



Full wwPDB X-ray Structure Validation Report ⓘ

May 24, 2020 – 09:13 am BST

PDB ID : 4MZ8
Title : Crystal Structure of the Inosine 5'-monophosphate Dehydrogenase, with an Internal Deletion of CBS Domain from *Campylobacter jejuni* complexed with inhibitor compound 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 : 2013-09-29
Resolution : 2.50 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

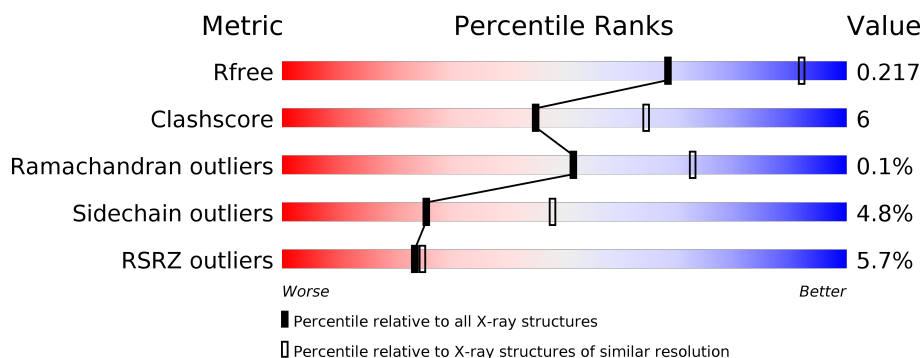
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	4661 (2.50-2.50)
Clashscore	141614	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)
RSRZ outliers	127900	4559 (2.50-2.50)

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	406	<div> <div>%</div> <div> <div></div> <div>78%</div> <div>9%</div> <div>13%</div> </div> </div>
1	B	406	<div> <div>%</div> <div> <div></div> <div>79%</div> <div>8%</div> <div>13%</div> </div> </div>
1	C	406	<div> <div>6%</div> <div> <div></div> <div>68%</div> <div>16%</div> <div>•</div> <div>14%</div> </div> </div>
1	D	406	<div> <div>12%</div> <div> <div></div> <div>63%</div> <div>20%</div> <div>•</div> <div>14%</div> </div> </div>

2 Entry composition

There are 11 unique types of molecules in this entry. The entry contains 11176 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	353	Total	C	N	O	S	0	5	0
			2669	1680	473	502	14			
1	B	354	Total	C	N	O	S	0	7	0
			2691	1691	476	509	15			
1	C	349	Total	C	N	O	S	0	3	0
			2619	1643	463	500	13			
1	D	350	Total	C	N	O	S	0	1	0
			2607	1642	460	491	14			

There are 100 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-23	MET	-	EXPRESSION TAG	UNP Q0P9J4
A	-22	HIS	-	EXPRESSION TAG	UNP Q0P9J4
A	-21	HIS	-	EXPRESSION TAG	UNP Q0P9J4
A	-20	HIS	-	EXPRESSION TAG	UNP Q0P9J4
A	-19	HIS	-	EXPRESSION TAG	UNP Q0P9J4
A	-18	HIS	-	EXPRESSION TAG	UNP Q0P9J4
A	-17	HIS	-	EXPRESSION TAG	UNP Q0P9J4
A	-16	SER	-	EXPRESSION TAG	UNP Q0P9J4
A	-15	SER	-	EXPRESSION TAG	UNP Q0P9J4
A	-14	GLY	-	EXPRESSION TAG	UNP Q0P9J4
A	-13	VAL	-	EXPRESSION TAG	UNP Q0P9J4
A	-12	ASP	-	EXPRESSION TAG	UNP Q0P9J4
A	-11	LEU	-	EXPRESSION TAG	UNP Q0P9J4
A	-10	GLY	-	EXPRESSION TAG	UNP Q0P9J4
A	-9	THR	-	EXPRESSION TAG	UNP Q0P9J4
A	-8	GLU	-	EXPRESSION TAG	UNP Q0P9J4
A	-7	ASN	-	EXPRESSION TAG	UNP Q0P9J4
A	-6	LEU	-	EXPRESSION TAG	UNP Q0P9J4
A	-5	TYR	-	EXPRESSION TAG	UNP Q0P9J4
A	-4	PHE	-	EXPRESSION TAG	UNP Q0P9J4
A	-3	GLN	-	EXPRESSION TAG	UNP Q0P9J4

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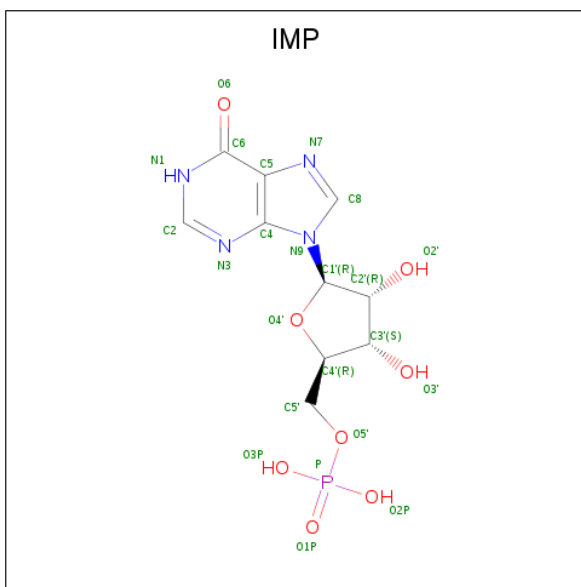
Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	SER	-	EXPRESSION TAG	UNP Q0P9J4
A	-1	ASN	-	EXPRESSION TAG	UNP Q0P9J4
A	0	ALA	-	EXPRESSION TAG	UNP Q0P9J4
A	92	GLY	-	LINKER	UNP Q0P9J4
B	-23	MET	-	EXPRESSION TAG	UNP Q0P9J4
B	-22	HIS	-	EXPRESSION TAG	UNP Q0P9J4
B	-21	HIS	-	EXPRESSION TAG	UNP Q0P9J4
B	-20	HIS	-	EXPRESSION TAG	UNP Q0P9J4
B	-19	HIS	-	EXPRESSION TAG	UNP Q0P9J4
B	-18	HIS	-	EXPRESSION TAG	UNP Q0P9J4
B	-17	HIS	-	EXPRESSION TAG	UNP Q0P9J4
B	-16	SER	-	EXPRESSION TAG	UNP Q0P9J4
B	-15	SER	-	EXPRESSION TAG	UNP Q0P9J4
B	-14	GLY	-	EXPRESSION TAG	UNP Q0P9J4
B	-13	VAL	-	EXPRESSION TAG	UNP Q0P9J4
B	-12	ASP	-	EXPRESSION TAG	UNP Q0P9J4
B	-11	LEU	-	EXPRESSION TAG	UNP Q0P9J4
B	-10	GLY	-	EXPRESSION TAG	UNP Q0P9J4
B	-9	THR	-	EXPRESSION TAG	UNP Q0P9J4
B	-8	GLU	-	EXPRESSION TAG	UNP Q0P9J4
B	-7	ASN	-	EXPRESSION TAG	UNP Q0P9J4
B	-6	LEU	-	EXPRESSION TAG	UNP Q0P9J4
B	-5	TYR	-	EXPRESSION TAG	UNP Q0P9J4
B	-4	PHE	-	EXPRESSION TAG	UNP Q0P9J4
B	-3	GLN	-	EXPRESSION TAG	UNP Q0P9J4
B	-2	SER	-	EXPRESSION TAG	UNP Q0P9J4
B	-1	ASN	-	EXPRESSION TAG	UNP Q0P9J4
B	0	ALA	-	EXPRESSION TAG	UNP Q0P9J4
B	92	GLY	-	LINKER	UNP Q0P9J4
C	-23	MET	-	EXPRESSION TAG	UNP Q0P9J4
C	-22	HIS	-	EXPRESSION TAG	UNP Q0P9J4
C	-21	HIS	-	EXPRESSION TAG	UNP Q0P9J4
C	-20	HIS	-	EXPRESSION TAG	UNP Q0P9J4
C	-19	HIS	-	EXPRESSION TAG	UNP Q0P9J4
C	-18	HIS	-	EXPRESSION TAG	UNP Q0P9J4
C	-17	HIS	-	EXPRESSION TAG	UNP Q0P9J4
C	-16	SER	-	EXPRESSION TAG	UNP Q0P9J4
C	-15	SER	-	EXPRESSION TAG	UNP Q0P9J4
C	-14	GLY	-	EXPRESSION TAG	UNP Q0P9J4
C	-13	VAL	-	EXPRESSION TAG	UNP Q0P9J4
C	-12	ASP	-	EXPRESSION TAG	UNP Q0P9J4
C	-11	LEU	-	EXPRESSION TAG	UNP Q0P9J4

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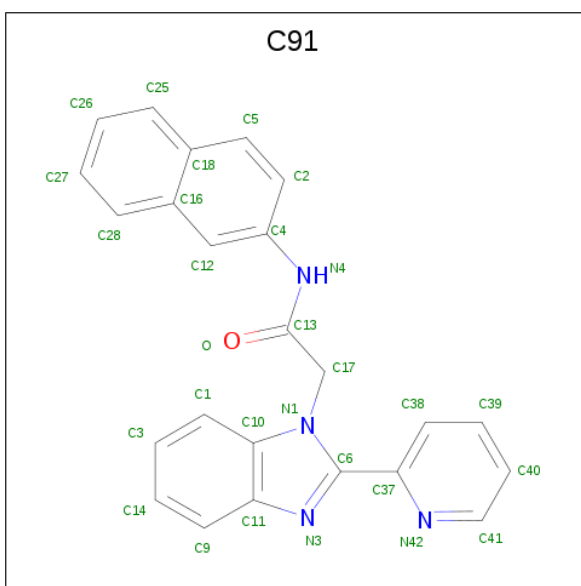
Chain	Residue	Modelled	Actual	Comment	Reference
C	-10	GLY	-	EXPRESSION TAG	UNP Q0P9J4
C	-9	THR	-	EXPRESSION TAG	UNP Q0P9J4
C	-8	GLU	-	EXPRESSION TAG	UNP Q0P9J4
C	-7	ASN	-	EXPRESSION TAG	UNP Q0P9J4
C	-6	LEU	-	EXPRESSION TAG	UNP Q0P9J4
C	-5	TYR	-	EXPRESSION TAG	UNP Q0P9J4
C	-4	PHE	-	EXPRESSION TAG	UNP Q0P9J4
C	-3	GLN	-	EXPRESSION TAG	UNP Q0P9J4
C	-2	SER	-	EXPRESSION TAG	UNP Q0P9J4
C	-1	ASN	-	EXPRESSION TAG	UNP Q0P9J4
C	0	ALA	-	EXPRESSION TAG	UNP Q0P9J4
C	92	GLY	-	LINKER	UNP Q0P9J4
D	-23	MET	-	EXPRESSION TAG	UNP Q0P9J4
D	-22	HIS	-	EXPRESSION TAG	UNP Q0P9J4
D	-21	HIS	-	EXPRESSION TAG	UNP Q0P9J4
D	-20	HIS	-	EXPRESSION TAG	UNP Q0P9J4
D	-19	HIS	-	EXPRESSION TAG	UNP Q0P9J4
D	-18	HIS	-	EXPRESSION TAG	UNP Q0P9J4
D	-17	HIS	-	EXPRESSION TAG	UNP Q0P9J4
D	-16	SER	-	EXPRESSION TAG	UNP Q0P9J4
D	-15	SER	-	EXPRESSION TAG	UNP Q0P9J4
D	-14	GLY	-	EXPRESSION TAG	UNP Q0P9J4
D	-13	VAL	-	EXPRESSION TAG	UNP Q0P9J4
D	-12	ASP	-	EXPRESSION TAG	UNP Q0P9J4
D	-11	LEU	-	EXPRESSION TAG	UNP Q0P9J4
D	-10	GLY	-	EXPRESSION TAG	UNP Q0P9J4
D	-9	THR	-	EXPRESSION TAG	UNP Q0P9J4
D	-8	GLU	-	EXPRESSION TAG	UNP Q0P9J4
D	-7	ASN	-	EXPRESSION TAG	UNP Q0P9J4
D	-6	LEU	-	EXPRESSION TAG	UNP Q0P9J4
D	-5	TYR	-	EXPRESSION TAG	UNP Q0P9J4
D	-4	PHE	-	EXPRESSION TAG	UNP Q0P9J4
D	-3	GLN	-	EXPRESSION TAG	UNP Q0P9J4
D	-2	SER	-	EXPRESSION TAG	UNP Q0P9J4
D	-1	ASN	-	EXPRESSION TAG	UNP Q0P9J4
D	0	ALA	-	EXPRESSION TAG	UNP Q0P9J4
D	92	GLY	-	LINKER	UNP Q0P9J4

