



Full wwPDB X-ray Structure Validation Report ⓘ

May 17, 2020 – 04:51 am BST

PDB ID : 4IQQ
Title : Crystal Structure of C.elegans Thymidylate Synthase in Complex with dUMP and Tomudex
Authors : Wilk, P.; Dowiercial, A.; Banaszak, K.; Jarmula, A.; Rypniewski, W.; Rode, W.
Deposited on : 2013-01-13
Resolution : 2.90 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

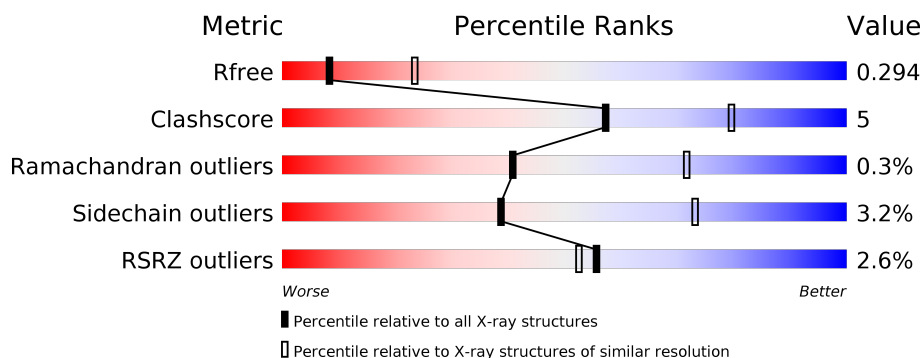
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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	1957 (2.90-2.90)
Clashscore	141614	2172 (2.90-2.90)
Ramachandran outliers	138981	2115 (2.90-2.90)
Sidechain outliers	138945	2117 (2.90-2.90)
RSRZ outliers	127900	1906 (2.90-2.90)

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	315	<div> <div>2%</div> <div>83% 8% 8%</div> </div>
1	B	315	<div> <div>3%</div> <div>78% 13% 8%</div> </div>
1	C	315	<div> <div>2%</div> <div>79% 13% 9%</div> </div>
1	D	315	<div> <div>3%</div> <div>76% 14% 9%</div> </div>

2 Entry composition [i](#)

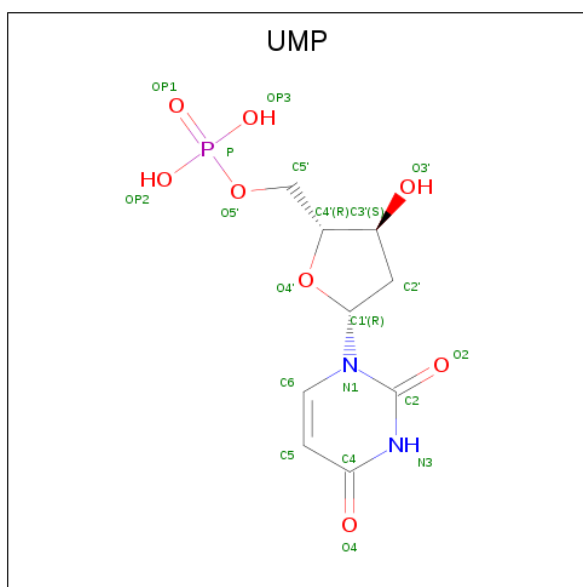
There are 4 unique types of molecules in this entry. The entry contains 9791 atoms, of which 0 are hydrogens and 0 are deuteriums.

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.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	289	Total	C	N	O	S	0	0	0
			2316	1468	401	431	16			
1	B	289	Total	C	N	O	S	0	3	0
			2336	1481	404	435	16			
1	C	288	Total	C	N	O	S	0	2	0
			2319	1471	402	430	16			
1	D	288	Total	C	N	O	S	0	0	0
			2307	1463	399	429	16			

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



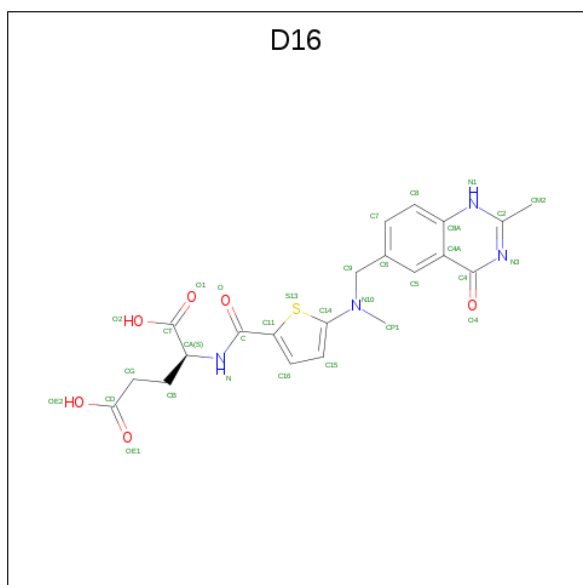
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			20	9	2	8	1		
2	B	1	Total	C	N	O	P	0	0
			20	9	2	8	1		

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

- Molecule 3 is TOMUDEX (three-letter code: D16) (formula: C₂₁H₂₂N₄O₆S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	S	0	0
			32	21	4	6	1		
3	B	1	Total	C	N	O	S	0	0
			32	21	4	6	1		
3	C	1	Total	C	N	O	S	0	0
			32	21	4	6	1		
3	D	1	Total	C	N	O	S	0	0
			32	21	4	6	1		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	80	Total	O	0	0
			80	80		
4	B	84	Total	O	0	0
			84	84		
4	C	105	Total	O	0	0
			105	105		

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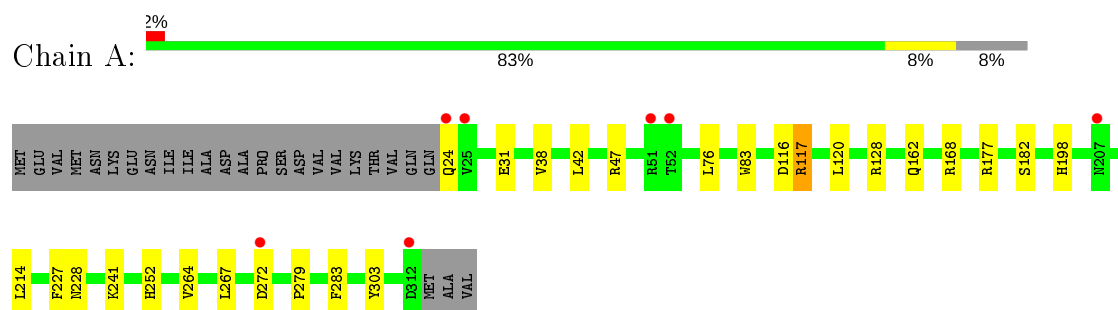
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	D	36	Total	O	0	0
			36	36		

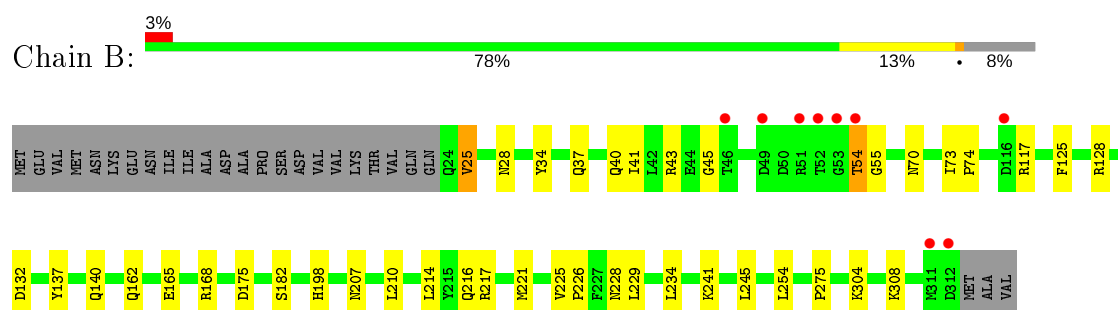
3 Residue-property plots [i](#)

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.

