



# Full wwPDB X-ray Structure Validation Report ⓘ

Jun 19, 2020 – 07:36 pm BST

PDB ID : 4KAT  
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](mailto: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.

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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

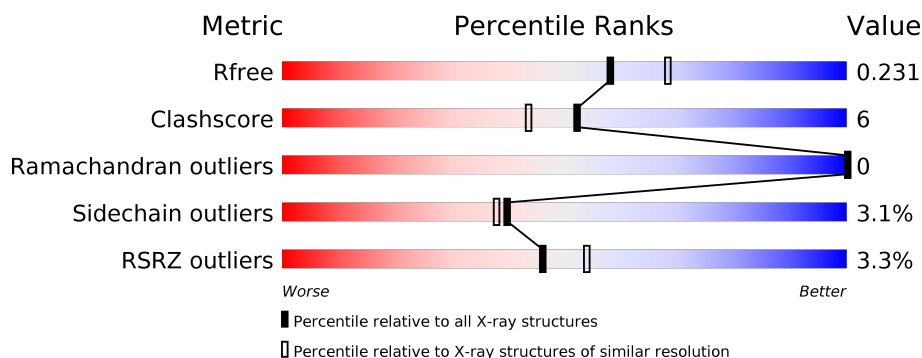
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.14 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))
$R_{free}$	130704	2523 (2.16-2.12)
Clashscore	141614	2653 (2.16-2.12)
Ramachandran outliers	138981	2618 (2.16-2.12)
Sidechain outliers	138945	2617 (2.16-2.12)
RSRZ outliers	127900	2485 (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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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	<div> <div>3%</div> <div> <div></div> <div>78%</div> <div>16%</div> <div>• 6%</div> </div> </div>
1	B	232	<div> <div>5%</div> <div> <div></div> <div>81%</div> <div>13%</div> <div>• 5%</div> </div> </div>
1	C	232	<div> <div>2%</div> <div> <div></div> <div>81%</div> <div>9%</div> <div>• 8%</div> </div> </div>
1	D	232	<div> <div>2%</div> <div> <div></div> <div>76%</div> <div>14%</div> <div>10%</div> </div> </div>

## 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:

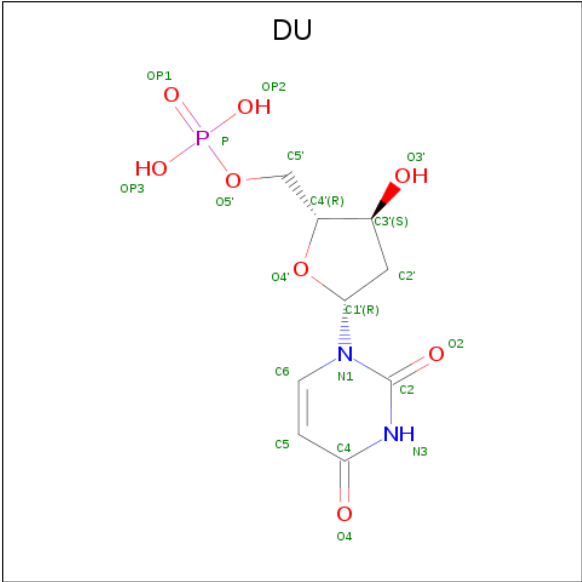
Chain	Residue	Modelled	Actual	Comment	Reference
A	-11	MET	-	INITIATING METHIONINE	UNP Q9WYT0
A	-10	GLY	-	EXPRESSION TAG	UNP Q9WYT0
A	-9	SER	-	EXPRESSION TAG	UNP Q9WYT0
A	-8	ASP	-	EXPRESSION TAG	UNP Q9WYT0
A	-7	LYS	-	EXPRESSION TAG	UNP Q9WYT0
A	-6	ILE	-	EXPRESSION TAG	UNP Q9WYT0
A	-5	HIS	-	EXPRESSION 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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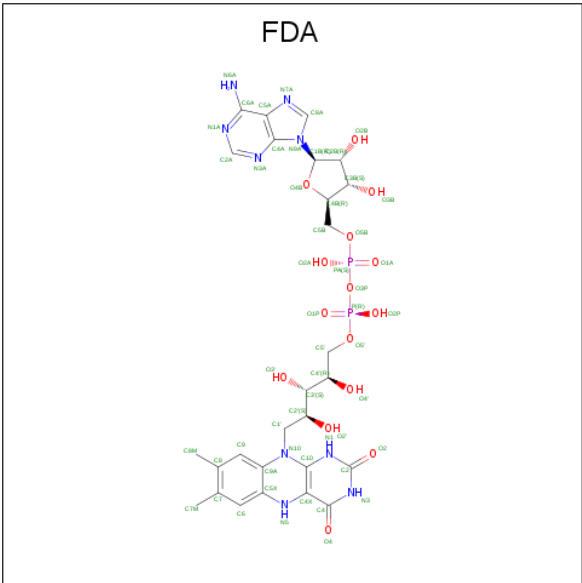
Chain	Residue	Modelled	Actual	Comment	Reference
B	-3	HIS	-	EXPRESSION TAG	UNP Q9WYT0
B	-2	HIS	-	EXPRESSION TAG	UNP Q9WYT0
B	-1	HIS	-	EXPRESSION TAG	UNP Q9WYT0
B	0	HIS	-	EXPRESSION TAG	UNP Q9WYT0
B	174	LYS	ARG	ENGINEERED MUTATION	UNP Q9WYT0
C	-11	MET	-	INITIATING METHIONINE	UNP Q9WYT0
C	-10	GLY	-	EXPRESSION 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 (three-letter code: DU) (formula: C<sub>9</sub>H<sub>13</sub>N<sub>2</sub>O<sub>8</sub>P).



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 (three-letter code: 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		

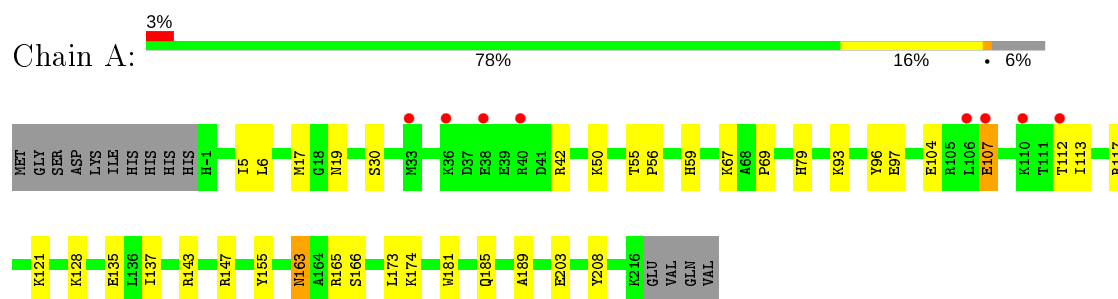
- Molecule 4 is water.

