



# wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 4, 2025 – 12:13 PM JST

PDB ID : 9LRJ  
Title : Glucosyl transferase NbUGT72AY1 co-crystallized with Scopoletin and UDP2Fglucose in the presence of retinol  
Authors : Arold, S.T.; Hameed, U.F.S.  
Deposited on : 2025-01-31  
Resolution : 3.11 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](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.

The types of validation reports are described at

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

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

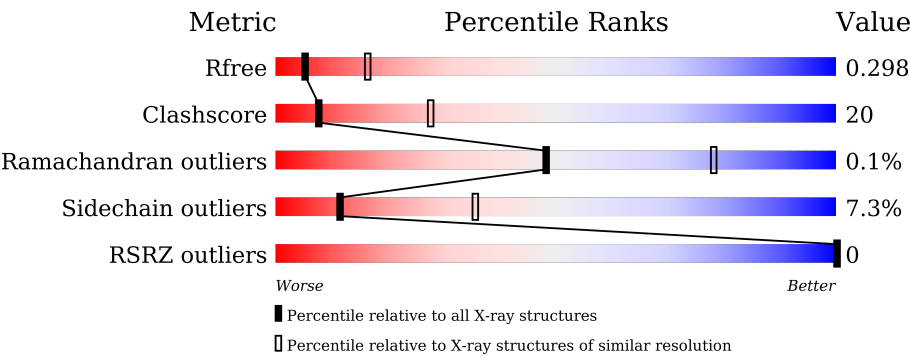
MolProbity : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.21  
EDS : 3.0  
buster-report : 1.1.7 (2018)  
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)  
CCP4 : 9.0.004 (Gargrove)  
Density-Fitness : 1.0.11  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.40

# 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 3.11 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))
$R_{free}$	164625	1668 (3.14-3.10)
Clashscore	180529	1788 (3.14-3.10)
Ramachandran outliers	177936	1696 (3.14-3.10)
Sidechain outliers	177891	1696 (3.14-3.10)
RSRZ outliers	164620	1668 (3.14-3.10)

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  $\geq 3$ , 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	479	
1	B	479	
1	C	479	
1	D	479	
1	E	479	
1	F	479	

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Mol	Chain	Length	Quality of chain
1	G	479	 <div>52% 37% 7% . .</div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	U2F	C	501	-	-	X	-
2	U2F	E	501	-	-	X	-

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 25741 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 Glycosyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	455	Total	C	N	O	S	0	0	0
			3600	2302	605	677	16			
1	B	459	Total	C	N	O	S	0	0	0
			3633	2321	611	685	16			
1	C	460	Total	C	N	O	S	0	0	0
			3640	2325	612	687	16			
1	D	460	Total	C	N	O	S	0	0	0
			3640	2325	612	687	16			
1	E	459	Total	C	N	O	S	0	0	0
			3633	2321	611	685	16			
1	F	459	Total	C	N	O	S	0	0	0
			3633	2321	611	685	16			
1	G	460	Total	C	N	O	S	0	0	0
			3640	2325	612	687	16			

There are 28 discrepancies between the modelled and reference sequences:

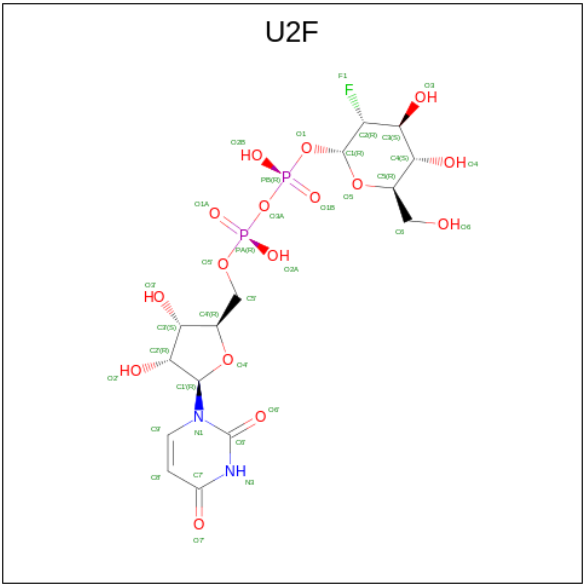
Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	expression tag	UNP A0A8K1ZRH3
A	0	PRO	-	expression tag	UNP A0A8K1ZRH3
A	1	LEU	-	expression tag	UNP A0A8K1ZRH3
A	2	GLY	-	expression tag	UNP A0A8K1ZRH3
B	-1	GLY	-	expression tag	UNP A0A8K1ZRH3
B	0	PRO	-	expression tag	UNP A0A8K1ZRH3
B	1	LEU	-	expression tag	UNP A0A8K1ZRH3
B	2	GLY	-	expression tag	UNP A0A8K1ZRH3
C	-1	GLY	-	expression tag	UNP A0A8K1ZRH3
C	0	PRO	-	expression tag	UNP A0A8K1ZRH3
C	1	LEU	-	expression tag	UNP A0A8K1ZRH3
C	2	GLY	-	expression tag	UNP A0A8K1ZRH3
D	-1	GLY	-	expression tag	UNP A0A8K1ZRH3
D	0	PRO	-	expression tag	UNP A0A8K1ZRH3
D	1	LEU	-	expression tag	UNP A0A8K1ZRH3

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Chain	Residue	Modelled	Actual	Comment	Reference
D	2	GLY	-	expression tag	UNP A0A8K1ZRH3
E	-1	GLY	-	expression tag	UNP A0A8K1ZRH3
E	0	PRO	-	expression tag	UNP A0A8K1ZRH3
E	1	LEU	-	expression tag	UNP A0A8K1ZRH3
E	2	GLY	-	expression tag	UNP A0A8K1ZRH3
F	-1	GLY	-	expression tag	UNP A0A8K1ZRH3
F	0	PRO	-	expression tag	UNP A0A8K1ZRH3
F	1	LEU	-	expression tag	UNP A0A8K1ZRH3
F	2	GLY	-	expression tag	UNP A0A8K1ZRH3
G	-1	GLY	-	expression tag	UNP A0A8K1ZRH3
G	0	PRO	-	expression tag	UNP A0A8K1ZRH3
G	1	LEU	-	expression tag	UNP A0A8K1ZRH3
G	2	GLY	-	expression tag	UNP A0A8K1ZRH3

- Molecule 2 is URIDINE-5'-DIPHOSPHATE-2-DEOXY-2-FLUORO-ALPHA-D-GLUCOS E (three-letter code: U2F) (formula: C<sub>15</sub>H<sub>23</sub>FN<sub>2</sub>O<sub>16</sub>P<sub>2</sub>) (labeled as "Ligand of Interest" by depositor).



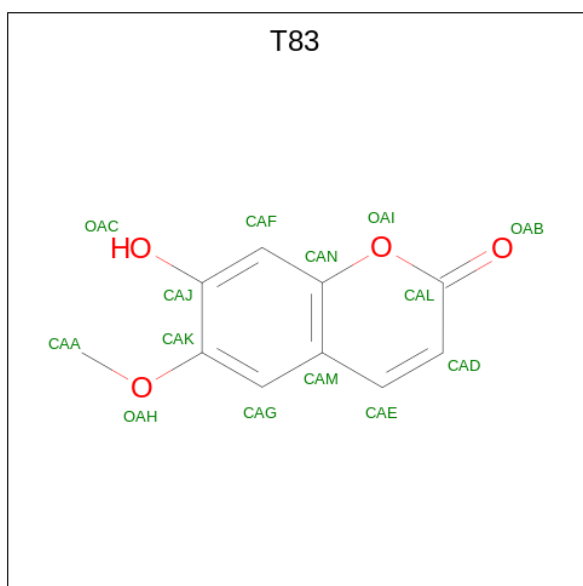
Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	A	1	Total	C	F	N	O	P	0	0
			36	15	1	2	16	2		
2	B	1	Total	C	F	N	O	P	0	0
			36	15	1	2	16	2		
2	C	1	Total	C	F	N	O	P	0	0
			36	15	1	2	16	2		
2	D	1	Total	C	F	N	O	P	0	0
			36	15	1	2	16	2		

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Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	E	1	Total	C	F	N	O	P	0	0
			36	15	1	2	16	2		
2	F	1	Total	C	F	N	O	P	0	0
			36	15	1	2	16	2		
2	G	1	Total	C	F	N	O	P	0	0
			36	15	1	2	16	2		

- Molecule 3 is 7-hydroxy-6-methoxy-2H-1-benzopyran-2-one (three-letter code: T83) (formula: C<sub>10</sub>H<sub>8</sub>O<sub>4</sub>) (labeled as "Ligand of Interest" by depositor).



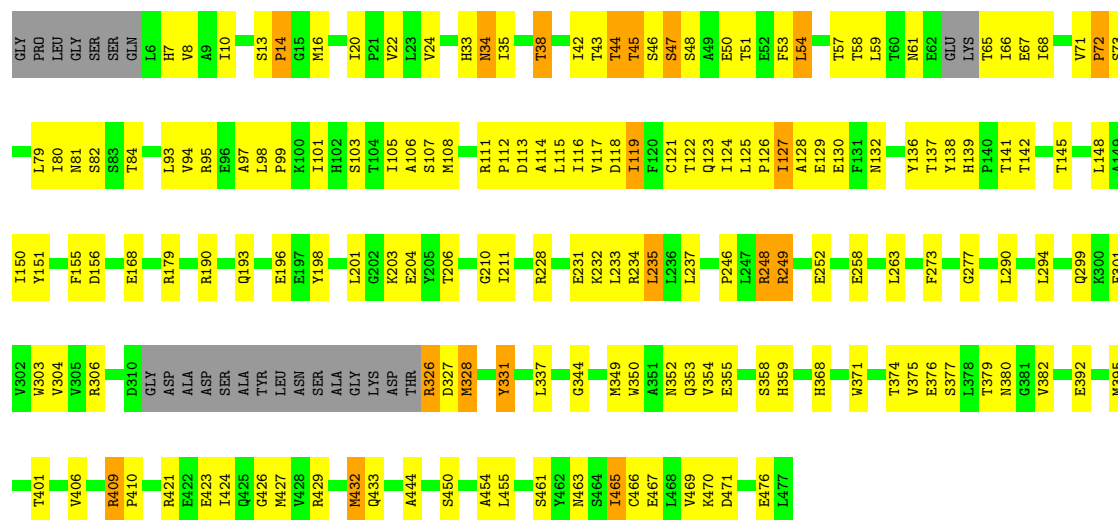
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			14	10	4		
3	B	1	Total	C	O	0	0
			14	10	4		
3	C	1	Total	C	O	0	0
			14	10	4		
3	D	1	Total	C	O	0	0
			14	10	4		
3	F	1	Total	C	O	0	0
			14	10	4		

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

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.