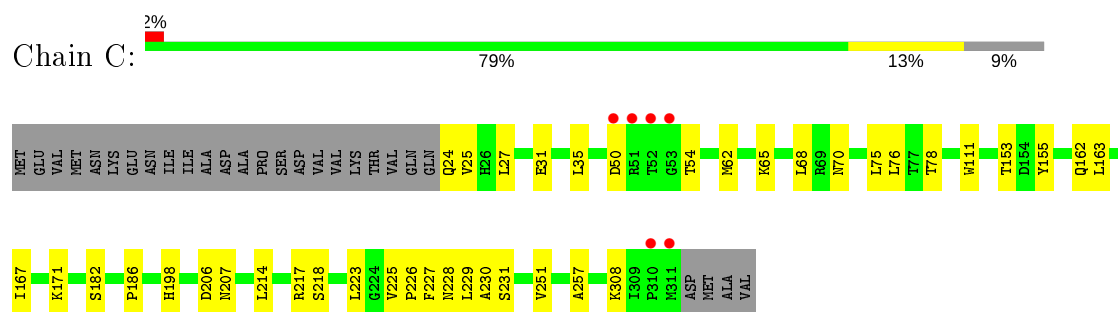
• Molecule 1: Thymidylate synthase



• Molecule 1: Thymidylate synthase

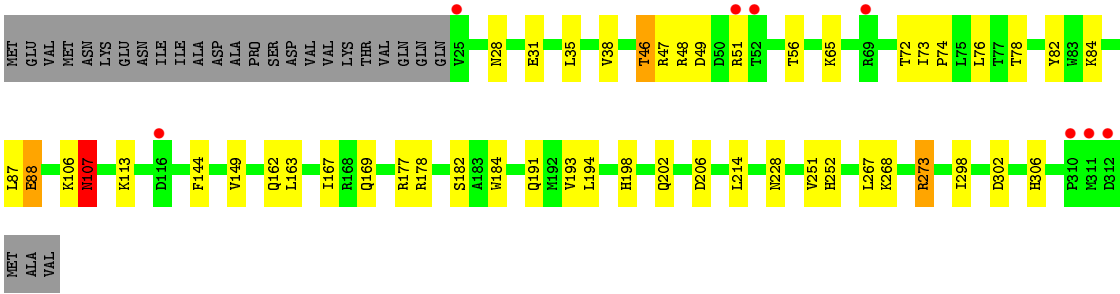


• Molecule 1: Thymidylate synthase



• Molecule 1: Thymidylate synthase





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	70.08Å 94.31Å 122.05Å 90.00° 105.36° 90.00°	Depositor
Resolution (Å)	14.70 – 2.90 14.71 – 2.90	Depositor EDS
% Data completeness (in resolution range)	96.0 (14.70-2.90) 96.8 (14.71-2.90)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.14	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.78 (at 2.90Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.218 , 0.295 0.217 , 0.294	Depositor DCC
R_{free} test set	1666 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å ²)	42.9	Xtriage
Anisotropy	0.112	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 47.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.035 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	9791	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 4.26% 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: D16, 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	0.46	1/2370 (0.0%)	0.53	0/3204
1	B	0.46	0/2399	0.54	0/3242
1	C	0.47	1/2379 (0.0%)	0.52	0/3216
1	D	0.47	1/2361 (0.0%)	0.50	0/3192
All	All	0.46	3/9509 (0.0%)	0.52	0/12854

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	111	TRP	CD2-CE2	5.24	1.47	1.41
1	A	83	TRP	CD2-CE2	5.04	1.47	1.41
1	D	184	TRP	CD2-CE2	5.03	1.47	1.41

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	2316	0	2270	22	0
1	B	2336	0	2295	24	0
1	C	2319	0	2280	27	0
1	D	2307	0	2262	28	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	20	0	11	1	0
2	B	20	0	11	2	0
2	C	20	0	11	1	0
2	D	20	0	11	1	0
3	A	32	0	20	2	0
3	B	32	0	20	1	0
3	C	32	0	20	2	0
3	D	32	0	20	1	0
4	A	80	0	0	0	0
4	B	84	0	0	2	0
4	C	105	0	0	1	0
4	D	36	0	0	2	0
All	All	9791	0	9231	91	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 5.

