Full wwPDB X-ray Structure Validation Report (i)

Jun 13, 2024 – 07:32 AM EDT

PDB ID	:	4DPD
Title	:	WILD TYPE PLASMODIUM FALCIPARUM DIHYDROFOLATE REDUCTASE-THYMIDYLATE SYNTHASE (PfDHFR-TS), DHF COMPLEX, NADP+, dUMP
Authors	:	Yuthavong, Y.; Vilaivan, T.; Kamchonwongpaisan, S.; Charman, S.A.; McLennan, D.N.; White, K.L.; Vivas, L.; Bongard, E.; Chitnumsub, P.; Tarnchompoo, B.; Thongphanchang, C.; Taweechai, S.; Vanichtanakul, J.; Arwon, U.; Fantauzzi, P.; Yuvaniyama, J.; Charman, W.N.; Matthews, D.
Deposited on	:	2012-02-13
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 (i) symbol.

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

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

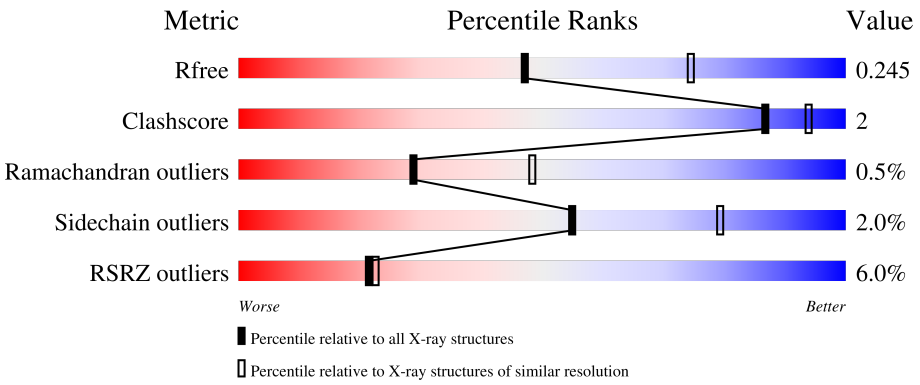
MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	2.36.2
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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 of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=5% The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	608	2% 80% 7% 13%
1	B	608	8% 83% 5% 11%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard

Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.36.2

residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	DHF	A	701	-	-	-	X

2 Entry composition (i)

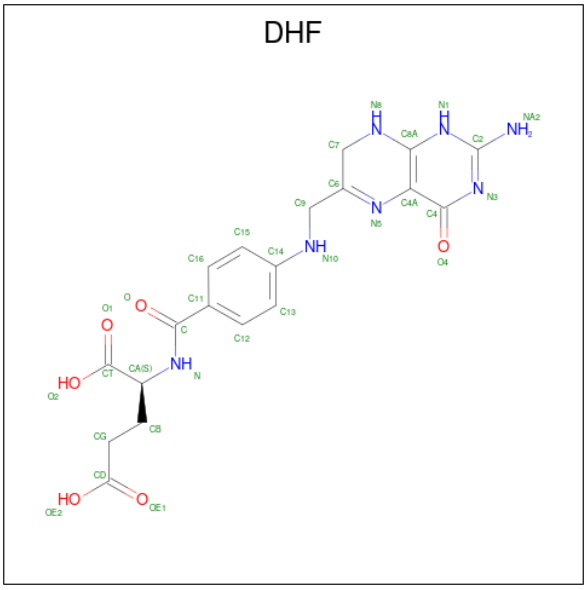
There are 5 unique types of molecules in this entry. The entry contains 9454 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Bifunctional dihydrofolate reductase-thymidylate synthase.

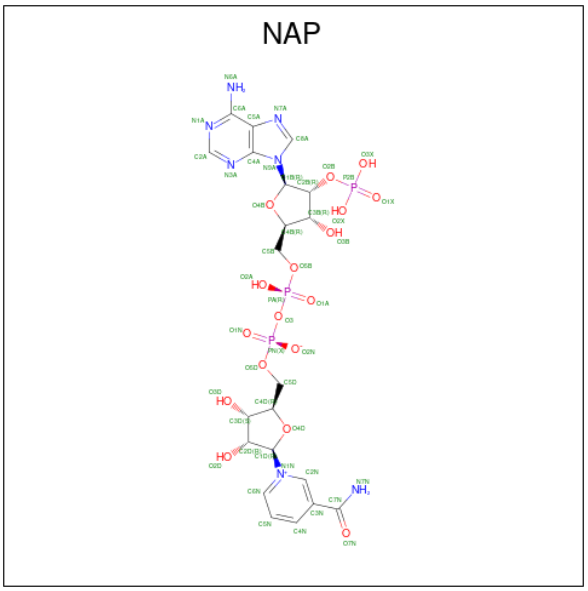
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	530	4413	2857	725	804	27	0	0	0
1	B	543	4514	2914	745	829	26	0	0	0

- Molecule 2 is DIHYDROFOLIC ACID (three-letter code: DHF) (formula: $C_{19}H_{21}N_7O_6$).



