



Full wwPDB X-ray Structure Validation Report ⓘ

May 21, 2020 – 01:19 am BST

PDB ID : 6OJV
Title : Crystal structure of human thymidylate synthase delta(7-29) in complex with dUMP and 2-amino-4-oxo-4,7-dihydro-pyrrolo[2,3-d]pyrimidine-methyl-phenyl-L-glutamic acid
Authors : Czyzyk, D.J.; Anderson, K.S.; Valhondo, M.; Jorgensen, W.L.
Deposited on : 2019-04-12
Resolution : 2.59 Å(reported)

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

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

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

with specific help available everywhere you see the ⓘ 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

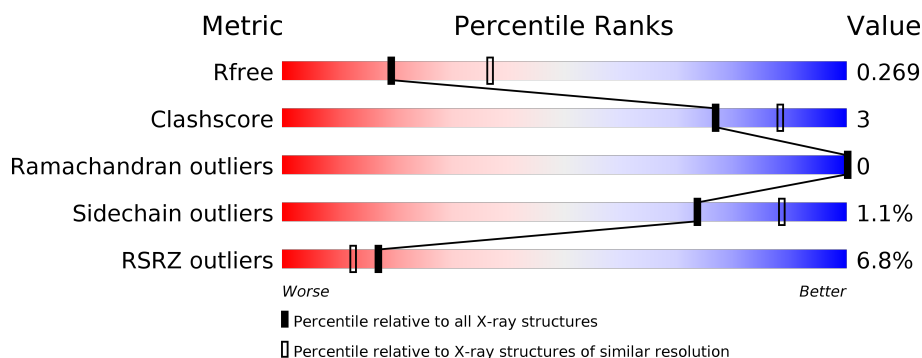
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.59 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	3163 (2.60-2.60)
Clashscore	141614	3518 (2.60-2.60)
Ramachandran outliers	138981	3455 (2.60-2.60)
Sidechain outliers	138945	3455 (2.60-2.60)
RSRZ outliers	127900	3104 (2.60-2.60)

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	290	<div> <div>5%</div> <div>88%</div> <div>10%</div> <div>•</div> </div>
1	B	290	<div> <div>5%</div> <div>92%</div> <div>6%</div> <div>•</div> </div>
1	C	290	<div> <div>5%</div> <div>89%</div> <div>9%</div> <div>•</div> </div>
1	D	290	<div> <div>11%</div> <div>91%</div> <div>7%</div> <div>••</div> </div>

2 Entry composition [i](#)

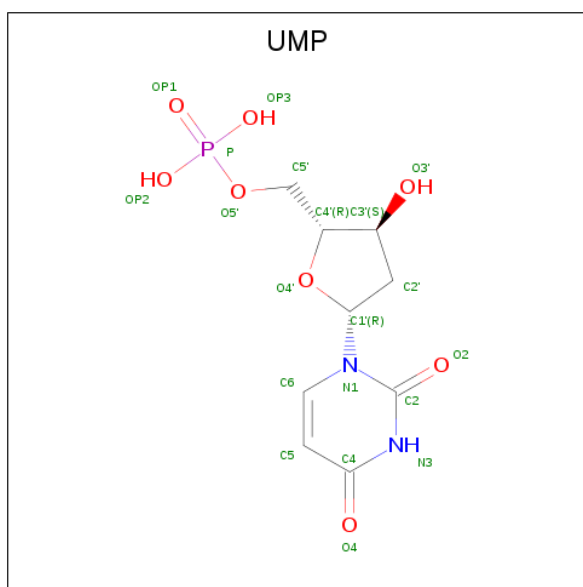
There are 4 unique types of molecules in this entry. The entry contains 9190 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Thymidylate synthase,Thymidylate synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	285	Total	C	N	O	S	0	0	0
			2262	1445	388	416	13			
1	A	285	Total	C	N	O	S	0	0	0
			2247	1435	385	414	13			
1	C	285	Total	C	N	O	S	0	0	0
			2225	1427	380	405	13			
1	D	285	Total	C	N	O	S	0	0	0
			2223	1422	376	412	13			

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



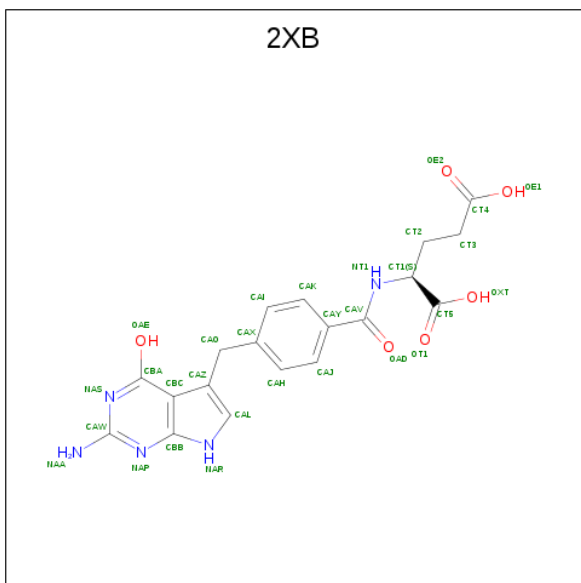
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	B	1	Total	C	N	O	P	0	0
			20	9	2	8	1		
2	A	1	Total	C	N	O	P	0	0
			20	9	2	8	1		

Continued on next page...

Continued from previous page...

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

- Molecule 3 is N-{4-[(2-amino-4-hydroxy-7H-pyrrolo[2,3-d]pyrimidin-5-yl)methyl]benzoyl}-L-glutamic acid (three-letter code: 2XB) (formula: C₁₉H₁₉N₅O₆).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	B	1	Total 30	C 19	N 5	O 6	0	0
3	A	1	Total 30	C 19	N 5	O 6	0	0
3	C	1	Total 30	C 19	N 5	O 6	0	0
3	D	1	Total 30	C 19	N 5	O 6	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	B	13	Total O 13 13	0	0
4	A	6	Total O 6 6	0	0
4	C	7	Total O 7 7	0	0

Continued on next page...

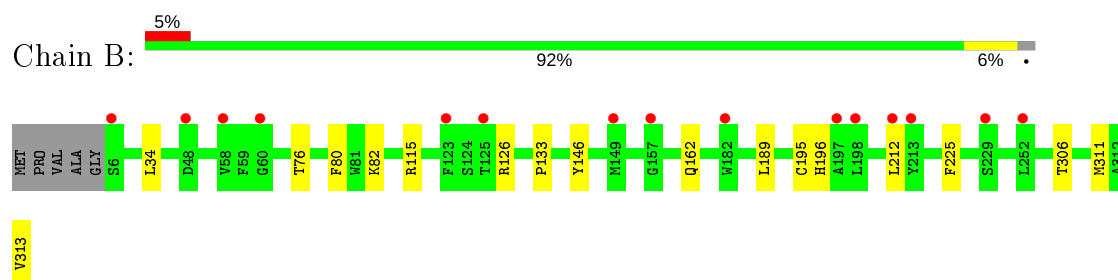
Continued from previous page...

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	D	7	Total	O	0	0
			7	7		

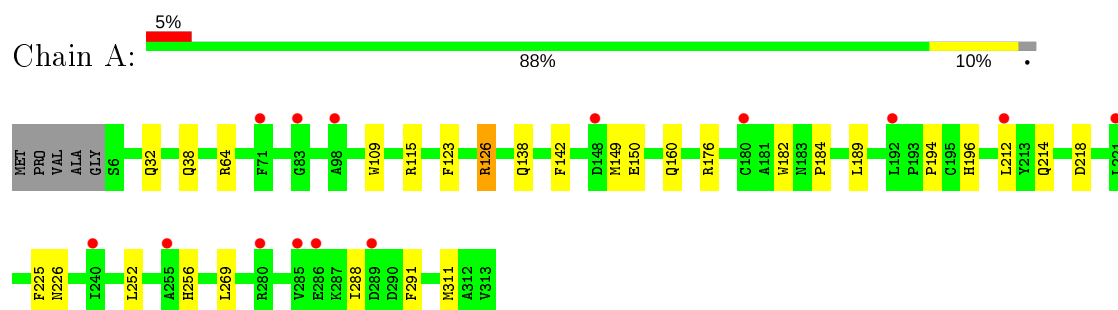
3 Residue-property plots [i](#)

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.

