



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2026 – 06:20 AM UTC

PDB ID : 4KAT / pdb_00004kat
Title : Crystal structure of FDTS from *T. maritima* mutant (R174K) with FAD and dUMP
Authors : Mathews, I.I.
Deposited on : 2013-04-22
Resolution : 2.14 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 types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

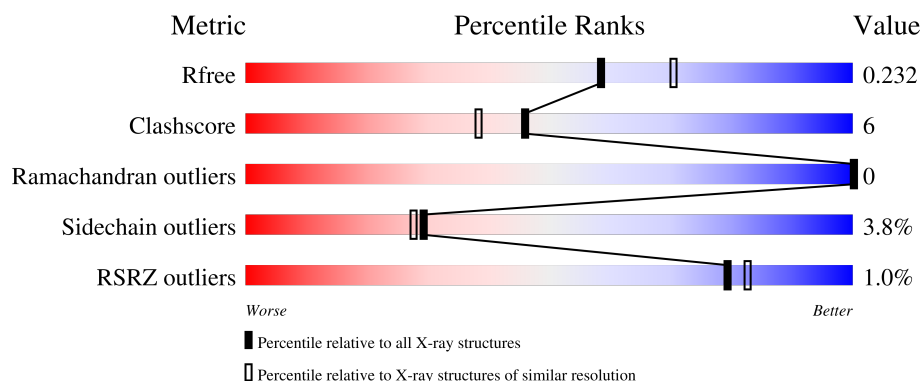
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.14 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}	180053	3689 (2.16-2.12)
Clashscore	190562	3812 (2.16-2.12)
Ramachandran outliers	187476	3773 (2.16-2.12)
Sidechain outliers	187428	3772 (2.16-2.12)
RSRZ outliers	180081	3691 (2.16-2.12)

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 $\leq 5\%$. 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	232	
1	B	232	
1	C	232	
1	D	232	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7606 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	218	Total	C	N	O	S	0	1	0
			1834	1194	316	318	6			
1	B	220	Total	C	N	O	S	0	0	0
			1841	1198	314	323	6			
1	C	213	Total	C	N	O	S	0	0	0
			1783	1161	304	313	5			
1	D	209	Total	C	N	O	S	0	0	0
			1753	1145	298	305	5			

There are 52 discrepancies between the modelled and reference sequences:

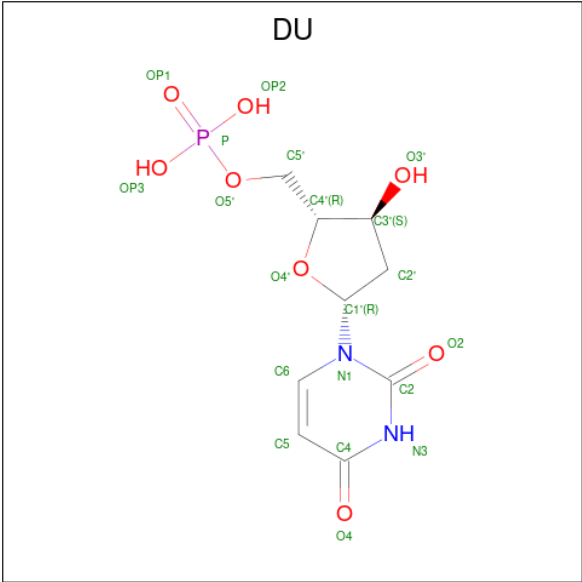
Chain	Residue	Modelled	Actual	Comment	Reference
A	-11	MET	-	initiating methionine	UNP Q9WYT0
A	-10	GLY	-	expression tag	UNP Q9WYT0
A	-9	SER	-	expression tag	UNP Q9WYT0
A	-8	ASP	-	expression tag	UNP Q9WYT0
A	-7	LYS	-	expression tag	UNP Q9WYT0
A	-6	ILE	-	expression tag	UNP Q9WYT0
A	-5	HIS	-	expression tag	UNP Q9WYT0
A	-4	HIS	-	expression tag	UNP Q9WYT0
A	-3	HIS	-	expression tag	UNP Q9WYT0
A	-2	HIS	-	expression tag	UNP Q9WYT0
A	-1	HIS	-	expression tag	UNP Q9WYT0
A	0	HIS	-	expression tag	UNP Q9WYT0
A	174	LYS	ARG	engineered mutation	UNP Q9WYT0
B	-11	MET	-	initiating methionine	UNP Q9WYT0
B	-10	GLY	-	expression tag	UNP Q9WYT0
B	-9	SER	-	expression tag	UNP Q9WYT0
B	-8	ASP	-	expression tag	UNP Q9WYT0
B	-7	LYS	-	expression tag	UNP Q9WYT0
B	-6	ILE	-	expression tag	UNP Q9WYT0
B	-5	HIS	-	expression tag	UNP Q9WYT0
B	-4	HIS	-	expression tag	UNP Q9WYT0

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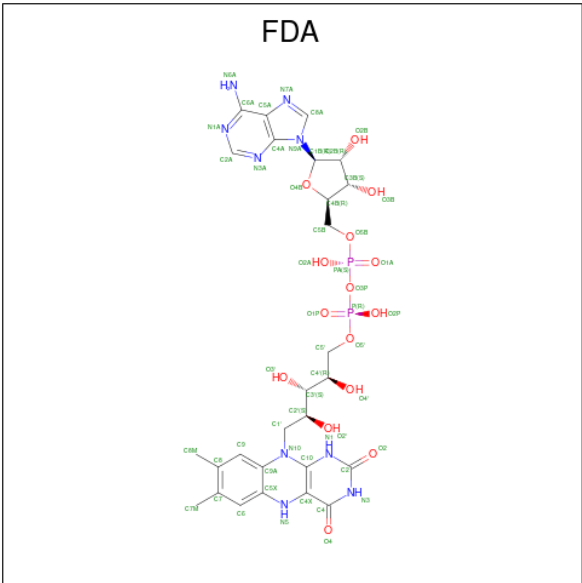
Chain	Residue	Modelled	Actual	Comment	Reference
B	-3	HIS	-	expression tag	UNP Q9WYT0
B	-2	HIS	-	expression tag	UNP Q9WYT0
B	-1	HIS	-	expression tag	UNP Q9WYT0
B	0	HIS	-	expression tag	UNP Q9WYT0
B	174	LYS	ARG	engineered mutation	UNP Q9WYT0
C	-11	MET	-	initiating methionine	UNP Q9WYT0
C	-10	GLY	-	expression tag	UNP Q9WYT0
C	-9	SER	-	expression tag	UNP Q9WYT0
C	-8	ASP	-	expression tag	UNP Q9WYT0
C	-7	LYS	-	expression tag	UNP Q9WYT0
C	-6	ILE	-	expression tag	UNP Q9WYT0
C	-5	HIS	-	expression tag	UNP Q9WYT0
C	-4	HIS	-	expression tag	UNP Q9WYT0
C	-3	HIS	-	expression tag	UNP Q9WYT0
C	-2	HIS	-	expression tag	UNP Q9WYT0
C	-1	HIS	-	expression tag	UNP Q9WYT0
C	0	HIS	-	expression tag	UNP Q9WYT0
C	174	LYS	ARG	engineered mutation	UNP Q9WYT0
D	-11	MET	-	initiating methionine	UNP Q9WYT0
D	-10	GLY	-	expression tag	UNP Q9WYT0
D	-9	SER	-	expression tag	UNP Q9WYT0
D	-8	ASP	-	expression tag	UNP Q9WYT0
D	-7	LYS	-	expression tag	UNP Q9WYT0
D	-6	ILE	-	expression tag	UNP Q9WYT0
D	-5	HIS	-	expression tag	UNP Q9WYT0
D	-4	HIS	-	expression tag	UNP Q9WYT0
D	-3	HIS	-	expression tag	UNP Q9WYT0
D	-2	HIS	-	expression tag	UNP Q9WYT0
D	-1	HIS	-	expression tag	UNP Q9WYT0
D	0	HIS	-	expression tag	UNP Q9WYT0
D	174	LYS	ARG	engineered mutation	UNP Q9WYT0

- Molecule 2 is 2'-DEOXYURIDINE-5'-MONOPHOSPHATE (CCD ID: DU) (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		
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 DIHYDROFLAVINE-ADENINE DINUCLEOTIDE (CCD ID: FDA) (formula: $C_{27}H_{35}N_9O_{15}P_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
3	B	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
3	C	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
3	D	1	Total	C	N	O	P	0	0
			53	27	9	15	2		

- Molecule 4 is water.

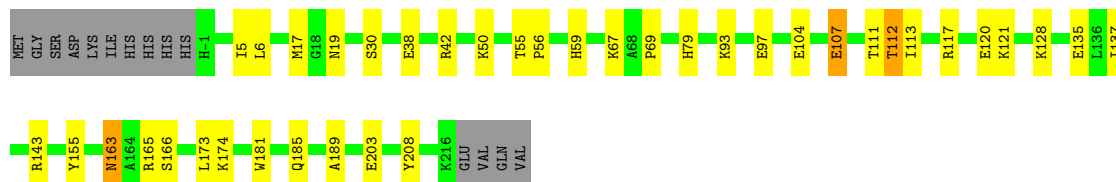
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	27	Total	O	0	0
			27	27		
4	B	22	Total	O	0	0
			22	22		
4	C	32	Total	O	0	0
			32	32		
4	D	22	Total	O	0	0
			22	22		

3 Residue-property plots

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.

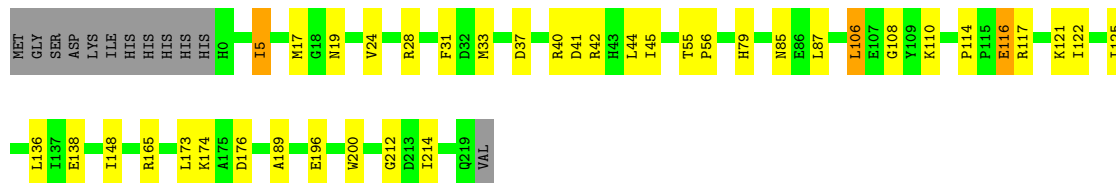
• Molecule 1: Thymidylate synthase

Chain A: 



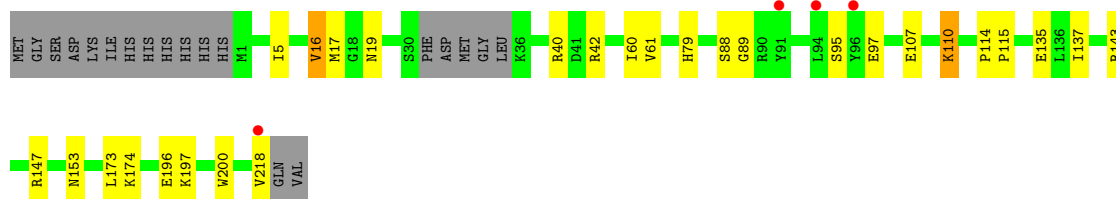
• Molecule 1: Thymidylate synthase

Chain B: 