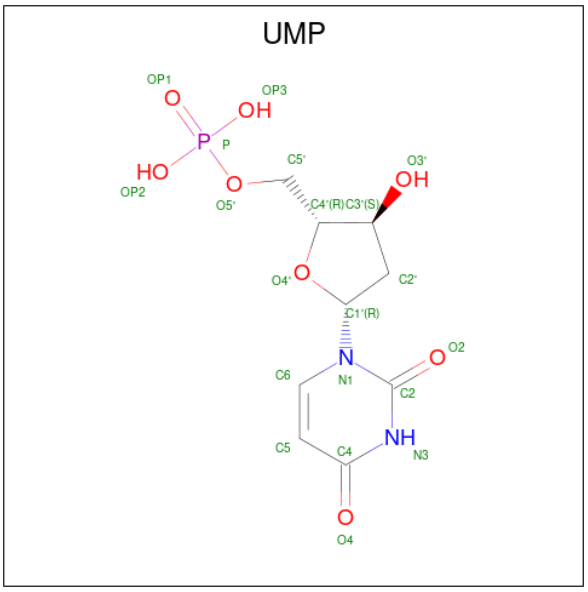
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
2	A	1	32	19	7	6	0	0

- Molecule 3 is NADP NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (three-letter code: NAP) (formula: $C_{21}H_{28}N_7O_{17}P_3$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
3	A	1	48	21	7	17	3	0	0

- Molecule 4 is 2'-DEOXYURIDINE 5'-MONOPHOSPHATE (three-letter code: UMP) (formula: $C_9H_{13}N_2O_8P$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
4	A	1	20	9	2	8	1	0	0
			Total	C	N	O	P		
4	B	1	20	9	2	8	1	0	0

- Molecule 5 is water.

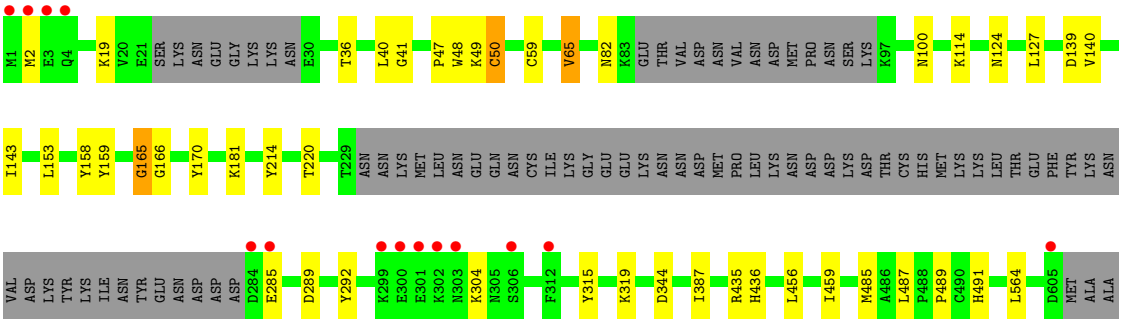
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
5	A	222	222	222	0	0
			Total	O		
5	B	185	185	185	0	0

3 Residue-property plots (i)

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

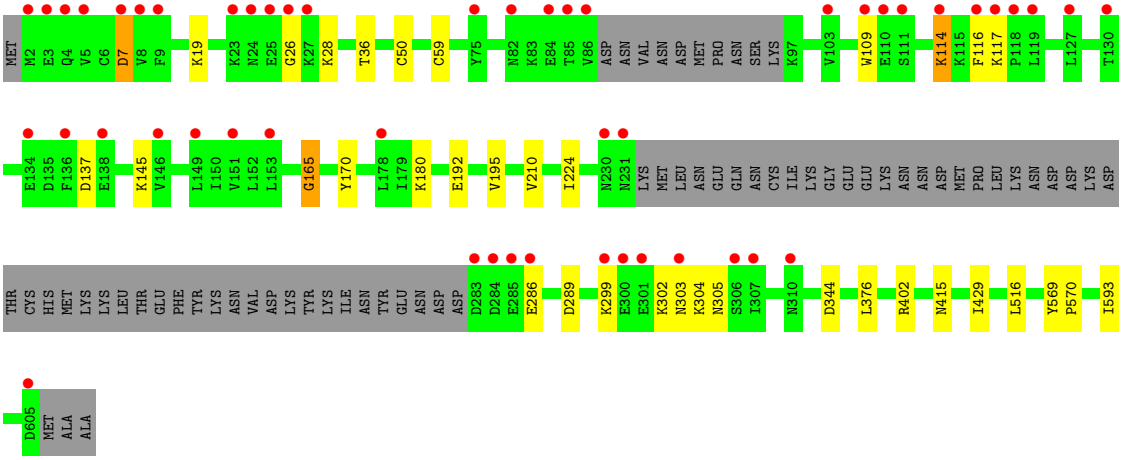
• Molecule 1: Bifunctional dihydrofolate reductase-thymidylate synthase

Chain A:



• Molecule 1: Bifunctional dihydrofolate reductase-thymidylate synthase

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	58.06Å 156.78Å 165.03Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.82 – 2.50 49.82 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.7 (49.82-2.50) 99.6 (49.82-2.50)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.09 (at 2.48Å)	Xtriage
Refinement program	CNS, REFMAC 5.6.0117	Depositor
R, R_{free}	0.212 , 0.266 0.199 , 0.245	Depositor DCC
R_{free} test set	2112 reflections (3.97%)	wwPDB-VP
Wilson B-factor (Å^2)	36.8	Xtriage
Anisotropy	0.553	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.33 , 36.0	EDS
L-test for twinning[2]	$< \|L\| > = 0.52$, $< L^2 > = 0.36$	Xtriage
Estimated twinning fraction	0.002 for -h,l,k	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	9454	wwPDB-VP
Average B, all atoms (Å^2)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.57% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

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

5 Model quality (i)

5.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: UMP, NAP, DHF

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	0.54	0/4517	0.65	0/6099
1	B	0.52	0/4619	0.63	0/6237
All	All	0.53	0/9136	0.64	0/12336

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand.A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2
1	B	0	1
All	All	0	3

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	165	GLY	Peptide
1	A	304	LYS	Peptide
1	B	165	GLY	Peptide

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	4413	0	4370	24	0
1	B	4514	0	4462	15	0
2	A	32	0	19	1	0
3	A	48	0	25	4	0
4	A	20	0	11	1	0
4	B	20	0	11	0	0
5	A	222	0	0	2	0
5	B	185	0	0	3	0
All	All	9454	0	8898	38	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 2.

