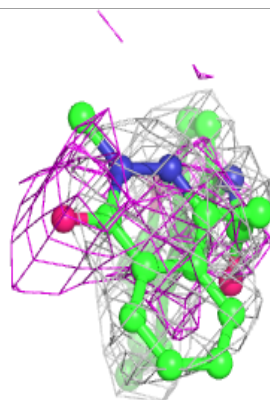
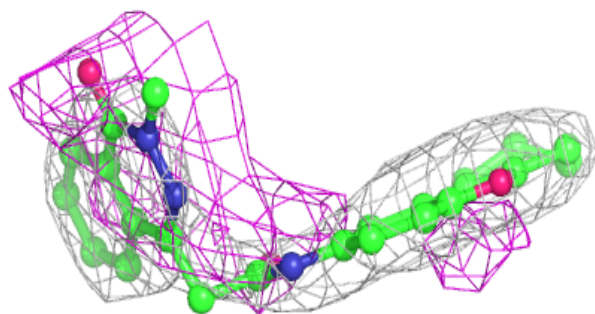
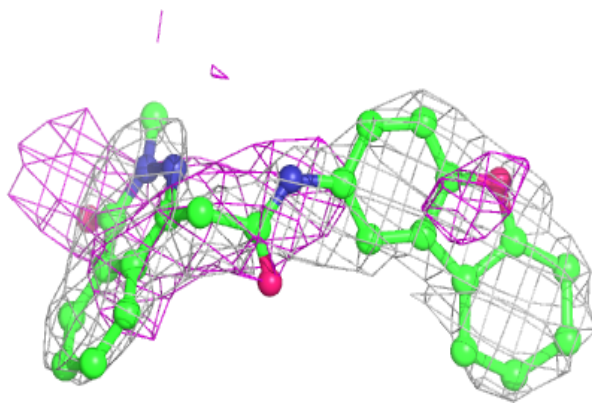
**Electron density around 39H C 503:**

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

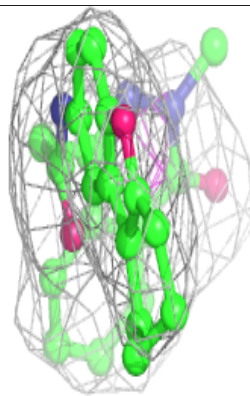
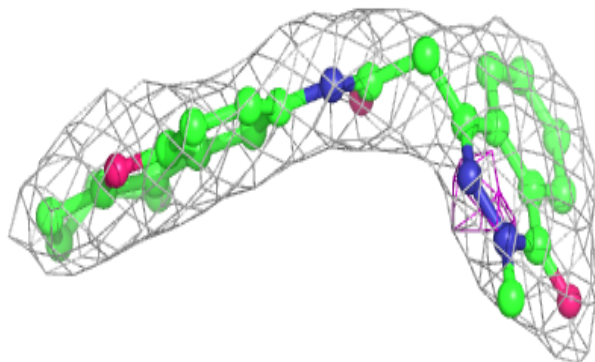
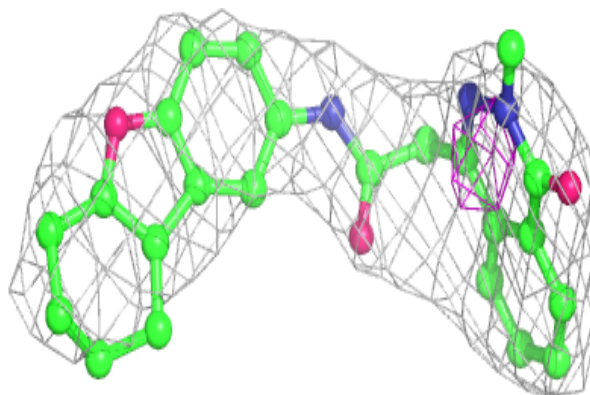


Electron density around 39H A 501:

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

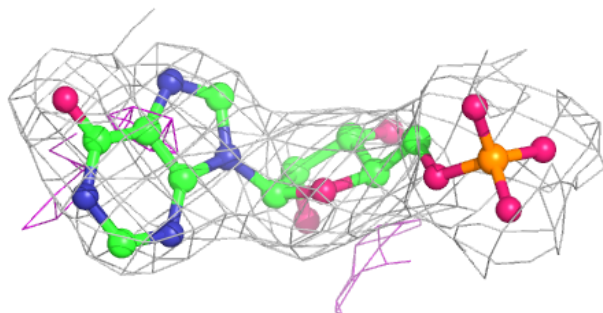
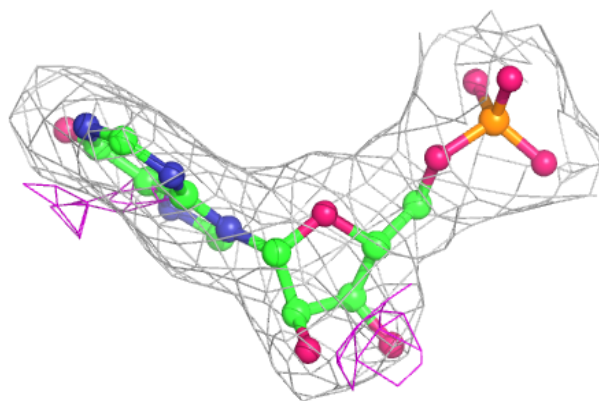
**Electron density around 39H C 501:**