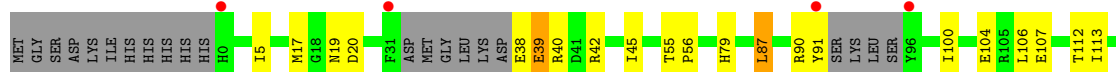
• Molecule 1: Thymidylate synthase

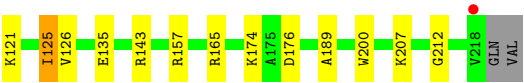
Chain C: 



• Molecule 1: Thymidylate synthase

Chain D: 





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	53.68Å 115.74Å 140.46Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	38.58 – 2.14 38.58 – 2.14	Depositor EDS
% Data completeness (in resolution range)	99.9 (38.58-2.14) 99.9 (38.58-2.14)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.01 (at 2.14Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.183 , 0.233 0.181 , 0.232	Depositor DCC
R_{free} test set	2457 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	40.0	Xtriage
Anisotropy	0.180	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 33.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	7606	wwPDB-VP
Average B, all atoms (Å ²)	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.94% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: FDA

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	0.88	0/1888	1.00	2/2551 (0.1%)
1	B	0.90	0/1891	1.03	2/2556 (0.1%)
1	C	0.92	0/1830	1.02	0/2473
1	D	0.87	0/1800	1.03	3/2433 (0.1%)
All	All	0.90	0/7409	1.02	7/10013 (0.1%)

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	C	0	1

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	108	GLY	N-CA-C	-7.02	106.13	114.48
1	B	214	ILE	N-CA-C	6.82	119.61	111.09
1	A	111	THR	CA-C-N	6.15	128.81	120.38
1	A	111	THR	C-N-CA	6.15	128.81	120.38
1	D	100	ILE	CA-C-N	6.00	126.02	119.90
1	D	100	ILE	C-N-CA	6.00	126.02	119.90
1	D	39	GLU	N-CA-C	-5.28	105.58	112.23

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	88	SER	Peptide

5.2 Too-close contacts ⓘ

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	1834	0	1834	29	0
1	B	1841	0	1837	24	0
1	C	1783	0	1785	17	0
1	D	1753	0	1744	18	0
2	A	20	0	11	1	0
2	B	20	0	11	0	0
2	C	20	0	11	3	0
2	D	20	0	11	1	0
3	A	53	0	33	2	0
3	B	53	0	33	4	0
3	C	53	0	33	3	0
3	D	53	0	33	1	0
4	A	27	0	0	1	0
4	B	22	0	0	1	0
4	C	32	0	0	2	0
4	D	22	0	0	1	0
All	All	7606	0	7376	86	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 6.

All (86) 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:79:HIS:NE2	1:A:174:LYS:NZ	2.25	0.85
1:B:31:PHE:HB2	1:B:33:MET:HE2	1.60	0.84
1:C:17:MET:HB2	1:D:17:MET:HB2	1.64	0.79
1:A:17:MET:HB2	1:B:17:MET:HB2	1.66	0.76
1:C:89:GLY:HA3	2:C:302:DU:OP3	1.87	0.75
3:B:302:FDA:N7A	4:C:401:HOH:O	2.22	0.72
1:A:5:ILE:HD11	1:A:189:ALA:HB2	1.72	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:107:GLU:HA	1:A:107:GLU:OE1	1.92	0.68
1:C:19:ASN:HB2	4:C:404:HOH:O	1.95	0.66
1:A:121:LYS:HE3	1:D:135:GLU:OE2	1.99	0.63
1:B:31:PHE:CB	1:B:33:MET:HE2	2.29	0.63
1:A:163:ASN:C	1:A:163:ASN:HD22	2.07	0.62
1:B:121:LYS:HE3	1:C:135:GLU:OE2	2.00	0.62
1:B:42:ARG:HE	1:B:200:TRP:CD1	2.19	0.61
1:B:19:ASN:HB2	4:B:416:HOH:O	2.02	0.59
1:D:91:TYR:HA	1:D:143:ARG:HH21	1.67	0.58
1:B:79:HIS:NE2	1:B:174:LYS:NZ	2.40	0.58
1:D:90:ARG:HG3	2:D:302:DU:OP2	2.04	0.58
1:B:165:ARG:CZ	3:D:301:FDA:H2B	2.35	0.57
1:C:218:VAL:HG12	1:C:218:VAL:O	2.04	0.57
1:D:42:ARG:HG2	1:D:200:TRP:CD2	2.40	0.56
1:C:16:VAL:HG11	1:C:197:LYS:HD2	1.88	0.56
1:D:19:ASN:HB2	4:D:422:HOH:O	2.05	0.56
2:C:302:DU:H6	2:C:302:DU:O5'	2.07	0.55
1:A:137:ILE:HD11	1:A:143:ARG:HA	1.88	0.55
1:B:28:ARG:HD3	1:B:33:MET:HE3	1.88	0.54
1:A:135:GLU:OE2	1:D:121:LYS:NZ	2.41	0.53
1:A:42:ARG:NH2	1:A:203:GLU:OE1	2.41	0.53
1:D:104:GLU:HA	1:D:107:GLU:HG2	1.91	0.51
1:B:114:PRO:HG2	1:B:117:ARG:HG2	1.93	0.51
1:A:113:ILE:HD12	1:A:117:ARG:CB	2.41	0.51
1:A:93:LYS:HB2	1:A:143:ARG:NH1	2.26	0.50
1:D:106:LEU:HD22	1:D:106:LEU:N	2.27	0.49
1:B:122:ILE:O	1:B:125:ILE:HG22	2.12	0.49
1:D:79:HIS:CD2	1:D:174:LYS:HZ3	2.27	0.49
1:A:67:LYS:NZ	1:A:97:GLU:OE2	2.44	0.49
1:C:79:HIS:NE2	1:C:174:LYS:NZ	2.44	0.49
1:A:5:ILE:HG22	1:A:6:LEU:HG	1.94	0.49
1:A:104:GLU:O	1:A:107:GLU:HG2	2.13	0.49
1:D:176:ASP:OD1	1:D:212:GLY:HA2	2.13	0.48
1:A:55:THR:OG1	1:A:56:PRO:HD3	2.13	0.48
1:A:113:ILE:HD12	1:A:117:ARG:HB2	1.95	0.48
1:A:112:THR:HG22	1:A:113:ILE:HG23	1.96	0.48
1:B:117:ARG:HH21	1:B:121:LYS:NZ	2.12	0.48
1:C:173:LEU:HD13	3:C:301:FDA:O4'	2.14	0.48
1:A:163:ASN:ND2	1:A:166:SER:H	2.13	0.47
1:B:5:ILE:HD11	1:B:189:ALA:HB2	1.97	0.47
1:A:107:GLU:OE1	1:A:107:GLU:CA	2.63	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:302:FDA:H2B	1:D:165:ARG:CZ	2.45	0.46
1:A:181:TRP:O	1:A:185:GLN:HG2	2.16	0.46
1:A:163:ASN:C	1:A:163:ASN:ND2	2.74	0.46
1:B:148:ILE:HB	1:C:153:ASN:HB3	1.96	0.46
1:D:20:ASP:HB3	1:D:45:ILE:CD1	2.45	0.46
1:A:19:ASN:HB2	4:A:419:HOH:O	2.15	0.46
1:B:176:ASP:OD1	1:B:212:GLY:HA2	2.16	0.45
1:D:87:LEU:HD22	1:D:157:ARG:HB2	1.98	0.45
1:D:5:ILE:HD11	1:D:189:ALA:HB2	1.99	0.45
1:B:173:LEU:HB3	3:B:302:FDA:HM83	1.98	0.44
1:C:110:LYS:HB2	1:C:110:LYS:HE2	1.68	0.44
1:B:114:PRO:HG2	1:B:117:ARG:CG	2.48	0.43
1:B:44:LEU:O	1:B:45:ILE:C	2.62	0.43
2:C:302:DU:H6	2:C:302:DU:C5'	2.48	0.43
1:A:50:LYS:HE3	1:A:208:TYR:CZ	2.54	0.43
1:C:218:VAL:O	1:C:218:VAL:CG1	2.67	0.43
3:A:302:FDA:H1'2	3:A:302:FDA:H9	1.77	0.43
1:C:143:ARG:O	1:C:147:ARG:HG3	2.20	0.42
1:A:59:HIS:NE2	1:B:85:ASN:ND2	2.62	0.42
1:A:69:PRO:HD3	1:A:155:TYR:CE1	2.54	0.42
1:B:55:THR:N	1:B:56:PRO:CD	2.82	0.42
1:B:116:GLU:H	1:B:116:GLU:CD	2.27	0.42
1:C:42:ARG:HG2	1:C:200:TRP:CD2	2.55	0.42
1:C:137:ILE:HD13	1:C:137:ILE:HA	1.83	0.42
1:D:125:ILE:HD13	1:D:125:ILE:HA	1.78	0.42
1:B:136:LEU:HD23	1:B:136:LEU:HA	1.86	0.42
1:A:165:ARG:CZ	3:C:301:FDA:H2B	2.50	0.42
1:C:114:PRO:HA	1:C:115:PRO:HD2	1.76	0.42
3:B:302:FDA:H9	3:B:302:FDA:H1'2	1.82	0.41
1:B:24:VAL:HG21	1:B:41:ASP:HB3	2.01	0.41
1:C:95:SER:HB3	1:C:97:GLU:HG3	2.03	0.41
1:A:173:LEU:HD13	3:A:302:FDA:O4'	2.20	0.41
1:D:55:THR:OG1	1:D:56:PRO:HD3	2.21	0.41
1:A:30:SER:HG	3:C:301:FDA:C4	2.32	0.41
1:A:128:LYS:HG3	1:D:125:ILE:HD11	2.03	0.41
2:A:301:DU:H6	2:A:301:DU:O5'	2.21	0.41
1:B:106:LEU:HD13	1:B:106:LEU:HA	1.87	0.41
1:C:60:ILE:C	1:C:61:VAL:HG23	2.45	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	217/232 (94%)	208 (96%)	9 (4%)	0	100	100
1	B	218/232 (94%)	212 (97%)	6 (3%)	0	100	100
1	C	209/232 (90%)	201 (96%)	8 (4%)	0	100	100
1	D	203/232 (88%)	199 (98%)	4 (2%)	0	100	100
All	All	847/928 (91%)	820 (97%)	27 (3%)	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	195/207 (94%)	190 (97%)	5 (3%)	40	42
1	B	196/207 (95%)	187 (95%)	9 (5%)	24	20
1	C	190/207 (92%)	184 (97%)	6 (3%)	34	35
1	D	185/207 (89%)	176 (95%)	9 (5%)	22	18
All	All	766/828 (92%)	737 (96%)	29 (4%)	29	28

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	38	GLU
1	A	107	GLU

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Mol	Chain	Res	Type
1	A	112	THR
1	A	120	GLU
1	A	163	ASN
1	B	5	ILE
1	B	37	ASP
1	B	40	ARG
1	B	87	LEU
1	B	106	LEU
1	B	110	LYS
1	B	116	GLU
1	B	138	GLU
1	B	196	GLU
1	C	5	ILE
1	C	16	VAL
1	C	40	ARG
1	C	107	GLU
1	C	110	LYS
1	C	196	GLU
1	D	38	GLU
1	D	39	GLU
1	D	40	ARG
1	D	87	LEU
1	D	112	THR
1	D	113	ILE
1	D	125	ILE
1	D	126	VAL
1	D	207	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	51	HIS
1	A	75	GLN
1	A	163	ASN
1	B	53	HIS
1	C	51	HIS
1	C	185	GLN
1	D	85	ASN
1	D	169	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