All (38) 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:A:50:CYS:SG	5:A:975:HOH:O	2.30	0.88
1:B:59:CYS:SG	5:B:984:HOH:O	2.34	0.85
1:A:59:CYS:SG	5:A:931:HOH:O	2.16	0.66
1:B:210:VAL:HG12	1:B:224:ILE:HG22	1.80	0.64
1:A:40:LEU:O	3:A:702:NAP:H2N	2.03	0.58
1:A:166:GLY:HA3	3:A:702:NAP:PA	2.44	0.57
1:A:127:LEU:HD23	1:A:143:ILE:HG13	1.88	0.54
1:A:19:LYS:HG2	1:A:36:THR:HG22	1.94	0.50
1:B:303:ASN:O	1:B:305:ASN:N	2.47	0.47
1:B:19:LYS:HG2	1:B:36:THR:HG22	1.96	0.46
1:A:181:LYS:NZ	1:B:289:ASP:OD2	2.46	0.45
1:A:315:TYR:HB2	1:A:564:LEU:O	2.16	0.45
1:A:319:LYS:HD3	1:B:286:GLU:HG3	1.98	0.45
1:A:456:LEU:O	1:A:459:ILE:HG13	2.17	0.45
1:B:192:GLU:O	5:B:850:HOH:O	2.21	0.44
1:B:109:TRP:CE2	1:B:117:LYS:HD2	2.53	0.44
1:B:415:ASN:ND2	5:B:829:HOH:O	2.51	0.43
1:B:114:LYS:HA	1:B:117:LYS:HB3	2.00	0.43
1:A:214:TYR:O	1:A:220:THR:HA	2.19	0.42
1:A:100:ASN:OD1	1:A:159:TYR:HB3	2.20	0.42
1:A:153:LEU:HD22	1:A:158:TYR:CZ	2.55	0.42
2:A:701:DHF:H72	3:A:702:NAP:C5N	2.50	0.42
1:B:165:GLY:HA3	1:B:170:TYR:CZ	2.55	0.42
1:A:166:GLY:HA3	3:A:702:NAP:O1A	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:41:GLY:HA2	1:A:47:PRO:HD3	2.02	0.41
1:B:569:TYR:HB3	1:B:570:PRO:HD2	2.01	0.41
1:A:491:HIS:CE1	4:A:703:UMP:O4	2.73	0.41
1:A:165:GLY:HA3	1:A:170:TYR:CZ	2.55	0.41
1:A:435:ARG:HB2	1:A:436:HIS:CD2	2.56	0.41
1:A:65:VAL:HG22	1:A:159:TYR:CB	2.51	0.41
1:A:124:ASN:HB2	1:A:140:VAL:HG12	2.02	0.41
1:A:387:ILE:O	1:A:435:ARG:NH1	2.50	0.41
1:B:165:GLY:CA	1:B:170:TYR:CZ	3.04	0.41
1:B:376:LEU:HD12	1:B:593:ILE:HG13	2.03	0.40
1:A:48:TRP:O	1:A:49:LYS:HB2	2.22	0.40
1:A:485:MET:SD	1:A:489:PRO:HD3	2.62	0.40
1:B:7:ASP:HA	1:B:180:LYS:HE2	2.03	0.40
1:A:289:ASP:HA	1:A:292:TYR:CD2	2.57	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	522/608 (86%)	495 (95%)	25 (5%)	2 (0%)	34	54
1	B	537/608 (88%)	506 (94%)	28 (5%)	3 (1%)	25	43
All	All	1059/1216 (87%)	1001 (94%)	53 (5%)	5 (0%)	29	48

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	304	LYS
1	A	139	ASP
1	B	26	GLY
1	A	82	ASN

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Mol	Chain	Res	Type
1	B	429	ILE

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	496/570 (87%)	489 (99%)	7 (1%)	67	86
1	B	508/570 (89%)	495 (97%)	13 (3%)	46	72
All	All	1004/1140 (88%)	984 (98%)	20 (2%)	55	79

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	MET
1	A	50	CYS
1	A	65	VAL
1	A	114	LYS
1	A	285	GLU
1	A	344	ASP
1	A	487	LEU
1	B	7	ASP
1	B	28	LYS
1	B	50	CYS
1	B	114	LYS
1	B	116	PHE
1	B	137	ASP
1	B	145	LYS
1	B	195	VAL
1	B	299	LYS
1	B	302	LYS
1	B	344	ASP
1	B	402	ARG
1	B	516	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

4 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	DHF	A	701	-	30,34,34	1.04	2 (6%)	36,47,47	1.58	4 (11%)
4	UMP	B	701	-	21,21,21	1.35	2 (9%)	30,31,31	1.96	9 (30%)
3	NAP	A	702	-	46,52,52	0.83	0	61,80,80	1.21	5 (8%)
4	UMP	A	703	-	21,21,21	1.34	4 (19%)	30,31,31	1.96	9 (30%)

