



Full wwPDB X-ray Structure Validation Report ⓘ

May 14, 2020 – 08:08 pm BST

PDB ID : 3GWC
Title : Crystal structure of Mycobacterium tuberculosis thymidylate synthase X bound to FdUMP and FAD
Authors : Seattle Structural Genomics Center for Infectious Disease (SSGCID)
Deposited on : 2009-03-31
Resolution : 1.90 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.11
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

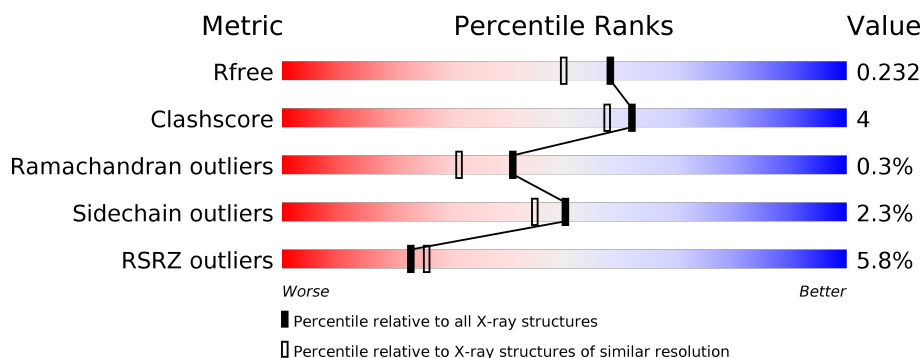
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.





Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	6207 (1.90-1.90)
Clashscore	141614	6847 (1.90-1.90)
Ramachandran outliers	138981	6760 (1.90-1.90)
Sidechain outliers	138945	6760 (1.90-1.90)
RSRZ outliers	127900	6082 (1.90-1.90)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	258	
1	B	258	
1	C	258	
1	D	258	
1	E	258	
1	F	258	

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Mol	Chain	Length	Quality of chain
1	G	258	
1	H	258	

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
3	UFP	D	260	X	-	-	-
3	UFP	F	260	X	-	-	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 16612 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 thyX.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	246	Total	C	N	O	S	0	0	0
			1912	1203	349	355	5			
1	B	239	Total	C	N	O	S	0	0	0
			1855	1166	339	345	5			
1	C	243	Total	C	N	O	S	0	0	0
			1882	1183	342	352	5			
1	D	242	Total	C	N	O	S	0	0	0
			1877	1180	341	351	5			
1	E	244	Total	C	N	O	S	0	0	0
			1887	1186	343	353	5			
1	F	242	Total	C	N	O	S	0	0	0
			1884	1183	345	351	5			
1	G	245	Total	C	N	O	S	0	0	0
			1897	1194	344	354	5			
1	H	243	Total	C	N	O	S	0	0	0
			1883	1185	341	352	5			

There are 64 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	251	LEU	-	EXPRESSION TAG	UNP P66930
A	252	GLU	-	EXPRESSION TAG	UNP P66930
A	253	HIS	-	EXPRESSION TAG	UNP P66930
A	254	HIS	-	EXPRESSION TAG	UNP P66930
A	255	HIS	-	EXPRESSION TAG	UNP P66930
A	256	HIS	-	EXPRESSION TAG	UNP P66930
A	257	HIS	-	EXPRESSION TAG	UNP P66930
A	258	HIS	-	EXPRESSION TAG	UNP P66930
B	251	LEU	-	EXPRESSION TAG	UNP P66930
B	252	GLU	-	EXPRESSION TAG	UNP P66930
B	253	HIS	-	EXPRESSION TAG	UNP P66930
B	254	HIS	-	EXPRESSION TAG	UNP P66930
B	255	HIS	-	EXPRESSION TAG	UNP P66930

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Chain	Residue	Modelled	Actual	Comment	Reference
B	256	HIS	-	EXPRESSION TAG	UNP P66930
B	257	HIS	-	EXPRESSION TAG	UNP P66930
B	258	HIS	-	EXPRESSION TAG	UNP P66930
C	251	LEU	-	EXPRESSION TAG	UNP P66930
C	252	GLU	-	EXPRESSION TAG	UNP P66930
C	253	HIS	-	EXPRESSION TAG	UNP P66930
C	254	HIS	-	EXPRESSION TAG	UNP P66930
C	255	HIS	-	EXPRESSION TAG	UNP P66930
C	256	HIS	-	EXPRESSION TAG	UNP P66930
C	257	HIS	-	EXPRESSION TAG	UNP P66930
C	258	HIS	-	EXPRESSION TAG	UNP P66930
D	251	LEU	-	EXPRESSION TAG	UNP P66930
D	252	GLU	-	EXPRESSION TAG	UNP P66930
D	253	HIS	-	EXPRESSION TAG	UNP P66930
D	254	HIS	-	EXPRESSION TAG	UNP P66930
D	255	HIS	-	EXPRESSION TAG	UNP P66930
D	256	HIS	-	EXPRESSION TAG	UNP P66930
D	257	HIS	-	EXPRESSION TAG	UNP P66930
D	258	HIS	-	EXPRESSION TAG	UNP P66930
E	251	LEU	-	EXPRESSION TAG	UNP P66930
E	252	GLU	-	EXPRESSION TAG	UNP P66930
E	253	HIS	-	EXPRESSION TAG	UNP P66930
E	254	HIS	-	EXPRESSION TAG	UNP P66930
E	255	HIS	-	EXPRESSION TAG	UNP P66930
E	256	HIS	-	EXPRESSION TAG	UNP P66930
E	257	HIS	-	EXPRESSION TAG	UNP P66930
E	258	HIS	-	EXPRESSION TAG	UNP P66930
F	251	LEU	-	EXPRESSION TAG	UNP P66930
F	252	GLU	-	EXPRESSION TAG	UNP P66930
F	253	HIS	-	EXPRESSION TAG	UNP P66930
F	254	HIS	-	EXPRESSION TAG	UNP P66930
F	255	HIS	-	EXPRESSION TAG	UNP P66930
F	256	HIS	-	EXPRESSION TAG	UNP P66930
F	257	HIS	-	EXPRESSION TAG	UNP P66930
F	258	HIS	-	EXPRESSION TAG	UNP P66930
G	251	LEU	-	EXPRESSION TAG	UNP P66930
G	252	GLU	-	EXPRESSION TAG	UNP P66930
G	253	HIS	-	EXPRESSION TAG	UNP P66930
G	254	HIS	-	EXPRESSION TAG	UNP P66930
G	255	HIS	-	EXPRESSION TAG	UNP P66930
G	256	HIS	-	EXPRESSION TAG	UNP P66930
G	257	HIS	-	EXPRESSION TAG	UNP P66930

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Chain	Residue	Modelled	Actual	Comment	Reference
G	258	HIS	-	EXPRESSION TAG	UNP P66930
H	251	LEU	-	EXPRESSION TAG	UNP P66930
H	252	GLU	-	EXPRESSION TAG	UNP P66930
H	253	HIS	-	EXPRESSION TAG	UNP P66930
H	254	HIS	-	EXPRESSION TAG	UNP P66930
H	255	HIS	-	EXPRESSION TAG	UNP P66930
H	256	HIS	-	EXPRESSION TAG	UNP P66930
H	257	HIS	-	EXPRESSION TAG	UNP P66930
H	258	HIS	-	EXPRESSION TAG	UNP P66930

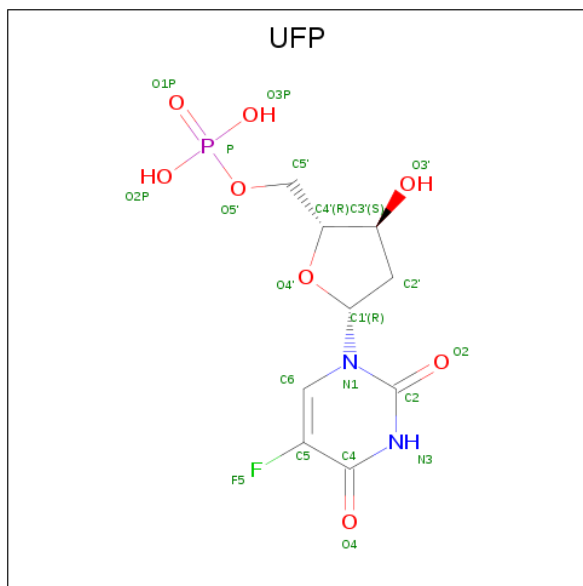
- # FAD

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

- Molecule 3 is 5-FLUORO-2'-DEOXYURIDINE-5'-MONOPHOSPHATE (three-letter code: UFP) (formula: $C_9H_{12}FN_2O_8P$).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
3	A	1	Total 21	C 9	F 1	N 2	O 8	P 1	0	0
3	B	1	Total 21	C 9	F 1	N 2	O 8	P 1	0	0
3	C	1	Total 21	C 9	F 1	N 2	O 8	P 1	0	0
3	D	1	Total 21	C 9	F 1	N 2	O 8	P 1	0	0
3	E	1	Total 21	C 9	F 1	N 2	O 8	P 1	0	0
3	F	1	Total 21	C 9	F 1	N 2	O 8	P 1	0	0
3	G	1	Total 21	C 9	F 1	N 2	O 8	P 1	0	0
3	H	1	Total 21	C 9	F 1	N 2	O 8	P 1	0	0

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	C	1	Total	C	O	0	0
			6	3	3		

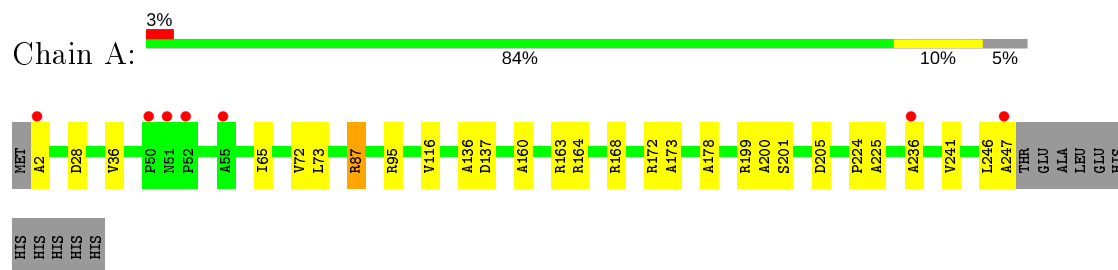
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	135	Total	O	0	0
			135	135		
5	B	145	Total	O	0	0
			145	145		
5	C	116	Total	O	0	0
			116	116		
5	D	119	Total	O	0	0
			119	119		
5	E	116	Total	O	0	0
			116	116		
5	F	94	Total	O	0	0
			94	94		
5	G	85	Total	O	0	0
			85	85		
5	H	127	Total	O	0	0
			127	127		

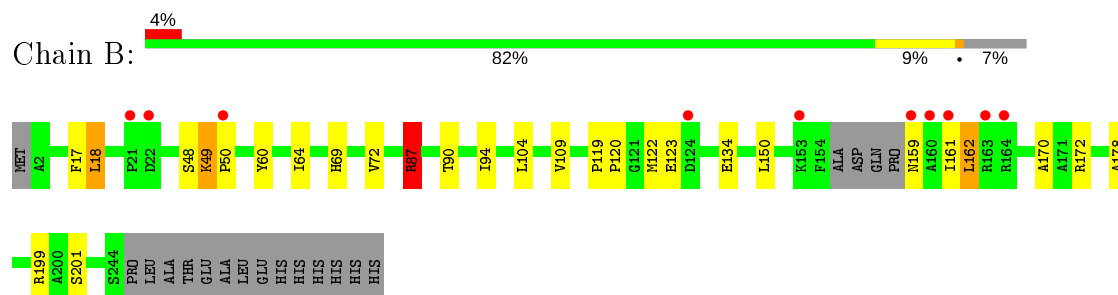
3 Residue-property plots [i](#)

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.

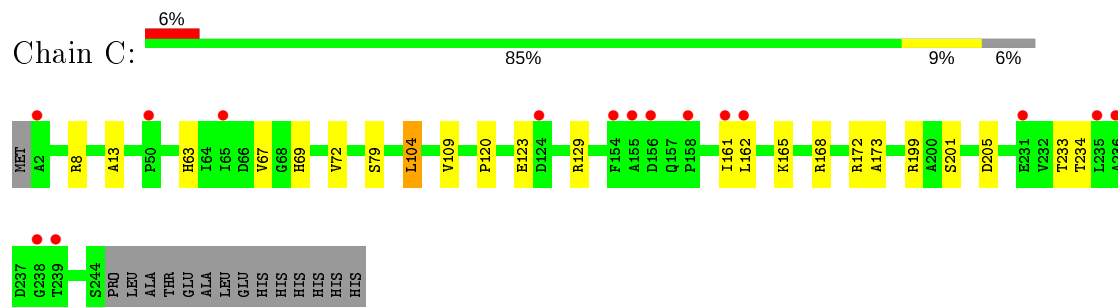
- Molecule 1: Thymidylate synthase thyX



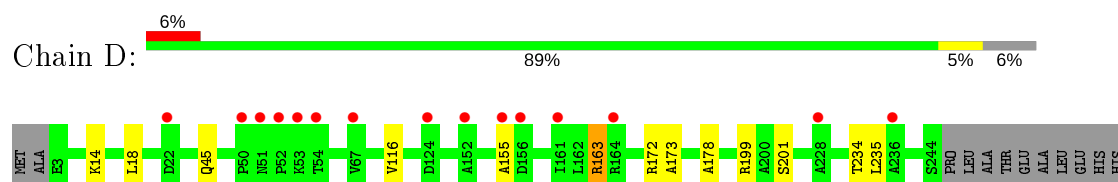
- Molecule 1: Thymidylate synthase thyX



- Molecule 1: Thymidylate synthase thyX




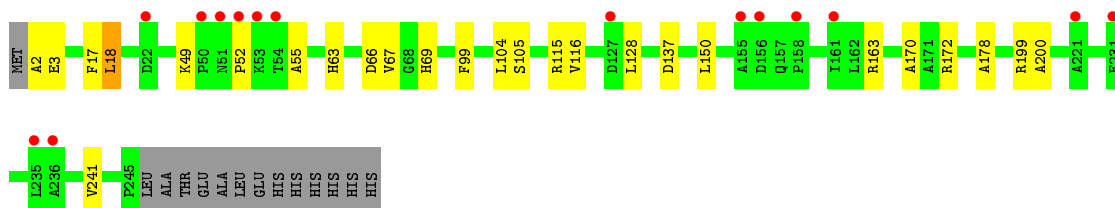
- Molecule 1: Thymidylate synthase thyX




HIS
HIS
HIS
HIS

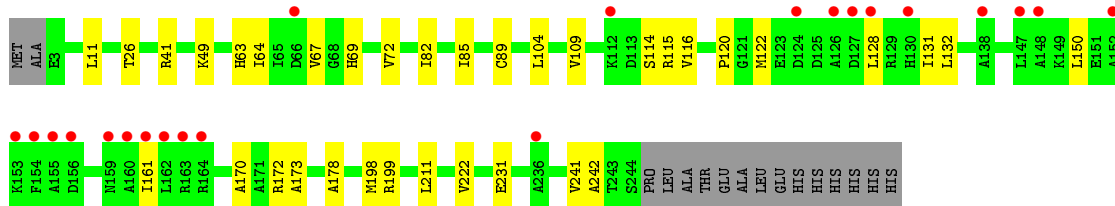
- Molecule 1: Thymidylate synthase thyX

Chain E: 




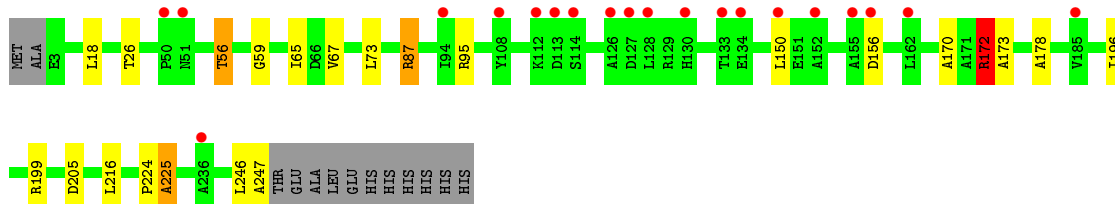
- Molecule 1: Thymidylate synthase thyX

Chain F: 

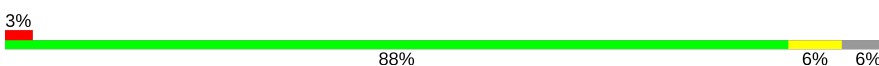


- Molecule 1: Thymidylate synthase thyX

Chain G: 



- Molecule 1: Thymidylate synthase thyX

Chain H: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	81.04Å 78.45Å 168.96Å 90.00° 96.80° 90.00°	Depositor
Resolution (Å)	48.70 – 1.90 48.71 – 1.90	Depositor EDS
% Data completeness (in resolution range)	(Not available) (48.70-1.90) 91.6 (48.71-1.90)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.88 (at 1.90Å)	Xtriage
Refinement program	REFMAC 5.5.0088	Depositor
R, R_{free}	0.187 , 0.229 0.191 , 0.232	Depositor DCC
R_{free} test set	7619 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	21.2	Xtriage
Anisotropy	0.119	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 55.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	16612	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.16% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: UFP, GOL, 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	0.62	0/1956	0.71	1/2667 (0.0%)
1	B	0.65	0/1896	0.71	2/2584 (0.1%)
1	C	0.60	0/1925	0.70	1/2626 (0.0%)
1	D	0.62	0/1920	0.68	1/2619 (0.0%)
1	E	0.56	0/1930	0.67	0/2633
1	F	0.55	0/1927	0.66	0/2626
1	G	0.55	0/1941	0.68	3/2649 (0.1%)
1	H	0.59	0/1926	0.68	0/2627
All	All	0.59	0/15421	0.69	8/21031 (0.0%)

There are no bond length outliers.