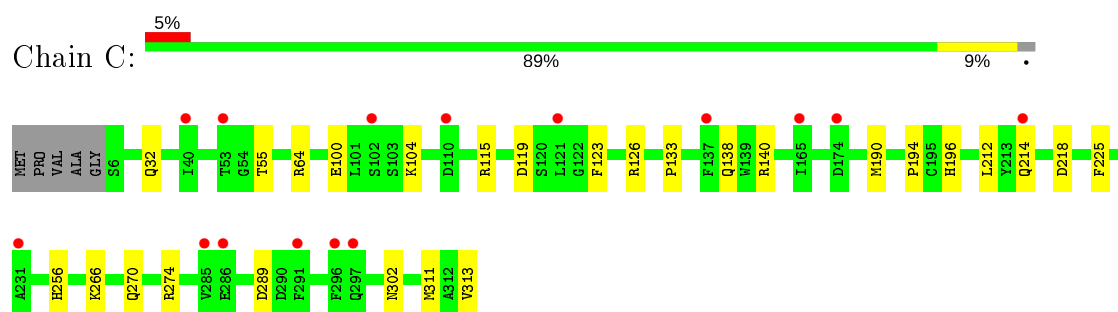
- Molecule 1: Thymidylate synthase,Thymidylate synthase



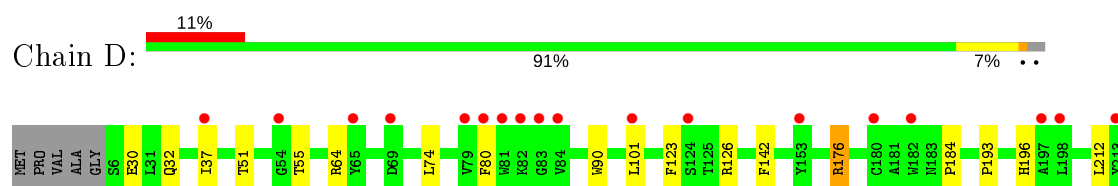
- Molecule 1: Thymidylate synthase,Thymidylate synthase

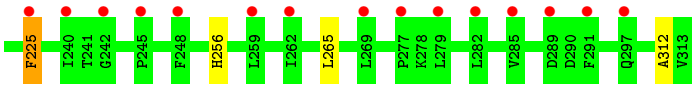


- Molecule 1: Thymidylate synthase,Thymidylate synthase



- Molecule 1: Thymidylate synthase,Thymidylate synthase





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	154.95Å 103.03Å 109.21Å 90.00° 128.92° 90.00°	Depositor
Resolution (Å)	48.56 – 2.59 48.55 – 2.59	Depositor EDS
% Data completeness (in resolution range)	98.9 (48.56-2.59) 99.2 (48.55-2.59)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.14	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.99 (at 2.58Å)	Xtriage
Refinement program	PHENIX (1.10.1_2155)	Depositor
R, R_{free}	0.240 , 0.269 0.240 , 0.269	Depositor DCC
R_{free} test set	2058 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	64.5	Xtriage
Anisotropy	0.582	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 41.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	9190	wwPDB-VP
Average B, all atoms (Å ²)	67.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.40% 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: 2XB, UMP

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	0.25	0/2304	0.41	0/3130
1	B	0.25	0/2319	0.42	0/3146
1	C	0.26	0/2282	0.43	0/3101
1	D	0.24	0/2280	0.41	0/3099
All	All	0.25	0/9185	0.42	0/12476

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	2247	0	2150	18	0
1	B	2262	0	2185	9	0
1	C	2225	0	2132	15	0
1	D	2223	0	2109	14	0
2	A	20	0	11	3	0
2	B	20	0	11	1	0
2	C	20	0	11	1	0
2	D	20	0	11	0	0
3	A	30	0	16	1	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	30	0	16	1	0
3	C	30	0	16	1	0
3	D	30	0	16	0	0
4	A	6	0	0	0	0
4	B	13	0	0	1	0
4	C	7	0	0	0	0
4	D	7	0	0	0	0
All	All	9190	0	8684	53	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 3.

All (53) 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)
2:A:600:UMP:OP1	1:D:176:ARG:NH1	2.21	0.73
1:D:123:PHE:HB3	1:D:126:ARG:HG3	1.71	0.72
1:B:115:ARG:NH1	1:B:126:ARG:O	2.29	0.66
1:A:126:ARG:NH2	1:A:189:LEU:O	2.31	0.62
1:B:196:HIS:HB2	1:B:212:LEU:HD11	1.83	0.60
1:A:218:ASP:HB2	2:A:600:UMP:H1'	1.85	0.59
1:B:162:GLN:NE2	4:B:702:HOH:O	2.34	0.59
1:C:140:ARG:NH2	1:C:289:ASP:OD1	2.35	0.59
1:A:196:HIS:HB2	1:A:212:LEU:HD11	1.85	0.58
1:A:123:PHE:HB3	1:A:126:ARG:HG3	1.85	0.57
1:A:32:GLN:NE2	1:A:64:ARG:O	2.39	0.56
1:D:37:ILE:HD13	1:D:265:LEU:HD22	1.88	0.55
1:C:196:HIS:HB2	1:C:212:LEU:HD11	1.90	0.54
1:C:32:GLN:NE2	1:C:64:ARG:O	2.37	0.52
1:D:196:HIS:HB2	1:D:212:LEU:HD11	1.91	0.51
1:B:133:PRO:HG3	1:B:146:TYR:CG	2.46	0.51
1:D:80:PHE:HB3	1:D:225:PHE:HE2	1.76	0.51
1:B:311:MET:HE2	3:B:601:2XB:H7	1.93	0.50
1:A:115:ARG:NH2	1:A:126:ARG:O	2.44	0.50
1:C:190:MET:SD	1:C:194:PRO:HD3	2.51	0.49
1:A:214:GLN:NE2	1:A:226:ASN:OD1	2.45	0.49
1:C:212:LEU:HD21	1:C:214:GLN:HB2	1.95	0.48
1:C:311:MET:CE	3:C:601:2XB:H8	2.43	0.48
1:B:34:LEU:HD11	1:B:76:THR:HG21	1.97	0.47
1:C:266:LYS:O	1:C:270:GLN:HG2	2.15	0.47
1:C:218:ASP:HB2	2:C:600:UMP:H1'	1.95	0.47

