WORLDWIDE PDB PROTEIN DATA BANK

Full wwPDB X-ray Structure Validation Report ⓘ

Oct 7, 2023 – 11:41 PM EDT

PDB ID	:	4QFC
Title	:	Co-crystal structure of compound 3 (4-hydroxy-6-[2-(7-hydroxy-2-oxo-4-phenyl-2h-chromen-6-yl)ethyl]pyridazin-3(2h)-one) and FAD bound to human DAAO at 2.4A
Authors	:	Lukacs, C.M.; Chun, L.
Deposited on	:	2014-05-20
Resolution	:	2.40 Å(reported)

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

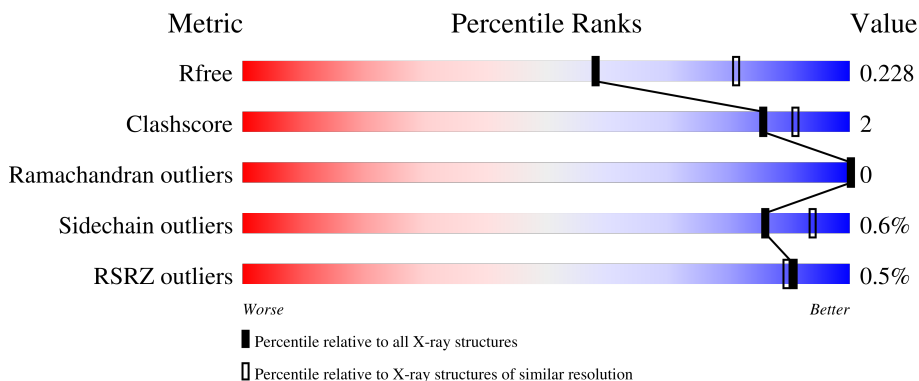|		
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.35.1
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35.1

1   Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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	Percentile Ranks	Value
Rfree		0.228
Clashscore		2
Ramachandran outliers		0
Sidechain outliers		0.6%
RSRZ outliers		0.5%

Worse   *Better*

▌ Percentile relative to all X-ray structures

▯ Percentile relative to X-ray structures of similar resolution

Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R$_{free}$	130704	3907 (2.40-2.40)
Clashscore	141614	4398 (2.40-2.40)
Ramachandran outliers	138981	4318 (2.40-2.40)
Sidechain outliers	138945	4319 (2.40-2.40)
RSRZ outliers	127900	3811 (2.40-2.40)

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 <=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	347	90% 5% 5%
1	B	347	.% 87% 6% 7%

2 Entry composition ⓘ

There are 5 unique types of molecules in this entry. The entry contains 5373 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 D-amino-acid oxidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	330	Total 2588	C 1665	N 443	O 471	S 9	0	1	0
1	B	323	Total 2524	C 1628	N 436	O 451	S 9	0	0	0

- 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
2	A	1	Total 53	C 27	N 9	O 15	P 2	0	0
2	B	1	Total 53	C 27	N 9	O 15	P 2	0	0

- Molecule 3 is 4-hydroxy-6-[2-(7-hydroxy-2-oxo-4-phenyl-2H-chromen-6-yl)ethyl]pyridazin-3(2H)-one (three-letter code: 31T) (formula: $C_{21}H_{16}N_2O_5$).
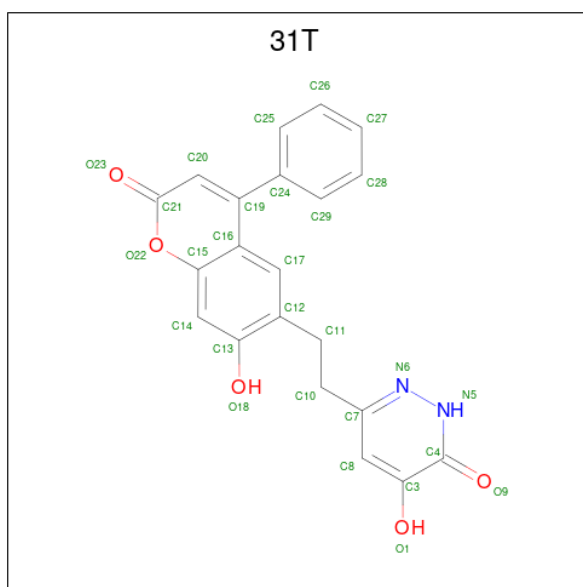
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total 28	C 21	N 2	O 5	0	0
3	B	1	Total 28	C 21	N 2	O 5	0	0

- Molecule 4 is POTASSIUM ION (three-letter code: K) (formula: K).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total 1	K 1	0	0

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	60	Total 60	O 60	0	0
5	B	38	Total 38	O 38	0	0

