



wwPDB X-ray Structure Validation Summary Report ⓘ

Aug 6, 2024 – 02:19 AM EDT

PDB ID : 3JSX
Title : X-ray Crystal structure of NAD(P)H: Quinone Oxidoreductase-1 (NQO1)
bound to the coumarin-based inhibitor AS1
Authors : Dunstan, M.S.; Levy, C.; Leys, D.
Deposited on : 2009-09-11
Resolution : 2.45 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.37.1
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) : 2.37.1

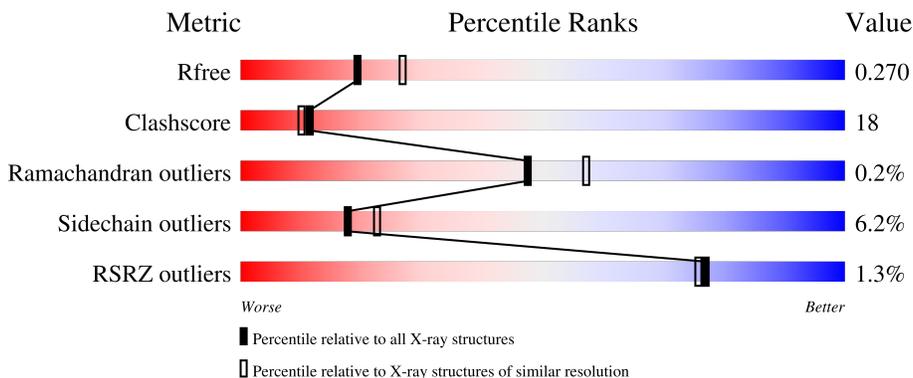
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.45 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	1544 (2.48-2.44)
Clashscore	141614	1613 (2.48-2.44)
Ramachandran outliers	138981	1598 (2.48-2.44)
Sidechain outliers	138945	1598 (2.48-2.44)
RSRZ outliers	127900	1523 (2.48-2.44)

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

Mol	Chain	Length	Quality of chain
1	A	273	 62% 32% ..
1	B	273	 70% 27% ..
1	C	273	 66% 29% ..
1	D	273	 69% 27% ..
1	E	273	 71% 24% ..

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Mol	Chain	Length	Quality of chain	
1	F	273	1%	
1	G	273	2%	
1	H	273	3%	

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	CC2	B	547	-	-	X	-

2 Entry composition

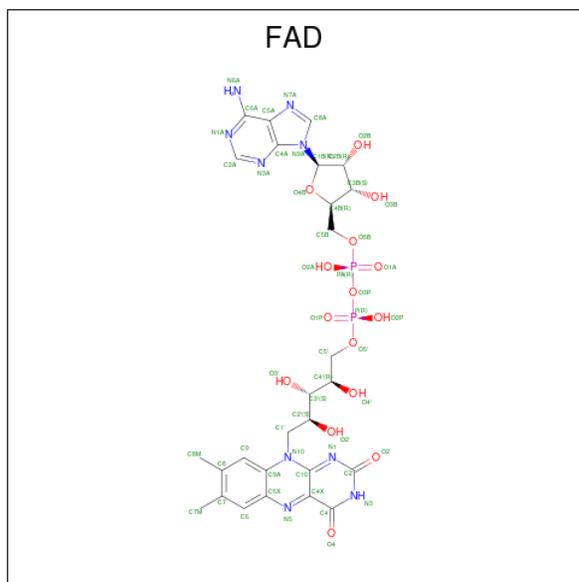
There are 4 unique types of molecules in this entry. The entry contains 18234 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 NAD(P)H dehydrogenase [quinone] 1.

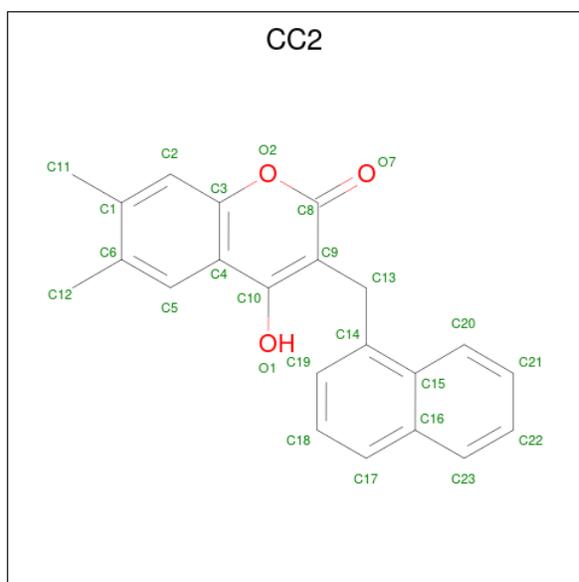
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	270	Total	C	N	O	S	0	0	0
			2147	1397	358	385	7			
1	B	270	Total	C	N	O	S	0	0	0
			2147	1397	358	385	7			
1	C	270	Total	C	N	O	S	0	0	0
			2147	1397	358	385	7			
1	D	270	Total	C	N	O	S	0	0	0
			2147	1397	358	385	7			
1	E	270	Total	C	N	O	S	0	0	0
			2147	1397	358	385	7			
1	F	270	Total	C	N	O	S	0	0	0
			2147	1397	358	385	7			
1	G	270	Total	C	N	O	S	0	0	0
			2147	1397	358	385	7			
1	H	270	Total	C	N	O	S	0	0	0
			2147	1397	358	385	7			

- Molecule 2 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula: $C_{27}H_{33}N_9O_{15}P_2$).



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

- Molecule 3 is 4-hydroxy-6,7-dimethyl-3-(naphthalen-1-ylmethyl)-2H-chromen-2-one (three-letter code: CC2) (formula: C₂₂H₁₈O₃).



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	C O	0	0
			25	22 3		
3	B	1	Total	C O	0	0
			25	22 3		
3	C	1	Total	C O	0	0
			25	22 3		
3	D	1	Total	C O	0	0
			25	22 3		
3	E	1	Total	C O	0	0
			25	22 3		
3	F	1	Total	C O	0	0
			25	22 3		
3	G	1	Total	C O	0	0
			25	22 3		
3	H	1	Total	C O	0	0
			25	22 3		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	66	Total	O	0	0
			66	66		
4	B	66	Total	O	0	0
			66	66		
4	C	70	Total	O	0	0
			70	70		
4	D	59	Total	O	0	0
			59	59		

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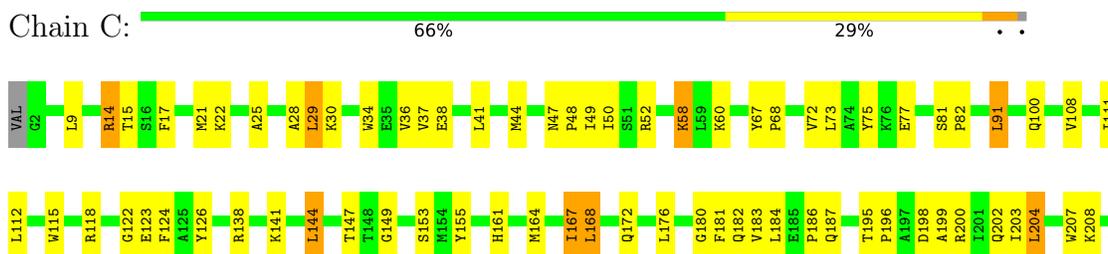
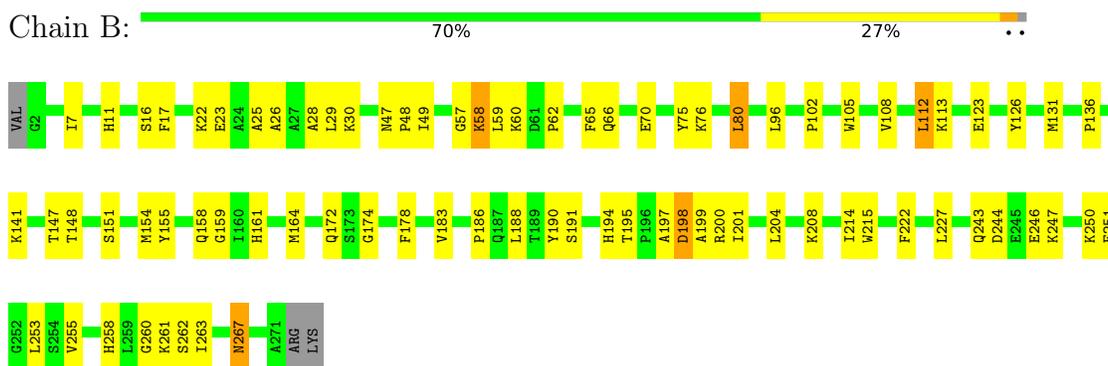
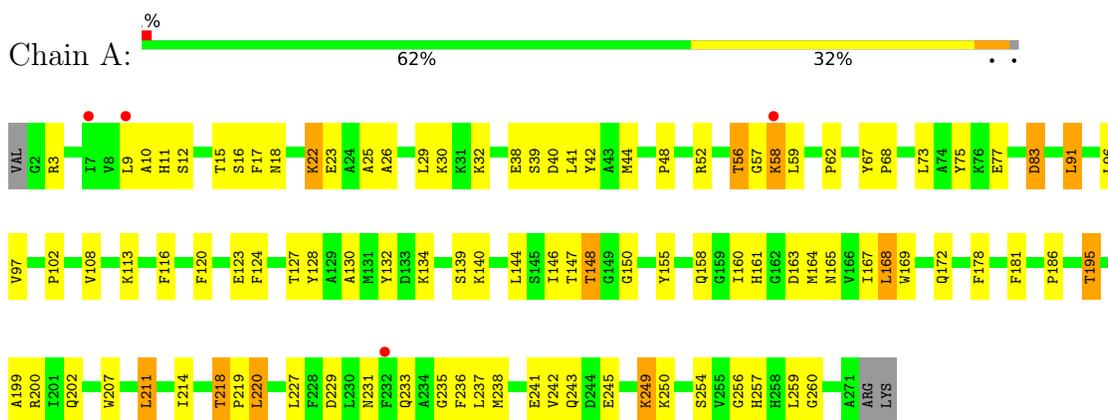
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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	E	37	Total O 37 37	0	0
4	F	41	Total O 41 41	0	0
4	G	54	Total O 54 54	0	0
4	H	41	Total O 41 41	0	0

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: NAD(P)H dehydrogenase [quinone] 1

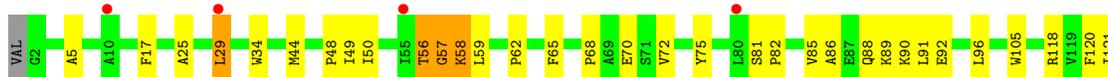




• Molecule 1: NAD(P)H dehydrogenase [quinone] 1



• Molecule 1: NAD(P)H dehydrogenase [quinone] 1

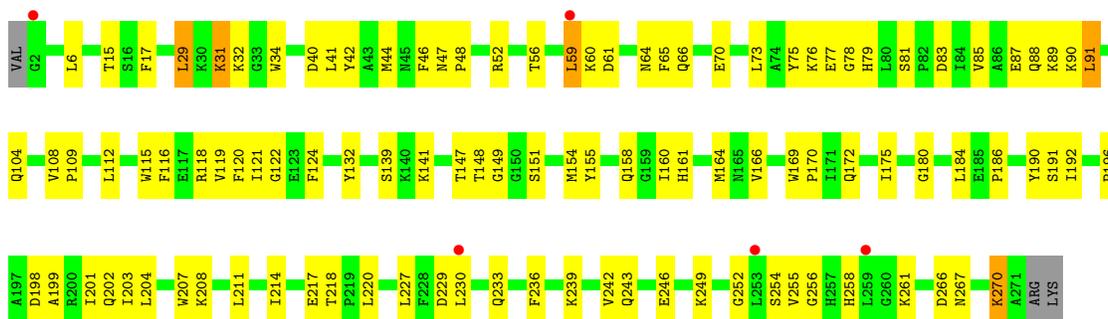


• Molecule 1: NAD(P)H dehydrogenase [quinone] 1

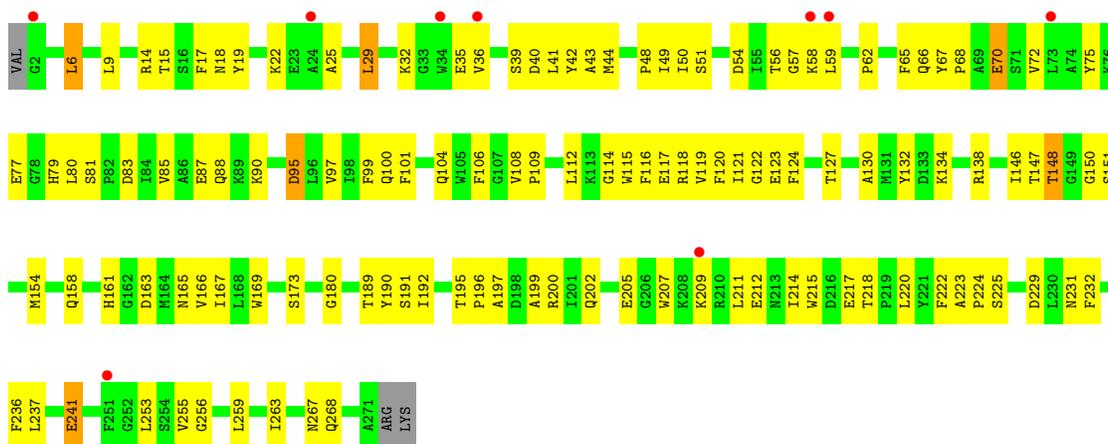


• Molecule 1: NAD(P)H dehydrogenase [quinone] 1





● Molecule 1: NAD(P)H dehydrogenase [quinone] 1



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	70.04Å 209.94Å 102.08Å 90.00° 109.93° 90.00°	Depositor
Resolution (Å)	47.95 – 2.45 47.95 – 2.45	Depositor EDS
% Data completeness (in resolution range)	99.7 (47.95-2.45) 99.1 (47.95-2.45)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.37 (at 2.45Å)	Xtrriage
Refinement program	PHENIX (phenix.refine)	Depositor
R, R_{free}	0.206 , 0.269 0.206 , 0.270	Depositor DCC
R_{free} test set	5065 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	53.9	Xtrriage
Anisotropy	0.646	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 46.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtrriage
Estimated twinning fraction	0.199 for h,-k,-h-l	Xtrriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	18234	wwPDB-VP
Average B, all atoms (Å ²)	64.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 10.51% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: FAD, CC2