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:176:ARG:HD2	1:D:193:PRO:HG2	1.97	0.46
1:B:126:ARG:NH2	1:B:189:LEU:O	2.42	0.46
1:D:55:THR:HB	1:D:256:HIS:HB2	1.97	0.46
1:C:115:ARG:NH1	1:C:119:ASP:OD2	2.49	0.46
1:D:80:PHE:HB3	1:D:225:PHE:CE2	2.51	0.46
1:A:184:PRO:HD2	1:D:142:PHE:CE1	2.50	0.46
1:D:32:GLN:NE2	1:D:64:ARG:O	2.45	0.46
1:A:214:GLN:HB3	1:A:252:LEU:HD23	1.98	0.45
1:B:80:PHE:CE1	1:B:82:LYS:HB3	2.52	0.45
1:C:100:GLU:OE2	1:C:104:LYS:NZ	2.50	0.45
1:A:256:HIS:CE1	2:A:600:UMP:HO3'	2.28	0.44
1:C:196:HIS:CB	1:C:212:LEU:HD11	2.47	0.43
1:A:142:PHE:CE1	1:D:184:PRO:HD2	2.54	0.43
1:C:55:THR:HB	1:C:256:HIS:HB2	2.01	0.42
1:D:51:THR:HG21	1:D:312:ALA:HB1	2.01	0.42
1:D:90:TRP:CE3	1:D:101:LEU:HD22	2.55	0.42
1:C:123:PHE:HB3	1:C:126:ARG:HG3	2.01	0.42
1:C:274:ARG:HD2	1:C:302:ASN:O	2.19	0.42
1:A:182:TRP:HE3	1:A:194:PRO:HG2	1.85	0.41
1:A:288:ILE:HD12	1:A:291:PHE:CD2	2.56	0.41
1:D:30:GLU:HG3	1:D:74:LEU:HD22	2.02	0.41
1:A:38:GLN:HB2	1:A:269:LEU:HD21	2.02	0.41
1:C:133:PRO:HB2	1:C:138:GLN:NE2	2.36	0.41
1:A:109:TRP:CZ3	3:A:601:2XB:H17	2.56	0.41
1:B:195:CYS:SG	2:B:600:UMP:C6	3.14	0.40
1:A:138:GLN:O	1:A:160:GLN:NE2	2.50	0.40
1:A:149:MET:HG3	1:A:150:GLU:HG3	2.03	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	283/290 (98%)	272 (96%)	11 (4%)	0	100	100
1	B	283/290 (98%)	272 (96%)	11 (4%)	0	100	100
1	C	283/290 (98%)	271 (96%)	12 (4%)	0	100	100
1	D	283/290 (98%)	270 (95%)	13 (5%)	0	100	100
All	All	1132/1160 (98%)	1085 (96%)	47 (4%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	236/252 (94%)	233 (99%)	3 (1%)	69	86
1	B	240/252 (95%)	237 (99%)	3 (1%)	69	86
1	C	231/252 (92%)	229 (99%)	2 (1%)	78	91
1	D	231/252 (92%)	229 (99%)	2 (1%)	78	91
All	All	938/1008 (93%)	928 (99%)	10 (1%)	73	88

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	225	PHE
1	B	306	THR
1	B	313	VAL
1	A	126	ARG
1	A	225	PHE
1	A	311	MET
1	C	225	PHE
1	C	313	VAL
1	D	176	ARG
1	D	225	PHE

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

Mol	Chain	Res	Type
1	B	196	HIS
1	C	196	HIS
1	D	196	HIS
1	D	297	GLN

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	UMP	A	600	-	18,21,21	1.47	2 (11%)	21,31,31	1.46	2 (9%)
3	2XB	C	601	-	26,32,32	1.50	4 (15%)	27,45,45	1.71	7 (25%)
3	2XB	A	601	-	26,32,32	1.45	4 (15%)	27,45,45	1.79	7 (25%)
3	2XB	B	601	-	26,32,32	1.48	4 (15%)	27,45,45	1.81	7 (25%)
2	UMP	C	600	-	18,21,21	1.72	3 (16%)	21,31,31	1.62	4 (19%)
3	2XB	D	601	-	26,32,32	1.50	4 (15%)	27,45,45	1.85	7 (25%)
2	UMP	B	600	-	18,21,21	1.52	3 (16%)	21,31,31	1.56	3 (14%)
2	UMP	D	600	-	18,21,21	1.49	3 (16%)	21,31,31	1.52	2 (9%)

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	UMP	A	600	-	-	2/7/22/22	0/2/2/2
3	2XB	C	601	-	-	2/15/21/21	0/3/3/3
3	2XB	A	601	-	-	3/15/21/21	0/3/3/3
3	2XB	B	601	-	-	2/15/21/21	0/3/3/3
2	UMP	C	600	-	-	2/7/22/22	0/2/2/2
3	2XB	D	601	-	-	5/15/21/21	0/3/3/3
2	UMP	B	600	-	-	1/7/22/22	0/2/2/2
2	UMP	D	600	-	-	3/7/22/22	0/2/2/2

All (27) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	600	UMP	C4-N3	4.17	1.40	1.33
2	B	600	UMP	C4-N3	4.03	1.40	1.33
2	A	600	UMP	C4-N3	4.01	1.40	1.33
2	D	600	UMP	C4-N3	4.01	1.40	1.33
3	B	601	2XB	CBA-CBC	3.87	1.48	1.43
2	C	600	UMP	P-OP1	3.78	1.62	1.50
3	C	601	2XB	CBA-CBC	3.74	1.48	1.43
3	D	601	2XB	CBA-CBC	3.71	1.48	1.43
3	A	601	2XB	CBA-CBC	3.47	1.48	1.43
3	D	601	2XB	CAW-NAS	3.06	1.40	1.35
3	B	601	2XB	CAW-NAS	3.05	1.40	1.35
3	C	601	2XB	CAW-NAS	2.99	1.40	1.35
2	B	600	UMP	C6-N1	2.94	1.39	1.35
3	A	601	2XB	CAW-NAS	2.93	1.40	1.35
2	D	600	UMP	C6-N1	2.79	1.39	1.35
2	C	600	UMP	C6-N1	2.72	1.39	1.35
2	A	600	UMP	C6-N1	2.70	1.39	1.35
3	D	601	2XB	CAW-NAA	2.26	1.38	1.33
3	A	601	2XB	CAW-NAA	2.24	1.38	1.33
3	C	601	2XB	CAW-NAA	2.21	1.38	1.33
3	A	601	2XB	CBC-CBB	-2.18	1.37	1.43
3	B	601	2XB	CAW-NAA	2.15	1.38	1.33
3	C	601	2XB	CBC-CBB	-2.09	1.37	1.43
2	B	600	UMP	P-OP3	2.05	1.62	1.54
3	B	601	2XB	CBC-CBB	-2.03	1.37	1.43

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	601	2XB	CBC-CBB	-2.02	1.37	1.43
2	D	600	UMP	P-OP3	2.00	1.62	1.54

All (39) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	601	2XB	CT1-NT1-CAV	4.26	127.84	122.34
2	C	600	UMP	C5-C4-N3	-4.08	114.33	123.31
3	A	601	2XB	NAA-CAW-NAP	4.02	124.33	117.79
2	D	600	UMP	C5-C4-N3	-3.98	114.56	123.31
2	B	600	UMP	C5-C4-N3	-3.95	114.62	123.31
3	B	601	2XB	CT1-NT1-CAV	3.95	127.43	122.34
2	A	600	UMP	C5-C4-N3	-3.93	114.65	123.31
3	D	601	2XB	NAA-CAW-NAP	3.91	124.17	117.79
3	C	601	2XB	NAA-CAW-NAP	3.91	124.16	117.79
3	D	601	2XB	CBC-CBA-NAS	-3.87	115.25	124.05
3	A	601	2XB	CBC-CBA-NAS	-3.87	115.26	124.05
3	B	601	2XB	CBC-CBA-NAS	-3.86	115.27	124.05
3	B	601	2XB	NAA-CAW-NAP	3.85	124.07	117.79
3	A	601	2XB	NAA-CAW-NAS	-3.80	111.33	117.25
3	B	601	2XB	NAA-CAW-NAS	-3.71	111.48	117.25
3	C	601	2XB	NAA-CAW-NAS	-3.69	111.51	117.25
3	C	601	2XB	CBC-CBA-NAS	-3.69	115.67	124.05
3	D	601	2XB	NAA-CAW-NAS	-3.66	111.57	117.25
2	B	600	UMP	O5'-P-OP1	3.64	116.69	106.47
2	C	600	UMP	OP2-P-O5'	3.45	115.91	106.73
2	D	600	UMP	O5'-P-OP1	3.44	116.12	106.47
2	A	600	UMP	O5'-P-OP1	3.22	115.50	106.47
3	A	601	2XB	CAZ-CAO-CAX	-2.76	107.58	114.21
3	A	601	2XB	CT1-NT1-CAV	2.61	125.71	122.34
3	D	601	2XB	CAW-NAS-CBA	2.53	122.81	116.43
3	D	601	2XB	CT2-CT1-NT1	2.52	113.86	110.19
3	B	601	2XB	CAW-NAS-CBA	2.47	122.65	116.43
3	A	601	2XB	CAW-NAS-CBA	2.46	122.64	116.43
3	C	601	2XB	CT2-CT1-NT1	2.40	113.69	110.19
3	C	601	2XB	CAW-NAS-CBA	2.36	122.38	116.43
3	C	601	2XB	CAZ-CAO-CAX	-2.31	108.68	114.21
3	D	601	2XB	NAP-CAW-NAS	-2.22	124.26	127.22
3	A	601	2XB	NAP-CAW-NAS	-2.17	124.33	127.22
3	C	601	2XB	NAP-CAW-NAS	-2.17	124.33	127.22
3	B	601	2XB	CAZ-CAO-CAX	-2.13	109.10	114.21
2	C	600	UMP	OP3-P-OP2	2.12	115.74	107.64

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	600	UMP	OP3-P-OP1	-2.10	102.47	110.68
3	B	601	2XB	NAP-CAW-NAS	-2.08	124.45	127.22
2	B	600	UMP	OP2-P-O5'	-2.04	101.31	106.73

There are no chirality outliers.

