



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 8, 2026 – 06:10 AM UTC

PDB ID : 5MOG / pdb_00005mog
Title : Oryza sativa phytoene desaturase inhibited by norflurazon
Authors : Brausemann, A.; Gemmecker, S.; Koschmieder, J.; Beyer, P.; Einsle, O.
Deposited on : 2016-12-14
Resolution : 2.77 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	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

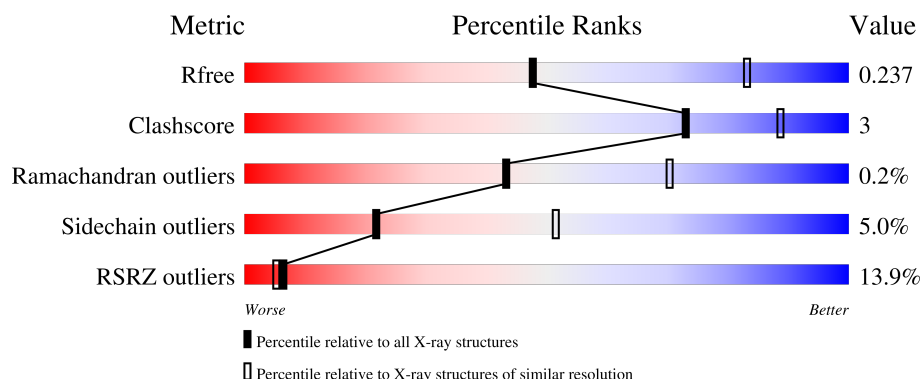
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.77 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}	180053	5248 (2.80-2.76)
Clashscore	190562	5693 (2.80-2.76)
Ramachandran outliers	187476	5590 (2.80-2.76)
Sidechain outliers	187428	5592 (2.80-2.76)
RSRZ outliers	180081	5251 (2.80-2.76)

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	497	3% 82% 11% • 6%
1	B	497	3% 83% 11% • 5%
1	C	497	5% 83% 11% • 5%
1	D	497	10% 84% 10% 5%
1	E	497	44% 79% 13% • 6%

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
5	IMD	D	603	-	-	-	X

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 19520 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 Phytoene dehydrogenase, chloroplastic/chromoplastic.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	467	Total	C	N	O	S	0	0	0
			3703	2388	618	677	20			
1	B	472	Total	C	N	O	S	0	1	0
			3757	2419	631	687	20			
1	C	470	Total	C	N	O	S	0	1	0
			3736	2408	626	682	20			
1	D	470	Total	C	N	O	S	0	0	0
			3728	2403	623	682	20			
1	E	466	Total	C	N	O	S	0	0	0
			3692	2380	616	676	20			

There are 30 discrepancies between the modelled and reference sequences:

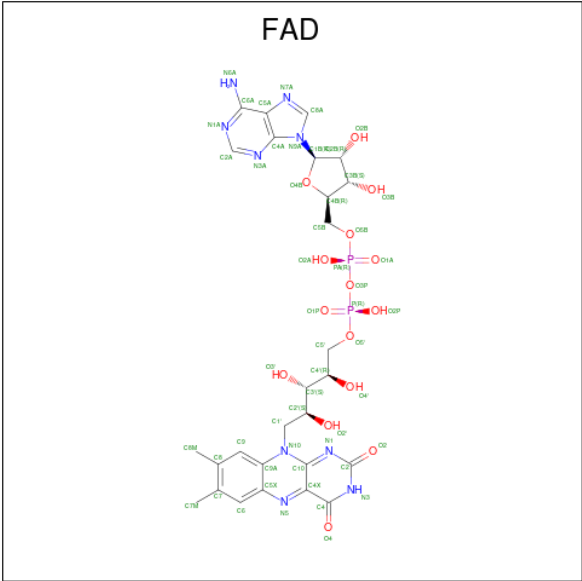
Chain	Residue	Modelled	Actual	Comment	Reference
A	579	HIS	-	expression tag	UNP A2XDA1
A	580	HIS	-	expression tag	UNP A2XDA1
A	581	HIS	-	expression tag	UNP A2XDA1
A	582	HIS	-	expression tag	UNP A2XDA1
A	583	HIS	-	expression tag	UNP A2XDA1
A	584	HIS	-	expression tag	UNP A2XDA1
B	579	HIS	-	expression tag	UNP A2XDA1
B	580	HIS	-	expression tag	UNP A2XDA1
B	581	HIS	-	expression tag	UNP A2XDA1
B	582	HIS	-	expression tag	UNP A2XDA1
B	583	HIS	-	expression tag	UNP A2XDA1
B	584	HIS	-	expression tag	UNP A2XDA1
C	579	HIS	-	expression tag	UNP A2XDA1
C	580	HIS	-	expression tag	UNP A2XDA1
C	581	HIS	-	expression tag	UNP A2XDA1
C	582	HIS	-	expression tag	UNP A2XDA1
C	583	HIS	-	expression tag	UNP A2XDA1
C	584	HIS	-	expression tag	UNP A2XDA1
D	579	HIS	-	expression tag	UNP A2XDA1