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	87	ARG	NE-CZ-NH2	-8.51	116.04	120.30
1	B	87	ARG	NE-CZ-NH2	-8.06	116.27	120.30
1	G	172	ARG	NE-CZ-NH2	-7.45	116.58	120.30
1	G	172	ARG	NE-CZ-NH1	6.70	123.65	120.30
1	C	168	ARG	NE-CZ-NH2	-6.09	117.25	120.30
1	G	87	ARG	NE-CZ-NH2	-5.48	117.56	120.30
1	D	163	ARG	NE-CZ-NH1	5.41	123.00	120.30
1	B	87	ARG	NE-CZ-NH1	5.25	122.92	120.30

There are no chirality outliers.

There are no planarity outliers.

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	1912	0	1885	20	0
1	B	1855	0	1815	25	0
1	C	1882	0	1840	17	0
1	D	1877	0	1835	7	0
1	E	1887	0	1841	20	0
1	F	1884	0	1848	27	0
1	G	1897	0	1858	19	0
1	H	1883	0	1848	12	0
2	A	53	0	31	0	0
2	B	53	0	31	0	0
2	C	53	0	31	0	0
2	D	53	0	31	0	0
2	E	53	0	31	0	0
2	F	53	0	31	0	0
2	G	53	0	31	0	0
2	H	53	0	31	0	0
3	A	21	0	10	2	0
3	B	21	0	9	2	0
3	C	21	0	10	1	0
3	D	21	0	9	3	0
3	E	21	0	9	2	0
3	F	21	0	9	2	0
3	G	21	0	9	4	0
3	H	21	0	10	1	0
4	C	6	0	8	0	0
5	A	135	0	0	5	0
5	B	145	0	0	2	0
5	C	116	0	0	1	0
5	D	119	0	0	0	0
5	E	116	0	0	4	0
5	F	94	0	0	4	0
5	G	85	0	0	1	0
5	H	127	0	0	1	0
All	All	16612	0	15101	134	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 4.

All (134) 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:B:161:ILE:HG23	1:B:162:LEU:HD13	1.42	1.02
1:A:199:ARG:HH12	3:D:260:UFP:HN3	1.14	0.95
1:C:161:ILE:HG22	1:C:162:LEU:HD23	1.50	0.93
3:A:260:UFP:HN3	1:D:199:ARG:HH12	1.11	0.93
1:F:199:ARG:HH12	3:G:260:UFP:HN3	1.15	0.89
3:F:260:UFP:HN3	1:G:199:ARG:HH12	1.18	0.88
3:E:260:UFP:HN3	1:H:199:ARG:HH12	1.21	0.88
1:C:123:GLU:O	1:C:129:ARG:NH1	2.07	0.87
3:B:260:UFP:HN3	1:C:199:ARG:HH12	1.23	0.84
1:B:199:ARG:HH12	3:C:260:UFP:HN3	1.22	0.84
1:E:199:ARG:HH12	3:H:260:UFP:HN3	1.25	0.84
1:F:69:HIS:O	1:F:72:VAL:HG12	1.78	0.83
1:E:67:VAL:HG13	1:E:69:HIS:CD2	2.14	0.82
1:F:41:ARG:CZ	5:F:898:HOH:O	2.26	0.81
1:F:41:ARG:NH1	5:F:898:HOH:O	2.17	0.78
1:C:69:HIS:O	1:C:72:VAL:HG12	1.85	0.76
1:A:168:ARG:CZ	5:A:645:HOH:O	2.35	0.73
1:C:67:VAL:HG12	1:C:69:HIS:NE2	2.04	0.73
1:G:56:THR:HG23	1:G:59:GLY:H	1.53	0.73
1:E:2:ALA:HB2	1:E:137:ASP:OD1	1.90	0.71
1:G:87:ARG:NH2	3:G:260:UFP:O3P	2.25	0.68
1:B:134:GLU:OE1	5:B:905:HOH:O	2.11	0.68
1:E:52:PRO:HA	1:E:55:ALA:HB2	1.76	0.67
1:B:49:LYS:CB	5:B:708:HOH:O	2.43	0.66
1:B:161:ILE:CG2	1:B:162:LEU:HD13	2.22	0.66
1:A:95:ARG:HG3	3:D:260:UFP:H2'2	1.77	0.66
1:A:2:ALA:HB2	1:A:137:ASP:OD1	1.95	0.65
1:B:104:LEU:HD13	1:D:45:GLN:HG2	1.76	0.65
1:G:224:PRO:O	1:G:225:ALA:HB3	1.98	0.64
1:A:87:ARG:NH2	3:A:260:UFP:O3P	2.31	0.64
1:F:104:LEU:CD1	1:F:109:VAL:HG21	2.29	0.62
1:C:63:HIS:CE1	1:C:67:VAL:HG21	2.36	0.61
1:E:99:PHE:CZ	5:E:284:HOH:O	2.52	0.59
1:F:173:ALA:HB1	1:G:178:ALA:HB2	1.83	0.58
1:B:69:HIS:O	1:B:72:VAL:HG22	2.03	0.58
1:A:205:ASP:OD2	5:A:835:HOH:O	2.17	0.58
1:A:65:ILE:HG22	1:A:246:LEU:HD13	1.86	0.57
1:C:162:LEU:HD22	1:C:165:LYS:HE3	1.86	0.57
1:F:199:ARG:NH1	3:G:260:UFP:HN3	1.96	0.56
1:D:18:LEU:HD22	1:H:112:LYS:HE2	1.88	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:82:ILE:HD13	1:F:211:LEU:HD11	1.89	0.55
1:B:161:ILE:HG23	1:B:162:LEU:CD1	2.28	0.55
1:B:104:LEU:CD1	1:B:109:VAL:HG21	2.36	0.54
1:B:159:ASN:OD1	1:B:161:ILE:HG22	2.07	0.54
1:A:36:VAL:HG13	1:A:72:VAL:HG12	1.89	0.54
1:F:122:MET:HE2	1:G:150:LEU:HD11	1.90	0.54
1:C:162:LEU:HD22	1:C:165:LYS:CE	2.38	0.53
1:B:87:ARG:NH2	3:B:260:UFP:O3P	2.42	0.53
1:A:178:ALA:HB2	1:D:173:ALA:HB1	1.90	0.53
1:F:104:LEU:HD13	1:H:45:GLN:HG2	1.91	0.53
1:C:104:LEU:HD13	1:C:109:VAL:HG21	1.90	0.53
1:E:2:ALA:HB3	5:E:272:HOH:O	2.08	0.53
1:H:200:ALA:O	1:H:241:VAL:HG13	2.09	0.53
1:F:64:ILE:HD13	1:F:72:VAL:HG11	1.92	0.52
1:B:104:LEU:HD12	1:B:109:VAL:HG21	1.92	0.52
1:B:120:PRO:HA	1:B:123:GLU:HG3	1.92	0.51
1:G:205:ASP:OD2	5:G:581:HOH:O	2.19	0.51
1:G:67:VAL:O	1:G:67:VAL:HG22	2.11	0.51
1:B:119:PRO:HG2	1:B:122:MET:HE3	1.93	0.50
1:C:67:VAL:CG1	1:C:69:HIS:NE2	2.73	0.50
3:E:260:UFP:HN3	1:H:199:ARG:NH1	2.00	0.50
1:B:60:TYR:CE2	1:B:64:ILE:HD11	2.46	0.50
1:H:104:LEU:C	1:H:104:LEU:HD23	2.31	0.50
1:C:63:HIS:CE1	1:C:67:VAL:CG2	2.95	0.50
1:A:160:ALA:HB1	1:A:164:ARG:NH2	2.27	0.49
1:E:3:GLU:HB3	1:E:115:ARG:HD3	1.95	0.49
1:G:65:ILE:HG22	1:G:246:LEU:HD13	1.94	0.49
1:E:52:PRO:HA	1:E:55:ALA:CB	2.42	0.48
1:B:48:SER:O	1:B:50:PRO:HD3	2.13	0.48
1:G:56:THR:HG23	1:G:59:GLY:N	2.25	0.48
1:B:17:PHE:C	1:B:18:LEU:HD13	2.34	0.48
1:F:104:LEU:HD12	1:F:109:VAL:HG21	1.95	0.48
1:H:120:PRO:HA	1:H:123:GLU:HG3	1.95	0.48
1:C:205:ASP:OD2	5:C:881:HOH:O	2.20	0.47
1:F:161:ILE:HG22	5:F:847:HOH:O	2.13	0.47
1:E:99:PHE:CE2	5:E:284:HOH:O	2.68	0.47
1:B:17:PHE:O	1:B:18:LEU:HD13	2.16	0.46
1:B:60:TYR:HE2	1:B:64:ILE:HD11	1.79	0.46
1:E:104:LEU:HD23	1:E:104:LEU:C	2.35	0.46
1:G:224:PRO:O	1:G:225:ALA:CB	2.62	0.46
1:H:3:GLU:N	5:H:575:HOH:O	2.48	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:90:THR:O	1:B:94:ILE:HG12	2.16	0.46
1:D:116:VAL:HG13	1:D:178:ALA:HB1	1.98	0.46
1:E:200:ALA:O	1:E:241:VAL:HG13	2.16	0.45
1:G:172:ARG:NH2	3:G:260:UFP:O3P	2.49	0.45
1:A:200:ALA:O	1:A:241:VAL:HG13	2.17	0.45
1:B:119:PRO:HG2	1:B:122:MET:CE	2.46	0.45
1:F:67:VAL:HG22	1:F:67:VAL:O	2.17	0.45
1:F:11:LEU:HD11	1:F:222:VAL:HG21	1.99	0.45
1:F:128:LEU:HA	1:F:131:ILE:HD12	1.98	0.45
1:A:247:ALA:HB2	5:A:539:HOH:O	2.17	0.45
1:C:13:ALA:HB3	1:C:79:SER:HB2	1.99	0.45
1:F:122:MET:HE3	1:G:170:ALA:HB1	1.99	0.45
1:F:122:MET:CE	1:G:170:ALA:HB1	2.47	0.45
3:F:260:UFP:H2'2	1:G:95:ARG:HG3	1.99	0.44
1:A:160:ALA:HB1	1:A:164:ARG:HH21	1.82	0.44
1:B:161:ILE:HG23	1:B:162:LEU:N	2.31	0.44
1:G:65:ILE:HG22	1:G:246:LEU:CD1	2.47	0.44
1:B:161:ILE:CG2	1:B:162:LEU:N	2.80	0.44
1:C:120:PRO:HA	1:C:123:GLU:HG3	2.00	0.44
1:D:234:THR:HG22	1:D:235:LEU:O	2.18	0.44
1:H:18:LEU:N	1:H:18:LEU:HD12	2.32	0.44
1:E:128:LEU:HD21	1:H:149:LYS:HB3	2.00	0.44
1:A:116:VAL:HG13	5:A:264:HOH:O	2.18	0.43
1:C:104:LEU:HD12	1:C:104:LEU:C	2.38	0.43
1:F:115:ARG:NE	5:F:611:HOH:O	2.52	0.43
1:B:178:ALA:HB2	1:C:173:ALA:HB1	2.00	0.43
1:F:122:MET:HE3	1:F:122:MET:HB2	1.82	0.43
1:A:116:VAL:HG11	1:A:136:ALA:CB	2.48	0.43
1:A:173:ALA:HB1	1:D:178:ALA:HB2	2.00	0.43
1:F:26:THR:O	1:F:49:LYS:NZ	2.44	0.42
1:G:196:ILE:HD13	1:G:216:LEU:HD13	2.02	0.42
1:F:231:GLU:O	1:F:242:ALA:HA	2.20	0.42
1:G:246:LEU:O	1:G:247:ALA:C	2.57	0.42
1:B:150:LEU:HD13	1:B:170:ALA:HB3	2.01	0.41
1:A:236:ALA:HB3	5:A:850:HOH:O	2.20	0.41
1:E:163:ARG:HD2	5:E:845:HOH:O	2.21	0.41
1:A:36:VAL:HG11	1:A:73:LEU:HD23	2.02	0.41
1:F:63:HIS:O	1:F:67:VAL:HG12	2.19	0.41
1:E:104:LEU:HD23	1:E:105:SER:N	2.36	0.41
1:F:120:PRO:C	1:F:122:MET:H	2.24	0.41
1:C:233:THR:HG22	1:C:234:THR:N	2.36	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:95:ARG:CG	3:D:260:UFP:H2'2	2.48	0.41
1:H:20:PRO:HA	1:H:21:PRO:HD3	1.90	0.41
1:E:150:LEU:HD13	1:E:170:ALA:HB3	2.01	0.41
1:E:17:PHE:C	1:E:18:LEU:HD13	2.41	0.41
1:F:178:ALA:HB2	1:G:173:ALA:HB1	2.02	0.41
1:E:67:VAL:CG1	1:E:69:HIS:CD2	2.94	0.41
1:F:85:ILE:HG12	1:F:89:CYS:HB3	2.03	0.41
1:E:178:ALA:HB2	1:H:173:ALA:HB1	2.03	0.40
1:A:224:PRO:O	1:A:225:ALA:HB3	2.21	0.40
1:F:150:LEU:HD13	1:F:170:ALA:HB3	2.03	0.40
1:E:116:VAL:HG23	1:E:178:ALA:CB	2.52	0.40
1:E:63:HIS:O	1:E:67:VAL:HG12	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	244/258 (95%)	241 (99%)	2 (1%)	1 (0%)	34	24
1	B	235/258 (91%)	232 (99%)	2 (1%)	1 (0%)	34	24
1	C	241/258 (93%)	234 (97%)	7 (3%)	0	100	100
1	D	240/258 (93%)	236 (98%)	3 (1%)	1 (0%)	34	24
1	E	242/258 (94%)	237 (98%)	4 (2%)	1 (0%)	34	24
1	F	240/258 (93%)	232 (97%)	8 (3%)	0	100	100
1	G	243/258 (94%)	232 (96%)	10 (4%)	1 (0%)	34	24
1	H	241/258 (93%)	238 (99%)	3 (1%)	0	100	100
All	All	1926/2064 (93%)	1882 (98%)	39 (2%)	5 (0%)	41	31

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	28	ASP
1	D	155	ALA
1	G	225	ALA
1	E	49	LYS
1	B	49	LYS

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	197/209 (94%)	194 (98%)	3 (2%)	65	62
1	B	190/209 (91%)	185 (97%)	5 (3%)	46	39
1	C	193/209 (92%)	189 (98%)	4 (2%)	53	48
1	D	193/209 (92%)	189 (98%)	4 (2%)	53	48
1	E	193/209 (92%)	190 (98%)	3 (2%)	62	60
1	F	194/209 (93%)	188 (97%)	6 (3%)	40	32
1	G	195/209 (93%)	189 (97%)	6 (3%)	40	32
1	H	194/209 (93%)	190 (98%)	4 (2%)	53	48
All	All	1549/1672 (93%)	1514 (98%)	35 (2%)	50	45

All (35) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	163	ARG
1	A	172	ARG
1	A	201	SER
1	B	18	LEU
1	B	87	ARG
1	B	162	LEU
1	B	172	ARG
1	B	201	SER
1	C	8	ARG
1	C	104	LEU
1	C	172	ARG

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Mol	Chain	Res	Type
1	C	201	SER
1	D	14	LYS
1	D	163	ARG
1	D	172	ARG
1	D	201	SER
1	E	18	LEU
1	E	66	ASP
1	E	172	ARG
1	F	114	SER
1	F	116	VAL
1	F	132	LEU
1	F	172	ARG
1	F	198	MET
1	F	241	VAL
1	G	18	LEU
1	G	26	THR
1	G	56	THR
1	G	73	LEU
1	G	156	ASP
1	G	172	ARG
1	H	3	GLU
1	H	172	ARG
1	H	201	SER
1	H	231	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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