All (20) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	C	601	2XB	CT5-CT1-CT2-CT3
3	A	601	2XB	CT5-CT1-CT2-CT3
3	B	601	2XB	CT5-CT1-CT2-CT3
3	D	601	2XB	CT1-CT2-CT3-CT4
3	D	601	2XB	CT5-CT1-CT2-CT3
3	D	601	2XB	CT2-CT1-NT1-CAV
3	D	601	2XB	CT5-CT1-NT1-CAV
3	A	601	2XB	NT1-CT1-CT2-CT3
2	D	600	UMP	O4'-C4'-C5'-O5'
3	C	601	2XB	NT1-CT1-CT2-CT3
3	B	601	2XB	NT1-CT1-CT2-CT3
3	D	601	2XB	NT1-CT1-CT2-CT3
2	D	600	UMP	C3'-C4'-C5'-O5'
2	C	600	UMP	C5'-O5'-P-OP1
3	A	601	2XB	CT1-CT2-CT3-CT4
2	B	600	UMP	O4'-C4'-C5'-O5'
2	C	600	UMP	O4'-C4'-C5'-O5'
2	A	600	UMP	C4'-C5'-O5'-P
2	D	600	UMP	C5'-O5'-P-OP2
2	A	600	UMP	O4'-C4'-C5'-O5'

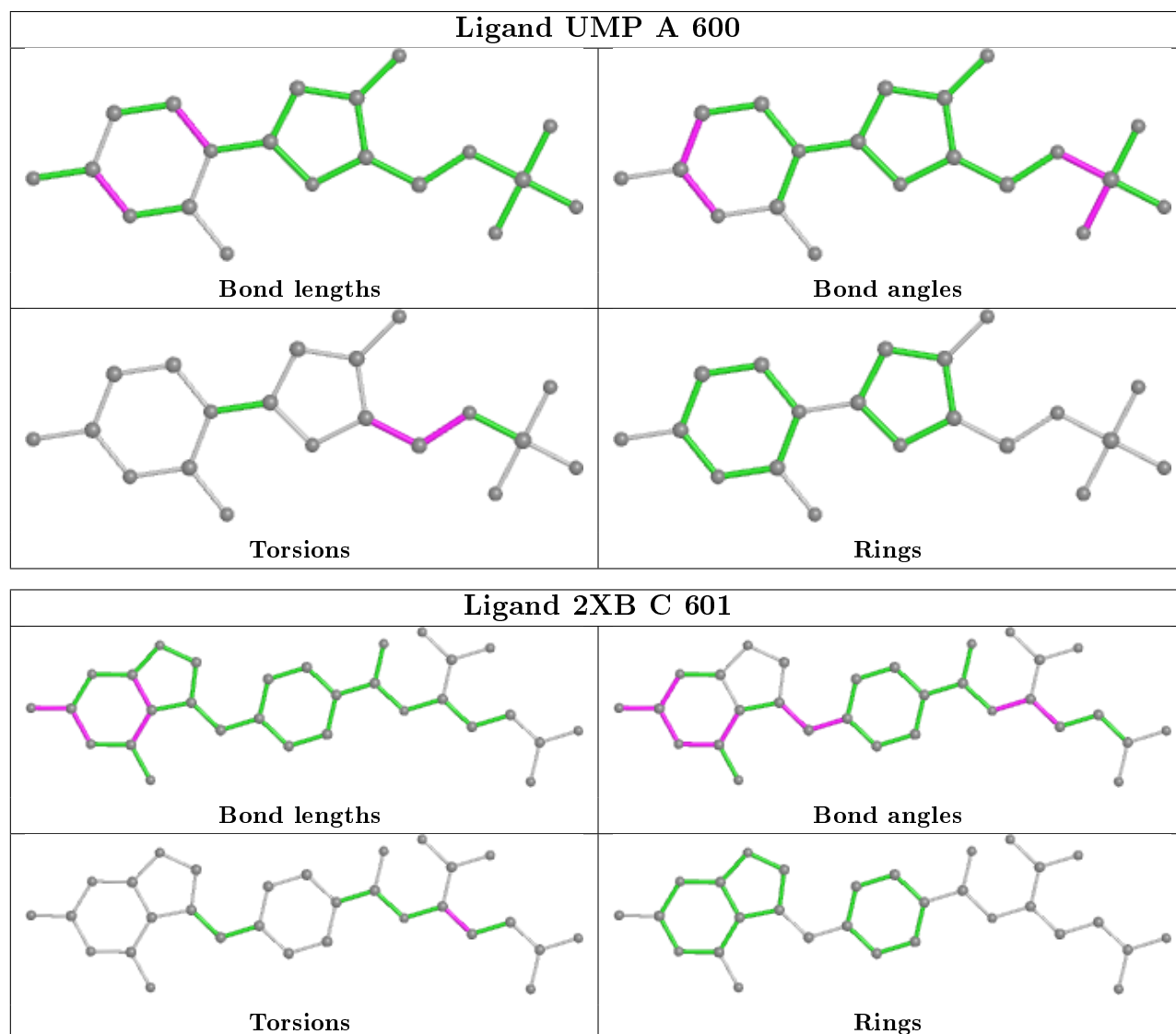
There are no ring outliers.

6 monomers are involved in 8 short contacts:

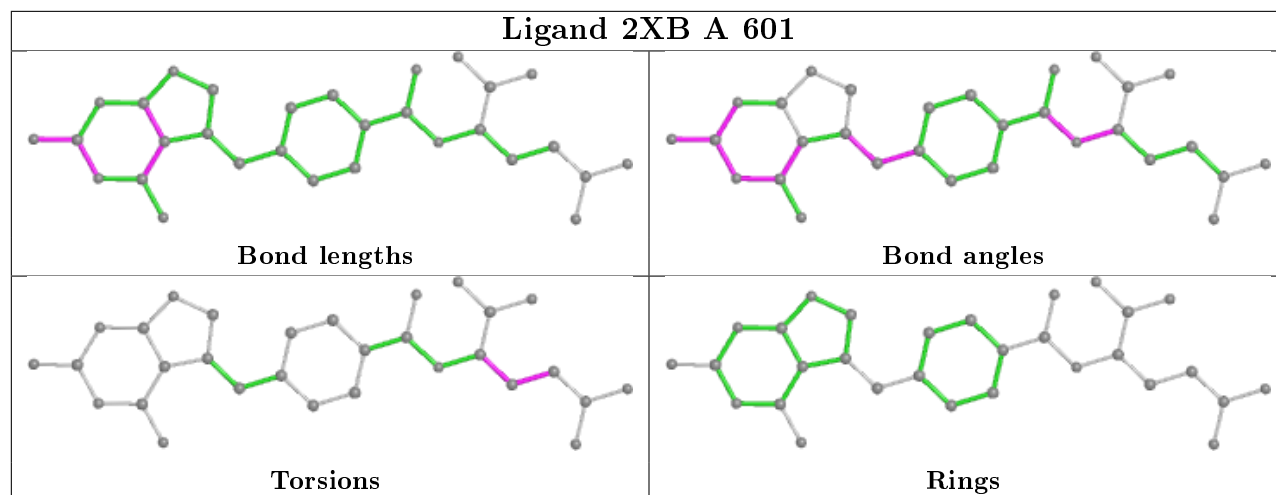
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	600	UMP	3	0
3	C	601	2XB	1	0
3	A	601	2XB	1	0
3	B	601	2XB	1	0
2	C	600	UMP	1	0
2	B	600	UMP	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths,