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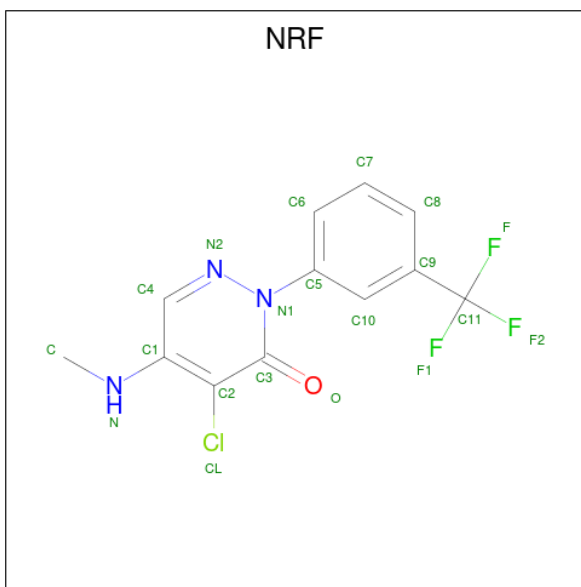
Chain	Residue	Modelled	Actual	Comment	Reference
D	580	HIS	-	expression tag	UNP A2XDA1
D	581	HIS	-	expression tag	UNP A2XDA1
D	582	HIS	-	expression tag	UNP A2XDA1
D	583	HIS	-	expression tag	UNP A2XDA1
D	584	HIS	-	expression tag	UNP A2XDA1
E	579	HIS	-	expression tag	UNP A2XDA1
E	580	HIS	-	expression tag	UNP A2XDA1
E	581	HIS	-	expression tag	UNP A2XDA1
E	582	HIS	-	expression tag	UNP A2XDA1
E	583	HIS	-	expression tag	UNP A2XDA1
E	584	HIS	-	expression tag	UNP A2XDA1

- Molecule 2 is FLAVIN-ADENINE DINUCLEOTIDE (CCD ID: FAD) (formula: C₂₇H₃₃N₉O₁₅P₂).



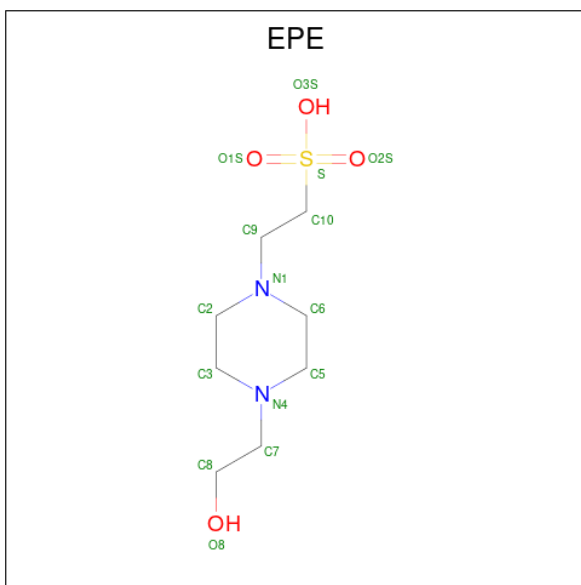
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
2	B	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
2	C	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
2	D	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
2	E	1	Total	C	N	O	P	0	0
			53	27	9	15	2		

- Molecule 3 is Norflurazon (CCD ID: NRF) (formula: C₁₂H₉ClF₃N₃O).



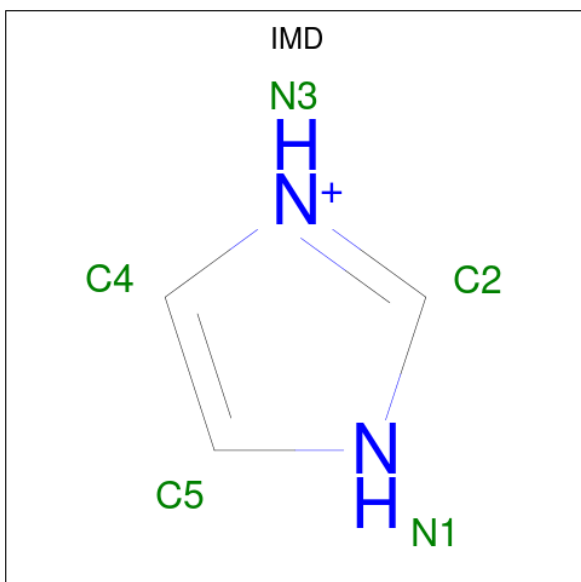
Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
3	A	1	Total	C	Cl	F	N	O	0	0
			20	12	1	3	3	1		
3	B	1	Total	C	Cl	F	N	O	0	0
			20	12	1	3	3	1		
3	C	1	Total	C	Cl	F	N	O	0	0
			20	12	1	3	3	1		
3	D	1	Total	C	Cl	F	N	O	0	0
			20	12	1	3	3	1		
3	E	1	Total	C	Cl	F	N	O	0	0
			20	12	1	3	3	1		

- Molecule 4 is 4-(2-HYDROXYETHYL)-1-PIPERAZINE ETHANESULFONIC ACID (CCD ID: EPE) (formula: C₈H₁₈N₂O₄S).



