



wwPDB X-ray Structure Validation Summary Report ⓘ

May 14, 2020 – 06:28 pm BST

PDB ID : 3UKK
Title : Crystal structure of R182K-UDP-galactopuranose mutase from *Aspergillus fumigatus* in complex with UDPgalp
Authors : Van Straaten, K.E.; Sanders, D.A.R.
Deposited on : 2011-11-09
Resolution : 2.75 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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

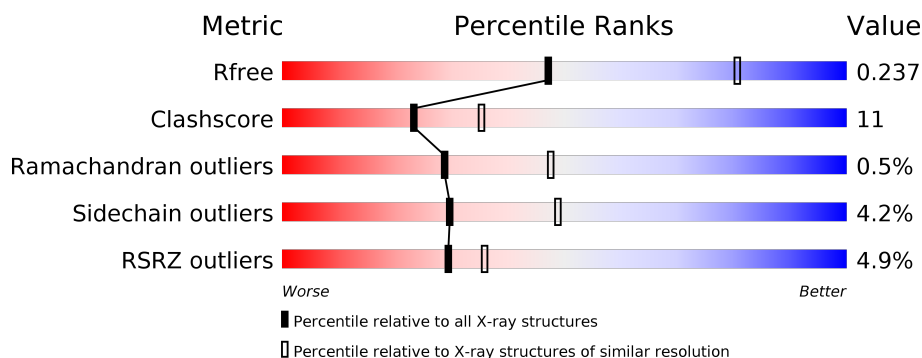
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.75 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	1235 (2.78-2.74)
Clashscore	141614	1277 (2.78-2.74)
Ramachandran outliers	138981	1257 (2.78-2.74)
Sidechain outliers	138945	1257 (2.78-2.74)
RSRZ outliers	127900	1207 (2.78-2.74)

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	517	<div> <div>4%</div> <div>78%</div> <div>21%</div> <div>.</div> </div>
1	B	517	<div> <div>4%</div> <div>74%</div> <div>23%</div> <div>..</div> </div>
1	C	517	<div> <div>8%</div> <div>72%</div> <div>26%</div> <div>.</div> </div>
1	D	517	<div> <div>4%</div> <div>73%</div> <div>24%</div> <div>..</div> </div>

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 16649 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 UDP-galactopyranose mutase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	517	Total	C	N	O	S	0	0	0
			4071	2581	704	765	21			
1	B	510	Total	C	N	O	S	0	0	0
			4002	2540	685	756	21			
1	C	516	Total	C	N	O	S	0	0	0
			4061	2575	701	764	21			
1	D	513	Total	C	N	O	S	0	0	0
			4031	2557	692	761	21			

There are 36 discrepancies between the modelled and reference sequences:

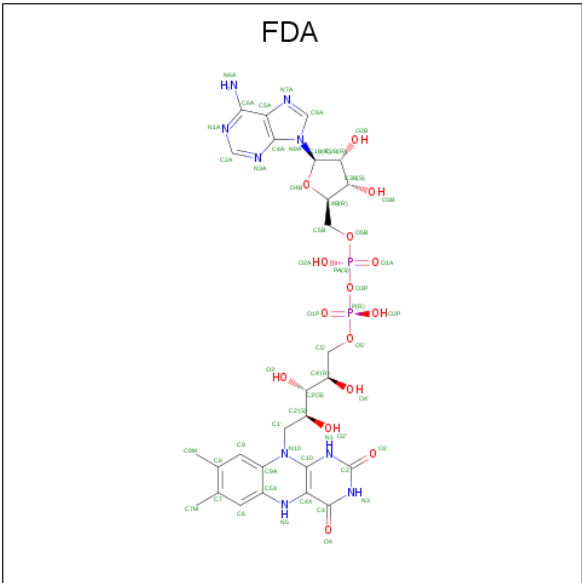
Chain	Residue	Modelled	Actual	Comment	Reference
A	182	LYS	ARG	ENGINEERED MUTATION	UNP Q4W1X2
A	511	LEU	-	EXPRESSION TAG	UNP Q4W1X2
A	512	GLU	-	EXPRESSION TAG	UNP Q4W1X2
A	513	HIS	-	EXPRESSION TAG	UNP Q4W1X2
A	514	HIS	-	EXPRESSION TAG	UNP Q4W1X2
A	515	HIS	-	EXPRESSION TAG	UNP Q4W1X2
A	516	HIS	-	EXPRESSION TAG	UNP Q4W1X2
A	517	HIS	-	EXPRESSION TAG	UNP Q4W1X2
A	518	HIS	-	EXPRESSION TAG	UNP Q4W1X2
B	182	LYS	ARG	ENGINEERED MUTATION	UNP Q4W1X2
B	511	LEU	-	EXPRESSION TAG	UNP Q4W1X2
B	512	GLU	-	EXPRESSION TAG	UNP Q4W1X2
B	513	HIS	-	EXPRESSION TAG	UNP Q4W1X2
B	514	HIS	-	EXPRESSION TAG	UNP Q4W1X2
B	515	HIS	-	EXPRESSION TAG	UNP Q4W1X2
B	516	HIS	-	EXPRESSION TAG	UNP Q4W1X2
B	517	HIS	-	EXPRESSION TAG	UNP Q4W1X2
B	518	HIS	-	EXPRESSION TAG	UNP Q4W1X2
C	182	LYS	ARG	ENGINEERED MUTATION	UNP Q4W1X2
C	511	LEU	-	EXPRESSION TAG	UNP Q4W1X2
C	512	GLU	-	EXPRESSION TAG	UNP Q4W1X2

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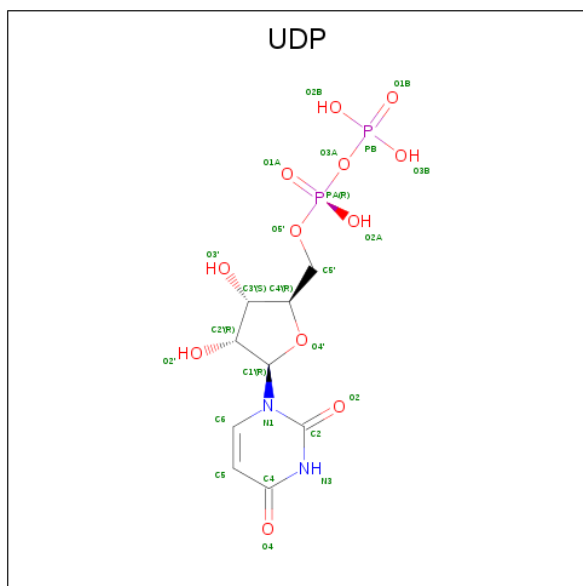
Chain	Residue	Modelled	Actual	Comment	Reference
C	513	HIS	-	EXPRESSION TAG	UNP Q4W1X2
C	514	HIS	-	EXPRESSION TAG	UNP Q4W1X2
C	515	HIS	-	EXPRESSION TAG	UNP Q4W1X2
C	516	HIS	-	EXPRESSION TAG	UNP Q4W1X2
C	517	HIS	-	EXPRESSION TAG	UNP Q4W1X2
C	518	HIS	-	EXPRESSION TAG	UNP Q4W1X2
D	182	LYS	ARG	ENGINEERED MUTATION	UNP Q4W1X2
D	511	LEU	-	EXPRESSION TAG	UNP Q4W1X2
D	512	GLU	-	EXPRESSION TAG	UNP Q4W1X2
D	513	HIS	-	EXPRESSION TAG	UNP Q4W1X2
D	514	HIS	-	EXPRESSION TAG	UNP Q4W1X2
D	515	HIS	-	EXPRESSION TAG	UNP Q4W1X2
D	516	HIS	-	EXPRESSION TAG	UNP Q4W1X2
D	517	HIS	-	EXPRESSION TAG	UNP Q4W1X2
D	518	HIS	-	EXPRESSION TAG	UNP Q4W1X2

- Molecule 2 is DIHYDROFLAVINE-ADENINE DINUCLEOTIDE (three-letter code: FDA) (formula: C₂₇H₃₅N₉O₁₅P₂).