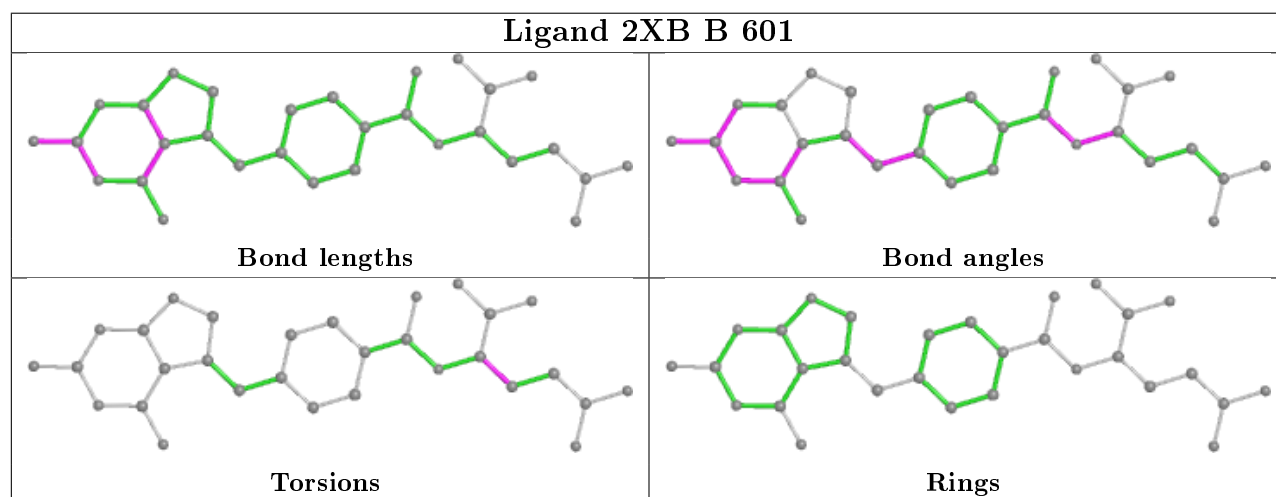
bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 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.