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	FDA	B	302	-	57,58,58	3.33	18 (31%)	78,89,89	1.87	18 (23%)
3	FDA	A	302	-	57,58,58	3.30	16 (28%)	78,89,89	1.82	17 (21%)
2	DU	C	302	-	21,21,21	1.21	2 (9%)	30,31,31	2.30	9 (30%)
3	FDA	C	301	-	57,58,58	3.29	19 (33%)	78,89,89	1.82	19 (24%)
2	DU	D	302	-	21,21,21	1.26	3 (14%)	30,31,31	1.83	6 (20%)
2	DU	B	301	-	21,21,21	1.22	3 (14%)	30,31,31	1.68	6 (20%)
3	FDA	D	301	-	57,58,58	3.23	18 (31%)	78,89,89	2.03	26 (33%)
2	DU	A	301	-	21,21,21	1.19	2 (9%)	30,31,31	2.08	6 (20%)

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	FDA	B	302	-	-	4/34/50/50	0/6/6/6
3	FDA	A	302	-	-	3/34/50/50	0/6/6/6
2	DU	C	302	-	-	0/10/22/22	0/2/2/2
3	FDA	C	301	-	-	4/34/50/50	0/6/6/6
2	DU	D	302	-	-	0/10/22/22	0/2/2/2
2	DU	B	301	-	-	0/10/22/22	0/2/2/2
3	FDA	D	301	-	-	7/34/50/50	0/6/6/6
2	DU	A	301	-	-	3/10/22/22	0/2/2/2

All (81) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	302	FDA	C5X-C9A	11.75	1.53	1.40
3	A	302	FDA	C5X-C9A	11.43	1.53	1.40
3	C	301	FDA	C5X-C9A	11.31	1.53	1.40
3	D	301	FDA	C5X-C9A	11.02	1.52	1.40
3	A	302	FDA	C6-C7	9.53	1.52	1.39
3	B	302	FDA	C6-C7	9.49	1.52	1.39
3	D	301	FDA	C6-C7	9.31	1.52	1.39
3	B	302	FDA	C6-C5X	9.27	1.54	1.39
3	C	301	FDA	C6-C7	9.10	1.52	1.39
3	D	301	FDA	C6-C5X	8.85	1.53	1.39
3	A	302	FDA	C6-C5X	8.64	1.53	1.39
3	C	301	FDA	C6-C5X	8.58	1.53	1.39
3	D	301	FDA	C9A-N10	-6.61	1.29	1.41
3	B	302	FDA	C9A-N10	-6.50	1.29	1.41
3	C	301	FDA	C9A-N10	-6.45	1.30	1.41
3	C	301	FDA	C5A-C4A	6.39	1.50	1.39
3	A	302	FDA	C10-N1	6.22	1.48	1.37
3	D	301	FDA	C9-C9A	6.15	1.49	1.39
3	B	302	FDA	C9-C9A	6.12	1.49	1.39
3	A	302	FDA	C9A-N10	-6.12	1.30	1.41
3	B	302	FDA	C10-N1	6.02	1.47	1.37
3	C	301	FDA	C2-N1	-6.00	1.27	1.37
3	A	302	FDA	C9-C9A	5.98	1.49	1.39
3	C	301	FDA	C10-N1	5.96	1.47	1.37
3	A	302	FDA	C5A-C4A	5.92	1.49	1.39
3	D	301	FDA	C5A-C4A	5.89	1.49	1.39
3	A	302	FDA	C2-N1	-5.77	1.28	1.37
3	D	301	FDA	C2-N1	-5.74	1.28	1.37
3	B	302	FDA	C5A-C4A	5.62	1.49	1.39
3	C	301	FDA	C9-C9A	5.59	1.48	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	302	FDA	C10-N10	5.25	1.47	1.38
3	D	301	FDA	C10-N1	5.21	1.46	1.37
3	B	302	FDA	C10-N10	5.19	1.47	1.38
3	B	302	FDA	C2-N1	-5.05	1.29	1.37
3	C	301	FDA	C10-N10	4.90	1.46	1.38
3	A	302	FDA	C4X-N5	4.86	1.45	1.35
3	B	302	FDA	C8-C7	4.71	1.52	1.40
3	C	301	FDA	C8-C7	4.70	1.52	1.40
3	A	302	FDA	C8-C7	4.64	1.52	1.40
3	C	301	FDA	C4X-N5	4.60	1.45	1.35
3	D	301	FDA	C10-N10	4.51	1.46	1.38
3	D	301	FDA	C8-C7	4.29	1.51	1.40
3	B	302	FDA	C4X-N5	4.26	1.44	1.35
3	D	301	FDA	C4X-N5	3.84	1.43	1.35
3	C	301	FDA	C5X-N5	3.54	1.45	1.39
3	A	302	FDA	C5X-N5	3.54	1.45	1.39
3	B	302	FDA	C5X-N5	3.40	1.45	1.39
3	A	302	FDA	C9-C8	-3.35	1.35	1.39
3	D	301	FDA	C4-N3	-3.08	1.33	1.38
2	D	302	DU	C2-N1	2.93	1.43	1.38
2	D	302	DU	C6-C5	2.90	1.41	1.35
2	C	302	DU	C2-N1	2.84	1.42	1.38
3	A	302	FDA	C4X-C4	2.84	1.50	1.41
3	D	301	FDA	O2-C2	-2.83	1.17	1.23
2	C	302	DU	C5-C4	-2.82	1.37	1.43
3	C	301	FDA	PA-O1A	2.80	1.60	1.50
3	B	302	FDA	C4X-C4	2.79	1.50	1.41
3	D	301	FDA	C4X-C4	2.76	1.50	1.41
3	C	301	FDA	C4X-C4	2.73	1.50	1.41
3	D	301	FDA	C5X-N5	2.70	1.44	1.39
2	D	302	DU	C4-N3	-2.58	1.34	1.38
3	B	302	FDA	C9-C8	-2.55	1.36	1.39
3	C	301	FDA	C4-N3	-2.48	1.34	1.38
2	B	301	DU	O2-C2	2.48	1.27	1.23
3	C	301	FDA	C8A-N7A	2.40	1.36	1.31
3	B	302	FDA	C4-N3	-2.39	1.34	1.38
3	B	302	FDA	P-O3P	2.39	1.62	1.59
2	B	301	DU	C6-C5	2.37	1.40	1.35
3	C	301	FDA	C9-C8	-2.36	1.36	1.39
3	D	301	FDA	C9-C8	-2.30	1.36	1.39
3	D	301	FDA	PA-O1A	2.25	1.58	1.50
2	A	301	DU	C5-C4	-2.24	1.38	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	302	FDA	PA-O1A	2.21	1.58	1.50
3	C	301	FDA	O2-C2	-2.20	1.18	1.23
2	B	301	DU	C2-N1	2.17	1.41	1.38
3	B	302	FDA	PA-O1A	2.17	1.58	1.50
3	D	301	FDA	PA-O3P	2.15	1.61	1.59
2	A	301	DU	C2-N1	2.14	1.41	1.38
3	C	301	FDA	C5A-N7A	-2.06	1.35	1.39
3	A	302	FDA	C4A-N9A	-2.02	1.33	1.37
3	B	302	FDA	PA-O3P	2.01	1.61	1.59

All (107) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	301	FDA	N3-C2-N1	6.87	126.54	115.74
3	D	301	FDA	C5A-C4A-N3A	-6.14	118.27	126.72
3	B	302	FDA	C5A-C4A-N3A	-5.80	118.74	126.72
2	C	302	DU	O4-C4-C5	-5.78	115.19	125.16
3	C	301	FDA	N3-C2-N1	5.72	124.73	115.74
2	A	301	DU	C4-N3-C2	-5.44	119.85	126.61
3	C	301	FDA	C5A-C4A-N3A	-5.28	119.45	126.72
2	A	301	DU	O4-C4-C5	-5.21	116.18	125.16
2	C	302	DU	C4-N3-C2	-5.12	120.26	126.61
3	B	302	FDA	N3-C2-N1	5.07	123.71	115.74
3	D	301	FDA	N3A-C4A-N9A	5.05	135.76	127.17
3	B	302	FDA	N3A-C4A-N9A	4.95	135.58	127.17
2	D	302	DU	N3-C2-N1	4.93	121.31	114.89
2	D	302	DU	C4-N3-C2	-4.91	120.52	126.61
2	A	301	DU	C5-C4-N3	4.90	121.67	114.80
2	C	302	DU	C5-C4-N3	4.76	121.47	114.80
3	A	302	FDA	C5A-C4A-N3A	-4.65	120.32	126.72
3	C	301	FDA	N3A-C4A-N9A	4.59	134.98	127.17
3	A	302	FDA	N3-C2-N1	4.53	122.86	115.74
3	A	302	FDA	N3A-C2A-N1A	-4.46	121.83	128.58
2	C	302	DU	C1'-N1-C2	4.37	126.21	117.66
2	B	301	DU	C5-C4-N3	4.21	120.70	114.80
3	A	302	FDA	N3A-C4A-N9A	4.18	134.28	127.17
3	A	302	FDA	C5X-C9A-N10	4.13	122.69	118.01
2	D	302	DU	C5-C4-N3	4.11	120.56	114.80
3	B	302	FDA	C4A-C5A-N7A	-3.85	106.18	110.58
2	B	301	DU	O4-C4-C5	-3.71	118.76	125.16
3	B	302	FDA	C5X-C9A-N10	3.69	122.18	118.01
2	A	301	DU	N3-C2-N1	3.57	119.53	114.89