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	DHF	A	701	-	-	2/20/31/31	0/3/3/3
4	UMP	B	701	-	-	4/10/22/22	0/2/2/2
3	NAP	A	702	-	-	3/31/67/67	0/5/5/5
4	UMP	A	703	-	-	3/10/22/22	0/2/2/2

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	703	UMP	C2-N1	3.40	1.43	1.38
4	B	701	UMP	C2-N1	3.06	1.43	1.38
2	A	701	DHF	C4A-C8A	2.71	1.48	1.41
4	A	703	UMP	C2-N3	-2.35	1.33	1.38
4	A	703	UMP	C5-C4	-2.16	1.39	1.43
4	B	701	UMP	C4-N3	-2.15	1.34	1.38
2	A	701	DHF	C7-C6	2.09	1.52	1.49
4	A	703	UMP	C4-N3	-2.09	1.35	1.38

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	701	DHF	C2-N3-C4	4.85	122.71	115.96
2	A	701	DHF	C4A-C4-N3	-4.38	117.56	123.42
4	A	703	UMP	C5-C4-N3	4.37	120.92	114.80
4	A	703	UMP	O4-C4-C5	-3.82	118.57	125.16
4	B	701	UMP	C4-N3-C2	-3.73	121.98	126.61
4	A	703	UMP	C4-N3-C2	-3.58	122.17	126.61
4	B	701	UMP	N3-C2-N1	3.58	119.55	114.89
4	B	701	UMP	O4-C4-C5	-3.54	119.06	125.16
4	B	701	UMP	C1'-N1-C2	3.51	124.52	117.66
3	A	702	NAP	C3N-C7N-N7N	3.48	122.02	117.74
4	B	701	UMP	C5-C4-N3	3.43	119.60	114.80
4	A	703	UMP	C1'-N1-C2	3.26	124.03	117.66
3	A	702	NAP	N3A-C2A-N1A	-3.16	124.38	128.67
4	A	703	UMP	N3-C2-N1	2.95	118.74	114.89
2	A	701	DHF	C2-N1-C8A	2.92	121.54	114.59
2	A	701	DHF	C8A-C4A-C4	2.67	116.26	114.53
4	B	701	UMP	C2'-C1'-N1	2.63	120.39	113.81
4	A	703	UMP	O2-C2-N3	-2.63	116.64	121.49
3	A	702	NAP	O2A-PA-O1A	2.55	124.31	112.44
4	B	701	UMP	O2-C2-N3	-2.54	116.81	121.49
4	A	703	UMP	C2'-C1'-N1	2.47	119.98	113.81
4	B	701	UMP	C1'-N1-C6	-2.42	116.77	121.53
3	A	702	NAP	C4A-C5A-N7A	-2.40	106.80	109.34
3	A	702	NAP	O7N-C7N-N7N	-2.38	119.18	122.62
4	A	703	UMP	C1'-N1-C6	-2.28	117.06	121.53
4	A	703	UMP	OP3-P-OP2	2.27	116.31	107.80
4	B	701	UMP	O4'-C4'-C3'	-2.24	100.54	105.65

There are no chirality outliers.

All (12) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	701	DHF	CT-CA-CB-CG
3	A	702	NAP	O4D-C1D-N1N-C6N
4	A	703	UMP	C5'-O5'-P-OP2
4	A	703	UMP	C5'-O5'-P-OP3
2	A	701	DHF	N-CA-CB-CG
4	B	701	UMP	C3'-C4'-C5'-O5'
4	B	701	UMP	O4'-C4'-C5'-O5'
4	A	703	UMP	C5'-O5'-P-OP1
4	B	701	UMP	C5'-O5'-P-OP1
3	A	702	NAP	C2B-O2B-P2B-O2X
4	B	701	UMP	C4'-C5'-O5'-P
3	A	702	NAP	O4D-C1D-N1N-C2N