3 Residue-property plots ⓘ

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.
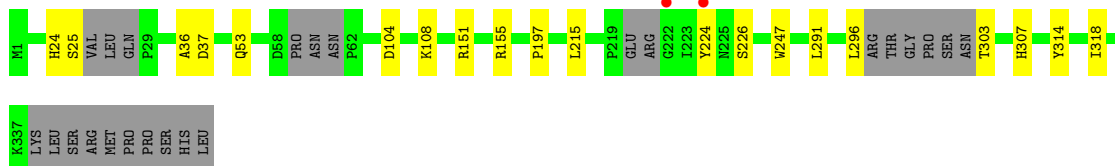
• Molecule 1: D-amino-acid oxidase

Chain A:
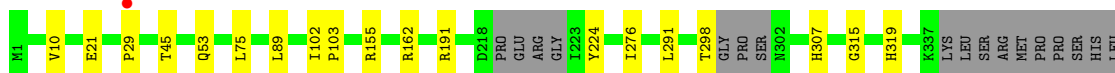
• Molecule 1: D-amino-acid oxidase

Chain B:

4 Data and refinement statistics ⓘ

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	44.39Å 60.93Å 258.45Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	43.79 – 2.40 43.75 – 2.40	Depositor EDS
% Data completeness (in resolution range)	(Not available) (43.79-2.40) 99.7 (43.75-2.40)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.86 (at 2.39Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, R_{free}	0.176 , 0.227 0.180 , 0.228	Depositor DCC
R_{free} test set	1437 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å^2)	34.0	Xtriage
Anisotropy	0.047	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.34 , 37.4	EDS
L-test for twinning [2]	$< |L| > = 0.48, < L^2 > = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	5373	wwPDB-VP
Average B, all atoms (Å^2)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.45% of the height of the origin peak. No significant pseudotranslation is detected.*

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[1]Intensities estimated from amplitudes.
[2]Theoretical values of $< |L| >$, $< L^2 >$ 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: FAD, K, 31T

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	0.69	0/2661	0.83	2/3637 (0.1%)
1	B	0.61	0/2595	0.76	2/3538 (0.1%)
All	All	0.65	0/5256	0.80	4/7175 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	191	ARG	NE-CZ-NH2	-6.99	116.81	120.30
1	A	191	ARG	NE-CZ-NH1	6.90	123.75	120.30
1	B	37	ASP	CB-CG-OD1	5.16	122.94	118.30
1	B	155	ARG	NE-CZ-NH2	-5.13	117.74	120.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	29	PRO	Peptide

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2588	0	2453	9	0
1	B	2524	0	2390	13	0
2	A	53	0	31	1	0
2	B	53	0	31	1	0
3	A	28	0	15	2	0
3	B	28	0	14	5	0
4	B	1	0	0	0	0
5	A	60	0	0	0	0
5	B	38	0	0	0	0
All	All	5373	0	4934	22	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 2.

All (22) 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:53:GLN:HB3	3:B:403:31T:O18	1.84	0.77
1:A:21:GLU:OE1	1:A:155:ARG:NH2	2.28	0.67
1:B:224:TYR:O	3:B:403:31T:C27	2.42	0.66
1:B:224:TYR:CD1	3:B:403:31T:H11	2.30	0.66
1:A:291:LEU:HA	1:A:307:HIS:O	1.99	0.63
1:A:102:ILE:HB	1:A:103:PRO:HD2	1.81	0.63
1:B:224:TYR:O	3:B:403:31T:H14	2.05	0.56
1:A:315:GLY:O	1:A:319:HIS:HB3	2.08	0.53
1:B:296:LEU:HD12	1:B:303:THR:CB	2.42	0.50
1:B:215:LEU:HD21	3:B:403:31T:H7	1.97	0.45
1:A:10:VAL:HB	1:A:45:THR:HG21	1.99	0.45
1:B:314:TYR:O	1:B:318:ILE:HG22	2.18	0.43
1:A:53:GLN:HB3	3:A:402:31T:O18	2.19	0.42
1:B:24:HIS:O	1:B:25:SER:C	2.58	0.42
1:A:224:TYR:O	3:A:402:31T:C27	2.68	0.41
1:B:36:ALA:HA	2:B:402:FAD:N3A	2.35	0.41
1:B:151:ARG:HG2	1:B:151:ARG:HH11	1.86	0.41
1:B:197:PRO:HG3	1:B:247:TRP:CE2	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:291:LEU:HA	1:B:307:HIS:O	2.20	0.41
1:A:75:LEU:O	1:A:89:LEU:HD21	2.21	0.41
1:A:162:ARG:O	2:A:401:FAD:H2A	2.22	0.40
1:B:104:ASP:HB3	1:B:108:LYS:HD3	2.04	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	325/347 (94%)	313 (96%)	12 (4%)	0	100	100
1	B	313/347 (90%)	302 (96%)	11 (4%)	0	100	100
All	All	638/694 (92%)	615 (96%)	23 (4%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	266/299 (89%)	264 (99%)	2 (1%)	81	91
1	B	254/299 (85%)	253 (100%)	1 (0%)	91	96
All	All	520/598 (87%)	517 (99%)	3 (1%)	86	94

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	276	ILE
1	A	298	THR
1	B	226	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	78	HIS
1	A	180	ASN
1	B	180	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no monosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 5 ligands modelled in this entry, 1 is monoatomic - leaving 4 for Mogul analysis.