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	0.43	0/2205	0.59	0/2980
1	B	0.46	0/2205	0.58	0/2980
1	C	0.45	0/2205	0.58	0/2980
1	D	0.46	0/2205	0.58	0/2980
1	E	0.40	0/2205	0.54	0/2980
1	F	0.41	0/2205	0.57	0/2980
1	G	0.41	0/2205	0.56	0/2980
1	H	0.40	0/2205	0.55	0/2980
All	All	0.43	0/17640	0.57	0/23840

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2147	0	2141	90	0
1	B	2147	0	2141	79	0
1	C	2147	0	2141	80	0
1	D	2147	0	2141	67	0
1	E	2147	0	2141	70	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	F	2147	0	2141	89	0
1	G	2147	0	2141	75	0
1	H	2147	0	2141	99	0
2	A	53	0	31	7	0
2	B	53	0	31	10	0
2	C	53	0	31	6	0
2	D	53	0	31	6	0
2	E	53	0	31	4	0
2	F	53	0	31	4	0
2	G	53	0	31	2	0
2	H	53	0	31	9	0
3	A	25	0	17	1	0
3	B	25	0	18	10	0
3	C	25	0	17	4	0
3	D	25	0	17	2	0
3	E	25	0	18	3	0
3	F	25	0	18	3	0
3	G	25	0	18	2	0
3	H	25	0	17	3	0
4	A	66	0	0	2	0
4	B	66	0	0	0	0
4	C	70	0	0	2	0
4	D	59	0	0	4	0
4	E	37	0	0	4	0
4	F	41	0	0	2	0
4	G	54	0	0	2	0
4	H	41	0	0	2	0
All	All	18234	0	17516	626	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 18.

The worst 5 of 626 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:E:228:PHE:HA	4:E:361:HOH:O	1.37	1.22
1:F:151:SER:H	1:F:154:MET:HE3	1.08	1.12
1:F:17:PHE:HB2	2:F:601:FAD:H52A	1.13	1.11
1:B:17:PHE:HB2	2:B:601:FAD:H52A	1.13	1.07
1:E:238:MET:HA	4:E:361:HOH:O	1.54	1.05

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	268/273 (98%)	247 (92%)	20 (8%)	1 (0%)	34	41
1	B	268/273 (98%)	242 (90%)	25 (9%)	1 (0%)	34	41
1	C	268/273 (98%)	246 (92%)	22 (8%)	0	100	100
1	D	268/273 (98%)	249 (93%)	19 (7%)	0	100	100
1	E	268/273 (98%)	243 (91%)	23 (9%)	2 (1%)	22	25
1	F	268/273 (98%)	249 (93%)	18 (7%)	1 (0%)	34	41
1	G	268/273 (98%)	240 (90%)	28 (10%)	0	100	100
1	H	268/273 (98%)	246 (92%)	22 (8%)	0	100	100
All	All	2144/2184 (98%)	1962 (92%)	177 (8%)	5 (0%)	47	57

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	83	ASP
1	E	56	THR
1	E	57	GLY
1	B	57	GLY
1	F	197	ALA

5.3.2 Protein sidechains [i](#)

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	224/227 (99%)	205 (92%)	19 (8%)	10	12
1	B	224/227 (99%)	216 (96%)	8 (4%)	35	46
1	C	224/227 (99%)	207 (92%)	17 (8%)	13	15
1	D	224/227 (99%)	212 (95%)	12 (5%)	22	28
1	E	224/227 (99%)	213 (95%)	11 (5%)	25	32
1	F	224/227 (99%)	206 (92%)	18 (8%)	12	14
1	G	224/227 (99%)	211 (94%)	13 (6%)	20	25
1	H	224/227 (99%)	210 (94%)	14 (6%)	18	22
All	All	1792/1816 (99%)	1680 (94%)	112 (6%)	18	22

5 of 112 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	E	58	LYS
1	H	231	ASN
1	F	44	MET
1	H	218	THR
1	H	15	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 62 such sidechains are listed below:

Mol	Chain	Res	Type
1	D	268	GLN
1	H	79	HIS
1	E	243	GLN
1	H	18	ASN
1	H	257	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

16 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
2	FAD	C	601	-	53,58,58	1.25	5 (9%)	68,89,89	1.66	12 (17%)
2	FAD	B	601	-	53,58,58	1.22	5 (9%)	68,89,89	1.78	17 (25%)
3	CC2	A	547	-	28,28,28	1.30	2 (7%)	38,41,41	1.40	6 (15%)
2	FAD	H	601	-	53,58,58	1.19	5 (9%)	68,89,89	1.63	12 (17%)
3	CC2	D	547	-	28,28,28	1.18	2 (7%)	38,41,41	1.47	8 (21%)
2	FAD	F	601	-	53,58,58	1.13	4 (7%)	68,89,89	1.67	12 (17%)
3	CC2	G	547	-	28,28,28	1.22	3 (10%)	38,41,41	1.38	5 (13%)
3	CC2	C	547	-	28,28,28	1.24	2 (7%)	38,41,41	1.23	5 (13%)
2	FAD	G	601	-	53,58,58	1.14	4 (7%)	68,89,89	1.72	13 (19%)
2	FAD	E	601	-	53,58,58	1.12	4 (7%)	68,89,89	1.51	11 (16%)
3	CC2	H	547	-	28,28,28	1.27	3 (10%)	38,41,41	1.32	6 (15%)
3	CC2	B	547	-	28,28,28	1.20	1 (3%)	38,41,41	1.35	8 (21%)
2	FAD	D	601	-	53,58,58	1.24	5 (9%)	68,89,89	1.72	16 (23%)
3	CC2	E	547	-	28,28,28	1.19	2 (7%)	38,41,41	1.10	2 (5%)
2	FAD	A	601	-	53,58,58	1.21	4 (7%)	68,89,89	1.98	15 (22%)
3	CC2	F	547	-	28,28,28	1.24	3 (10%)	38,41,41	1.21	5 (13%)

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
2	FAD	C	601	-	-	5/30/50/50	0/5/6/6
2	FAD	B	601	-	-	9/30/50/50	0/5/6/6
3	CC2	A	547	-	-	0/4/4/4	0/4/4/4
2	FAD	H	601	-	-	5/30/50/50	0/5/6/6
3	CC2	D	547	-	-	0/4/4/4	0/4/4/4
2	FAD	F	601	-	-	11/30/50/50	0/5/6/6
3	CC2	G	547	-	-	0/4/4/4	0/4/4/4
3	CC2	C	547	-	-	1/4/4/4	0/4/4/4
2	FAD	G	601	-	-	4/30/50/50	0/5/6/6
2	FAD	E	601	-	-	5/30/50/50	0/5/6/6
3	CC2	H	547	-	-	0/4/4/4	0/4/4/4
3	CC2	B	547	-	-	0/4/4/4	0/4/4/4
2	FAD	D	601	-	-	6/30/50/50	0/5/6/6
3	CC2	E	547	-	-	1/4/4/4	0/4/4/4
2	FAD	A	601	-	-	7/30/50/50	0/5/6/6
3	CC2	F	547	-	-	1/4/4/4	0/4/4/4

The worst 5 of 54 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	601	FAD	C2A-N3A	4.28	1.39	1.32
2	D	601	FAD	C4X-N5	4.21	1.38	1.30
2	A	601	FAD	C2A-N3A	4.07	1.38	1.32
2	H	601	FAD	C2A-N3A	4.02	1.38	1.32
2	E	601	FAD	C4X-N5	3.87	1.38	1.30

The worst 5 of 153 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	601	FAD	C1'-N10-C9A	-9.00	105.50	120.51
2	G	601	FAD	C1'-N10-C9A	-7.37	108.22	120.51
2	F	601	FAD	C1'-N10-C9A	-6.28	110.05	120.51
2	H	601	FAD	C1'-N10-C9A	-6.09	110.36	120.51
2	C	601	FAD	C1'-N10-C9A	-5.68	111.03	120.51

There are no chirality outliers.

5 of 55 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	601	FAD	C5'-O5'-P-O1P
2	B	601	FAD	C5B-O5B-PA-O3P

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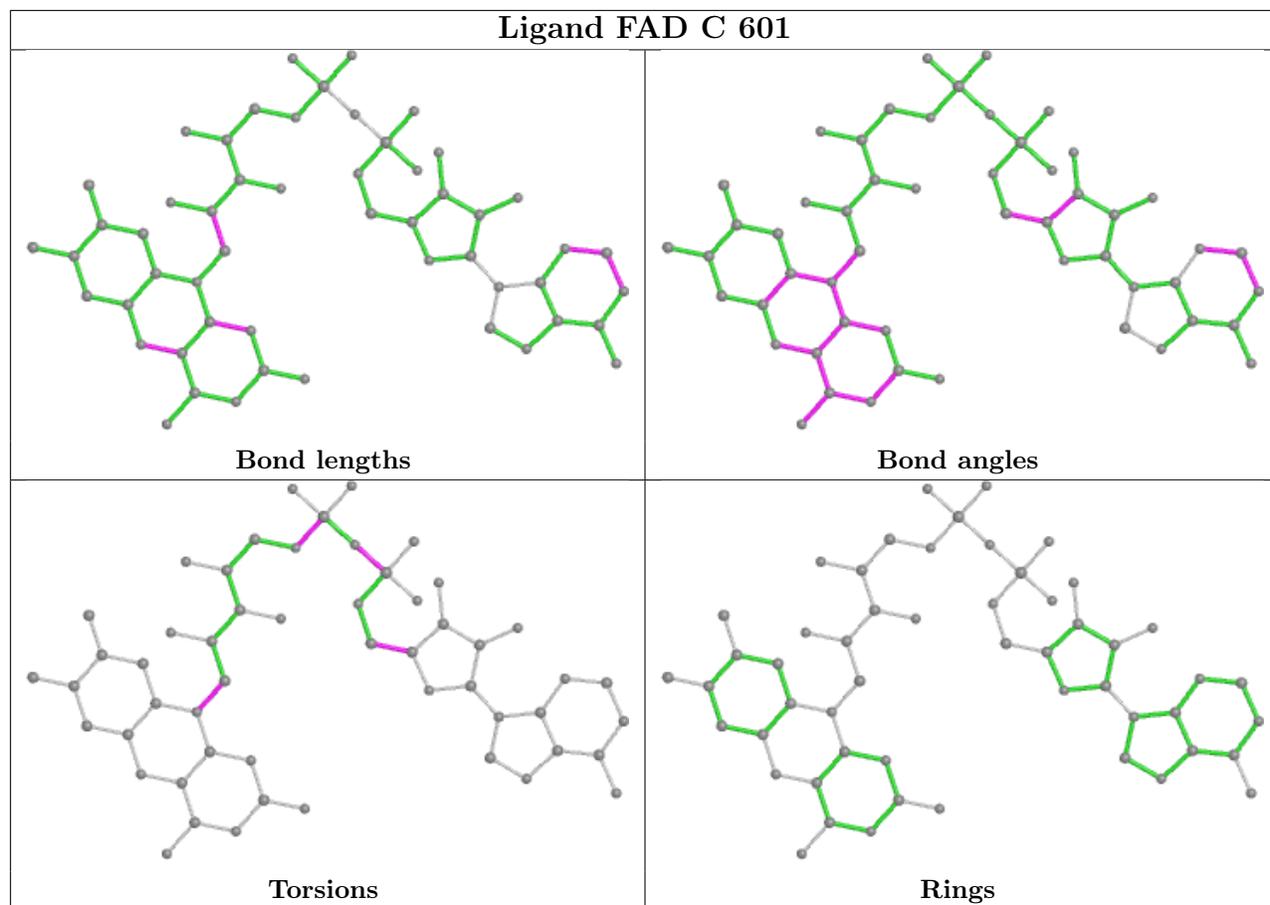
Mol	Chain	Res	Type	Atoms
2	B	601	FAD	C2'-C1'-N10-C10
2	B	601	FAD	C5'-O5'-P-O3P
2	C	601	FAD	C2'-C1'-N10-C10