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



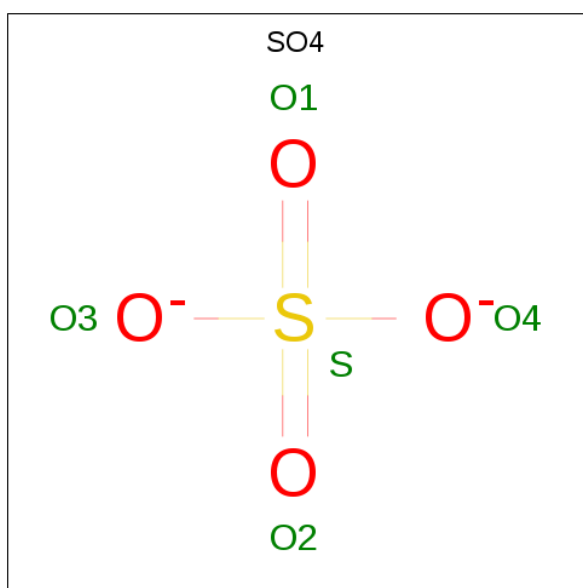
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			23	10	4	8	1		
2	B	1	Total	C	N	O	P	0	0
			23	10	4	8	1		
2	C	1	Total	C	N	O	P	0	0
			23	10	4	8	1		
2	D	1	Total	C	N	O	P	0	0
			23	10	4	8	1		

- Molecule 3 is N-(naphthalen-2-yl)-2-[2-(pyridin-2-yl)-1H-benzimidazol-1-yl]acetamide (three-letter code: C91) (formula: C₂₄H₁₈N₄O).



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

- Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



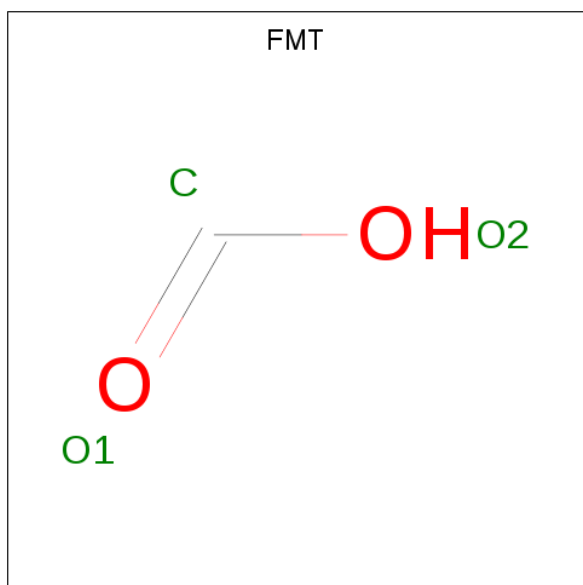
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	O	S	0	0
			5	4	1		
4	B	1	Total	O	S	0	0
			5	4	1		
4	B	1	Total	O	S	0	0
			5	4	1		
4	C	1	Total	O	S	0	0
			5	4	1		
4	C	1	Total	O	S	0	0
			5	4	1		
4	D	1	Total	O	S	0	0
			5	4	1		

- Molecule 5 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



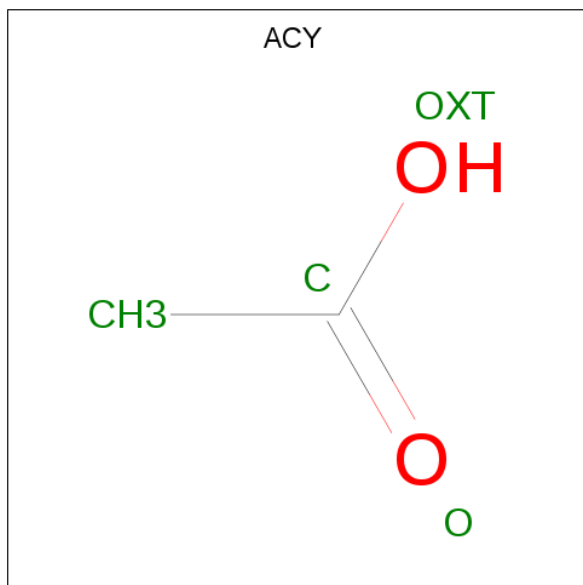
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			6	3	3		
5	B	1	Total	C	O	0	0
			6	3	3		
5	B	1	Total	C	O	0	0
			6	3	3		
5	D	1	Total	C	O	0	0
			6	3	3		

- Molecule 6 is FORMIC ACID (three-letter code: FMT) (formula: CH_2O_2).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			3	1	2		
6	B	1	Total	C	O	0	0
			3	1	2		

- Molecule 7 is ACETIC ACID (three-letter code: ACY) (formula: $C_2H_4O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total	C	O	0	0
			4	2	2		
7	B	1	Total	C	O	0	0
			4	2	2		

- Molecule 8 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: $C_2H_6O_2$).

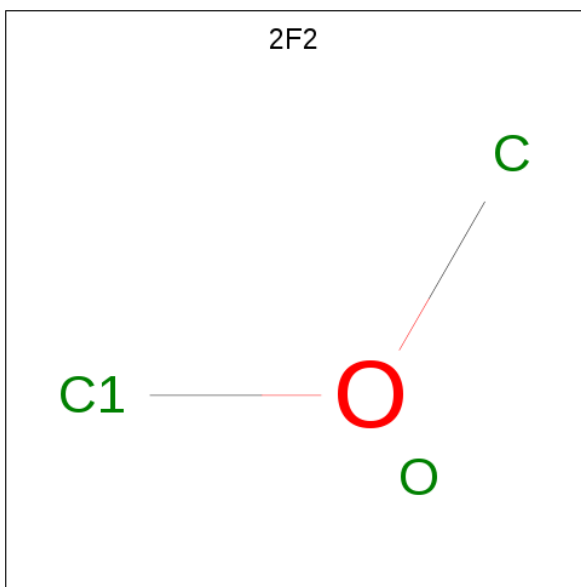


Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	A	1	Total C O 4 2 2	0	0
8	B	1	Total C O 4 2 2	0	0
8	B	1	Total C O 4 2 2	0	0

- Molecule 9 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
9	B	1	Total Cl 1 1	0	0
9	A	1	Total Cl 1 1	0	0
9	C	3	Total Cl 3 3	0	0

- Molecule 10 is dimethyl ether (three-letter code: 2F2) (formula: C₂H₆O).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
10	A	1	Total	C	O	0	0
			3	2	1		

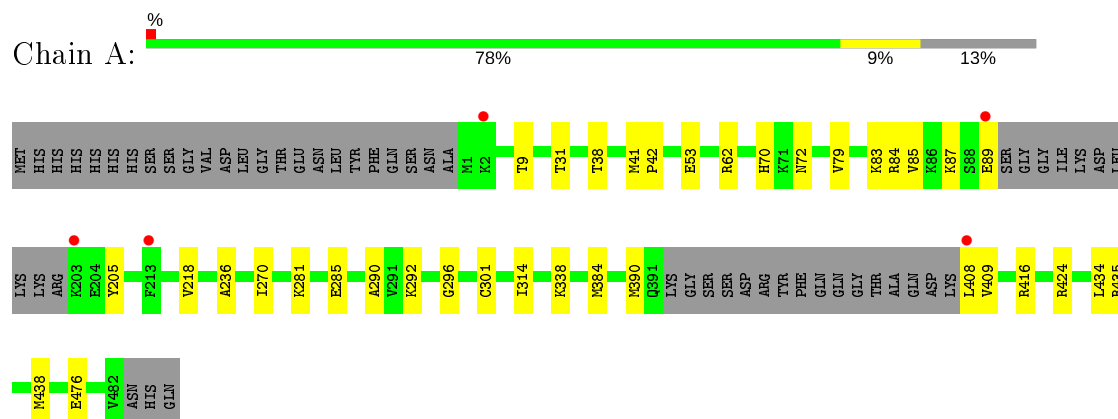
- Molecule 11 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
11	A	107	Total	O	0	0
			107	107		
11	B	114	Total	O	0	0
			114	114		
11	C	46	Total	O	0	0
			46	46		
11	D	27	Total	O	0	0
			27	27		

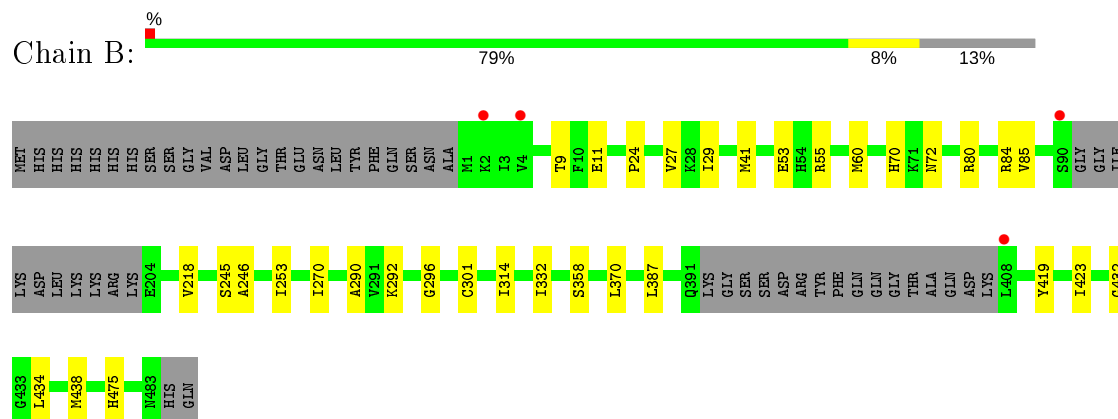
3 Residue-property plots [i](#)

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.