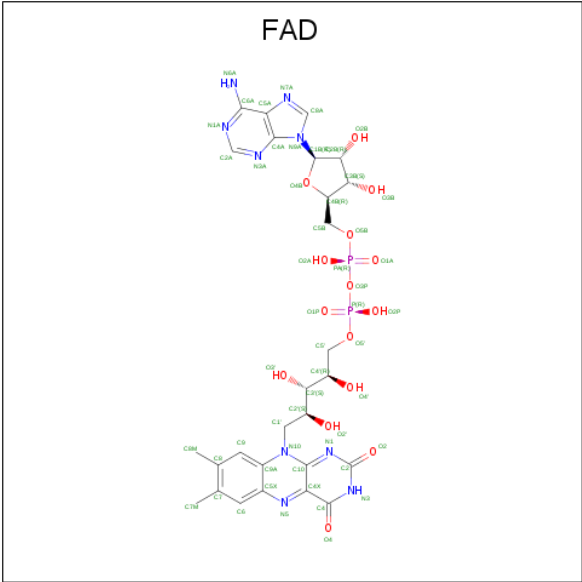
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
2	B	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
2	D	1	Total	C	N	O	P	0	0
			53	27	9	15	2		

- Molecule 3 is URIDINE-5'-DIPHOSPHATE (three-letter code: UDP) (formula: $C_9H_{14}N_2O_{12}P_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			25	9	2	12	2		
3	B	1	Total	C	N	O	P	0	0
			25	9	2	12	2		
3	C	1	Total	C	N	O	P	0	0
			25	9	2	12	2		
3	D	1	Total	C	N	O	P	0	0
			25	9	2	12	2		

- Molecule 4 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula: $C_{27}H_{33}N_9O_{15}P_2$).

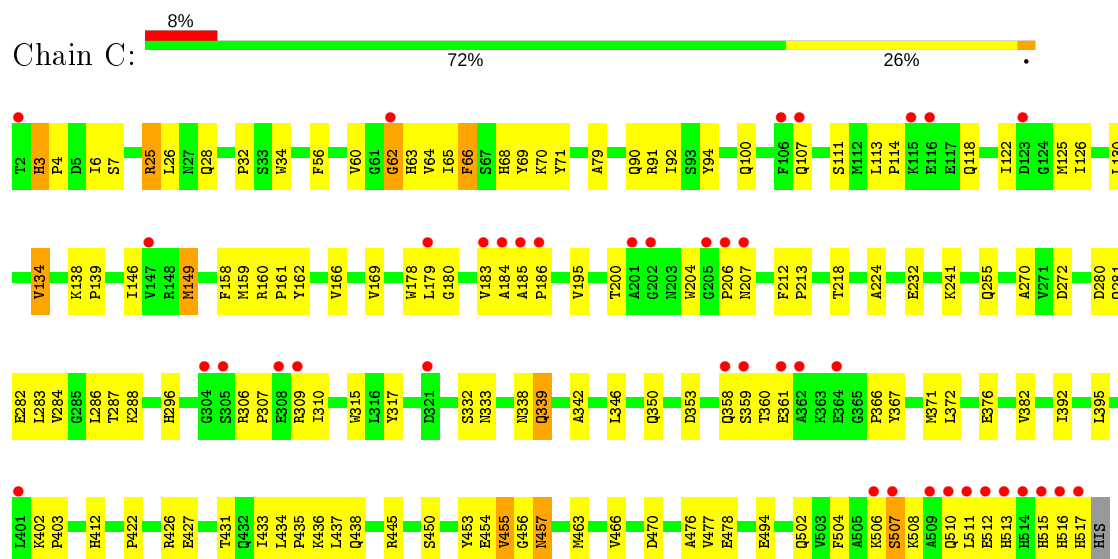


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	C	1	Total	C	N	O	P	0	0
			53	27	9	15	2		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	54	Total	O	0	0
			54	54		
5	B	53	Total	O	0	0
			53	53		
5	C	37	Total	O	0	0
			37	37		
5	D	28	Total	O	0	0
			28	28		

• Molecule 1: UDP-galactopyranose mutase



• Molecule 1: UDP-galactopyranose mutase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	70.66Å 123.66Å 156.37Å 90.00° 103.26° 90.00°	Depositor
Resolution (Å)	47.99 – 2.75 47.99 – 2.75	Depositor EDS
% Data completeness (in resolution range)	93.4 (47.99-2.75) 93.5 (47.99-2.75)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.30 (at 2.77Å)	Xtriage
Refinement program	PHENIX 1.7.1_743	Depositor
R, R_{free}	0.183 , 0.241 0.180 , 0.237	Depositor DCC
R_{free} test set	3182 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	46.5	Xtriage
Anisotropy	0.034	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 54.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.031 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	16649	wwPDB-VP
Average B, all atoms (Å ²)	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.04% 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: FDA, UDP, FAD

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	0.26	0/4176	0.43	0/5675
1	B	0.26	0/4101	0.43	0/5573
1	C	0.27	0/4165	0.44	0/5660
1	D	0.25	0/4132	0.43	0/5615
All	All	0.26	0/16574	0.43	0/22523

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4071	0	3973	87	0
1	B	4002	0	3925	92	0
1	C	4061	0	3966	104	0
1	D	4031	0	3945	96	0
2	A	53	0	30	2	0
2	B	53	0	30	5	0
2	D	53	0	30	3	0
3	A	25	0	11	1	0
3	B	25	0	11	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	C	25	0	11	2	0
3	D	25	0	11	2	0
4	C	53	0	30	3	0
5	A	54	0	0	1	0
5	B	53	0	0	6	0
5	C	37	0	0	1	0
5	D	28	0	0	4	0
All	All	16649	0	15973	366	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 11.

The worst 5 of 366 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:C:66:PHE:CE2	1:C:206:PRO:HB3	2.00	0.95
1:A:456:GLY:O	1:A:457:ASN:HB2	1.71	0.90
1:C:456:GLY:O	1:C:457:ASN:HB2	1.72	0.88
1:D:63:HIS:HB2	1:D:218:THR:HG21	1.54	0.87
1:C:91:ARG:HD3	1:C:207:ASN:HB2	1.55	0.85

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	515/517 (100%)	494 (96%)	19 (4%)	2 (0%)	34 53
1	B	508/517 (98%)	487 (96%)	18 (4%)	3 (1%)	25 42
1	C	514/517 (99%)	486 (95%)	25 (5%)	3 (1%)	25 42
1	D	511/517 (99%)	479 (94%)	29 (6%)	3 (1%)	25 42

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	2048/2068 (99%)	1946 (95%)	91 (4%)	11 (0%)	29	47

5 of 11 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	457	ASN
1	B	457	ASN
1	C	457	ASN
1	D	457	ASN
1	B	208	ALA

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	439/439 (100%)	425 (97%)	14 (3%)	39	59
1	B	432/439 (98%)	415 (96%)	17 (4%)	32	52
1	C	438/439 (100%)	416 (95%)	22 (5%)	24	42
1	D	435/439 (99%)	414 (95%)	21 (5%)	25	44
All	All	1744/1756 (99%)	1670 (96%)	74 (4%)	30	49

5 of 74 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	25	ARG
1	C	200	THR
1	D	450	SER
1	C	66	PHE
1	C	134	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 55 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	412	HIS
1	C	118	GLN
1	D	384	GLN
1	B	471	ASN
1	C	28	GLN

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	UDP	D	602	-	20,26,26	1.17	1 (5%)	25,40,40	1.14	2 (8%)
2	FDA	A	601	-	51,58,58	2.24	22 (43%)	60,89,89	2.47	17 (28%)
3	UDP	C	602	-	20,26,26	1.15	1 (5%)	25,40,40	1.04	2 (8%)
3	UDP	A	602	-	20,26,26	1.16	1 (5%)	25,40,40	1.10	3 (12%)
3	UDP	B	602	-	20,26,26	1.16	1 (5%)	25,40,40	1.00	2 (8%)
2	FDA	D	601	-	51,58,58	2.24	23 (45%)	60,89,89	2.31	19 (31%)
2	FDA	B	601	-	51,58,58	2.24	24 (47%)	60,89,89	2.34	17 (28%)
4	FAD	C	601	-	51,58,58	2.45	15 (29%)	60,89,89	1.81	16 (26%)