17 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
4	GOL	C	261	-	5,5,5	0.69	0	5,5,5	0.38	0
3	UFP	H	260	-	19,22,22	1.35	3 (15%)	24,33,33	1.93	4 (16%)
2	FAD	D	259	-	51,58,58	1.00	2 (3%)	60,89,89	1.68	9 (15%)
2	FAD	F	259	-	51,58,58	0.98	2 (3%)	60,89,89	1.58	9 (15%)
3	UFP	C	260	-	19,22,22	1.16	2 (10%)	24,33,33	2.18	5 (20%)
3	UFP	A	260	-	19,22,22	1.41	3 (15%)	24,33,33	2.09	3 (12%)
2	FAD	B	259	-	51,58,58	1.09	3 (5%)	60,89,89	1.80	11 (18%)
3	UFP	G	260	-	19,22,22	1.28	2 (10%)	24,33,33	2.41	5 (20%)
3	UFP	E	260	-	19,22,22	1.19	1 (5%)	24,33,33	1.94	4 (16%)
3	UFP	D	260	-	19,22,22	1.12	1 (5%)	24,33,33	2.79	10 (41%)
2	FAD	H	259	-	51,58,58	1.08	4 (7%)	60,89,89	1.86	12 (20%)
2	FAD	E	259	-	51,58,58	1.04	4 (7%)	60,89,89	1.65	11 (18%)
3	UFP	B	260	-	19,22,22	1.17	1 (5%)	24,33,33	2.31	6 (25%)
2	FAD	G	259	-	51,58,58	1.02	2 (3%)	60,89,89	1.76	10 (16%)
2	FAD	A	259	-	51,58,58	1.07	4 (7%)	60,89,89	1.52	9 (15%)
3	UFP	F	260	-	19,22,22	1.10	1 (5%)	24,33,33	1.90	6 (25%)
2	FAD	C	259	-	51,58,58	1.06	4 (7%)	60,89,89	1.55	10 (16%)

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
4	GOL	C	261	-	-	1/4/4/4	-
3	UFP	H	260	-	-	0/7/22/22	0/2/2/2
2	FAD	D	259	-	-	5/30/50/50	0/6/6/6

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	FAD	F	259	-	-	7/30/50/50	0/6/6/6
3	UFP	C	260	-	-	1/7/22/22	0/2/2/2
3	UFP	A	260	-	-	1/7/22/22	0/2/2/2
2	FAD	B	259	-	-	7/30/50/50	0/6/6/6
3	UFP	G	260	-	-	0/7/22/22	0/2/2/2
3	UFP	E	260	-	-	0/7/22/22	0/2/2/2
3	UFP	D	260	-	1/1/4/4	0/7/22/22	0/2/2/2
2	FAD	H	259	-	-	8/30/50/50	0/6/6/6
2	FAD	E	259	-	-	6/30/50/50	0/6/6/6
3	UFP	B	260	-	-	0/7/22/22	0/2/2/2
2	FAD	G	259	-	-	7/30/50/50	0/6/6/6
2	FAD	A	259	-	-	7/30/50/50	0/6/6/6
3	UFP	F	260	-	1/1/4/4	1/7/22/22	0/2/2/2
2	FAD	C	259	-	-	6/30/50/50	0/6/6/6

All (39) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	H	259	FAD	C4X-N5	3.65	1.38	1.33
3	A	260	UFP	C4-N3	3.51	1.39	1.33
2	E	259	FAD	C4-N3	3.41	1.39	1.33
2	D	259	FAD	C10-N1	3.27	1.37	1.33
2	D	259	FAD	C4X-N5	3.27	1.38	1.33
2	E	259	FAD	C4X-N5	3.23	1.38	1.33
3	B	260	UFP	C4-N3	3.22	1.38	1.33
3	G	260	UFP	C4-N3	3.22	1.38	1.33
3	D	260	UFP	C4-N3	3.18	1.38	1.33
2	G	259	FAD	C10-N1	3.16	1.37	1.33
3	H	260	UFP	C4-N3	3.14	1.38	1.33
3	E	260	UFP	C4-N3	3.08	1.38	1.33
2	C	259	FAD	C10-N1	3.02	1.37	1.33
2	B	259	FAD	C4X-C10	3.00	1.41	1.38
3	C	260	UFP	C4-N3	2.99	1.38	1.33
2	B	259	FAD	O4B-C1B	2.85	1.45	1.41
2	C	259	FAD	O4B-C1B	2.80	1.45	1.41
2	A	259	FAD	C10-N1	2.75	1.36	1.33
2	C	259	FAD	C4-N3	2.73	1.37	1.33
2	H	259	FAD	C4X-C10	2.60	1.41	1.38
3	F	260	UFP	C4-N3	2.57	1.37	1.33
2	F	259	FAD	C4-N3	2.55	1.37	1.33
2	B	259	FAD	C4X-N5	2.49	1.36	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	H	260	UFP	O4'-C1'	-2.48	1.36	1.42
2	F	259	FAD	C10-N1	2.43	1.36	1.33
2	E	259	FAD	O4B-C1B	2.42	1.44	1.41
3	A	260	UFP	O4'-C1'	-2.40	1.37	1.42
2	G	259	FAD	C4X-N5	2.38	1.36	1.33
2	C	259	FAD	C4X-N5	2.33	1.36	1.33
2	H	259	FAD	O4B-C1B	2.27	1.44	1.41
3	C	260	UFP	C1'-N1	-2.26	1.42	1.49
3	A	260	UFP	C1'-N1	-2.25	1.42	1.49
3	G	260	UFP	P-O3P	-2.20	1.46	1.54
2	A	259	FAD	C4X-N5	2.15	1.36	1.33
2	H	259	FAD	C10-N1	2.14	1.36	1.33
2	E	259	FAD	C10-N1	2.13	1.36	1.33
2	A	259	FAD	C4-N3	2.10	1.36	1.33
3	H	260	UFP	C1'-N1	-2.07	1.43	1.49
2	A	259	FAD	C2A-N3A	2.00	1.35	1.32

All (124) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	260	UFP	C4-N3-C2	8.10	121.98	115.14
3	G	260	UFP	C4-N3-C2	8.04	121.93	115.14
3	C	260	UFP	C4-N3-C2	7.99	121.89	115.14
3	A	260	UFP	C4-N3-C2	7.96	121.86	115.14
3	D	260	UFP	C4-N3-C2	7.18	121.20	115.14
2	G	259	FAD	C4-N3-C2	7.03	121.07	115.14
3	E	260	UFP	C4-N3-C2	6.97	121.02	115.14
2	H	259	FAD	C4-N3-C2	6.84	120.92	115.14
3	H	260	UFP	C4-N3-C2	6.83	120.91	115.14
2	B	259	FAD	C4-N3-C2	6.21	120.38	115.14
2	C	259	FAD	C4-N3-C2	6.16	120.34	115.14
2	D	259	FAD	C4-N3-C2	5.99	120.20	115.14
3	F	260	UFP	C4-N3-C2	5.75	120.00	115.14
2	B	259	FAD	C1'-N10-C9A	5.65	122.74	118.29
2	F	259	FAD	C4-N3-C2	5.50	119.78	115.14
3	D	260	UFP	C4'-O4'-C1'	-5.46	96.28	109.45
2	F	259	FAD	N3A-C2A-N1A	-5.39	120.25	128.68
2	H	259	FAD	N3A-C2A-N1A	-5.26	120.45	128.68
2	A	259	FAD	N3A-C2A-N1A	-5.18	120.58	128.68
2	G	259	FAD	N3A-C2A-N1A	-5.10	120.71	128.68
2	E	259	FAD	C4-N3-C2	5.04	119.39	115.14
2	E	259	FAD	N3A-C2A-N1A	-4.98	120.90	128.68

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	259	FAD	C1'-N10-C9A	4.89	122.14	118.29
2	B	259	FAD	N3A-C2A-N1A	-4.67	121.39	128.68
3	D	260	UFP	O4'-C1'-C2'	4.60	114.94	106.25
3	G	260	UFP	O4'-C1'-C2'	4.59	114.93	106.25
2	A	259	FAD	C4-N3-C2	4.53	118.97	115.14
2	D	259	FAD	N3A-C2A-N1A	-4.49	121.67	128.68
2	C	259	FAD	N3A-C2A-N1A	-4.27	122.00	128.68
2	B	259	FAD	C5X-C9A-N10	4.25	120.80	117.72
3	D	260	UFP	C2'-C1'-N1	4.21	123.97	114.27
3	D	260	UFP	P-O5'-C5'	4.07	129.51	118.30
2	E	259	FAD	C4X-N5-C5X	4.06	120.83	116.77
2	A	259	FAD	C1'-N10-C9A	3.98	121.42	118.29
2	G	259	FAD	C5X-C9A-N10	3.98	120.60	117.72
3	H	260	UFP	O4'-C1'-C2'	3.96	113.73	106.25
2	D	259	FAD	C1'-N10-C9A	3.89	121.35	118.29
2	D	259	FAD	C4X-N5-C5X	3.87	120.64	116.77
3	G	260	UFP	C5-C4-N3	-3.85	118.33	122.39
3	C	260	UFP	C5-C4-N3	-3.84	118.33	122.39
3	B	260	UFP	O4'-C1'-C2'	3.81	113.45	106.25
2	G	259	FAD	C1'-N10-C9A	3.77	121.26	118.29
2	F	259	FAD	C4X-N5-C5X	3.75	120.52	116.77
2	E	259	FAD	C4-C4X-N5	3.67	122.79	118.60
3	G	260	UFP	C4'-O4'-C1'	-3.58	100.80	109.45
2	D	259	FAD	C5X-C9A-N10	3.58	120.31	117.72
3	B	260	UFP	C4'-O4'-C1'	-3.51	100.97	109.45
2	H	259	FAD	C4-C4X-N5	3.48	122.58	118.60
2	E	259	FAD	C1'-N10-C9A	3.47	121.02	118.29
3	A	260	UFP	C5-C4-N3	-3.46	118.74	122.39
3	F	260	UFP	O4'-C1'-C2'	3.37	112.61	106.25
3	D	260	UFP	C6-N1-C1'	-3.32	111.79	119.24
2	C	259	FAD	C10-C4X-N5	-3.31	118.97	121.26
3	A	260	UFP	F5-C5-C4	3.31	124.89	118.90
3	D	260	UFP	C3'-C2'-C1'	-3.31	94.25	102.54
2	C	259	FAD	C4-C4X-N5	3.30	122.36	118.60
3	D	260	UFP	F5-C5-C4	3.29	124.85	118.90
3	E	260	UFP	O4'-C1'-C2'	3.28	112.44	106.25
3	F	260	UFP	P-O5'-C5'	3.21	127.13	118.30
2	E	259	FAD	C10-C4X-N5	-3.19	119.06	121.26
2	H	259	FAD	C5X-C9A-N10	3.12	119.98	117.72
2	D	259	FAD	O2A-PA-O1A	3.12	127.66	112.24
2	F	259	FAD	C10-C4X-N5	-3.09	119.12	121.26
2	C	259	FAD	C4X-N5-C5X	3.05	119.82	116.77

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	259	FAD	C4X-N5-C5X	3.05	119.81	116.77
2	F	259	FAD	C1'-N10-C9A	3.02	120.67	118.29
2	B	259	FAD	C4-C4X-N5	2.99	122.02	118.60
3	B	260	UFP	F5-C5-C4	2.95	124.24	118.90
2	F	259	FAD	C4X-C4-N3	-2.90	119.46	123.43
2	H	259	FAD	C4-C4X-C10	-2.88	118.05	119.95
2	E	259	FAD	O2A-PA-O1A	2.83	126.22	112.24
2	B	259	FAD	C4-C4X-C10	-2.82	118.08	119.95
2	G	259	FAD	C3B-C2B-C1B	2.80	105.19	100.98
2	H	259	FAD	O2A-PA-O1A	2.75	125.85	112.24
3	C	260	UFP	O4'-C1'-C2'	2.73	111.41	106.25
2	H	259	FAD	C10-C4X-N5	-2.73	119.37	121.26
2	E	259	FAD	C4-C4X-C10	-2.72	118.15	119.95
3	B	260	UFP	C5-C4-N3	-2.71	119.53	122.39
3	E	260	UFP	C4'-O4'-C1'	-2.69	102.95	109.45
3	F	260	UFP	C4'-O4'-C1'	-2.66	103.04	109.45
2	H	259	FAD	C4X-C4-N3	-2.63	119.84	123.43
2	F	259	FAD	C3B-C2B-C1B	2.59	104.87	100.98
2	G	259	FAD	C4X-C4-N3	-2.59	119.90	123.43
3	F	260	UFP	C2'-C1'-N1	2.57	120.19	114.27
2	C	259	FAD	O2A-PA-O1A	2.56	124.92	112.24
3	G	260	UFP	C2'-C1'-N1	2.55	120.16	114.27
2	A	259	FAD	C5X-C9A-N10	2.55	119.56	117.72
2	C	259	FAD	C1'-N10-C9A	2.51	120.27	118.29
2	B	259	FAD	C4X-C4-N3	-2.50	120.01	123.43
2	B	259	FAD	O2A-PA-O1A	2.49	124.55	112.24
2	G	259	FAD	C9A-N10-C10	-2.46	118.69	121.91
2	D	259	FAD	C9A-C5X-N5	-2.43	118.56	122.36
2	B	259	FAD	C9A-N10-C10	-2.41	118.75	121.91
2	C	259	FAD	C3B-C2B-C1B	2.40	104.59	100.98
2	G	259	FAD	C4X-N5-C5X	2.38	119.15	116.77
3	F	260	UFP	C5-C4-N3	-2.36	119.90	122.39
2	E	259	FAD	C5X-C9A-N10	2.36	119.42	117.72
3	D	260	UFP	C2'-C3'-C4'	2.33	107.62	102.76
2	H	259	FAD	C4'-C3'-C2'	2.33	118.21	113.36
2	F	259	FAD	O2B-C2B-C3B	-2.31	104.36	111.82
3	D	260	UFP	C5-C4-N3	-2.30	119.97	122.39
3	B	260	UFP	C2'-C1'-N1	2.28	119.53	114.27
3	C	260	UFP	P-O5'-C5'	2.27	124.54	118.30
2	G	259	FAD	O2A-PA-O1A	2.25	123.38	112.24
2	B	259	FAD	C4X-N5-C5X	2.25	119.02	116.77
3	C	260	UFP	O5'-P-O1P	-2.24	100.19	106.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	259	FAD	C3B-C2B-C1B	2.23	104.33	100.98
2	F	259	FAD	C4-C4X-N5	2.20	121.11	118.60
2	A	259	FAD	C4-C4X-N5	2.19	121.10	118.60
2	G	259	FAD	C2A-N1A-C6A	2.16	122.45	118.75
2	E	259	FAD	C6-C5X-N5	2.13	121.39	119.05
2	A	259	FAD	C1B-N9A-C4A	-2.12	122.91	126.64
3	H	260	UFP	C4'-O4'-C1'	-2.12	104.34	109.45
2	A	259	FAD	O2A-PA-O1A	2.11	122.67	112.24
2	D	259	FAD	C4-C4X-N5	2.10	121.00	118.60
2	C	259	FAD	C4A-C5A-N7A	-2.09	107.22	109.40
3	E	260	UFP	O3P-P-O2P	2.08	115.60	107.64
2	C	259	FAD	C4X-C4-N3	-2.07	120.60	123.43
2	A	259	FAD	C10-C4X-N5	-2.05	119.84	121.26
2	D	259	FAD	C4-C4X-C10	-2.05	118.60	119.95
3	H	260	UFP	O5'-P-O1P	2.04	112.20	106.47
2	B	259	FAD	O3'-C3'-C4'	-2.04	103.89	108.81
2	E	259	FAD	C9A-C5X-N5	-2.02	119.20	122.36
2	A	259	FAD	C4X-C4-N3	-2.00	120.69	123.43

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	D	260	UFP	C1'
3	F	260	UFP	C1'

All (57) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	D	259	FAD	O4'-C4'-C5'-O5'
2	F	259	FAD	O4'-C4'-C5'-O5'
2	F	259	FAD	PA-O3P-P-O5'
2	B	259	FAD	C5B-O5B-PA-O2A
2	H	259	FAD	C5B-O5B-PA-O2A
2	H	259	FAD	C3'-C4'-C5'-O5'
2	H	259	FAD	O4'-C4'-C5'-O5'
2	E	259	FAD	O4'-C4'-C5'-O5'
2	G	259	FAD	C3'-C4'-C5'-O5'
2	G	259	FAD	O4'-C4'-C5'-O5'
2	A	259	FAD	C5B-O5B-PA-O2A
2	C	259	FAD	C3'-C4'-C5'-O5'
2	C	259	FAD	O4'-C4'-C5'-O5'
4	C	261	GOL	C1-C2-C3-O3