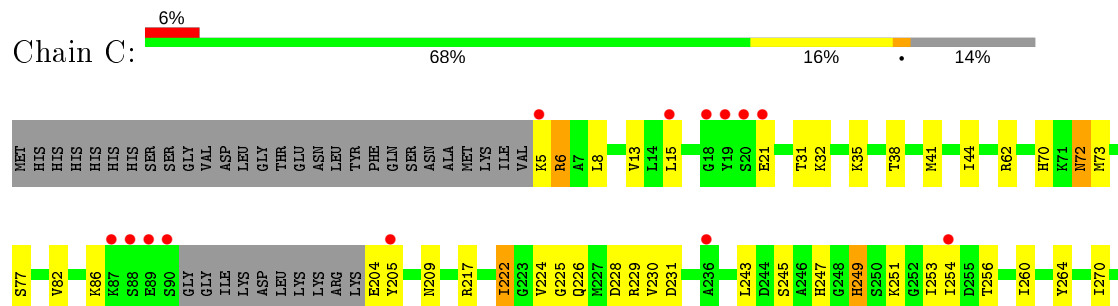
- Molecule 1: Inosine-5'-monophosphate dehydrogenase

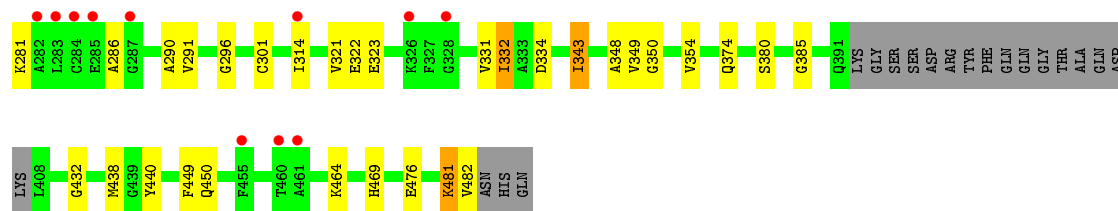


- Molecule 1: Inosine-5'-monophosphate dehydrogenase

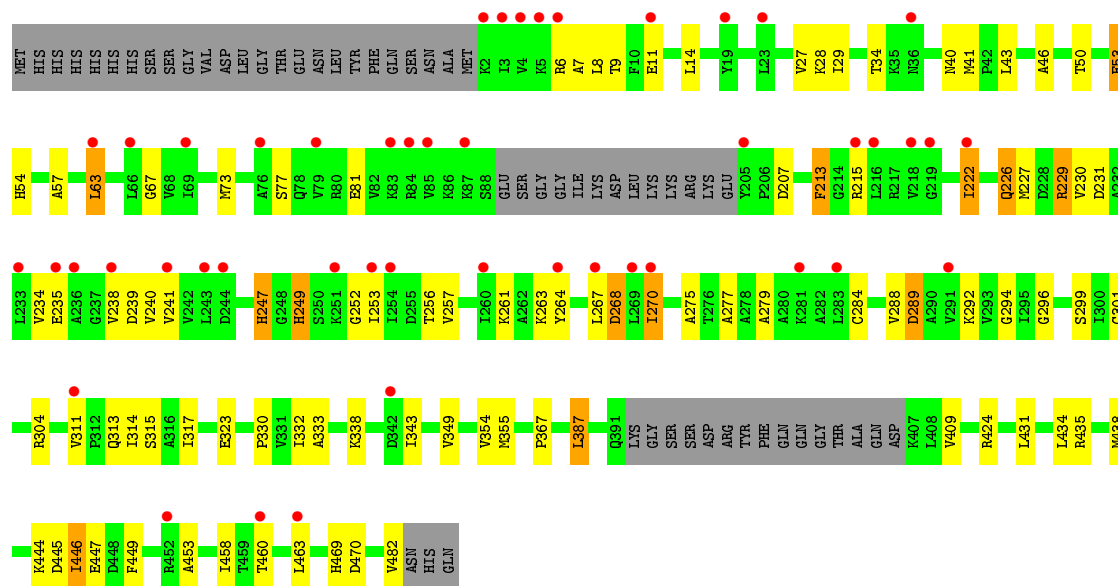


- Molecule 1: Inosine-5'-monophosphate dehydrogenase





● Molecule 1: Inosine-5'-monophosphate dehydrogenase



4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	114.46Å 114.46Å 256.31Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	39.23 – 2.50 47.55 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.7 (39.23-2.50) 99.7 (47.55-2.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.11	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.68 (at 2.51Å)	Xtriage
Refinement program	PHENIX (phenix.refine: dev_1367)	Depositor
R, R_{free}	0.176 , 0.217 0.180 , 0.217	Depositor DCC
R_{free} test set	3448 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å ²)	46.5	Xtriage
Anisotropy	0.255	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 51.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.018 for -h,-k,l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	11176	wwPDB-VP
Average B, all atoms (Å ²)	66.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.34% 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: 2F2, C91, GOL, CL, FMT, EDO, IMP, SO4, ACY

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	0.44	0/2704	0.59	0/3646
1	B	0.43	0/2726	0.59	0/3676
1	C	0.35	0/2653	0.53	0/3578
1	D	0.33	0/2641	0.53	0/3561
All	All	0.39	0/10724	0.56	0/14461

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 [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	2669	0	2759	22	0
1	B	2691	0	2770	21	0
1	C	2619	0	2681	37	0
1	D	2607	0	2694	61	0
2	A	23	0	11	1	0
2	B	23	0	11	2	0
2	C	23	0	11	3	0
2	D	23	0	11	2	0
3	A	29	0	18	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	29	0	18	3	0
3	C	29	0	18	0	0
3	D	29	0	18	1	0
4	A	5	0	0	0	0
4	B	10	0	0	0	0
4	C	10	0	0	0	0
4	D	5	0	0	0	0
5	A	6	0	8	0	0
5	B	12	0	16	0	0
5	D	6	0	8	0	0
6	A	3	0	1	0	0
6	B	3	0	1	0	0
7	A	4	0	3	0	0
7	B	4	0	3	1	0
8	A	4	0	6	1	0
8	B	8	0	12	2	0
9	A	1	0	0	0	0
9	B	1	0	0	0	0
9	C	3	0	0	0	0
10	A	3	0	6	0	0
11	A	107	0	0	1	0
11	B	114	0	0	0	0
11	C	46	0	0	1	0
11	D	27	0	0	2	0
All	All	11176	0	11084	142	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 6.

All (142) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:246:ALA:H	8:B:509:EDO:H12	1.50	0.76
1:D:41:MET:HG3	1:D:43:LEU:H	1.52	0.74
1:D:222:ILE:HD12	1:D:230:VAL:HG22	1.67	0.74
1:C:281:LYS:NZ	1:C:323:GLU:OE2	2.26	0.69
1:C:270:ILE:HG12	1:C:290:ALA:HB3	1.75	0.68
1:D:213:PHE:HB3	1:D:215:ARG:HG3	1.74	0.68
1:D:261:LYS:NZ	1:D:267:LEU:O	2.26	0.67
1:B:301:CYS:SG	2:B:501:IMP:H2	2.35	0.66
1:D:292:LYS:NZ	11:D:610:HOH:O	2.28	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:434:LEU:HG	1:A:438:MET:HE1	1.78	0.65
1:D:53:GLU:HB2	1:D:367:PRO:HG3	1.79	0.65
1:D:54:HIS:HB3	1:D:81:GLU:HG3	1.78	0.65
1:B:292:LYS:HG2	1:B:332:ILE:HB	1.79	0.64
1:A:89:GLU:HA	8:A:507:EDO:H12	1.80	0.64
1:D:277:ALA:HB1	1:D:323:GLU:HG3	1.81	0.63
2:B:501:IMP:H2'	7:B:506:ACY:H2	1.82	0.62
1:C:225:GLY:H	1:C:256:THR:HG21	1.64	0.62
1:D:296:GLY:HA2	1:D:301:CYS:SG	2.40	0.61
1:B:370:LEU:HD21	1:B:419:TYR:CD1	2.36	0.60
1:B:434:LEU:HG	1:B:438:MET:HE1	1.85	0.59
1:A:270:ILE:HG12	1:A:290:ALA:HB3	1.85	0.58
1:D:333:ALA:HB3	1:D:354:VAL:HG22	1.85	0.58
1:A:84[B]:ARG:HH11	1:A:84[B]:ARG:HB3	1.68	0.57
1:B:245:SER:HB2	1:B:253:ILE:HD11	1.86	0.57
1:D:40:ASN:ND2	1:D:207:ASP:O	2.37	0.56
1:C:296:GLY:HA2	1:C:301:CYS:SG	2.45	0.56
1:D:240:VAL:HG23	1:D:268:ASP:HB2	1.87	0.56
1:C:440:TYR:O	1:D:247:HIS:NE2	2.38	0.56
1:D:28:LYS:HD3	1:D:444:LYS:HE2	1.87	0.56
1:B:70:HIS:HD2	1:B:72:ASN:H	1.53	0.55
1:C:222:ILE:HD12	1:C:243:LEU:HD13	1.88	0.55
1:D:14:LEU:HD12	1:D:458:ILE:HG21	1.88	0.55
1:D:46:ALA:HA	1:D:355:MET:SD	2.47	0.54
1:C:321:VAL:HG11	1:C:350:GLY:HA3	1.89	0.54
1:D:268:ASP:N	1:D:268:ASP:OD1	2.39	0.54
1:D:41:MET:SD	1:D:43:LEU:HB2	2.49	0.53
1:C:21:GLU:OE1	1:C:21:GLU:N	2.38	0.53
1:D:252:GLY:O	1:D:256:THR:OG1	2.21	0.53
1:D:57:ALA:HA	1:D:67:GLY:HA3	1.91	0.52
1:A:435:ARG:HA	1:A:438:MET:HE2	1.91	0.52
1:D:275:ALA:HB1	1:D:311:VAL:HB	1.92	0.52
1:D:9:THR:HA	1:D:314:ILE:HB	1.93	0.51
1:D:447:GLU:N	1:D:447:GLU:OE1	2.39	0.51
1:A:9:THR:HA	1:A:314:ILE:HB	1.93	0.50
1:A:296:GLY:HA2	1:A:301:CYS:SG	2.51	0.50
1:C:222:ILE:HG22	1:C:229:ARG:HD2	1.93	0.50
1:D:28:LYS:HG2	1:D:444:LYS:O	2.12	0.50
1:D:296:GLY:HA3	1:D:304:ARG:HG3	1.93	0.50
1:A:301:CYS:SG	2:A:501:IMP:H2	2.52	0.49
1:C:82:VAL:O	1:C:86:LYS:HG2	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:73:MET:HG2	1:D:77:SER:HB2	1.95	0.48
1:B:270:ILE:HG12	1:B:290:ALA:HB3	1.94	0.48
1:D:226:GLN:HG2	1:D:229:ARG:HB2	1.96	0.48
1:C:343:ILE:CG2	1:C:354:VAL:HG21	2.43	0.48
1:C:8:LEU:O	1:C:314:ILE:HG22	2.13	0.48
1:A:390:MET:HG2	3:A:502:C91:C14	2.43	0.48
1:C:231:ASP:OD1	1:C:264:TYR:OH	2.21	0.48
1:C:70:HIS:HD2	1:C:72:ASN:H	1.62	0.48
1:D:27:VAL:HG11	1:D:438[A]:MET:SD	2.54	0.48
1:A:281:LYS:HE2	1:A:285:GLU:OE2	2.14	0.47
1:D:253:ILE:O	1:D:257:VAL:HG23	2.15	0.47
1:A:53:GLU:OE2	1:A:70:HIS:HE1	1.98	0.47
1:C:301:CYS:SG	2:C:501:IMP:H2	2.54	0.47
1:D:469:HIS:ND1	1:D:470:ASP:OD2	2.34	0.47
1:A:79:VAL:HG13	1:A:236:ALA:HB2	1.97	0.47
1:D:431:LEU:O	1:D:435:ARG:HG3	2.15	0.46
1:A:70:HIS:HD2	1:A:72:ASN:H	1.62	0.46
1:C:209:ASN:O	1:C:217[A]:ARG:HG3	2.16	0.46
1:D:29:ILE:HG12	11:D:627:HOH:O	2.16	0.46
1:D:11:GLU:HB3	1:D:463:LEU:HD22	1.98	0.46
1:A:62:ARG:HD3	1:A:205:TYR:CE2	2.50	0.46
1:D:270:ILE:HD11	1:D:292:LYS:HG3	1.97	0.46
1:D:317:ILE:HG23	1:D:349:VAL:HG11	1.97	0.46
1:A:409[A]:VAL:HG21	1:B:432:GLY:HA3	1.97	0.46
1:C:464:LYS:HB2	1:C:464:LYS:HE3	1.75	0.45
1:D:434:LEU:O	1:D:438[B]:MET:HG2	2.16	0.45
1:C:432:GLY:HA3	1:D:409:VAL:HG21	1.99	0.45
1:B:24:PRO:O	1:B:27:VAL:HG22	2.17	0.45
1:D:222:ILE:HD13	1:D:241:VAL:HG13	1.98	0.45
1:B:9:THR:HA	1:B:314:ILE:HB	1.98	0.44
1:D:289:ASP:N	1:D:289:ASP:OD1	2.50	0.44
1:D:314:ILE:HA	1:D:317:ILE:HG22	2.00	0.44
1:B:434:LEU:HG	1:B:438:MET:CE	2.47	0.44
1:D:301:CYS:SG	2:D:501:IMP:H2	2.57	0.44
1:B:296:GLY:HA2	1:B:301:CYS:SG	2.58	0.44
1:B:60:MET:CE	1:B:423:ILE:HD13	2.48	0.44
1:C:62:ARG:HD3	1:C:205:TYR:CZ	2.53	0.44
1:D:263:LYS:HG2	1:D:264:TYR:CD2	2.52	0.44
1:C:350:GLY:HA2	1:C:450:GLN:HG3	1.99	0.44
1:C:469:HIS:O	1:D:299:SER:HB2	2.17	0.44
1:C:438:MET:HB2	1:C:438:MET:HE3	1.84	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:502:C91:H9	3:D:502:C91:N42	2.33	0.44
1:B:29:ILE:HG22	1:B:438:MET:CE	2.48	0.43
1:A:72:ASN:HB2	1:A:384:MET:CE	2.48	0.43
3:B:502:C91:H8	3:B:502:C91:O	2.18	0.43
1:C:348:ALA:HB2	1:C:449:PHE:CE2	2.53	0.43
1:B:85:VAL:HG11	1:B:218:VAL:HB	2.00	0.43
1:C:476:GLU:HB3	1:C:481:LYS:HE3	2.01	0.43
1:D:313:GLN:O	1:D:317:ILE:HG22	2.18	0.43
1:C:6:ARG:HH21	1:D:6:ARG:HH12	1.66	0.43
1:D:43:LEU:HD11	1:D:343:ILE:HD13	1.99	0.43
3:B:502:C91:H9	3:B:502:C91:N42	2.34	0.43
1:C:226:GLN:O	1:C:230:VAL:HG23	2.18	0.43
1:C:291:VAL:HG23	1:C:331:VAL:HG22	2.01	0.43
1:D:289:ASP:O	1:D:330:PRO:HD2	2.19	0.43
1:D:387:LEU:HD22	1:D:387:LEU:HA	1.86	0.43
2:D:501:IMP:H8	2:D:501:IMP:H3'	2.00	0.43
1:A:476:GLU:HG2	11:A:690:HOH:O	2.18	0.43
1:D:34:THR:O	1:D:330:PRO:HG3	2.19	0.43
1:C:245:SER:HB2	1:C:253:ILE:HD11	2.00	0.42
1:B:29:ILE:HG22	1:B:438:MET:HE1	2.02	0.42
1:B:11[B]:GLU:CD	1:B:11[B]:GLU:H	2.22	0.42
1:A:85:VAL:HG11	1:A:218:VAL:HB	2.00	0.42
1:D:230:VAL:O	1:D:234:VAL:HG23	2.18	0.42
1:A:434:LEU:HG	1:A:438:MET:CE	2.47	0.42
1:D:446:ILE:O	1:D:449:PHE:HB3	2.19	0.42
1:C:247:HIS:ND1	11:C:635:HOH:O	2.37	0.42
1:D:446:ILE:H	1:D:446:ILE:HG12	1.52	0.42
1:A:31:THR:HG21	1:A:42:PRO:HB3	2.02	0.42
1:D:249:HIS:CE1	1:D:279:ALA:HA	2.55	0.42
1:D:28:LYS:HE3	1:D:445:ASP:HB3	2.00	0.42
1:B:53:GLU:OE2	1:B:70:HIS:HE1	2.02	0.41
1:C:31:THR:OG1	1:C:32:LYS:N	2.52	0.41
1:A:390:MET:HE1	1:A:408[A]:LEU:HD22	2.02	0.41
1:C:254:ILE:HG23	1:C:286:ALA:HB2	2.02	0.41
1:D:449:PHE:O	1:D:453:ALA:N	2.54	0.41
1:C:32:LYS:HA	1:C:38:THR:HA	2.02	0.41
1:B:80:ARG:NH1	1:B:84:ARG:HH12	2.19	0.41
1:C:224:VAL:HG23	1:C:253:ILE:HD13	2.02	0.41
1:D:7:ALA:HB1	1:D:315:SER:HB2	2.02	0.41
1:C:385:GLY:N	2:C:501:IMP:O6	2.41	0.41
3:B:502:C91:O	8:B:509:EDO:H11	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:334:ASP:OD2	2:C:501:IMP:O2'	2.35	0.41
1:D:226:GLN:HB2	1:D:226:GLN:HE21	1.66	0.41
1:D:284:CYS:HA	1:D:288:VAL:HG22	2.02	0.41
1:D:63:LEU:HD21	1:D:424:ARG:HA	2.03	0.41
1:B:387:LEU:HD12	1:B:387:LEU:HA	1.90	0.41
1:C:44:ILE:HD12	1:C:332:ILE:HD13	2.02	0.41
1:A:83:LYS:O	1:A:87:LYS:HG2	2.21	0.40
1:C:73:MET:HB3	1:C:77:SER:HB2	2.02	0.40
1:D:41:MET:HE1	1:D:63:LEU:HB3	2.02	0.40
1:D:231:ASP:O	1:D:235:GLU:HG2	2.21	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	351/406 (86%)	343 (98%)	8 (2%)	0	100	100
1	B	354/406 (87%)	344 (97%)	10 (3%)	0	100	100
1	C	346/406 (85%)	332 (96%)	13 (4%)	1 (0%)	41	61
1	D	345/406 (85%)	330 (96%)	14 (4%)	1 (0%)	41	61
All	All	1396/1624 (86%)	1349 (97%)	45 (3%)	2 (0%)	51	73