All (91) 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:D:228:ASN:HD21	2:D:401:UMP:HN3	1.15	0.94
1:B:228:ASN:HD21	2:B:401:UMP:HN3	1.20	0.87
1:A:168:ARG:HH11	1:A:168:ARG:HG2	1.46	0.79
1:C:228:ASN:HD21	2:C:401:UMP:HN3	1.32	0.74
1:A:228:ASN:HD21	2:A:401:UMP:HN3	1.39	0.71
1:B:165[B]:GLU:OE2	1:B:168[B]:ARG:CZ	2.42	0.68
1:C:163:LEU:O	1:C:167:ILE:HG12	1.96	0.66
1:A:42:LEU:HD13	1:C:25:VAL:HG11	1.77	0.65
1:A:264:VAL:HG21	1:C:27:LEU:HB2	1.78	0.64
1:C:62:MET:HE2	1:C:257:ALA:HB1	1.81	0.63
1:D:193:VAL:HG23	1:D:194:LEU:H	1.67	0.59
1:A:214:LEU:HD23	1:A:252:HIS:HD2	1.69	0.58
1:C:167:ILE:HG22	1:C:171:LYS:HE2	1.84	0.57
1:B:165[B]:GLU:OE2	1:B:168[B]:ARG:NH1	2.37	0.57
1:C:62:MET:HE2	1:C:257:ALA:CB	2.35	0.57
1:D:31:GLU:HG3	1:D:76:LEU:HD22	1.85	0.56
1:B:162:GLN:HE22	1:B:182:SER:H	1.54	0.54
1:D:162:GLN:HE22	1:D:182:SER:H	1.55	0.54
1:A:162:GLN:HE22	1:A:182:SER:H	1.56	0.53
1:A:31:GLU:HG3	1:A:76:LEU:HD22	1.89	0.53
1:C:217:ARG:HA	1:D:202:GLN:HE22	1.74	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:50:ASP:HB2	1:C:54:THR:H	1.72	0.53
1:A:117:ARG:HH21	1:A:120:LEU:HD23	1.74	0.52
1:C:24:GLN:HB2	4:C:602:HOH:O	2.09	0.52
1:A:168:ARG:HG2	1:A:168:ARG:NH1	2.20	0.52
1:C:223:LEU:HD11	3:C:402:D16:HG2	1.91	0.51
1:D:214:LEU:HD23	1:D:252:HIS:CD2	2.46	0.51
1:A:116:ASP:O	1:A:120:LEU:HB2	2.11	0.51
1:B:198:HIS:HB2	1:B:214:LEU:HD11	1.91	0.51
1:B:216:GLN:HB3	1:B:254:LEU:HD23	1.93	0.51
1:D:84:LYS:HE2	1:D:107:ASN:HB3	1.93	0.51
1:A:227:PHE:CG	3:A:402:D16:H15	2.46	0.50
1:D:163:LEU:O	1:D:167:ILE:HG12	2.11	0.50
1:C:198:HIS:HB2	1:C:214:LEU:HD11	1.92	0.50
3:D:402:D16:HB2	4:D:504:HOH:O	2.11	0.50
1:B:41:ILE:O	1:B:45:GLY:HA3	2.12	0.49
1:D:65:LYS:HG2	1:D:251:VAL:HG22	1.94	0.49
1:A:198:HIS:CB	1:A:214:LEU:HD11	2.42	0.49
1:B:210:LEU:HB2	1:B:245:LEU:HD23	1.94	0.49
1:D:88:GLU:HG2	1:D:106:LYS:HB3	1.95	0.48
1:A:279:PRO:HG3	1:A:303:TYR:CD1	2.47	0.48
1:C:308:LYS:H	1:C:308:LYS:HD2	1.78	0.48
1:B:128:ARG:HD3	1:B:132:ASP:HB3	1.96	0.48
1:A:162:GLN:NE2	1:A:182:SER:H	2.12	0.48
1:C:31:GLU:HG3	1:C:76:LEU:HD22	1.95	0.48
1:B:40:GLN:HG3	1:B:43:ARG:HH12	1.78	0.47
1:B:125:PHE:HB3	1:B:128:ARG:HG2	1.96	0.47
1:B:37:GLN:HG3	4:B:579:HOH:O	2.13	0.47
1:C:206:ASP:OD1	1:D:46:THR:HG21	2.13	0.47
1:C:35:LEU:HD11	1:C:78:THR:HG21	1.97	0.47
1:D:38:VAL:HG11	1:D:267:LEU:HD12	1.96	0.46
1:B:34:TYR:OH	1:B:221:MET:O	2.32	0.46
1:B:25:VAL:HG12	1:D:268:LYS:HG2	1.97	0.46
1:B:25:VAL:O	1:D:268:LYS:HD3	2.15	0.46
1:C:65:LYS:HG2	1:C:251:VAL:HG22	1.98	0.46
1:B:54:THR:HG22	1:B:55:GLY:H	1.81	0.46
1:D:169:GLN:HE22	1:D:178:ARG:HG2	1.81	0.45
1:B:73:ILE:HA	1:B:74:PRO:HD3	1.83	0.45
1:D:193:VAL:HG23	1:D:194:LEU:N	2.31	0.45
1:B:28:ASN:HB3	1:B:275:PRO:HB2	1.99	0.45
1:C:75:LEU:HD23	1:C:230:ALA:HB2	1.99	0.45
1:B:198:HIS:CB	1:B:214:LEU:HD11	2.47	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:35:LEU:HD11	1:D:78:THR:HG21	1.97	0.45
1:D:198:HIS:HB2	1:D:214:LEU:HD11	1.98	0.44
1:C:162:GLN:HE22	1:C:182:SER:H	1.64	0.44
3:B:402:D16:H16	4:B:522:HOH:O	2.17	0.44
1:A:227:PHE:CD2	3:A:402:D16:H15	2.52	0.44
1:A:214:LEU:HD23	1:A:252:HIS:CD2	2.49	0.44
1:C:186:PRO:HD2	1:D:144:PHE:CE2	2.53	0.44
1:B:207:ASN:H	1:C:70:ASN:HD21	1.66	0.43
1:B:70:ASN:H	1:C:207:ASN:ND2	2.16	0.43
1:C:227:PHE:CG	3:C:402:D16:H15	2.53	0.43
1:B:70:ASN:HD21	1:C:207:ASN:H	1.67	0.43
1:C:218:SER:OG	1:D:177:ARG:HD3	2.20	0.42
1:A:117:ARG:NH2	1:A:128:ARG:O	2.52	0.42
1:D:214:LEU:HD23	1:D:252:HIS:HD2	1.83	0.42
1:D:191:GLN:HG2	4:D:536:HOH:O	2.19	0.42
1:C:225:VAL:HB	1:C:226:PRO:HD3	2.01	0.42
1:D:87:LEU:HD11	1:D:298:ILE:HB	2.01	0.42
1:A:177:ARG:HD2	2:B:401:UMP:OP2	2.19	0.42
1:D:73:ILE:HA	1:D:74:PRO:HD3	1.89	0.42
1:A:241:LYS:HE3	1:A:283:PHE:HB2	2.03	0.41
1:D:273:ARG:HG2	1:D:306:HIS:CG	2.56	0.41
1:A:117:ARG:NH2	1:A:120:LEU:HD23	2.34	0.41
1:D:178:ARG:HG3	1:D:178:ARG:O	2.20	0.41
1:D:106:LYS:O	1:D:107:ASN:CB	2.68	0.41
1:A:177:ARG:O	1:B:217:ARG:HD2	2.20	0.41
1:D:178:ARG:CG	1:D:178:ARG:O	2.68	0.40
1:B:225:VAL:HB	1:B:226:PRO:HD3	2.03	0.40
1:A:38:VAL:HG11	1:A:267:LEU:HD12	2.03	0.40
1:C:153:THR:HB	1:C:155:TYR:CZ	2.57	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	287/315 (91%)	272 (95%)	15 (5%)	0	100	100
1	B	290/315 (92%)	274 (94%)	15 (5%)	1 (0%)	41	71
1	C	288/315 (91%)	273 (95%)	15 (5%)	0	100	100
1	D	286/315 (91%)	267 (93%)	17 (6%)	2 (1%)	22	54
All	All	1151/1260 (91%)	1086 (94%)	62 (5%)	3 (0%)	41	71

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	137	TYR
1	D	107	ASN
1	D	82	TYR

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	253/276 (92%)	249 (98%)	4 (2%)	62	86
1	B	256/276 (93%)	246 (96%)	10 (4%)	32	66
1	C	254/276 (92%)	251 (99%)	3 (1%)	71	91
1	D	252/276 (91%)	237 (94%)	15 (6%)	19	49
All	All	1015/1104 (92%)	983 (97%)	32 (3%)	39	73

All (32) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	24	GLN
1	A	47	ARG
1	A	117	ARG
1	A	272	ASP
1	B	25	VAL
1	B	54	THR
1	B	117	ARG
1	B	140	GLN

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Mol	Chain	Res	Type
1	B	175	ASP
1	B	229	LEU
1	B	234	LEU
1	B	241	LYS
1	B	304	LYS
1	B	308	LYS
1	C	68	LEU
1	C	229	LEU
1	C	231	SER
1	D	28	ASN
1	D	46	THR
1	D	47	ARG
1	D	48	ARG
1	D	49	ASP
1	D	51	ARG
1	D	56	THR
1	D	72	THR
1	D	88	GLU
1	D	107	ASN
1	D	113	LYS
1	D	149	VAL
1	D	206	ASP
1	D	273	ARG
1	D	302	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (26) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	63	GLN
1	A	143	HIS
1	A	162	GLN
1	A	191	GLN
1	A	213	GLN
1	A	228	ASN
1	A	262	ASN
1	B	24	GLN
1	B	70	ASN
1	B	162	GLN
1	B	202	GLN
1	B	228	ASN
1	C	40	GLN
1	C	70	ASN

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Mol	Chain	Res	Type
1	C	162	GLN
1	C	207	ASN
1	C	228	ASN
1	D	28	ASN
1	D	70	ASN
1	D	114	ASN
1	D	158	GLN
1	D	162	GLN
1	D	202	GLN
1	D	213	GLN
1	D	228	ASN
1	D	263	HIS