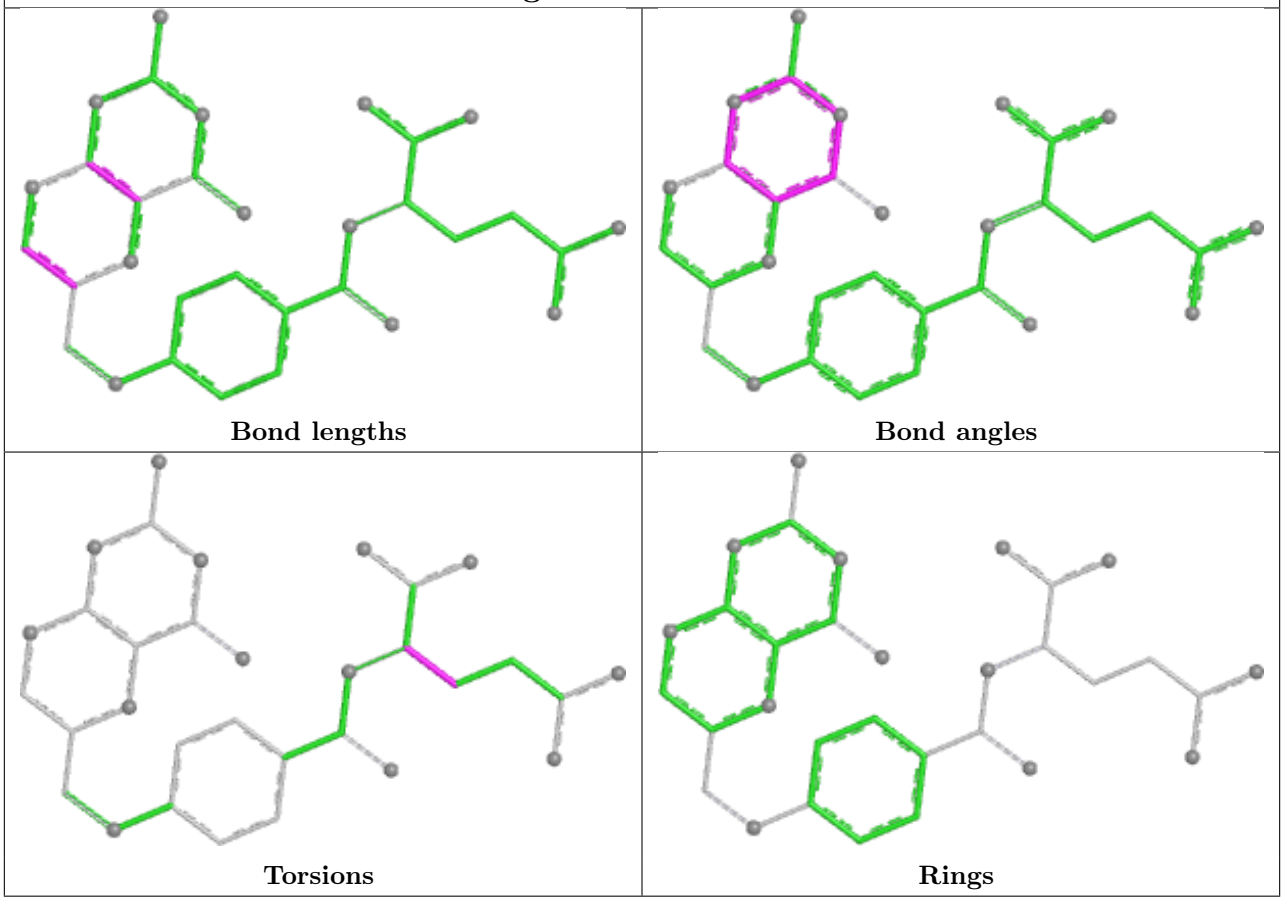
There are no ring outliers.

3 monomers are involved in 5 short contacts:

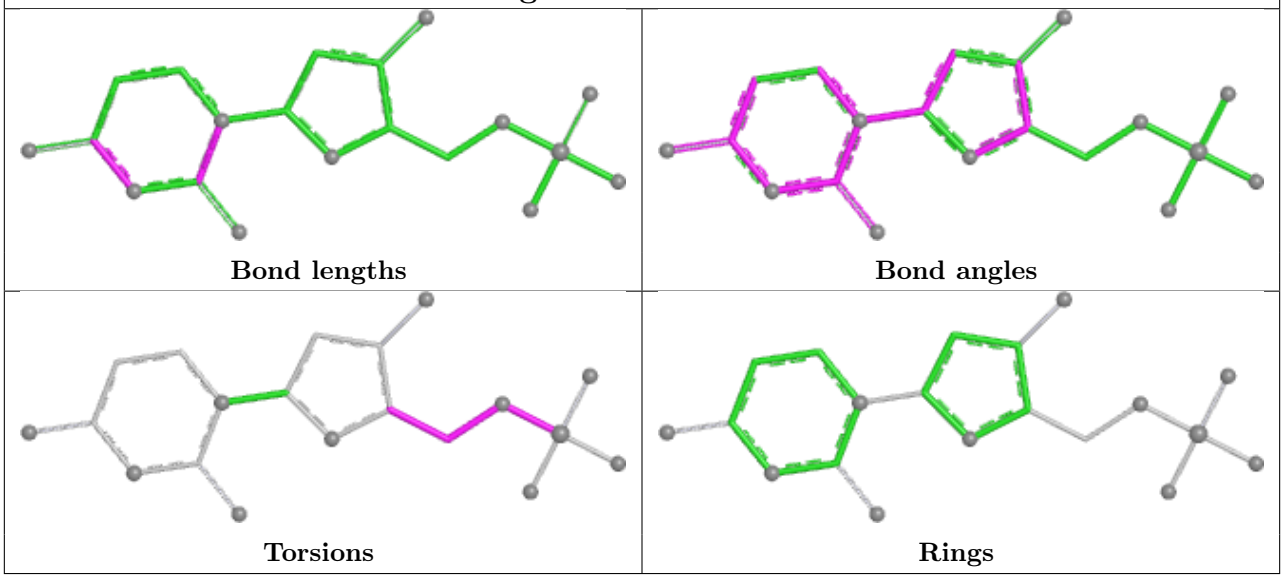
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	701	DHF	1	0
3	A	702	NAP	4	0
4	A	703	UMP	1	0