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	A	401	-	53,58,58	1.15	5 (9%)	68,89,89	1.48	13 (19%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	#$\|Z\|$ > 2	Counts	RMSZ	#$\|Z\|$ > 2
3	31T	B	403	-	30,31,31	2.76	9 (30%)	39,44,44	1.75	9 (23%)
3	31T	A	402	-	30,31,31	2.47	7 (23%)	39,44,44	1.59	7 (17%)
2	FAD	B	402	-	53,58,58	1.25	6 (11%)	68,89,89	1.44	13 (19%)

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	A	401	-	-	1/30/50/50	0/6/6/6
3	31T	B	403	-	-	3/9/9/9	0/4/4/4
3	31T	A	402	-	-	1/9/9/9	0/4/4/4
2	FAD	B	402	-	-	1/30/50/50	0/6/6/6

All (27) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	403	31T	C4-N5	6.43	1.45	1.36
3	B	403	31T	C24-C19	-6.40	1.39	1.49
3	B	403	31T	C10-C7	-5.77	1.37	1.49
3	A	402	31T	C10-C7	-5.74	1.37	1.49
3	A	402	31T	C24-C19	-5.72	1.40	1.49
3	B	403	31T	C11-C12	-5.67	1.40	1.51
3	A	402	31T	C20-C19	5.27	1.41	1.35
2	B	402	FAD	C9A-C5X	4.87	1.49	1.41
3	A	402	31T	N5-N6	-4.76	1.26	1.38
3	B	403	31T	N5-N6	-4.63	1.26	1.38
3	A	402	31T	C11-C12	-4.62	1.42	1.51
3	A	402	31T	C4-N5	4.16	1.42	1.36
3	B	403	31T	C16-C19	-3.96	1.39	1.47
2	A	401	FAD	C9A-C5X	3.54	1.47	1.41
3	B	403	31T	C20-C19	3.21	1.39	1.35
3	A	402	31T	C16-C19	-3.20	1.40	1.47
3	B	403	31T	C7-N6	3.13	1.34	1.31
3	B	403	31T	C20-C21	-2.96	1.38	1.44
2	A	401	FAD	O4B-C1B	2.87	1.45	1.41
2	B	402	FAD	C8-C7	2.50	1.47	1.40
2	A	401	FAD	C4-N3	-2.35	1.34	1.38
2	A	401	FAD	C8-C7	2.34	1.46	1.40
2	B	402	FAD	C5A-C4A	2.31	1.47	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	402	FAD	C4X-N5	2.26	1.35	1.30
2	B	402	FAD	C4-N3	-2.19	1.34	1.38
2	A	401	FAD	C5X-N5	-2.13	1.35	1.39
2	B	402	FAD	C5A-N7A	-2.12	1.32	1.39

All (42) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	A	402	31T	C7-N6-N5	4.80	120.91	117.06
3	B	403	31T	C4-N5-N6	-4.39	122.14	127.59
3	B	403	31T	C3-C8-C7	-3.90	119.26	125.19
3	A	402	31T	C4-N5-N6	-3.61	123.11	127.59
3	A	402	31T	C3-C8-C7	-3.49	119.88	125.19
2	A	401	FAD	N3A-C2A-N1A	-3.46	123.26	128.68
2	A	401	FAD	O4B-C1B-C2B	-3.42	101.93	106.93
2	B	402	FAD	N3A-C2A-N1A	-3.36	123.42	128.68
2	B	402	FAD	C4-C4X-N5	3.35	123.00	118.23
3	B	403	31T	O23-C21-C20	-3.32	119.06	125.84
3	B	403	31T	C19-C20-C21	-3.19	120.08	122.96
2	A	401	FAD	C4-C4X-N5	3.10	122.65	118.23
3	B	403	31T	C25-C24-C19	-3.04	117.07	120.98
2	A	401	FAD	C1'-N10-C9A	3.01	125.52	120.51
3	B	403	31T	O22-C21-O23	2.94	120.19	116.44
2	A	401	FAD	O2-C2-N1	-2.92	116.99	121.83
3	A	402	31T	O23-C21-C20	-2.82	120.09	125.84
2	A	401	FAD	O2'-C2'-C1'	-2.80	103.03	109.80
3	B	403	31T	O22-C21-C20	2.78	120.75	117.17
3	B	403	31T	C7-N6-N5	2.76	119.27	117.06
2	B	402	FAD	C4A-C5A-N7A	-2.73	106.55	109.40
3	B	403	31T	O9-C4-C3	-2.71	118.06	123.55
2	B	402	FAD	C4X-C4-N3	2.59	119.77	113.19
2	B	402	FAD	C4X-C10-N10	2.57	120.24	116.48
2	B	402	FAD	C2A-N1A-C6A	2.57	123.15	118.75
3	A	402	31T	O22-C21-O23	2.53	119.66	116.44
2	A	401	FAD	O4-C4-C4X	-2.47	120.05	126.60
2	A	401	FAD	C5X-C9A-N10	2.44	120.47	117.95
2	B	402	FAD	O4-C4-C4X	-2.41	120.19	126.60
3	A	402	31T	O22-C21-C20	2.39	120.25	117.17
2	A	401	FAD	C4A-C5A-N7A	-2.38	106.92	109.40
2	B	402	FAD	C10-N1-C2	2.37	121.64	116.90
2	A	401	FAD	C9A-C5X-N5	-2.31	119.92	122.43
3	A	402	31T	C19-C20-C21	-2.29	120.89	122.96