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
3	D16	A	402	-	23,34,34	1.21	1 (4%)	25,48,48	3.21	6 (24%)
2	UMP	B	401	-	18,21,21	0.71	0	21,31,31	0.99	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	D16	B	402	-	23,34,34	1.18	2 (8%)	25,48,48	3.26	5 (20%)
2	UMP	A	401	-	18,21,21	0.72	0	21,31,31	0.91	0
3	D16	D	402	-	23,34,34	1.19	1 (4%)	25,48,48	3.28	6 (24%)
2	UMP	C	401	-	18,21,21	0.74	0	21,31,31	0.89	0
3	D16	C	402	-	23,34,34	1.20	1 (4%)	25,48,48	3.32	8 (32%)
2	UMP	D	401	-	18,21,21	0.70	0	21,31,31	0.97	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
3	D16	A	402	-	-	1/13/25/25	0/3/3/3
2	UMP	B	401	-	-	1/7/22/22	0/2/2/2
3	D16	B	402	-	-	3/13/25/25	0/3/3/3
2	UMP	A	401	-	-	1/7/22/22	0/2/2/2
3	D16	D	402	-	-	7/13/25/25	0/3/3/3
2	UMP	C	401	-	-	1/7/22/22	0/2/2/2
3	D16	C	402	-	-	1/13/25/25	0/3/3/3
2	UMP	D	401	-	-	1/7/22/22	0/2/2/2

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	402	D16	C4A-C8A	3.88	1.48	1.41
3	D	402	D16	C4A-C8A	3.87	1.48	1.41
3	C	402	D16	C4A-C8A	3.86	1.48	1.41
3	B	402	D16	C4A-C8A	3.70	1.48	1.41
3	B	402	D16	C8A-N1	-2.00	1.34	1.37

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	402	D16	C4A-C4-N3	-9.90	117.50	124.40
3	D	402	D16	C4A-C8A-N1	-9.18	118.62	123.60
3	A	402	D16	C4A-C4-N3	-9.11	118.05	124.40
3	C	402	D16	C4A-C4-N3	-9.06	118.08	124.40
3	D	402	D16	C4A-C4-N3	-8.99	118.13	124.40
3	A	402	D16	C4A-C8A-N1	-8.91	118.76	123.60
3	C	402	D16	C4A-C8A-N1	-8.67	118.89	123.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	402	D16	C4A-C8A-N1	-8.58	118.94	123.60
3	C	402	D16	C2-N1-C8A	8.03	121.99	116.54
3	A	402	D16	C2-N1-C8A	7.96	121.94	116.54
3	D	402	D16	C2-N1-C8A	7.91	121.90	116.54
3	B	402	D16	C2-N1-C8A	7.82	121.84	116.54
3	C	402	D16	CA-N-C	4.08	127.60	122.34
3	D	402	D16	CA-N-C	3.05	126.27	122.34
3	D	402	D16	C8-C8A-N1	2.36	122.29	118.69
3	C	402	D16	CB-CG-CD	-2.26	108.73	113.59
3	C	402	D16	C6-C5-C4A	-2.19	119.37	122.65
3	A	402	D16	C8-C8A-N1	2.19	122.03	118.69
3	C	402	D16	C4-C4A-C8A	-2.18	116.70	118.59
3	C	402	D16	C8-C8A-N1	2.12	121.92	118.69
3	B	402	D16	CB-CG-CD	-2.10	109.08	113.59
3	B	402	D16	C8-C8A-N1	2.09	121.88	118.69
3	D	402	D16	C6-C5-C4A	-2.06	119.57	122.65
3	A	402	D16	C4-C4A-C8A	-2.04	116.81	118.59
3	A	402	D16	C6-C5-C4A	-2.03	119.61	122.65

There are no chirality outliers.

All (16) torsion outliers are listed below:

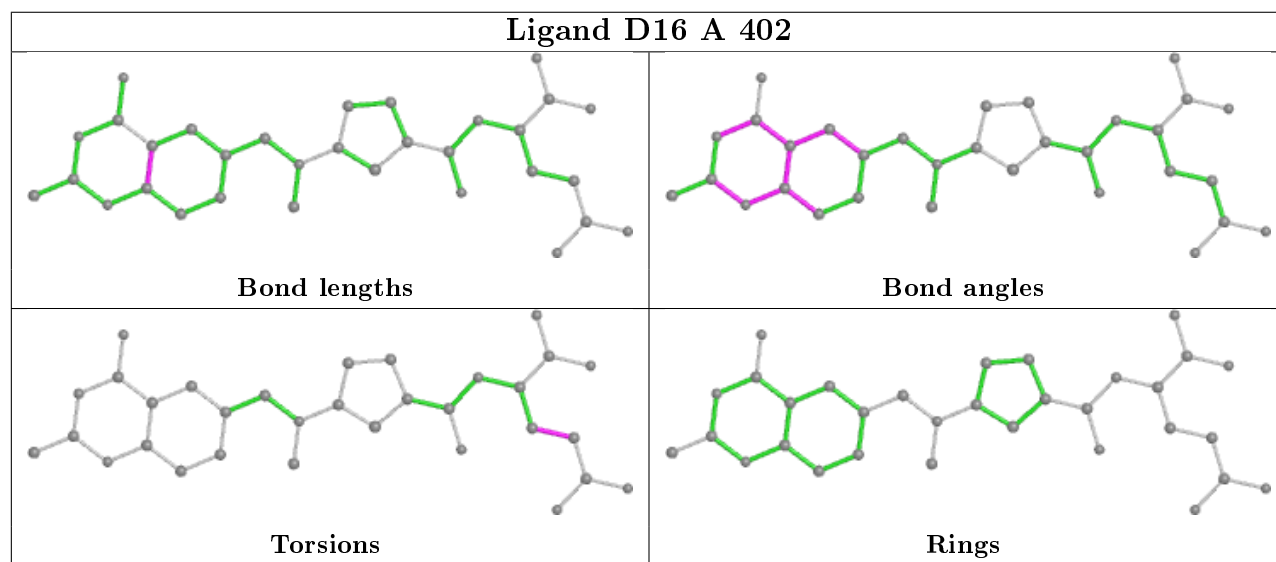
Mol	Chain	Res	Type	Atoms
3	B	402	D16	CT-CA-CB-CG
3	D	402	D16	C11-C-N-CA
3	D	402	D16	N-CA-CB-CG
3	D	402	D16	CT-CA-CB-CG
3	C	402	D16	CB-CA-N-C
3	D	402	D16	O-C-N-CA
3	B	402	D16	N-CA-CB-CG
3	D	402	D16	CA-CB-CG-CD
3	D	402	D16	CT-CA-N-C
3	A	402	D16	CA-CB-CG-CD
3	B	402	D16	CA-CB-CG-CD
3	D	402	D16	CB-CA-N-C
2	A	401	UMP	O4'-C4'-C5'-O5'
2	C	401	UMP	O4'-C4'-C5'-O5'
2	D	401	UMP	O4'-C4'-C5'-O5'
2	B	401	UMP	O4'-C4'-C5'-O5'