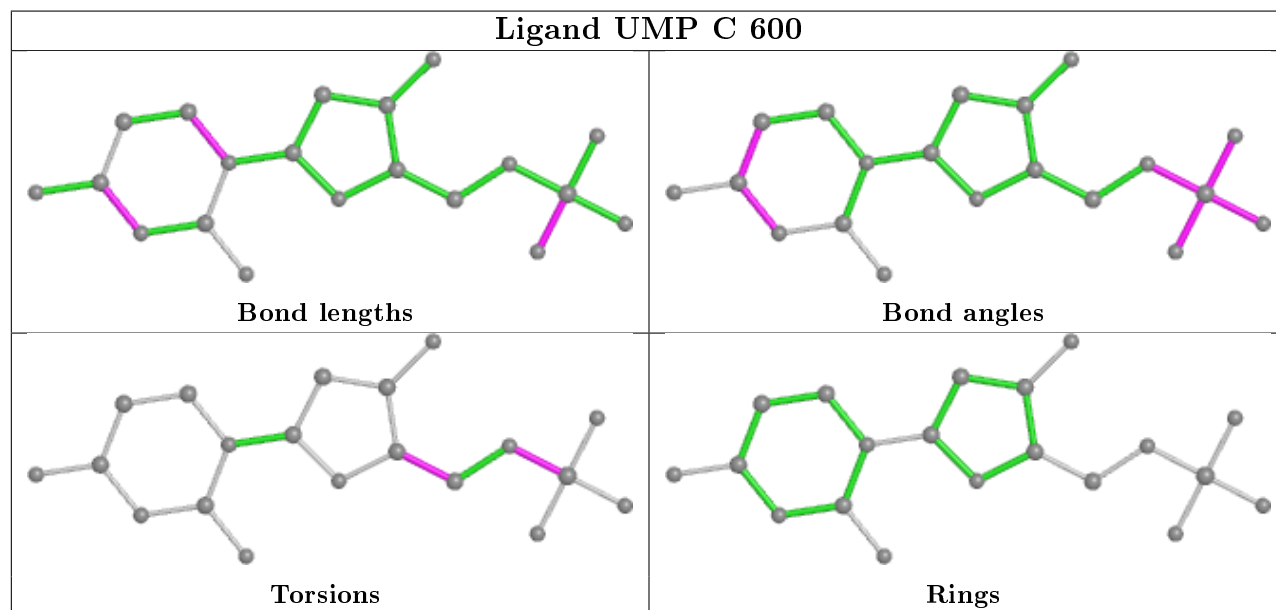
Ligand 2XB A 601



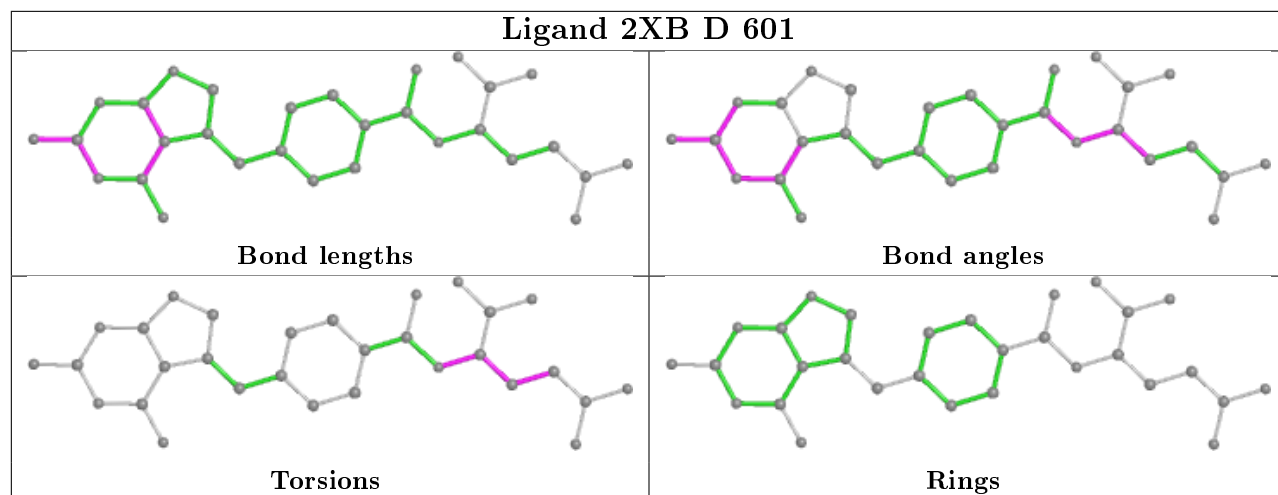
Ligand 2XB B 601



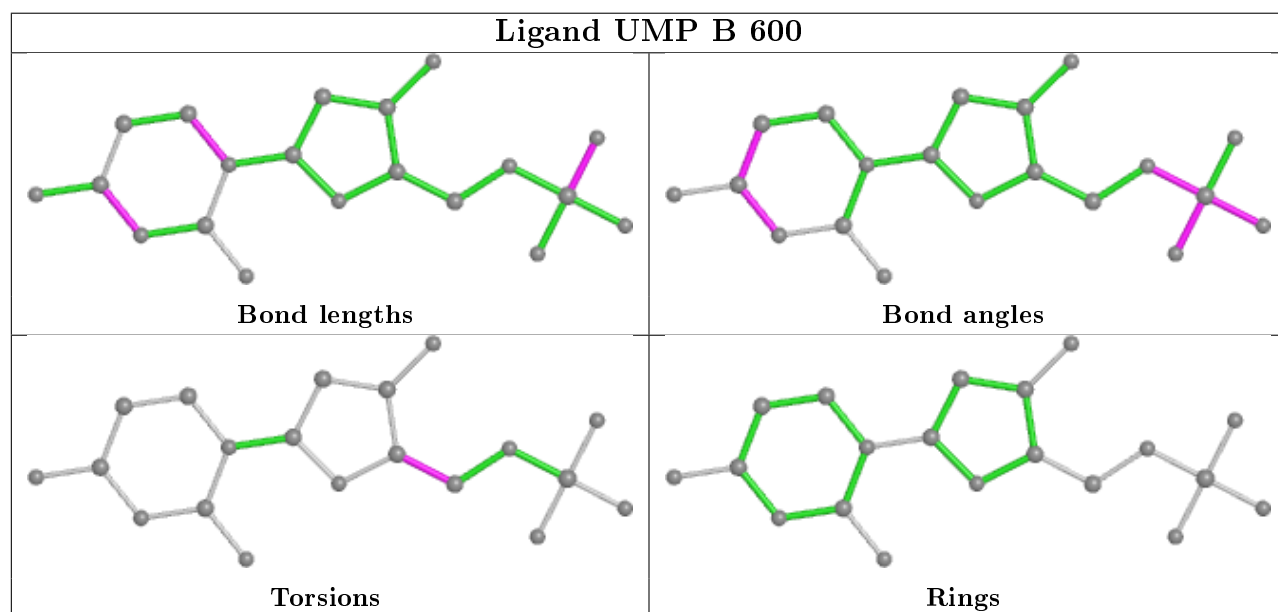
Ligand UMP C 600



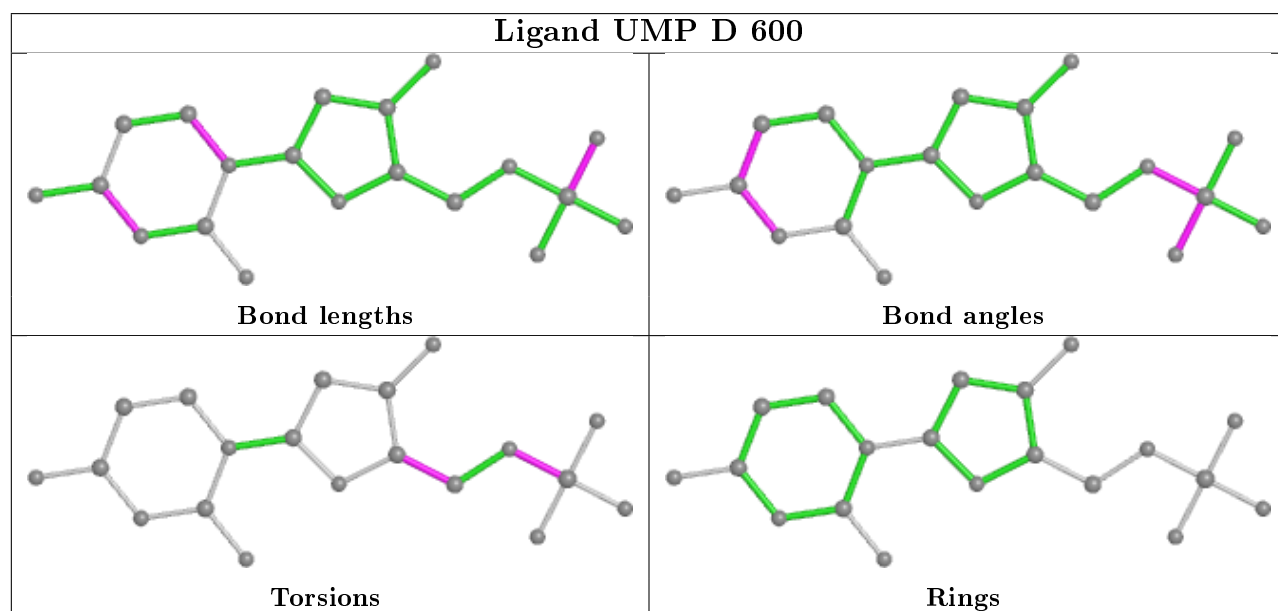
Ligand 2XB D 601



Ligand UMP B 600



Ligand UMP D 600



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	285/290 (98%)	0.76	14 (4%)	29 23	57, 65, 75, 83	0
1	B	285/290 (98%)	0.68	15 (5%)	26 20	51, 63, 82, 95	0
1	C	285/290 (98%)	0.69	15 (5%)	26 20	54, 67, 75, 84	0
1	D	285/290 (98%)	0.87	33 (11%)	4 3	62, 74, 85, 93	0
All	All	1140/1160 (98%)	0.75	77 (6%)	17 12	51, 67, 82, 95	0

All (77) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	282	LEU	5.0
1	D	297	GLN	4.7
1	D	79	VAL	3.7
1	C	285	VAL	3.5
1	D	279	LEU	3.4
1	A	83	GLY	3.4
1	A	285	VAL	3.4
1	D	262	ILE	3.3
1	B	213	TYR	3.3
1	C	121	LEU	3.2
1	D	80	PHE	3.2
1	A	180	CYS	3.1
1	A	212	LEU	3.0
1	D	54	GLY	3.0
1	A	240	ILE	3.0
1	D	83	GLY	3.0
1	C	110	ASP	2.9
1	B	212	LEU	2.9
1	A	98	ALA	2.9
1	B	197	ALA	2.8
1	D	84	VAL	2.8

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	B	157	GLY	2.7
1	A	289	ASP	2.7
1	C	291	PHE	2.7
1	C	286	GLU	2.7
1	D	197	ALA	2.6
1	C	297	GLN	2.6
1	B	182	TRP	2.6
1	B	125	THR	2.6
1	B	252	LEU	2.6
1	C	40	ILE	2.6
1	D	291	PHE	2.5
1	D	124	SER	2.5
1	C	165	ILE	2.5
1	C	137	PHE	2.5
1	D	180	CYS	2.5
1	B	123	PHE	2.5
1	B	6	SER	2.5
1	D	245	PRO	2.4
1	A	148	ASP	2.4
1	D	289	ASP	2.4
1	C	174	ASP	2.4
1	D	269	LEU	2.4
1	B	60	GLY	2.3
1	D	259	LEU	2.3
1	A	192	LEU	2.3
1	D	248	PHE	2.3
1	D	213	TYR	2.3
1	D	285	VAL	2.3
1	C	231	ALA	2.3
1	D	277	PRO	2.3
1	C	53	THR	2.3
1	A	280	ARG	2.2
1	D	182	TRP	2.2
1	D	37	ILE	2.2
1	B	229	SER	2.2
1	C	102	SER	2.2
1	D	82	LYS	2.1
1	B	48	ASP	2.1
1	D	242	GLY	2.1
1	A	286	GLU	2.1
1	B	58	VAL	2.1
1	A	71	PHE	2.1

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	C	214	GLN	2.1
1	D	198	LEU	2.1
1	D	153	TYR	2.1
1	D	225	PHE	2.1
1	B	198	LEU	2.1
1	A	221	LEU	2.1
1	B	149	MET	2.1
1	D	65	TYR	2.1
1	D	240	ILE	2.1
1	D	81	TRP	2.1
1	C	296	PHE	2.0
1	D	101	LEU	2.0
1	D	69	ASP	2.0
1	A	255	ALA	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](#)