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

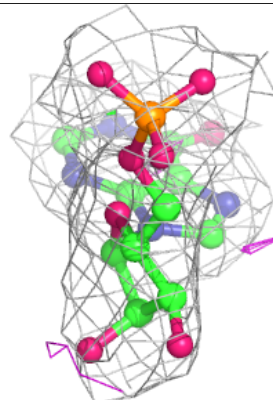
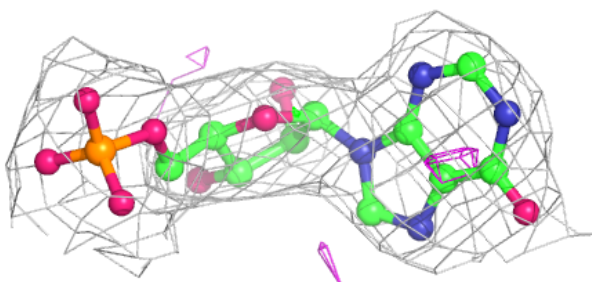
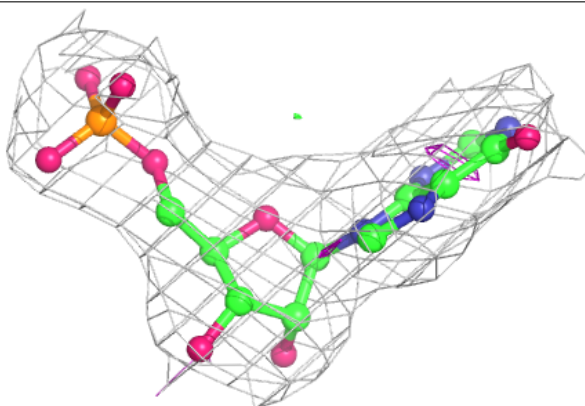


Electron density around IMP B 500:

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

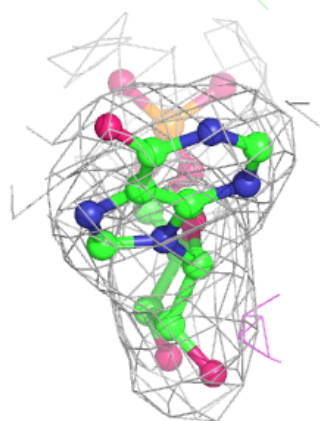
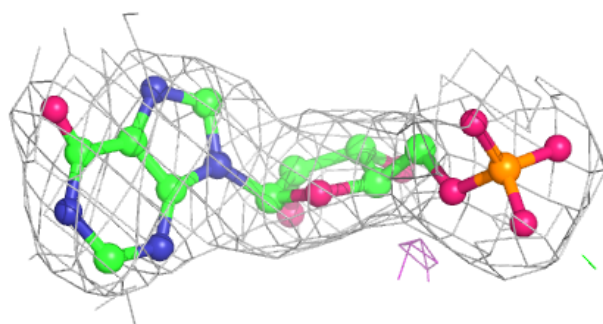
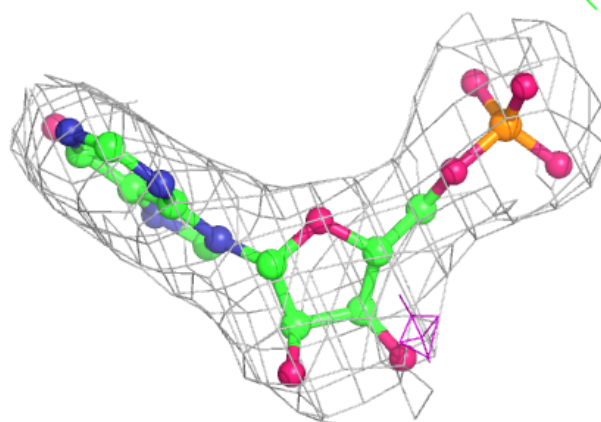
**Electron density around IMP C 502:**

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

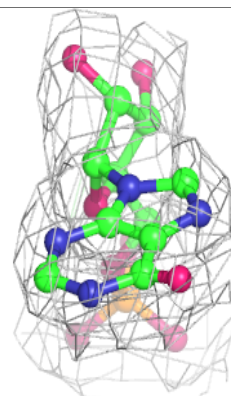
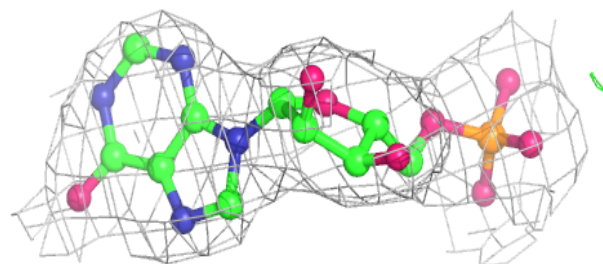
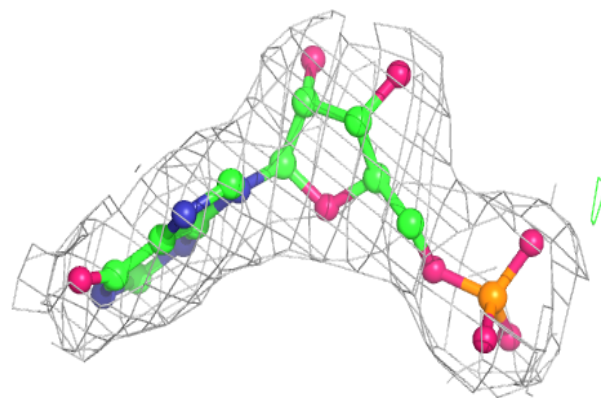


Electron density around IMP A 500:

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

**Electron density around IMP 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)



6.5 Other polymers

There are no such residues in this entry.