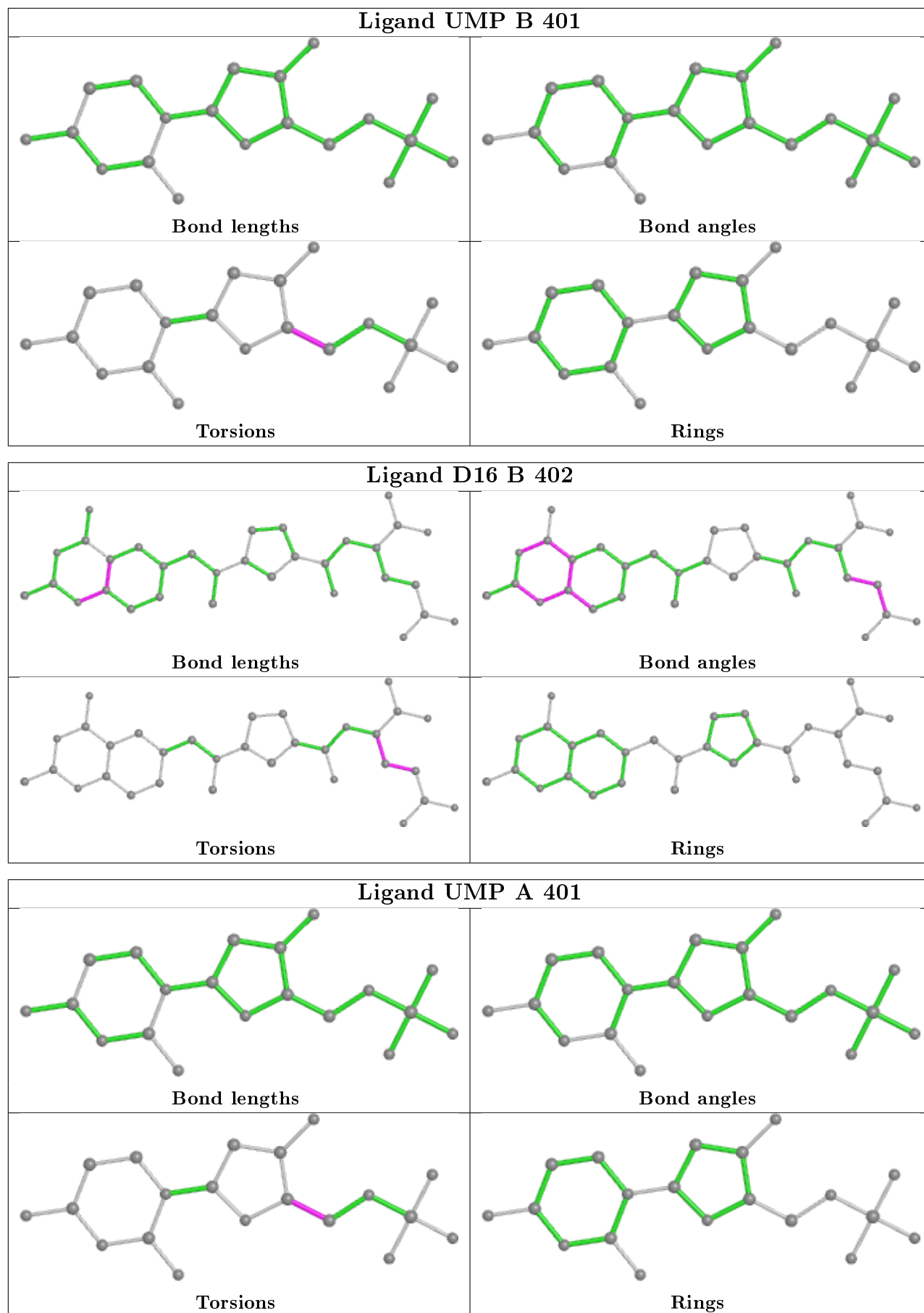
There are no ring outliers.

8 monomers are involved in 11 short contacts:

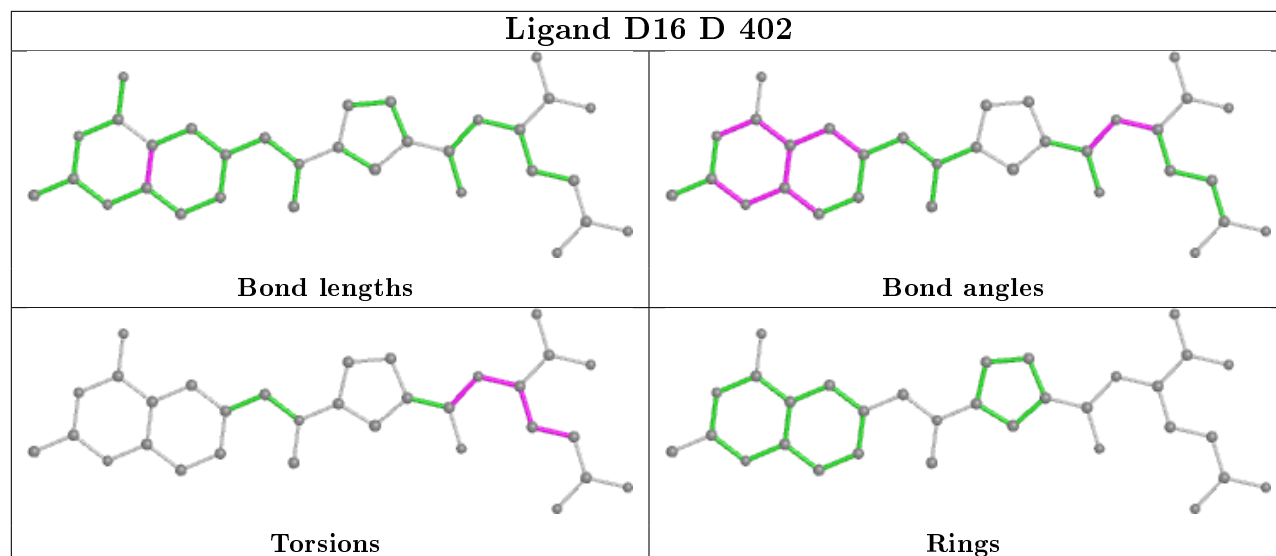
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	402	D16	2	0
2	B	401	UMP	2	0
3	B	402	D16	1	0
2	A	401	UMP	1	0
3	D	402	D16	1	0
2	C	401	UMP	1	0
3	C	402	D16	2	0
2	D	401	UMP	1	0

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.