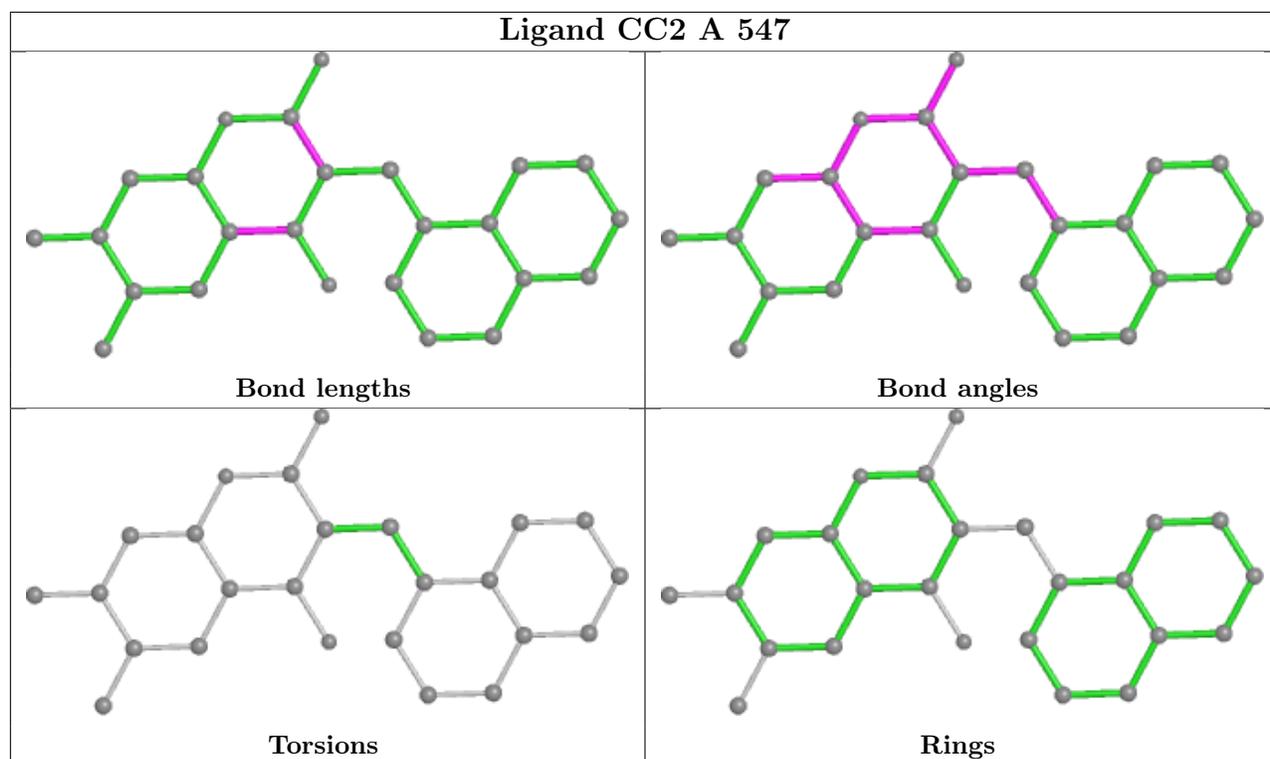
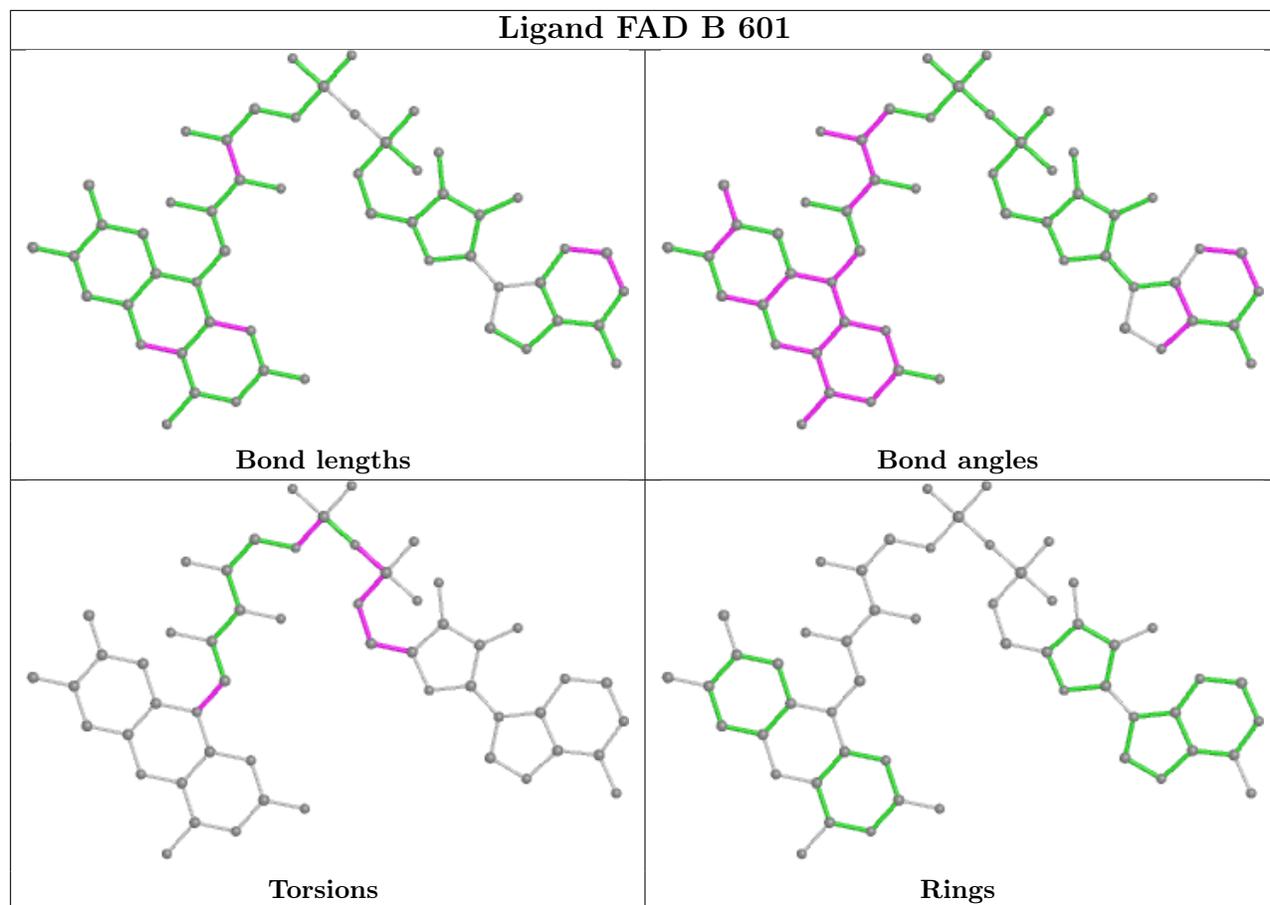
There are no ring outliers.

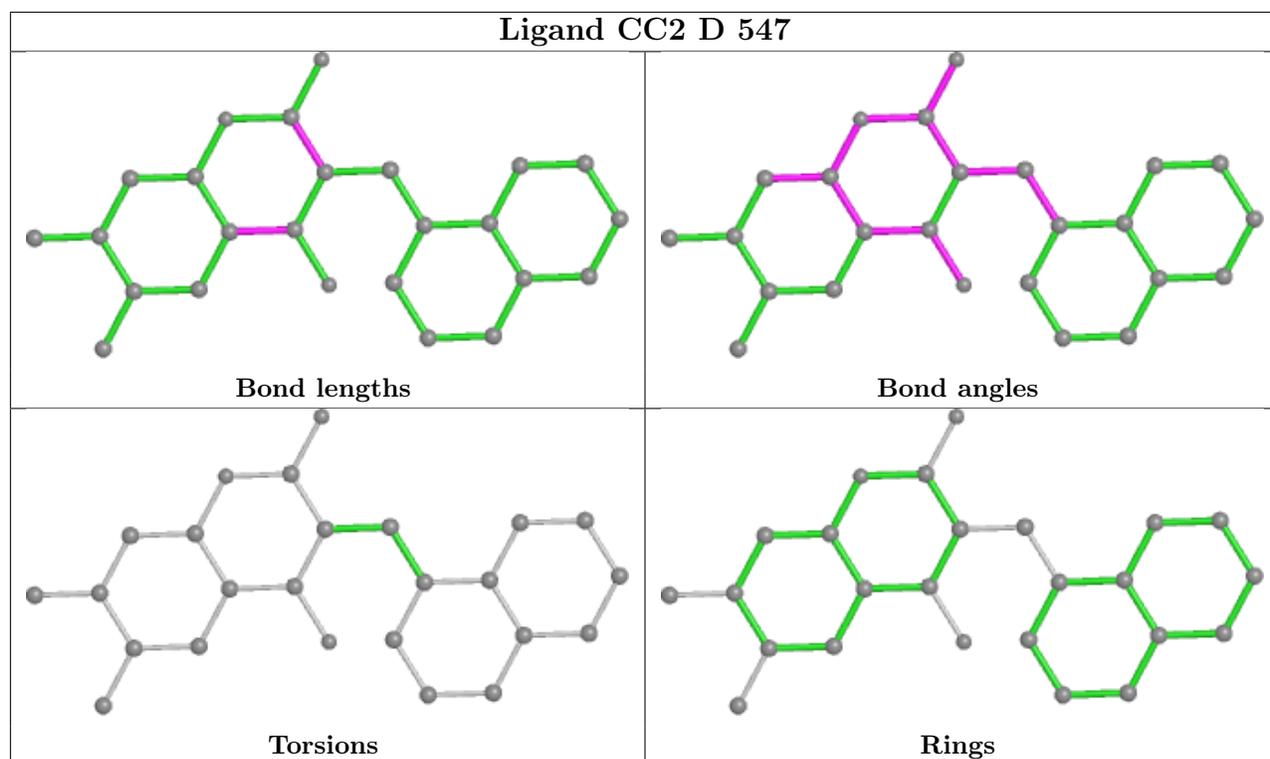
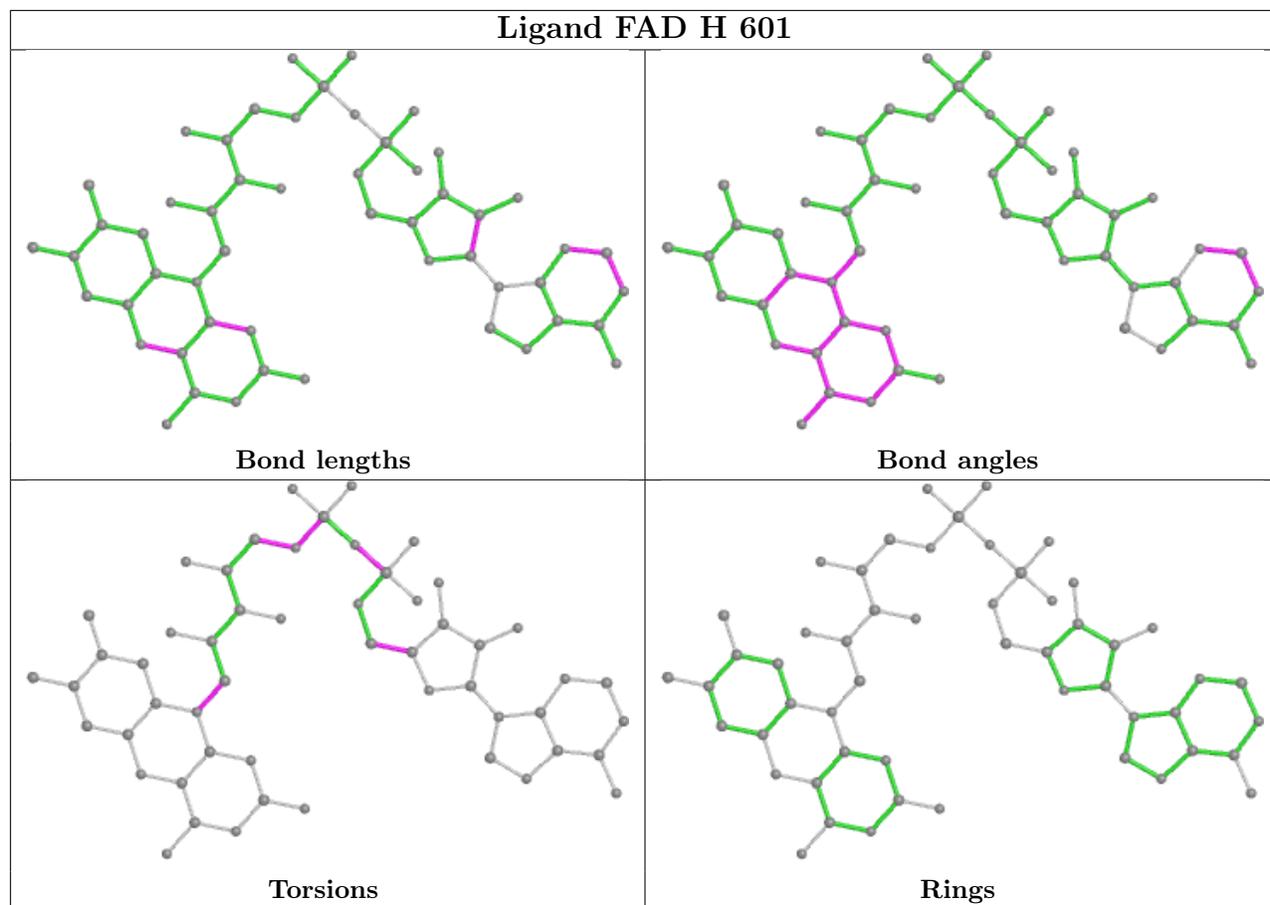
16 monomers are involved in 74 short contacts:

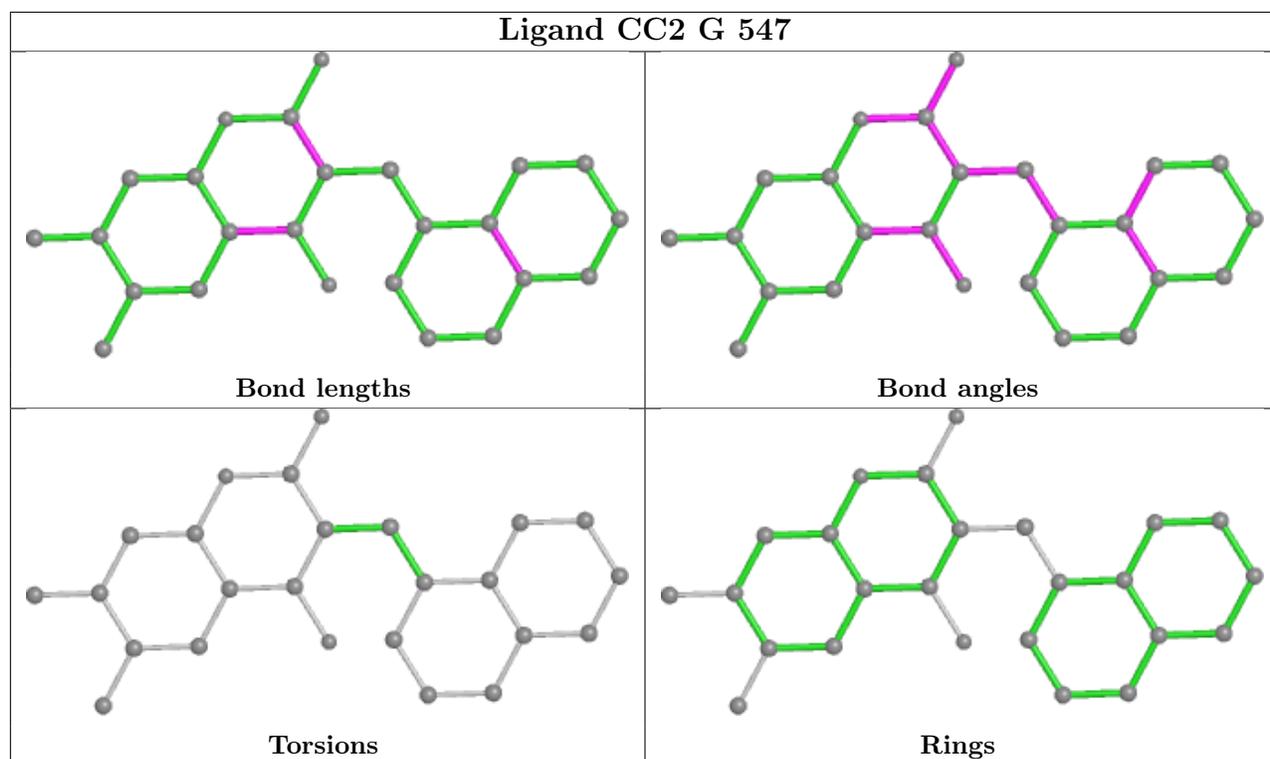
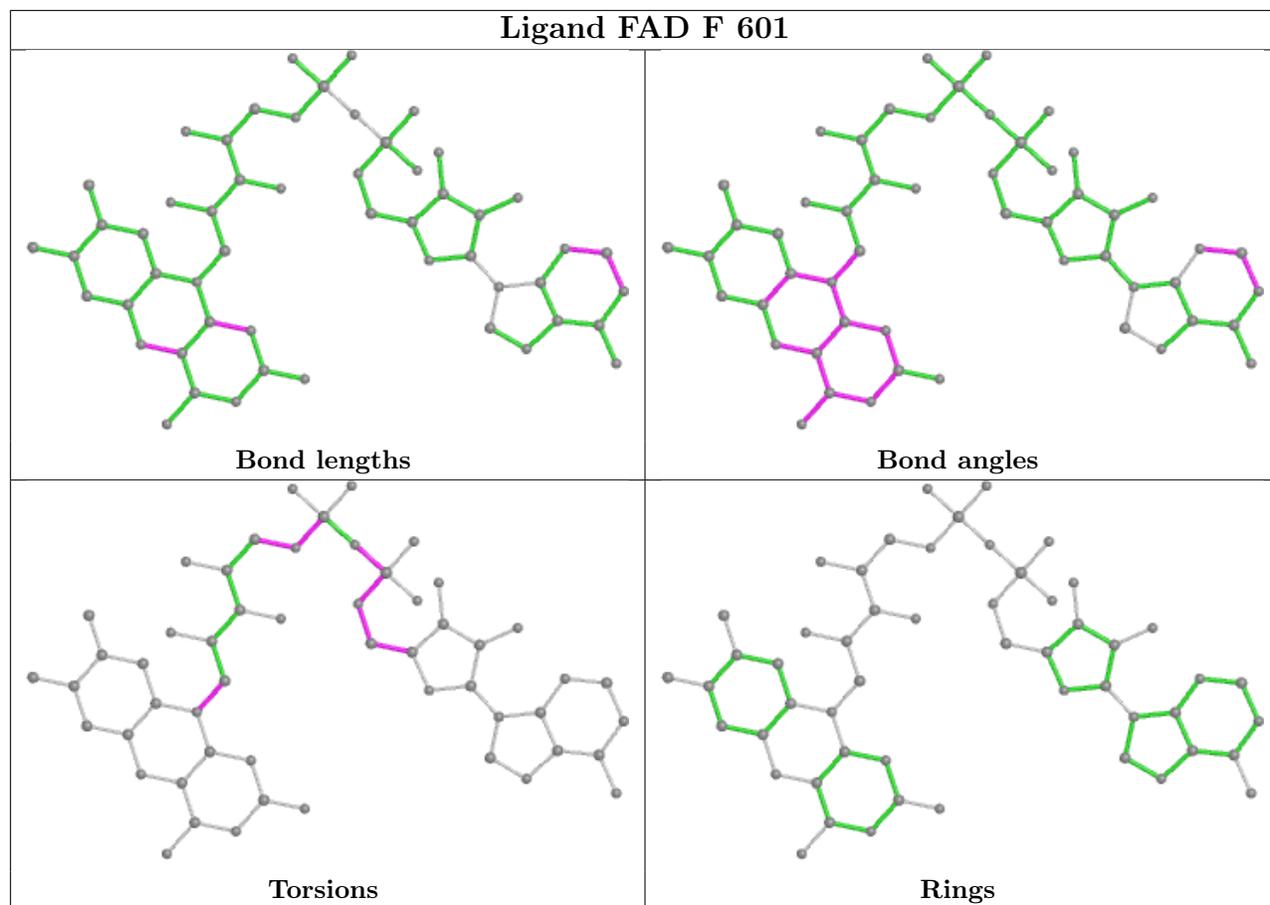
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	601	FAD	6	0
2	B	601	FAD	10	0
3	A	547	CC2	1	0
2	H	601	FAD	9	0
3	D	547	CC2	2	0
2	F	601	FAD	4	0
3	G	547	CC2	2	0
3	C	547	CC2	4	0
2	G	601	FAD	2	0
2	E	601	FAD	4	0
3	H	547	CC2	3	0
3	B	547	CC2	10	0
2	D	601	FAD	6	0
3	E	547	CC2	3	0
2	A	601	FAD	7	0
3	F	547	CC2	3	0

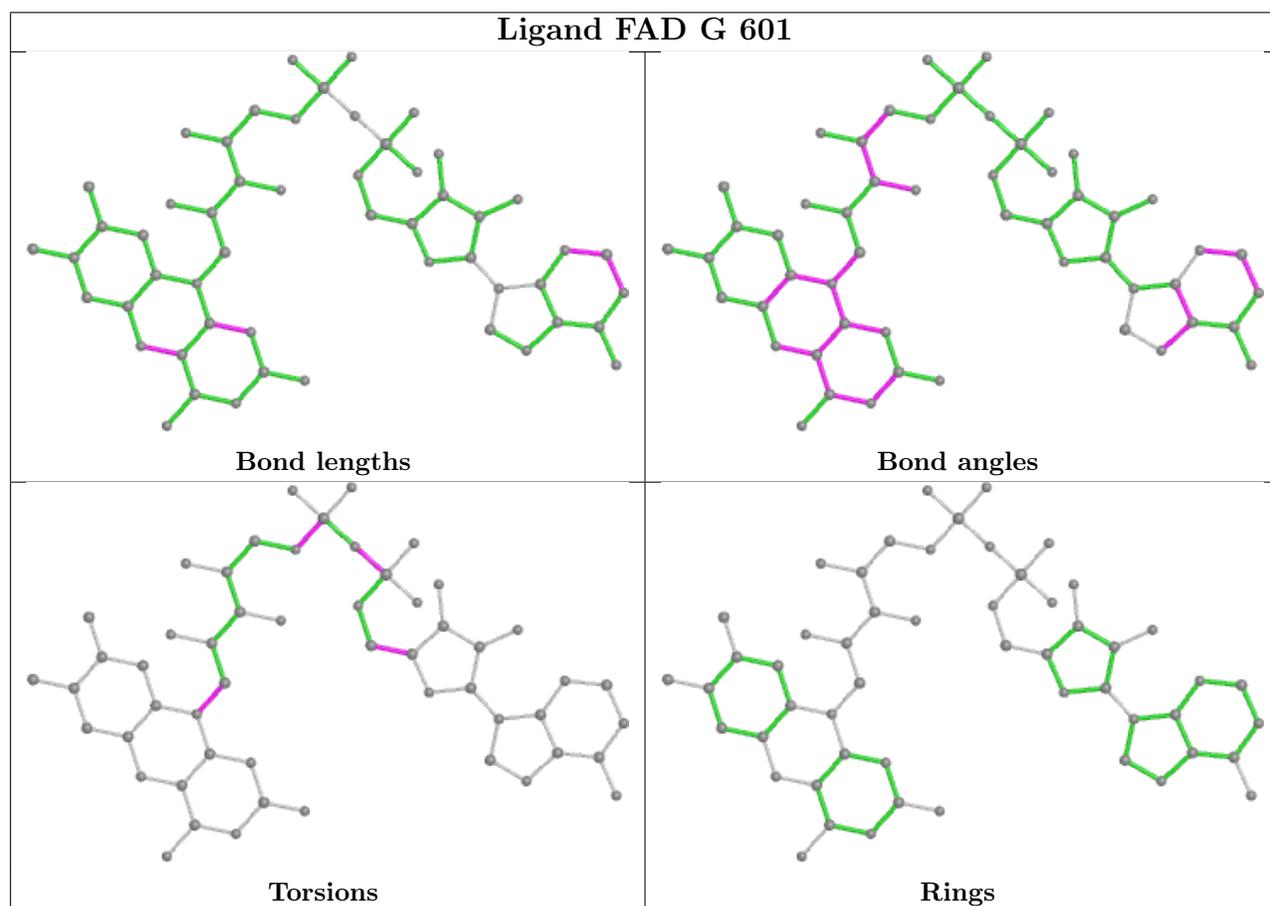
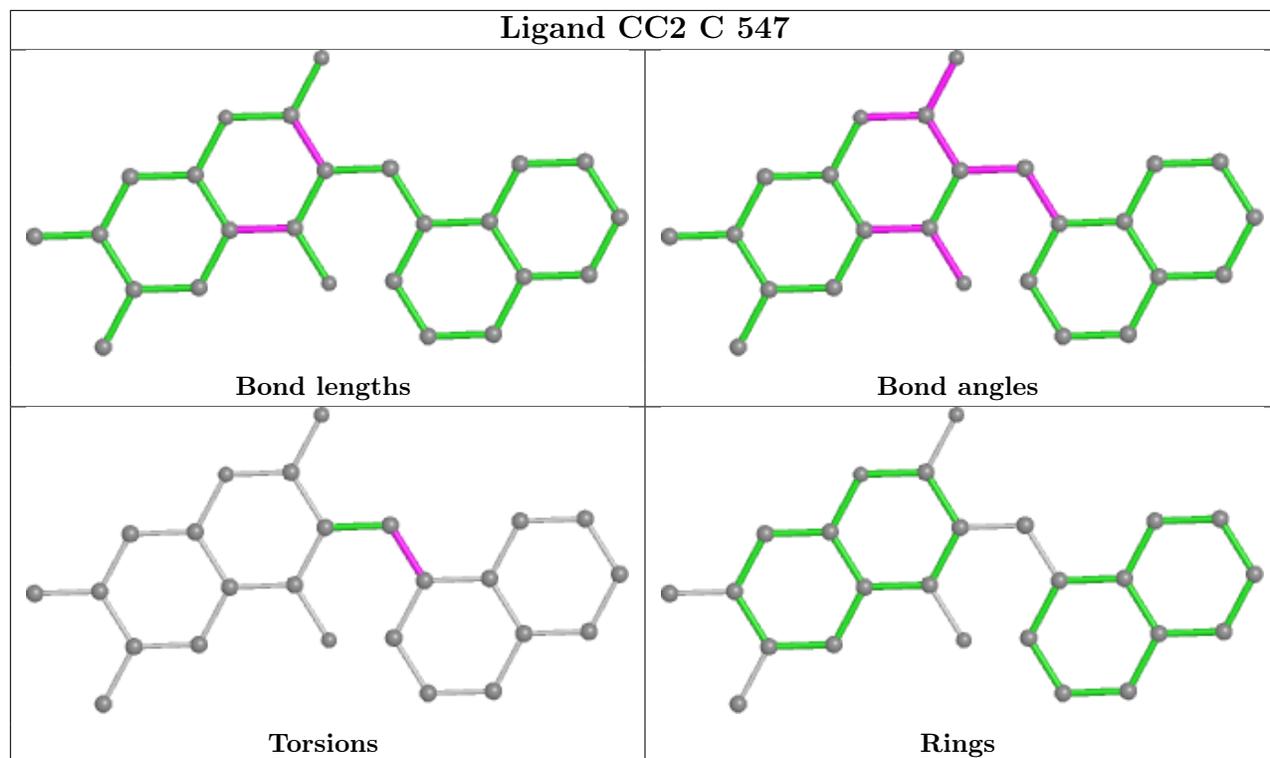
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