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		

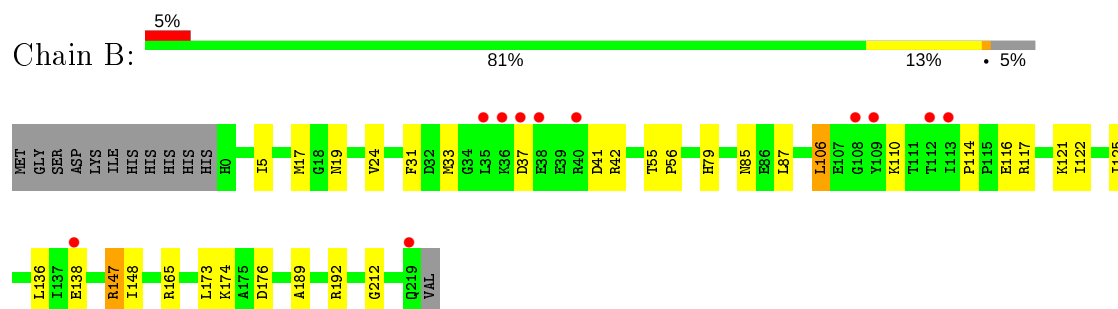
### 3 Residue-property plots [i](#)

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.

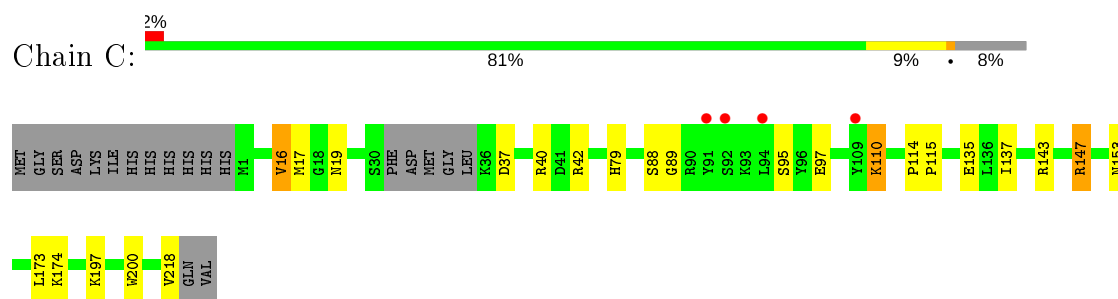
- Molecule 1: Thymidylate synthase



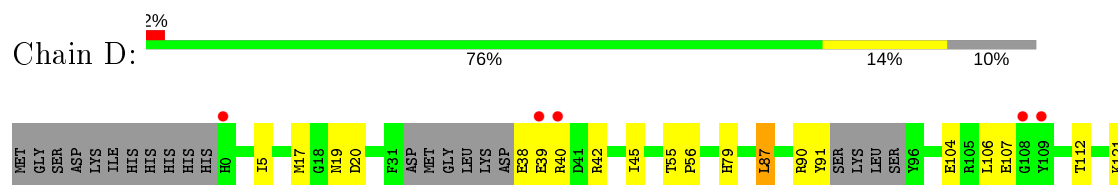
- Molecule 1: Thymidylate synthase



- Molecule 1: Thymidylate synthase



- Molecule 1: Thymidylate synthase



I125	E135	R143	R147	R157	R165	L173	K174	A175	D176	A189	W200	K207	G212	V218	GLN	VAL
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## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	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
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.01 (at 2.14Å)	Xtriage
Refinement program	REFMAC	Depositor
R, $R_{free}$	0.183 , 0.233 0.181 , 0.231	Depositor DCC
$R_{free}$ test set	2457 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	40.0	Xtriage
Anisotropy	0.180	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.37 , 48.7	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	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 <sup>2</sup> )	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.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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	0.70	0/1888	0.80	0/2551
1	B	0.69	0/1891	0.79	1/2556 (0.0%)
1	C	0.73	0/1830	0.81	1/2473 (0.0%)
1	D	0.67	0/1800	0.75	0/2433
All	All	0.70	0/7409	0.79	2/10013 (0.0%)

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 (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	147	ARG	NE-CZ-NH2	-5.26	117.67	120.30
1	C	37	ASP	CB-CG-OD1	5.22	123.00	118.30

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	19	0
1	C	1783	0	1785	17	0
1	D	1753	0	1744	19	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	3	0
3	B	53	0	33	4	0
3	C	53	0	33	4	0
3	D	53	0	33	2	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	82	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 (82) 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: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
1:A:107:GLU:OE1	1:A:107:GLU:HA	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:121:LYS:HE3	1:C:135:GLU:OE2	2.00	0.62
1:B:19:ASN:HB2	4:B:416:HOH:O	2.02	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:A:163:ASN:HD22	1:A:163:ASN:C	2.07	0.58
1:B:165:ARG:CZ	3:D:301:FDA:H2B	2.35	0.57
1:C:218:VAL:O	1:C:218:VAL:HG12	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:173:LEU:HB3	3:B:302:FDA:HM81	1.88	0.54
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:D:173:LEU:HB3	3:D:301:FDA:HM81	1.92	0.50
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:C:79:HIS:NE2	1:C:174:LYS:NZ	2.44	0.49
1:A:104:GLU:O	1:A:107:GLU:HG2	2.13	0.49
1:A:5:ILE:HG22	1:A:6:LEU:HG	1.94	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:CA	1:A:107:GLU:OE1	2.63	0.46
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:B:148:ILE:HB	1:C:153:ASN:HB3	1.96	0.46
1:A:173:LEU:HB3	3:A:302:FDA:HM81	1.97	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:31:PHE:HB2	1:B:33:MET:HE2	1.97	0.46
1:B:176:ASP:OD1	1:B:212:GLY:HA2	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:C:110:LYS:HE2	1:C:110:LYS:HB2	1.68	0.44
1:A:135:GLU:OE2	1:D:121:LYS:NZ	2.41	0.44
1:B:114:PRO:HG2	1:B:117:ARG:CG	2.48	0.43
2:C:302:DU:C5'	2:C:302:DU:H6	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:173:LEU:HB3	3:C:301:FDA:HM81	1.99	0.43
1:A:30:SER:HG	3:C:301:FDA:C4	2.32	0.42
1:C:143:ARG:O	1:C:147:ARG:HG3	2.20	0.42
1:A:69:PRO:HD3	1:A:155:TYR:CE1	2.54	0.42
1:A:59:HIS:NE2	1:B:85:ASN:ND2	2.62	0.42
1:B:55:THR:N	1:B:56:PRO:CD	2.82	0.42
1:C:137:ILE:HD13	1:C:137:ILE:HA	1.83	0.42
1:C:42:ARG:HG2	1:C:200:TRP:CD2	2.55	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:67:LYS:NZ	1:A:97:GLU:OE2	2.44	0.42
1:C:114:PRO:HA	1:C:115:PRO:HD2	1.76	0.42
1:A:165:ARG:CZ	3:C:301:FDA:H2B	2.50	0.42
3:B:302:FDA:H1'2	3:B:302:FDA:H9	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: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