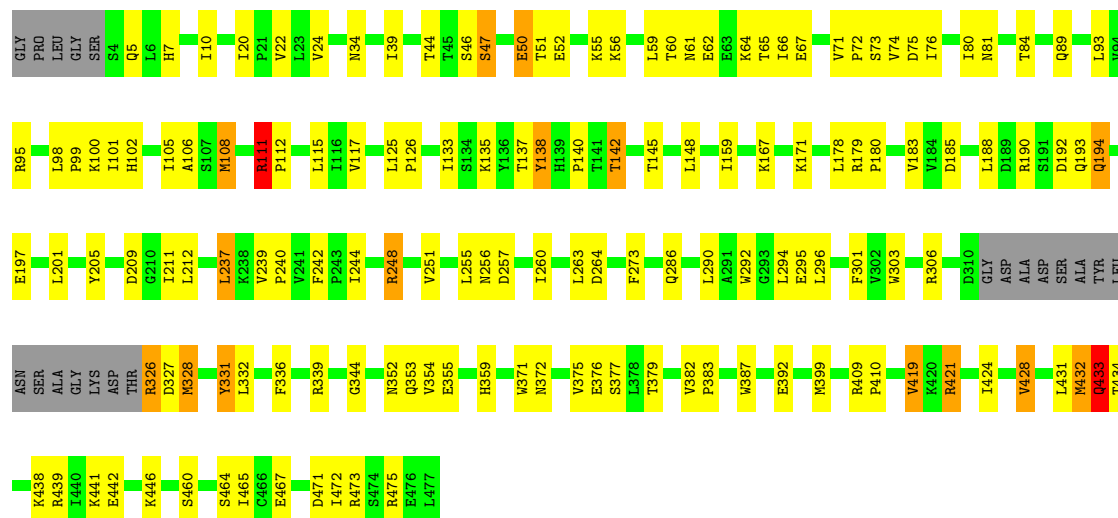
#### • Molecule 1: Glycosyltransferase

Chain A: 



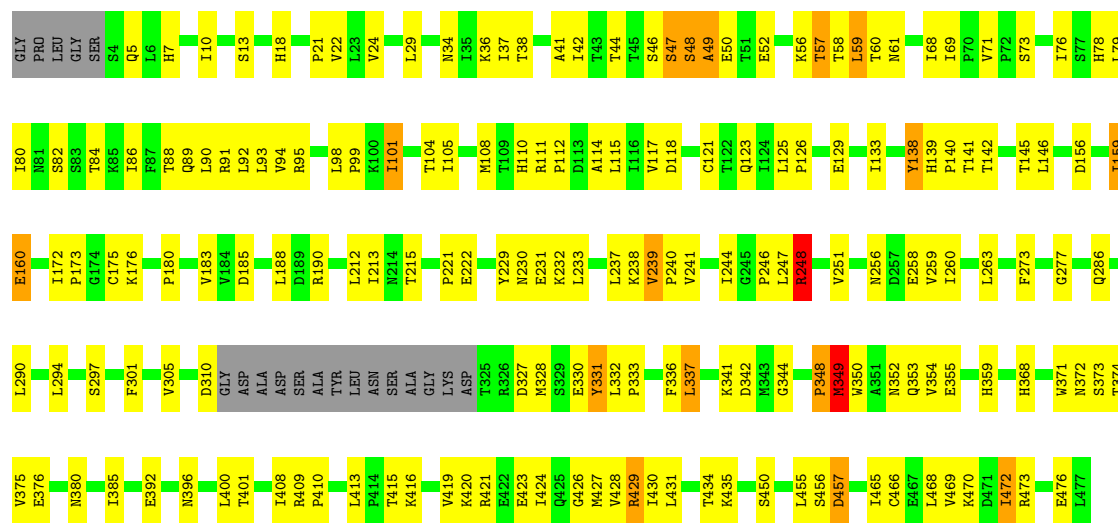
#### • Molecule 1: Glycosyltransferase

Chain B: 



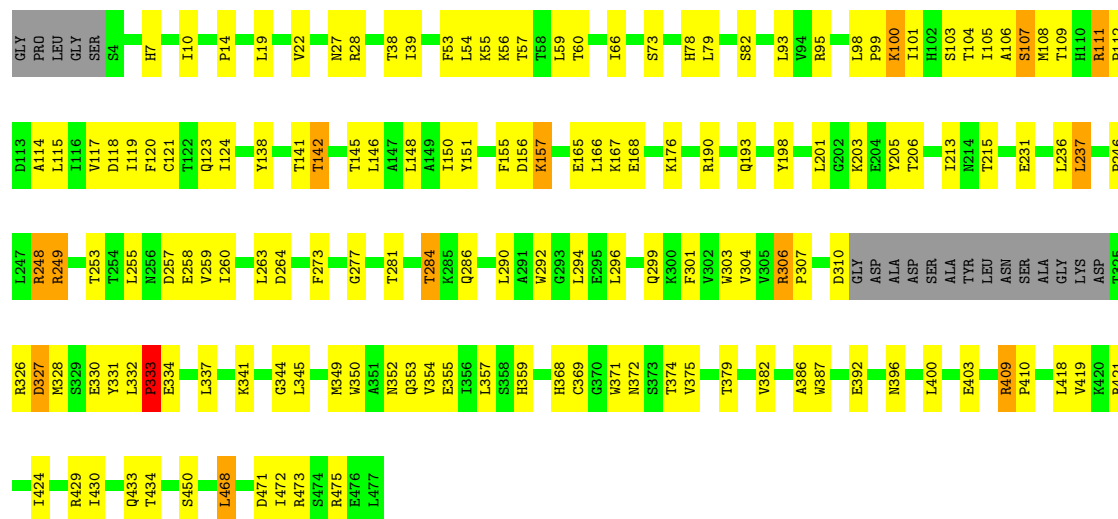
• Molecule 1: Glycosyltransferase

Chain C:  58% 34%



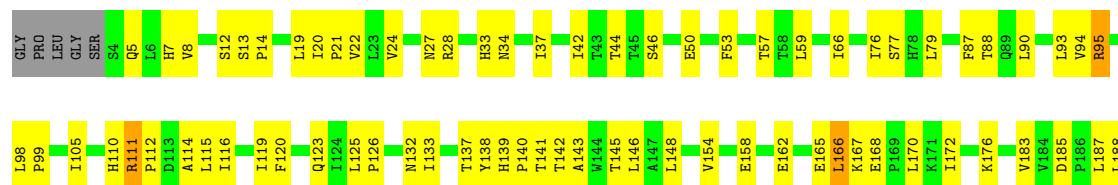
• Molecule 1: Glycosyltransferase

Chain D:  65% 28%



• Molecule 1: Glycosyltransferase

Chain E:  61% 31%









## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	87.72Å 171.78Å 153.38Å 90.00° 93.35° 90.00°	Depositor
Resolution (Å)	48.89 – 3.11 48.89 – 3.11	Depositor EDS
% Data completeness (in resolution range)	68.8 (48.89-3.11) 58.2 (48.89-3.11)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.55 (at 3.12Å)	Xtriage
Refinement program	REFMAC Refmac5	Depositor
R, $R_{free}$	0.268 , 0.312 0.264 , 0.298	Depositor DCC
$R_{free}$ test set	2878 reflections (5.13%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	58.5	Xtriage
Anisotropy	0.308	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 73.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.41$ , $\langle L^2 \rangle = 0.24$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.90	EDS
Total number of atoms	25741	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.94% 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: U2F, T83

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	0.44	0/3670	0.75	5/4978 (0.1%)
1	B	0.43	0/3704	0.85	16/5024 (0.3%)
1	C	0.46	1/3711 (0.0%)	0.85	10/5034 (0.2%)
1	D	0.43	0/3711	0.78	6/5034 (0.1%)
1	E	0.45	0/3704	0.75	6/5024 (0.1%)
1	F	0.44	0/3704	0.76	6/5024 (0.1%)
1	G	0.50	0/3711	0.96	24/5034 (0.5%)
All	All	0.45	1/25915 (0.0%)	0.82	73/35152 (0.2%)

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	A	0	9
1	B	0	8
1	C	0	6
1	D	0	8
1	E	0	10
1	F	0	7
1	G	0	9
All	All	0	57

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	231	GLU	C-O	-7.08	1.09	1.23

The worst 5 of 73 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	G	59	LEU	CB-CA-C	-19.66	72.85	110.20
1	C	57	THR	CB-CA-C	14.99	152.07	111.60
1	G	278	SER	CB-CA-C	14.54	137.74	110.10
1	C	348	PRO	N-CA-C	14.48	149.75	112.10
1	G	41	ALA	N-CA-CB	-13.55	91.13	110.10

There are no chirality outliers.

5 of 57 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	111	ARG	Sidechain
1	A	179	ARG	Sidechain
1	A	228	ARG	Sidechain
1	A	234	ARG	Sidechain
1	A	248	ARG	Sidechain

## 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	3600	0	3690	160	1
1	B	3633	0	3723	118	1
1	C	3640	0	3730	151	0
1	D	3640	0	3729	130	0
1	E	3633	0	3723	156	0
1	F	3633	0	3722	120	0
1	G	3640	0	3729	183	0
2	A	36	0	21	0	0
2	B	36	0	20	3	0
2	C	36	0	20	15	0
2	D	36	0	20	5	0
2	E	36	0	20	10	0
2	F	36	0	20	4	0
2	G	36	0	20	5	0
3	A	14	0	0	1	0
3	B	14	0	0	1	0
3	C	14	0	0	1	0
3	D	14	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	F	14	0	0	1	0
All	All	25741	0	26187	1025	1

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 20.