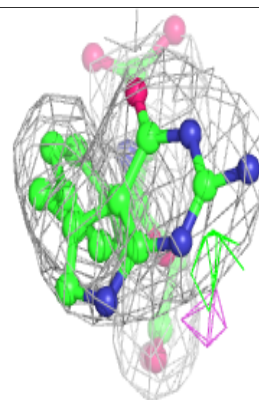
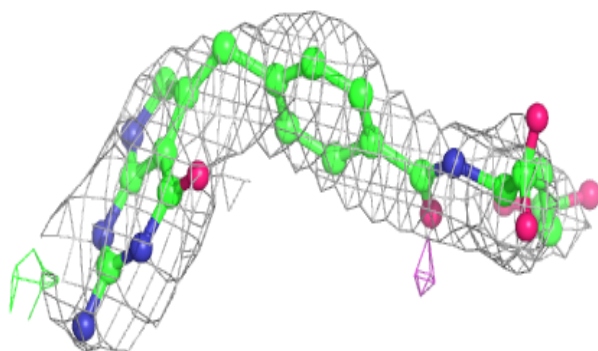
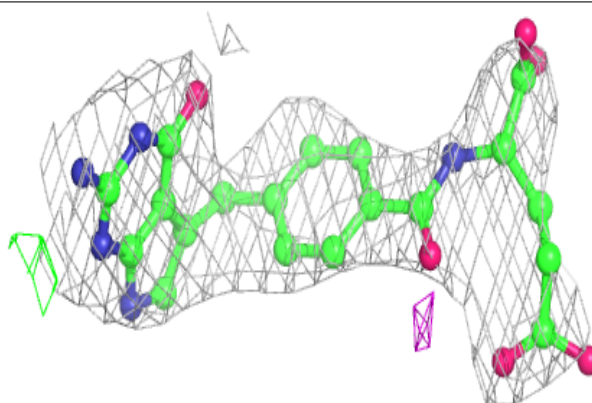
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	2XB	D	601	30/30	0.82	0.27	74,80,93,96	0
3	2XB	C	601	30/30	0.83	0.26	59,71,87,91	0
3	2XB	B	601	30/30	0.86	0.21	56,62,81,84	0
3	2XB	A	601	30/30	0.86	0.20	60,67,73,76	0
2	UMP	C	600	20/20	0.90	0.23	55,61,64,64	0
2	UMP	A	600	20/20	0.90	0.19	55,63,67,67	0
2	UMP	B	600	20/20	0.93	0.23	54,62,66,68	0
2	UMP	D	600	20/20	0.93	0.20	60,73,77,77	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.

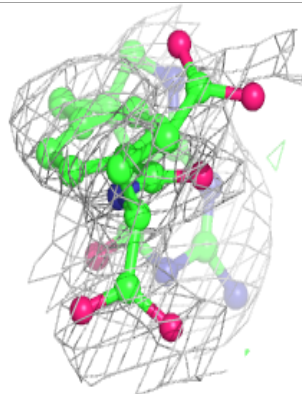
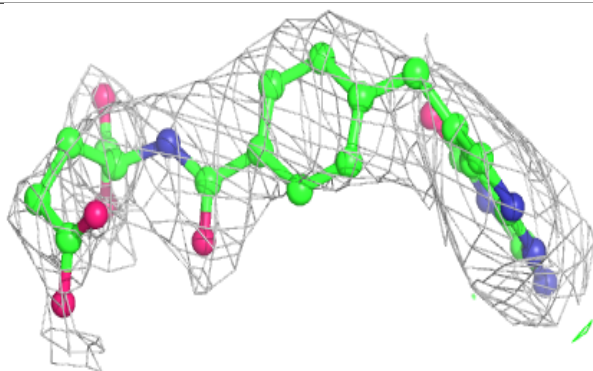
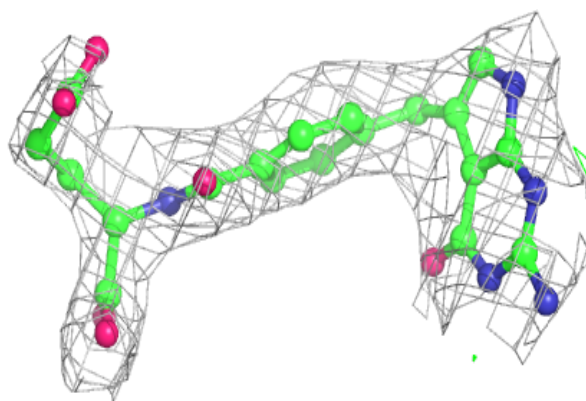
Electron density around 2XB D 601:

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

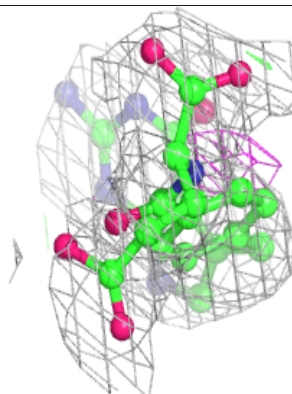
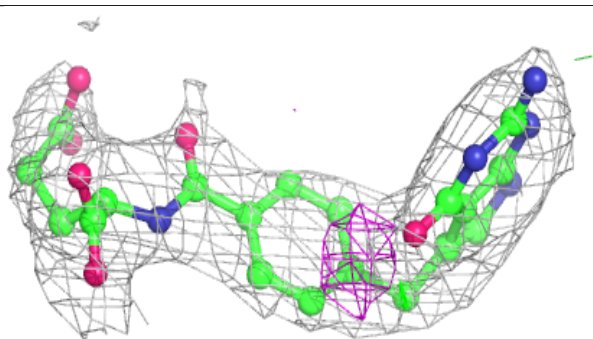
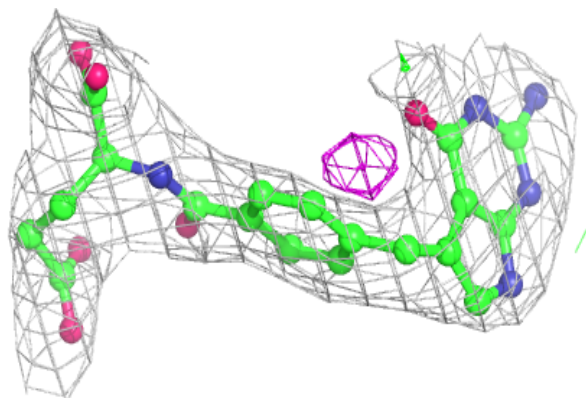


Electron density around 2XB C 601:

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

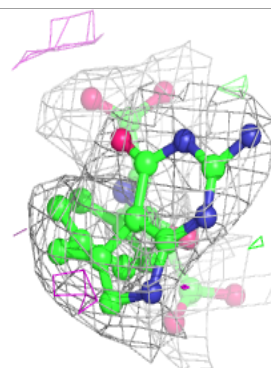
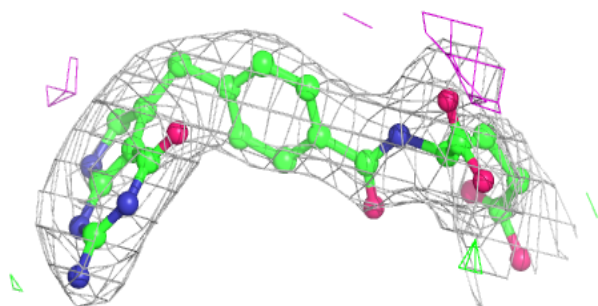
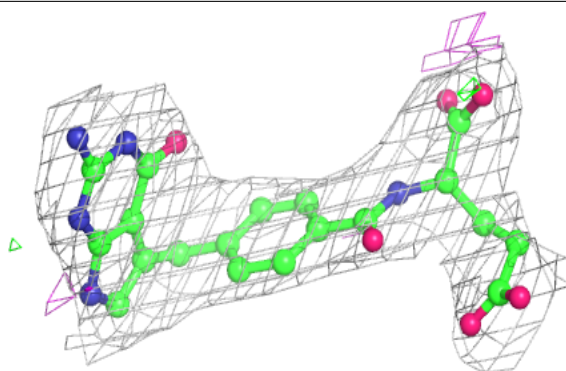
**Electron density around 2XB B 601:**

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

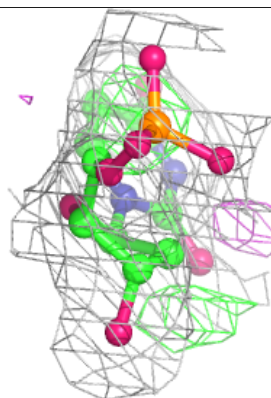
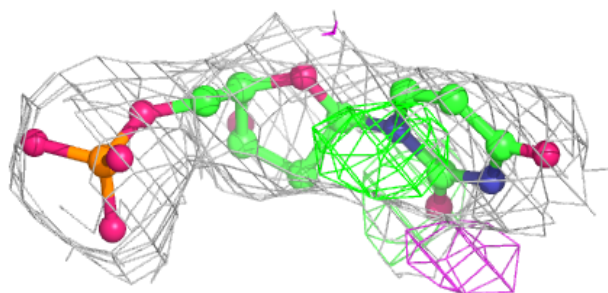
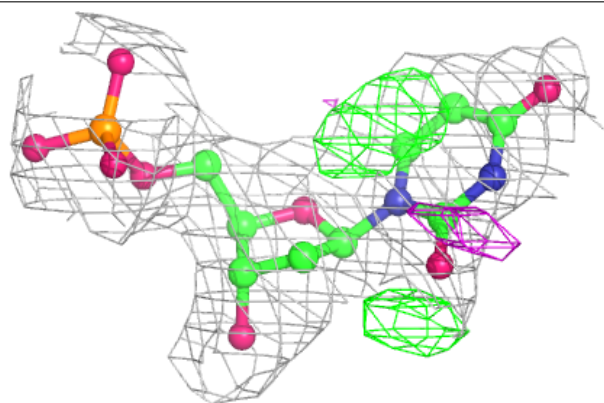