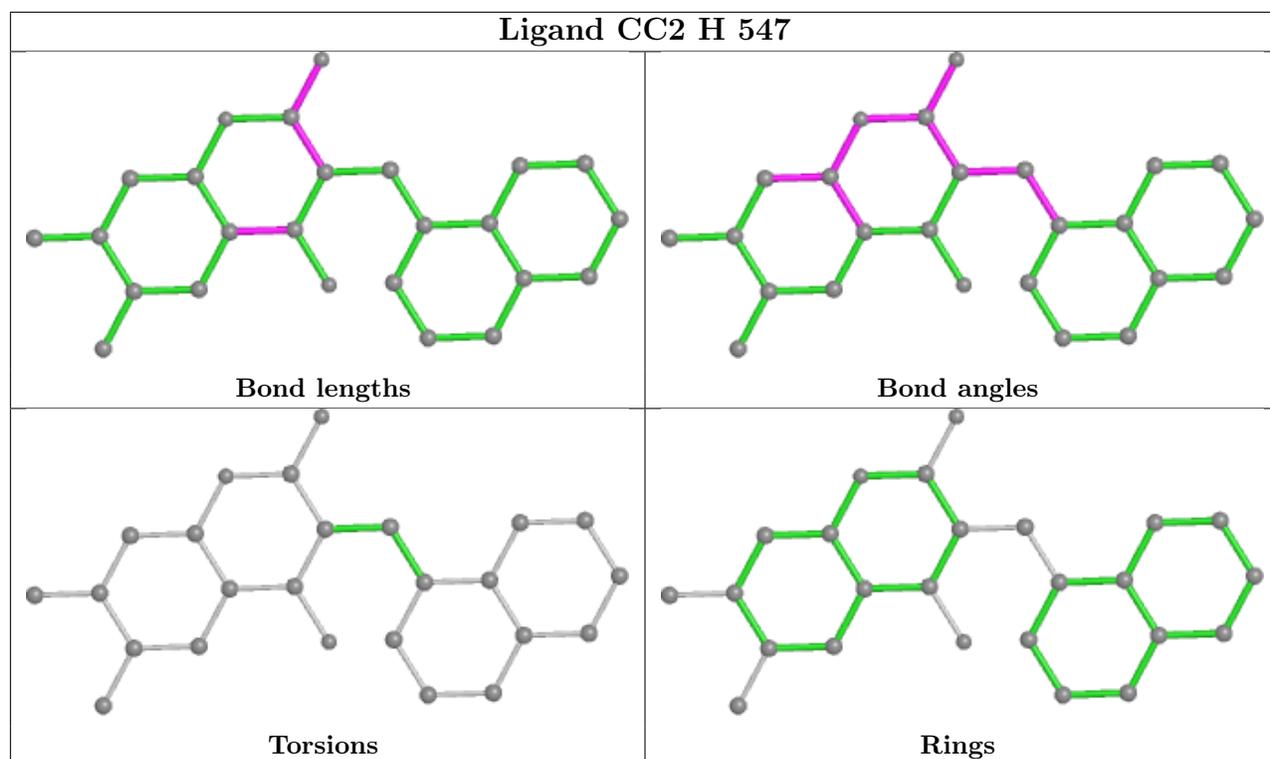
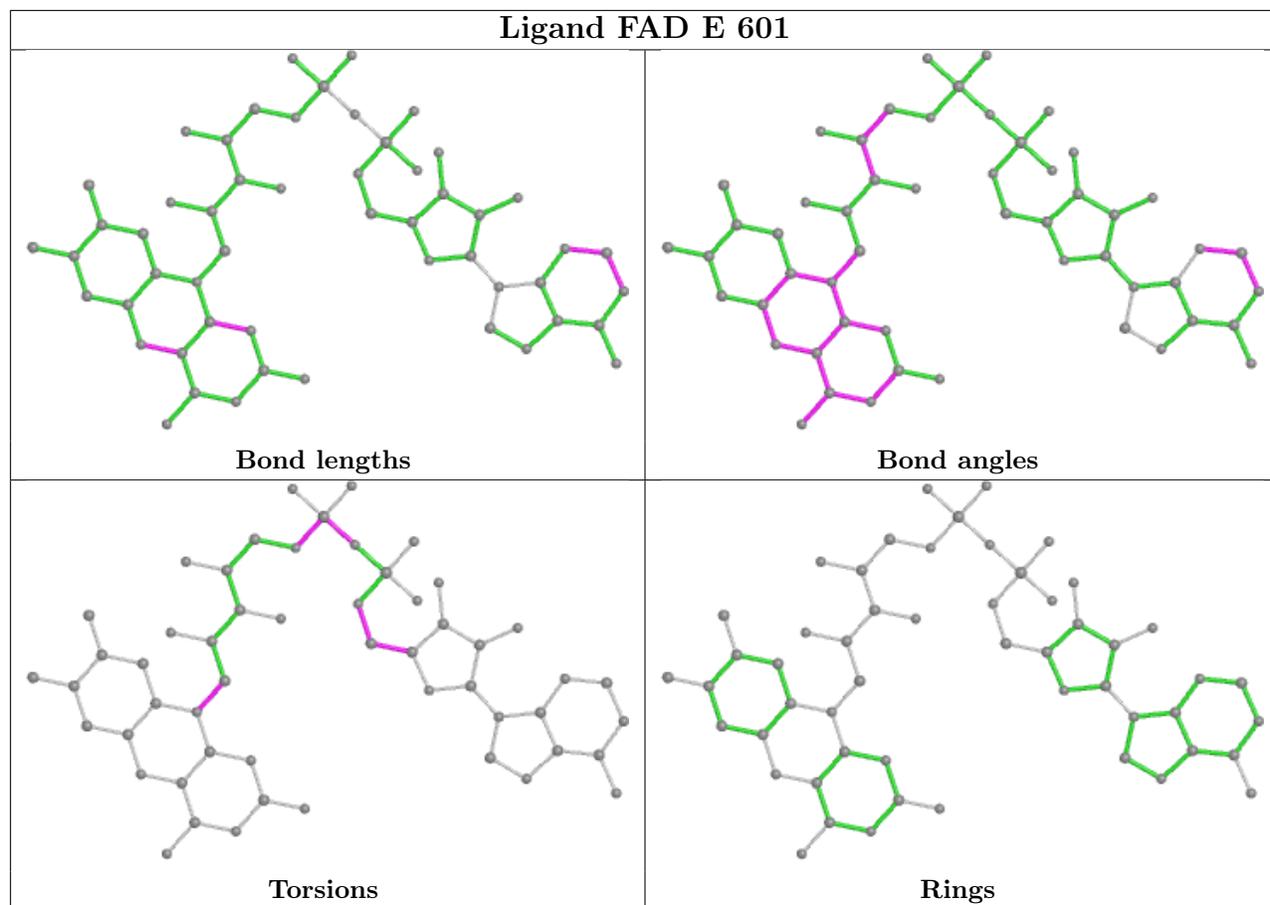


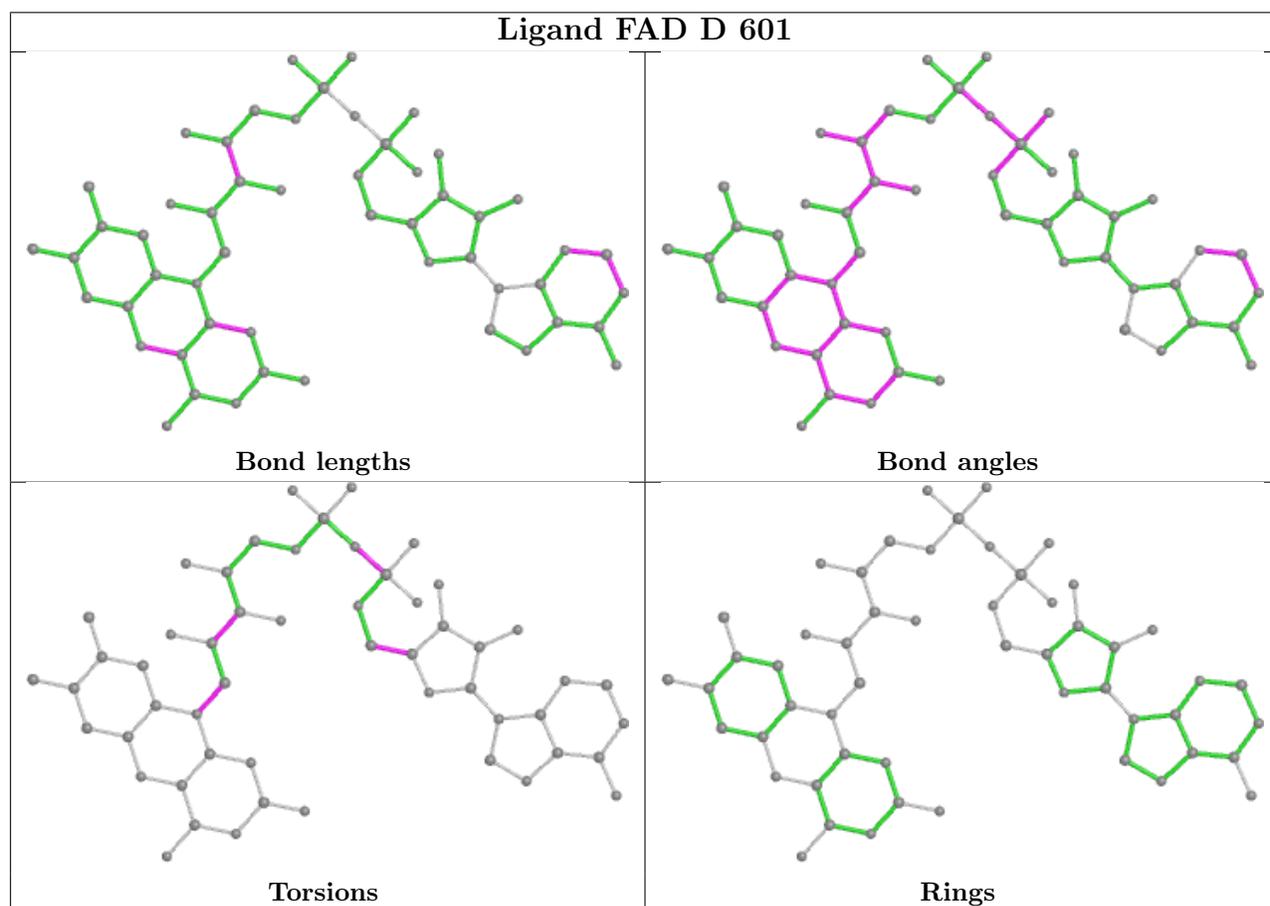
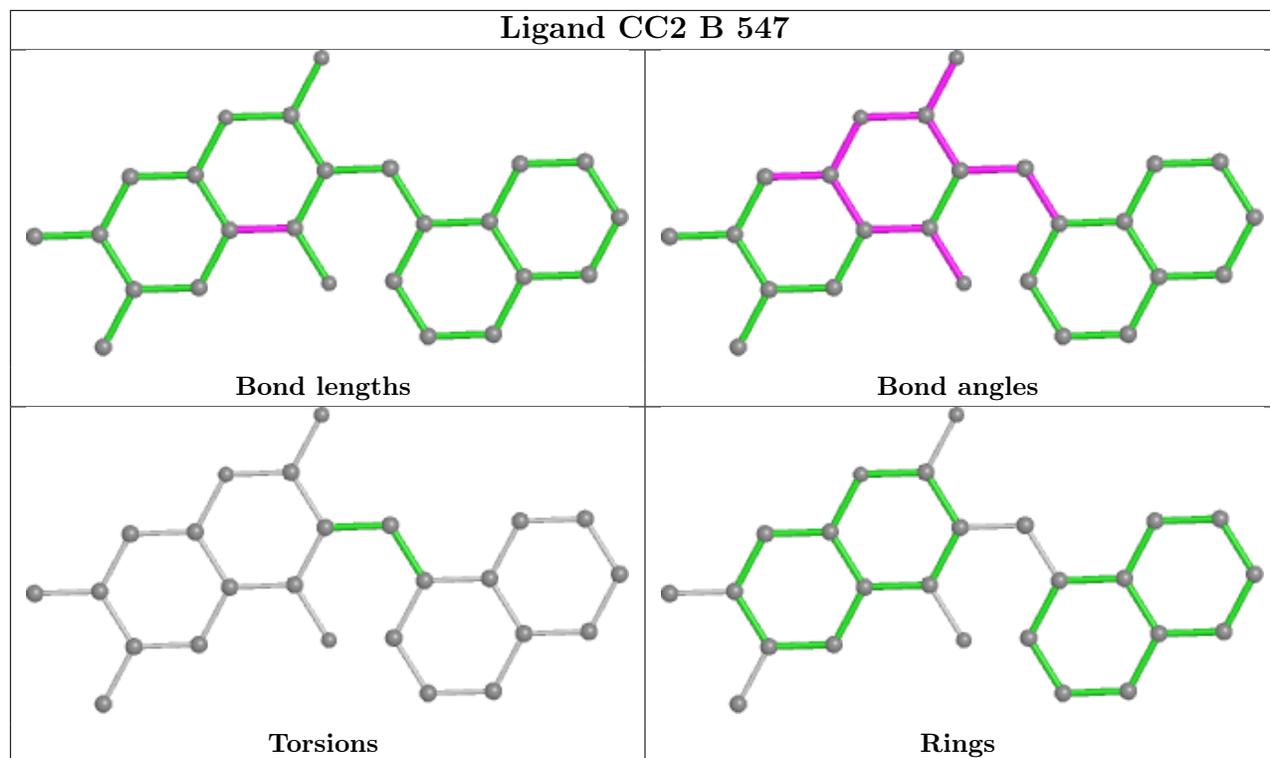