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	301	FDA	N3A-C2A-N1A	-3.48	123.31	128.58
2	B	301	DU	C4-N3-C2	-3.47	122.31	126.61
2	C	302	DU	N3-C2-N1	3.45	119.38	114.89
3	A	302	FDA	C4A-C5A-N7A	-3.45	106.64	110.58
3	C	301	FDA	N3A-C2A-N1A	-3.41	123.42	128.58
3	C	301	FDA	O2A-PA-O1A	3.39	128.19	112.44
3	D	301	FDA	C4A-C5A-N7A	-3.38	106.72	110.58
3	A	302	FDA	C4A-N9A-C8A	3.37	109.28	105.74
3	C	301	FDA	C5X-C9A-N10	3.36	121.81	118.01
3	A	302	FDA	C2A-N1A-C6A	3.31	124.17	118.73
3	B	302	FDA	C4-C4X-N5	3.25	124.69	116.37
3	D	301	FDA	C2A-N3A-C4A	3.23	119.72	111.83
3	A	302	FDA	C4-C4X-N5	3.19	124.53	116.37
2	C	302	DU	C1'-N1-C6	-3.17	115.29	121.53
3	B	302	FDA	N3A-C2A-N1A	-3.17	123.78	128.58
3	B	302	FDA	C2A-N3A-C4A	3.11	119.42	111.83
3	D	301	FDA	O2-C2-N3	-3.09	116.37	121.86
3	C	301	FDA	C4-C4X-N5	3.05	124.16	116.37
3	C	301	FDA	C4A-C5A-N7A	-3.00	107.15	110.58
3	D	301	FDA	O4B-C1B-N9A	3.00	113.85	108.09
3	D	301	FDA	C5X-C9A-N10	2.98	121.38	118.01
3	B	302	FDA	C5A-N7A-C8A	2.97	108.12	103.45
3	A	302	FDA	C2A-N3A-C4A	2.94	119.02	111.83
3	D	301	FDA	C4-N3-C2	-2.92	122.35	126.37
3	D	301	FDA	C4-C4X-N5	2.88	123.74	116.37
3	B	302	FDA	C4A-N9A-C8A	2.86	108.74	105.74
2	B	301	DU	O5'-P-OP1	-2.84	98.76	106.44
3	A	302	FDA	O4-C4-C4X	-2.83	120.43	127.26
3	D	301	FDA	O2A-PA-O1A	2.82	125.54	112.44
3	D	301	FDA	C3B-C2B-C1B	2.80	106.76	101.46
2	C	302	DU	O4-C4-N3	2.80	123.33	119.27
3	C	301	FDA	C2A-N1A-C6A	2.78	123.29	118.73
2	A	301	DU	O2-C2-N1	-2.76	119.21	122.80
3	B	302	FDA	O2-C2-N3	-2.75	116.97	121.86
3	D	301	FDA	O2-C2-N1	-2.69	117.08	121.86
3	B	302	FDA	C3B-C2B-C1B	2.68	106.54	101.46
3	A	302	FDA	C9A-C9-C8	2.67	124.59	119.22
2	B	301	DU	N3-C2-N1	2.65	118.35	114.89
3	D	301	FDA	O5'-C5'-C4'	-2.61	102.40	109.36
2	C	302	DU	OP3-P-O5'	-2.59	99.92	106.67
3	C	301	FDA	O2-C2-N3	-2.58	117.27	121.86
3	C	301	FDA	C2A-N3A-C4A	2.58	118.13	111.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	301	FDA	O4-C4-C4X	-2.57	121.06	127.26
3	D	301	FDA	O2B-C2B-C1B	-2.56	101.28	110.10
3	C	301	FDA	O4-C4-C4X	-2.53	121.17	127.26
3	C	301	FDA	C4A-N9A-C8A	2.52	108.39	105.74
2	D	302	DU	O4-C4-N3	-2.49	115.66	119.27
3	B	302	FDA	O5'-C5'-C4'	-2.45	102.83	109.36
3	B	302	FDA	C4X-C4-N3	2.43	118.80	112.13
3	A	302	FDA	O2A-PA-O1A	2.41	123.66	112.44
2	D	302	DU	O4'-C1'-N1	2.39	112.10	107.86
3	A	302	FDA	C5A-N7A-C8A	2.35	107.15	103.45
3	B	302	FDA	O4-C4-C4X	-2.34	121.61	127.26
3	D	301	FDA	C9-C9A-C5X	-2.29	116.75	119.95
3	B	302	FDA	C9-C8-C7	2.24	122.97	119.69
3	D	301	FDA	C4A-N9A-C8A	2.23	108.08	105.74
3	D	301	FDA	C2'-C1'-N10	2.22	120.70	110.20
3	B	302	FDA	C2B-C1B-N9A	2.22	118.82	113.30
3	C	301	FDA	C3B-C2B-C1B	2.20	105.63	101.46
2	D	302	DU	O2-C2-N3	-2.20	117.44	121.49
3	D	301	FDA	C4X-C4-N3	2.19	118.16	112.13
3	C	301	FDA	O2-C2-N1	-2.19	117.97	121.86
3	A	302	FDA	C8M-C8-C9	-2.19	115.72	119.57
3	A	302	FDA	C4X-C4-N3	2.17	118.09	112.13
3	A	302	FDA	C2B-C1B-N9A	2.16	118.67	113.30
3	D	301	FDA	O2P-P-O1P	2.15	122.44	112.44
3	C	301	FDA	C4X-C4-N3	2.14	118.02	112.13
3	C	301	FDA	O2A-PA-O3P	-2.14	101.48	107.27
3	D	301	FDA	O3'-C3'-C2'	2.14	113.78	108.93
3	C	301	FDA	C9A-C9-C8	2.08	123.39	119.22
2	A	301	DU	C2'-C1'-N1	-2.07	108.65	113.81
3	D	301	FDA	C5A-N7A-C8A	2.06	106.69	103.45
2	C	302	DU	C6-N1-C2	-2.05	118.50	121.00
3	D	301	FDA	C9-C8-C7	2.04	122.68	119.69
3	D	301	FDA	C9A-C9-C8	2.03	123.31	119.22
3	B	302	FDA	C9A-C9-C8	2.03	123.30	119.22
2	B	301	DU	C2'-C1'-N1	-2.02	108.77	113.81
3	C	301	FDA	C5X-N5-C4X	-2.02	116.42	121.08

There are no chirality outliers.

All (21) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	301	DU	C5'-O5'-P-OP1

Continued on next page...

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Mol	Chain	Res	Type	Atoms
3	A	302	FDA	C3'-C4'-C5'-O5'
3	A	302	FDA	O4'-C4'-C5'-O5'
3	B	302	FDA	C3'-C4'-C5'-O5'
3	B	302	FDA	O4'-C4'-C5'-O5'
3	C	301	FDA	P-O3P-PA-O5B
3	C	301	FDA	O4'-C4'-C5'-O5'
3	D	301	FDA	O4'-C4'-C5'-O5'
3	D	301	FDA	C3'-C4'-C5'-O5'
3	D	301	FDA	P-O3P-PA-O5B
3	D	301	FDA	C2'-C3'-C4'-O4'
3	A	302	FDA	C2'-C3'-C4'-O4'
3	D	301	FDA	C2'-C3'-C4'-C5'
3	C	301	FDA	C5'-O5'-P-O1P
3	D	301	FDA	C5'-O5'-P-O1P
3	B	302	FDA	O4B-C4B-C5B-O5B
2	A	301	DU	C5'-O5'-P-OP3
2	A	301	DU	C5'-O5'-P-OP2
3	D	301	FDA	O3'-C3'-C4'-C5'
3	B	302	FDA	PA-O3P-P-O2P
3	C	301	FDA	PA-O3P-P-O2P

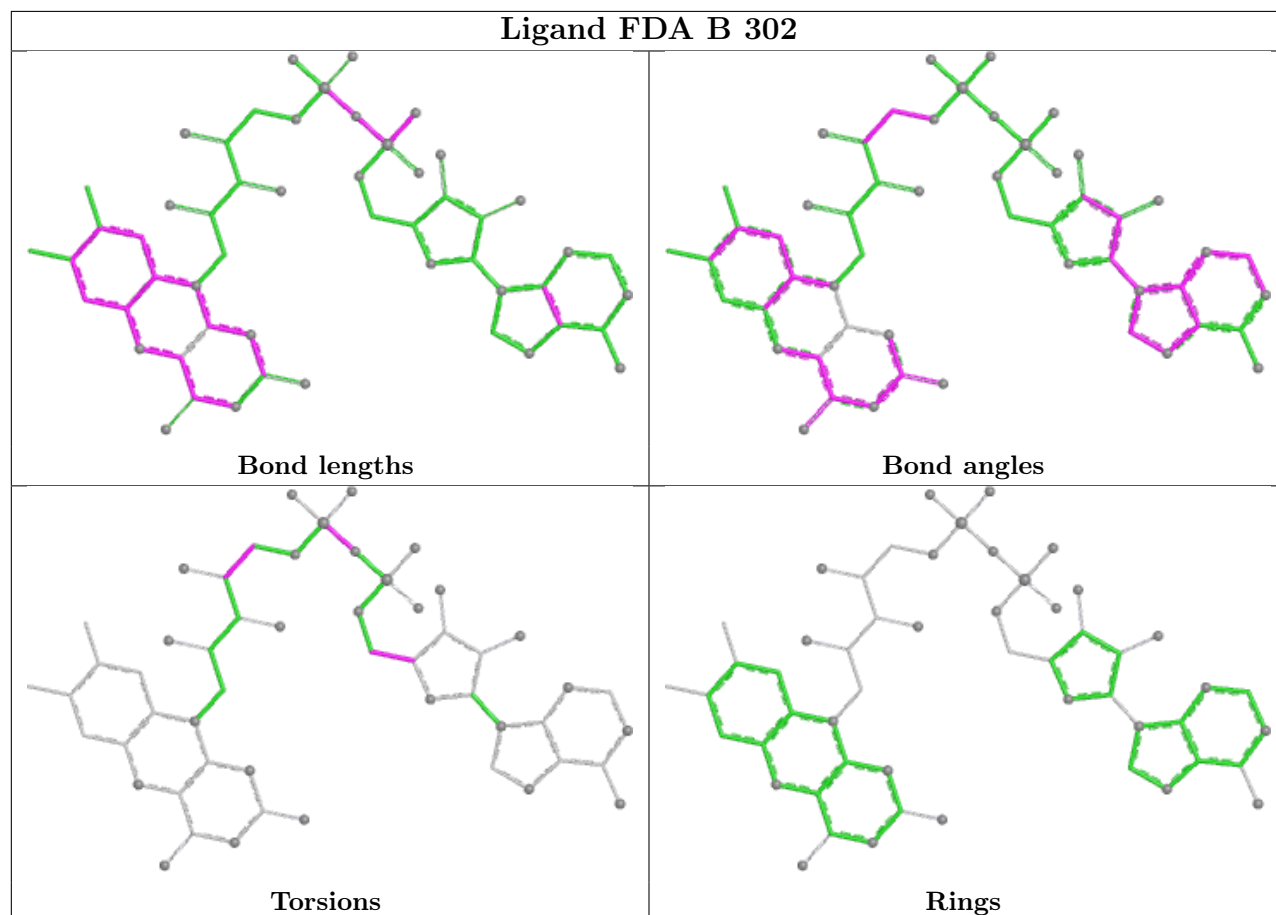
There are no ring outliers.