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	402	FAD	C5X-C9A-N10	2.26	120.29	117.95
2	B	402	FAD	O2-C2-N1	-2.25	118.10	121.83
2	A	401	FAD	C4X-C10-N10	2.25	119.77	116.48
2	B	402	FAD	C4X-C10-N1	-2.25	119.52	124.73
2	B	402	FAD	O2P-P-O1P	2.19	123.08	112.24
2	A	401	FAD	C9A-N10-C10	-2.12	117.47	120.77
2	B	402	FAD	C9A-N10-C10	-2.07	117.54	120.77
2	A	401	FAD	C4X-C4-N3	2.03	118.33	113.19

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	402	31T	C7-C10-C11-C12
3	B	403	31T	C7-C10-C11-C12
2	B	402	FAD	O4B-C4B-C5B-O5B
3	B	403	31T	C16-C19-C24-C25
2	A	401	FAD	O4B-C4B-C5B-O5B
3	B	403	31T	C20-C19-C24-C25

There are no ring outliers.

4 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	401	FAD	1	0
3	B	403	31T	5	0
3	A	402	31T	2	0
2	B	402	FAD	1	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 then 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 FAD A 401

Bond lengths

Bond angles

Torsions

Rings

Ligand 31T B 403

Bond lengths

Bond angles

Torsions

Rings

Ligand 31T A 402

Bond lengths

Bond angles

Torsions

Rings

Ligand FAD B 402