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	294	GLY
1	C	249	HIS

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	283/323 (88%)	277 (98%)	6 (2%)	53	78
1	B	286/323 (88%)	282 (99%)	4 (1%)	67	86
1	C	277/323 (86%)	255 (92%)	22 (8%)	12	24
1	D	276/323 (85%)	254 (92%)	22 (8%)	12	23
All	All	1122/1292 (87%)	1068 (95%)	54 (5%)	25	48

All (54) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	38	THR
1	A	41	MET
1	A	292	LYS
1	A	338	LYS
1	A	416	ARG
1	A	424	ARG
1	B	41	MET
1	B	55	ARG
1	B	358	SER
1	B	475	HIS
1	C	5	LYS
1	C	6	ARG
1	C	13	VAL
1	C	15	LEU
1	C	35	LYS
1	C	41	MET
1	C	72	ASN
1	C	204	GLU
1	C	222	ILE
1	C	228	ASP
1	C	249	HIS
1	C	251	LYS
1	C	260	ILE
1	C	322	GLU

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Mol	Chain	Res	Type
1	C	332	ILE
1	C	343	ILE
1	C	349	VAL
1	C	374[A]	GLN
1	C	374[B]	GLN
1	C	380	SER
1	C	481	LYS
1	C	482	VAL
1	D	8	LEU
1	D	50	THR
1	D	53	GLU
1	D	63	LEU
1	D	213	PHE
1	D	222	ILE
1	D	226	GLN
1	D	227	MET
1	D	229	ARG
1	D	238	VAL
1	D	239	ASP
1	D	247	HIS
1	D	249	HIS
1	D	268	ASP
1	D	270	ILE
1	D	289	ASP
1	D	332	ILE
1	D	338	LYS
1	D	387	LEU
1	D	446	ILE
1	D	460	THR
1	D	482	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	70	HIS
1	B	70	HIS
1	C	70	HIS
1	C	247	HIS
1	C	377	GLN
1	D	226	GLN
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 carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 31 ligands modelled in this entry, 5 are monoatomic - leaving 26 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
4	SO4	C	503	-	4,4,4	0.12	0	6,6,6	0.27	0
2	IMP	D	501	-	21,25,25	1.37	3 (14%)	23,38,38	1.42	2 (8%)
6	FMT	B	510	-	0,2,2	0.00	-	0,1,1	0.00	-
2	IMP	B	501	-	21,25,25	1.28	3 (14%)	23,38,38	1.61	4 (17%)
5	GOL	A	504	-	5,5,5	0.34	0	5,5,5	0.53	0
4	SO4	C	506	-	4,4,4	0.14	0	6,6,6	0.10	0
10	2F2	A	509	-	2,2,2	0.05	0	1,1,1	0.27	0
3	C91	A	502	-	30,33,33	1.73	10 (33%)	38,46,46	1.14	4 (10%)
8	EDO	A	507	-	3,3,3	0.51	0	2,2,2	0.30	0
5	GOL	D	504	-	5,5,5	0.39	0	5,5,5	0.30	0
4	SO4	B	505	-	4,4,4	0.15	0	6,6,6	0.09	0
8	EDO	B	509	-	3,3,3	0.54	0	2,2,2	0.26	0
3	C91	B	502	-	30,33,33	1.78	11 (36%)	38,46,46	1.20	5 (13%)
7	ACY	A	506	-	1,3,3	1.94	0	0,3,3	0.00	-
4	SO4	A	503	-	4,4,4	0.16	0	6,6,6	0.13	0
5	GOL	B	507	-	5,5,5	0.42	0	5,5,5	0.45	0
8	EDO	B	508	-	3,3,3	0.51	0	2,2,2	0.30	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	IMP	A	501	-	21,25,25	1.36	3 (14%)	23,38,38	1.78	5 (21%)
4	SO4	D	503	-	4,4,4	0.15	0	6,6,6	0.16	0
6	FMT	A	505	-	0,2,2	0.00	-	0,1,1	0.00	-
3	C91	D	502	-	30,33,33	1.80	8 (26%)	38,46,46	1.22	4 (10%)
3	C91	C	502	-	30,33,33	1.75	8 (26%)	38,46,46	1.18	4 (10%)
7	ACY	B	506	-	1,3,3	1.52	0	0,3,3	0.00	-
2	IMP	C	501	-	21,25,25	1.25	3 (14%)	23,38,38	1.59	4 (17%)
5	GOL	B	511	-	5,5,5	0.35	0	5,5,5	0.22	0
4	SO4	B	503	-	4,4,4	0.13	0	6,6,6	0.26	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
8	EDO	B	509	-	-	1/1/1/1	-
3	C91	B	502	-	-	0/12/12/12	0/5/5/5
3	C91	A	502	-	-	0/12/12/12	0/5/5/5
3	C91	D	502	-	-	0/12/12/12	0/5/5/5
3	C91	C	502	-	-	0/12/12/12	0/5/5/5
8	EDO	A	507	-	-	0/1/1/1	-
5	GOL	D	504	-	-	2/4/4/4	-
2	IMP	D	501	-	-	5/6/26/26	0/3/3/3
8	EDO	B	508	-	-	0/1/1/1	-
2	IMP	B	501	-	-	1/6/26/26	0/3/3/3
5	GOL	B	507	-	-	3/4/4/4	-
5	GOL	A	504	-	-	2/4/4/4	-
2	IMP	C	501	-	-	5/6/26/26	0/3/3/3
5	GOL	B	511	-	-	3/4/4/4	-
2	IMP	A	501	-	-	0/6/26/26	0/3/3/3