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	UDP	D	602	-	-	3/14/32/32	0/2/2/2
2	FDA	A	601	-	-	6/30/50/50	0/6/6/6
3	UDP	C	602	-	-	5/14/32/32	0/2/2/2
3	UDP	A	602	-	-	4/14/32/32	0/2/2/2
3	UDP	B	602	-	-	7/14/32/32	0/2/2/2
2	FDA	D	601	-	-	4/30/50/50	0/6/6/6
2	FDA	B	601	-	-	5/30/50/50	0/6/6/6
4	FAD	C	601	-	-	7/30/50/50	0/6/6/6

The worst 5 of 88 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	C	601	FAD	C10-N1	8.48	1.44	1.33
4	C	601	FAD	C1'-N10	-6.77	1.41	1.48
2	A	601	FDA	P-O1P	-5.70	1.30	1.50
2	B	601	FDA	P-O1P	-5.62	1.31	1.50
2	D	601	FDA	P-O1P	-5.59	1.31	1.50

The worst 5 of 78 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	601	FDA	C4-N3-C2	8.56	122.37	115.14
2	D	601	FDA	C4-N3-C2	8.41	122.24	115.14
2	B	601	FDA	C4-N3-C2	8.06	121.94	115.14
2	A	601	FDA	C3B-C2B-C1B	-6.08	91.82	100.98
4	C	601	FAD	C4-N3-C2	5.84	120.08	115.14

There are no chirality outliers.

5 of 41 torsion outliers are listed below:

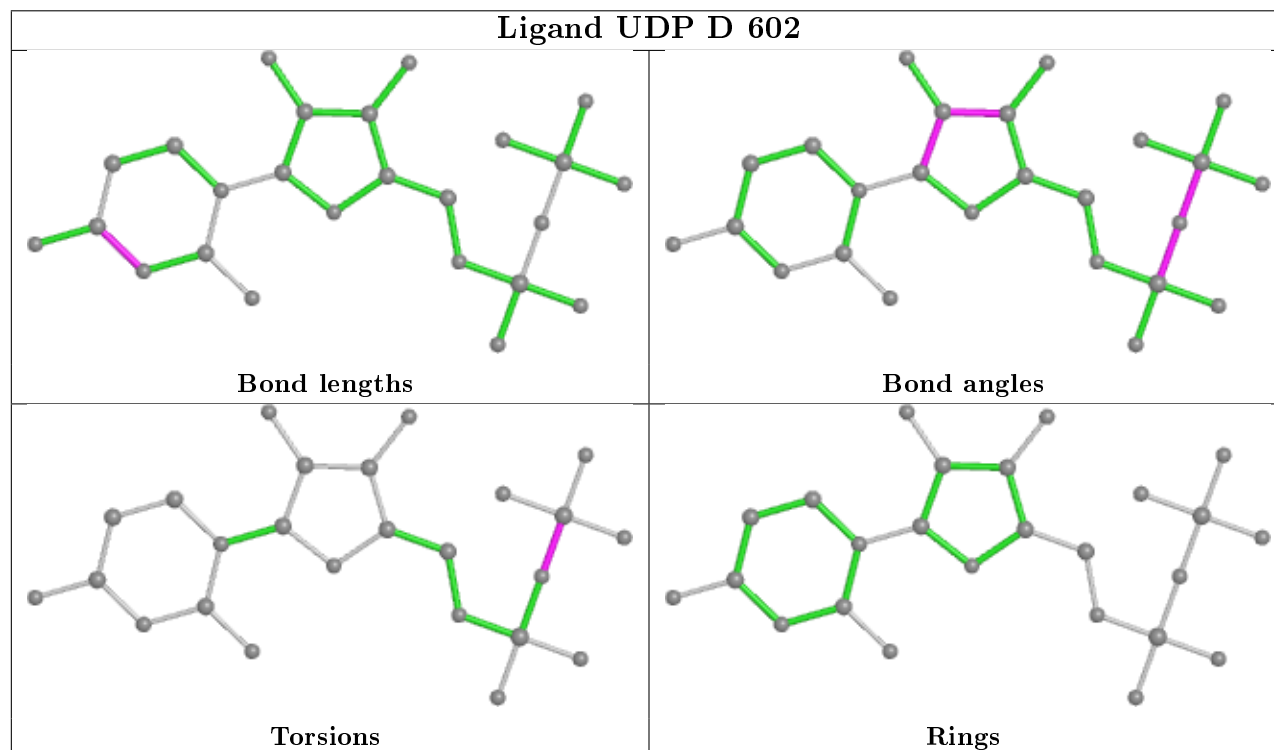
Mol	Chain	Res	Type	Atoms
3	D	602	UDP	PA-O3A-PB-O3B
2	A	601	FDA	N10-C1'-C2'-O2'
2	A	601	FDA	PA-O3P-P-O5'
3	C	602	UDP	C2'-C1'-N1-C6
3	C	602	UDP	O4'-C1'-N1-C6