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	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%)	191 (98%)	4 (2%)	53	54
1	B	196/207 (95%)	187 (95%)	9 (5%)	27	22
1	C	190/207 (92%)	186 (98%)	4 (2%)	53	54
1	D	185/207 (89%)	178 (96%)	7 (4%)	33	30
All	All	766/828 (92%)	742 (97%)	24 (3%)	40	38

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	96	TYR
1	A	107	GLU
1	A	147	ARG
1	A	163	ASN
1	B	37	ASP
1	B	42	ARG
1	B	87	LEU
1	B	106	LEU
1	B	110	LYS
1	B	116	GLU

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Mol	Chain	Res	Type
1	B	138	GLU
1	B	147	ARG
1	B	192	ARG
1	C	16	VAL
1	C	40	ARG
1	C	110	LYS
1	C	147	ARG
1	D	38	GLU
1	D	39	GLU
1	D	40	ARG
1	D	87	LEU
1	D	112	THR
1	D	147	ARG
1	D	207	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	75	GLN
1	A	163	ASN
1	B	85	ASN
1	C	51	HIS
1	C	185	GLN
1	D	85	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 carbohydrates 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	D	301	-	51,58,58	3.56	16 (31%)	60,89,89	2.33	18 (30%)
3	FDA	B	302	-	51,58,58	3.72	16 (31%)	60,89,89	2.43	13 (21%)
3	FDA	C	301	-	51,58,58	3.71	15 (29%)	60,89,89	2.38	14 (23%)
3	FDA	A	302	-	51,58,58	3.78	15 (29%)	60,89,89	2.56	12 (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	D	301	-	-	5/30/50/50	0/6/6/6
3	FDA	B	302	-	-	4/30/50/50	0/6/6/6
3	FDA	C	301	-	-	3/30/50/50	0/6/6/6
3	FDA	A	302	-	-	3/30/50/50	0/6/6/6

All (62) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	302	FDA	C4X-C10	14.06	1.52	1.38
3	B	302	FDA	C4X-C10	13.53	1.52	1.38
3	C	301	FDA	C4X-C10	13.39	1.52	1.38
3	D	301	FDA	C4X-C10	12.90	1.51	1.38
3	A	302	FDA	C10-N1	11.61	1.48	1.33
3	B	302	FDA	C10-N1	11.33	1.47	1.33
3	C	301	FDA	C10-N1	11.25	1.47	1.33
3	D	301	FDA	C10-N1	10.23	1.46	1.33
3	A	302	FDA	C4X-N5	8.51	1.45	1.33
3	C	301	FDA	C4X-N5	8.16	1.45	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	302	FDA	C6-C5X	7.90	1.54	1.41
3	B	302	FDA	C4X-N5	7.69	1.44	1.33
3	D	301	FDA	C6-C5X	7.47	1.53	1.41
3	A	302	FDA	C6-C5X	7.26	1.53	1.41
3	C	301	FDA	C6-C5X	7.20	1.53	1.41
3	D	301	FDA	C4X-N5	7.11	1.43	1.33
3	D	301	FDA	C9A-N10	-6.50	1.29	1.38
3	C	301	FDA	C5X-N5	6.38	1.45	1.35
3	A	302	FDA	C5X-N5	6.38	1.45	1.35
3	B	302	FDA	C9A-N10	-6.37	1.29	1.38
3	C	301	FDA	C9A-N10	-6.31	1.30	1.38
3	B	302	FDA	C5X-N5	6.23	1.45	1.35
3	A	302	FDA	C9A-N10	-5.87	1.30	1.38
3	A	302	FDA	C6-C7	5.83	1.52	1.37
3	B	302	FDA	C6-C7	5.81	1.52	1.37
3	D	301	FDA	C6-C7	5.71	1.52	1.37
3	C	301	FDA	C6-C7	5.60	1.52	1.37
3	D	301	FDA	C5X-N5	5.51	1.44	1.35
3	B	302	FDA	C9A-C5X	5.46	1.53	1.42
3	C	301	FDA	C2-N1	-5.33	1.27	1.38
3	A	302	FDA	C9A-C5X	5.29	1.53	1.42
3	C	301	FDA	C9A-C5X	5.22	1.53	1.42
3	A	302	FDA	C4-C4X	5.21	1.50	1.41
3	A	302	FDA	C2-N1	-5.15	1.28	1.38
3	B	302	FDA	C4-C4X	5.14	1.50	1.41
3	D	301	FDA	C2-N1	-5.12	1.28	1.38
3	D	301	FDA	C4-C4X	5.07	1.50	1.41
3	D	301	FDA	C9A-C5X	5.06	1.52	1.42
3	C	301	FDA	C4-C4X	5.03	1.50	1.41
3	B	302	FDA	C8-C7	4.57	1.52	1.40
3	C	301	FDA	C8-C7	4.57	1.52	1.40
3	B	302	FDA	C2-N1	-4.54	1.29	1.38
3	A	302	FDA	C8-C7	4.50	1.52	1.40
3	D	301	FDA	C9-C9A	4.43	1.49	1.40
3	B	302	FDA	C9-C9A	4.41	1.49	1.40
3	A	302	FDA	C9-C9A	4.29	1.49	1.40
3	D	301	FDA	C8-C7	4.17	1.51	1.40
3	C	301	FDA	C9-C9A	3.98	1.48	1.40
3	C	301	FDA	C5A-C4A	3.61	1.50	1.40
3	A	302	FDA	C5A-C4A	3.29	1.49	1.40
3	D	301	FDA	C5A-C4A	3.27	1.49	1.40
3	B	302	FDA	C5A-C4A	3.09	1.49	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	301	FDA	O4B-C1B	2.75	1.44	1.41
3	C	301	FDA	PA-O1A	2.71	1.60	1.50
3	D	301	FDA	C2A-N3A	2.58	1.36	1.32
3	B	302	FDA	O4B-C1B	2.42	1.44	1.41
3	C	301	FDA	O4B-C1B	2.28	1.44	1.41
3	D	301	FDA	PA-O1A	2.17	1.58	1.50
3	A	302	FDA	O4B-C1B	2.16	1.44	1.41
3	A	302	FDA	PA-O1A	2.14	1.58	1.50
3	B	302	FDA	PA-O1A	2.10	1.58	1.50
3	B	302	FDA	C2A-N3A	2.03	1.35	1.32