All (49) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	501	IMP	C2-N3	4.22	1.38	1.32
2	B	501	IMP	C2-N3	4.05	1.38	1.32
3	B	502	C91	C13-N4	3.96	1.44	1.35
3	C	502	C91	C37-C6	3.90	1.55	1.47
3	D	502	C91	C37-C6	3.80	1.55	1.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501	IMP	C2-N3	3.78	1.38	1.32
3	A	502	C91	C13-N4	3.71	1.43	1.35
3	D	502	C91	C13-N4	3.68	1.43	1.35
2	C	501	IMP	C2-N3	3.62	1.37	1.32
3	C	502	C91	C13-N4	3.44	1.43	1.35
2	A	501	IMP	C6-N1	3.36	1.38	1.33
2	D	501	IMP	C6-N1	3.17	1.38	1.33
2	C	501	IMP	C6-N1	3.11	1.38	1.33
3	B	502	C91	C26-C25	3.07	1.43	1.36
3	A	502	C91	C37-C6	3.01	1.53	1.47
3	D	502	C91	C5-C2	2.97	1.42	1.36
3	B	502	C91	C37-C6	2.95	1.53	1.47
3	B	502	C91	C5-C2	2.89	1.42	1.36
3	A	502	C91	C26-C25	2.89	1.43	1.36
3	A	502	C91	C5-C2	2.82	1.42	1.36
3	D	502	C91	C26-C25	2.80	1.43	1.36
3	C	502	C91	C26-C25	2.79	1.43	1.36
2	B	501	IMP	C6-N1	2.76	1.37	1.33
3	C	502	C91	C16-C18	-2.67	1.36	1.42
3	C	502	C91	C5-C2	2.65	1.42	1.36
3	A	502	C91	C16-C18	-2.56	1.36	1.42
3	D	502	C91	C16-C18	-2.52	1.36	1.42
2	A	501	IMP	C2-N1	2.51	1.38	1.33
2	D	501	IMP	C2-N1	2.47	1.38	1.33
3	B	502	C91	C28-C16	2.46	1.47	1.41
3	D	502	C91	C3-C1	2.46	1.42	1.36
3	B	502	C91	C3-C1	2.41	1.42	1.36
3	D	502	C91	C9-C11	2.39	1.46	1.41
3	B	502	C91	C16-C18	-2.38	1.36	1.42
3	C	502	C91	C3-C1	2.38	1.42	1.36
2	C	501	IMP	C2-N1	2.37	1.38	1.33
3	A	502	C91	C3-C1	2.32	1.42	1.36
3	D	502	C91	C28-C16	2.31	1.47	1.41
3	A	502	C91	C27-C28	2.30	1.41	1.36
3	A	502	C91	C28-C16	2.30	1.47	1.41
3	C	502	C91	C9-C11	2.28	1.45	1.41
2	B	501	IMP	C2-N1	2.27	1.38	1.33
3	C	502	C91	C28-C16	2.21	1.47	1.41
3	B	502	C91	C27-C28	2.21	1.41	1.36
3	B	502	C91	C9-C11	2.17	1.45	1.41
3	B	502	C91	C12-C4	2.09	1.42	1.37
3	B	502	C91	C4-N4	2.08	1.45	1.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	502	C91	C12-C4	2.07	1.42	1.37
3	A	502	C91	C9-C11	2.03	1.45	1.41

All (32) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	501	IMP	N3-C2-N1	-5.49	120.10	128.68
2	D	501	IMP	N3-C2-N1	-5.35	120.31	128.68
2	A	501	IMP	N3-C2-N1	-5.04	120.81	128.68
2	B	501	IMP	N3-C2-N1	-4.70	121.33	128.68
2	A	501	IMP	C1'-N9-C4	-3.64	120.25	126.64
3	D	502	C91	C6-C37-N42	3.13	122.13	116.35
2	A	501	IMP	C2-N1-C6	3.10	121.08	115.88
3	C	502	C91	C6-N3-C11	3.07	109.88	103.78
3	D	502	C91	C6-N3-C11	2.99	109.70	103.78
3	A	502	C91	C6-C37-N42	2.97	121.85	116.35
3	B	502	C91	C6-C37-N42	2.93	121.78	116.35
2	C	501	IMP	C2-N1-C6	2.93	120.79	115.88
3	C	502	C91	C6-C37-N42	2.91	121.74	116.35
2	B	501	IMP	C1'-N9-C4	-2.88	121.58	126.64
3	D	502	C91	C41-N42-C37	2.85	121.23	117.23
3	C	502	C91	N3-C6-N1	-2.85	107.97	115.11
2	B	501	IMP	C2-N1-C6	2.83	120.62	115.88
3	B	502	C91	C6-N3-C11	2.73	109.20	103.78
3	A	502	C91	C6-N3-C11	2.72	109.18	103.78
3	C	502	C91	C41-N42-C37	2.71	121.03	117.23
3	D	502	C91	N3-C6-N1	-2.66	108.43	115.11
2	D	501	IMP	C2-N1-C6	2.46	120.01	115.88
3	B	502	C91	O-C13-C17	-2.45	116.04	120.04
2	B	501	IMP	O3P-P-O5'	-2.37	100.41	106.73
3	B	502	C91	C41-N42-C37	2.37	120.55	117.23
3	B	502	C91	N3-C6-N1	-2.33	109.27	115.11
3	A	502	C91	N3-C6-N1	-2.28	109.38	115.11
3	A	502	C91	C41-N42-C37	2.22	120.34	117.23
2	A	501	IMP	C4-C5-N7	-2.13	107.18	109.40
2	C	501	IMP	O2P-P-O1P	2.07	118.79	110.68
2	C	501	IMP	C1'-N9-C4	-2.07	123.01	126.64
2	A	501	IMP	O5'-P-O1P	2.05	112.22	106.47

There are no chirality outliers.

All (22) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	D	501	IMP	C5'-O5'-P-O1P
2	D	501	IMP	C5'-O5'-P-O2P
2	D	501	IMP	C5'-O5'-P-O3P
2	D	501	IMP	C3'-C4'-C5'-O5'
5	A	504	GOL	O1-C1-C2-C3
5	D	504	GOL	O1-C1-C2-O2
5	D	504	GOL	O1-C1-C2-C3
5	B	507	GOL	C1-C2-C3-O3
2	C	501	IMP	C5'-O5'-P-O1P
2	C	501	IMP	C5'-O5'-P-O2P
2	C	501	IMP	C5'-O5'-P-O3P
2	C	501	IMP	C3'-C4'-C5'-O5'
5	B	511	GOL	O1-C1-C2-C3
2	D	501	IMP	O4'-C4'-C5'-O5'
2	C	501	IMP	O4'-C4'-C5'-O5'
5	B	507	GOL	O1-C1-C2-C3
5	A	504	GOL	O1-C1-C2-O2
5	B	507	GOL	O2-C2-C3-O3
5	B	511	GOL	O1-C1-C2-O2
8	B	509	EDO	O1-C1-C2-O2
2	B	501	IMP	C3'-C4'-C5'-O5'
5	B	511	GOL	C1-C2-C3-O3

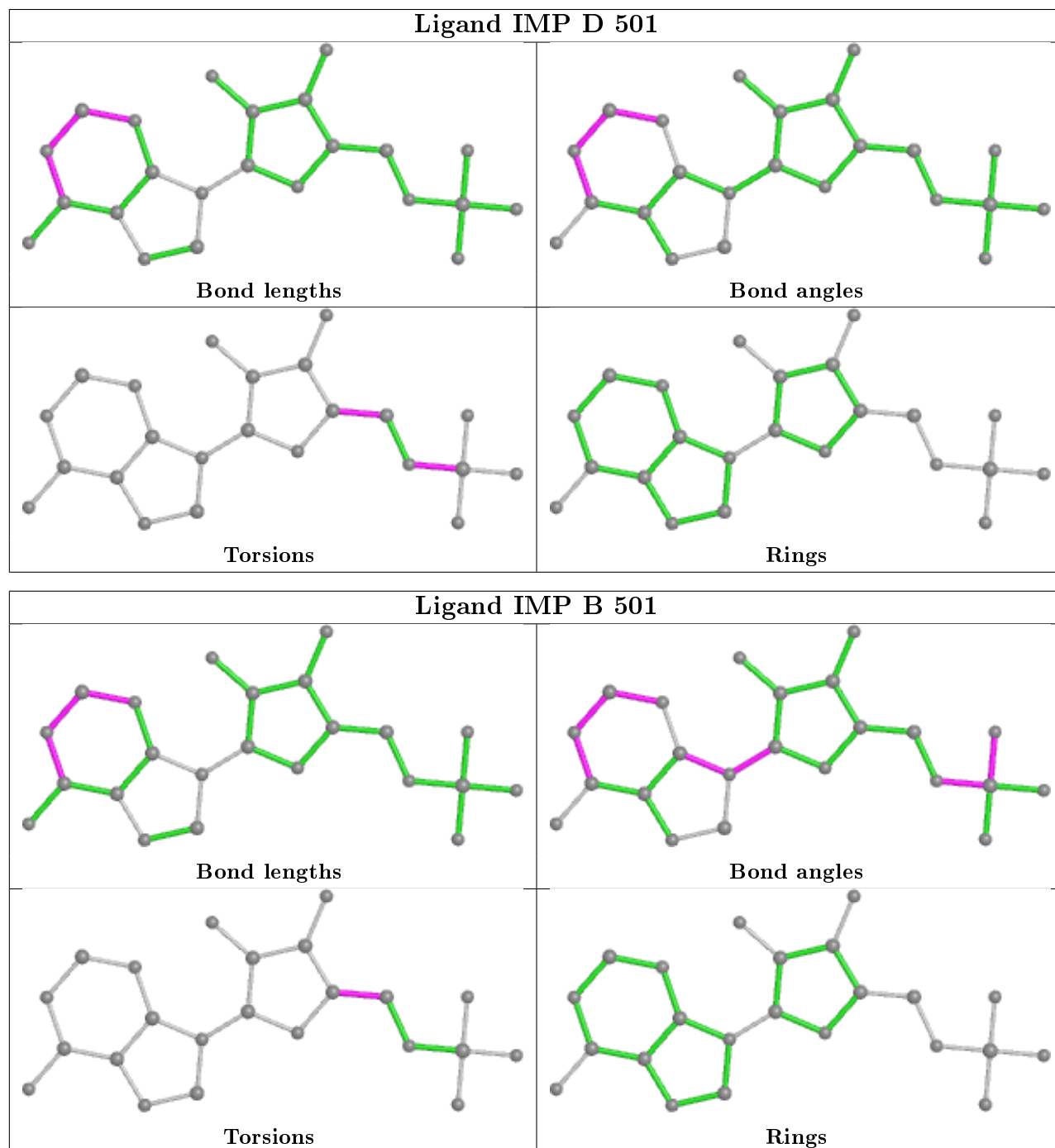
There are no ring outliers.

10 monomers are involved in 15 short contacts:

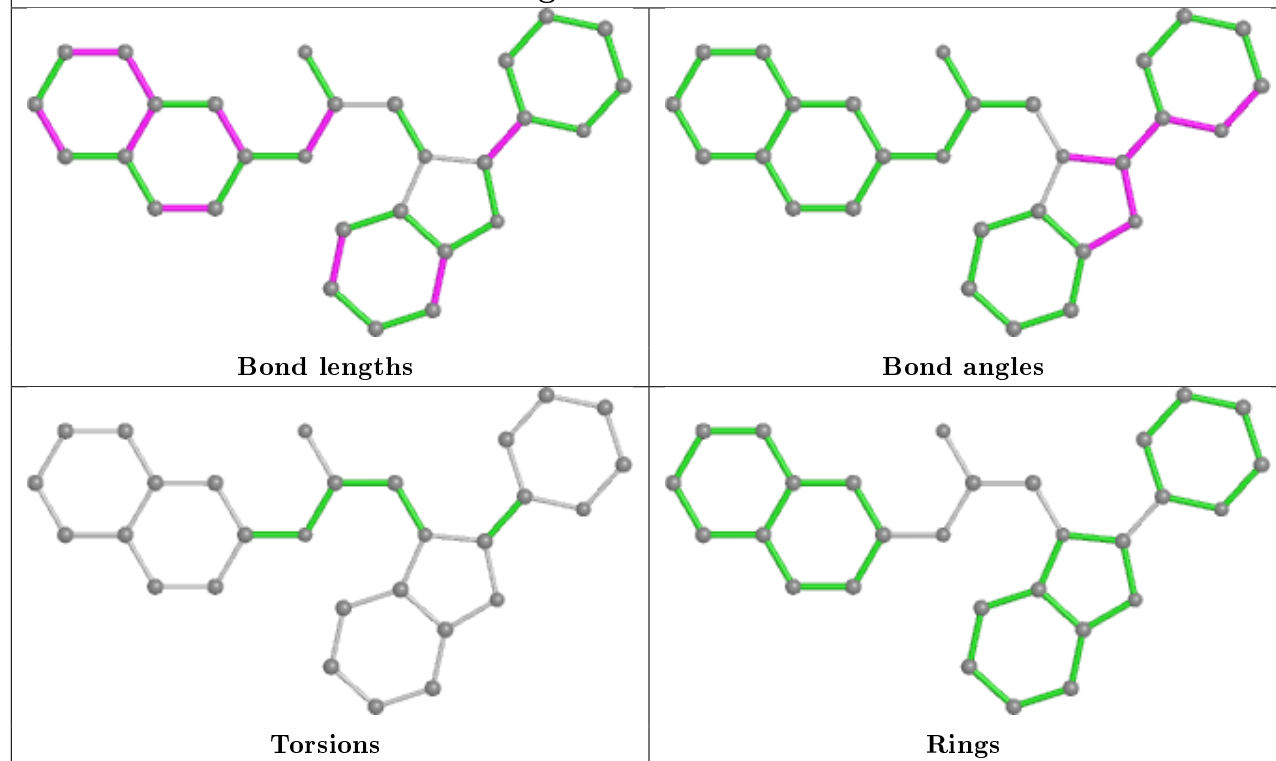
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	501	IMP	2	0
2	B	501	IMP	2	0
3	A	502	C91	1	0
8	A	507	EDO	1	0
8	B	509	EDO	2	0
3	B	502	C91	3	0
2	A	501	IMP	1	0
3	D	502	C91	1	0
7	B	506	ACY	1	0
2	C	501	IMP	3	0

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.