Electron density around 2XB A 601:

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

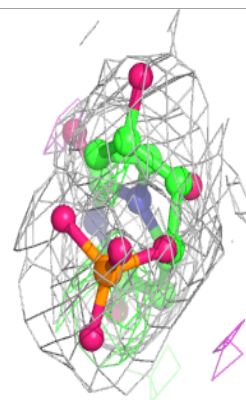
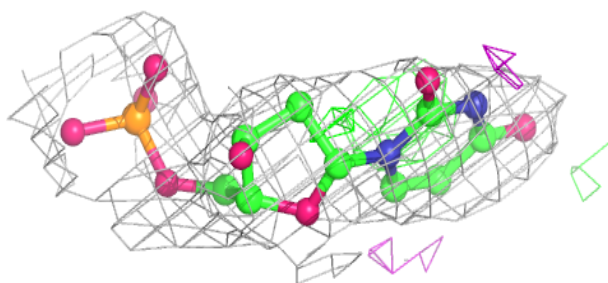
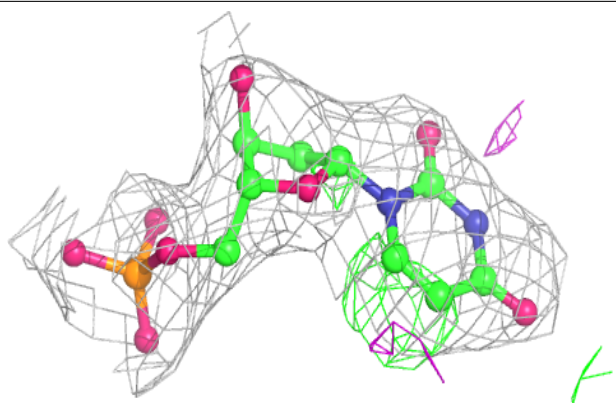
**Electron density around UMP C 600:**

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

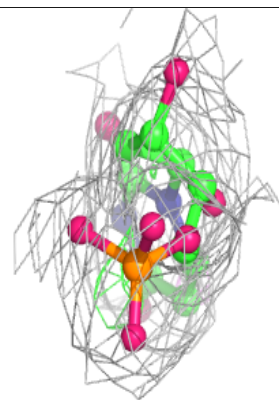
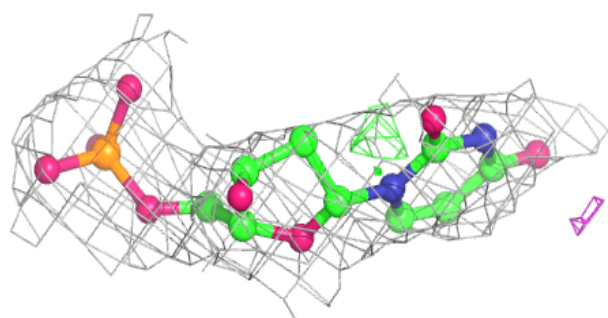
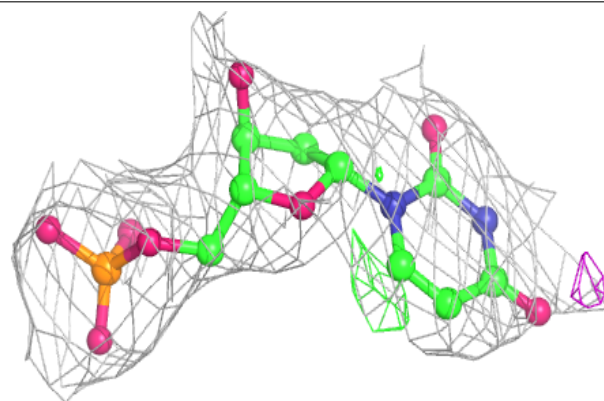


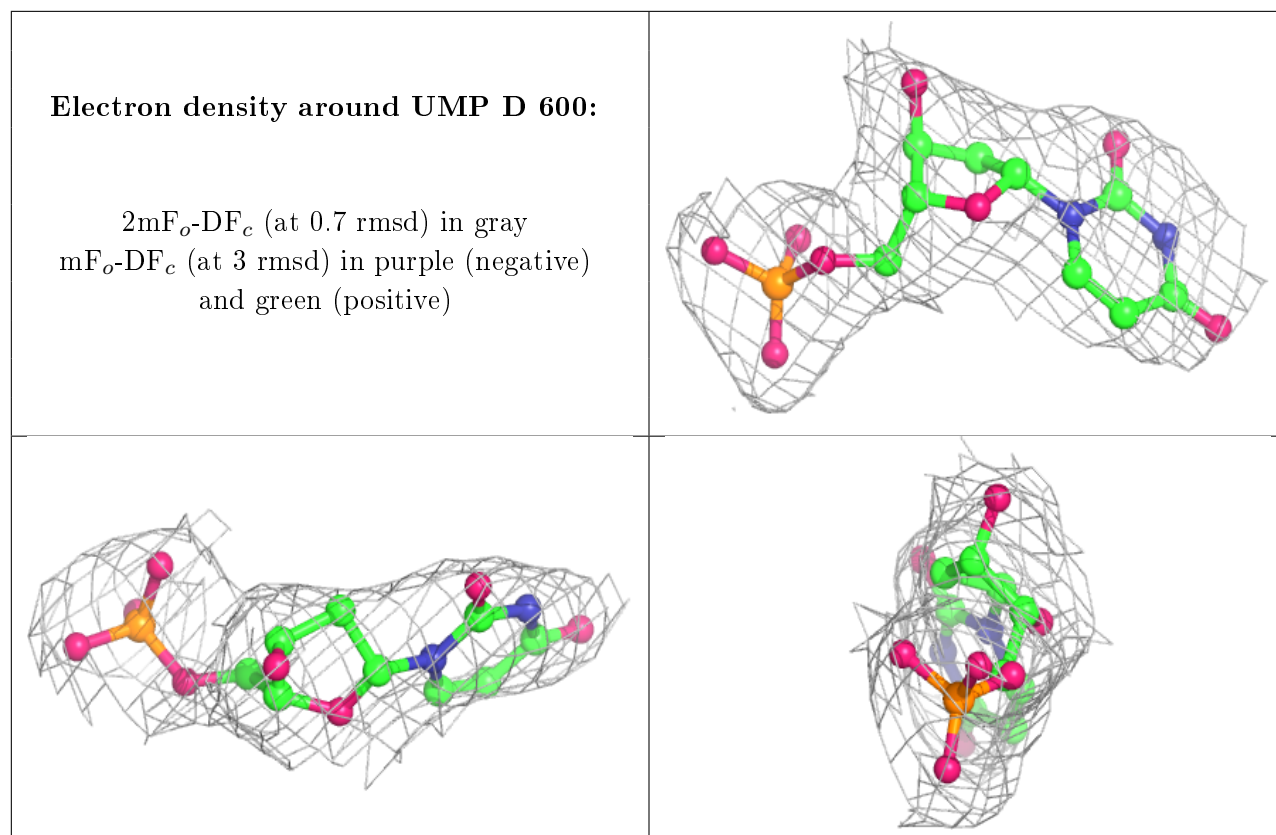
Electron density around UMP A 600:

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

**Electron density around UMP B 600:**

$2mF_o-DF_c$ (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.