All (57) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	302	FDA	C4-N3-C2	11.80	125.11	115.14
3	C	301	FDA	C4-N3-C2	10.58	124.07	115.14
3	B	302	FDA	C4-N3-C2	10.54	124.04	115.14
3	D	301	FDA	C4-N3-C2	8.54	122.35	115.14
3	A	302	FDA	C4-C4X-C10	-6.86	115.41	119.95
3	A	302	FDA	C5X-C9A-N10	6.86	122.69	117.72
3	D	301	FDA	C1'-N10-C9A	6.51	123.41	118.29
3	B	302	FDA	C5X-C9A-N10	6.16	122.18	117.72
3	B	302	FDA	C4-C4X-C10	-6.05	115.94	119.95
3	C	301	FDA	C4-C4X-C10	-5.70	116.18	119.95
3	C	301	FDA	C5X-C9A-N10	5.65	121.81	117.72
3	B	302	FDA	C4-C4X-N5	5.33	124.69	118.60
3	A	302	FDA	C4-C4X-N5	5.19	124.53	118.60
3	D	301	FDA	C5X-C9A-N10	5.06	121.38	117.72
3	C	301	FDA	C4-C4X-N5	4.87	124.16	118.60
3	D	301	FDA	C4-C4X-N5	4.50	123.74	118.60
3	A	302	FDA	N3A-C2A-N1A	-4.38	121.83	128.68
3	B	302	FDA	C1'-N10-C9A	4.35	121.71	118.29
3	C	301	FDA	C1'-N10-C9A	4.08	121.50	118.29
3	C	301	FDA	C4X-C4-N3	-3.96	118.02	123.43
3	A	302	FDA	C4X-C4-N3	-3.90	118.09	123.43
3	D	301	FDA	C4-C4X-C10	-3.89	117.38	119.95
3	D	301	FDA	C4X-C4-N3	-3.86	118.16	123.43
3	D	301	FDA	C3B-C2B-C1B	3.84	106.76	100.98
3	D	301	FDA	C7-C6-C5X	-3.80	115.85	121.22
3	B	302	FDA	C3B-C2B-C1B	3.69	106.54	100.98
3	D	301	FDA	C10-C4X-N5	-3.44	118.88	121.26
3	D	301	FDA	N3A-C2A-N1A	-3.44	123.31	128.68

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	302	FDA	C1'-N10-C9A	3.41	120.98	118.29
3	B	302	FDA	C4X-C4-N3	-3.39	118.80	123.43
3	C	301	FDA	N3A-C2A-N1A	-3.36	123.42	128.68
3	C	301	FDA	O2A-PA-O1A	3.23	128.19	112.24
3	B	302	FDA	C7-C6-C5X	-3.21	116.67	121.22
3	A	302	FDA	C2A-N1A-C6A	3.17	124.17	118.75
3	B	302	FDA	N3A-C2A-N1A	-3.13	123.78	128.68
3	C	301	FDA	C3B-C2B-C1B	3.09	105.63	100.98
3	B	302	FDA	C4A-C5A-N7A	-3.08	106.18	109.40
3	C	301	FDA	P-O3P-PA	-2.96	122.66	132.83
3	C	301	FDA	C7-C6-C5X	-2.89	117.13	121.22
3	A	302	FDA	P-O3P-PA	-2.81	123.18	132.83
3	B	302	FDA	C10-C4X-N5	-2.75	119.36	121.26
3	D	301	FDA	O2A-PA-O1A	2.69	125.54	112.24
3	C	301	FDA	C2A-N1A-C6A	2.65	123.29	118.75
3	D	301	FDA	C1'-N10-C10	-2.65	116.03	118.41
3	A	302	FDA	C4A-C5A-N7A	-2.65	106.64	109.40
3	D	301	FDA	O5'-C5'-C4'	-2.61	102.40	109.36
3	D	301	FDA	O2B-C2B-C1B	-2.59	101.28	110.85
3	D	301	FDA	C4A-C5A-N7A	-2.57	106.72	109.40
3	A	302	FDA	C7-C6-C5X	-2.50	117.68	121.22
3	B	302	FDA	P-O3P-PA	-2.46	124.37	132.83
3	B	302	FDA	O5'-C5'-C4'	-2.45	102.83	109.36
3	C	301	FDA	C10-C4X-N5	-2.32	119.66	121.26
3	A	302	FDA	O2A-PA-O1A	2.31	123.66	112.24
3	C	301	FDA	C4A-C5A-N7A	-2.16	107.15	109.40
3	D	301	FDA	O2P-P-O1P	2.06	122.44	112.24
3	D	301	FDA	O3'-C3'-C2'	2.06	113.78	108.81
3	D	301	FDA	C6-C5X-C9A	2.06	121.75	119.05

There are no chirality outliers.