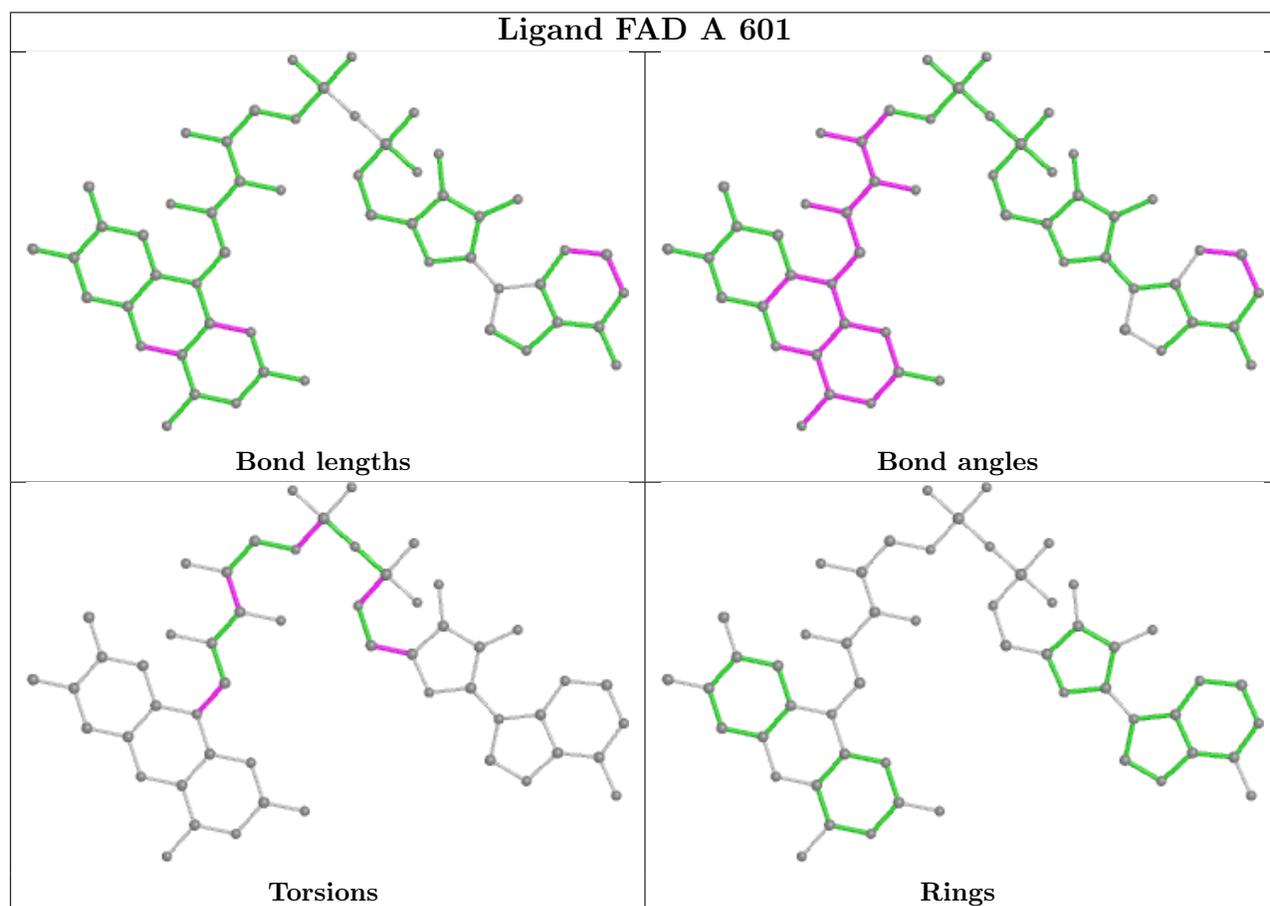
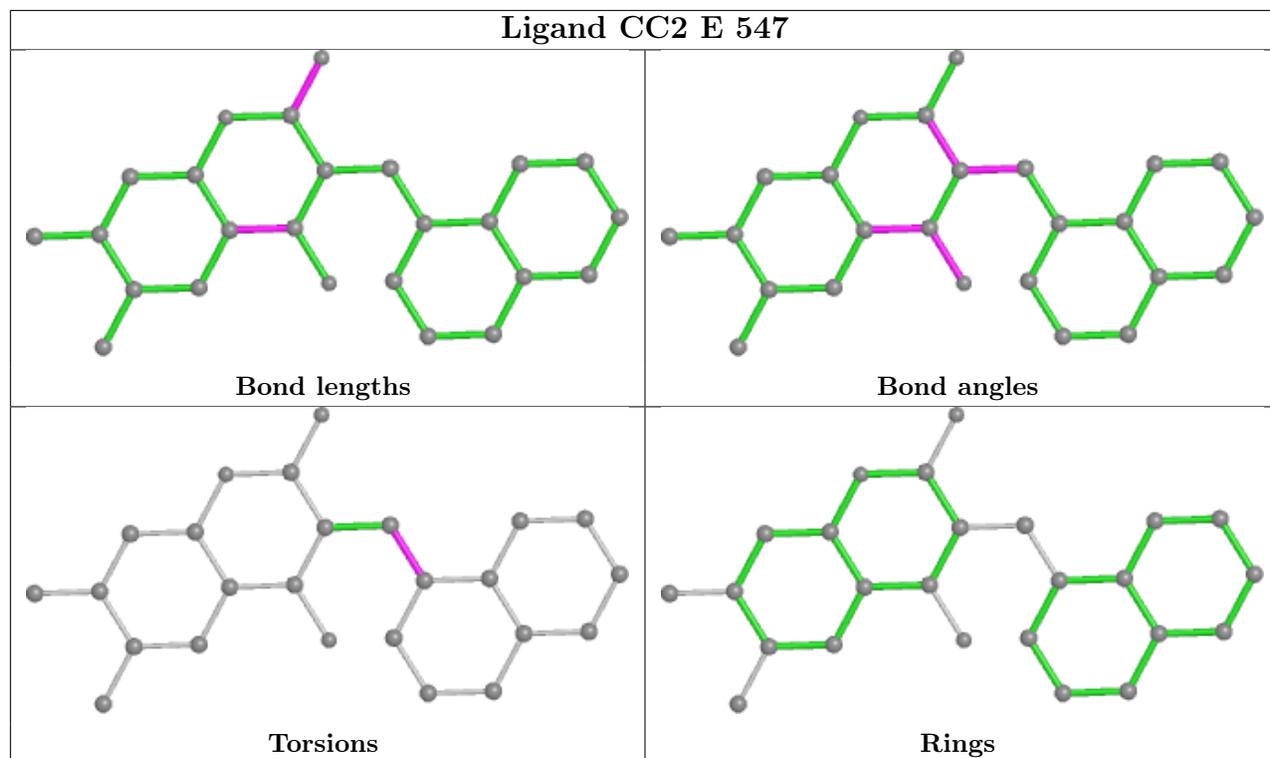


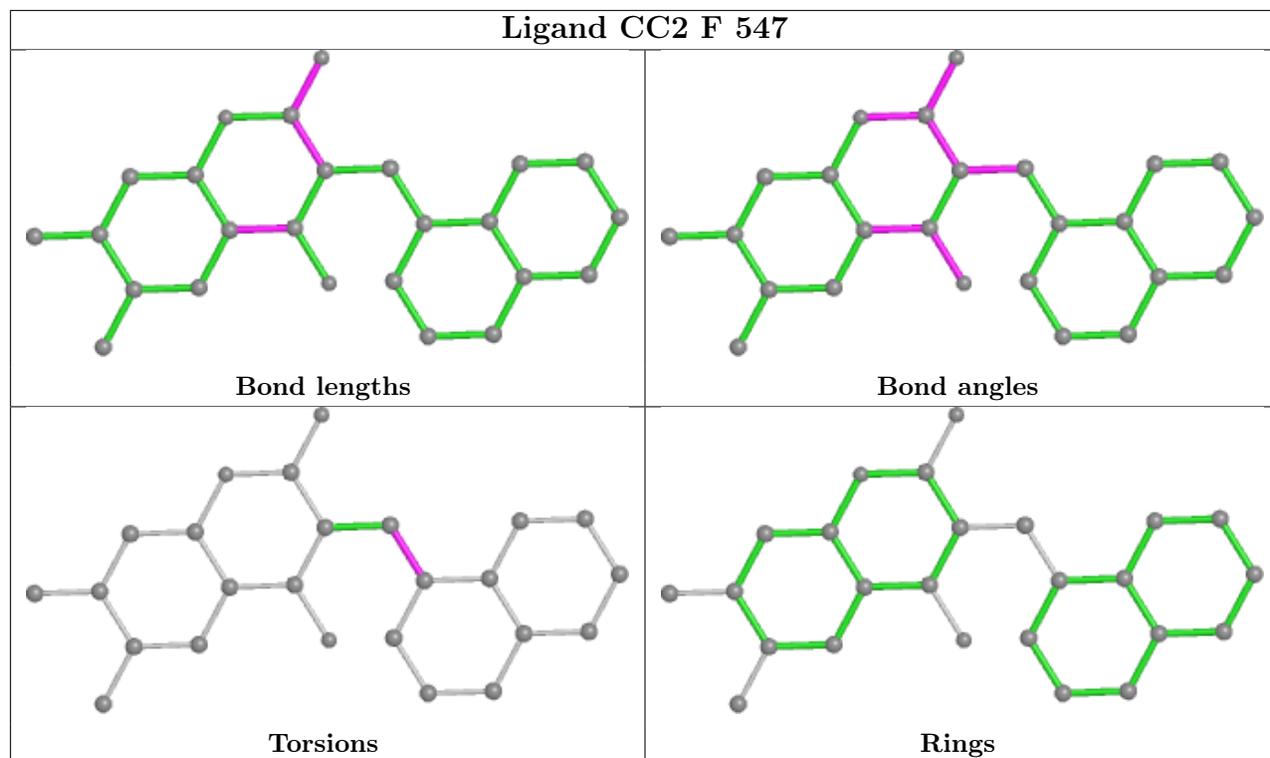












5.7 Other polymers [\(i\)](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [\(i\)](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	270/273 (98%)	0.09	4 (1%) 73 71	42, 61, 78, 92	0
1	B	270/273 (98%)	-0.04	0 100 100	42, 56, 75, 87	0
1	C	270/273 (98%)	-0.07	0 100 100	41, 57, 75, 91	0
1	D	270/273 (98%)	0.03	2 (0%) 87 88	42, 59, 78, 90	0
1	E	270/273 (98%)	0.09	5 (1%) 66 64	51, 67, 89, 98	0
1	F	270/273 (98%)	0.13	4 (1%) 73 71	50, 66, 87, 102	0
1	G	270/273 (98%)	0.15	5 (1%) 66 64	50, 69, 95, 105	0
1	H	270/273 (98%)	0.27	9 (3%) 46 43	51, 70, 93, 105	0
All	All	2160/2184 (98%)	0.08	29 (1%) 77 76	41, 63, 87, 105	0

The worst 5 of 29 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	H	2	GLY	10.3
1	F	2	GLY	5.0
1	G	2	GLY	4.5
1	H	59	LEU	3.7
1	A	232	PHE	3.1

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

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

6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

6.4 Ligands

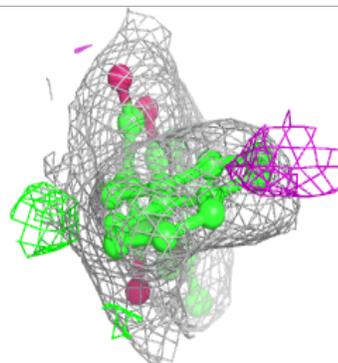
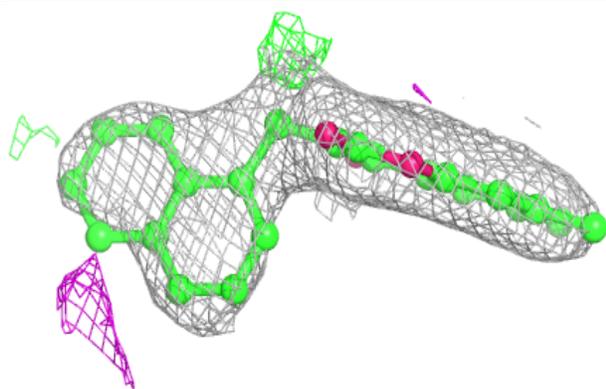
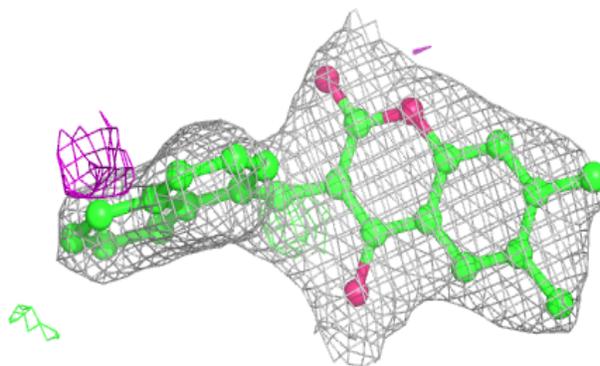
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	CC2	B	547	25/25	0.91	0.22	57,66,87,89	0
3	CC2	C	547	25/25	0.93	0.21	52,62,85,89	0
2	FAD	F	601	53/53	0.94	0.15	53,62,69,76	0
3	CC2	F	547	25/25	0.94	0.23	59,68,81,82	0
3	CC2	E	547	25/25	0.95	0.16	60,68,81,85	0
3	CC2	A	547	25/25	0.96	0.16	44,49,53,56	0
2	FAD	C	601	53/53	0.96	0.14	39,54,63,69	0
2	FAD	E	601	53/53	0.96	0.15	56,63,73,80	0
2	FAD	B	601	53/53	0.96	0.14	46,58,66,68	0
2	FAD	H	601	53/53	0.96	0.15	43,56,69,73	0
3	CC2	H	547	25/25	0.96	0.15	51,55,58,60	0
3	CC2	D	547	25/25	0.97	0.14	46,51,54,55	0
2	FAD	D	601	53/53	0.97	0.14	42,50,62,68	0
2	FAD	G	601	53/53	0.97	0.14	42,54,62,68	0
3	CC2	G	547	25/25	0.97	0.17	53,56,59,60	0
2	FAD	A	601	53/53	0.97	0.15	42,49,61,64	0