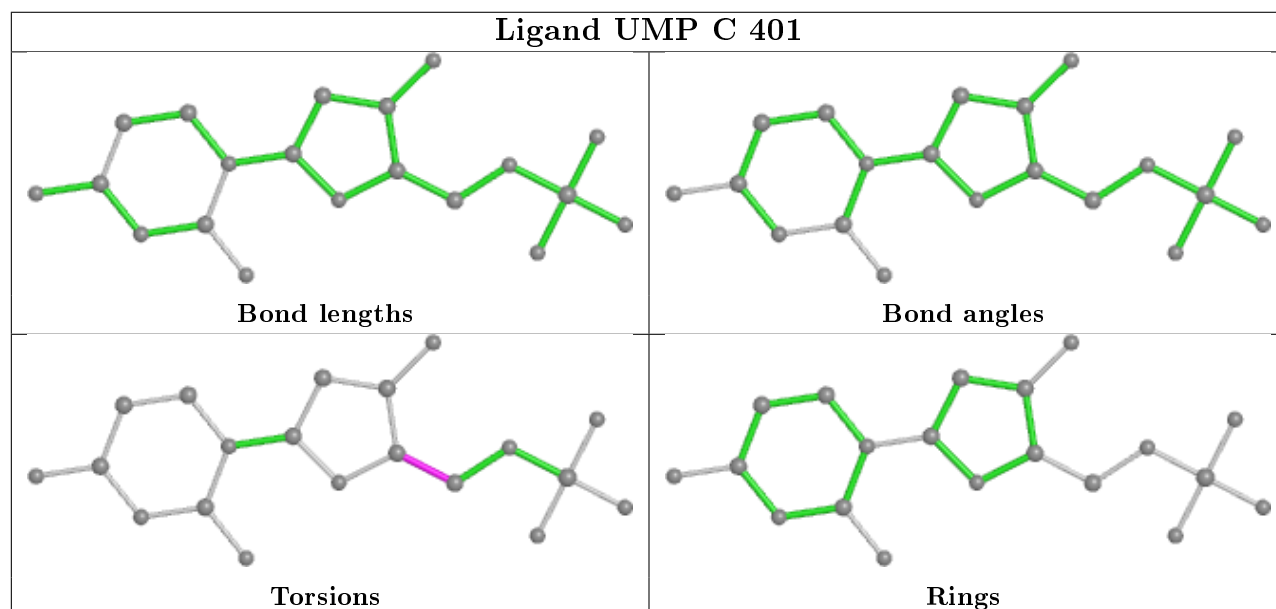




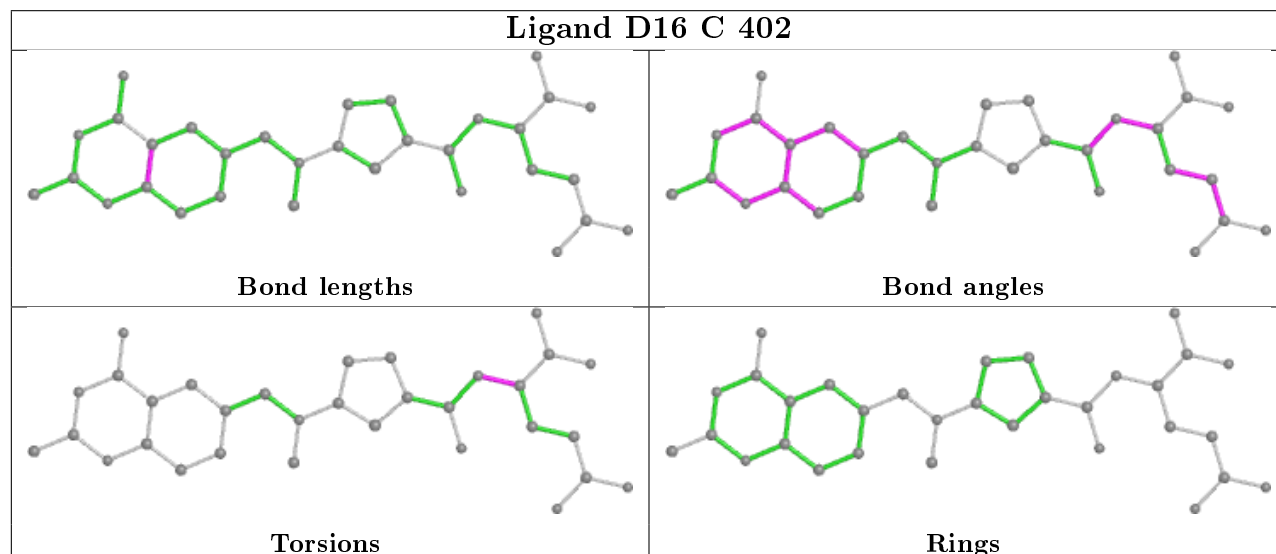
Ligand D16 D 402

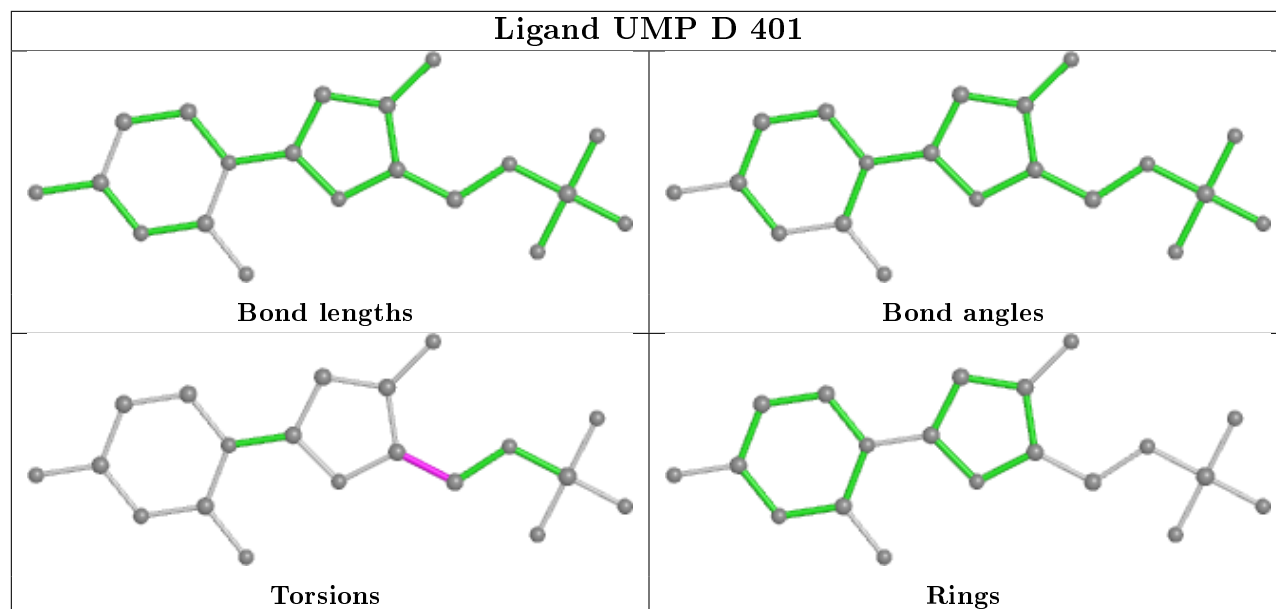


Ligand UMP C 401



Ligand D16 C 402





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	289/315 (91%)	-0.41	7 (2%) 59 56	21, 35, 56, 95	0
1	B	289/315 (91%)	-0.31	9 (3%) 49 44	17, 34, 77, 99	0
1	C	288/315 (91%)	-0.52	6 (2%) 63 61	20, 32, 45, 104	0
1	D	288/315 (91%)	-0.02	8 (2%) 53 49	29, 55, 81, 108	0
All	All	1154/1260 (91%)	-0.31	30 (2%) 56 52	17, 37, 77, 108	0

All (30) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	52	THR	6.5
1	D	25	VAL	4.7
1	A	51	ARG	4.5
1	C	52	THR	4.2
1	C	51	ARG	4.2
1	B	311	MET	4.1
1	D	52	THR	3.5
1	B	53	GLY	3.4
1	C	310	PRO	3.3
1	D	51	ARG	3.2
1	D	312	ASP	3.2
1	D	311	MET	3.0
1	A	24	GLN	2.9
1	B	312	ASP	2.9
1	A	272	ASP	2.8
1	A	312	ASP	2.7
1	A	52	THR	2.6
1	C	50	ASP	2.6
1	C	311	MET	2.5
1	B	116	ASP	2.4
1	D	310	PRO	2.4