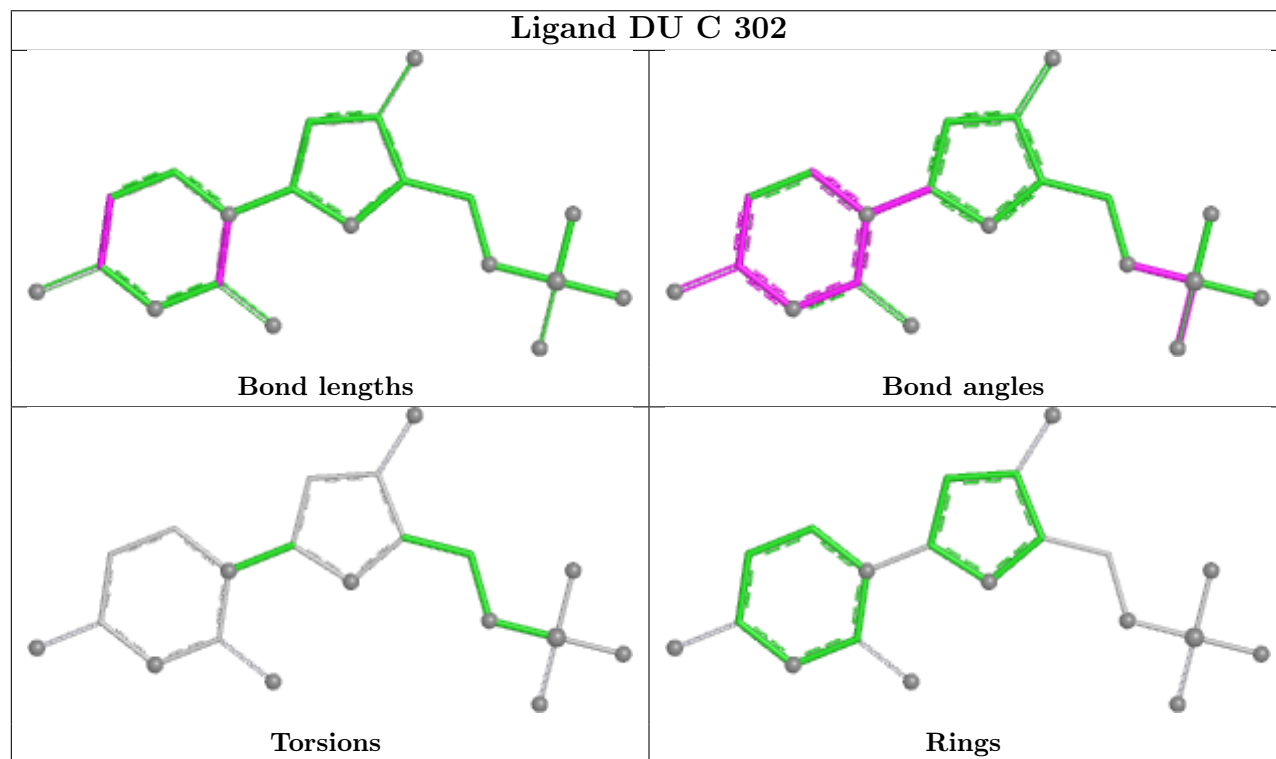
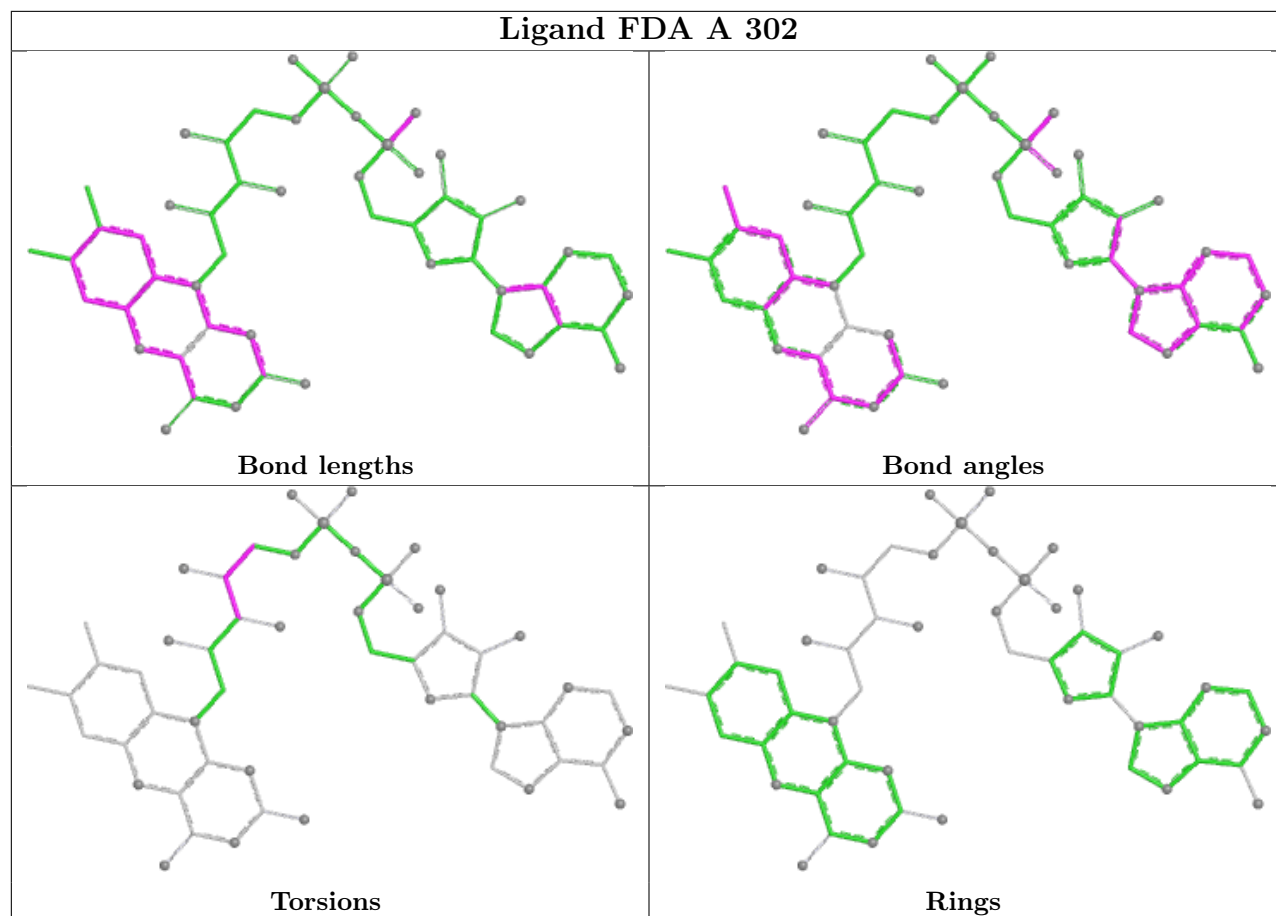
7 monomers are involved in 15 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	302	FDA	4	0
3	A	302	FDA	2	0
2	C	302	DU	3	0
3	C	301	FDA	3	0
2	D	302	DU	1	0
3	D	301	FDA	1	0
2	A	301	DU	1	0

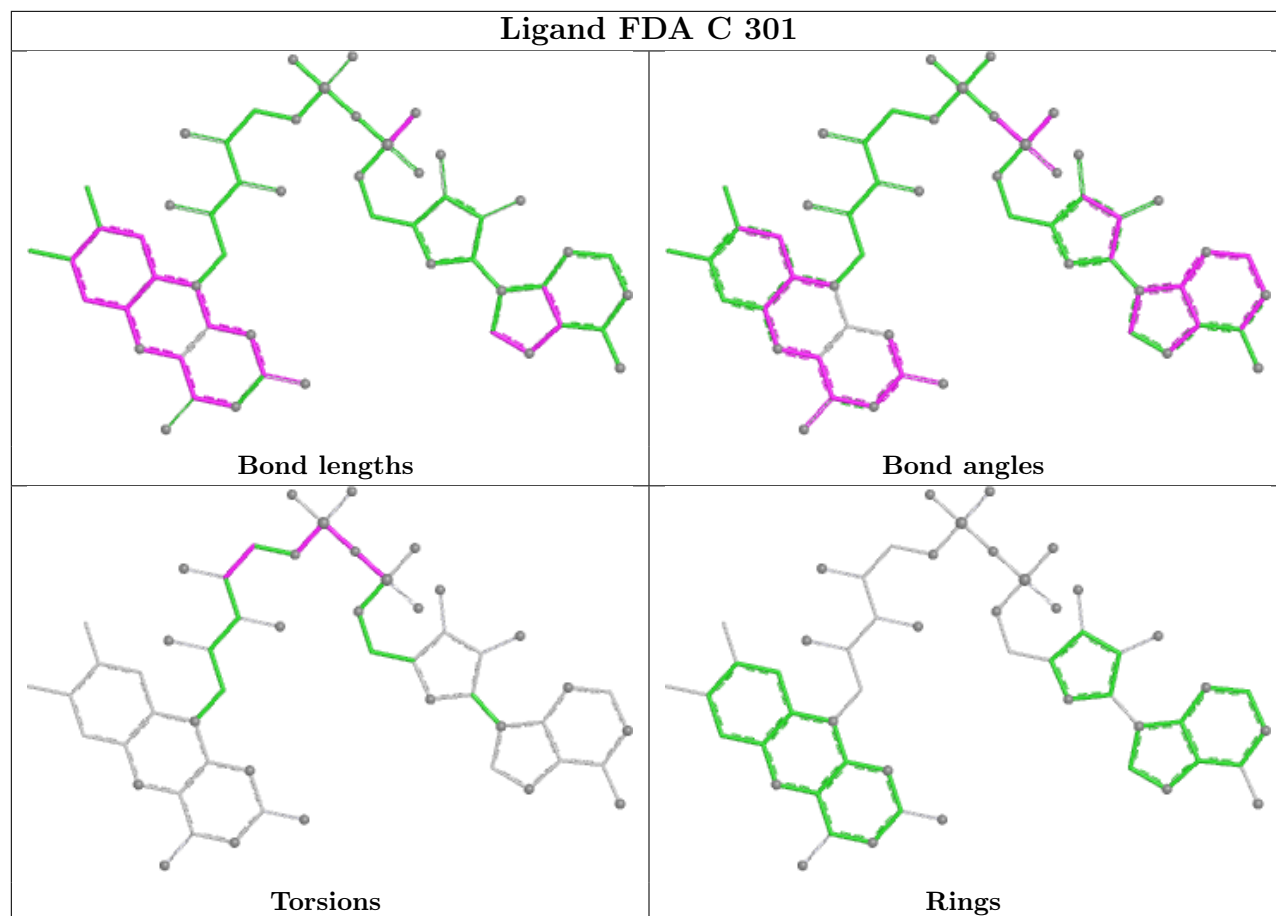
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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.