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

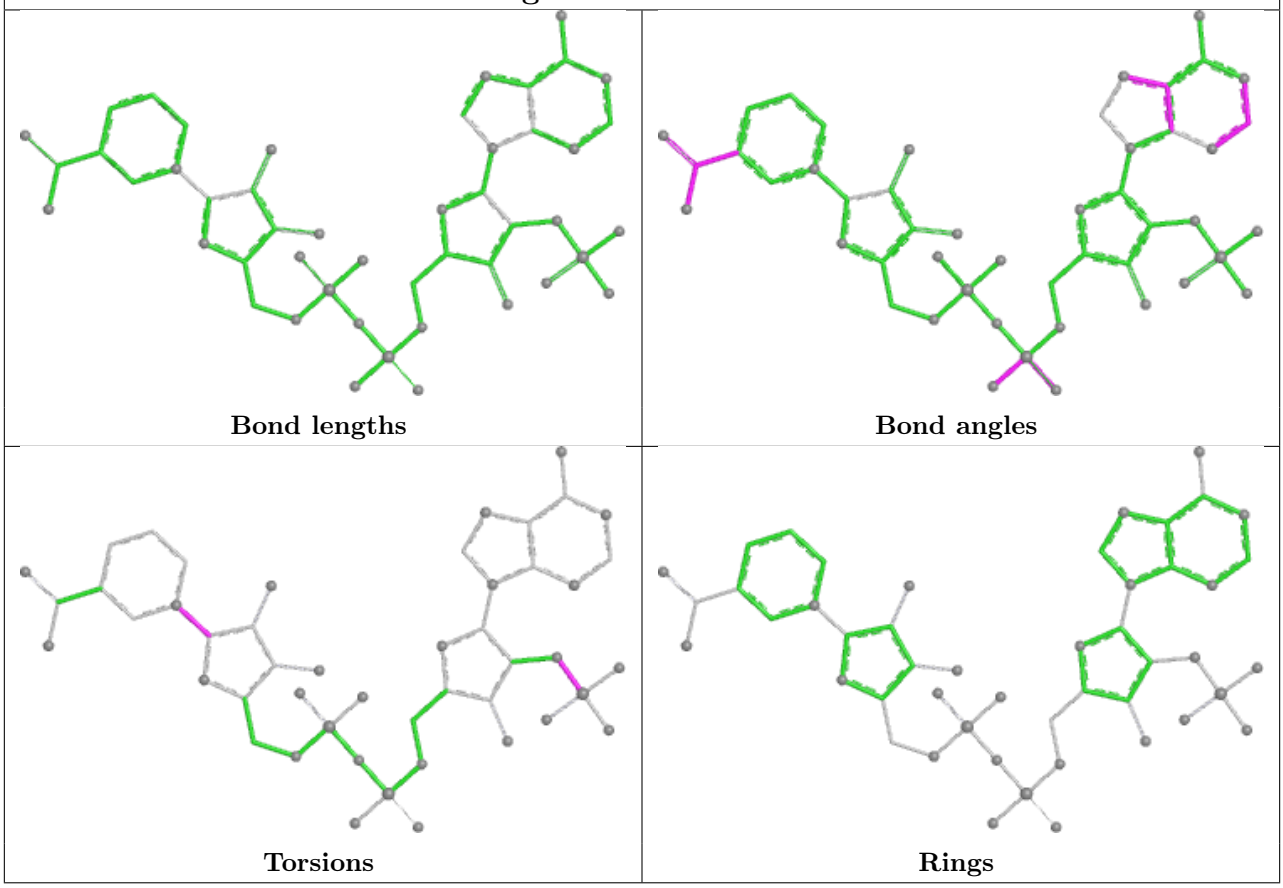
Ligand DHF A 701



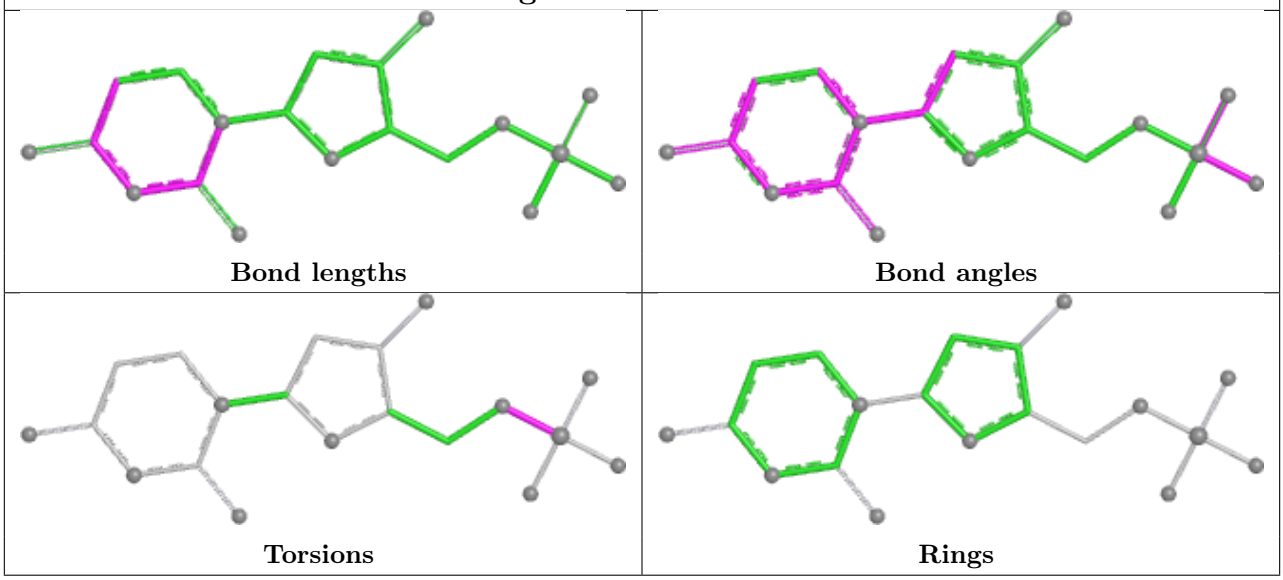
Ligand UMP B 701



Ligand NAP A 702



Ligand UMP A 703



5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

6 Fit of model and data (i)

6.1 Protein, DNA and RNA chains (i)

In the following table, the column labelled '#RSRZ> 2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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	530/608 (87%)	-0.18	14 (2%)	56	59	20, 31, 72, 90	0
1	B	543/608 (89%)	0.17	50 (9%)	9	9	20, 38, 90, 90	0
All	All	1073/1216 (88%)	-0.00	64 (5%)	21	22	20, 34, 89, 90	0