The worst 5 of 1025 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:G:501:U2F:O4'	2:G:501:U2F:C4'	1.67	1.30
2:E:501:U2F:O4'	2:E:501:U2F:C4'	1.65	1.28
2:B:501:U2F:O4'	2:B:501:U2F:C4'	1.66	1.28
2:F:501:U2F:C4'	2:F:501:U2F:O4'	1.66	1.25
2:C:501:U2F:C4'	2:C:501:U2F:O4'	1.65	1.23

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:433:GLN:NE2	1:B:264:ASP:OD2[2_646]	2.07	0.13

## 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	449/479 (94%)	418 (93%)	30 (7%)	1 (0%)	44 73
1	B	455/479 (95%)	424 (93%)	31 (7%)	0	100 100
1	C	456/479 (95%)	420 (92%)	36 (8%)	0	100 100
1	D	456/479 (95%)	422 (92%)	33 (7%)	1 (0%)	44 73
1	E	455/479 (95%)	423 (93%)	32 (7%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	F	455/479 (95%)	422 (93%)	33 (7%)	0	100	100
1	G	456/479 (95%)	405 (89%)	50 (11%)	1 (0%)	44	73
All	All	3182/3353 (95%)	2934 (92%)	245 (8%)	3 (0%)	48	78

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	46	SER
1	D	73	SER
1	G	77	SER

### 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	408/425 (96%)	379 (93%)	29 (7%)	12	37
1	B	412/425 (97%)	390 (95%)	22 (5%)	19	46
1	C	413/425 (97%)	387 (94%)	26 (6%)	15	41
1	D	413/425 (97%)	388 (94%)	25 (6%)	15	42
1	E	412/425 (97%)	389 (94%)	23 (6%)	17	44
1	F	412/425 (97%)	376 (91%)	36 (9%)	8	29
1	G	413/425 (97%)	364 (88%)	49 (12%)	4	16
All	All	2883/2975 (97%)	2673 (93%)	210 (7%)	11	35

5 of 210 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	E	296	LEU
1	F	237	LEU
1	G	347	VAL
1	E	331	TYR
1	F	100	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 51 such sidechains are listed below:

Mol	Chain	Res	Type
1	E	286	GLN
1	F	425	GLN
1	G	433	GLN
1	E	299	GLN
1	F	5	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 oligosaccharides in this entry.

## 5.6 Ligand geometry ⓘ

12 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	T83	A	502	-	15,15,15	0.59	0	21,21,21	1.02	2 (9%)
3	T83	C	502	-	15,15,15	0.49	0	21,21,21	0.96	2 (9%)
3	T83	B	502	-	15,15,15	0.47	0	21,21,21	0.95	2 (9%)
3	T83	D	502	-	15,15,15	0.52	0	21,21,21	0.93	2 (9%)
2	U2F	F	501	-	35,38,38	3.55	18 (51%)	51,58,58	1.94	13 (25%)
2	U2F	A	501	-	35,38,38	0.54	0	51,58,58	0.61	0



Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
2	U2F	E	501	-	35,38,38	3.36	19 (54%)	51,58,58	2.06	18 (35%)
2	U2F	C	501	-	35,38,38	3.60	18 (51%)	51,58,58	1.96	10 (19%)
3	T83	F	502	-	15,15,15	0.48	0	21,21,21	0.93	2 (9%)
2	U2F	B	501	-	35,38,38	3.39	19 (54%)	51,58,58	1.77	15 (29%)
2	U2F	G	501	-	35,38,38	3.45	20 (57%)	51,58,58	1.89	13 (25%)
2	U2F	D	501	-	35,38,38	3.44	19 (54%)	51,58,58	1.82	13 (25%)

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	T83	A	502	-	-	2/2/2/2	0/2/2/2
3	T83	C	502	-	-	2/2/2/2	0/2/2/2
3	T83	B	502	-	-	2/2/2/2	0/2/2/2
3	T83	D	502	-	-	2/2/2/2	0/2/2/2
2	U2F	F	501	-	-	7/22/59/59	0/3/3/3
2	U2F	A	501	-	-	8/22/59/59	0/3/3/3
2	U2F	E	501	-	-	14/22/59/59	0/3/3/3
2	U2F	C	501	-	-	8/22/59/59	0/3/3/3
3	T83	F	502	-	-	2/2/2/2	0/2/2/2
2	U2F	B	501	-	-	8/22/59/59	0/3/3/3
2	U2F	G	501	-	-	9/22/59/59	0/3/3/3
2	U2F	D	501	-	-	8/22/59/59	0/3/3/3

The worst 5 of 113 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	G	501	U2F	O4'-C4'	9.83	1.67	1.45
2	F	501	U2F	O4'-C4'	9.80	1.66	1.45
2	D	501	U2F	O4'-C4'	9.65	1.66	1.45
2	B	501	U2F	O4'-C4'	9.42	1.66	1.45
2	E	501	U2F	O4'-C4'	9.33	1.65	1.45

The worst 5 of 92 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	501	U2F	N3-C6'-N1	7.10	124.32	114.89
2	F	501	U2F	N3-C6'-N1	7.03	124.22	114.89
2	G	501	U2F	N3-C6'-N1	6.84	123.98	114.89
2	B	501	U2F	N3-C6'-N1	6.12	123.01	114.89
2	C	501	U2F	C7'-N3-C6'	-6.03	118.62	126.58

There are no chirality outliers.

5 of 72 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	501	U2F	O5-C1-O1-PB
2	B	501	U2F	C1-O1-PB-O3A
2	C	501	U2F	O5-C1-O1-PB
2	C	501	U2F	C4'-C5'-O5'-PA
2	D	501	U2F	O5-C1-O1-PB

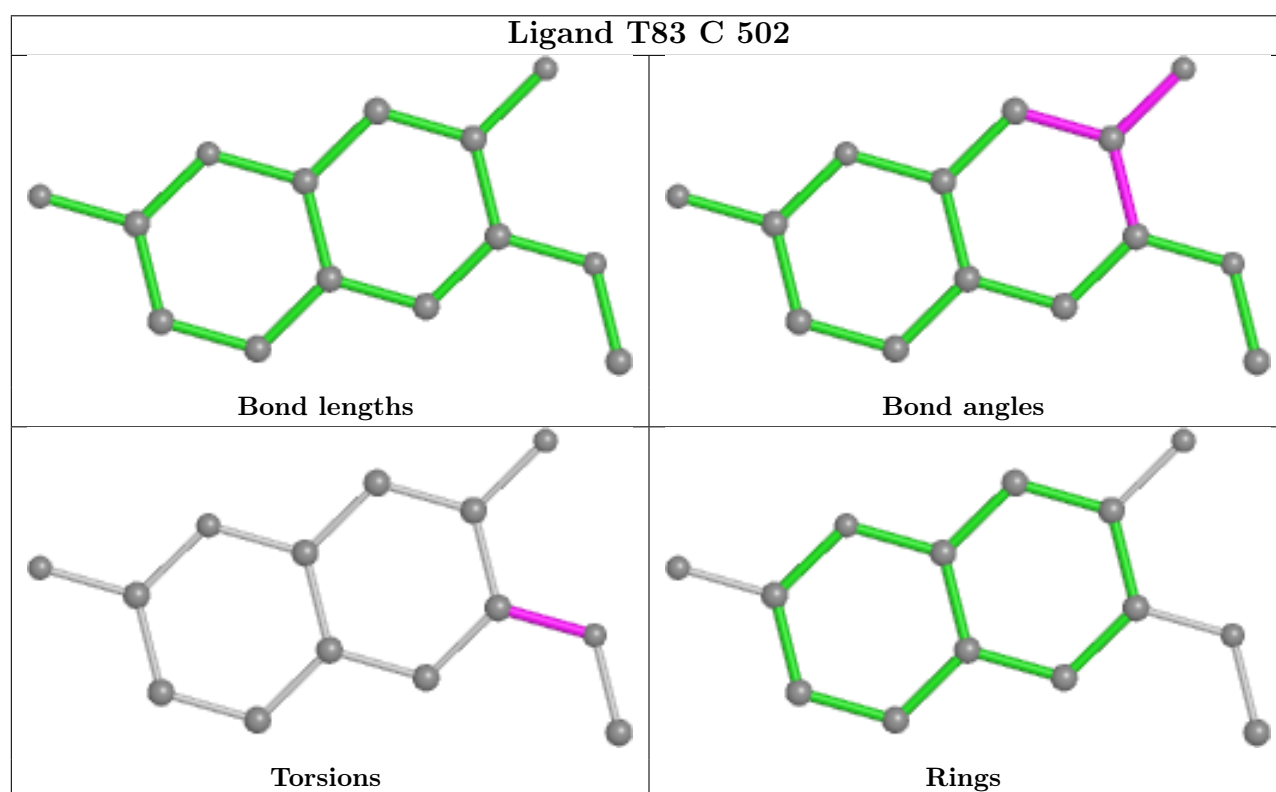
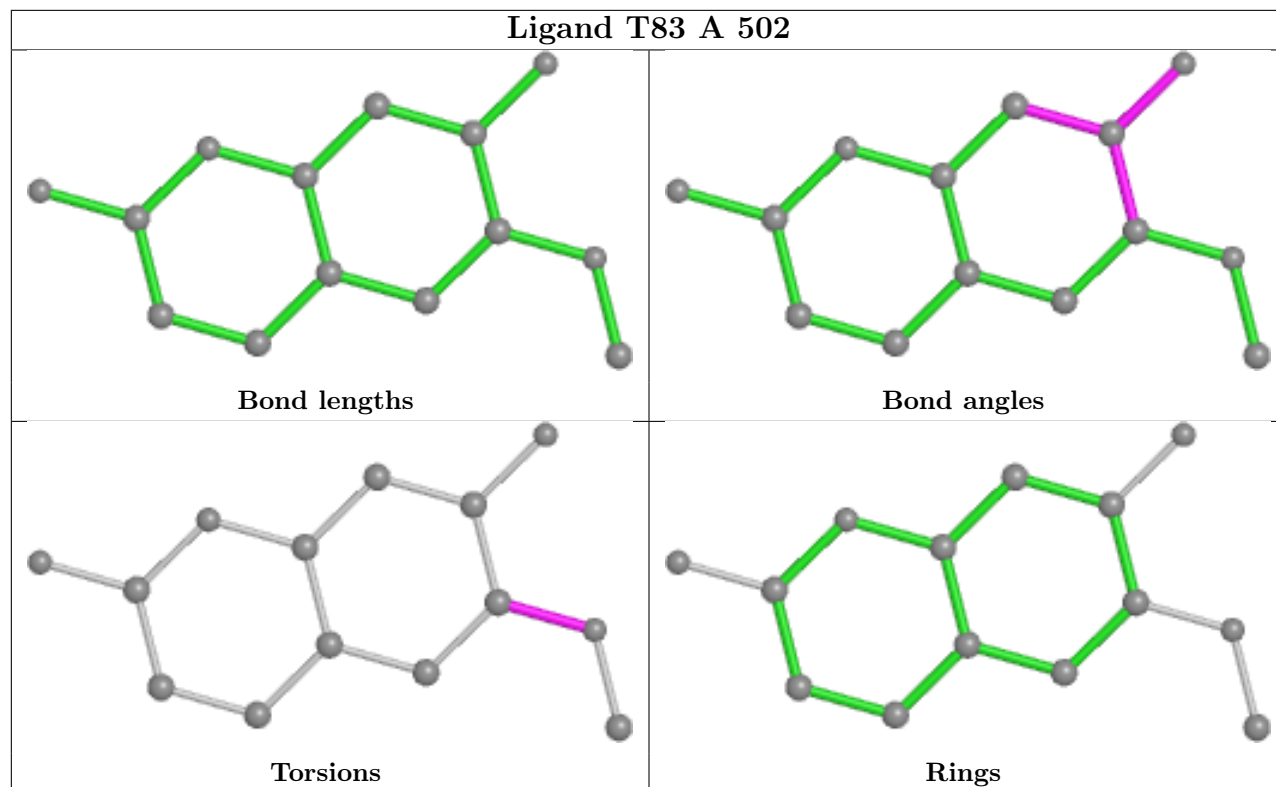
There are no ring outliers.