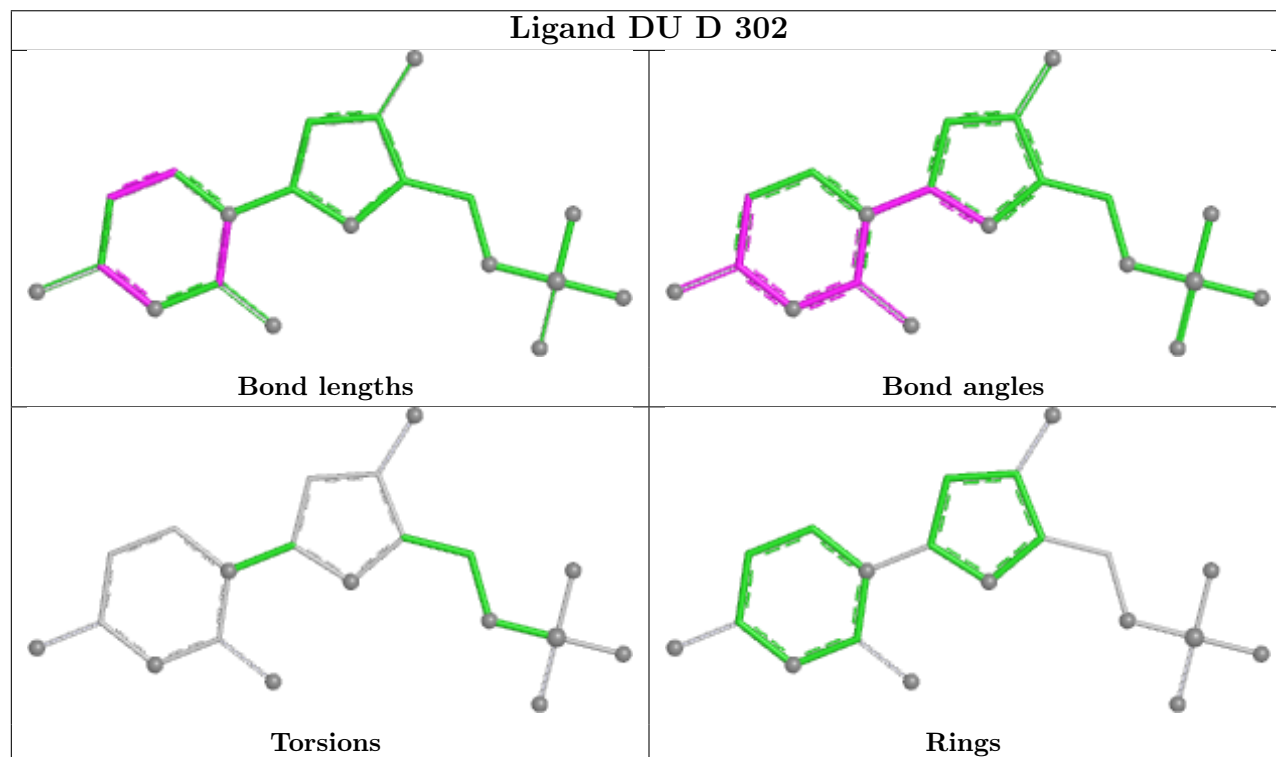




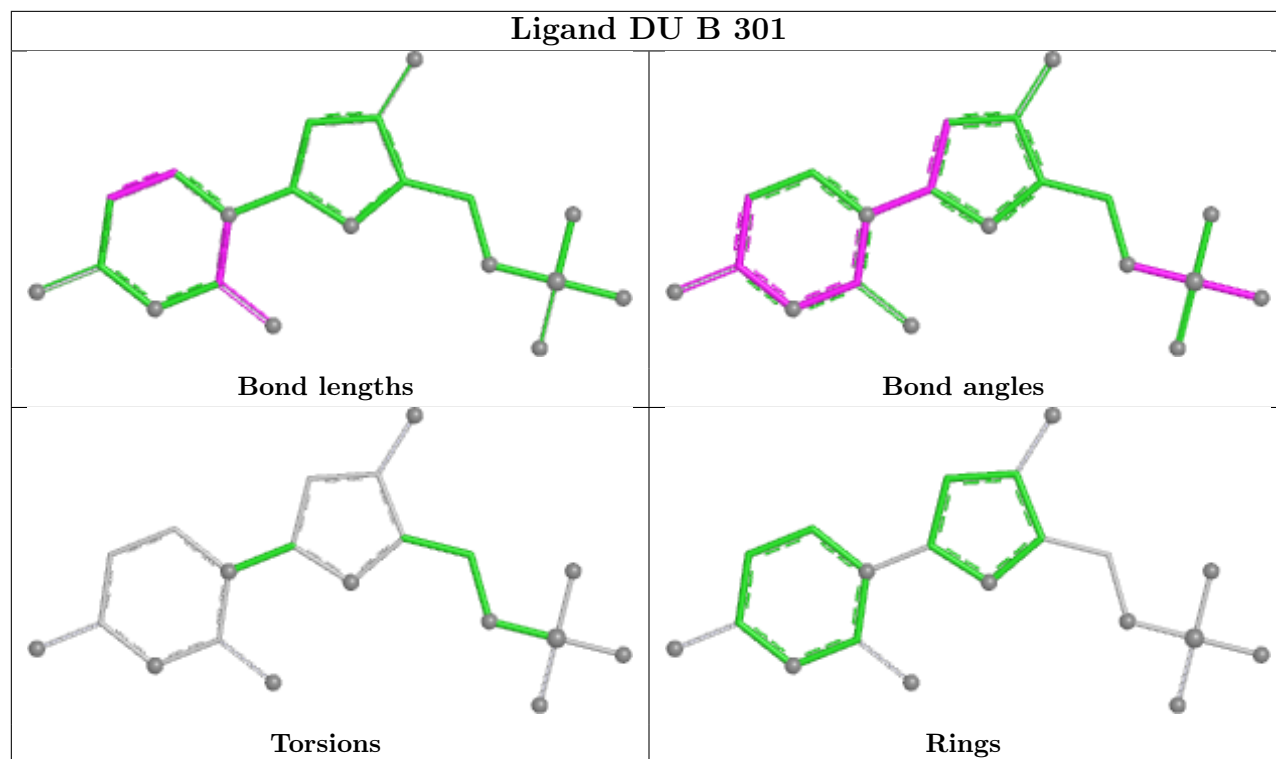
Ligand FDA C 301



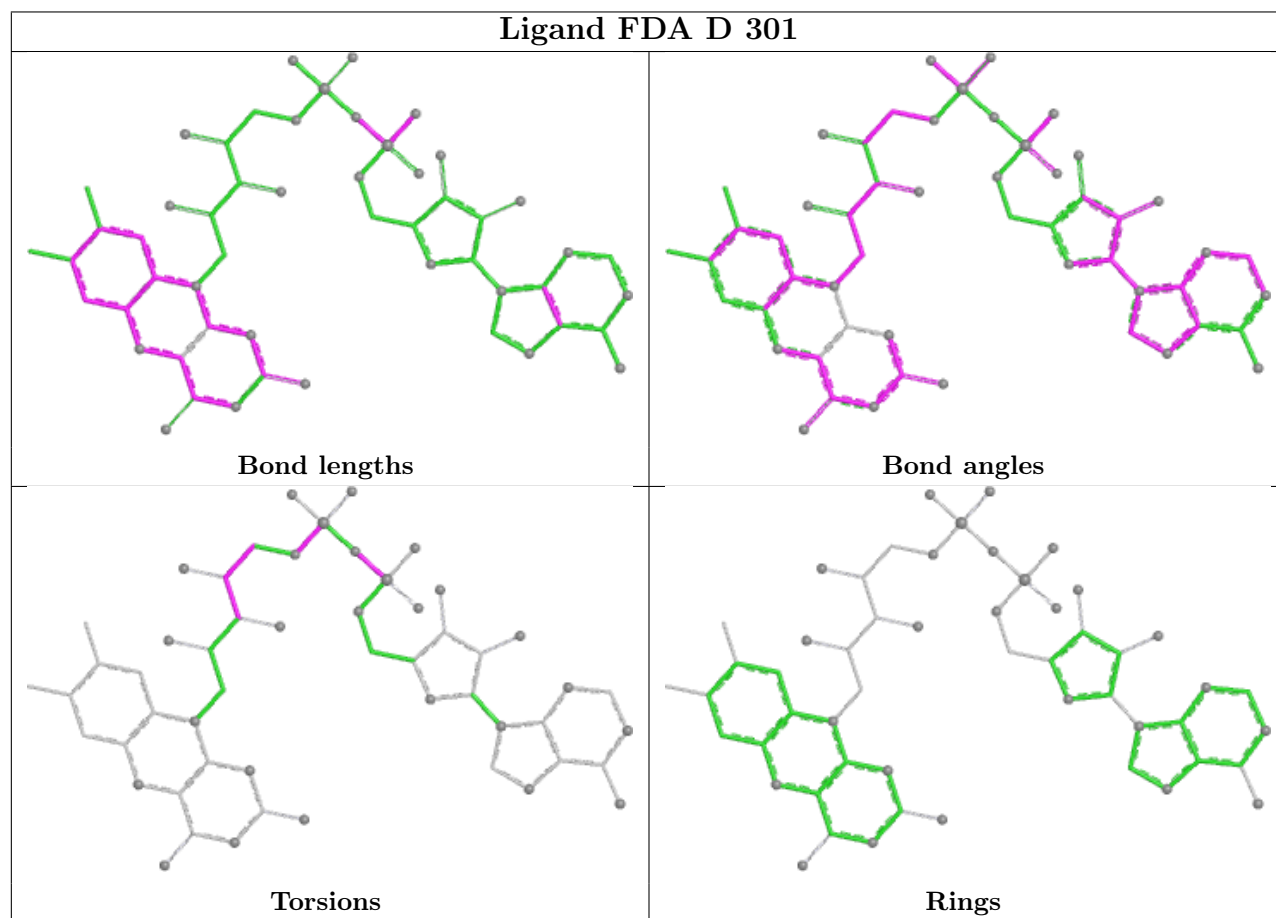
Ligand DU D 302

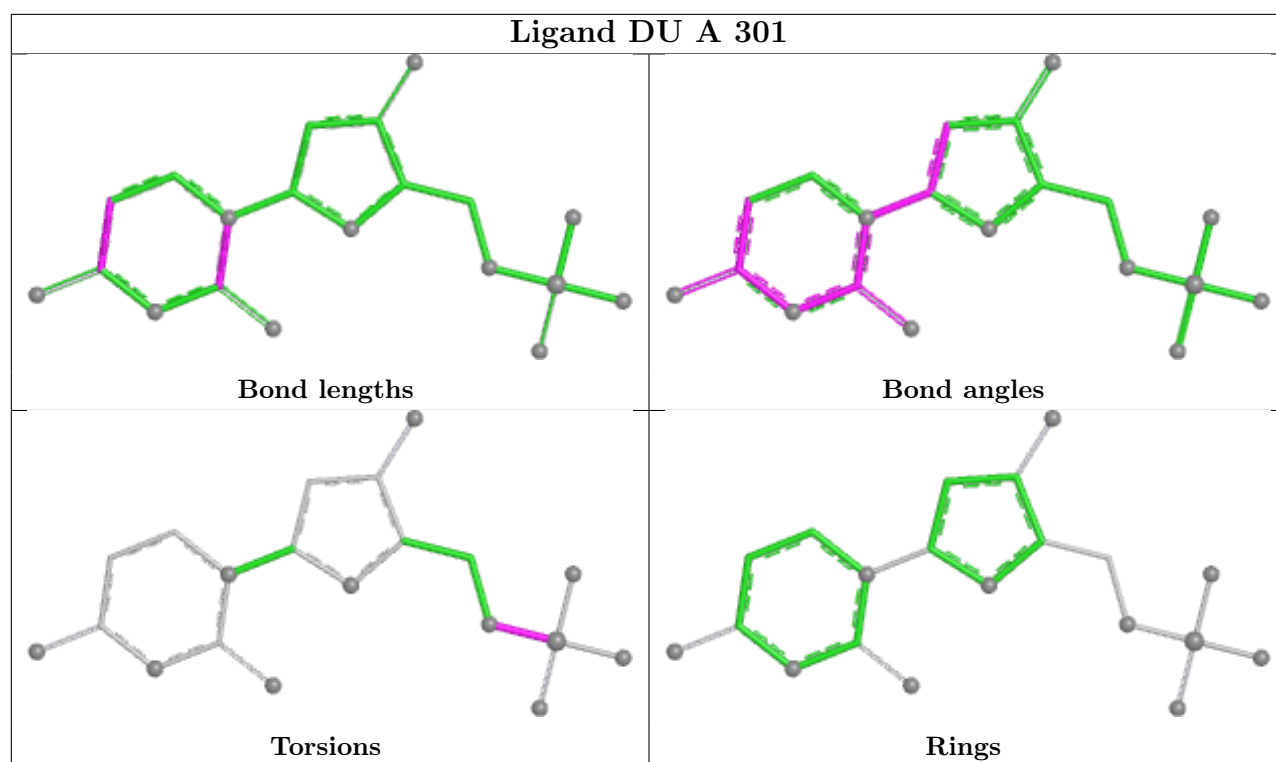


Ligand DU B 301



Ligand FDA D 301





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	218/232 (93%)	-0.23	0 100 100	23, 42, 65, 87	1 (0%)
1	B	220/232 (94%)	-0.21	0 100 100	31, 42, 74, 99	0
1	C	213/232 (91%)	-0.22	4 (1%) 66 70	29, 40, 68, 80	0
1	D	209/232 (90%)	-0.19	5 (2%) 59 63	30, 40, 68, 90	0
All	All	860/928 (92%)	-0.21	9 (1%) 79 83	23, 41, 68, 99	1 (0%)