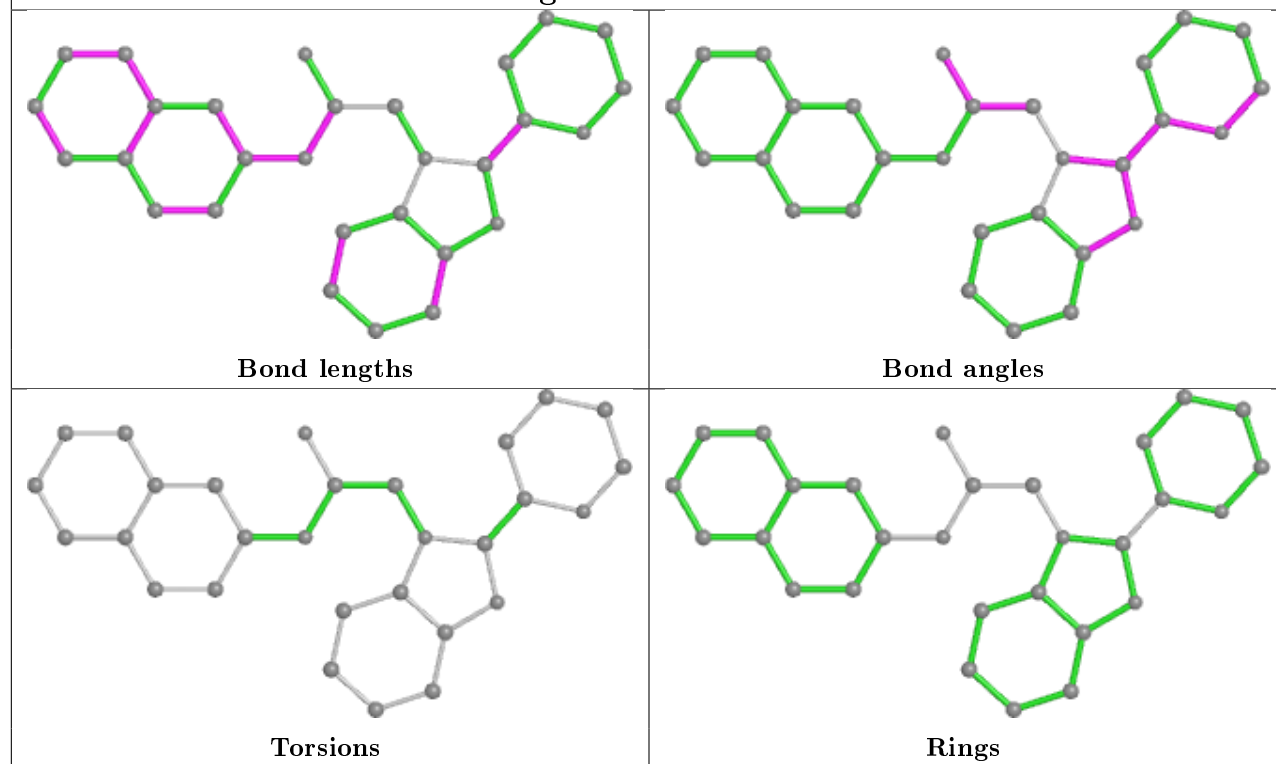
Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD 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.