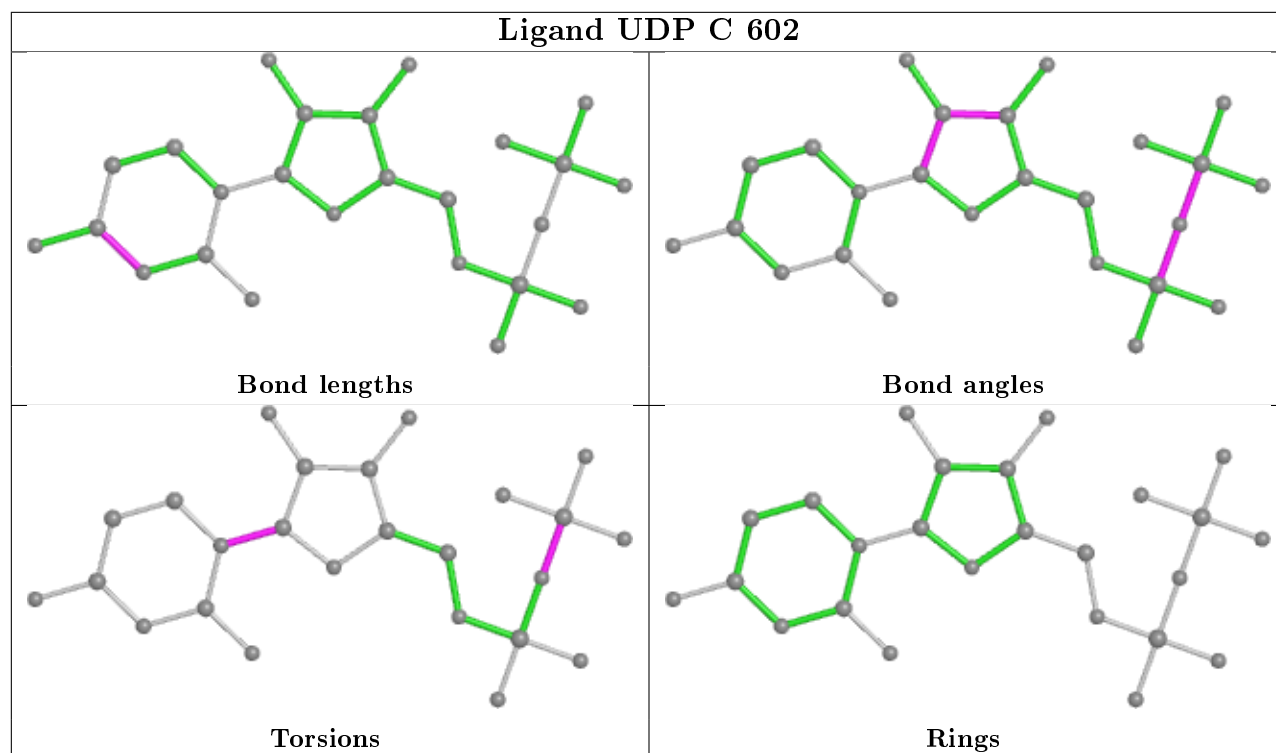
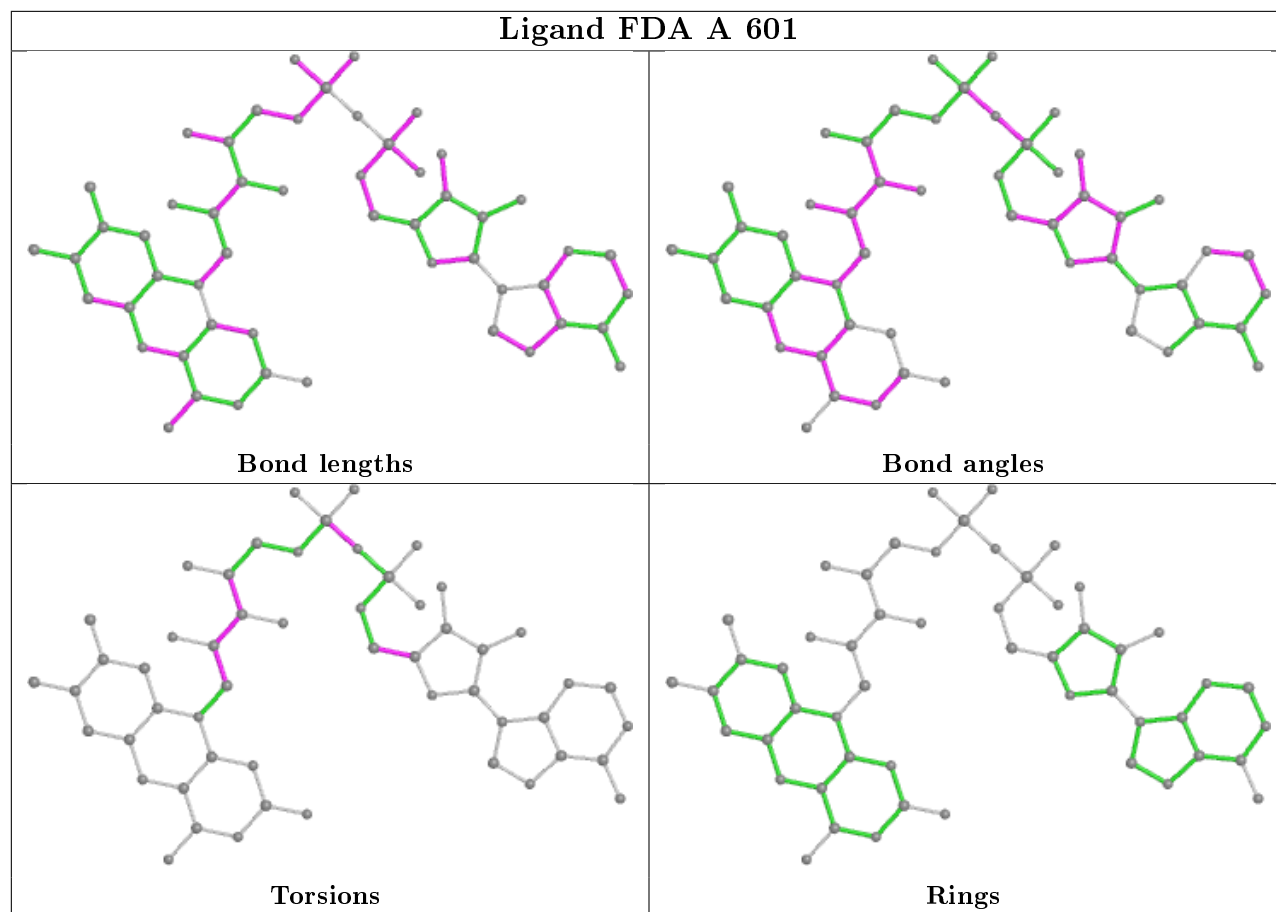
There are no ring outliers.

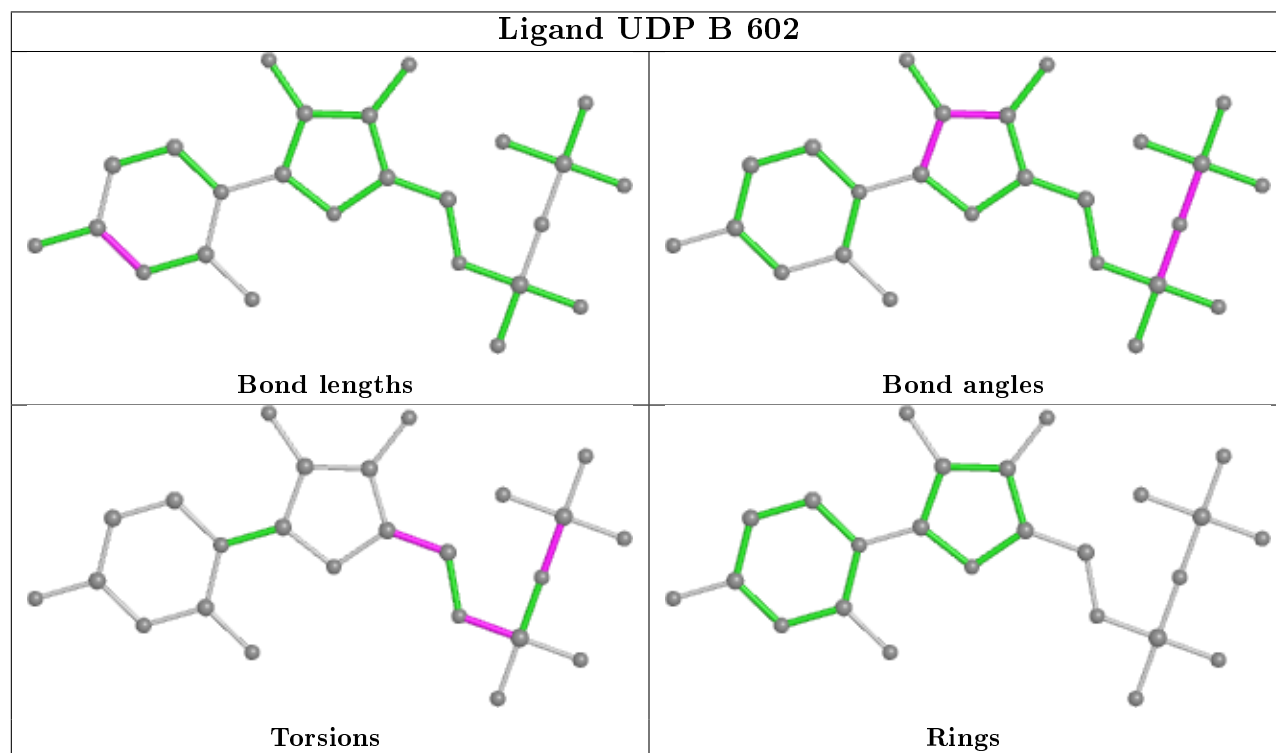
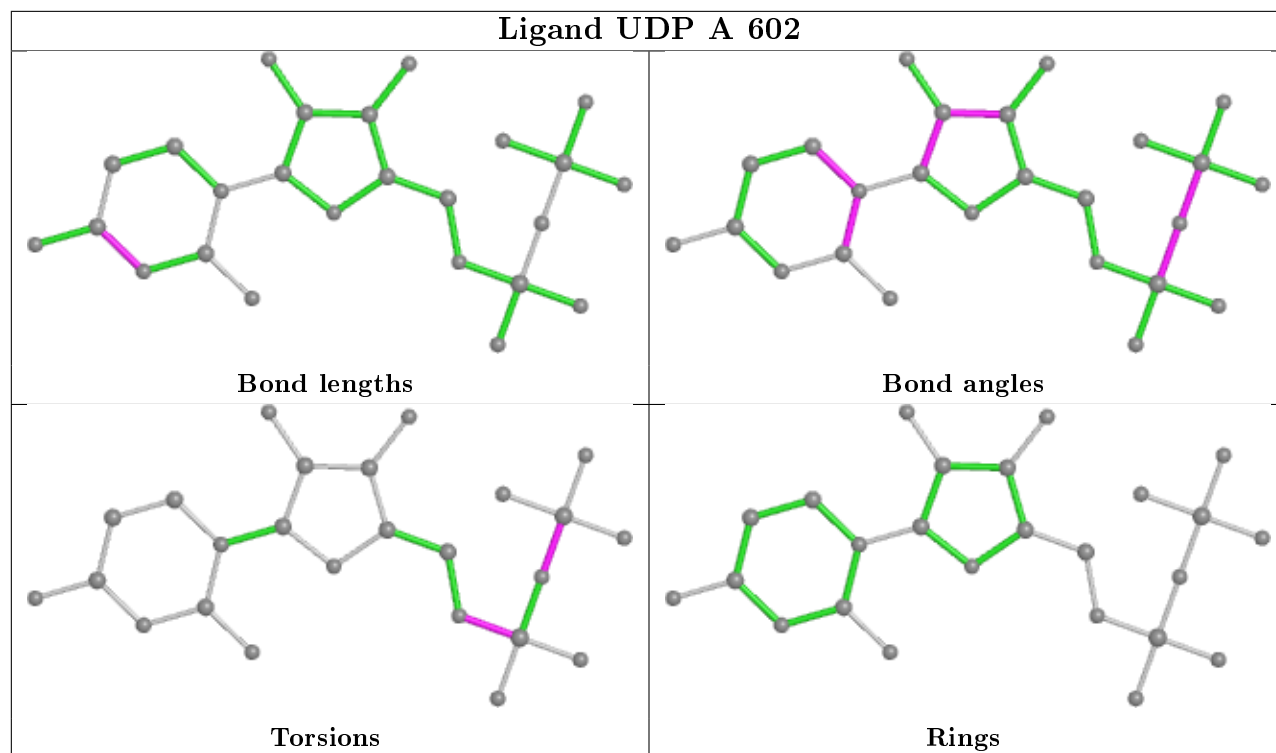
8 monomers are involved in 19 short contacts:

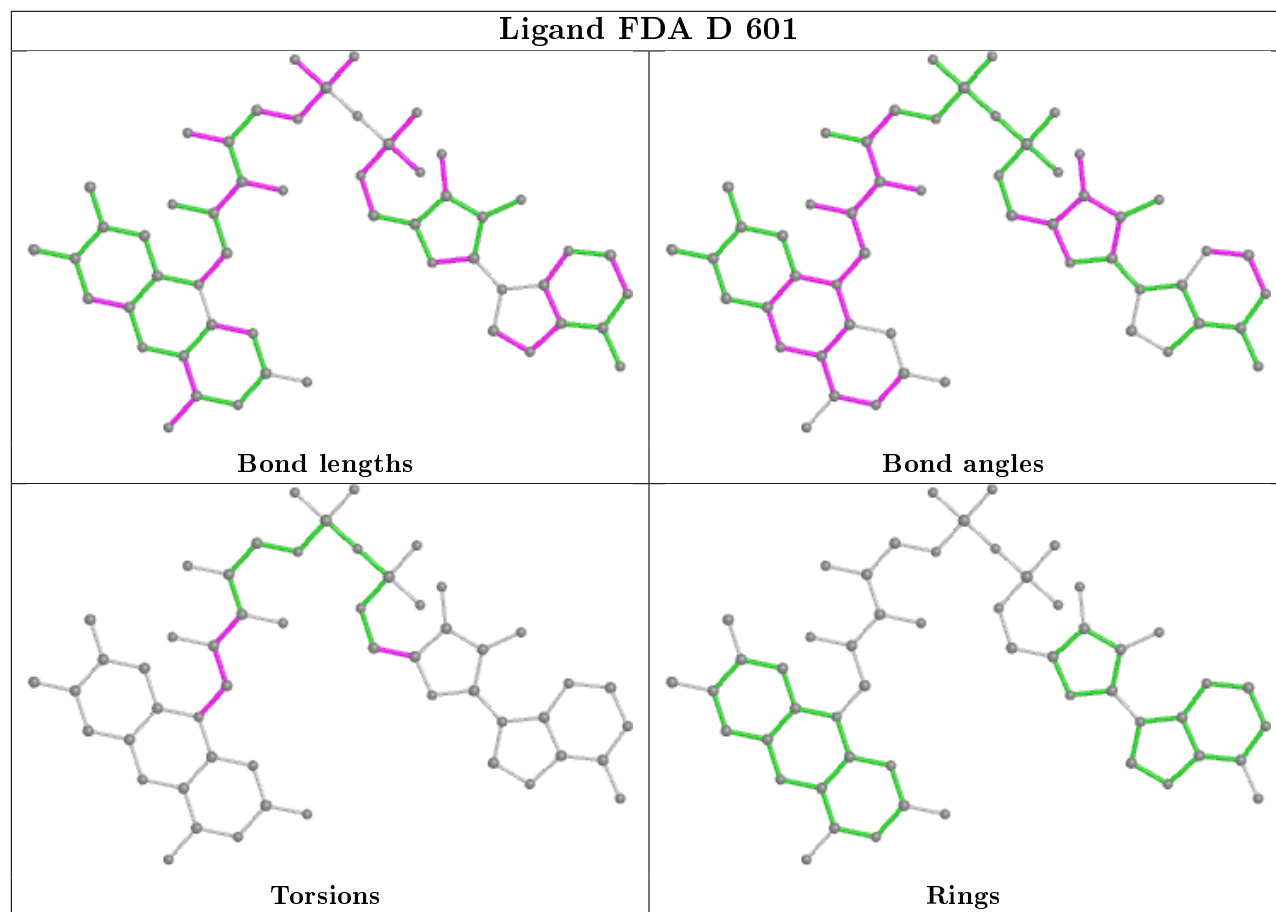
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	D	602	UDP	2	0
2	A	601	FDA	2	0
3	C	602	UDP	2	0
3	A	602	UDP	1	0
3	B	602	UDP	1	0
2	D	601	FDA	3	0
2	B	601	FDA	5	0
4	C	601	FAD	3	0

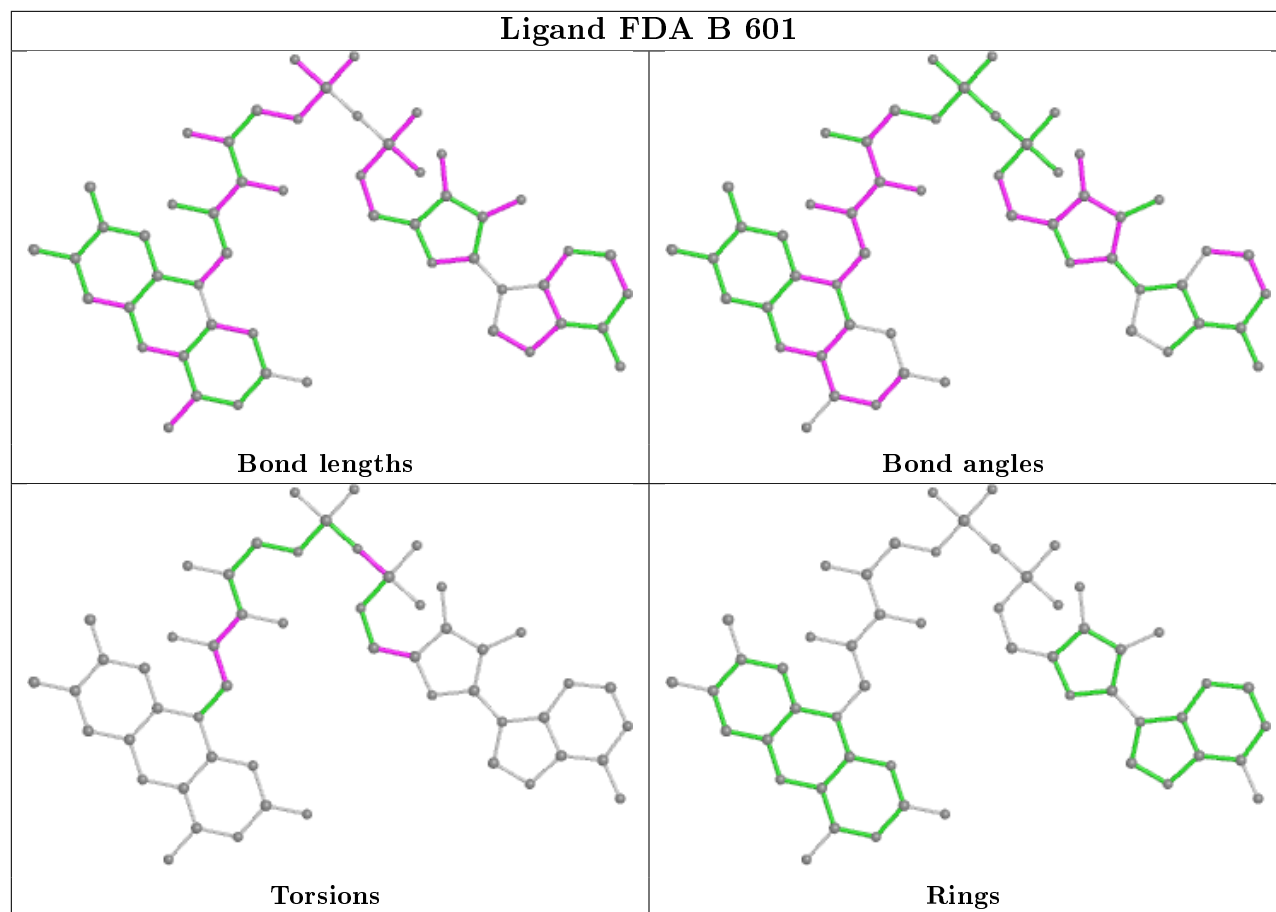
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