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Mol	Chain	Res	Type	Atoms
2	D	259	FAD	C3'-C4'-C5'-O5'
2	F	259	FAD	C3'-C4'-C5'-O5'
2	B	259	FAD	C3'-C4'-C5'-O5'
2	E	259	FAD	C3'-C4'-C5'-O5'
3	C	260	UFP	C5'-O5'-P-O1P
2	D	259	FAD	PA-O3P-P-O1P
2	B	259	FAD	P-O3P-PA-O1A
2	B	259	FAD	PA-O3P-P-O5'
2	H	259	FAD	PA-O3P-P-O5'
2	E	259	FAD	PA-O3P-P-O5'
2	G	259	FAD	PA-O3P-P-O5'
2	A	259	FAD	PA-O3P-P-O5'
2	C	259	FAD	PA-O3P-P-O5'
2	F	259	FAD	C5B-O5B-PA-O3P
2	B	259	FAD	C5B-O5B-PA-O3P
2	G	259	FAD	C5B-O5B-PA-O3P
2	A	259	FAD	C5B-O5B-PA-O3P
2	F	259	FAD	P-O3P-PA-O2A
2	G	259	FAD	P-O3P-PA-O2A
2	A	259	FAD	P-O3P-PA-O2A
2	C	259	FAD	P-O3P-PA-O2A
3	A	260	UFP	C5'-O5'-P-O1P
2	G	259	FAD	C4'-C5'-O5'-P
3	F	260	UFP	C5'-O5'-P-O1P
2	B	259	FAD	P-O3P-PA-O2A
2	H	259	FAD	P-O3P-PA-O2A
2	E	259	FAD	P-O3P-PA-O2A
2	F	259	FAD	C4'-C5'-O5'-P
2	E	259	FAD	C4'-C5'-O5'-P
2	A	259	FAD	C4'-C5'-O5'-P
2	B	259	FAD	O4'-C4'-C5'-O5'
2	A	259	FAD	O4'-C4'-C5'-O5'
2	F	259	FAD	P-O3P-PA-O1A
2	D	259	FAD	PA-O3P-P-O5'
2	H	259	FAD	C4'-C5'-O5'-P
2	H	259	FAD	C5B-O5B-PA-O3P
2	E	259	FAD	C5B-O5B-PA-O3P
2	D	259	FAD	C4'-C5'-O5'-P
2	H	259	FAD	P-O3P-PA-O1A
2	G	259	FAD	P-O3P-PA-O1A
2	A	259	FAD	P-O3P-PA-O1A
2	C	259	FAD	P-O3P-PA-O1A

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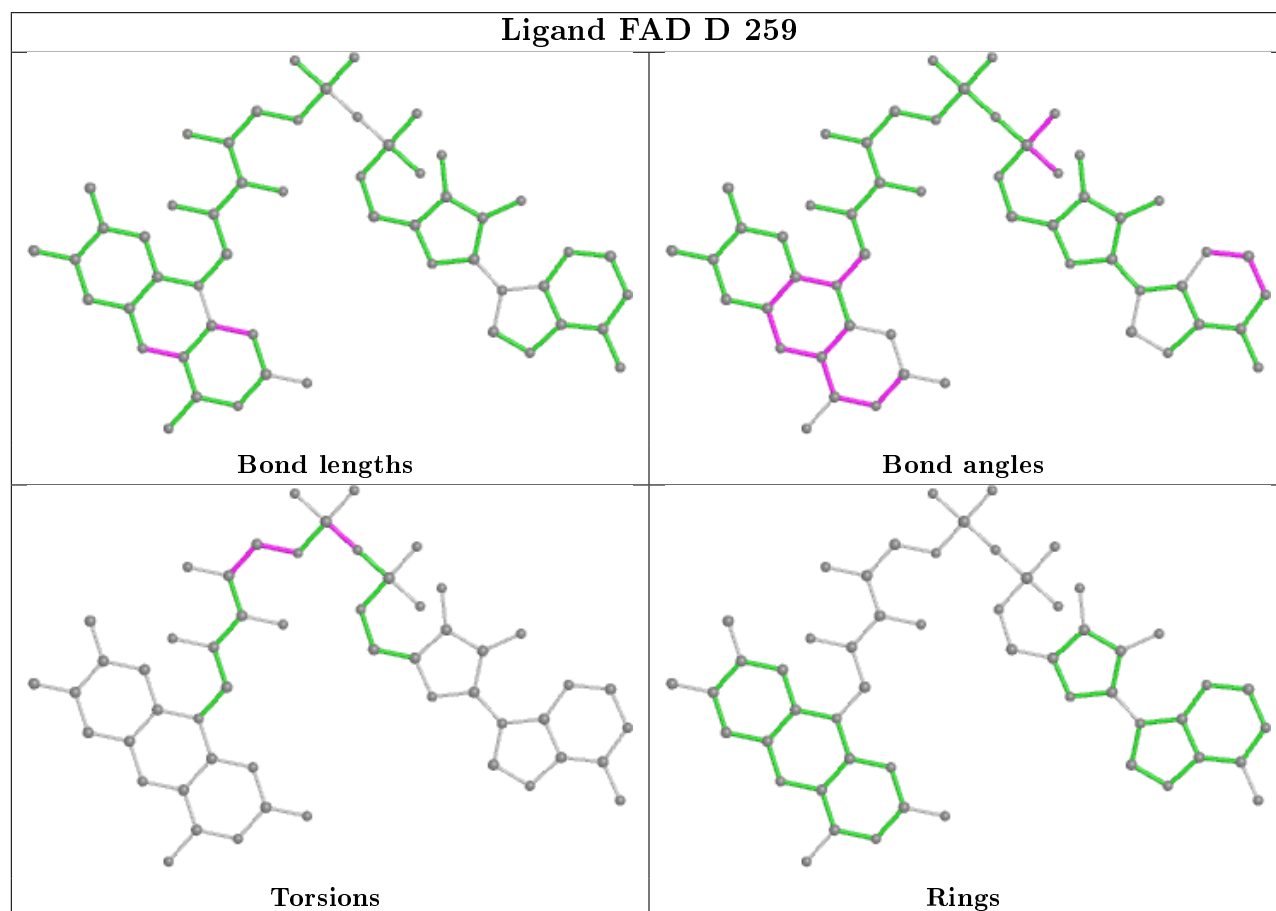
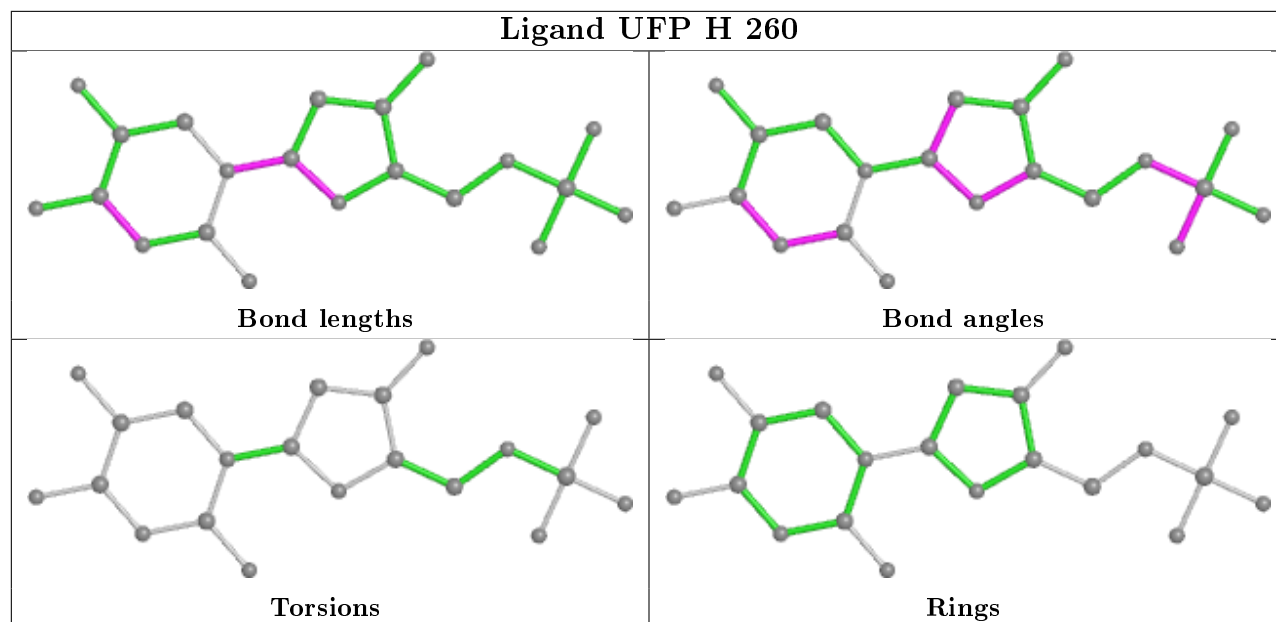
Mol	Chain	Res	Type	Atoms
2	C	259	FAD	C4'-C5'-O5'-P

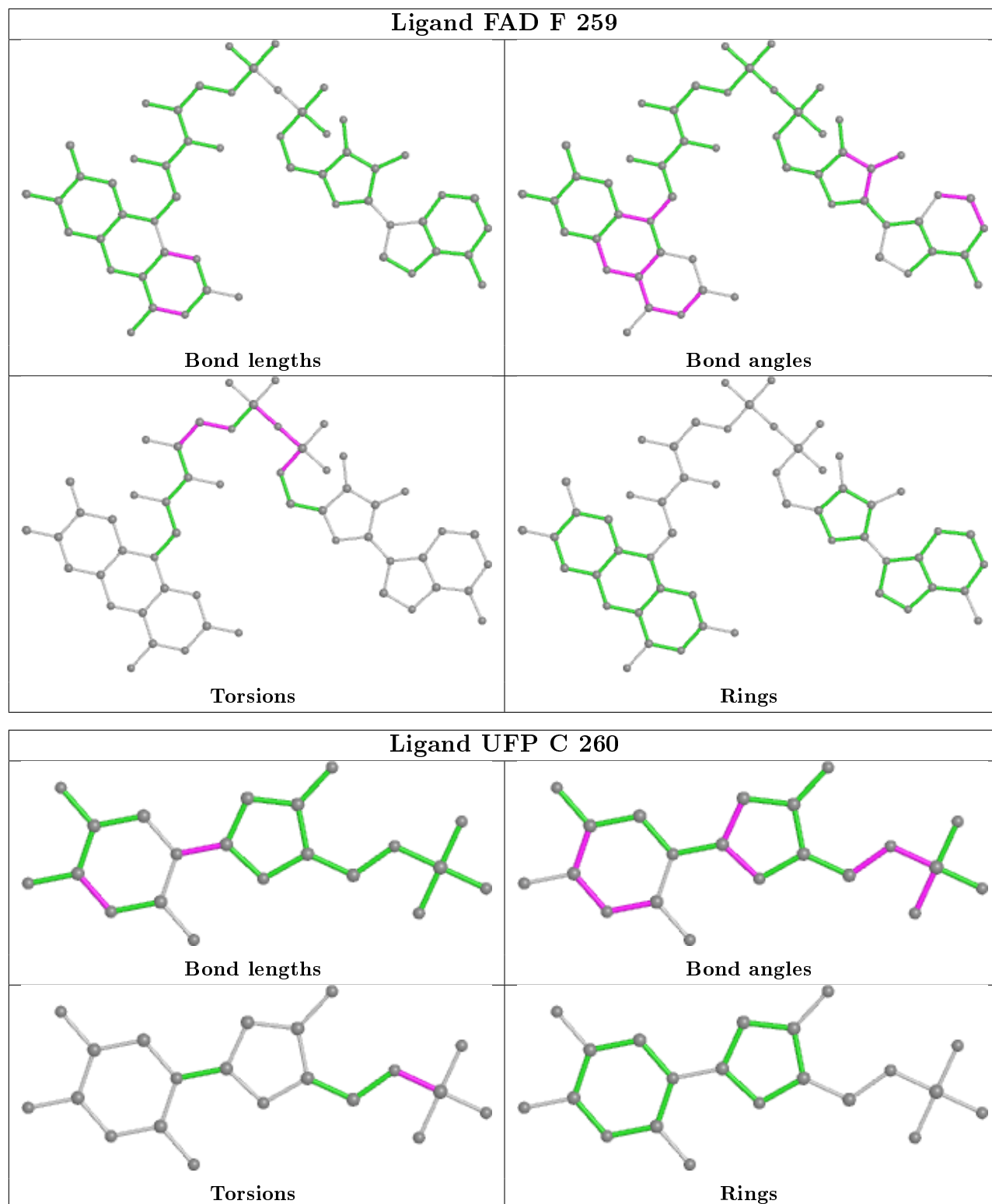
There are no ring outliers.

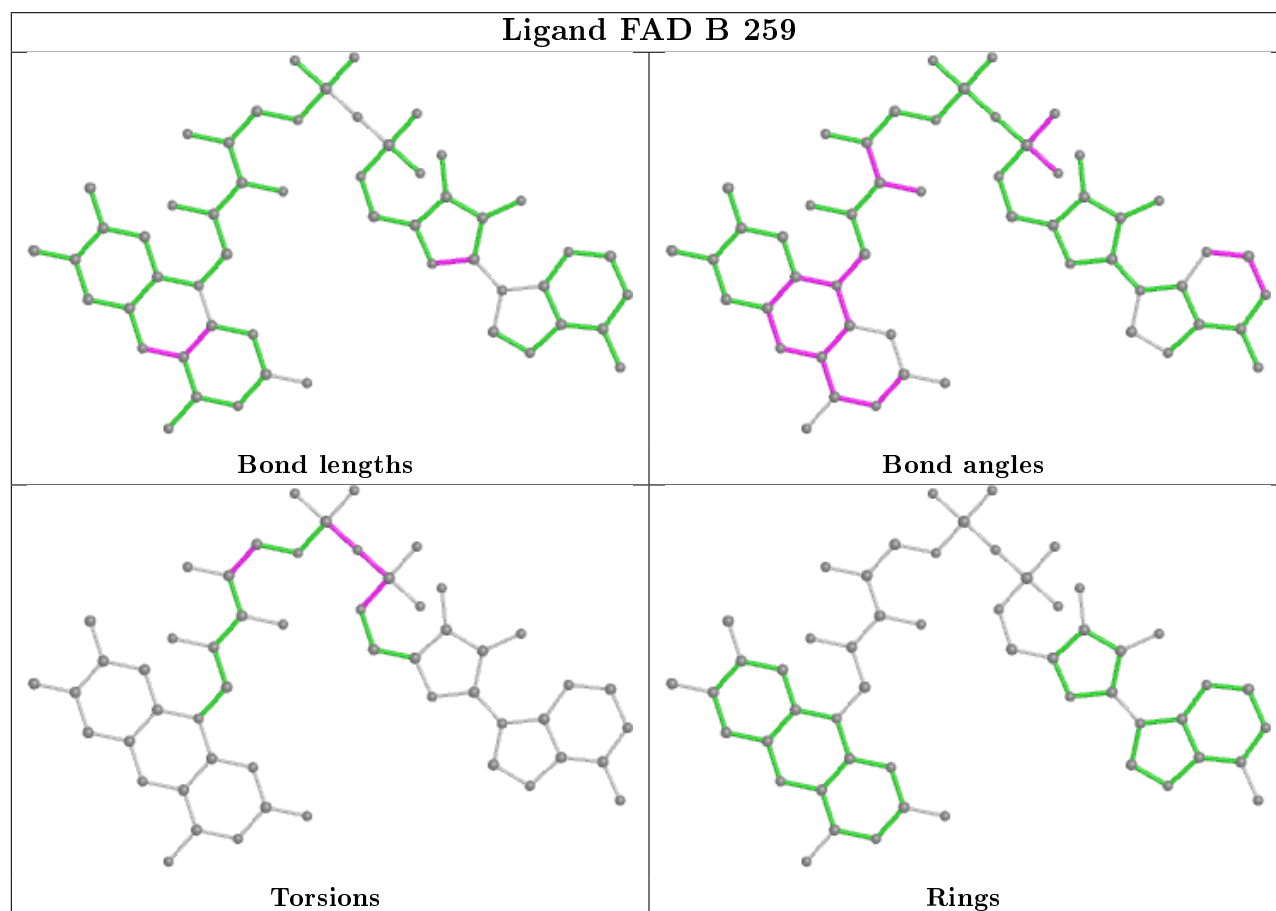
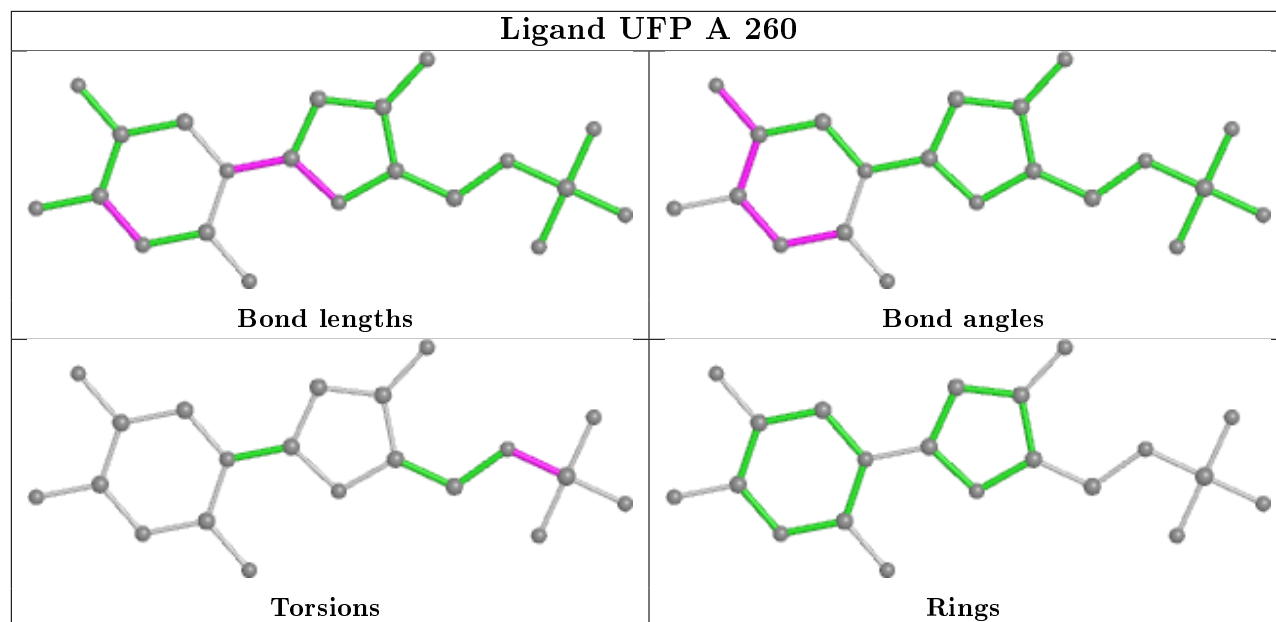
8 monomers are involved in 17 short contacts:

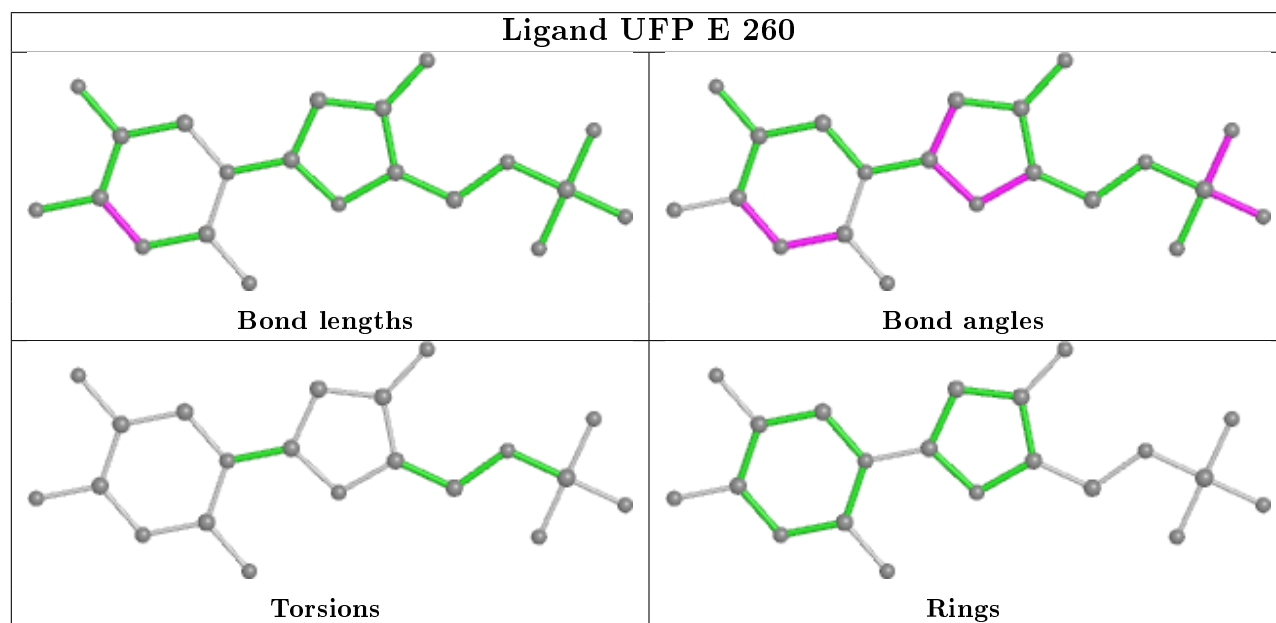
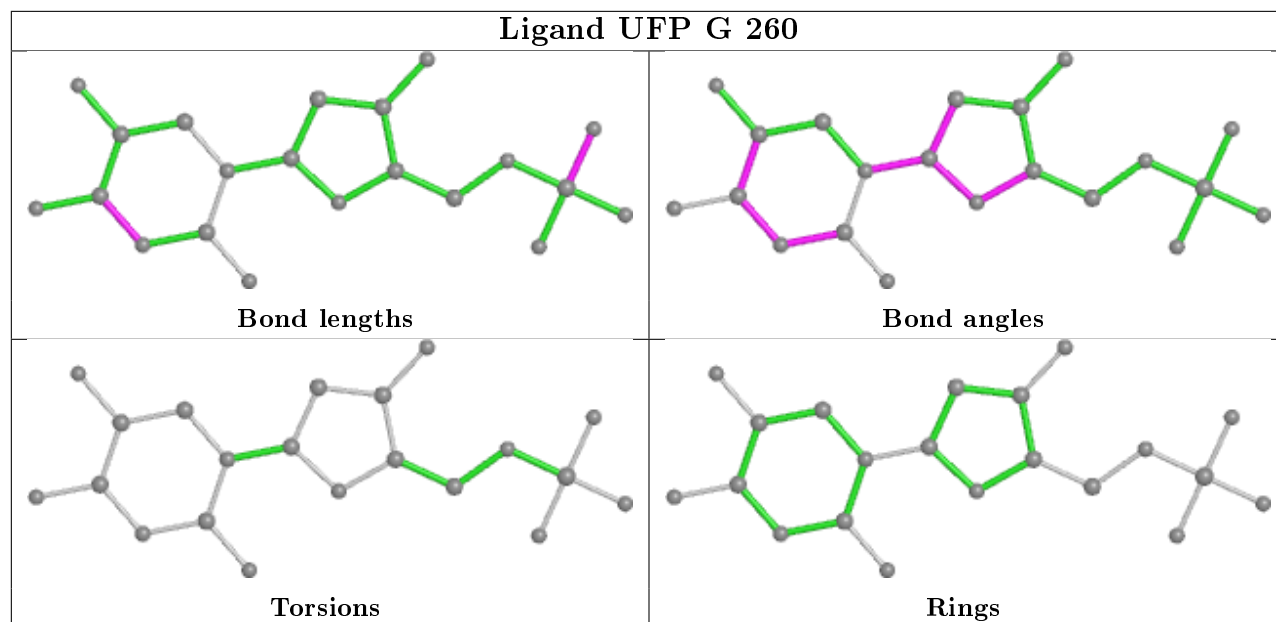
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	H	260	UFP	1	0
3	C	260	UFP	1	0
3	A	260	UFP	2	0
3	G	260	UFP	4	0
3	E	260	UFP	2	0
3	D	260	UFP	3	0
3	B	260	UFP	2	0
3	F	260	UFP	2	0

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.