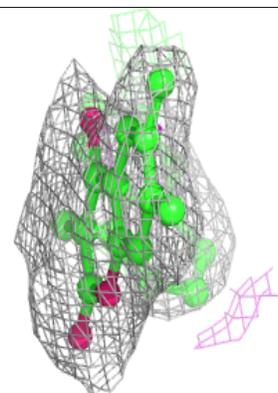
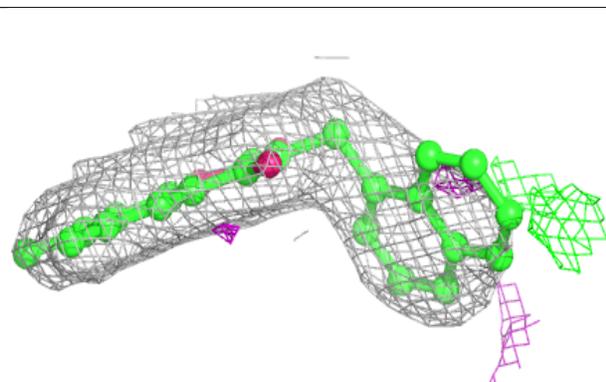
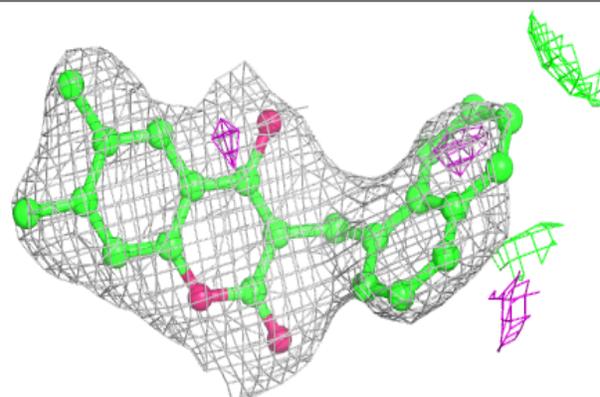
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around CC2 B 547:

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

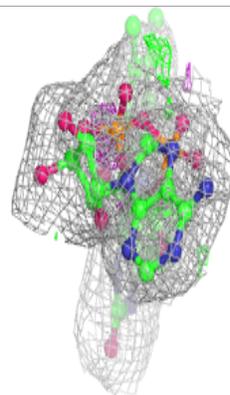
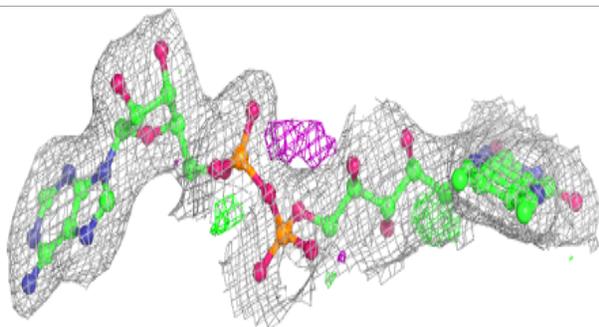
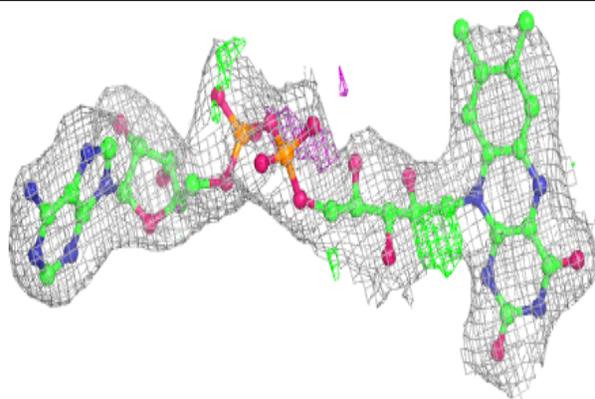
**Electron density around CC2 C 547:**

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

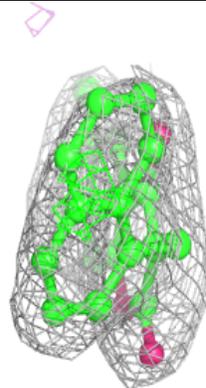
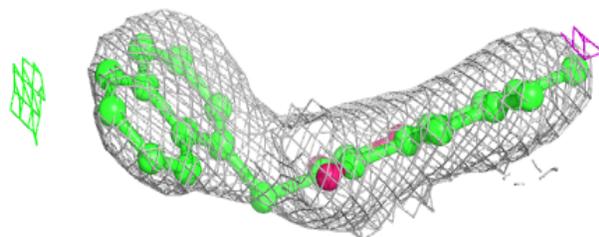
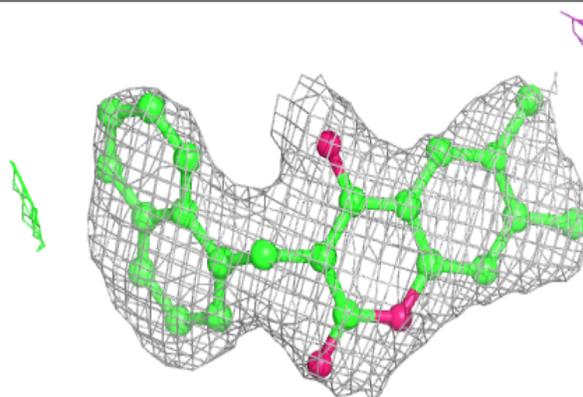


Electron density around FAD F 601:

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

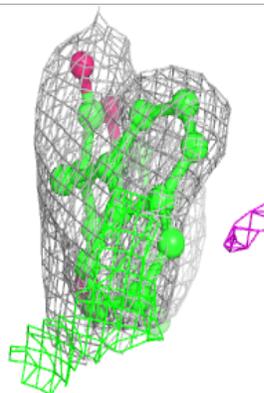
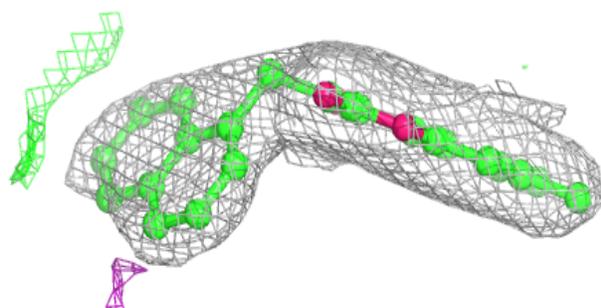
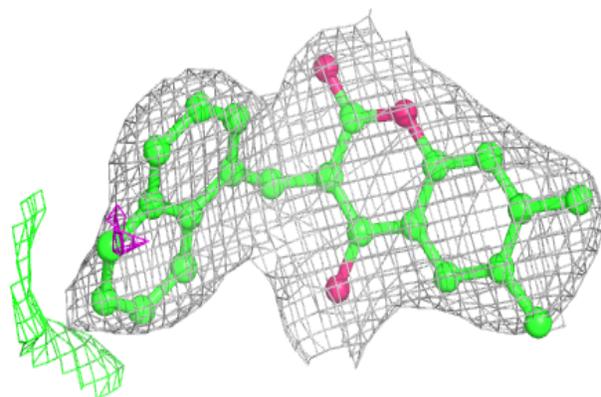
**Electron density around CC2 F 547:**

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

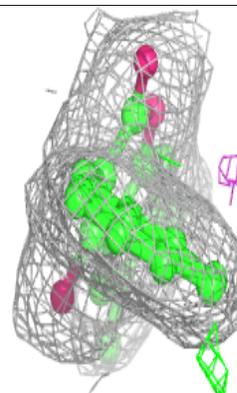
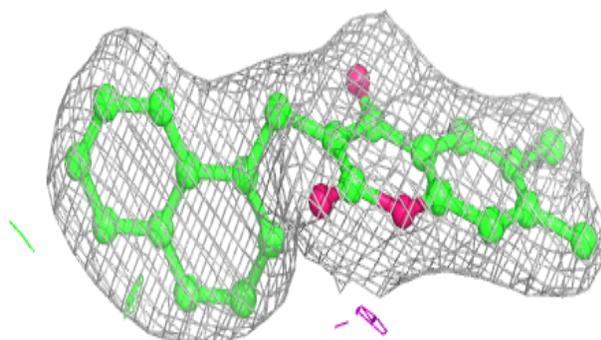
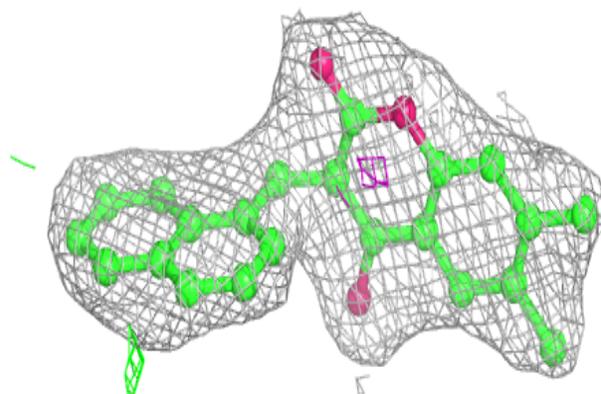


Electron density around CC2 E 547:

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and green (positive)

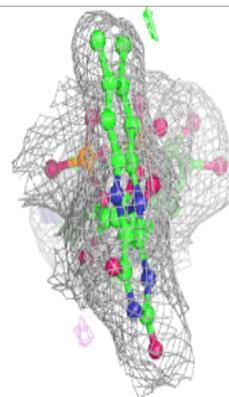
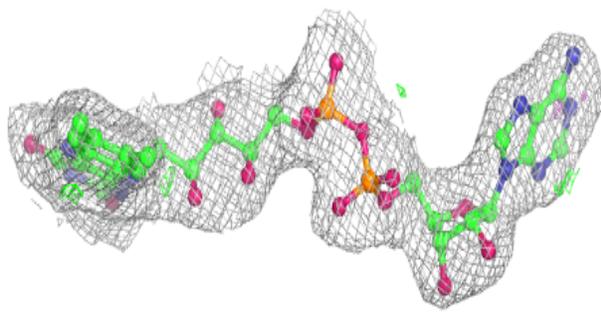
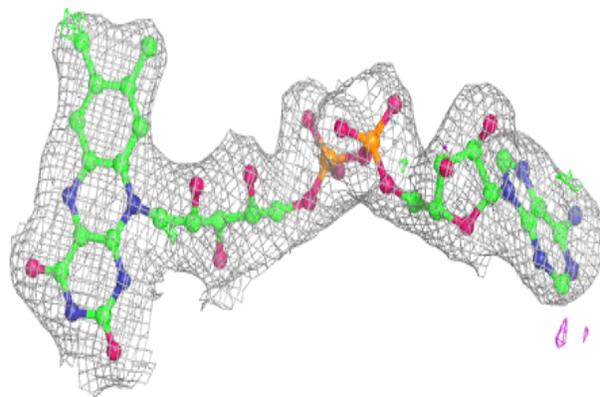
**Electron density around CC2 A 547:**

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and green (positive)

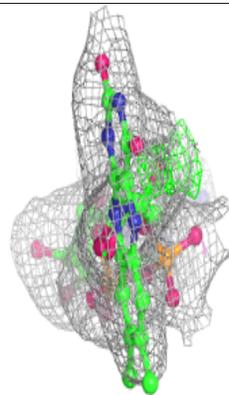
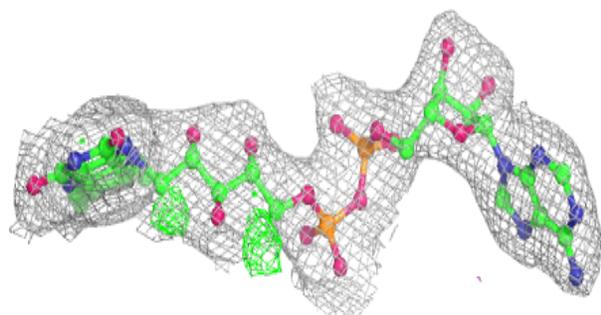
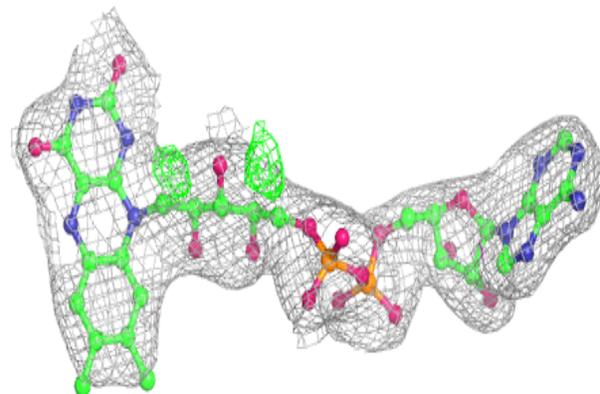


Electron density around FAD C 601:

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

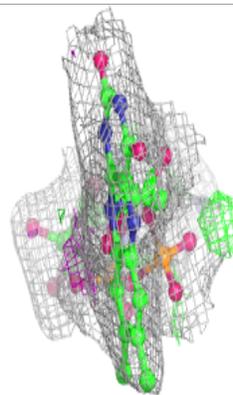
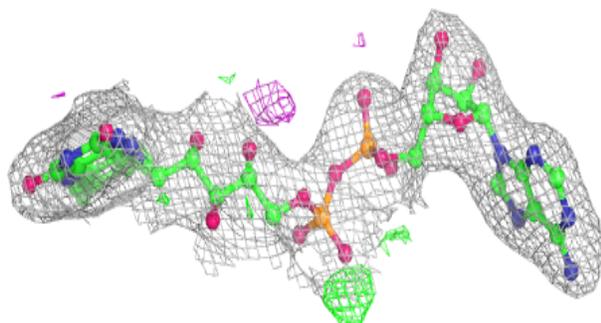
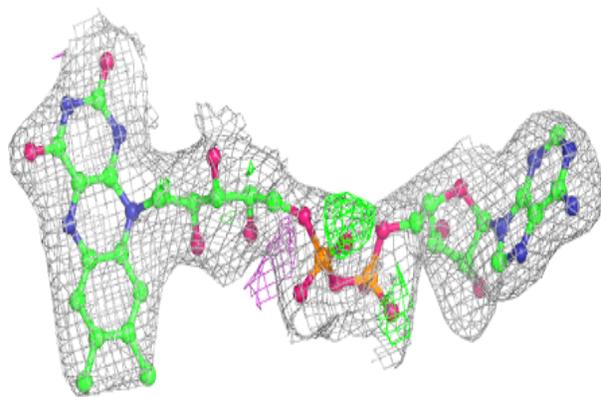
**Electron density around FAD E 601:**