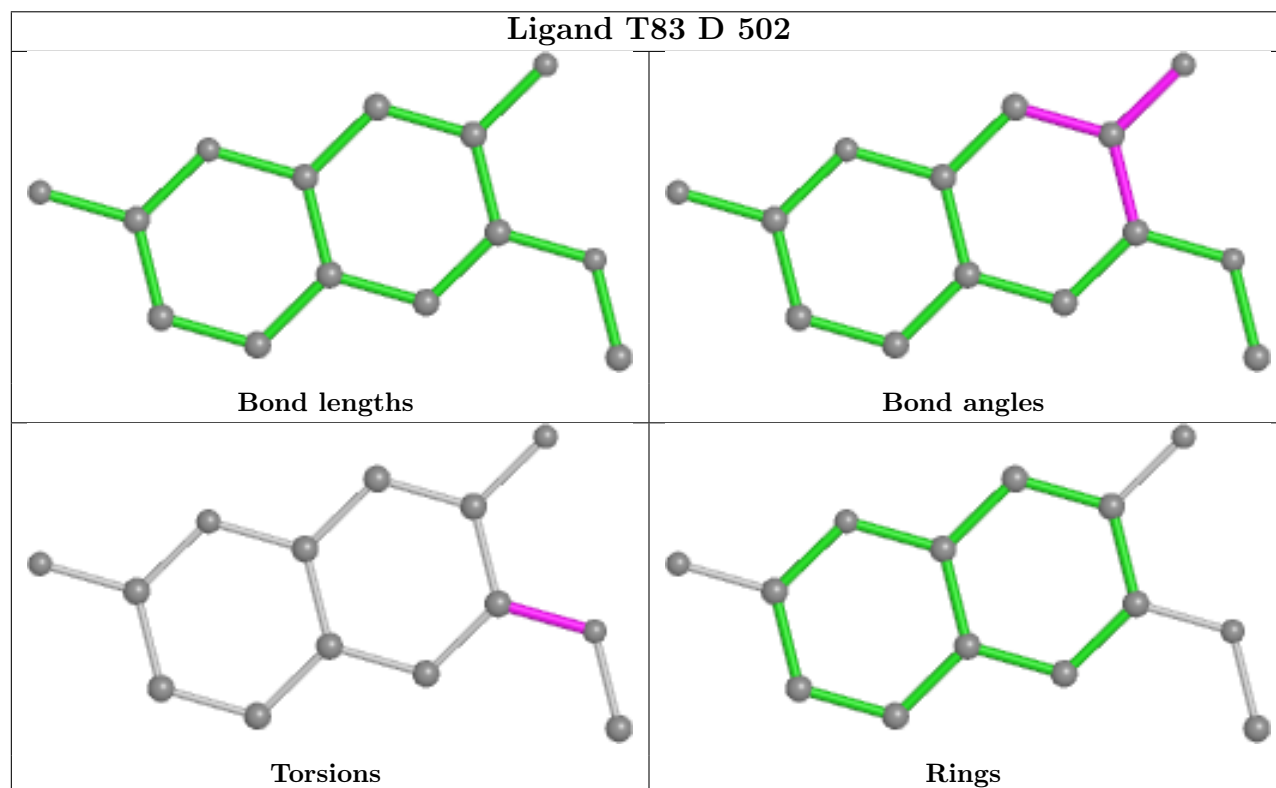
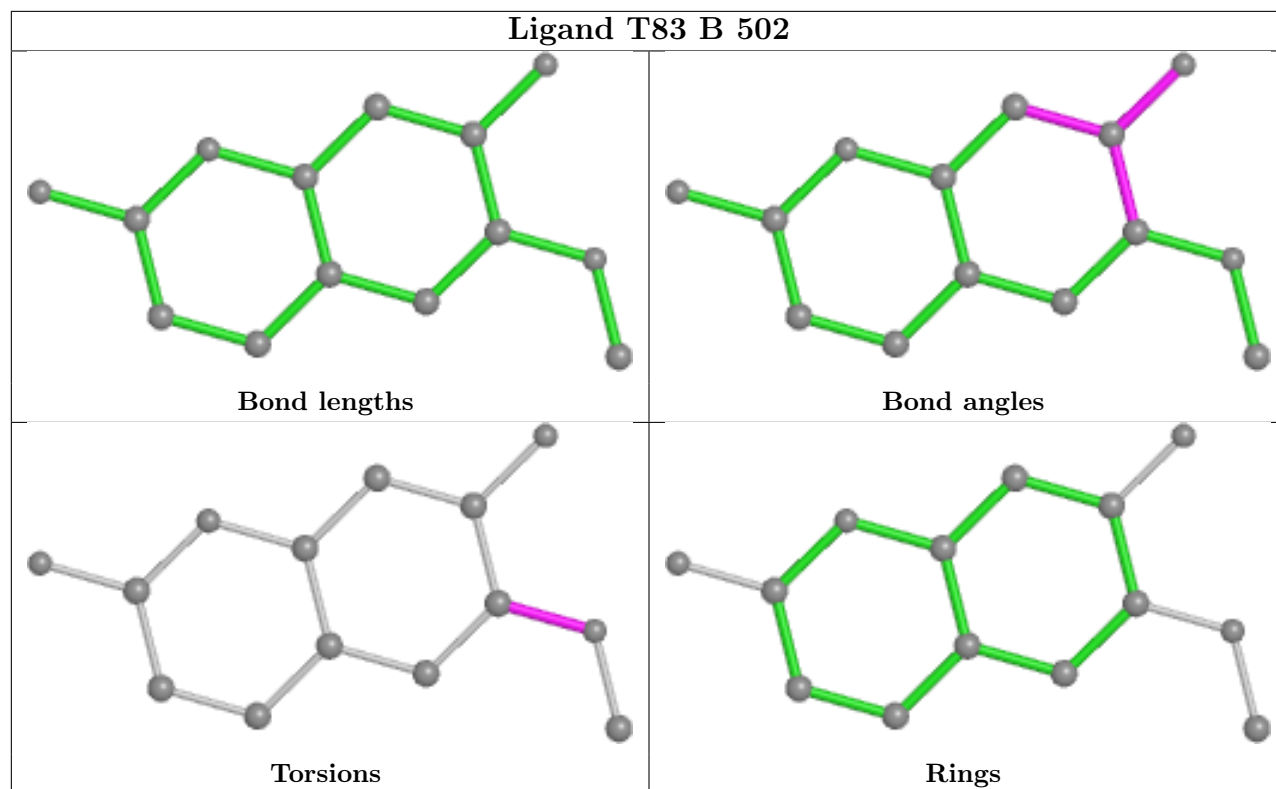
11 monomers are involved in 47 short contacts:

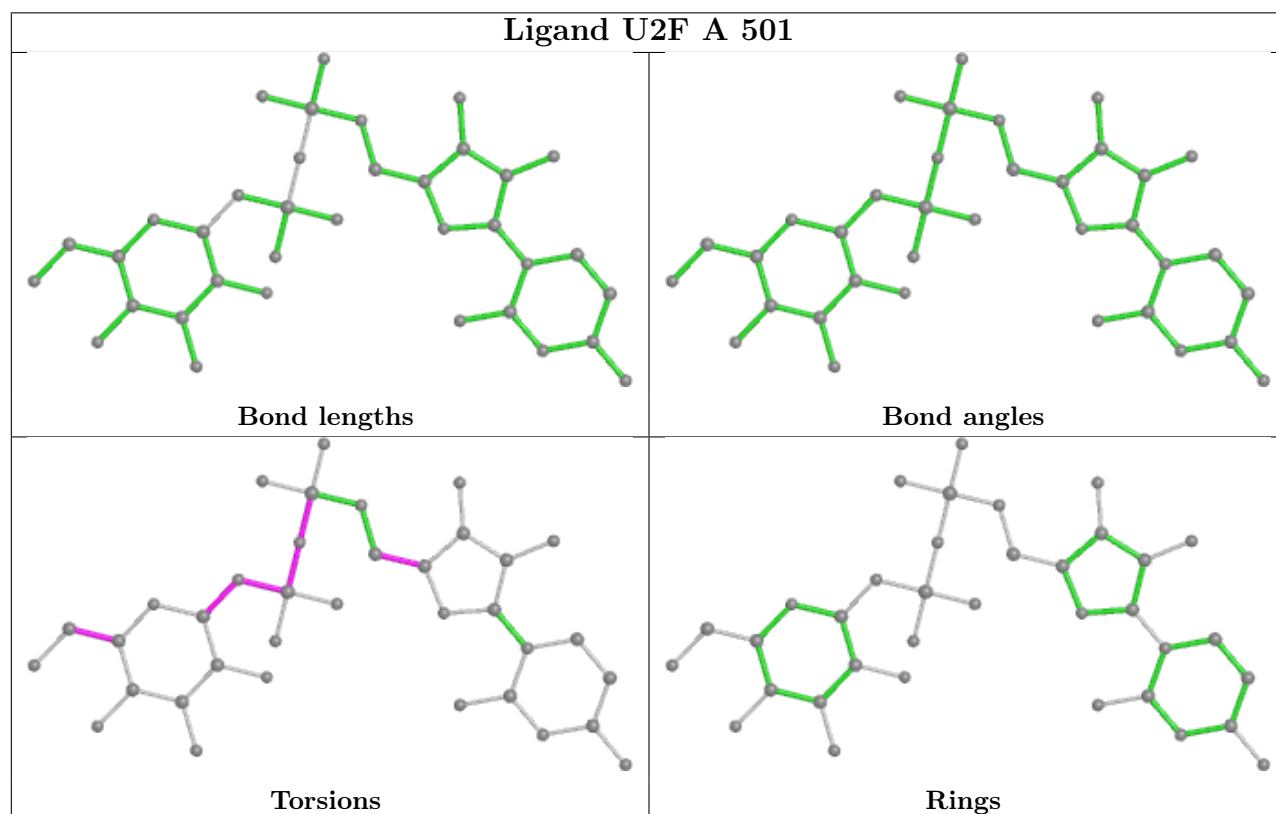
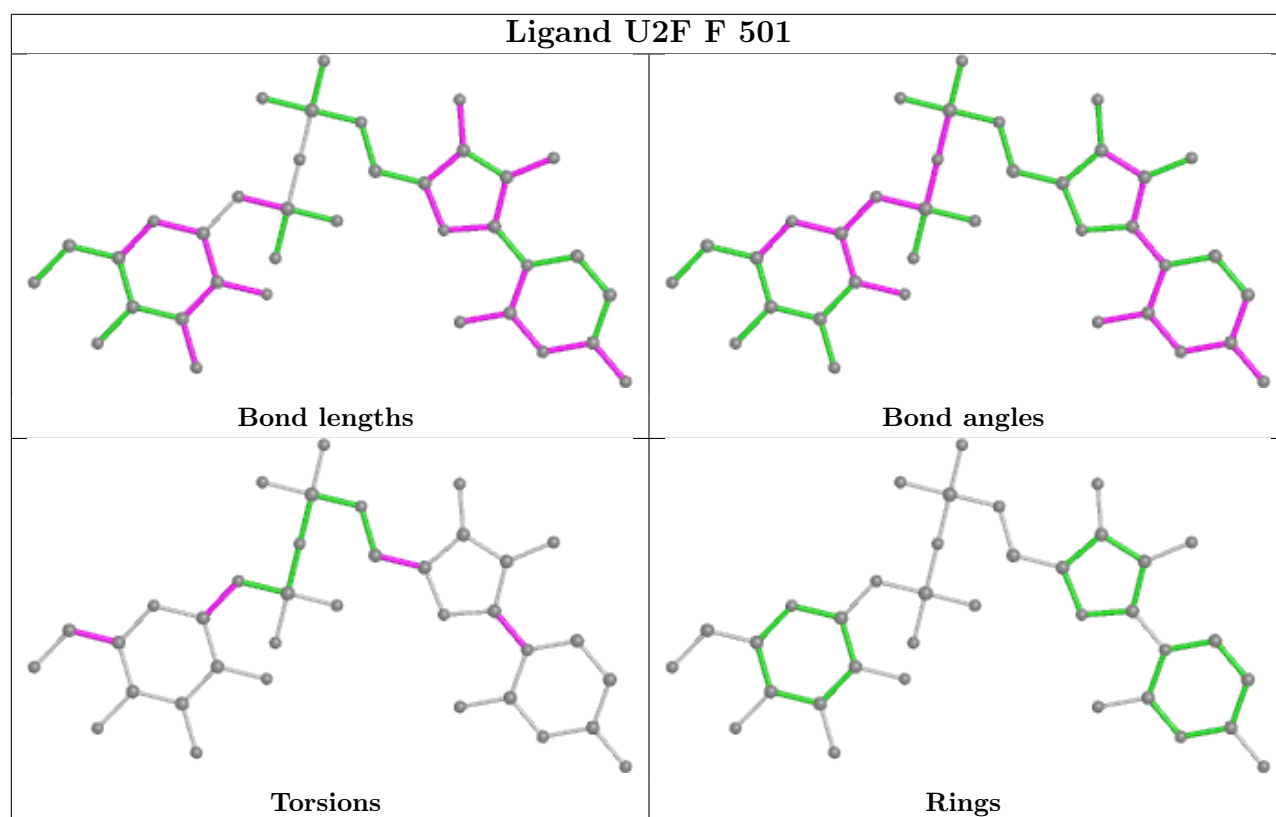
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	502	T83	1	0
3	C	502	T83	1	0
3	B	502	T83	1	0
3	D	502	T83	1	0
2	F	501	U2F	4	0
2	E	501	U2F	10	0
2	C	501	U2F	15	0
3	F	502	T83	1	0
2	B	501	U2F	3	0
2	G	501	U2F	5	0
2	D	501	U2F	5	0

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

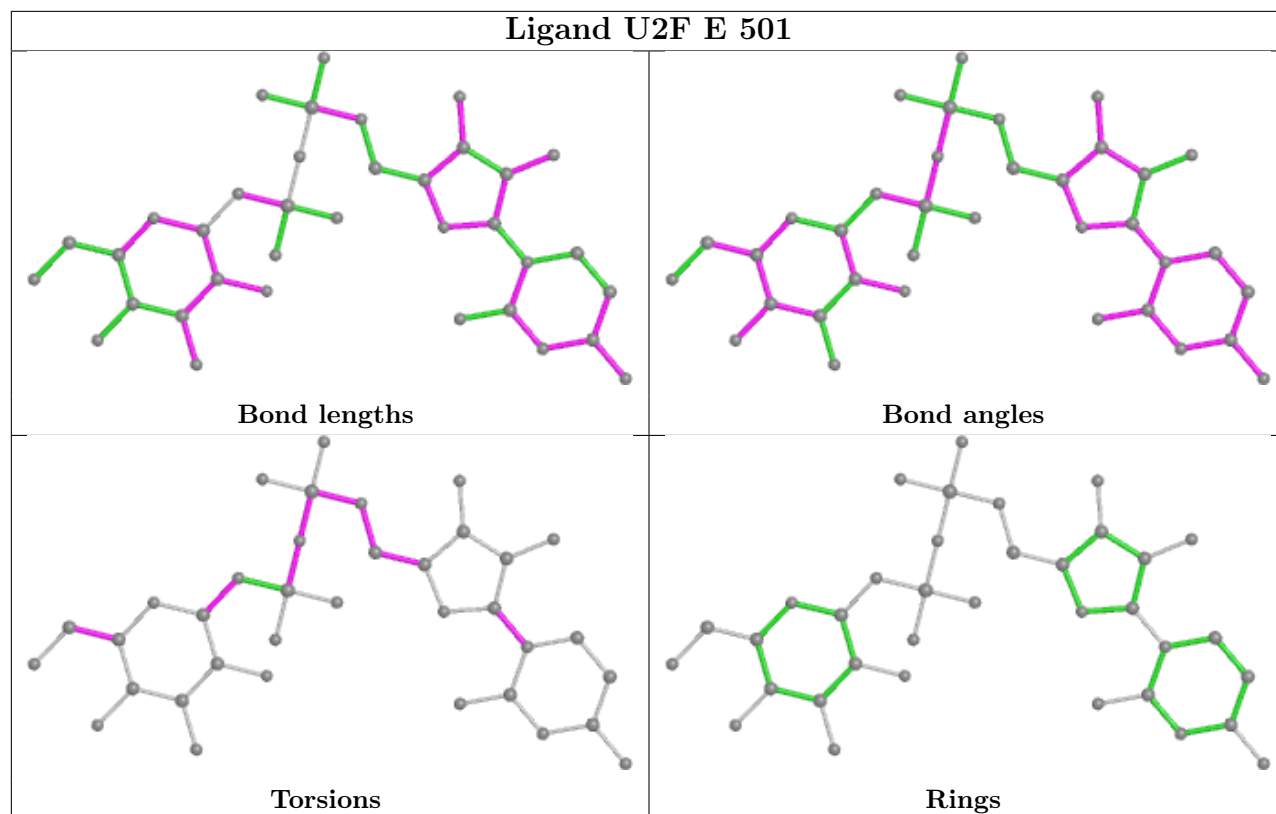
equivalents in the CSD to analyse the geometry.