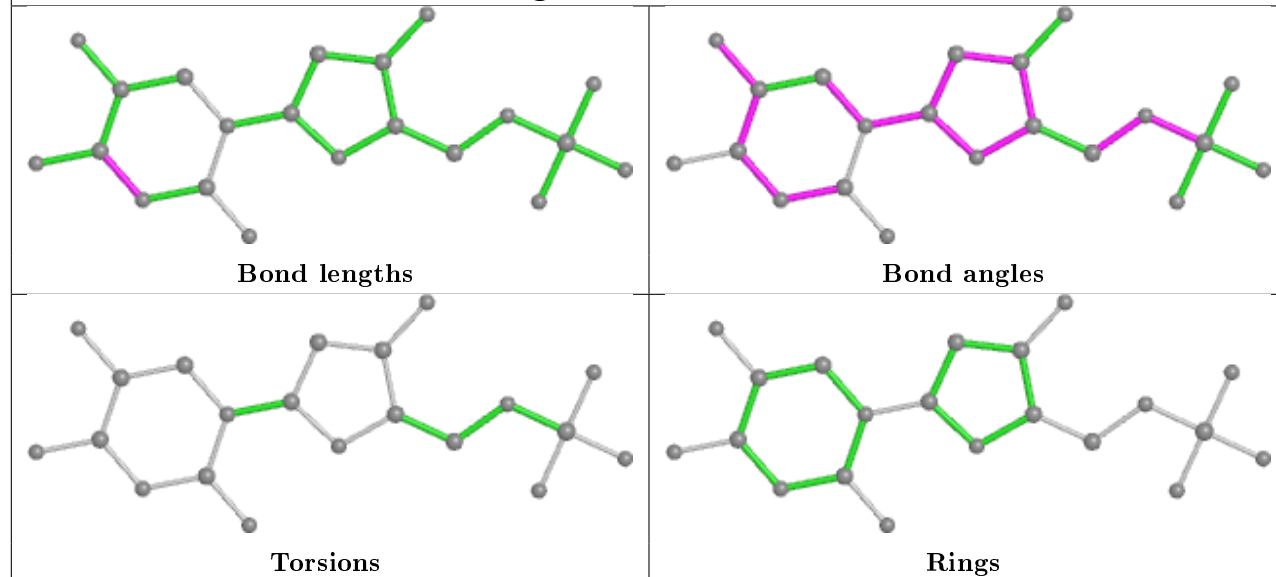




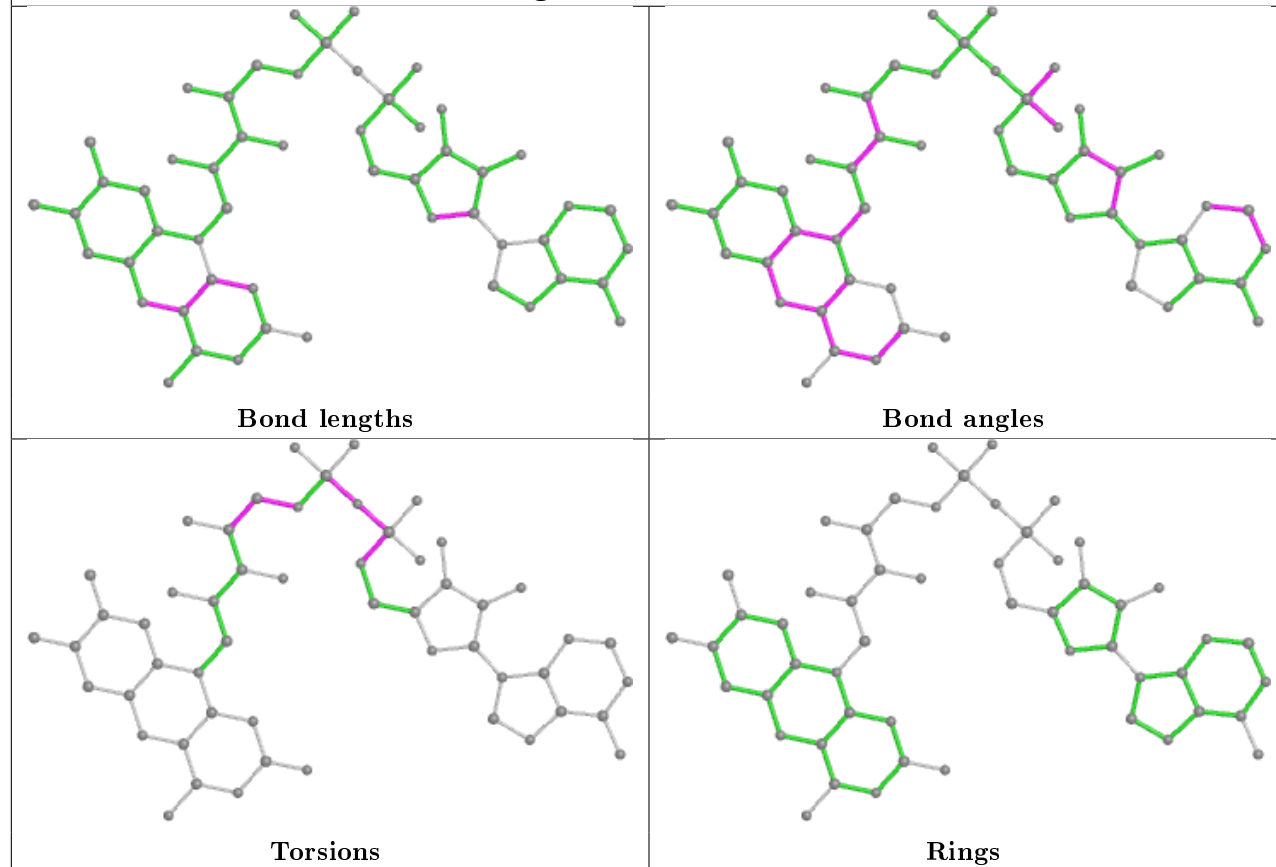




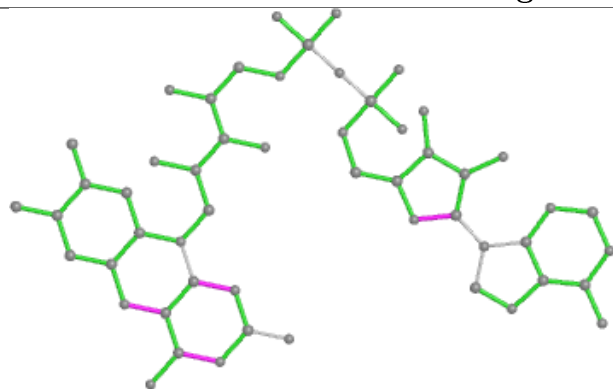
Ligand UFP D 260



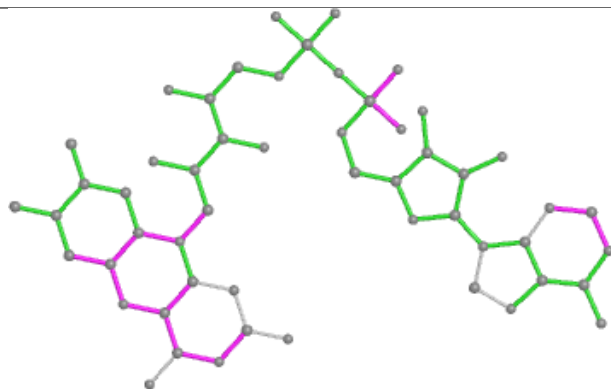
Ligand FAD H 259



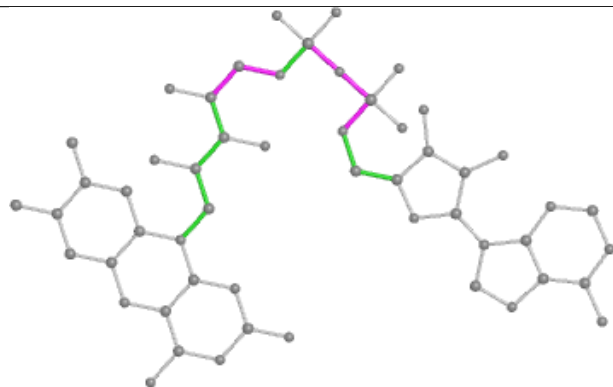
Ligand FAD E 259



Bond lengths



Bond angles

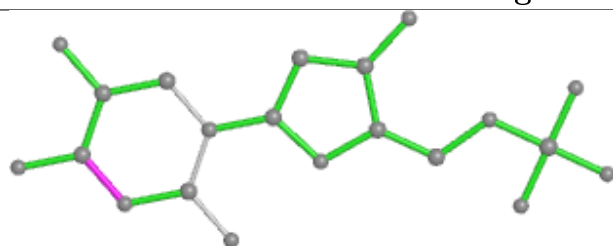


Torsions

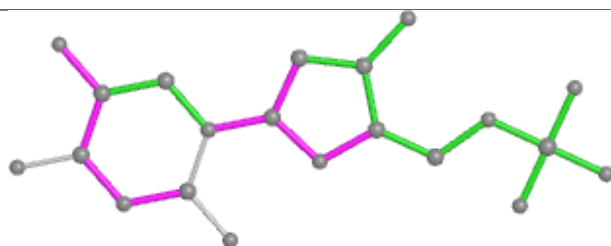


Rings

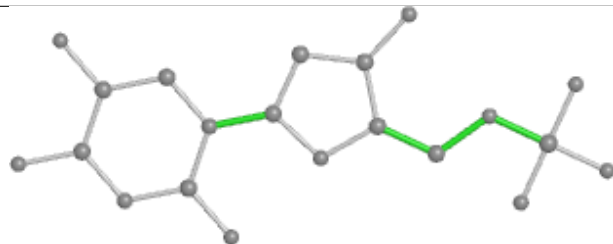
Ligand UFP B 260



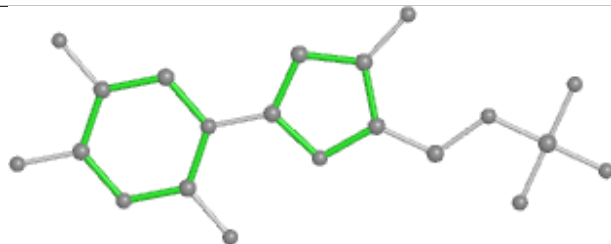
Bond lengths



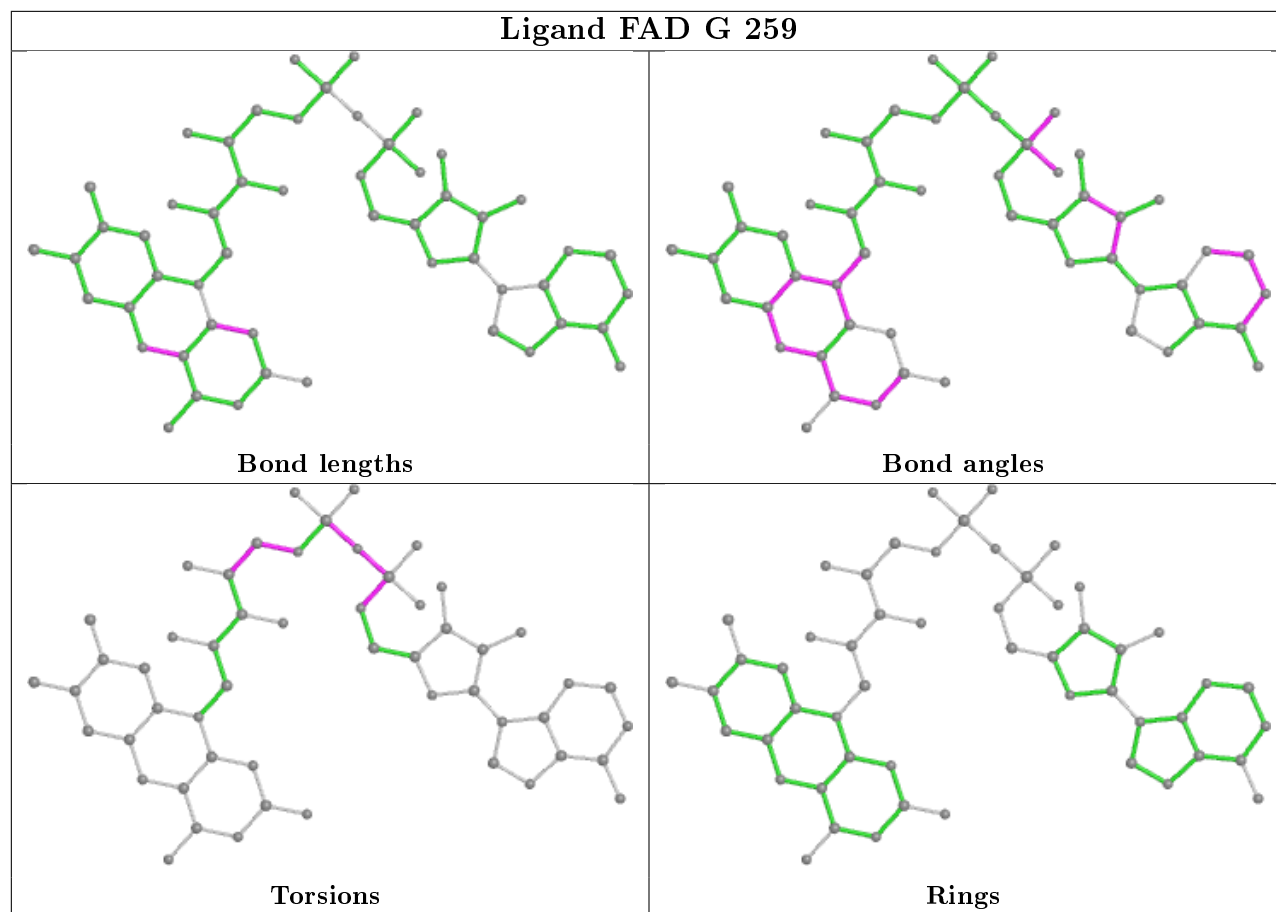
Bond angles



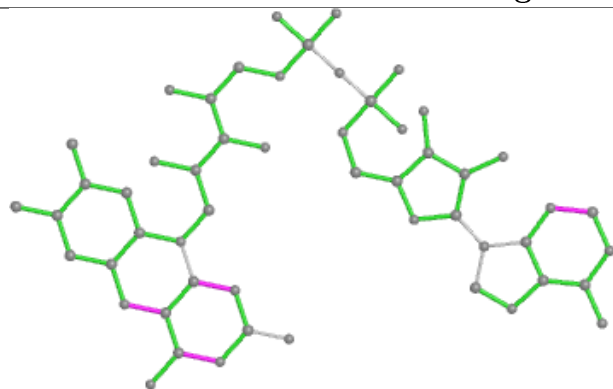
Torsions



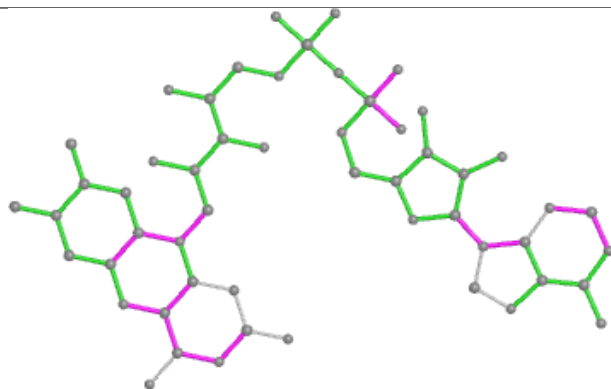
Rings



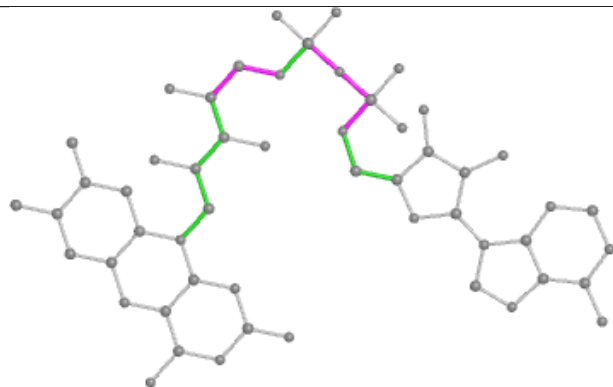
Ligand FAD A 259



Bond lengths



Bond angles

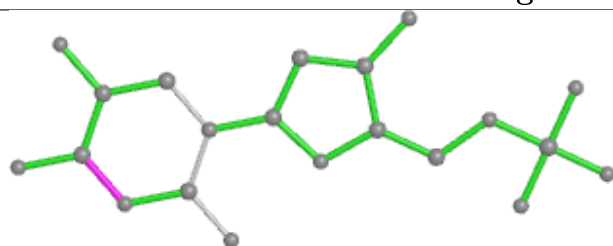


Torsions

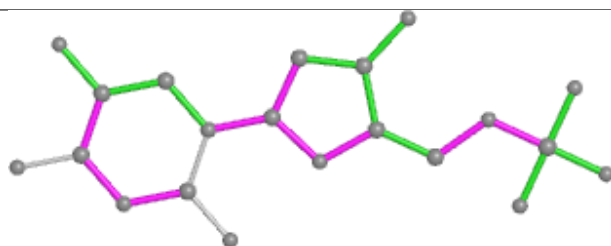


Rings

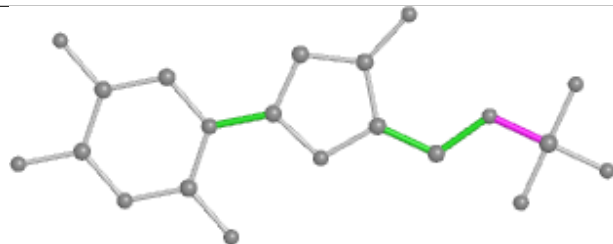
Ligand UFP F 260



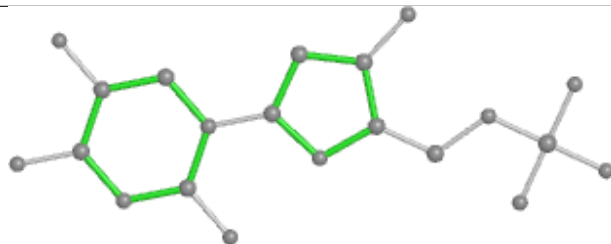
Bond lengths



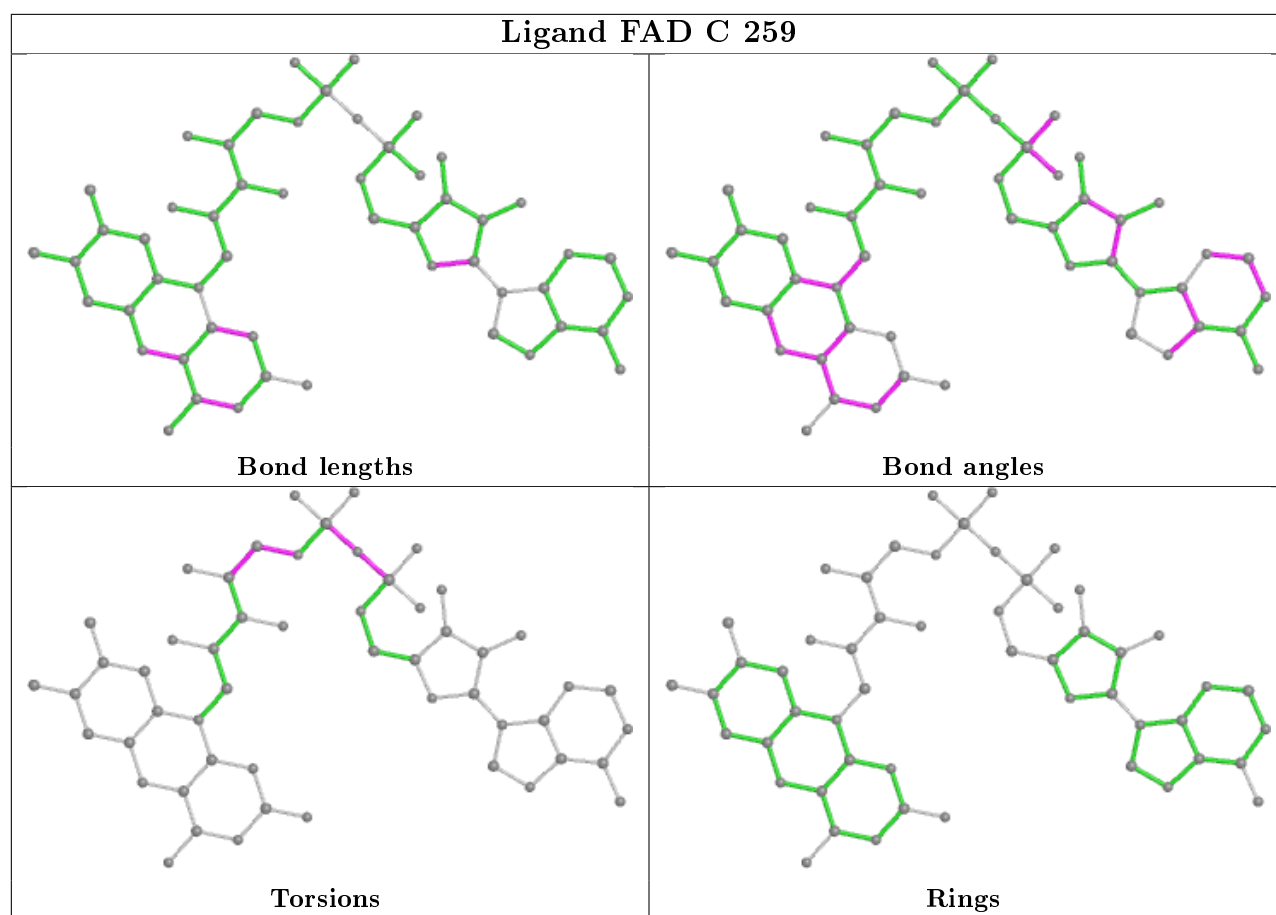
Bond angles



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 ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	246/258 (95%)	0.08	7 (2%) 53 56	9, 21, 36, 49	0
1	B	239/258 (92%)	0.16	10 (4%) 36 39	9, 19, 38, 49	1 (0%)
1	C	243/258 (94%)	0.13	15 (6%) 20 23	10, 23, 43, 51	0
1	D	242/258 (93%)	0.27	15 (6%) 20 23	10, 22, 41, 48	1 (0%)
1	E	244/258 (94%)	0.16	15 (6%) 21 24	12, 24, 44, 54	0
1	F	242/258 (93%)	0.43	22 (9%) 9 10	11, 26, 54, 63	1 (0%)
1	G	245/258 (94%)	0.41	20 (8%) 11 13	13, 27, 48, 53	1 (0%)
1	H	243/258 (94%)	-0.01	8 (3%) 46 49	10, 20, 40, 48	0
All	All	1944/2064 (94%)	0.20	112 (5%) 23 25	9, 23, 44, 63	4 (0%)

All (112) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	155	ALA	5.1
1	F	154	PHE	5.0
1	E	52	PRO	5.0
1	D	53	LYS	5.0
1	D	124	ASP	4.7
1	D	236	ALA	4.5
1	E	155	ALA	4.5
1	A	52	PRO	4.5
1	H	245	PRO	4.4
1	D	54	THR	4.4
1	A	247	ALA	4.1
1	D	155	ALA	4.1
1	C	235	LEU	3.8
1	C	238	GLY	3.7
1	F	161	ILE	3.6
1	G	236	ALA	3.5

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Mol	Chain	Res	Type	RSRZ
1	E	50	PRO	3.5
1	G	134	GLU	3.4
1	D	51	ASN	3.4
1	F	138	ALA	3.3
1	F	127	ASP	3.3
1	G	51	ASN	3.3
1	D	52	PRO	3.3
1	A	236	ALA	3.2
1	G	156	ASP	3.2
1	H	155	ALA	3.2
1	G	127	ASP	3.2
1	F	148	ALA	3.2
1	C	2	ALA	3.1
1	F	159	ASN	3.1
1	C	156	ASP	3.0
1	E	53	LYS	3.0
1	G	50	PRO	3.0
1	D	161	ILE	3.0
1	D	228	ALA	3.0
1	G	130	HIS	2.9
1	F	163	ARG	2.9
1	C	158	PRO	2.9
1	B	160	ALA	2.9
1	C	236	ALA	2.9
1	G	128	LEU	2.9
1	F	130	HIS	2.9
1	E	236	ALA	2.8
1	E	156	ASP	2.8
1	C	162	LEU	2.8
1	E	231	GLU	2.8
1	H	236	ALA	2.7
1	C	124	ASP	2.7
1	F	112	LYS	2.7
1	A	50	PRO	2.7
1	D	50	PRO	2.7
1	A	55	ALA	2.7
1	E	51	ASN	2.7
1	F	153	LYS	2.6
1	F	164	ARG	2.6
1	E	158	PRO	2.6
1	D	156	ASP	2.6
1	F	160	ALA	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	22	ASP	2.6
1	G	113	ASP	2.6
1	F	236	ALA	2.6
1	B	153	LYS	2.6
1	F	156	ASP	2.6
1	F	152	ALA	2.5
1	B	50	PRO	2.5
1	B	159	ASN	2.5
1	G	150	LEU	2.5
1	B	161	ILE	2.5
1	B	21	PRO	2.5
1	C	231	GLU	2.5
1	G	112	LYS	2.4
1	C	155	ALA	2.4
1	C	65	ILE	2.4
1	C	161	ILE	2.4
1	H	162	LEU	2.4
1	E	127	ASP	2.4
1	H	156	ASP	2.4
1	E	54	THR	2.3
1	C	50	PRO	2.3
1	F	126	ALA	2.3
1	F	147	LEU	2.3
1	G	155	ALA	2.3
1	B	124	ASP	2.3
1	E	235	LEU	2.3
1	A	51	ASN	2.3
1	A	2	ALA	2.3
1	D	164	ARG	2.3
1	D	152	ALA	2.2
1	E	221	ALA	2.2
1	G	162	LEU	2.2
1	B	164	ARG	2.2
1	C	239	THR	2.2
1	G	126	ALA	2.2
1	G	152	ALA	2.2
1	H	235	LEU	2.2
1	G	114	SER	2.2
1	G	108	TYR	2.2
1	D	67	VAL	2.2
1	E	22	ASP	2.2
1	H	244	SER	2.1