All (15) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	C	301	FDA	P-O3P-PA-O5B
3	C	301	FDA	O4'-C4'-C5'-O5'
3	B	302	FDA	C3'-C4'-C5'-O5'
3	B	302	FDA	O4'-C4'-C5'-O5'
3	A	302	FDA	C3'-C4'-C5'-O5'
3	A	302	FDA	O4'-C4'-C5'-O5'
3	D	301	FDA	O4'-C4'-C5'-O5'
3	D	301	FDA	C3'-C4'-C5'-O5'

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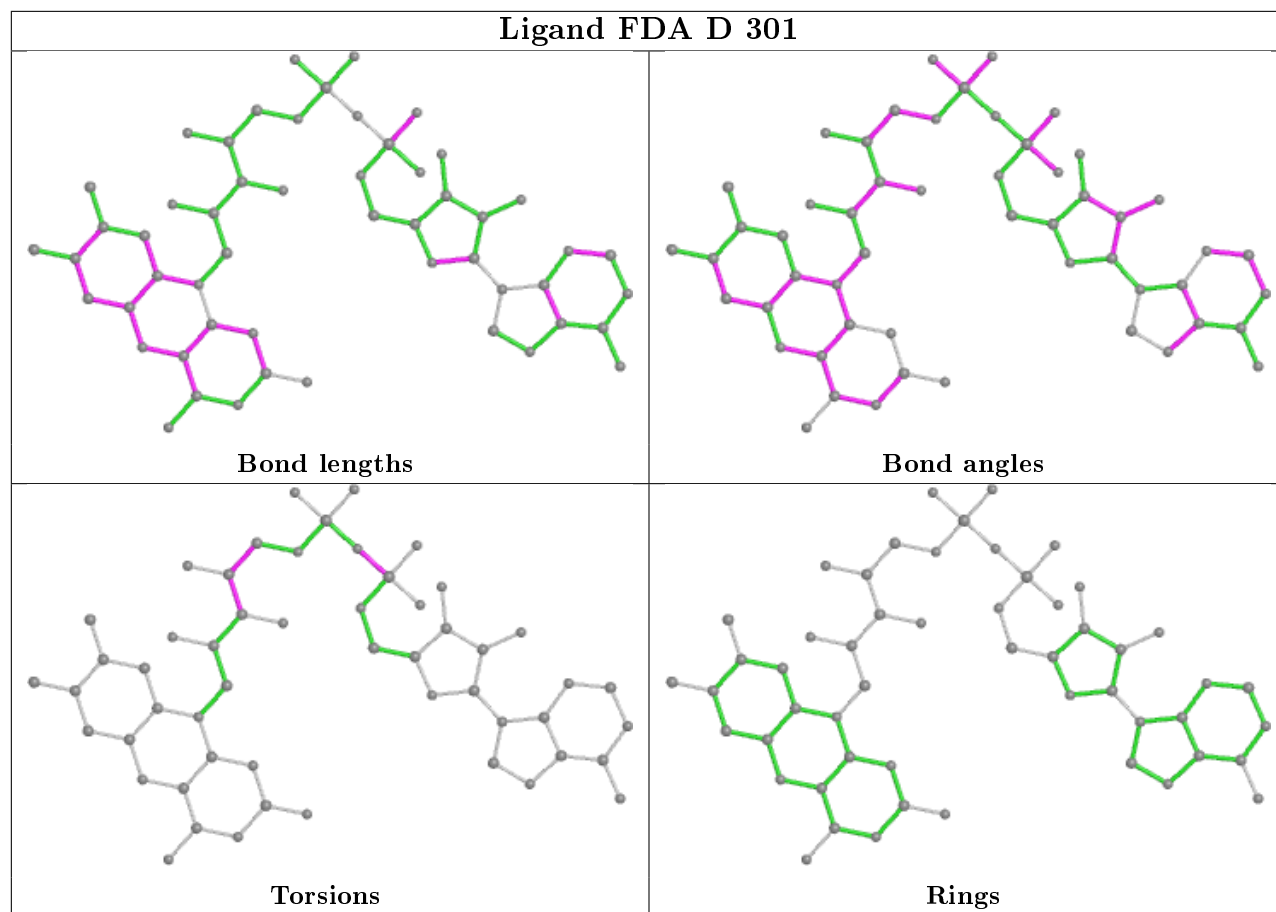
Mol	Chain	Res	Type	Atoms
3	D	301	FDA	P-O3P-PA-O5B
3	D	301	FDA	C2'-C3'-C4'-O4'
3	D	301	FDA	C2'-C3'-C4'-C5'
3	A	302	FDA	C2'-C3'-C4'-O4'
3	B	302	FDA	O4B-C4B-C5B-O5B
3	C	301	FDA	C5'-O5'-P-O1P
3	B	302	FDA	C3B-C4B-C5B-O5B

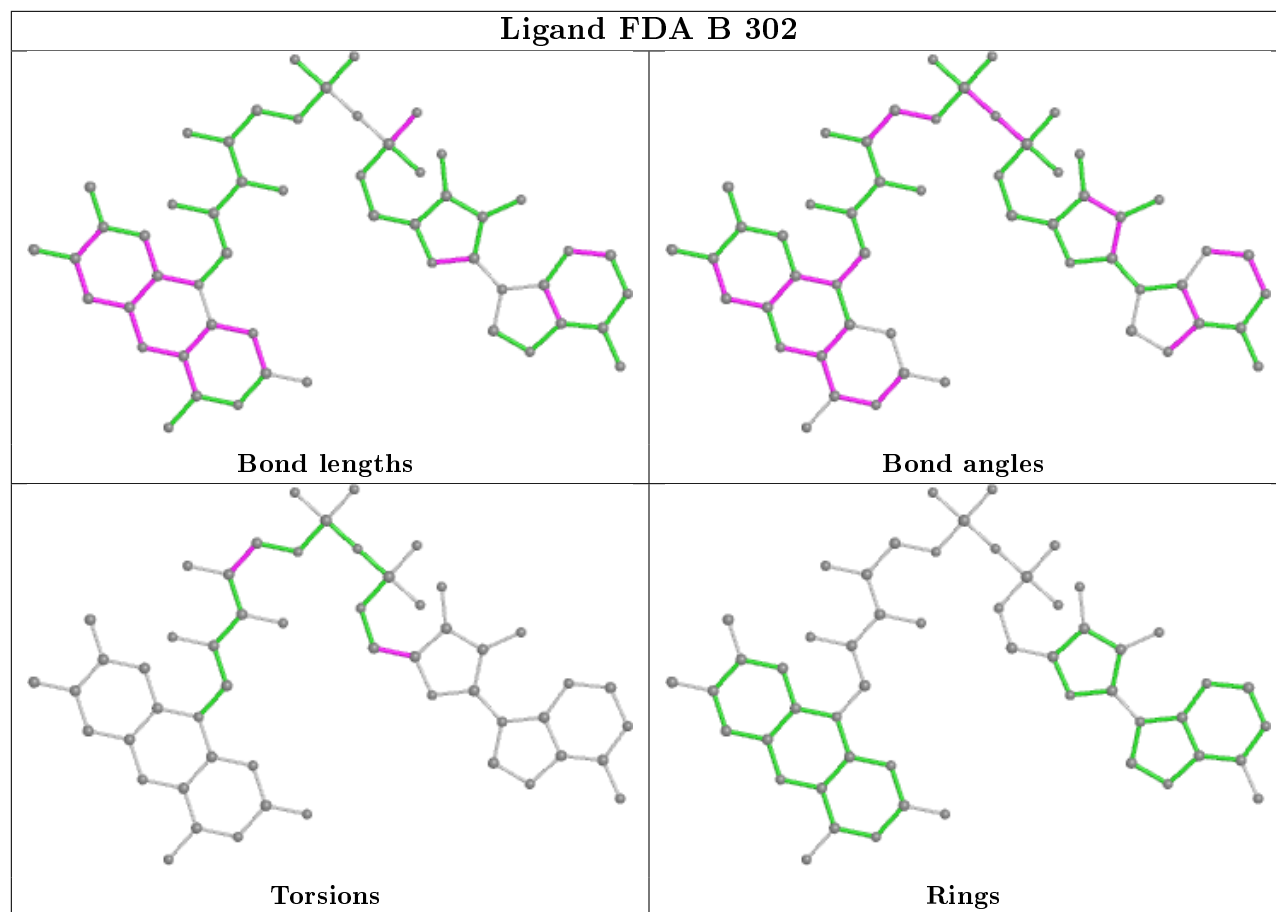
There are no ring outliers.