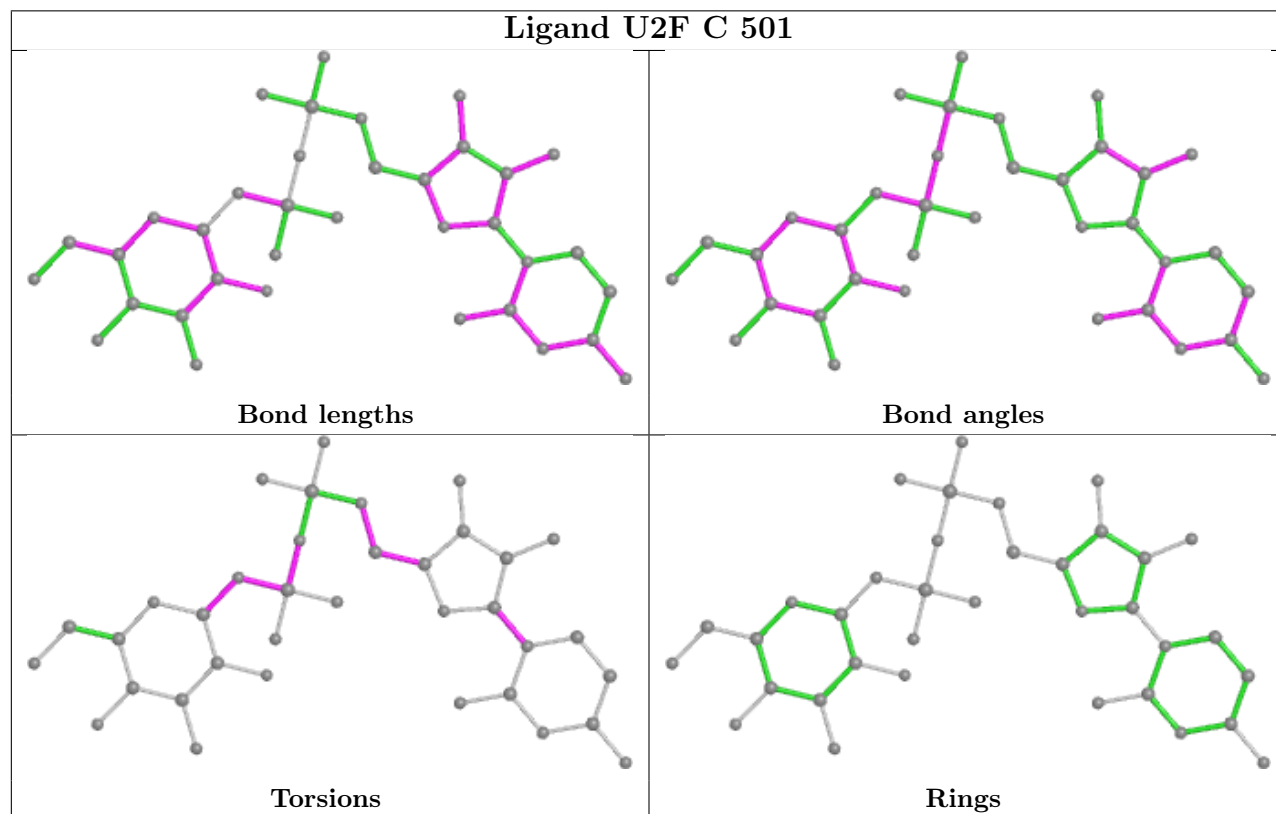


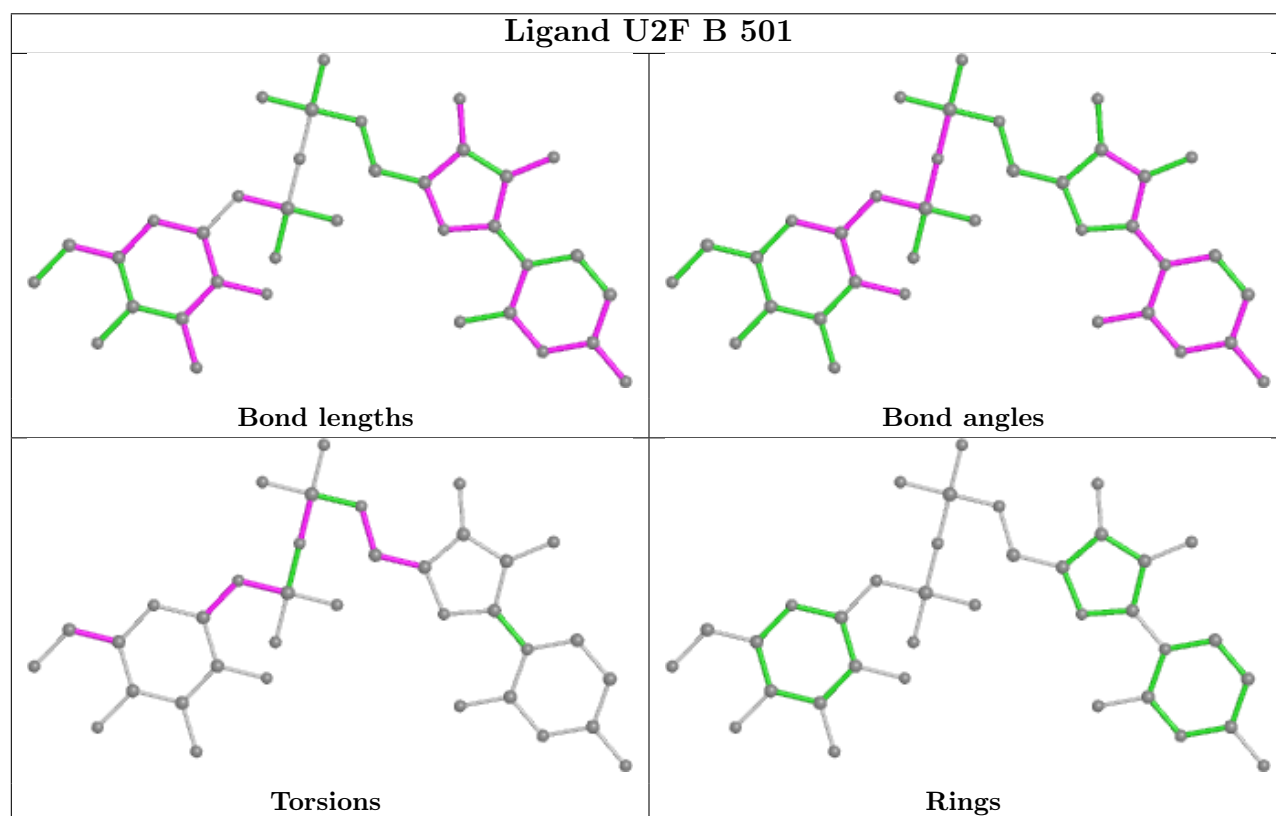
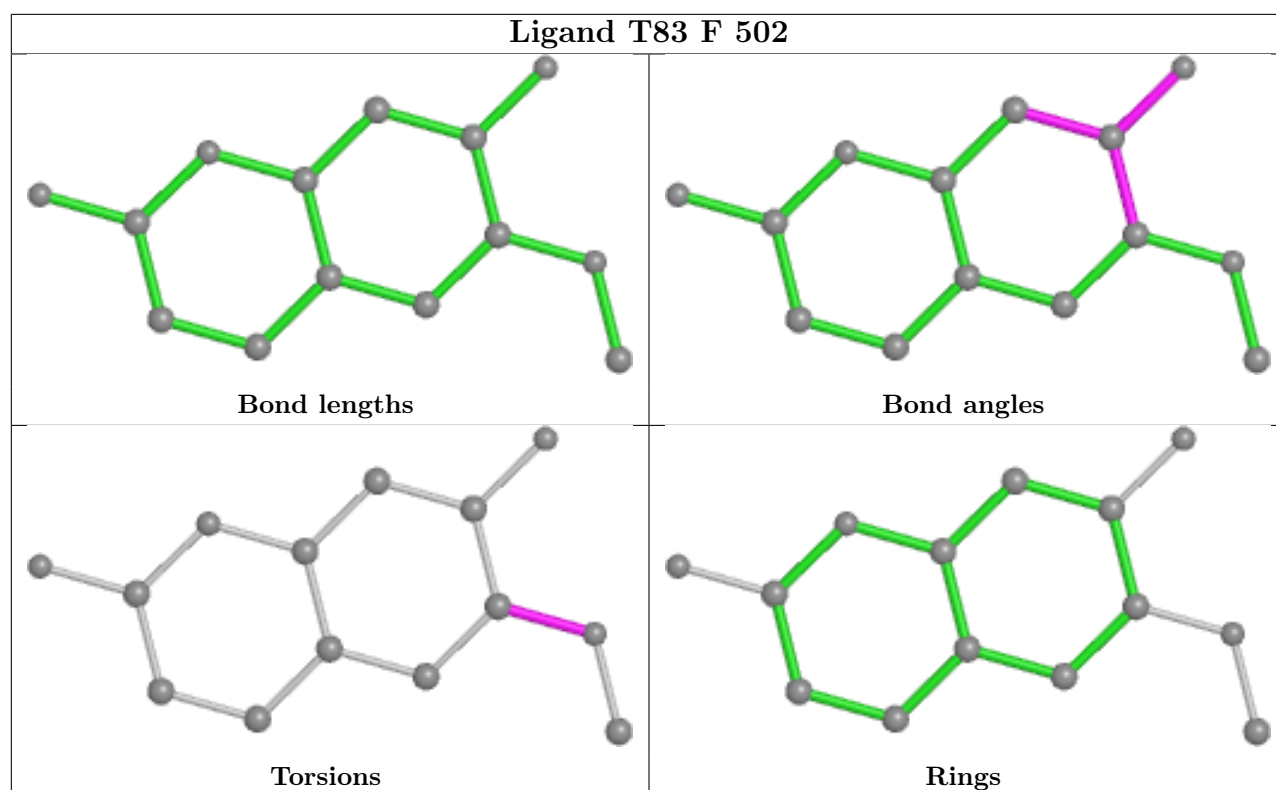