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Mol	Chain	Res	Type	RSRZ
1	C	53	GLY	2.3
1	A	207	ASN	2.3
1	B	54	THR	2.2
1	D	69	ARG	2.1
1	B	46	THR	2.1
1	A	25	VAL	2.1
1	B	51	ARG	2.1
1	B	49	ASP	2.0
1	D	116	ASP	2.0

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

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

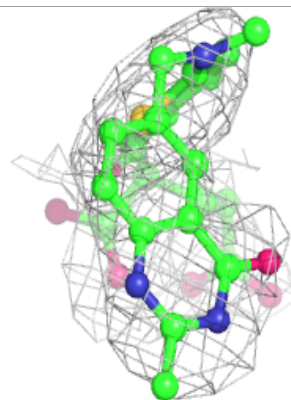
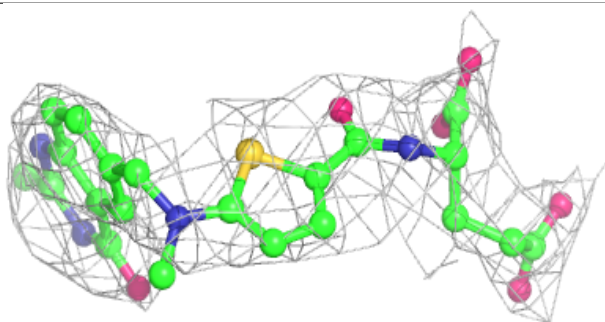
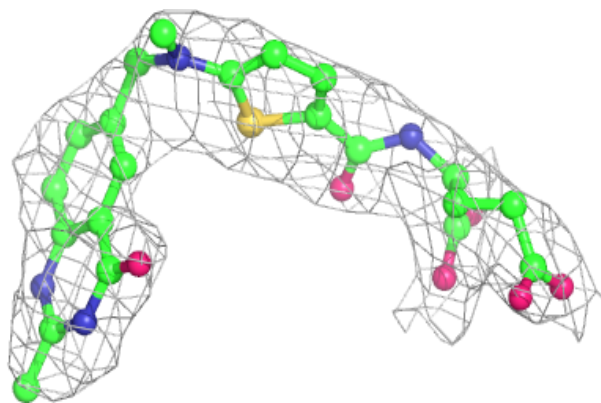
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	D16	D	402	32/32	0.89	0.23	52,58,68,69	0
3	D16	B	402	32/32	0.90	0.19	52,54,58,59	0
3	D16	C	402	32/32	0.92	0.20	37,39,42,42	0
3	D16	A	402	32/32	0.93	0.15	33,33,38,38	0
2	UMP	B	401	20/20	0.94	0.16	41,46,46,47	0
2	UMP	C	401	20/20	0.96	0.13	41,42,44,44	0
2	UMP	D	401	20/20	0.96	0.15	37,41,45,45	0
2	UMP	A	401	20/20	0.97	0.13	22,26,29,30	0

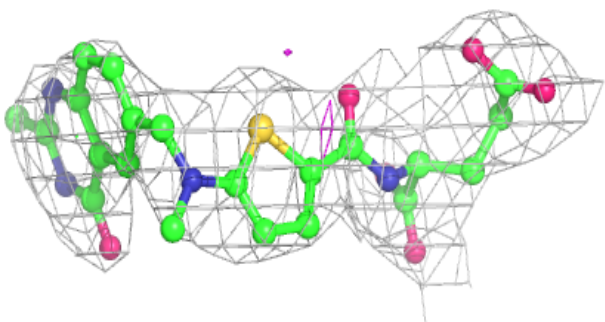
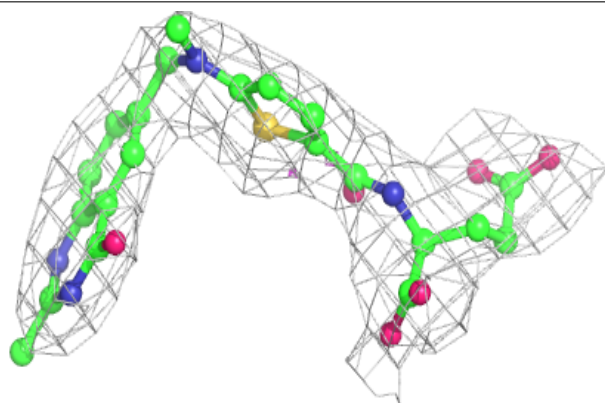
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around D16 D 402:

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

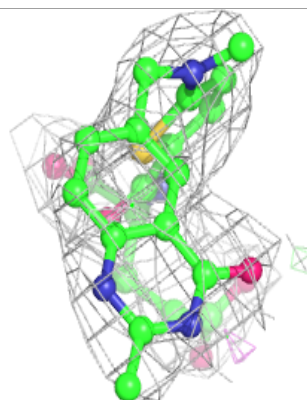
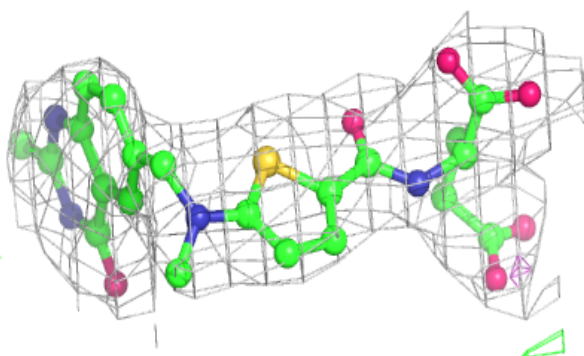
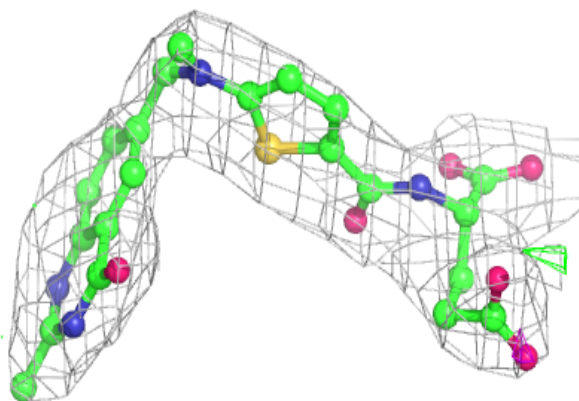
**Electron density around D16 B 402:**

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

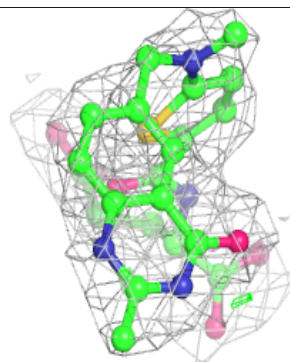
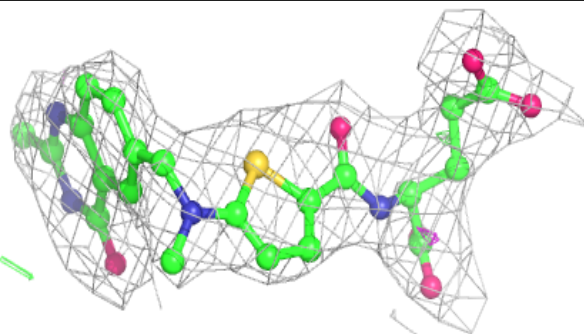
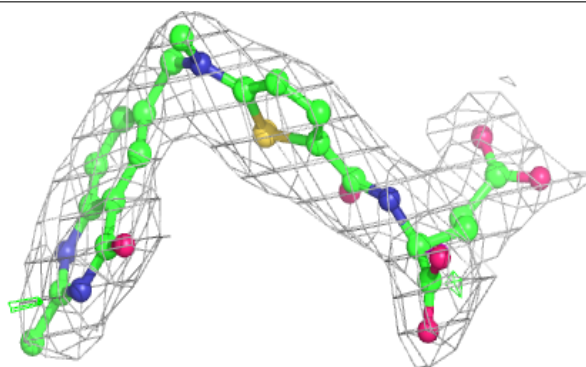