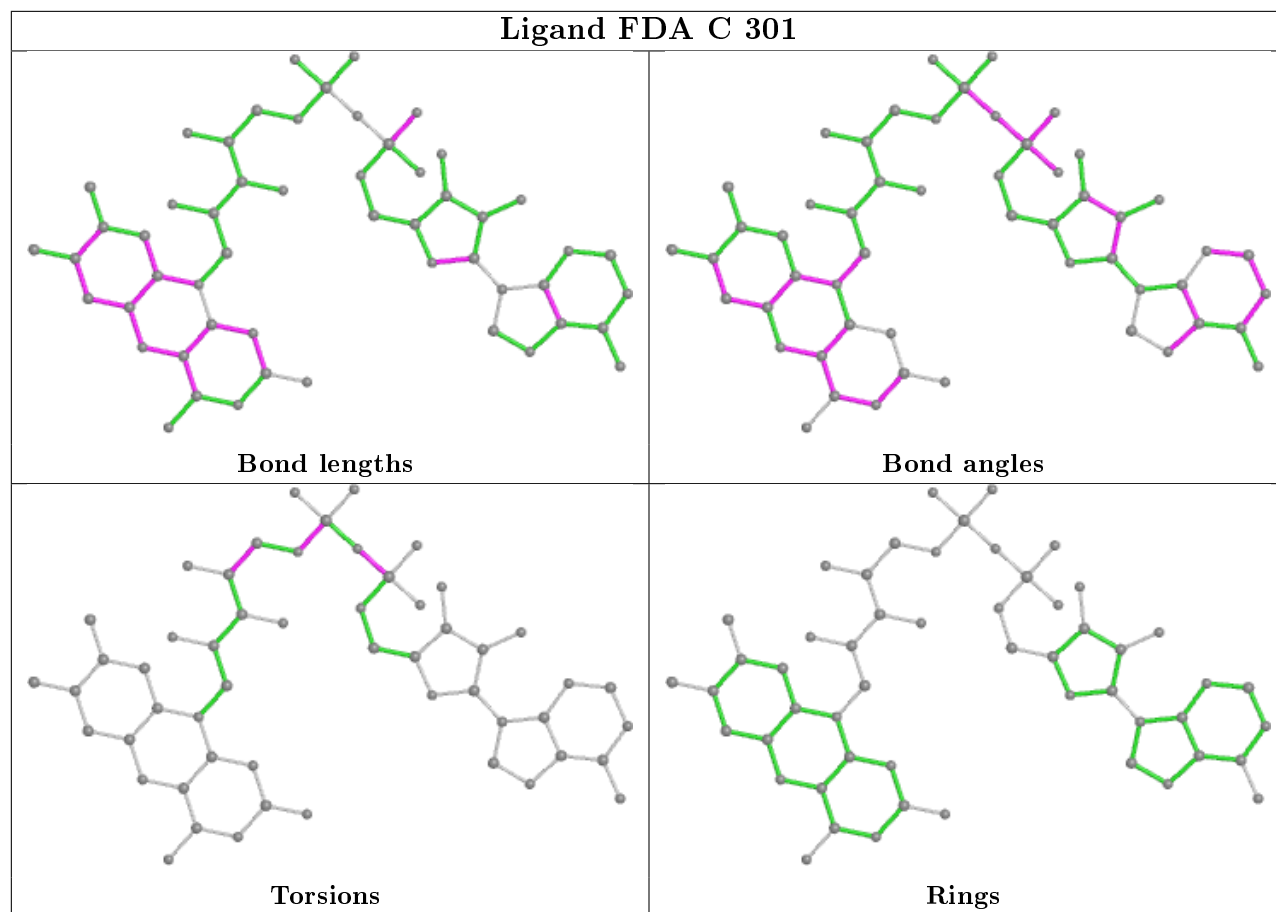
4 monomers are involved in 13 short contacts:

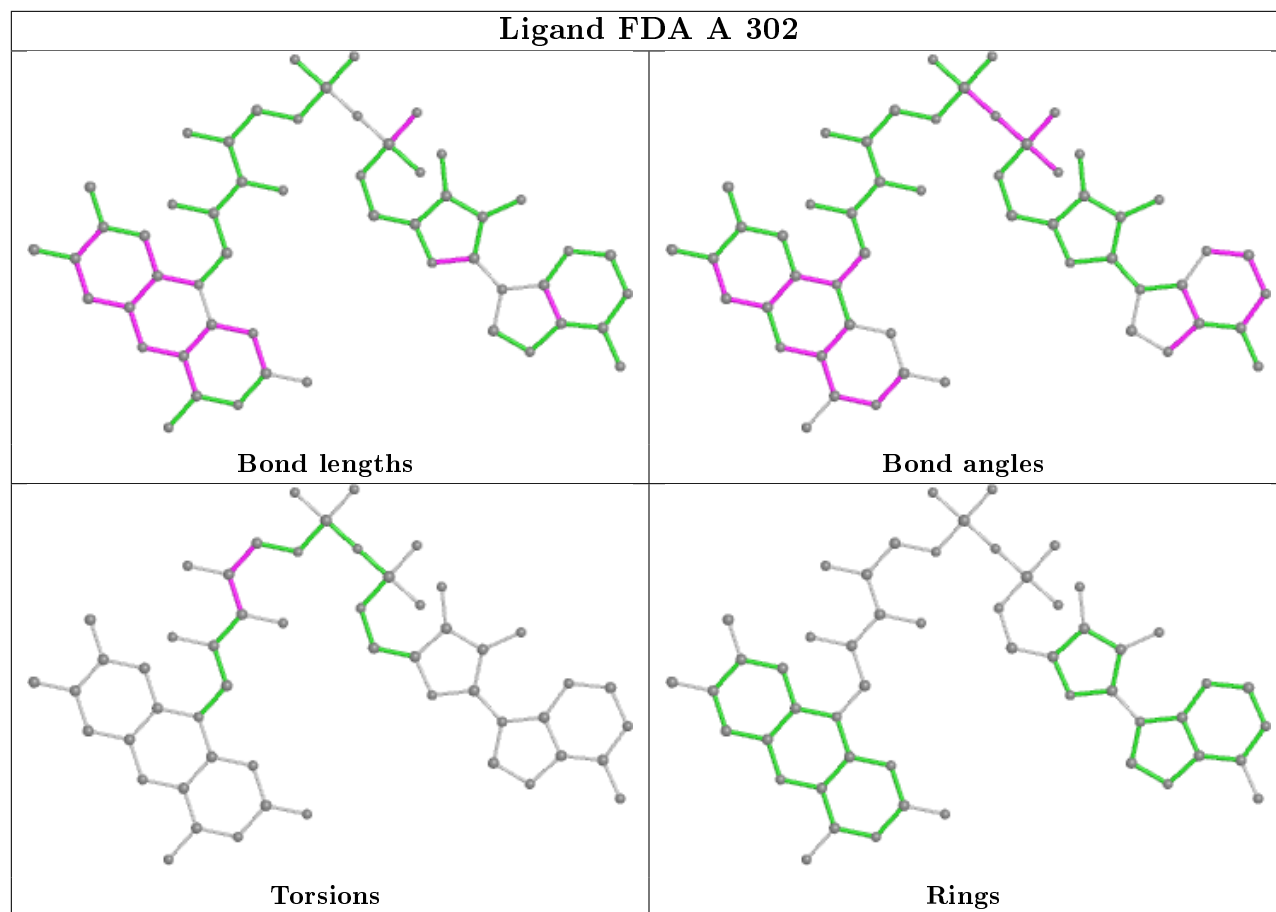
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	D	301	FDA	2	0
3	B	302	FDA	4	0
3	C	301	FDA	4	0
3	A	302	FDA	3	0

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









## 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	218/232 (93%)	0.07	8 (3%) 41 49	29, 42, 65, 87	0
1	B	220/232 (94%)	0.08	11 (5%) 28 35	31, 42, 74, 99	0
1	C	213/232 (91%)	-0.12	4 (1%) 66 72	29, 40, 68, 80	0
1	D	209/232 (90%)	-0.06	5 (2%) 59 65	30, 40, 68, 90	0
All	All	860/928 (92%)	-0.01	28 (3%) 46 54	29, 41, 68, 99	0

All (28) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	108	GLY	5.4
1	A	40	ARG	4.1
1	B	38	GLU	3.7
1	D	40	ARG	3.7
1	B	40	ARG	3.6
1	B	219	GLN	3.3
1	D	39	GLU	3.2
1	D	109	TYR	3.2
1	A	38	GLU	3.1
1	B	35	LEU	3.0
1	C	91	TYR	2.8
1	C	94	LEU	2.8
1	B	112	THR	2.7
1	A	107	GLU	2.6
1	C	92	SER	2.6
1	B	138	GLU	2.6
1	A	36	LYS	2.5
1	B	109	TYR	2.5
1	A	112	THR	2.4
1	B	36	LYS	2.4
1	B	108	GLY	2.4

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Mol	Chain	Res	Type	RSRZ
1	D	0	HIS	2.3
1	B	113	ILE	2.1
1	C	109	TYR	2.1
1	A	33	MET	2.1
1	A	106	LEU	2.1
1	A	110	LYS	2.1
1	B	37	ASP	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 carbohydrates in this entry.

## 6.4 Ligands [i](#)

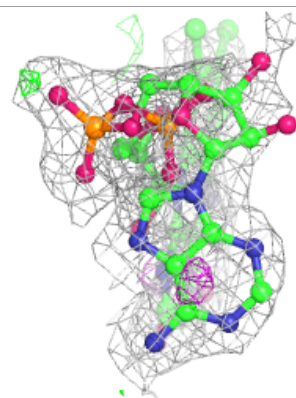
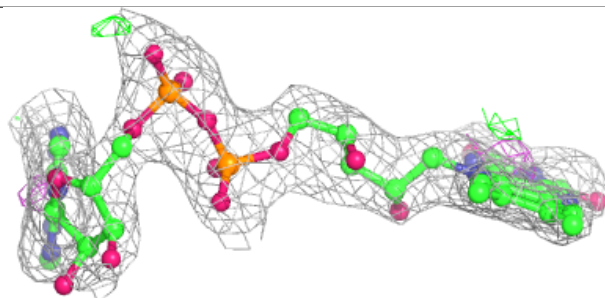
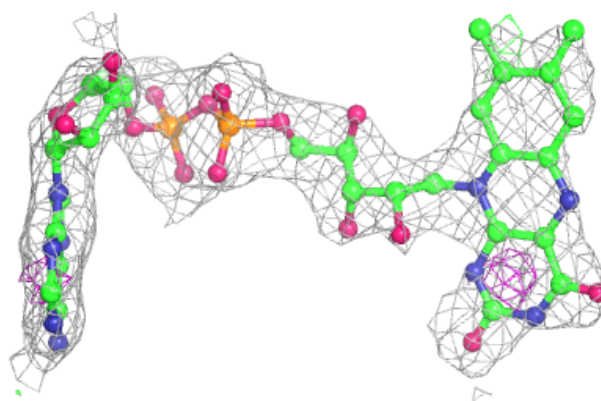
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
3	FDA	B	302	53/53	0.89	0.16	39,64,77,86	0
3	FDA	D	301	53/53	0.90	0.15	35,60,69,82	0
3	FDA	A	302	53/53	0.91	0.15	40,61,83,85	0
3	FDA	C	301	53/53	0.91	0.14	35,60,69,74	0
2	DU	D	302	20/20	0.93	0.13	44,51,71,77	0
2	DU	C	302	20/20	0.95	0.13	46,55,69,82	0
2	DU	A	301	20/20	0.97	0.12	37,43,58,60	0
2	DU	B	301	20/20	0.97	0.12	37,43,51,52	0