Bond lengths

Bond angles

Torsions

Rings
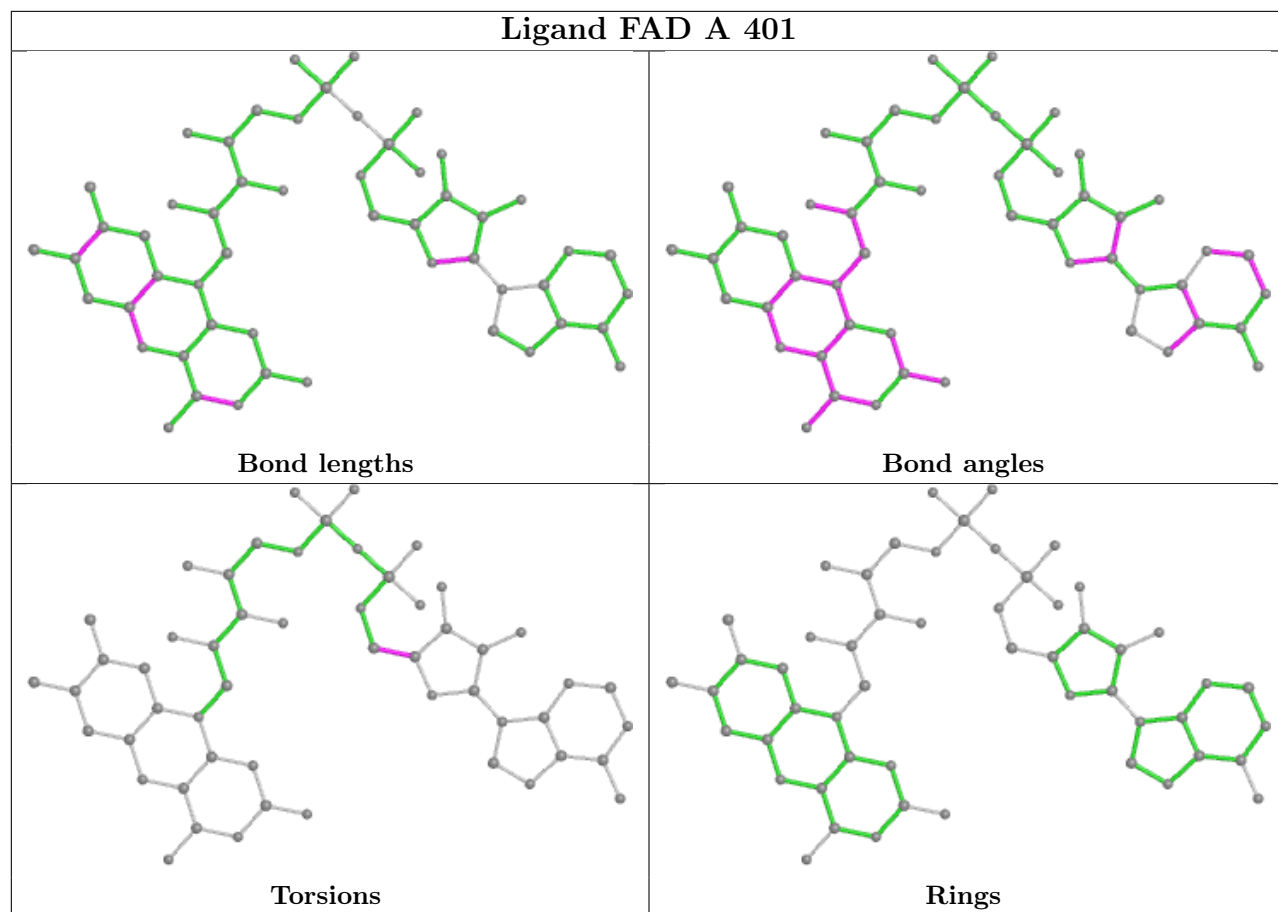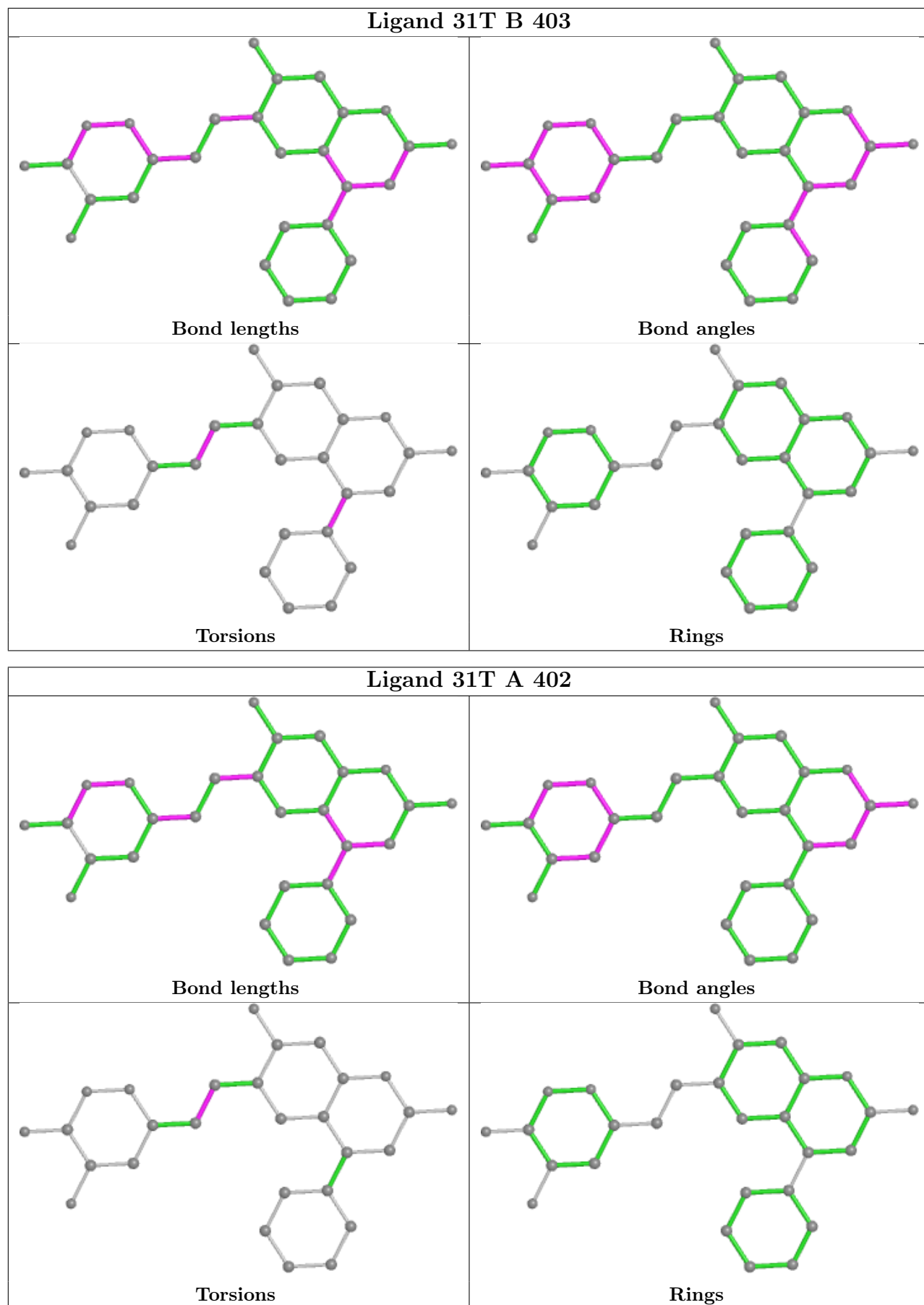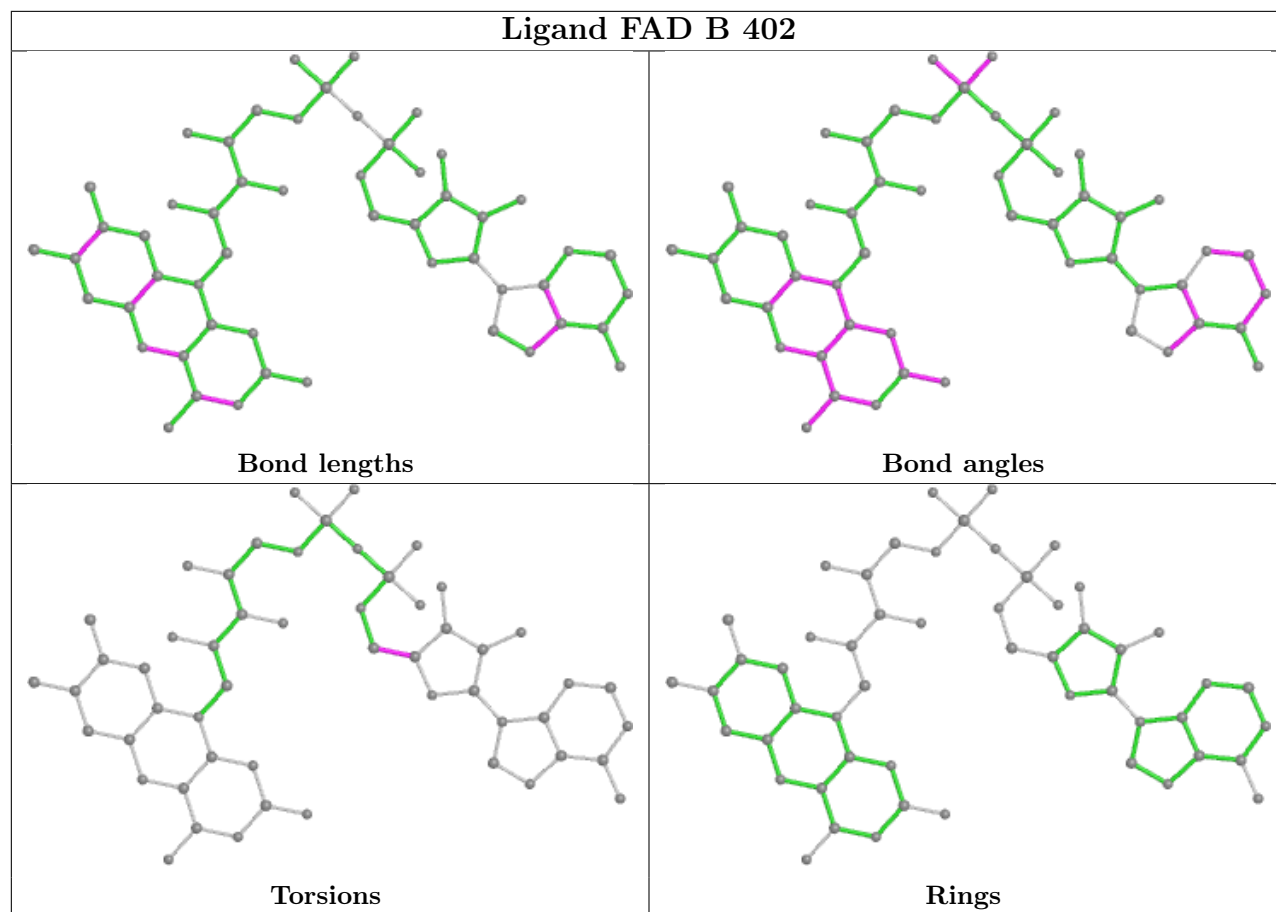
5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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	330/347 (95%)	-0.47	1 (0%)	94	93	17, 28, 53, 73	0
1	B	323/347 (93%)	-0.50	2 (0%)	89	88	22, 36, 58, 83	0
All	All	653/694 (94%)	-0.49	3 (0%)	91	89	17, 32, 55, 83	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	222	GLY	3.3
1	B	224	TYR	2.4
1	A	29	PRO	2.2