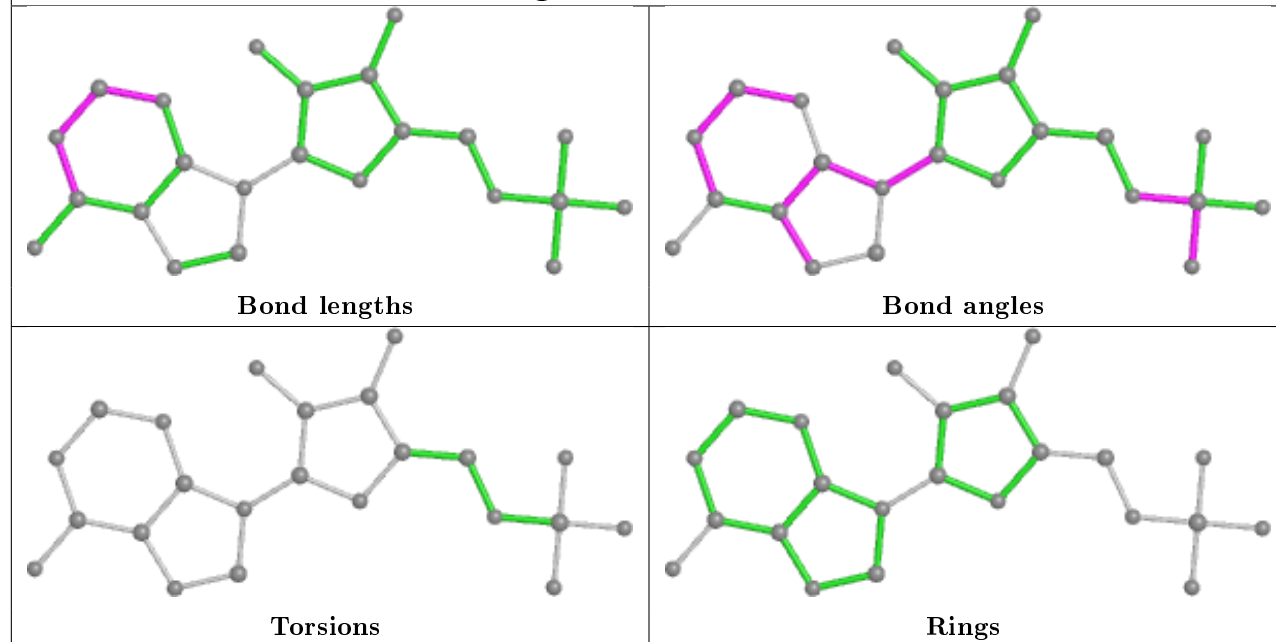
Ligand C91 A 502



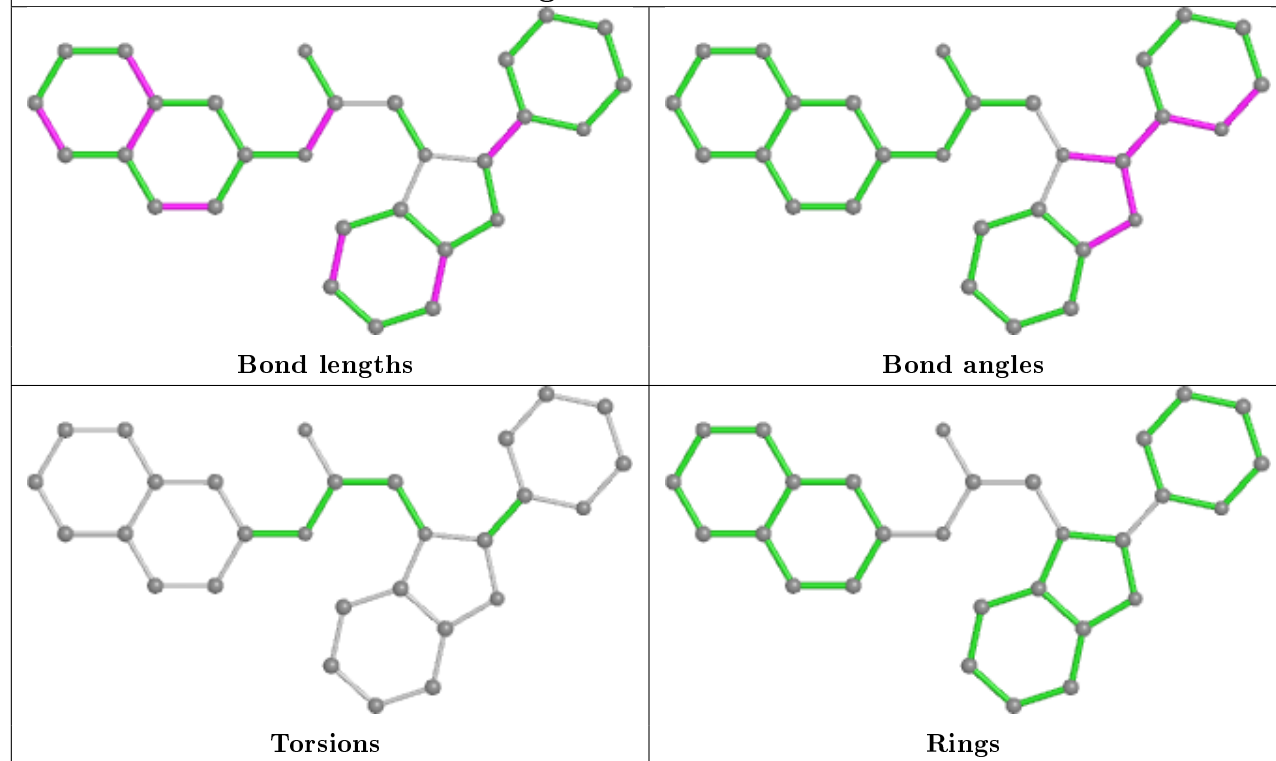
Ligand C91 B 502

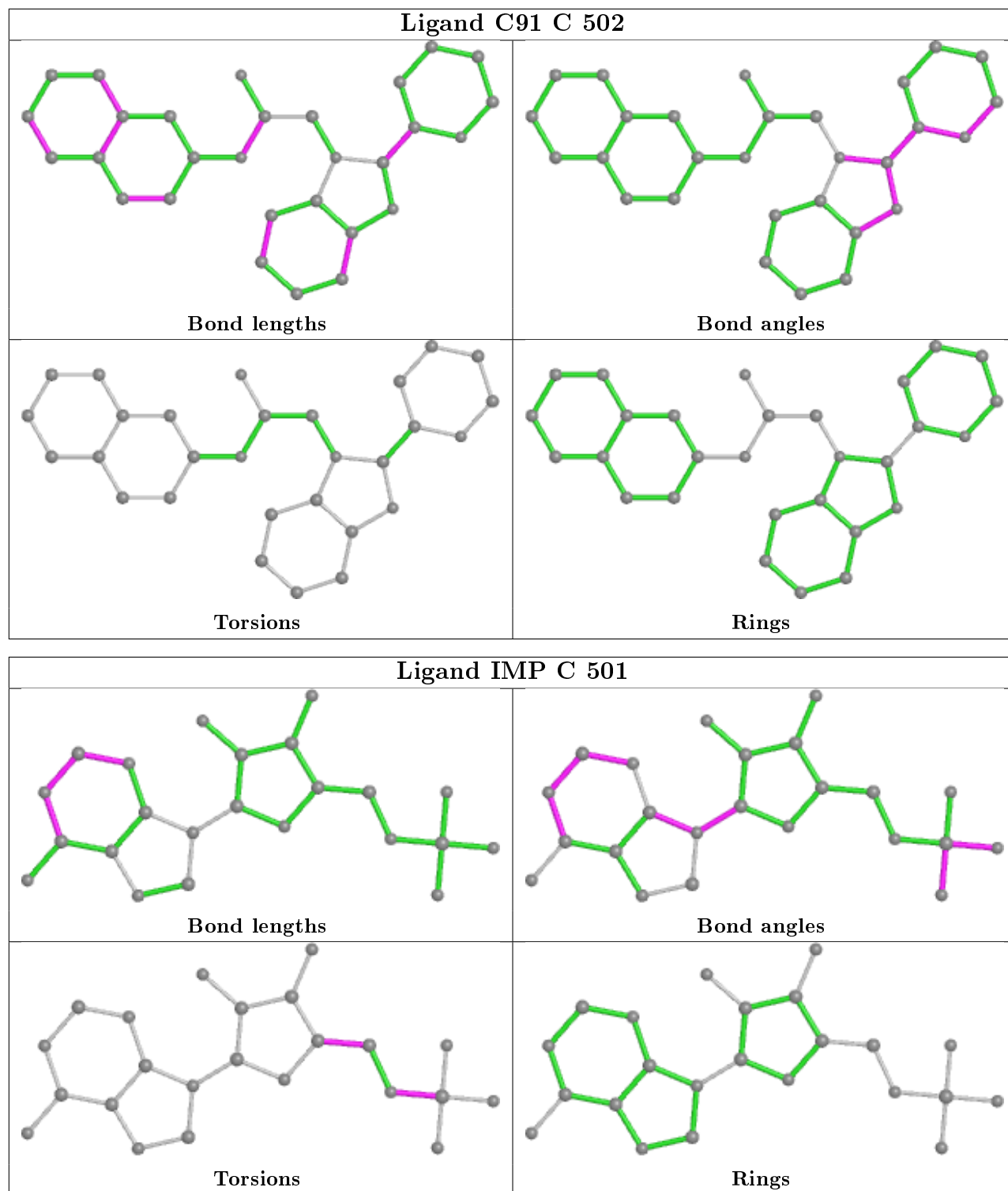


Ligand IMP A 501



Ligand C91 D 502





5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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	353/406 (86%)	-0.21	5 (1%) 75 77	31, 43, 69, 112	0
1	B	354/406 (87%)	-0.26	4 (1%) 80 82	31, 44, 67, 120	0
1	C	349/406 (85%)	0.42	24 (6%) 16 17	35, 78, 110, 134	0
1	D	350/406 (86%)	0.74	47 (13%) 3 2	39, 94, 124, 140	0
All	All	1406/1624 (86%)	0.17	80 (5%) 23 25	31, 60, 114, 140	0

All (80) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	408[A]	LEU	8.2
1	C	283	LEU	6.0
1	D	3	ILE	4.9
1	C	328	GLY	4.8
1	C	19	TYR	4.6
1	D	79	VAL	4.6
1	D	83	LYS	4.5
1	D	218	VAL	4.5
1	D	260	ILE	4.4
1	D	5	LYS	4.4
1	C	90	SER	4.3
1	C	326	LYS	3.9
1	D	244	ASP	3.8
1	C	88	SER	3.8
1	A	89	GLU	3.8
1	C	455	PHE	3.7
1	C	89	GLU	3.7
1	D	236	ALA	3.7
1	D	251	LYS	3.6
1	D	85	VAL	3.6
1	C	236	ALA	3.5

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Mol	Chain	Res	Type	RSRZ
1	D	235	GLU	3.4
1	C	21	GLU	3.4
1	C	254	ILE	3.3
1	D	4	VAL	3.3
1	D	460	THR	3.3
1	C	18	GLY	3.2
1	D	11	GLU	3.2
1	D	267	LEU	3.2
1	D	243	LEU	3.2
1	C	284	CYS	3.2
1	C	205	TYR	3.1
1	A	213	PHE	3.1
1	D	84	ARG	3.1
1	B	90	SER	3.1
1	C	287	GLY	3.1
1	C	285	GLU	3.0
1	D	254	ILE	3.0
1	D	291	VAL	3.0
1	A	2	LYS	3.0
1	B	4	VAL	2.9
1	A	408[A]	LEU	2.9
1	D	87	LYS	2.9
1	C	282	ALA	2.8
1	D	216	LEU	2.8
1	D	238	VAL	2.7
1	D	270	ILE	2.7
1	D	269	LEU	2.6
1	D	19	TYR	2.6
1	D	205	TYR	2.6
1	C	20	SER	2.6
1	C	461	ALA	2.6
1	D	283	LEU	2.6
1	A	203	LYS	2.5
1	D	36	ASN	2.5
1	D	264	TYR	2.5
1	D	311	VAL	2.5
1	D	463	LEU	2.4
1	D	2	LYS	2.3
1	D	233	LEU	2.3
1	D	253	ILE	2.3
1	C	5	LYS	2.2
1	D	66	LEU	2.2

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Mol	Chain	Res	Type	RSRZ
1	D	76	ALA	2.2
1	D	452	ARG	2.2
1	D	241	VAL	2.2
1	D	342	ASP	2.2
1	C	87	LYS	2.1
1	D	222	ILE	2.1
1	D	63	LEU	2.1
1	D	215	ARG	2.1
1	C	460	THR	2.1
1	C	15	LEU	2.1
1	D	23	LEU	2.1
1	C	314	ILE	2.0
1	D	281	LYS	2.0
1	B	2	LYS	2.0
1	D	6	ARG	2.0
1	D	219	GLY	2.0
1	D	69	ILE	2.0

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

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 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
8	EDO	A	507	4/4	0.80	0.27	78,79,79,80	0
3	C91	C	502	29/29	0.82	0.36	99,116,145,145	0
5	GOL	B	507	6/6	0.84	0.42	58,66,69,75	0
9	CL	B	504	1/1	0.86	0.26	85,85,85,85	0
7	ACY	A	506	4/4	0.86	0.19	77,80,80,82	0
9	CL	C	505	1/1	0.88	0.07	88,88,88,88	0

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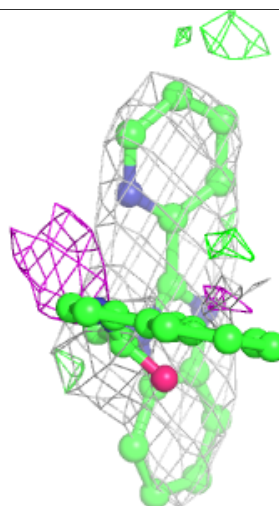
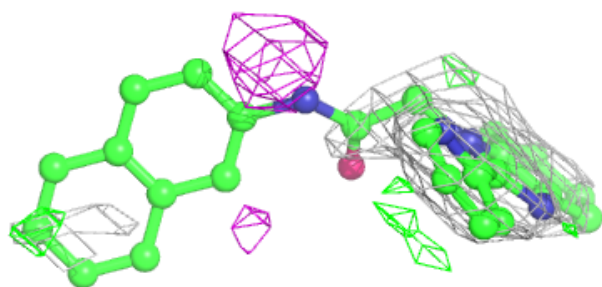
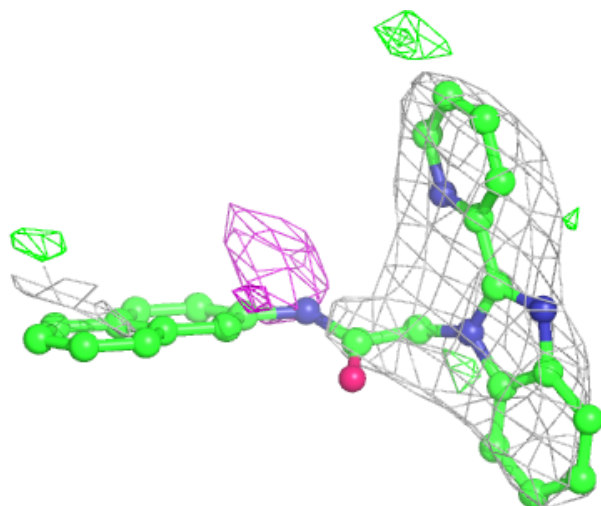
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
8	EDO	B	509	4/4	0.89	0.27	54,57,62,64	0
3	C91	D	502	29/29	0.89	0.36	92,94,96,97	0
7	ACY	B	506	4/4	0.90	0.34	83,84,84,84	0
5	GOL	D	504	6/6	0.91	0.13	79,83,85,86	0
8	EDO	B	508	4/4	0.92	0.11	55,57,59,63	0
9	CL	C	504	1/1	0.92	0.06	81,81,81,81	0
5	GOL	B	511	6/6	0.92	0.11	74,75,78,80	0
5	GOL	A	504	6/6	0.93	0.12	71,73,77,80	0
4	SO4	C	506	5/5	0.93	0.14	128,130,131,131	0
10	2F2	A	509	3/3	0.93	0.13	48,48,54,54	0
6	FMT	B	510	3/3	0.94	0.14	54,54,55,56	0
2	IMP	D	501	23/23	0.94	0.13	40,58,70,77	0
6	FMT	A	505	3/3	0.94	0.17	61,61,62,63	0
4	SO4	B	505	5/5	0.94	0.23	117,119,121,121	0
3	C91	A	502	29/29	0.95	0.17	50,63,76,78	0
3	C91	B	502	29/29	0.96	0.17	41,60,62,63	0
4	SO4	D	503	5/5	0.96	0.19	83,84,85,86	0
4	SO4	C	503	5/5	0.97	0.15	68,69,70,72	0
2	IMP	C	501	23/23	0.97	0.12	47,53,63,67	0
4	SO4	A	503	5/5	0.97	0.12	68,72,78,78	0
4	SO4	B	503	5/5	0.97	0.13	75,77,77,80	0
9	CL	C	507	1/1	0.98	0.36	79,79,79,79	1
2	IMP	A	501	23/23	0.98	0.15	28,33,38,41	0
9	CL	A	508	1/1	0.99	0.13	70,70,70,70	1
2	IMP	B	501	23/23	0.99	0.18	26,39,44,46	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.