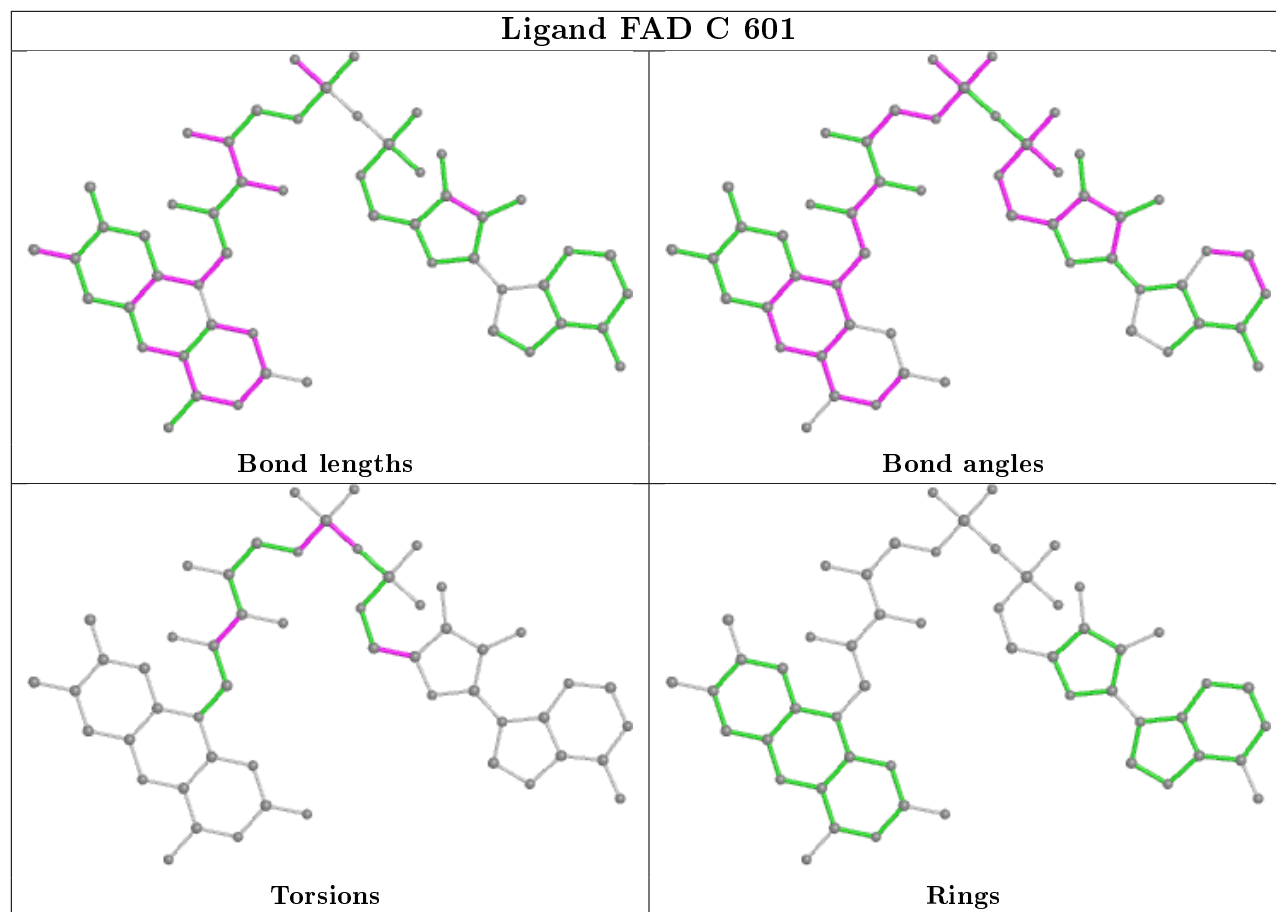












5.7 Other polymers [i](#)

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 [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	517/517 (100%)	0.29	21 (4%) 37 44	18, 44, 85, 112	0
1	B	510/517 (98%)	0.22	19 (3%) 41 49	17, 43, 83, 112	0
1	C	516/517 (99%)	0.29	40 (7%) 13 15	18, 45, 94, 152	0
1	D	513/517 (99%)	0.25	21 (4%) 37 44	19, 49, 88, 126	0
All	All	2056/2068 (99%)	0.26	101 (4%) 29 36	17, 46, 88, 152	0

The worst 5 of 101 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	507	SER	7.8
1	C	514	HIS	7.4
1	B	511	LEU	6.4
1	C	184	ALA	5.3
1	B	509	ALA	4.3

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

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

6.3 Carbohydrates [i](#)

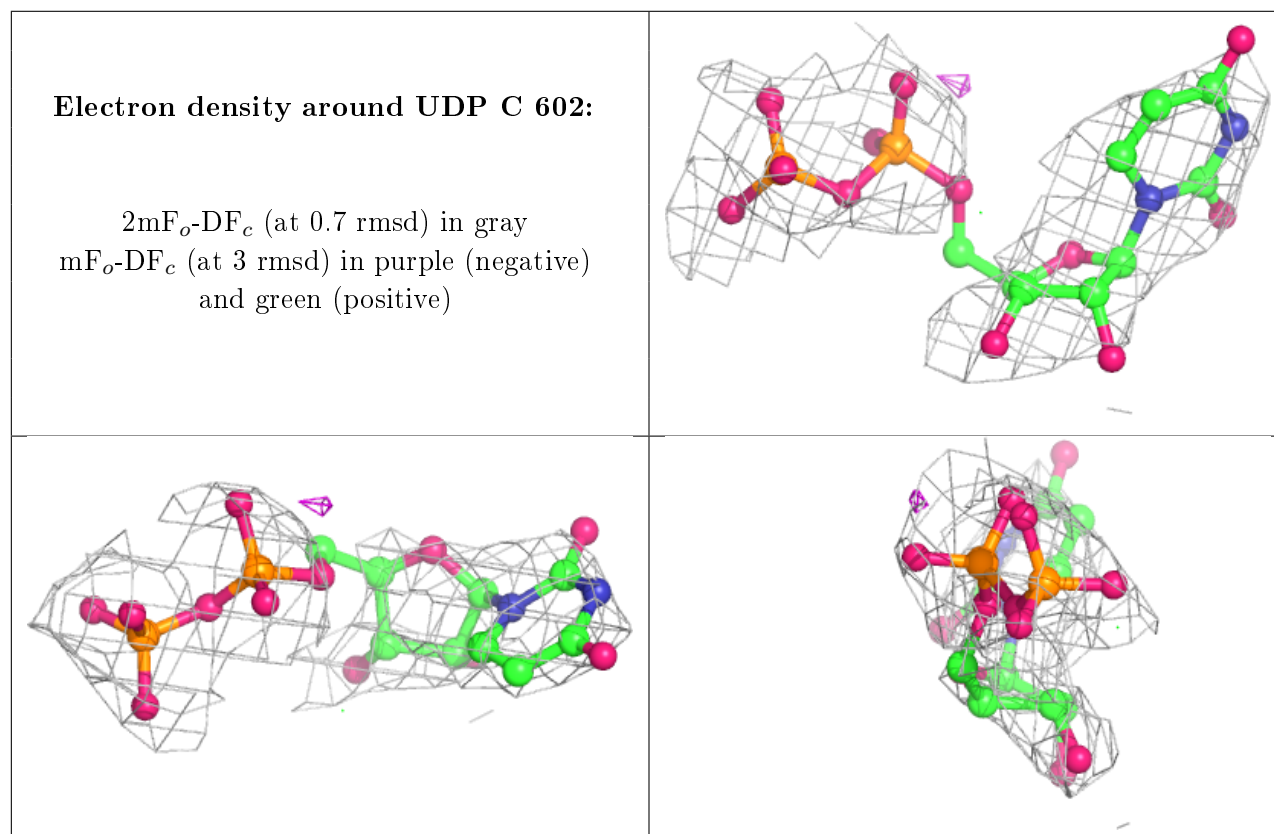
There are no carbohydrates in this entry.