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Mol	Chain	Res	Type	RSRZ
1	H	243	THR	2.1
1	C	154	PHE	2.1
1	F	66	ASP	2.1
1	F	124	ASP	2.1
1	F	128	LEU	2.1
1	D	22	ASP	2.1
1	G	185	VAL	2.1
1	F	162	LEU	2.1
1	B	163	ARG	2.0
1	G	133	THR	2.0
1	E	161	ILE	2.0
1	G	94	ILE	2.0

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

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	GOL	C	261	6/6	0.81	0.17	32,34,35,35	0
2	FAD	F	259	53/53	0.94	0.13	21,27,31,32	0
2	FAD	D	259	53/53	0.95	0.11	14,22,25,29	0
2	FAD	G	259	53/53	0.95	0.13	18,25,28,30	0
2	FAD	H	259	53/53	0.96	0.10	16,20,26,28	0
2	FAD	E	259	53/53	0.96	0.09	14,22,26,29	0
2	FAD	B	259	53/53	0.96	0.10	14,18,24,26	0
2	FAD	A	259	53/53	0.96	0.11	14,19,25,27	0
2	FAD	C	259	53/53	0.96	0.10	14,20,25,27	0
3	UFP	G	260	21/21	0.97	0.09	17,21,24,27	0
3	UFP	A	260	21/21	0.97	0.09	11,15,19,20	0

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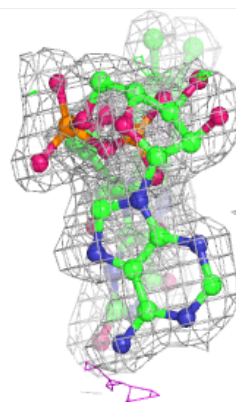
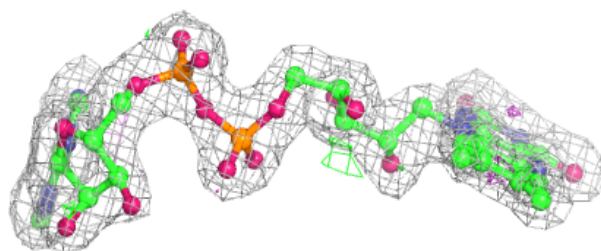
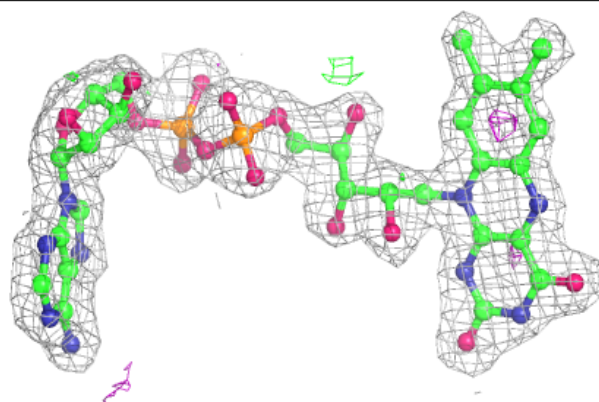
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	UFP	H	260	21/21	0.97	0.09	11,15,19,24	0
3	UFP	B	260	21/21	0.98	0.08	11,16,17,17	0
3	UFP	D	260	21/21	0.98	0.08	11,13,14,16	0
3	UFP	C	260	21/21	0.98	0.09	10,12,15,15	0
3	UFP	F	260	21/21	0.98	0.08	17,20,21,26	0
3	UFP	E	260	21/21	0.98	0.07	11,16,20,20	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.

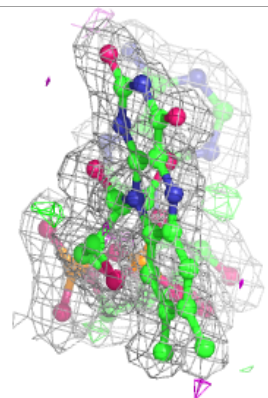
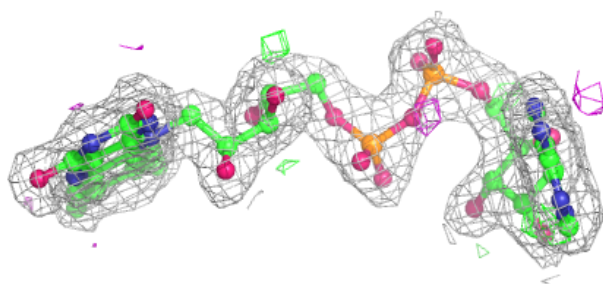
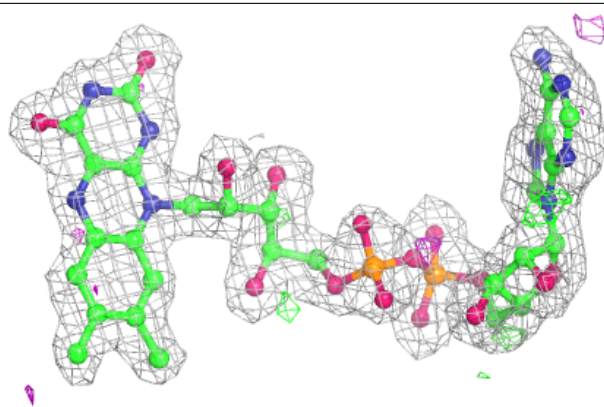
Electron density around FAD F 259:

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

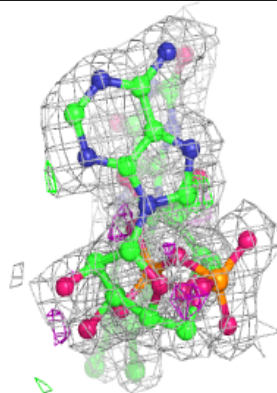
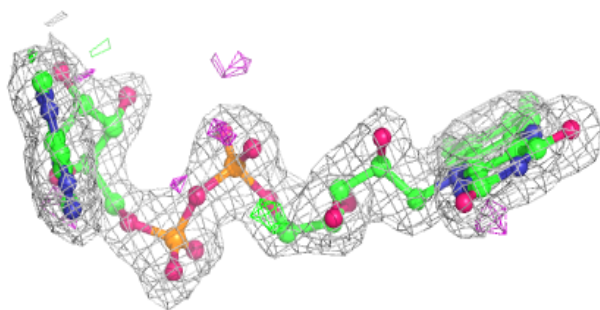
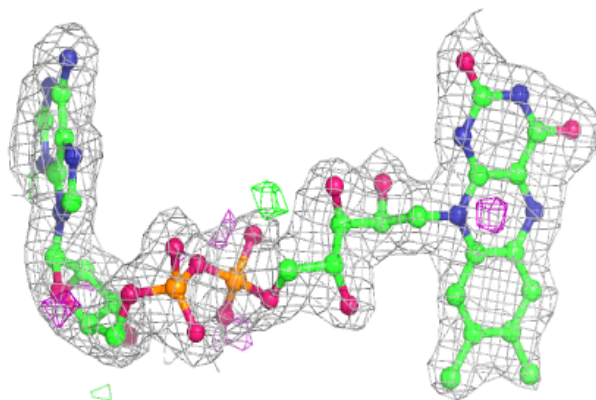


Electron density around FAD D 259:

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

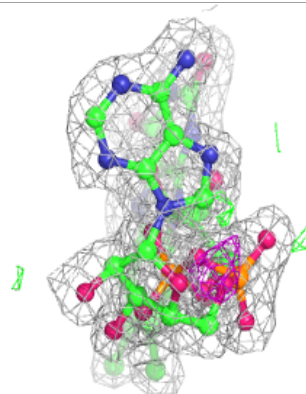
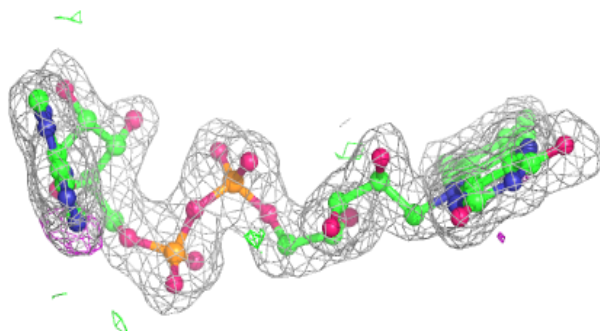
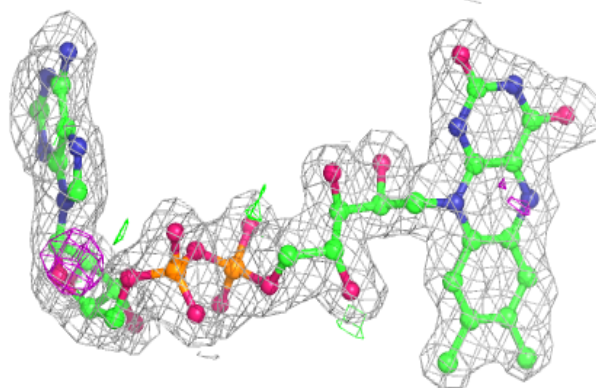
**Electron density around FAD G 259:**

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

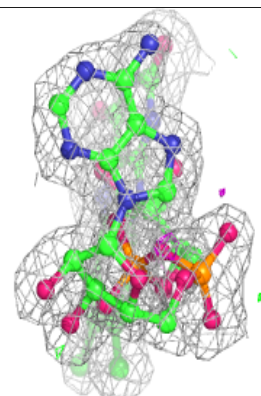
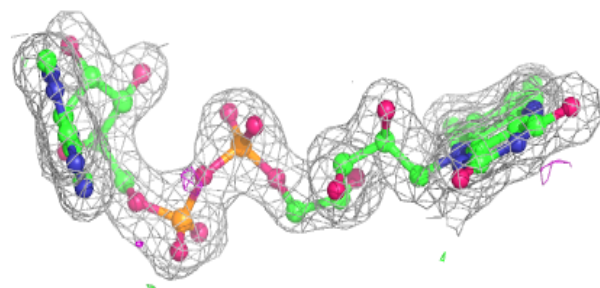
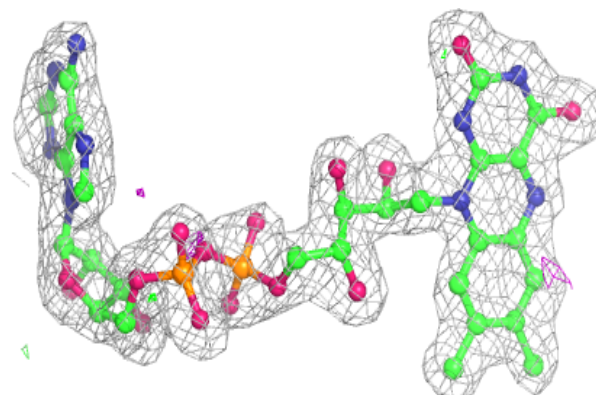


Electron density around FAD H 259:

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

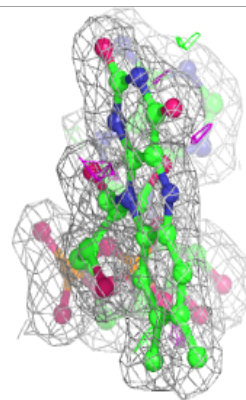
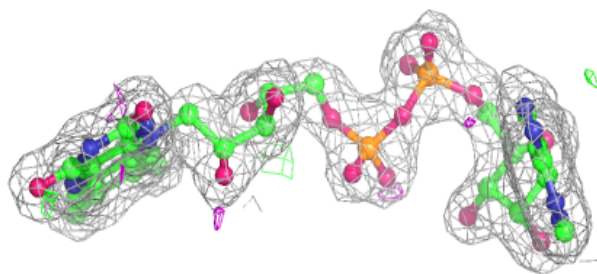
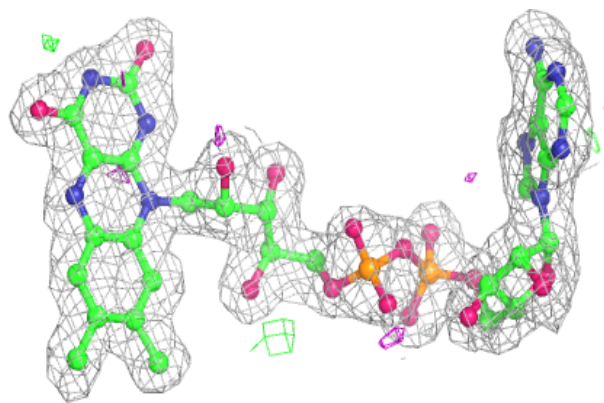
**Electron density around FAD E 259:**

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

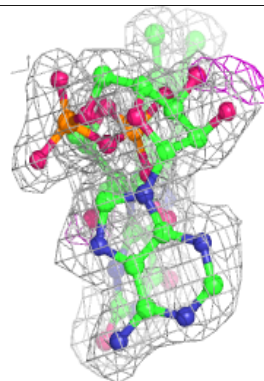
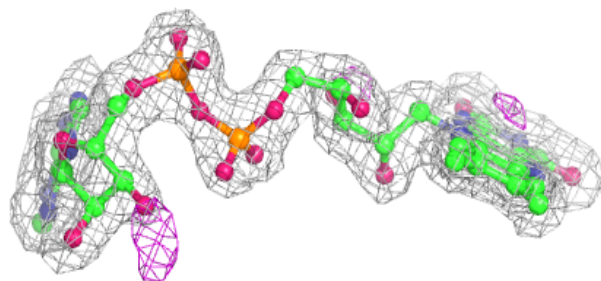
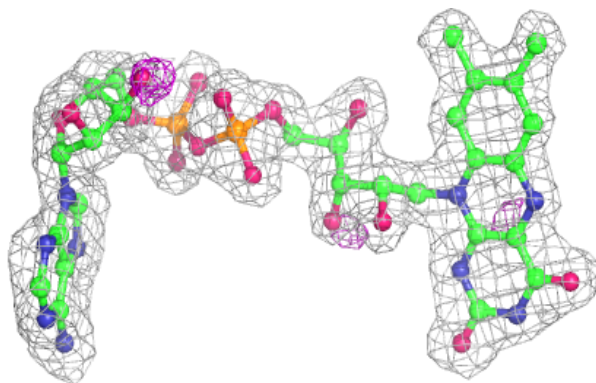


Electron density around FAD B 259:

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

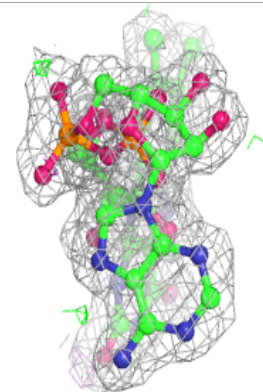
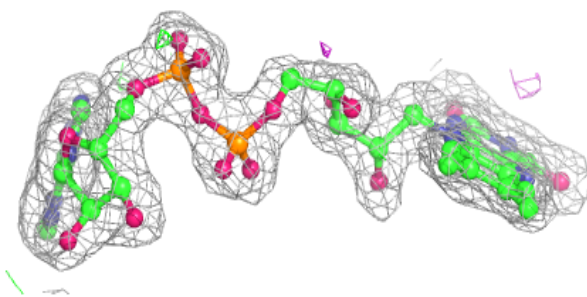
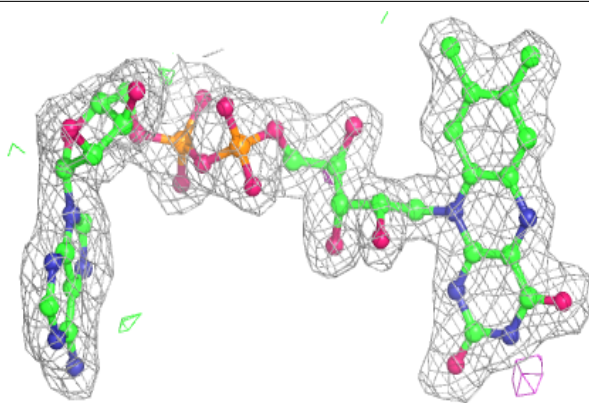
**Electron density around FAD A 259:**