Electron density around D16 C 402:

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

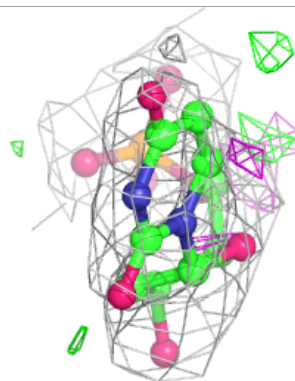
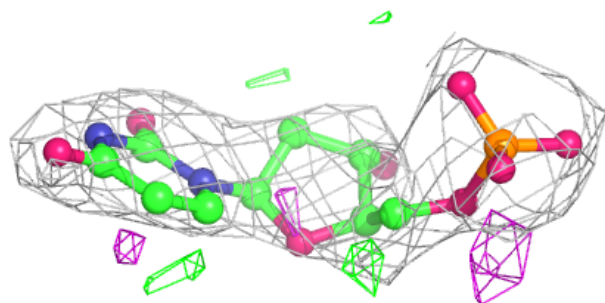
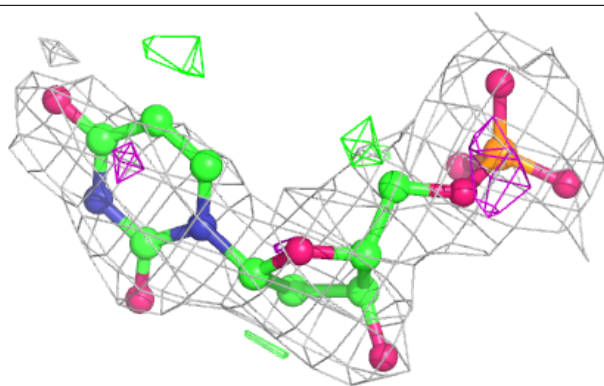
**Electron density around D16 A 402:**

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

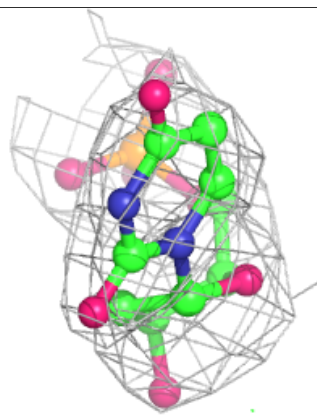
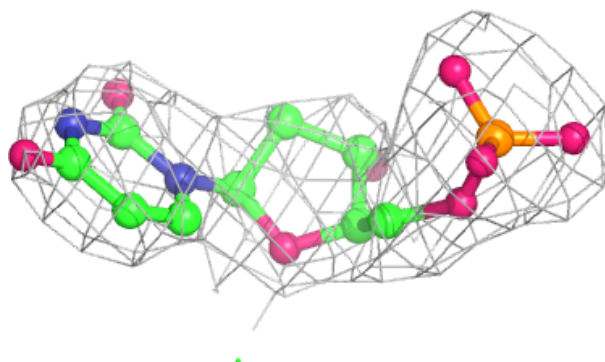
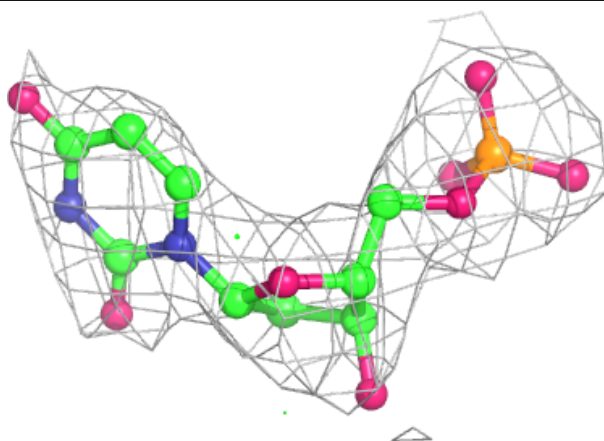


Electron density around UMP B 401:

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

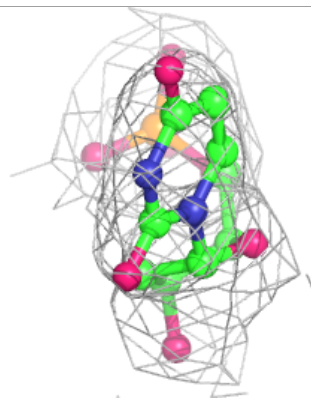
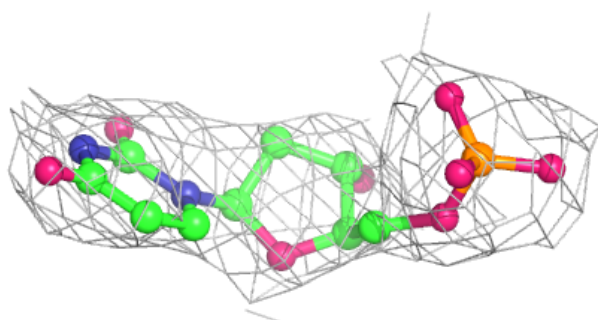
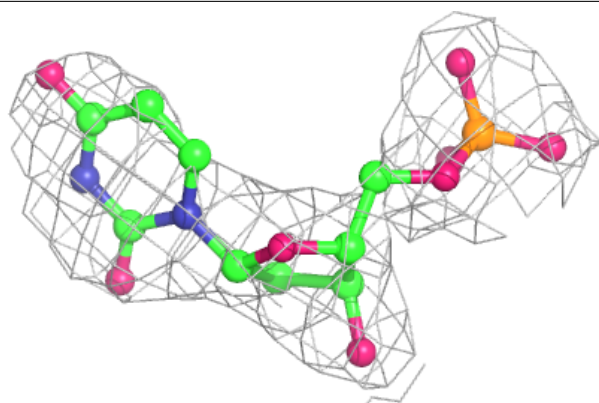
**Electron density around UMP C 401:**

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

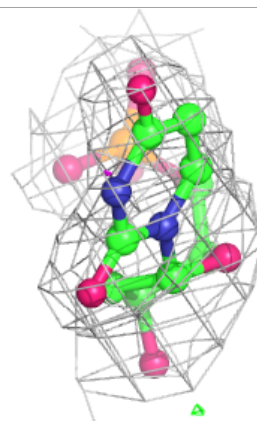
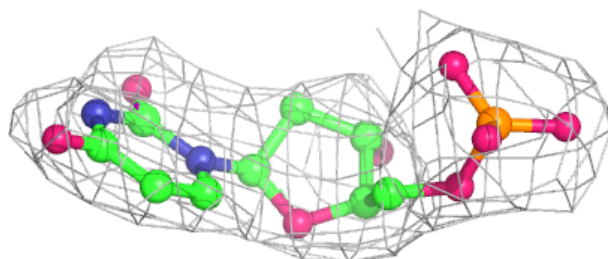


Electron density around UMP D 401:

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

**Electron density around UMP A 401:**

$2mF_o - DF_c$ (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.