6.4 Ligands [i](#)

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

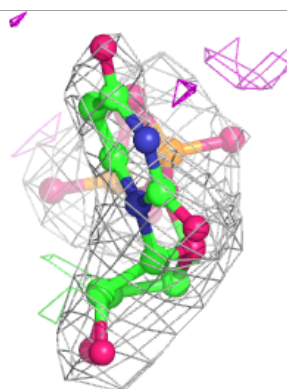
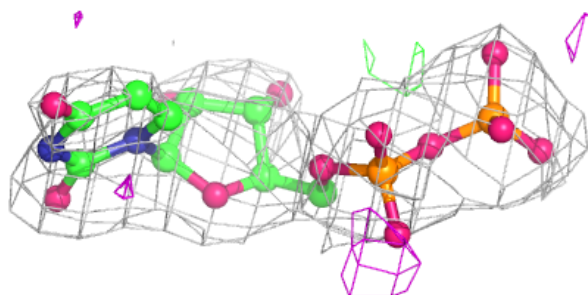
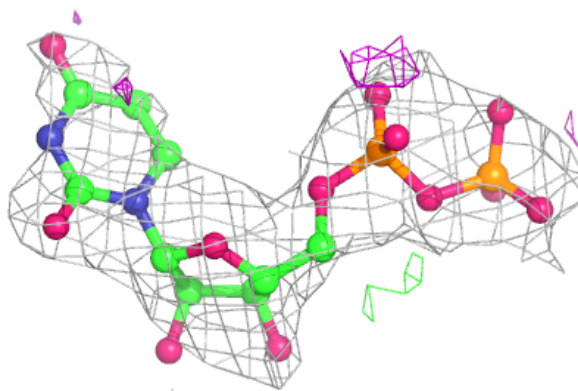
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	UDP	C	602	25/25	0.80	0.26	68,99,141,156	0
3	UDP	D	602	25/25	0.91	0.19	55,95,117,134	0
3	UDP	B	602	25/25	0.92	0.18	45,79,109,134	0
3	UDP	A	602	25/25	0.94	0.17	23,66,98,123	0
2	FDA	A	601	53/53	0.97	0.19	17,38,65,71	0
2	FDA	D	601	53/53	0.97	0.16	32,53,78,99	0
2	FDA	B	601	53/53	0.98	0.18	13,33,56,67	0
4	FAD	C	601	53/53	0.98	0.18	16,39,65,74	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.

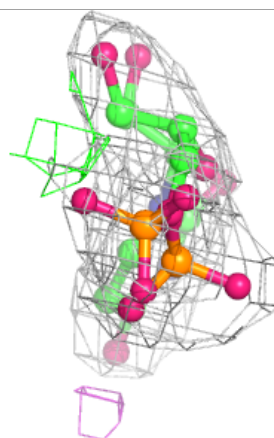
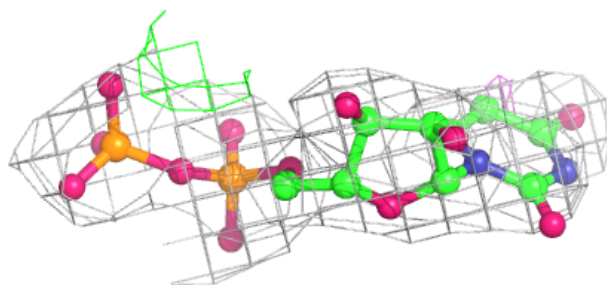
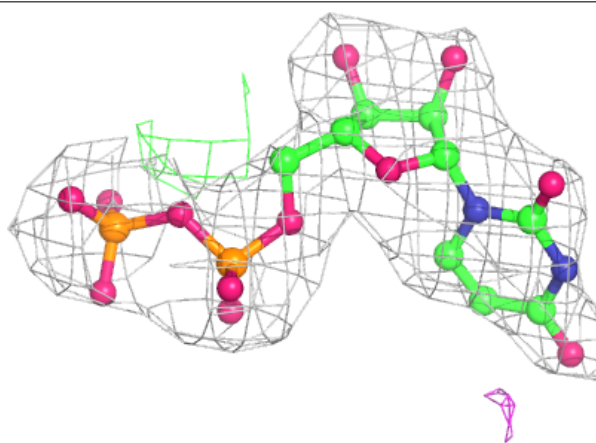


Electron density around UDP D 602:

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

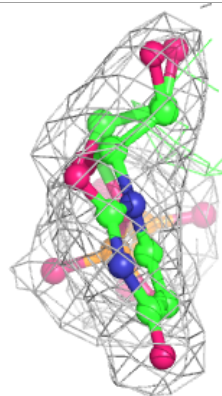
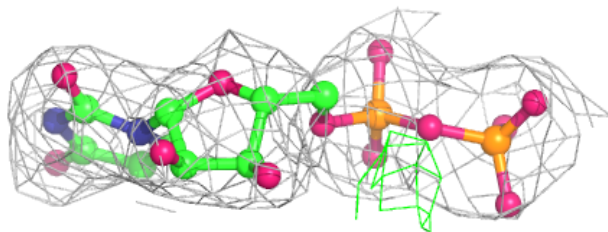
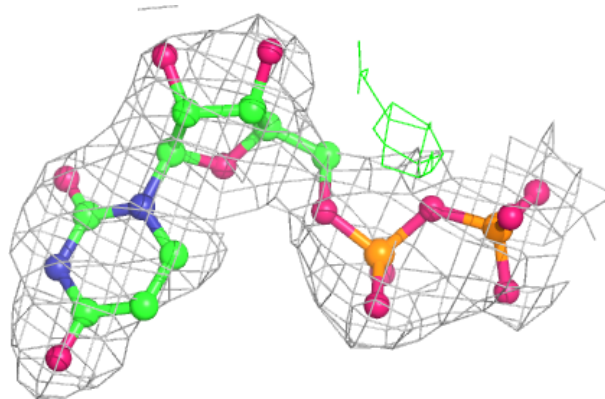
**Electron density around UDP B 602:**