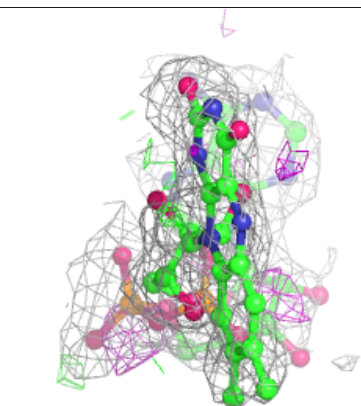
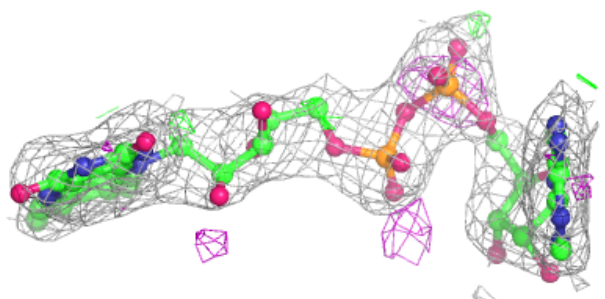
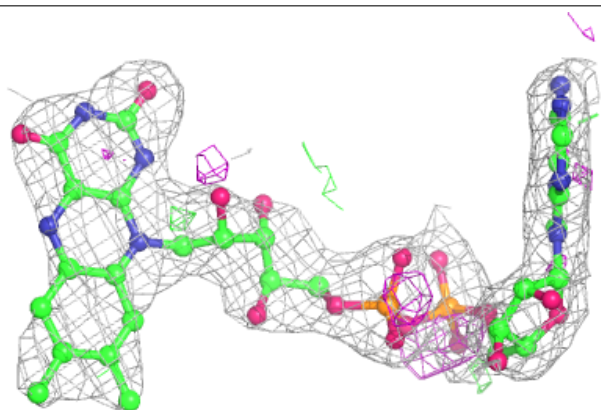
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

**Electron density around 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)

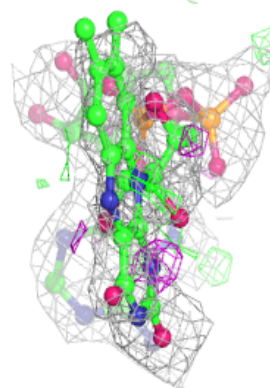
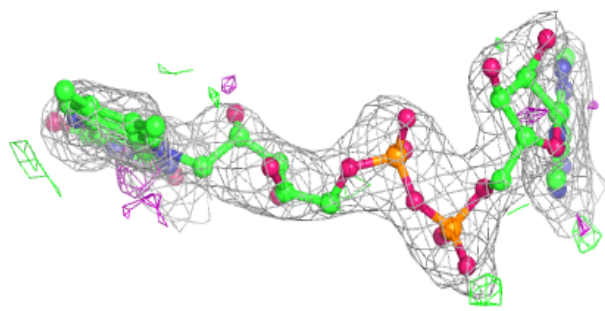
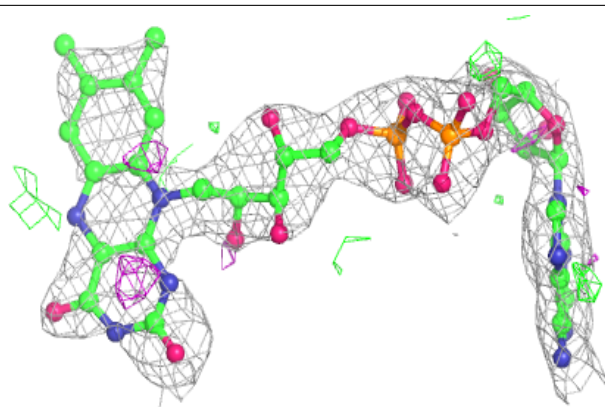
**Electron density around FDA D 301:**

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

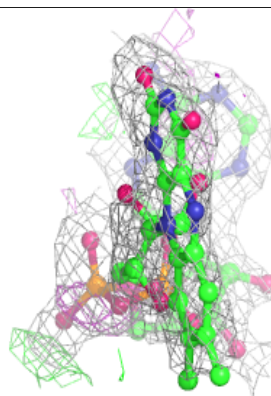
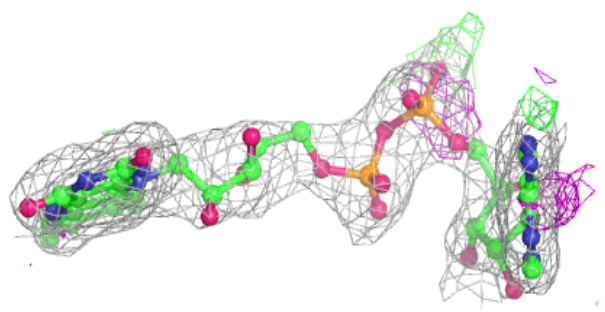
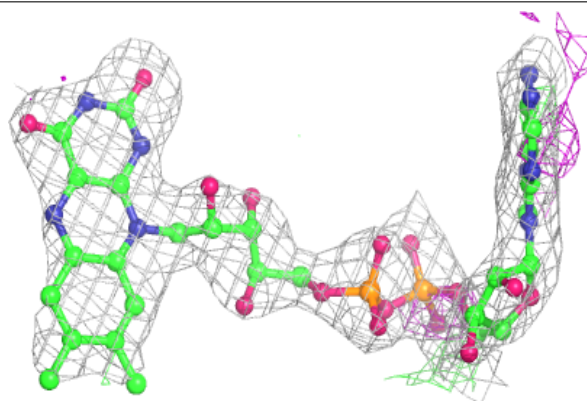


**Electron density around FDA A 302:**

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

**Electron density around FDA 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)



## 6.5 Other polymers [i](#)

There are no such residues in this entry.