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



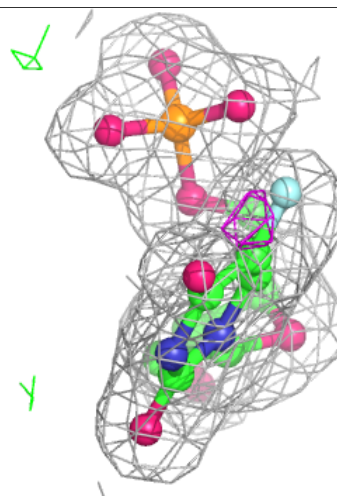
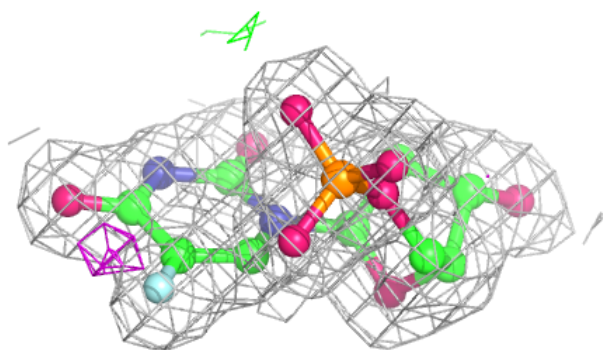
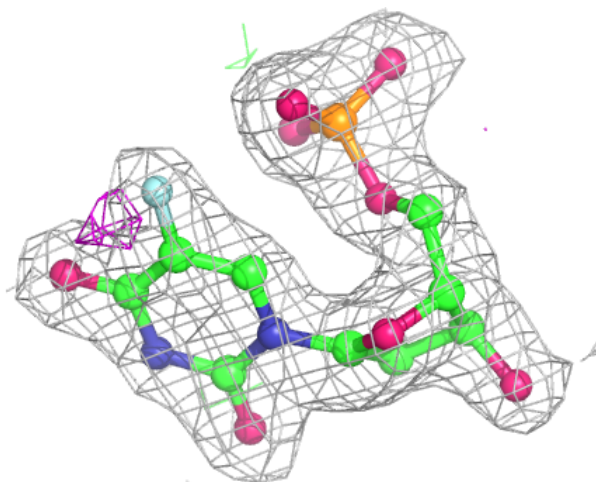
Electron density around FAD C 259:

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



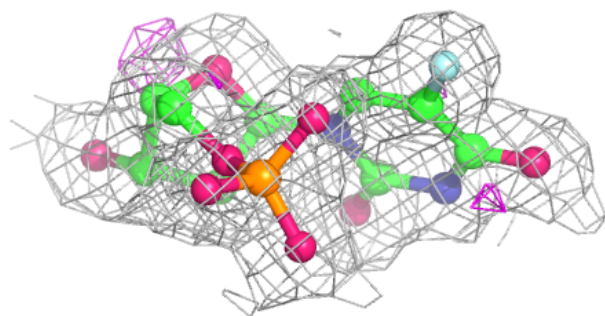
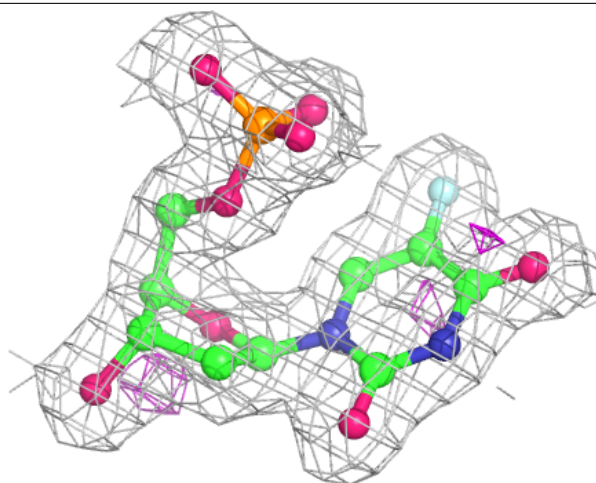
Electron density around UFP G 260:

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



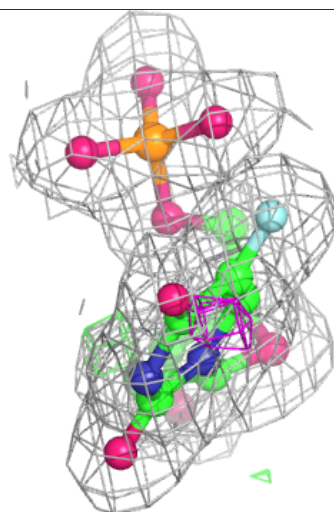
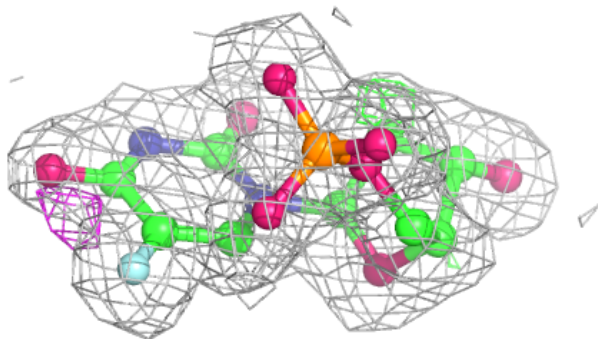
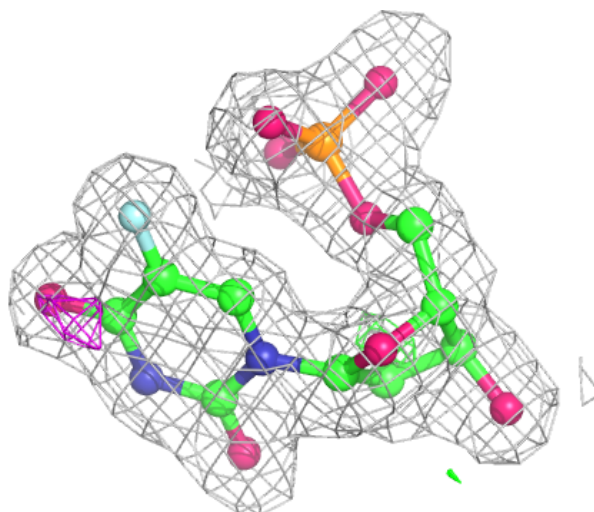
Electron density around UFP A 260:

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



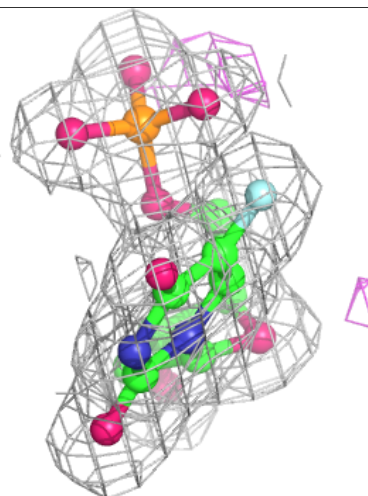
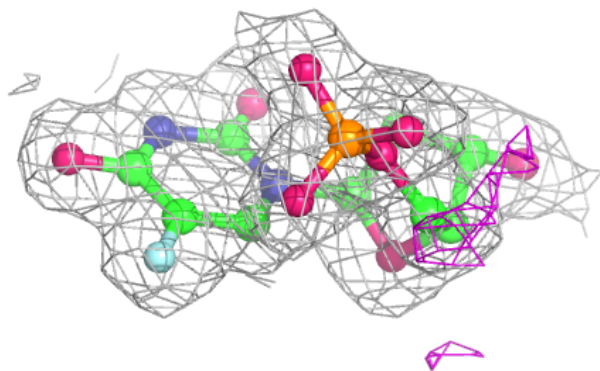
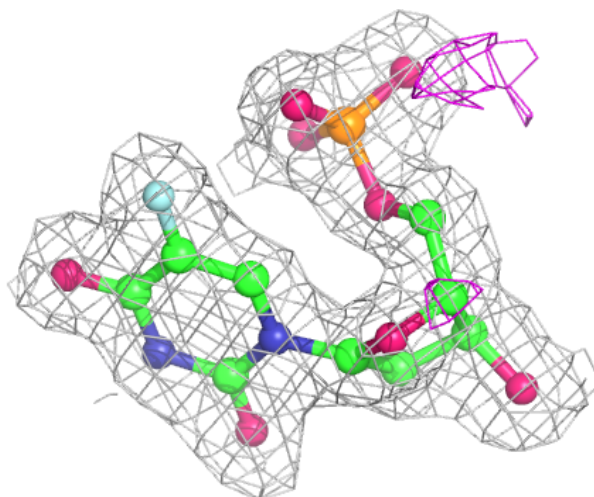
Electron density around UFP H 260:

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



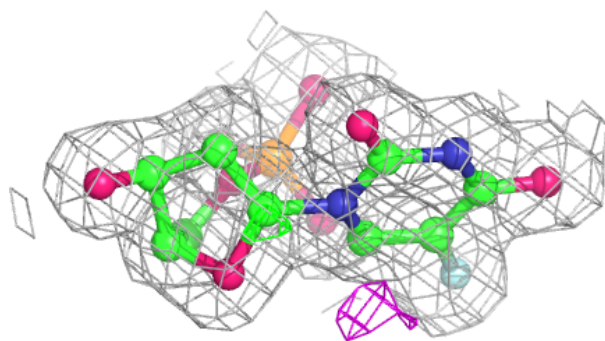
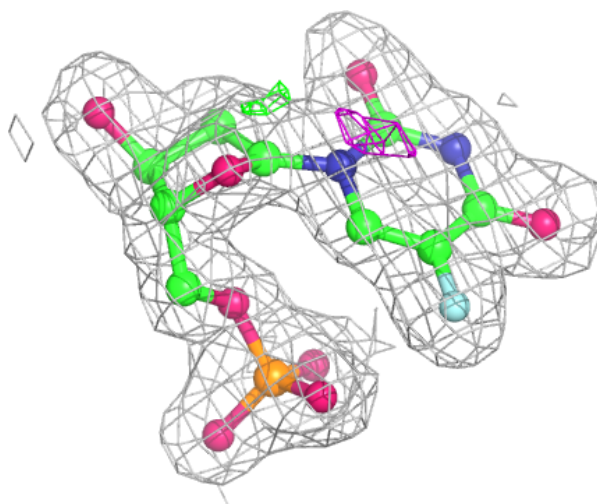
Electron density around UFP B 260:

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



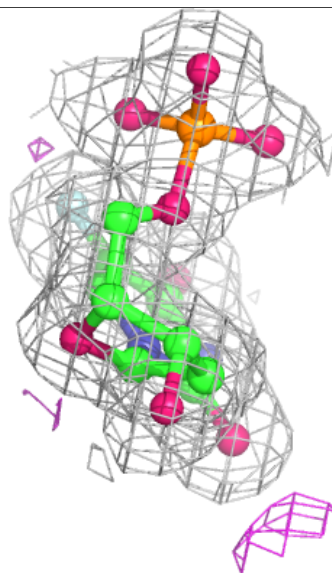
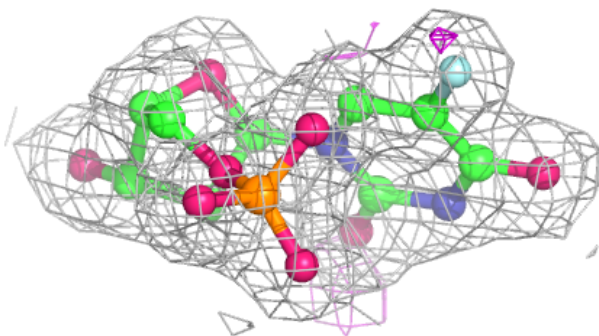
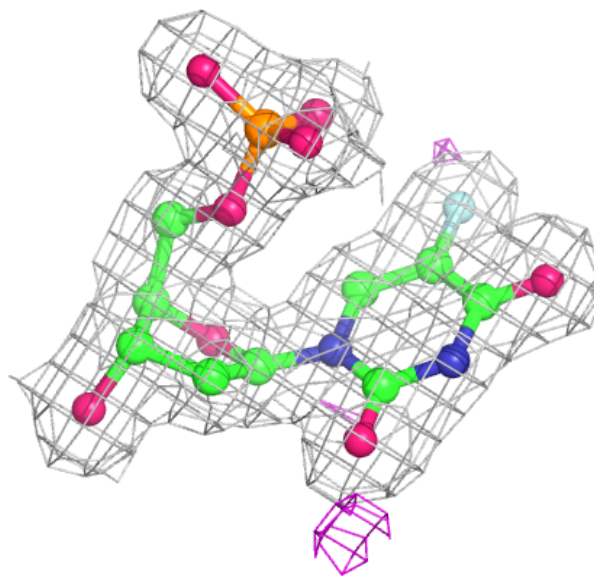
Electron density around UFP D 260:

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



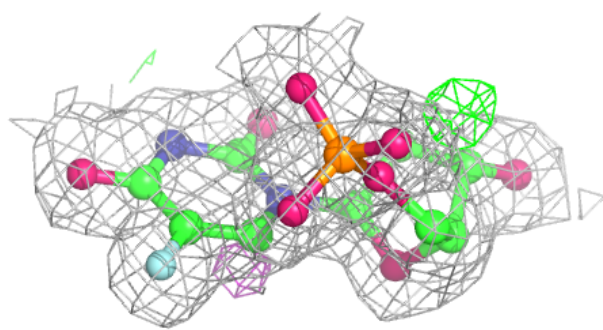
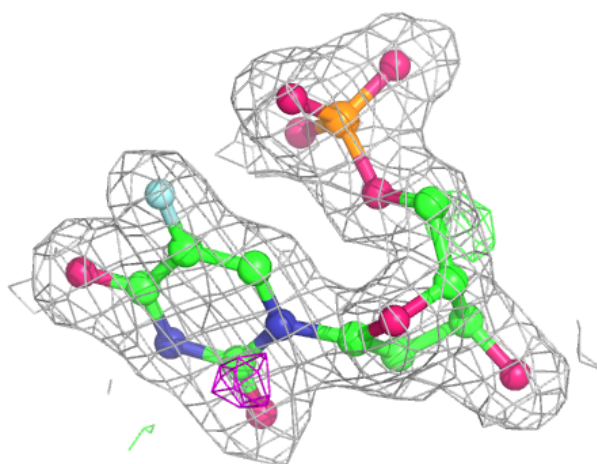
Electron density around UFP C 260:

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



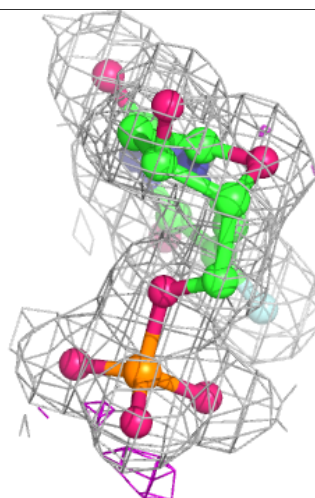
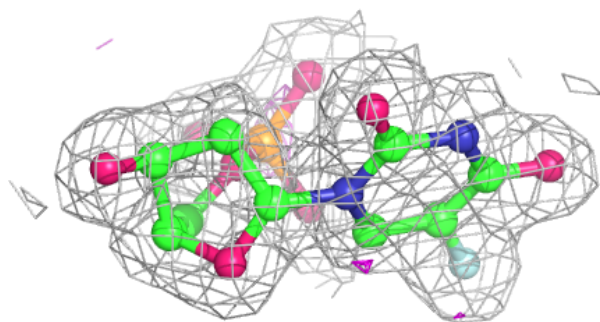
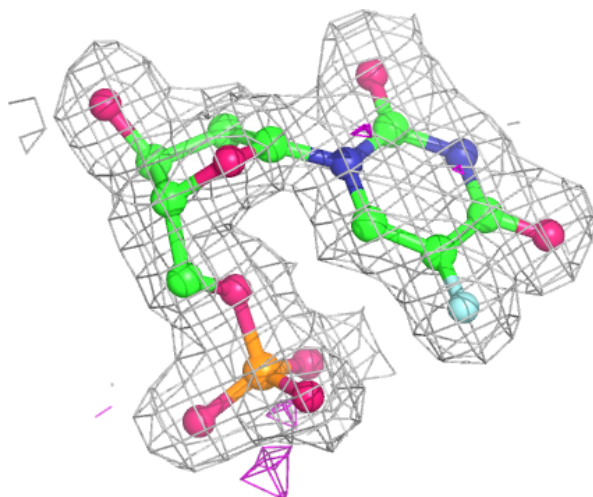
Electron density around UFP F 260:

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



Electron density around UFP E 260:

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



6.5 Other polymers ⓘ

There are no such residues in this entry.