All (64) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1	MET	7.9
1	B	2	MET	7.4
1	A	2	MET	7.0
1	B	85	THR	6.9
1	B	116	PHE	6.2
1	B	231	ASN	5.7
1	A	302	LYS	5.3
1	A	299	LYS	5.1
1	B	24	ASN	5.1
1	A	301	GLU	4.9
1	A	300	GLU	4.7
1	B	75	TYR	4.7
1	B	4	GLN	4.7
1	B	27	LYS	4.6
1	B	299	LYS	4.6
1	B	151	VAL	4.5
1	B	138	GLU	4.5
1	B	3	GLU	4.4
1	B	118	PRO	4.1
1	B	23	LYS	3.8
1	B	300	GLU	3.7
1	B	86	VAL	3.7
1	B	136	PHE	3.7
1	B	301	GLU	3.6

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Mol	Chain	Res	Type	RSRZ
1	B	9	PHE	3.6
1	B	605	ASP	3.5
1	B	230	ASN	3.5
1	B	306	SER	3.4
1	B	26	GLY	3.4
1	B	8	VAL	3.4
1	B	119	LEU	3.4
1	B	127	LEU	3.3
1	B	117	LYS	3.1
1	B	7	ASP	3.0
1	A	3	GLU	2.9
1	B	114	LYS	2.8
1	B	25	GLU	2.7
1	B	178	LEU	2.7
1	B	307	ILE	2.6
1	B	5	VAL	2.6
1	B	284	ASP	2.5
1	A	312	PHE	2.5
1	A	306	SER	2.5
1	B	82	ASN	2.4
1	B	286	GLU	2.4
1	A	303	ASN	2.4
1	B	103	VAL	2.4
1	B	109	TRP	2.4
1	A	4	GLN	2.3
1	B	285	GLU	2.3
1	B	153	LEU	2.3
1	A	284	ASP	2.2
1	A	605	ASP	2.2
1	B	303	ASN	2.2
1	A	285	GLU	2.2
1	B	134	GLU	2.2
1	B	110	GLU	2.2
1	B	310	ASN	2.1
1	B	149	LEU	2.1
1	B	130	THR	2.1
1	B	111	SER	2.1
1	B	84	GLU	2.0
1	B	283	ASP	2.0
1	B	146	VAL	2.0

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

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

6.3 Carbohydrates (i)

There are no monosaccharides in this entry.

6.4 Ligands (i)

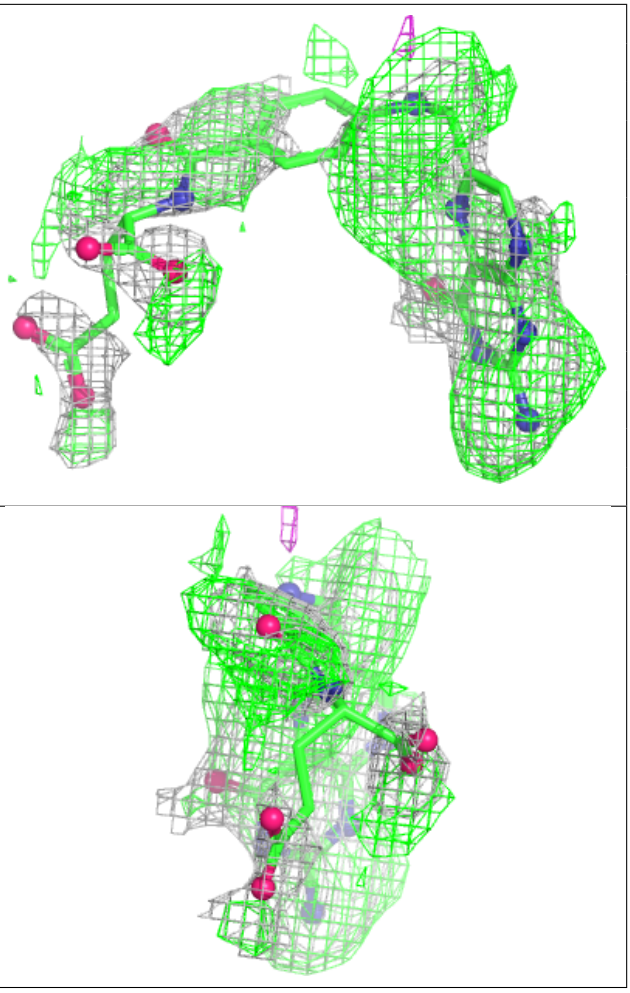
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	DHF	A	701	32/32	0.59	0.58	25,34,36,37	32
4	UMP	A	703	20/20	0.97	0.13	29,36,41,42	0
4	UMP	B	701	20/20	0.97	0.17	28,39,50,51	0
3	NAP	A	702	48/48	0.98	0.10	24,33,41,45	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.