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

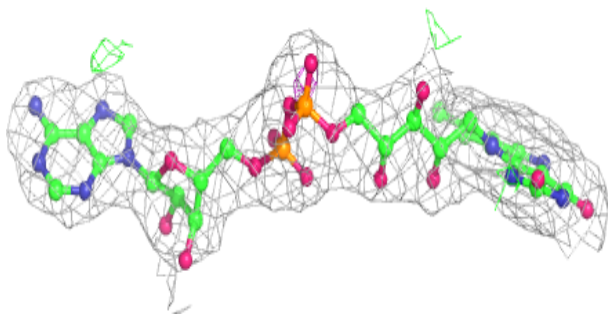
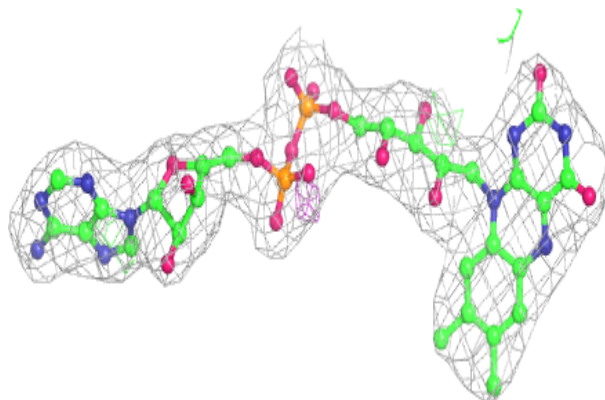


Electron density around UDP A 602:

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

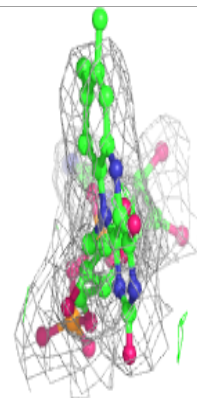
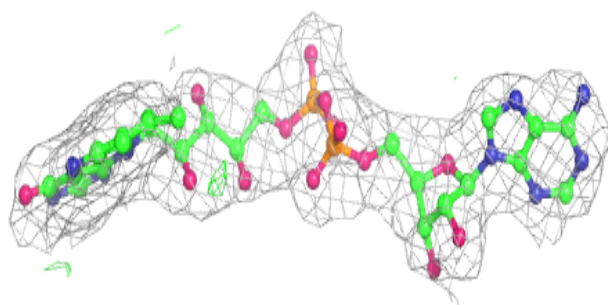
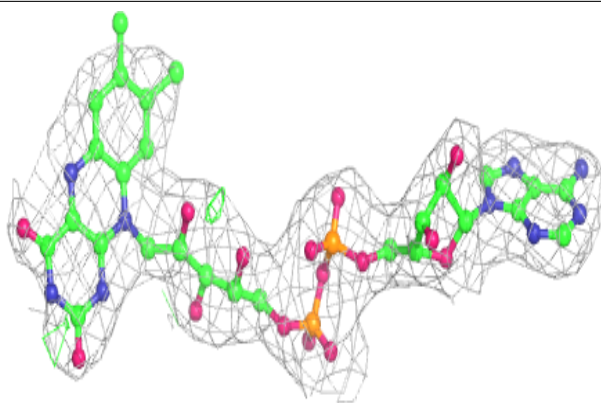
**Electron density around FDA A 601:**

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

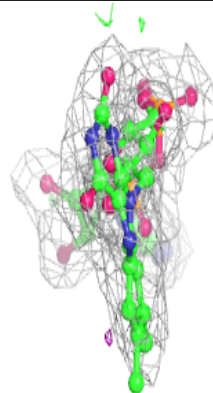
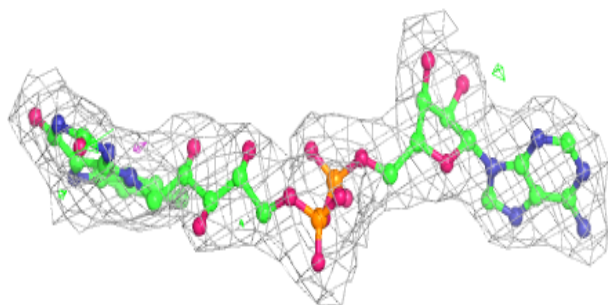
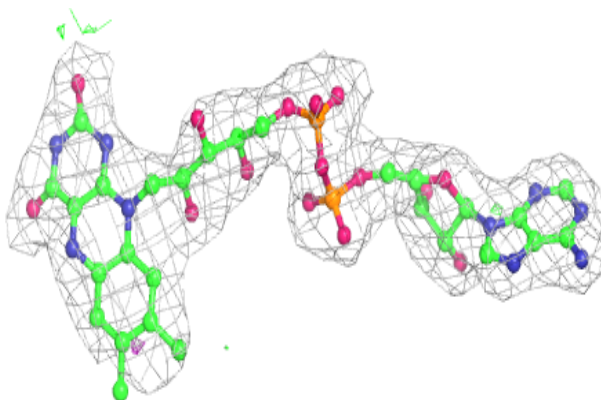


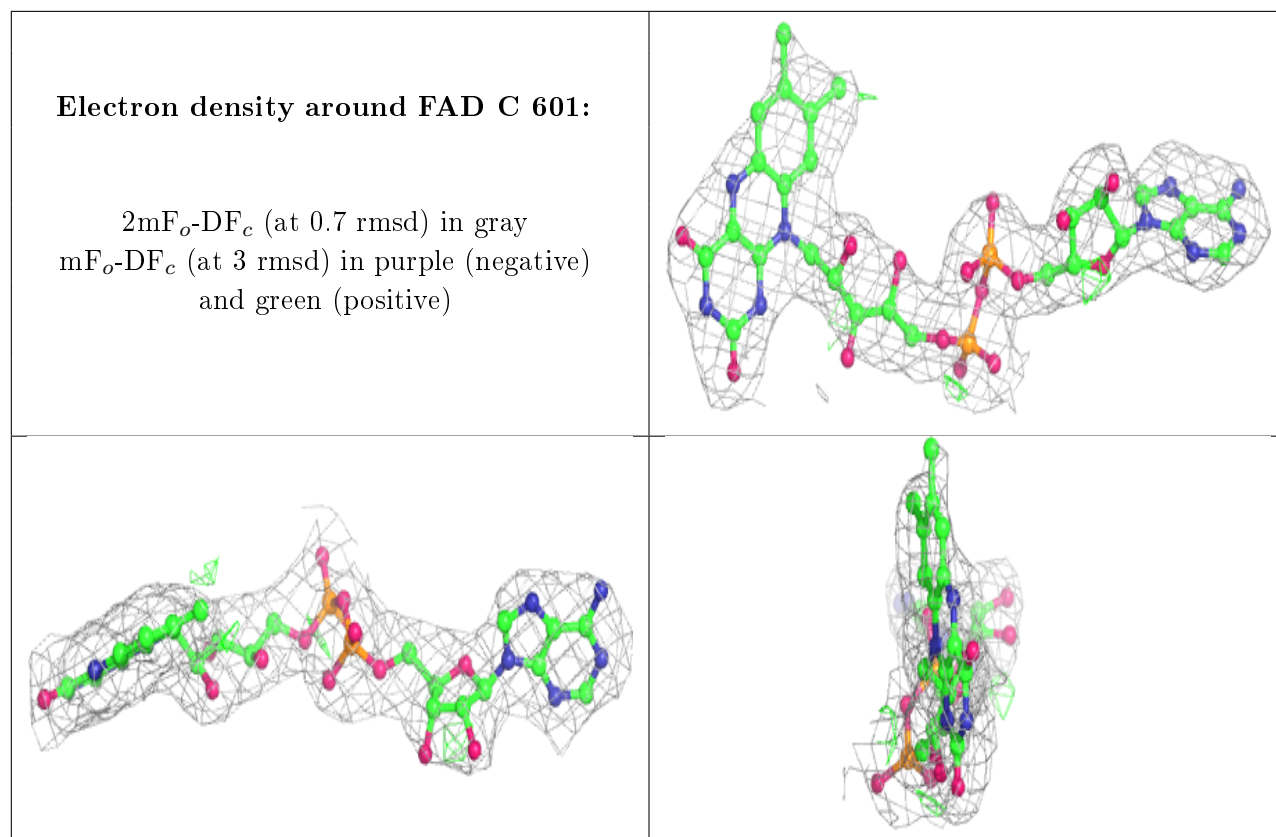
Electron density around FDA D 601:

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

**Electron density around FDA B 601:**

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





6.5 Other polymers [i](#)

There are no such residues in this entry.