## Ligand U2F E 501

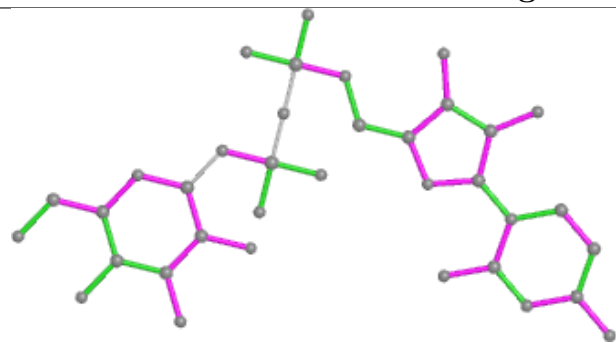


## Ligand U2F C 501

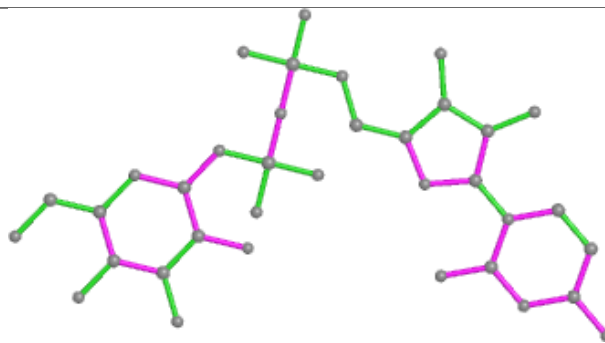




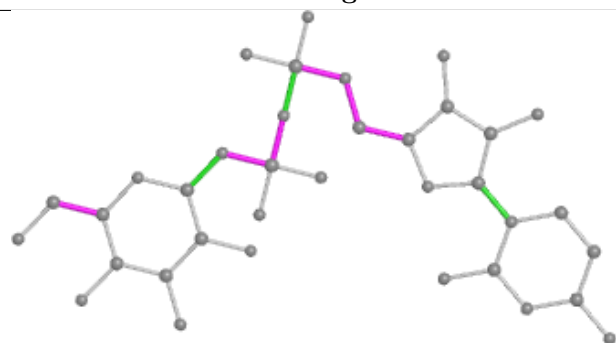
## Ligand U2F G 501



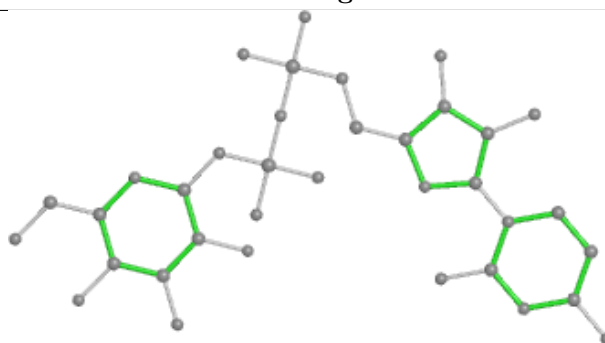
Bond lengths



Bond angles

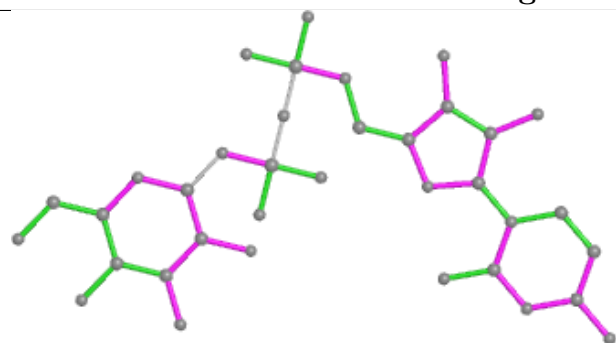


Torsions

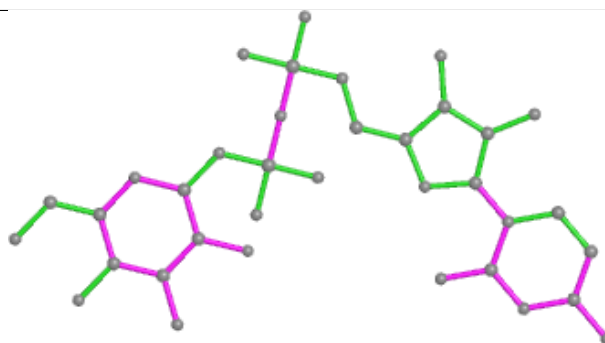


Rings

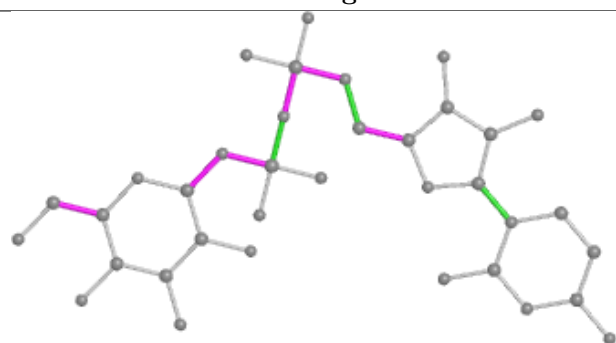
## Ligand U2F D 501



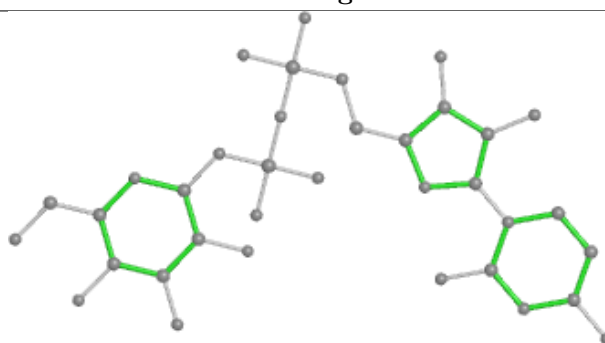
Bond lengths



Bond angles



Torsions



Rings



## 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, 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	455/479 (94%)	-1.63	0 100 100	11, 30, 65, 138	0
1	B	459/479 (95%)	-1.65	0 100 100	12, 28, 57, 121	0
1	C	460/479 (96%)	-1.65	0 100 100	10, 28, 61, 92	0
1	D	460/479 (96%)	-1.65	0 100 100	13, 29, 56, 97	0
1	E	459/479 (95%)	-1.63	0 100 100	10, 31, 59, 81	0
1	F	459/479 (95%)	-1.61	0 100 100	14, 34, 62, 126	0
1	G	460/479 (96%)	-1.59	0 100 100	13, 35, 67, 94	0
All	All	3212/3353 (95%)	-1.63	0 100 100	10, 30, 61, 138	0

There are no RSRZ outliers to report.

### 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 monosaccharides in this entry.

### 6.4 Ligands [i](#)

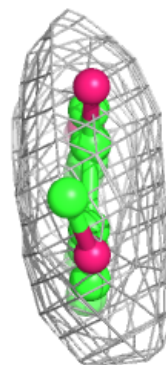
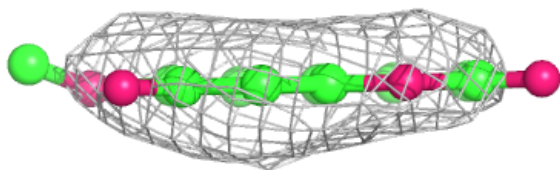
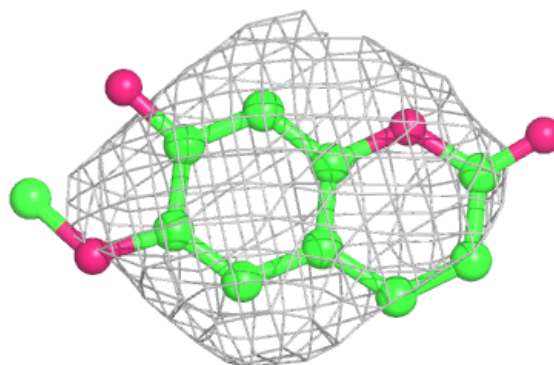
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
3	T83	F	502	14/14	0.98	0.06	62,68,75,76	0
2	U2F	D	501	36/36	0.99	0.03	34,49,69,84	0
2	U2F	F	501	36/36	0.99	0.03	33,49,79,80	0
2	U2F	G	501	36/36	0.99	0.04	36,63,119,125	0
3	T83	A	502	14/14	0.99	0.04	52,62,85,95	0
3	T83	B	502	14/14	0.99	0.04	49,54,62,66	0
3	T83	C	502	14/14	0.99	0.04	56,67,80,83	0
3	T83	D	502	14/14	0.99	0.04	35,56,75,87	0
2	U2F	C	501	36/36	0.99	0.03	30,69,117,141	0
2	U2F	B	501	36/36	1.00	0.03	22,27,60,62	0
2	U2F	E	501	36/36	1.00	0.03	39,65,105,117	0
2	U2F	A	501	36/36	1.00	0.02	29,38,78,92	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.

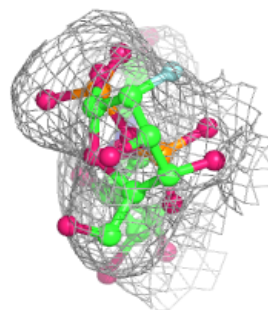
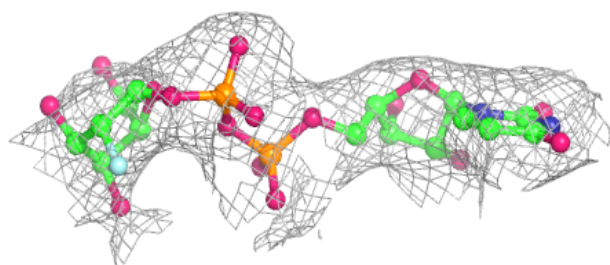
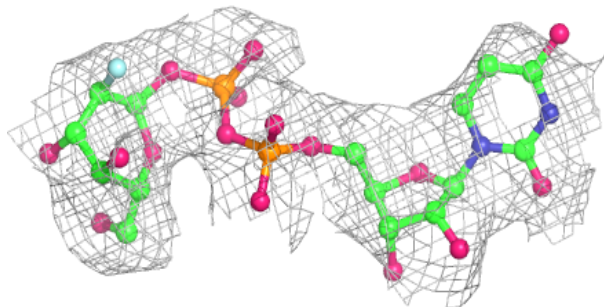
**Electron density around T83 F 502:**

2mF<sub>o</sub>-DF<sub>c</sub> (at 0.7 rmsd) in gray  
 mF<sub>o</sub>-DF<sub>c</sub> (at 3 rmsd) in purple (negative)  
 and green (positive)

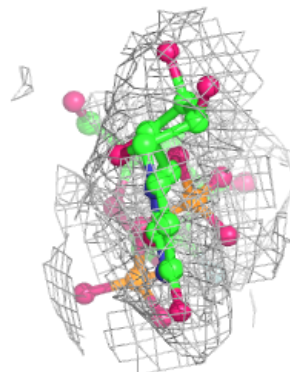
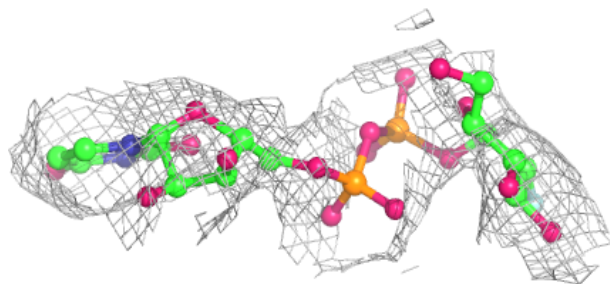
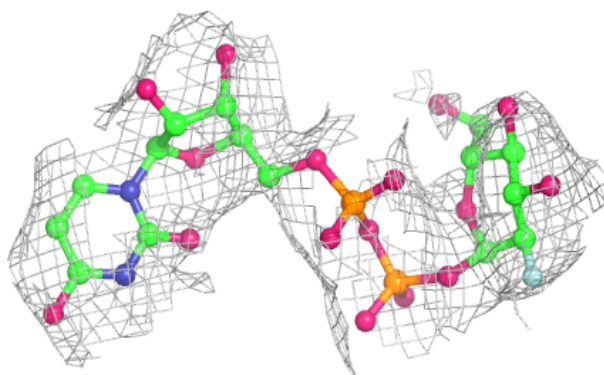


**Electron density around U2F D 501:**

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

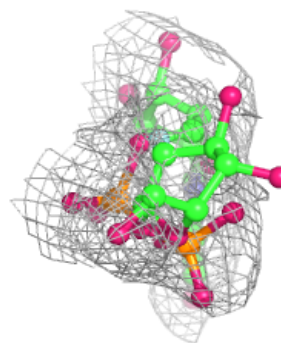
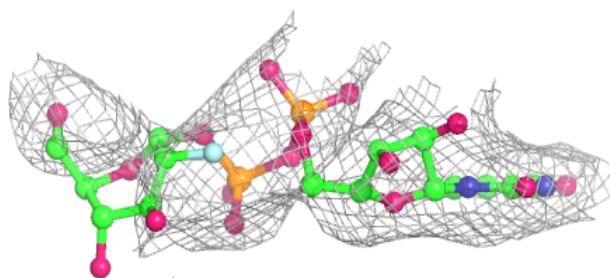
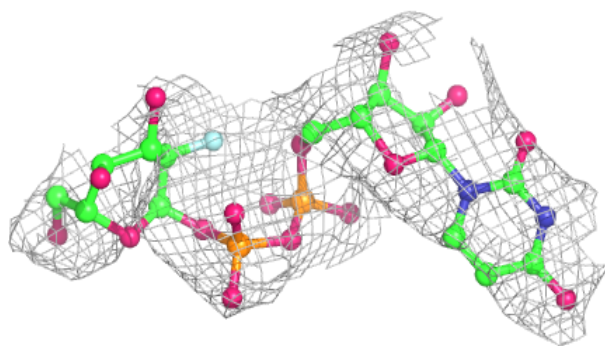
**Electron density around U2F F 501:**