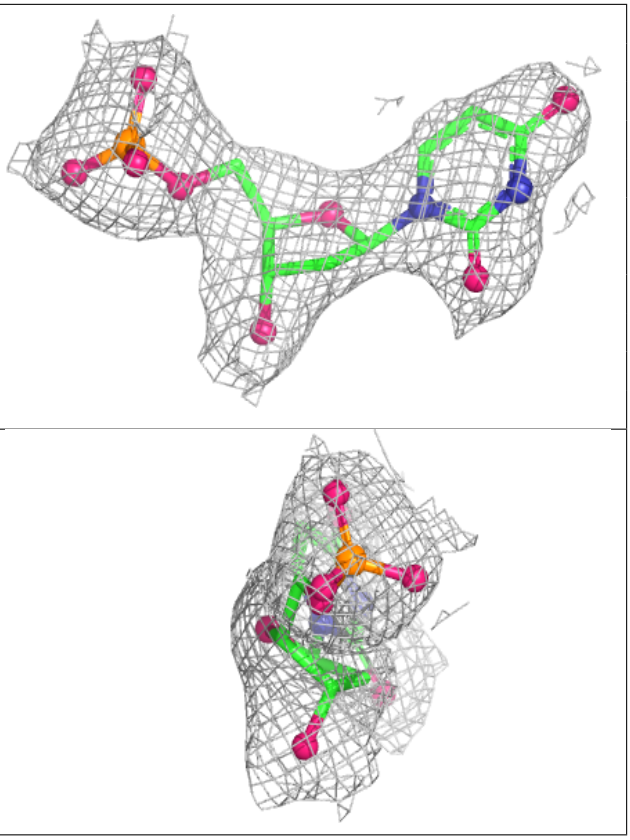
Electron density around DHF A 701:

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



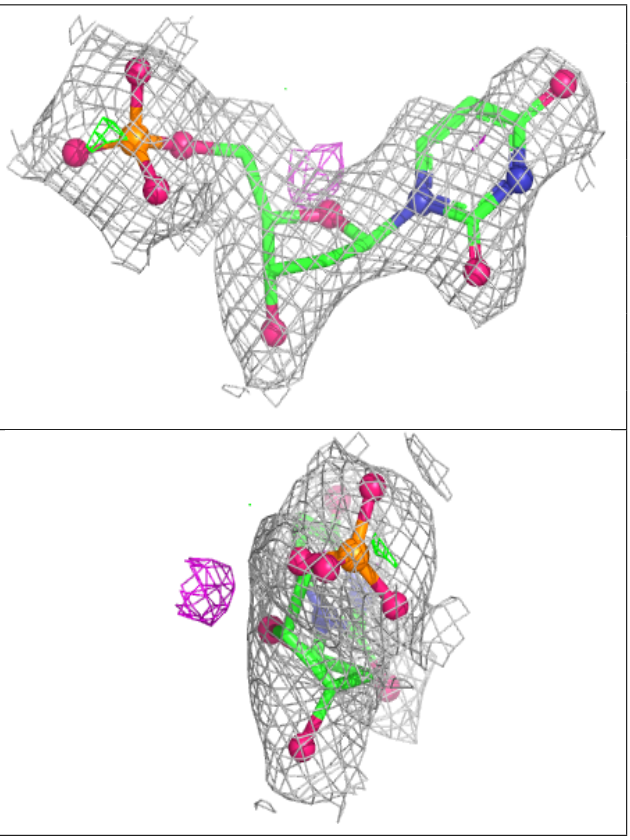
Electron density around UMP A 703:

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



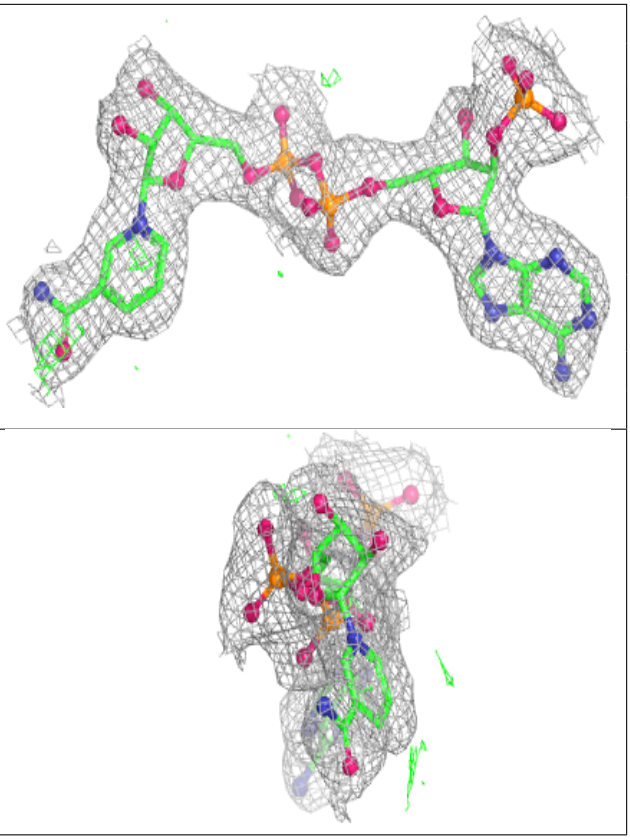
Electron density around UMP B 701:

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



Electron density around NAP A 702:

$2mF_o$-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.