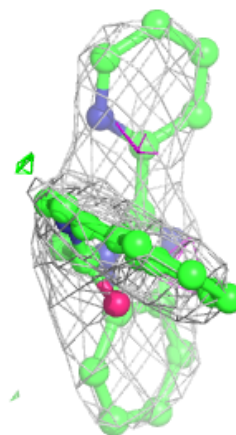
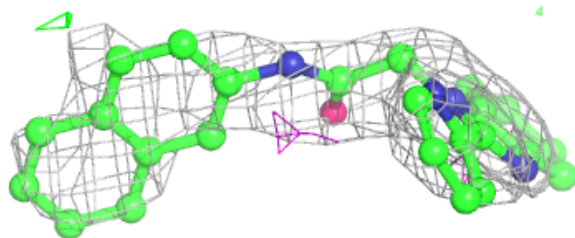
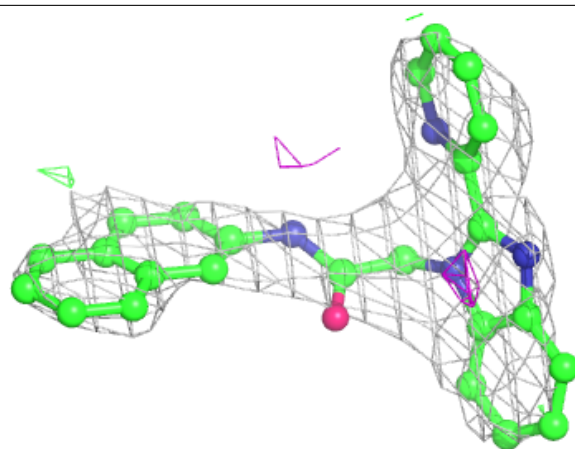
Electron density around C91 C 502:

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

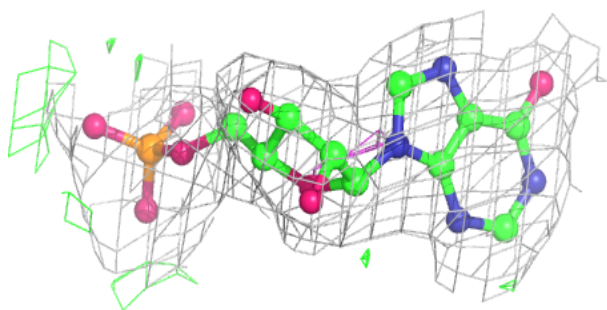
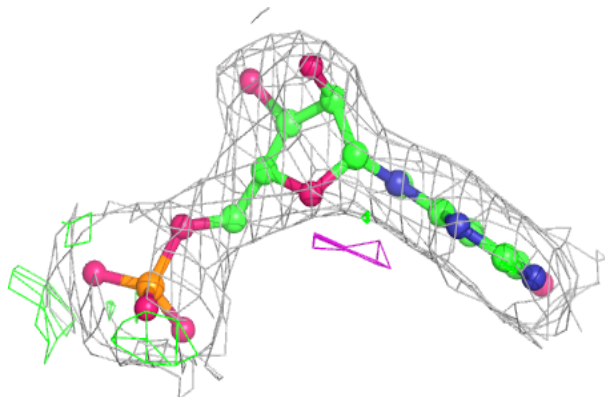


Electron density around C91 D 502:

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

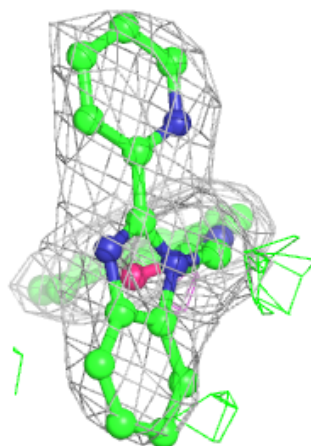
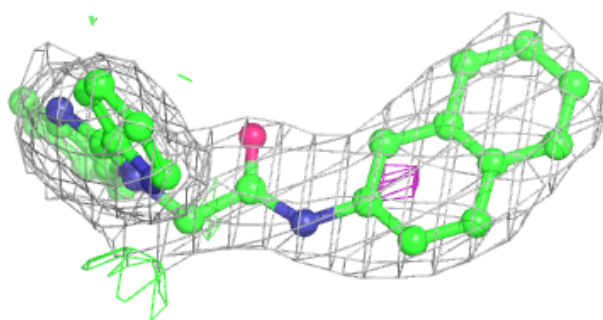
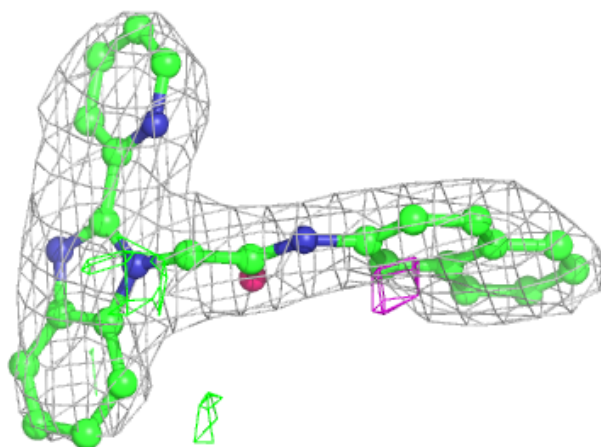
**Electron density around IMP D 501:**

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



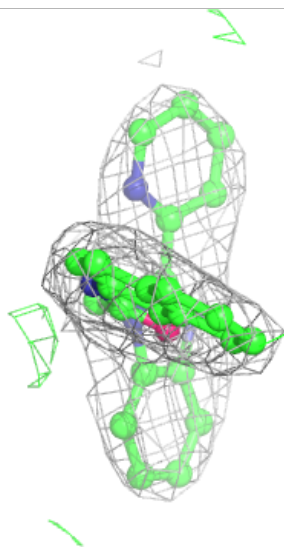
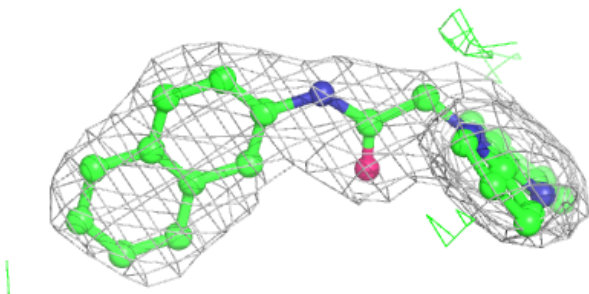
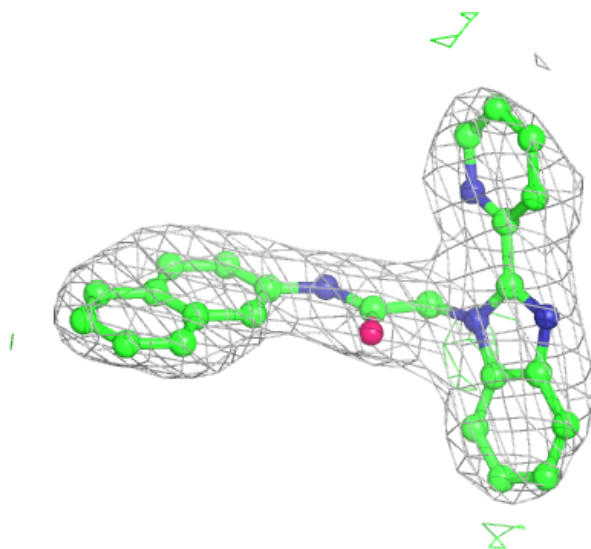
Electron density around C91 A 502:

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



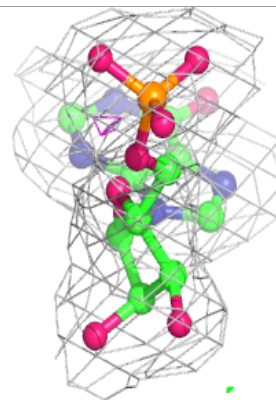
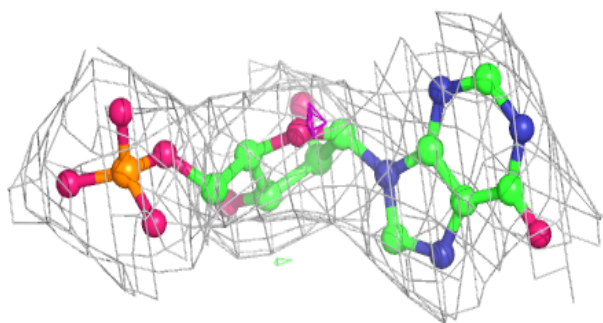
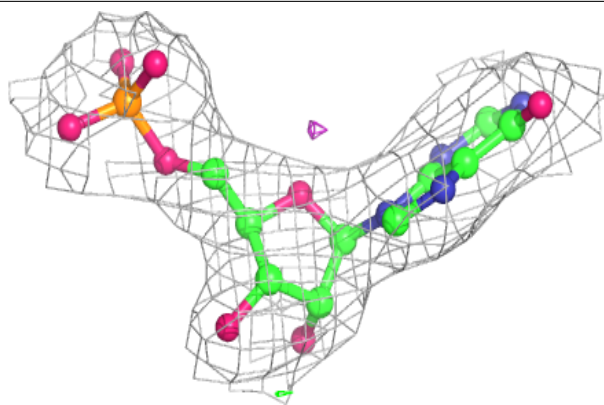
Electron density around C91 B 502:

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

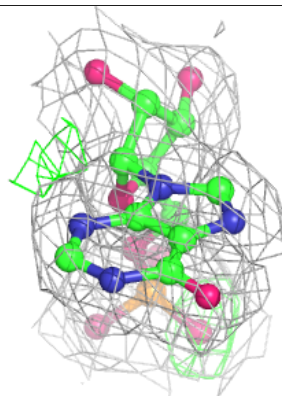
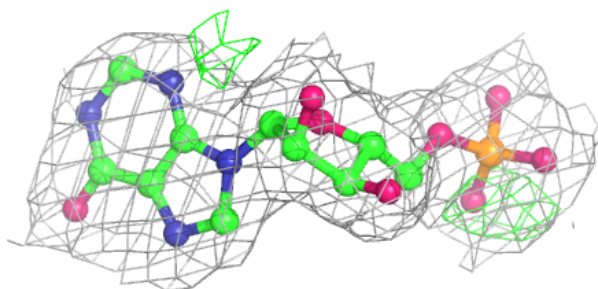
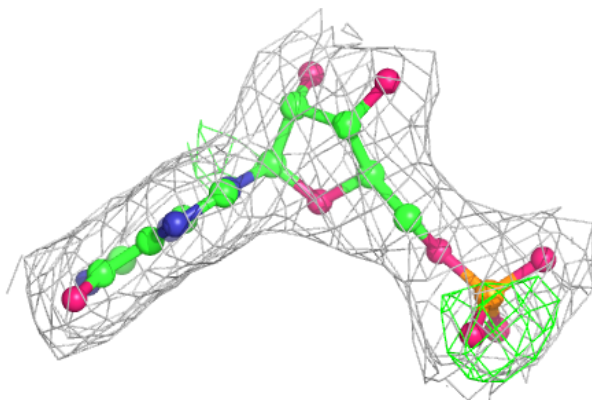


Electron density around IMP C 501:

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

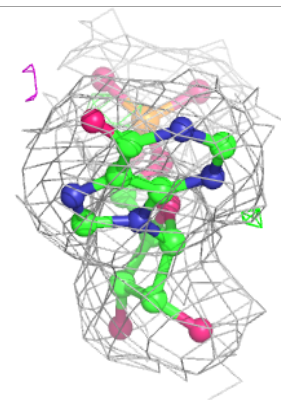
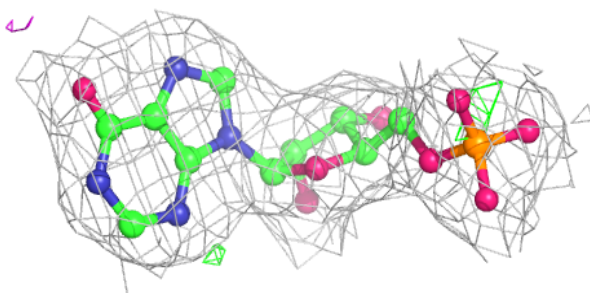
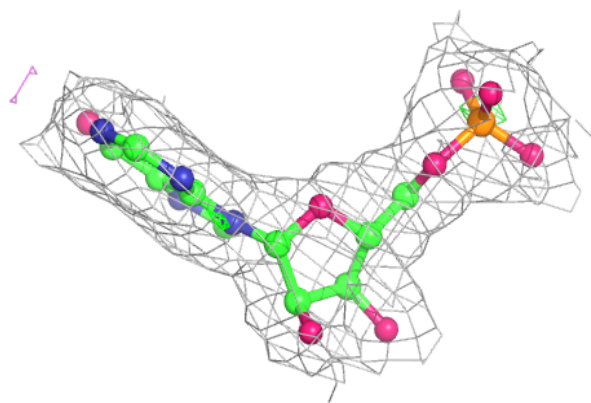
**Electron density around IMP A 501:**

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



Electron density around IMP 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)



6.5 Other polymers [i](#)

There are no such residues in this entry.