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

6.3 Carbohydrates ⓘ

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	31T	B	403	28/28	0.86	0.25	32,38,42,44	13
3	31T	A	402	28/28	0.93	0.15	21,35,38,41	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	FAD	B	402	53/53	0.97	0.10	25,29,31,34	0
2	FAD	A	401	53/53	0.98	0.10	15,17,20,21	0
4	K	B	401	1/1	0.98	0.03	37,37,37,37	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 31T B 403:

2mF$_o$-DF$_c$ (at 0.7 rmsd) in gray
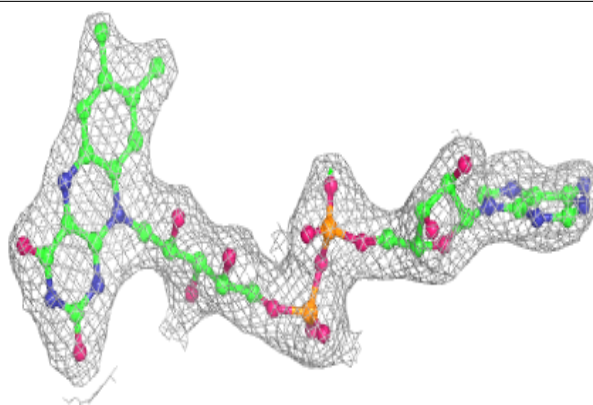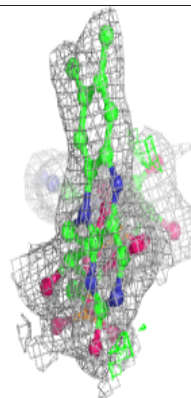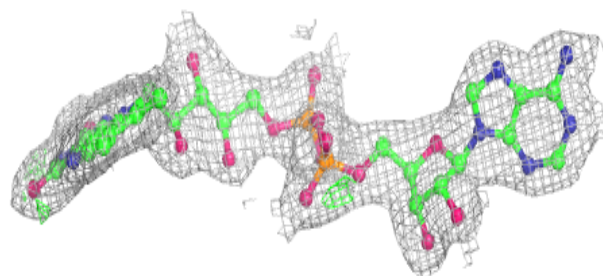mF$_o$-DF$_c$ (at 3 rmsd) in purple (negative)
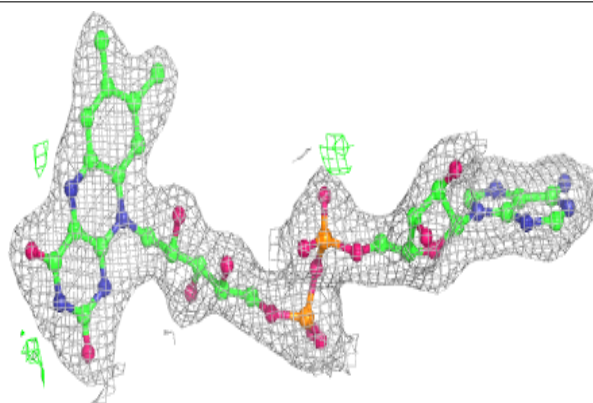and green (positive)