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	94	LEU	3.4
1	C	218	VAL	3.0
1	D	218	VAL	2.7
1	D	91	TYR	2.3
1	D	0	HIS	2.2
1	C	96	TYR	2.2
1	D	96	TYR	2.2
1	C	91	TYR	2.1
1	D	31	PHE	2.1

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](#)

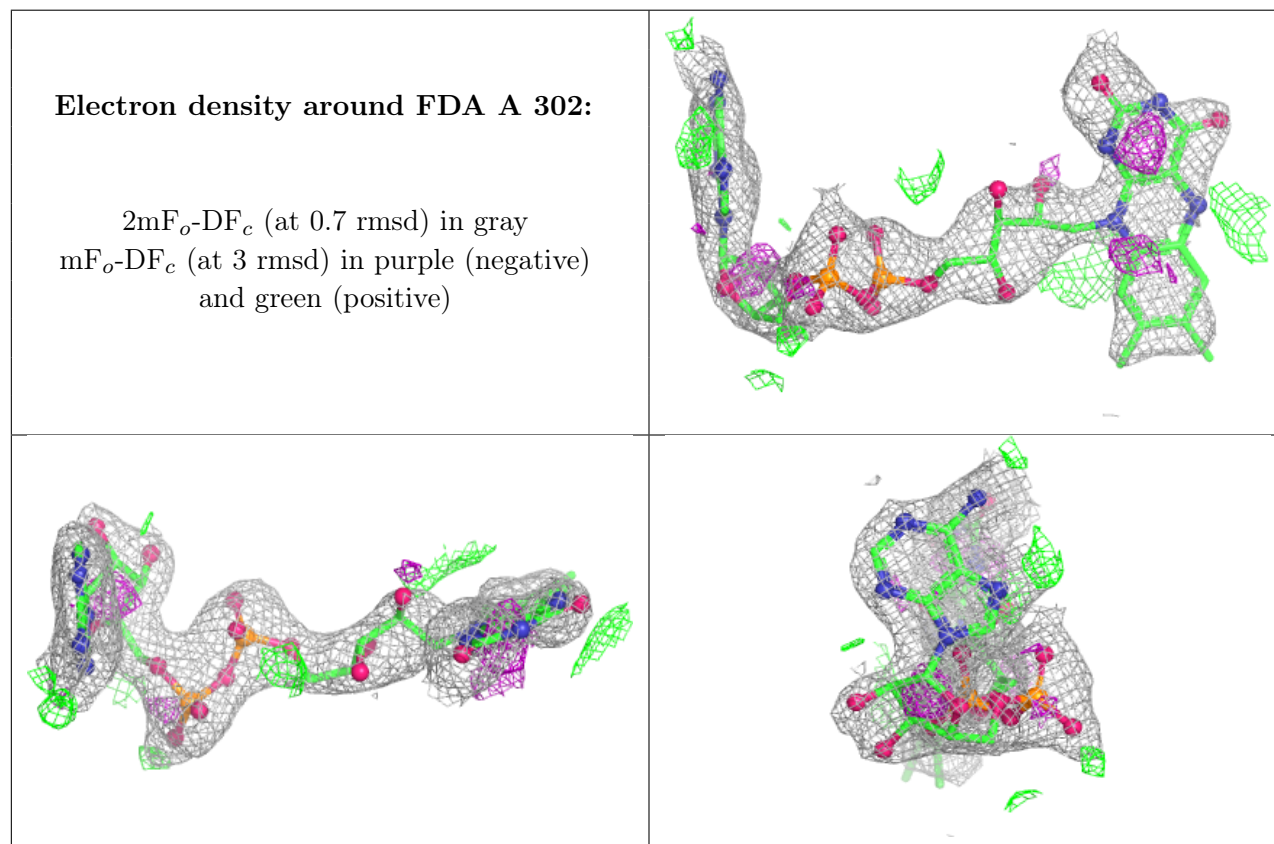
There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

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

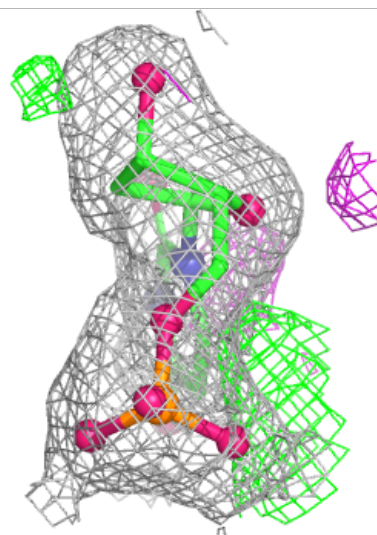
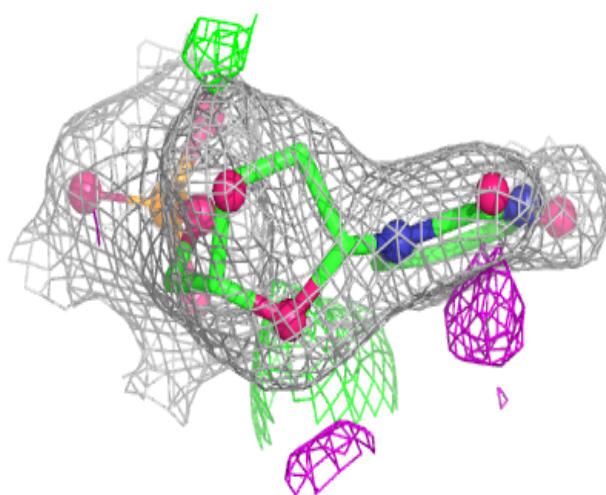
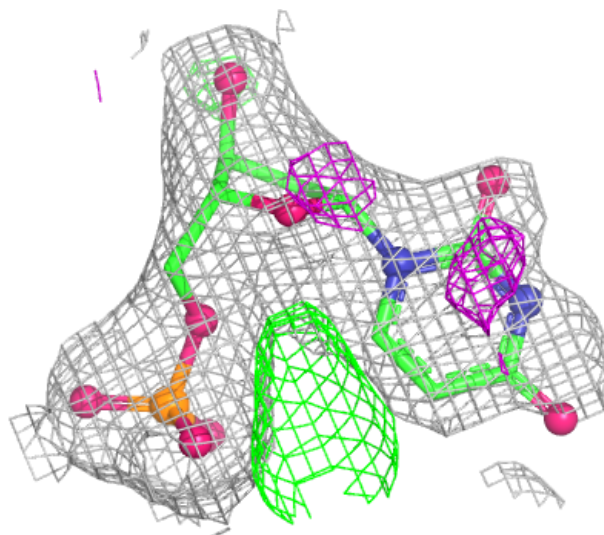
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	FDA	A	302	53/53	0.88	0.12	40,61,83,85	0
2	DU	D	302	20/20	0.89	0.11	44,51,71,77	0
3	FDA	B	302	53/53	0.89	0.11	39,64,77,86	0
3	FDA	D	301	53/53	0.89	0.11	35,60,69,82	0
3	FDA	C	301	53/53	0.91	0.10	35,60,69,74	0
2	DU	C	302	20/20	0.92	0.09	46,55,69,82	0
2	DU	A	301	20/20	0.96	0.07	37,43,58,60	0
2	DU	B	301	20/20	0.96	0.06	37,43,51,52	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.



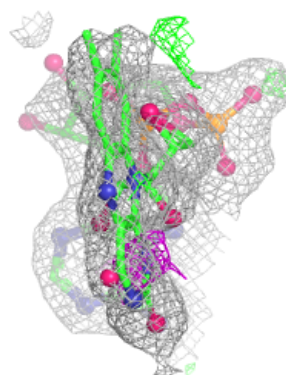
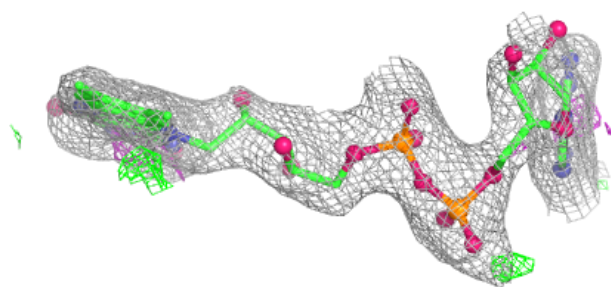
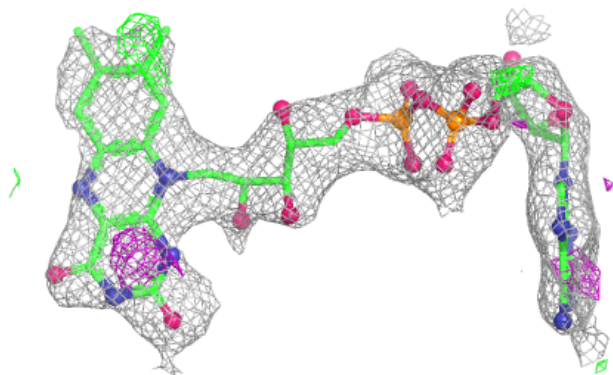
Electron density around DU D 302:

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

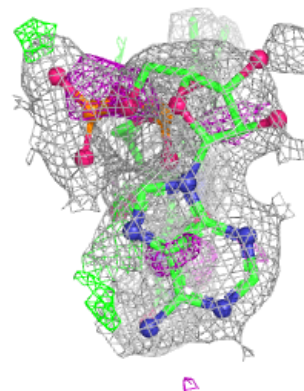
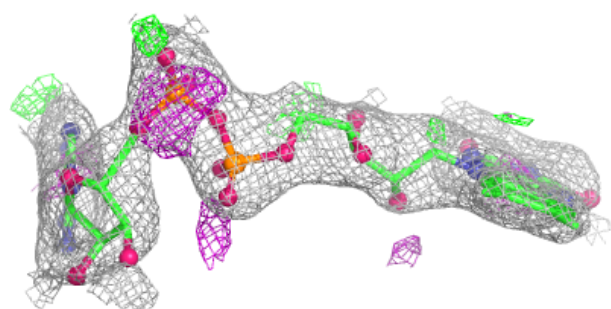
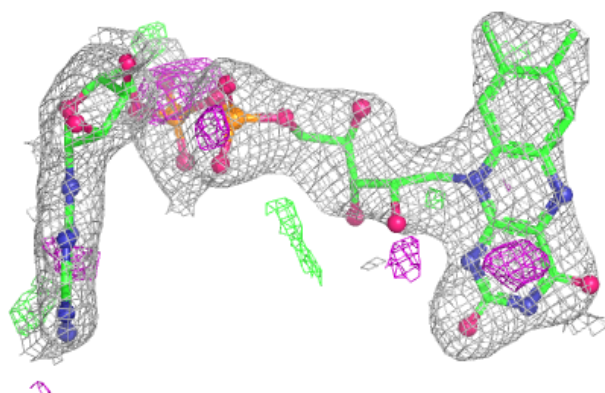


Electron density around FDA B 302:

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

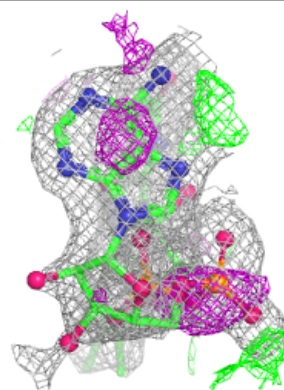
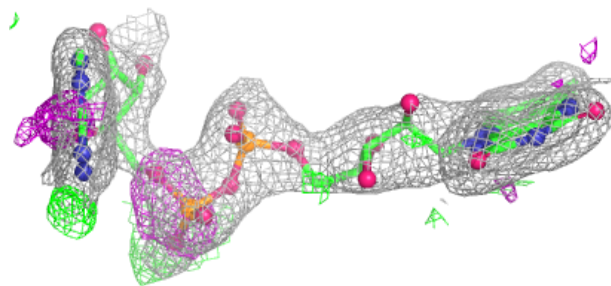
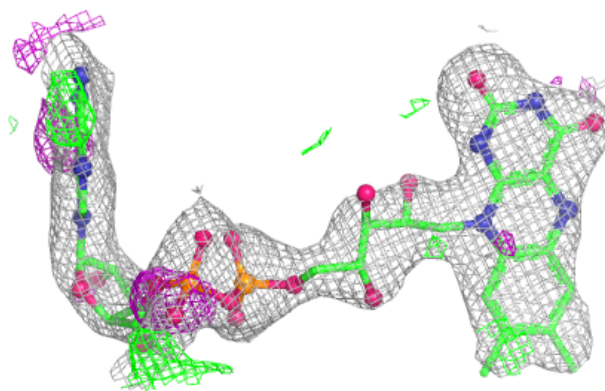
**Electron density around FDA D 301:**

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



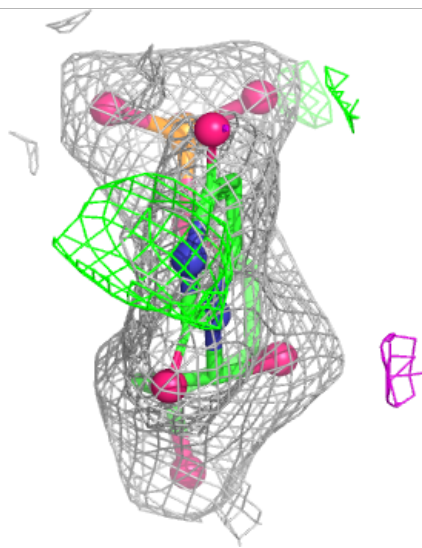
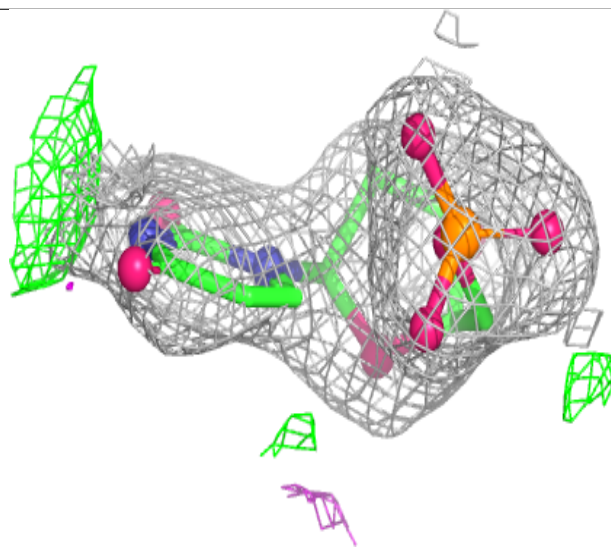
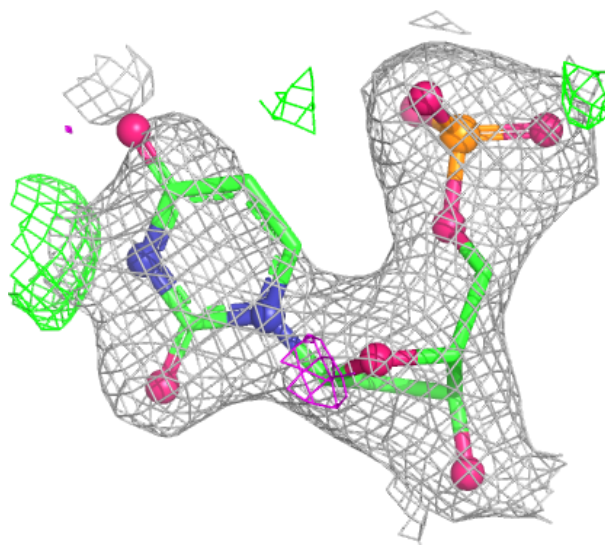
Electron density around FDA C 301:

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



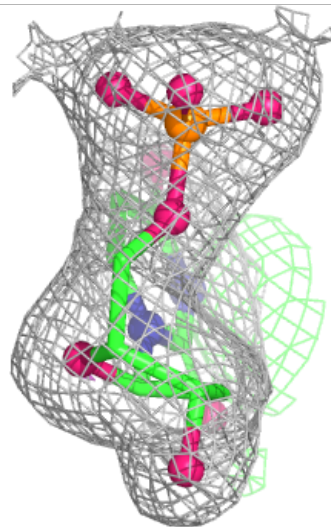
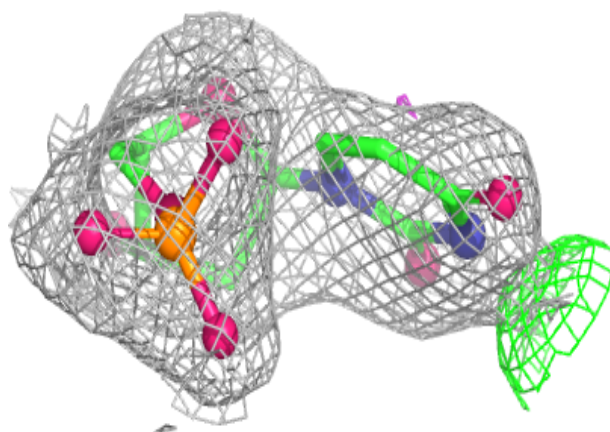
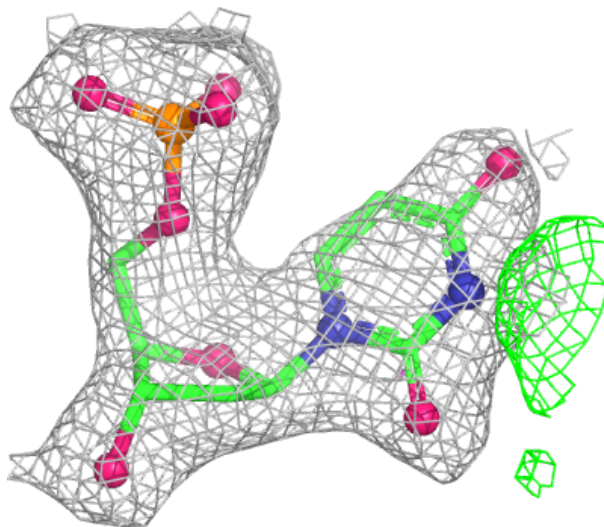
Electron density around DU C 302:

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



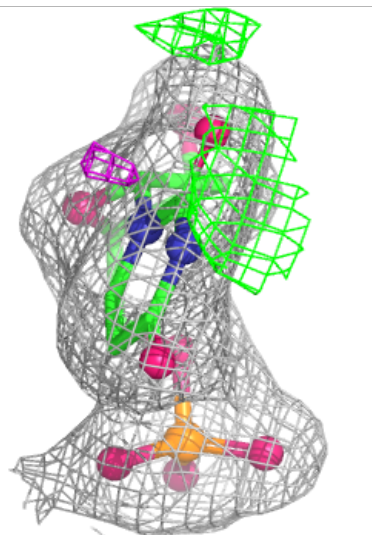
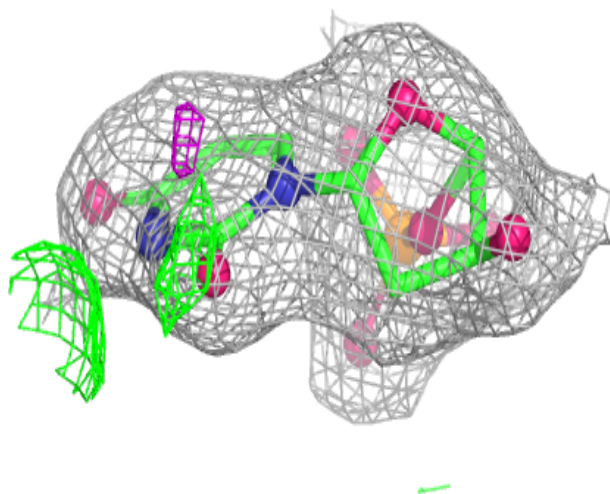
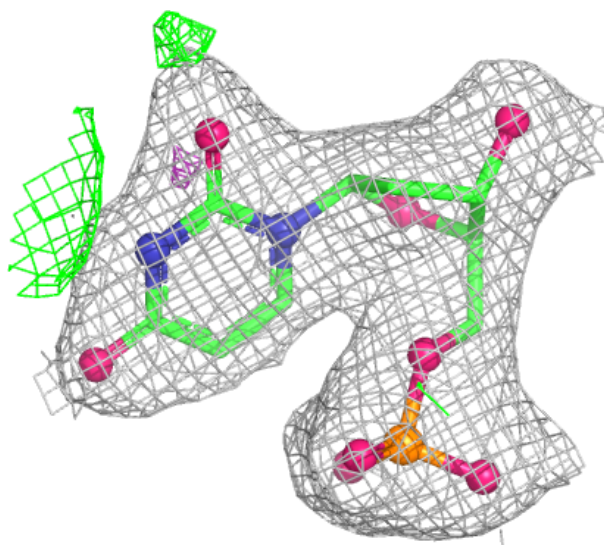
Electron density around DU A 301:

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



Electron density around DU B 301:

$2mF_o-DF_c$ (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.