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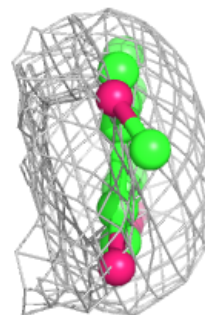
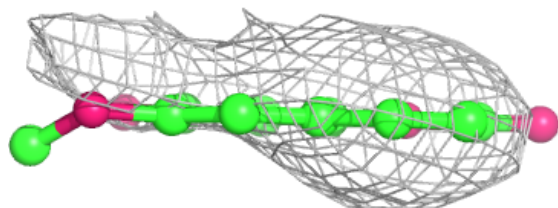
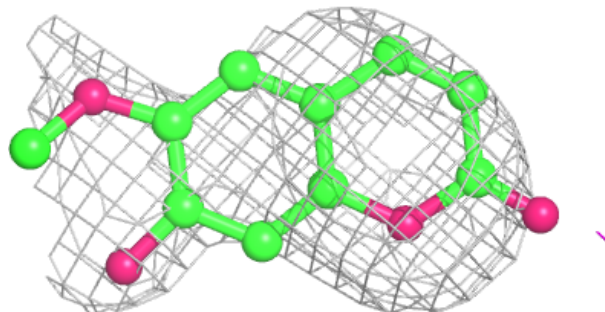


**Electron density around U2F G 501:**

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

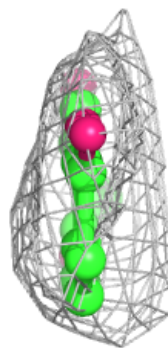
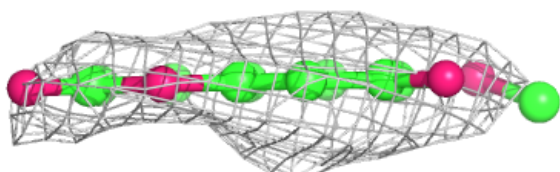
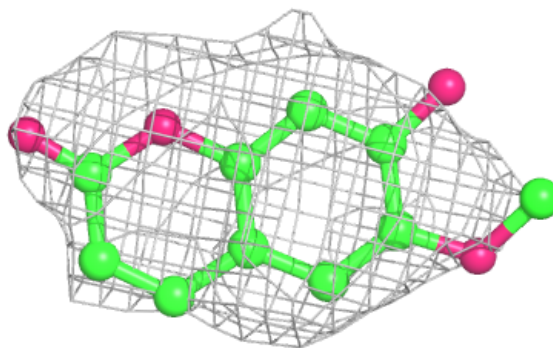
**Electron density around T83 A 502:**

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

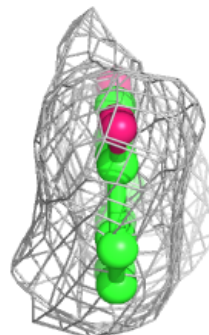
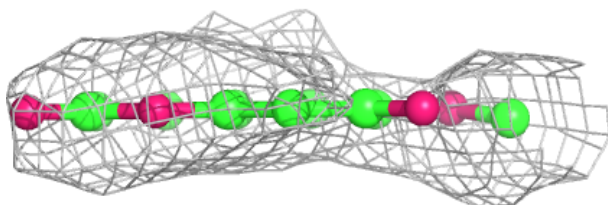
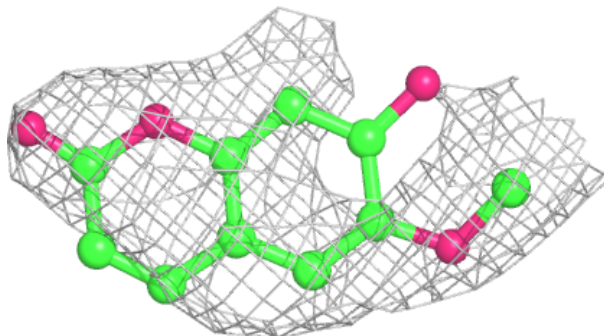


**Electron density around T83 B 502:**

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**Electron density around T83 C 502:**

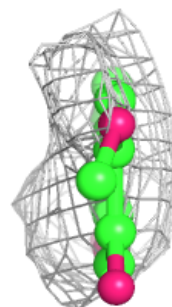
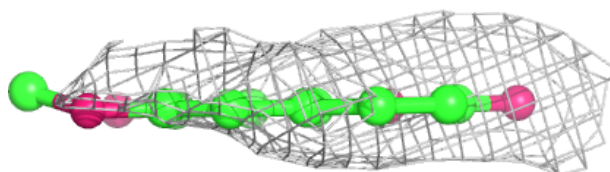
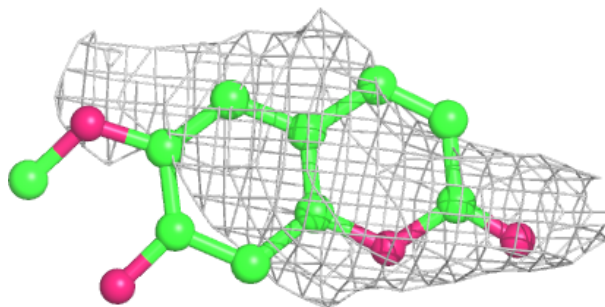
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 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



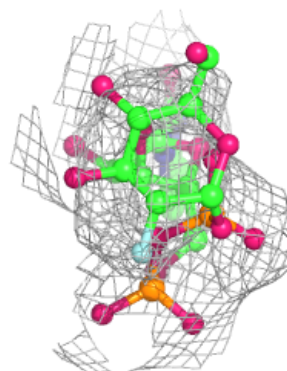
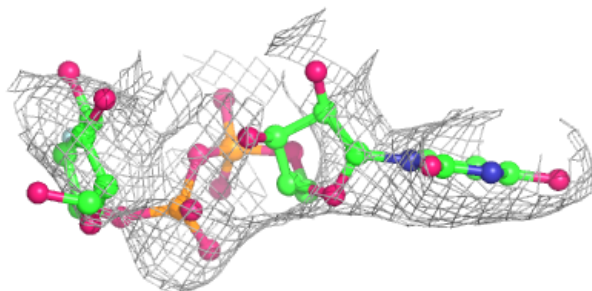
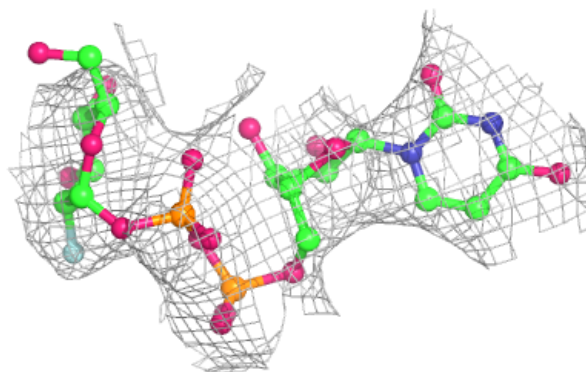


**Electron density around T83 D 502:**

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and green (positive)

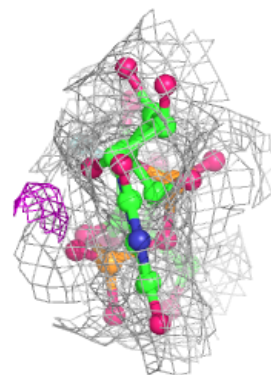
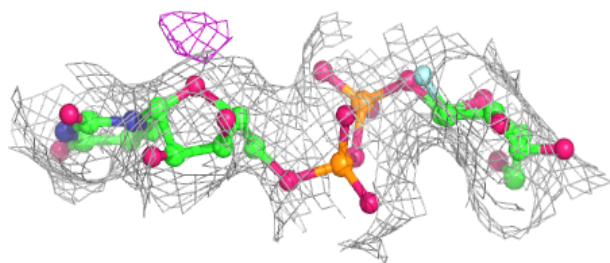
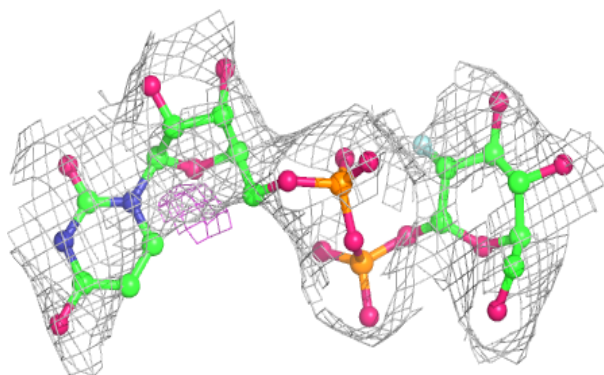
**Electron density around U2F C 501:**

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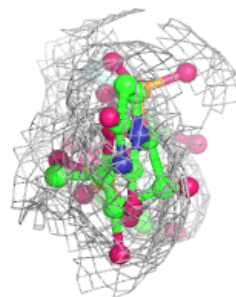
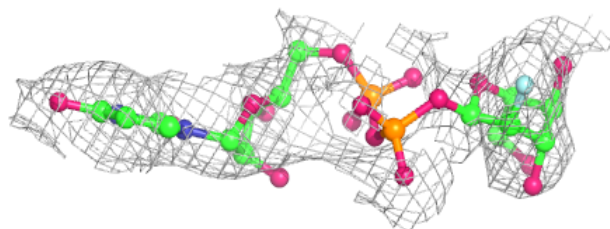
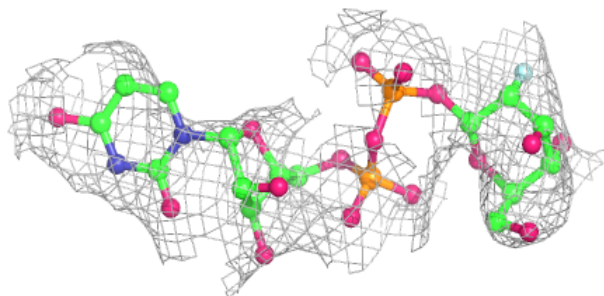


**Electron density around U2F B 501:**

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

**Electron density around U2F E 501:**

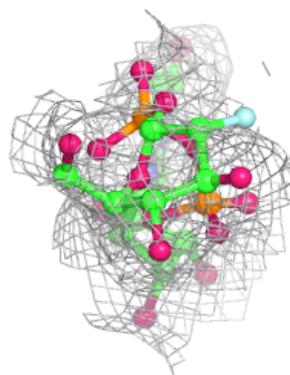
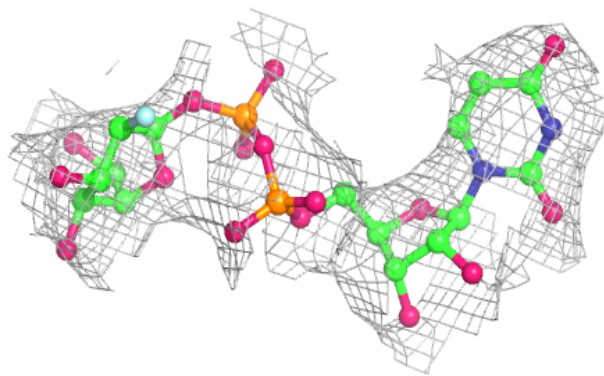
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





**Electron density around U2F A 501:**

$2mF_o-DF_c$  (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.