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and green (positive)

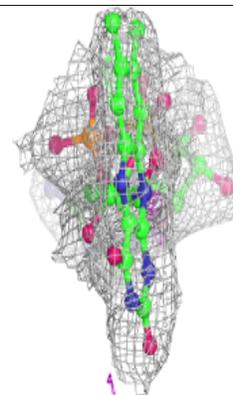
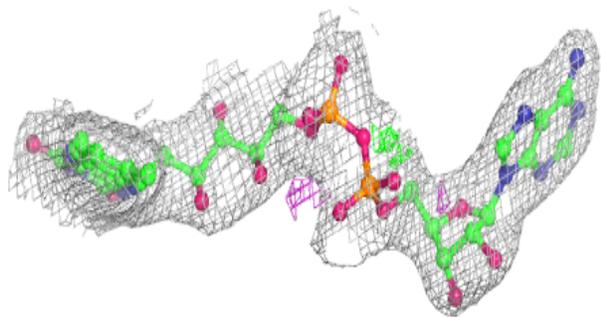
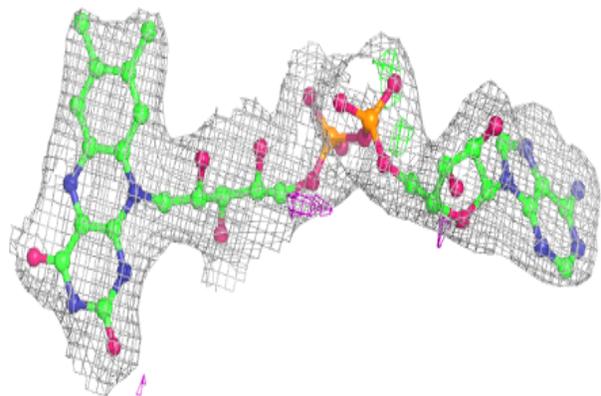


Electron density around FAD B 601:

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and green (positive)

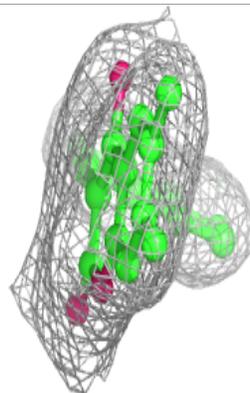
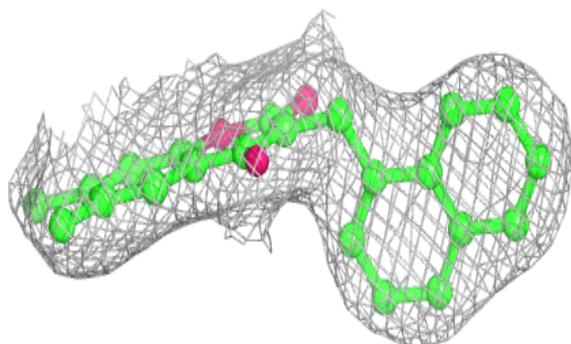
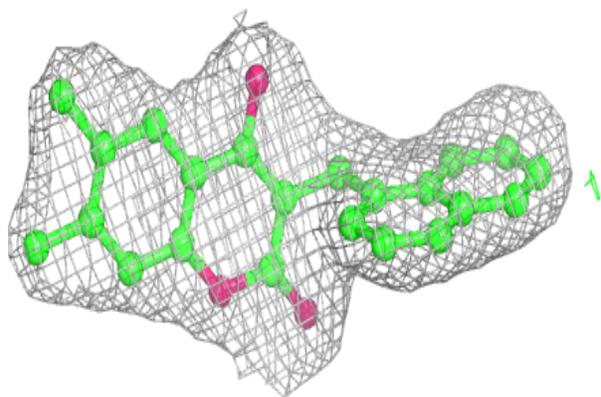
**Electron density around FAD H 601:**

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

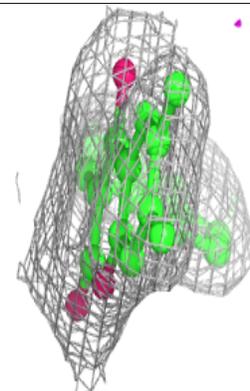
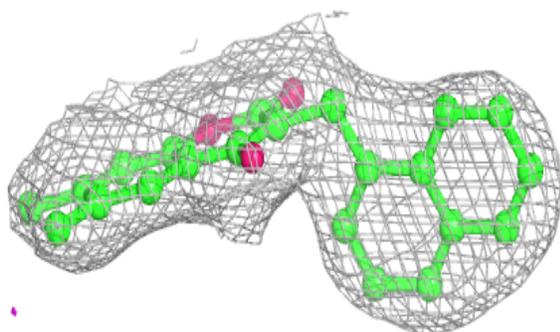
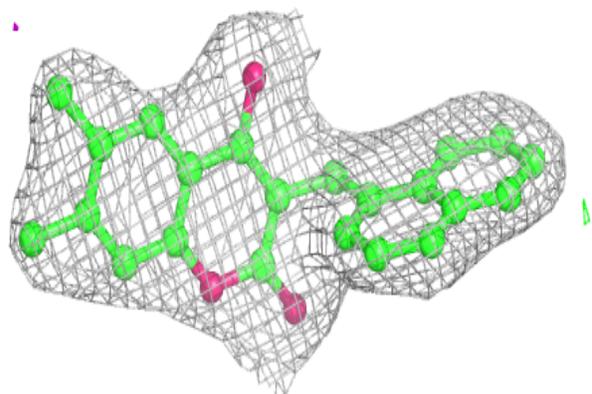


Electron density around CC2 H 547:

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

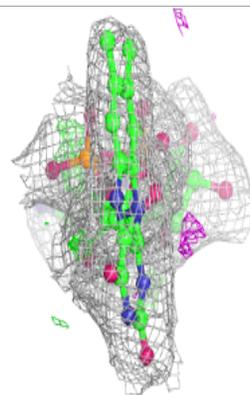
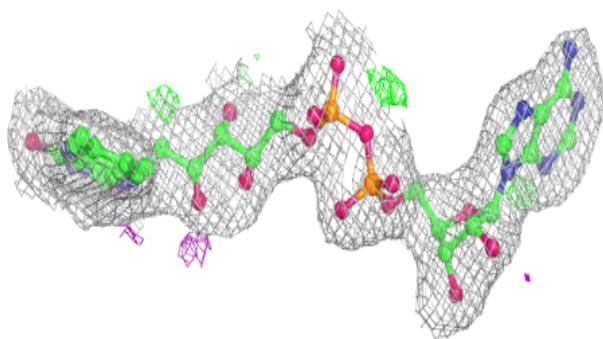
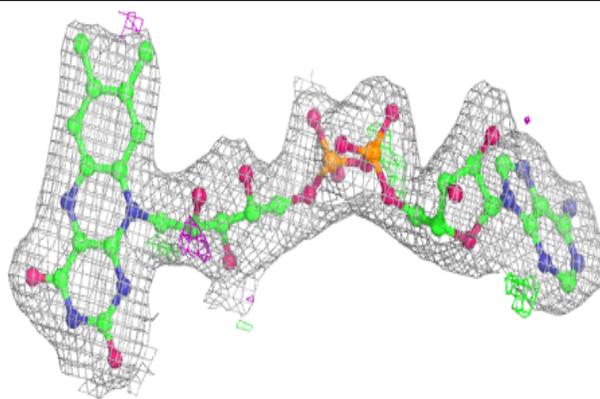
**Electron density around CC2 D 547:**

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
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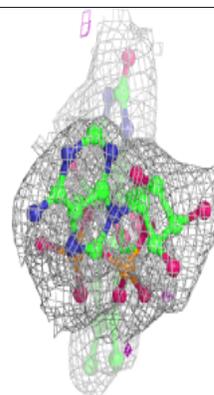
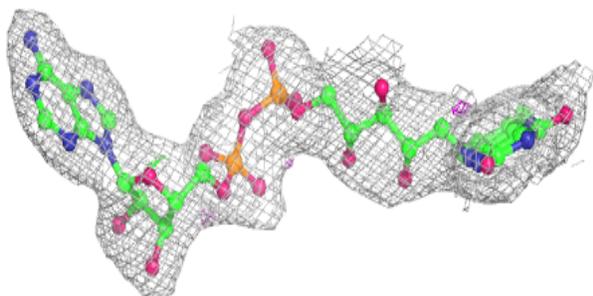
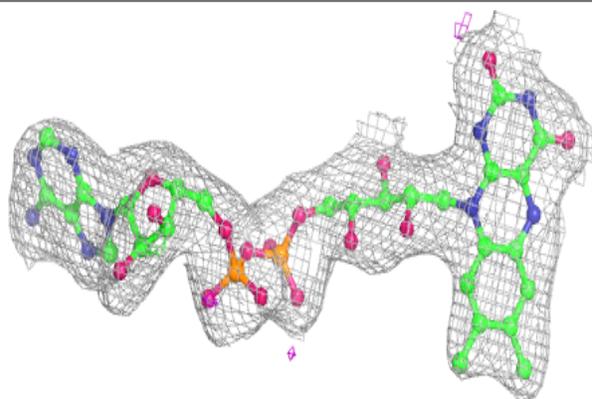


Electron density around FAD D 601:

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and green (positive)

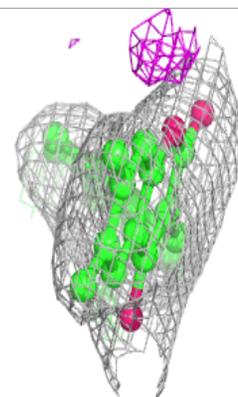
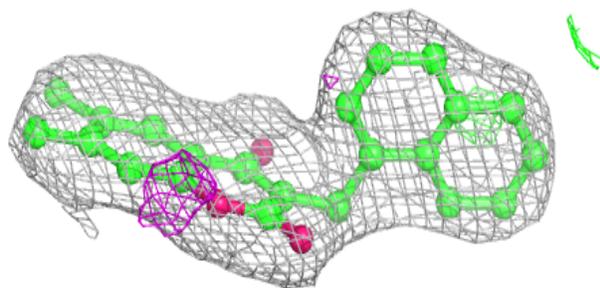
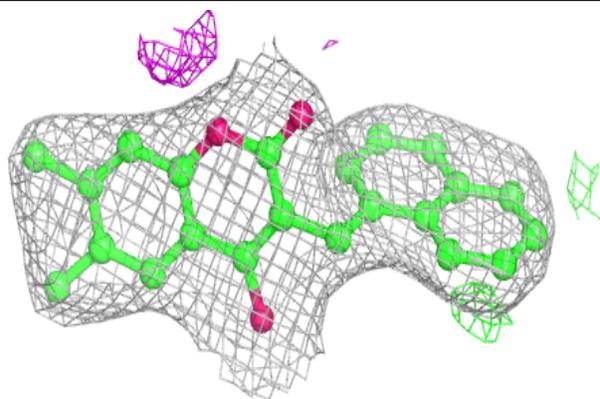
**Electron density around FAD G 601:**

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and green (positive)

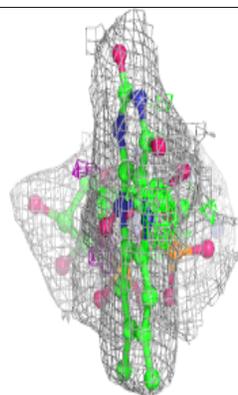
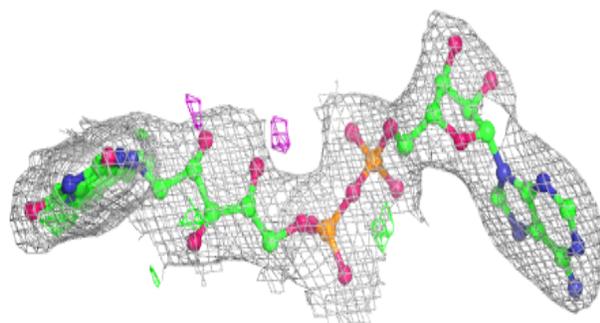
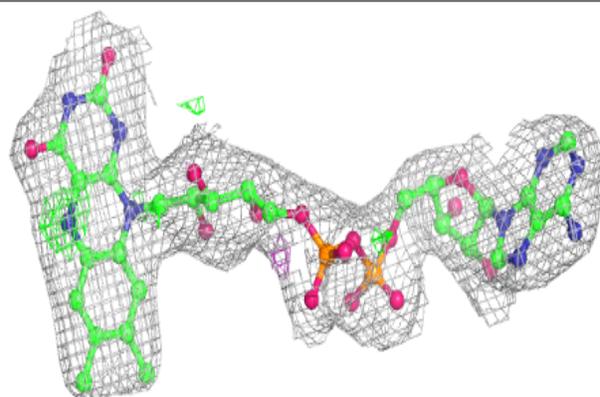


Electron density around CC2 G 547:

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

**Electron density around FAD A 601:**

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



6.5 Other polymers [i](#)

There are no such residues in this entry.