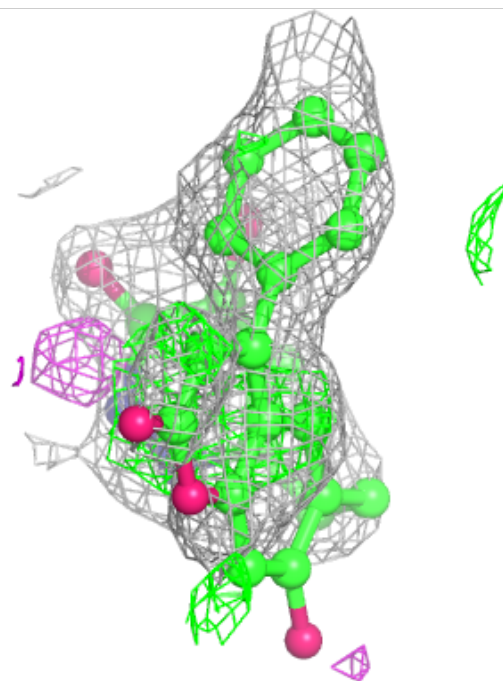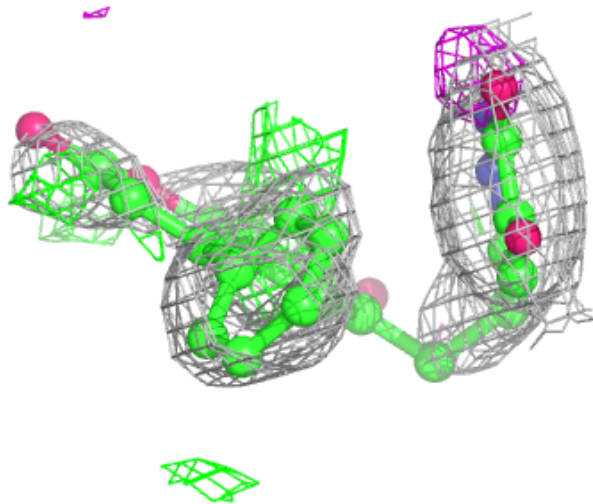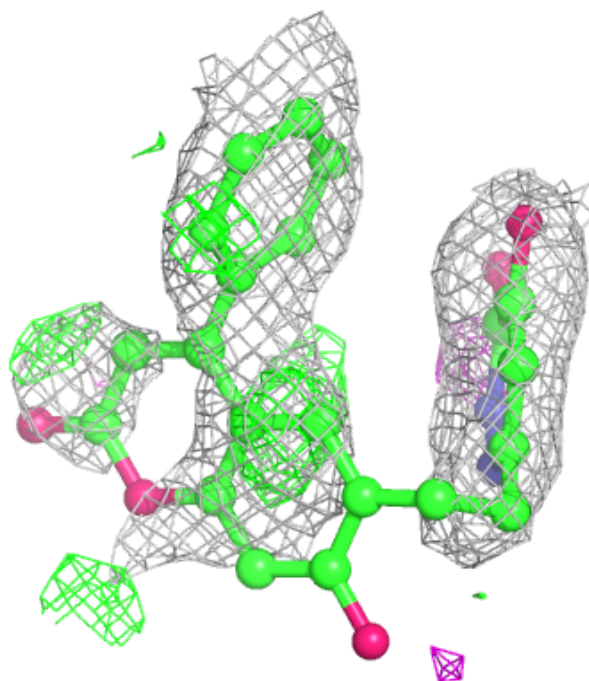
Electron density around 31T A 402:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)
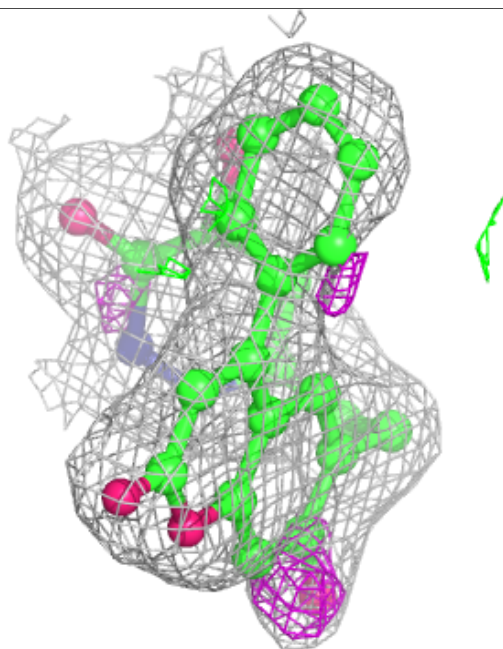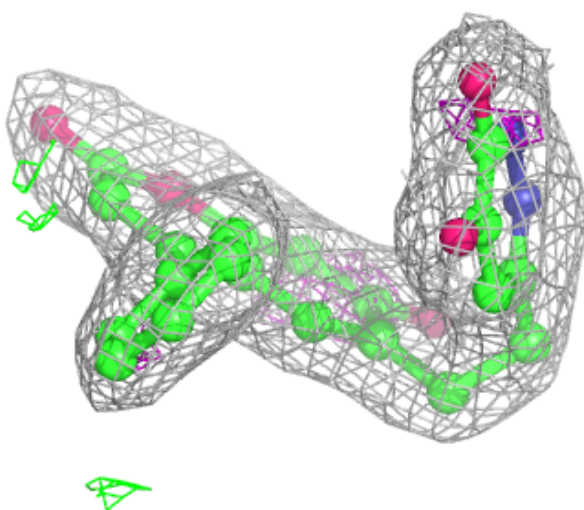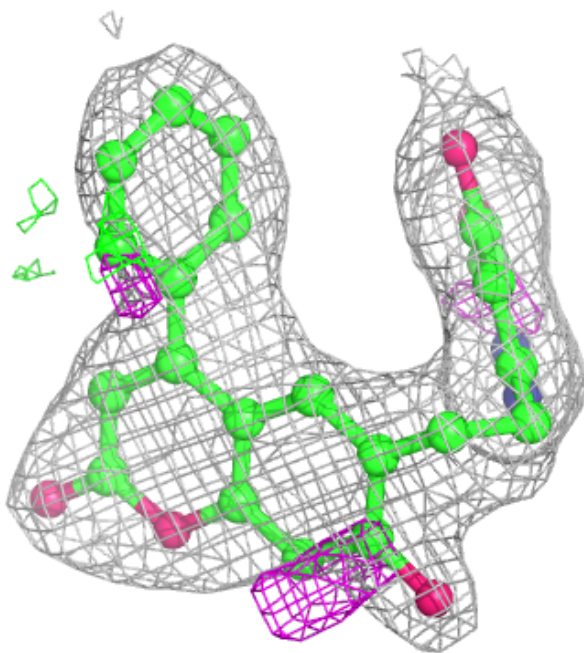
Electron density around FAD B 402:

$2mF_o\text{-}DF_c$ (at 0.7 rmsd) in gray
$mF_o\text{-}DF_c$ (at 3 rmsd) in purple (negative)
and green (positive)
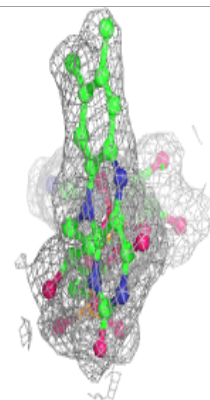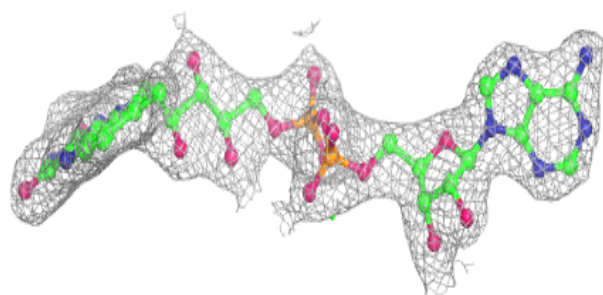
Electron density around FAD A 401:

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

6.5 Other polymers ⓘ

There are no such residues in this entry.