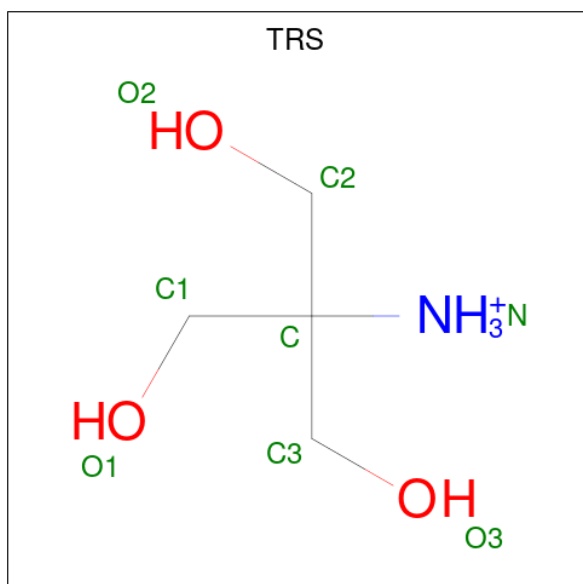
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	S	0	0
			15	8	2	4	1		
4	A	1	Total	C	N	O	S	0	0
			15	8	2	4	1		
4	B	1	Total	C	N	O	S	0	0
			15	8	2	4	1		
4	B	1	Total	C	N	O	S	0	0
			15	8	2	4	1		
4	E	1	Total	C	N	O	S	0	0
			15	8	2	4	1		

- Molecule 5 is IMIDAZOLE (CCD ID: IMD) (formula: $C_3H_5N_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	N	0	0
			5	3	2		
5	B	1	Total	C	N	0	0
			5	3	2		
5	B	1	Total	C	N	0	0
			5	3	2		
5	D	1	Total	C	N	0	0
			5	3	2		
5	D	1	Total	C	N	0	0
			5	3	2		
5	D	1	Total	C	N	0	0
			5	3	2		

- Molecule 6 is 2-AMINO-2-HYDROXYMETHYL-PROPANE-1,3-DIOL (CCD ID: TRS) (formula: $C_4H_{12}NO_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	B	1	Total	C	N	O	0	0
			8	4	1	3		
6	D	1	Total	C	N	O	0	0
			8	4	1	3		

- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	144	Total	O	0	0
			144	144		

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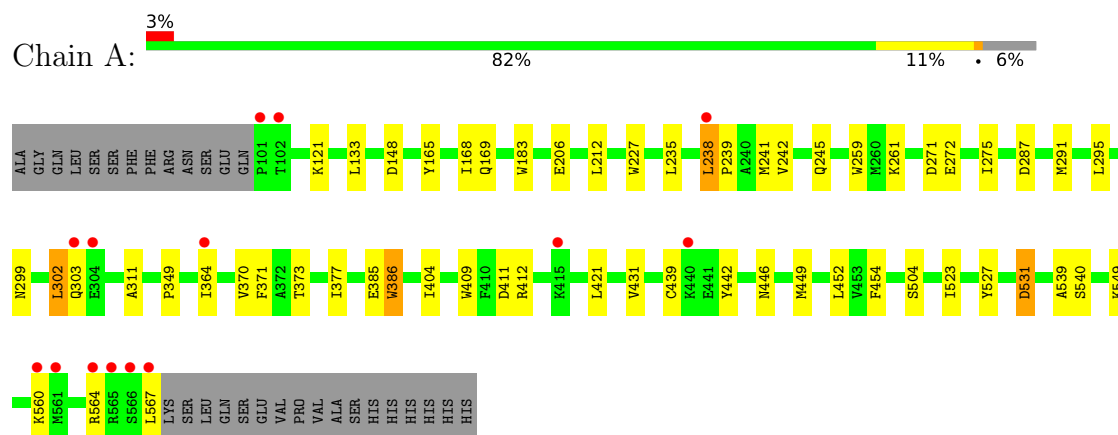
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	B	139	Total 139	O 139	0	0
7	C	62	Total 62	O 62	0	0
7	D	62	Total 62	O 62	0	0
7	E	11	Total 11	O 11	0	0

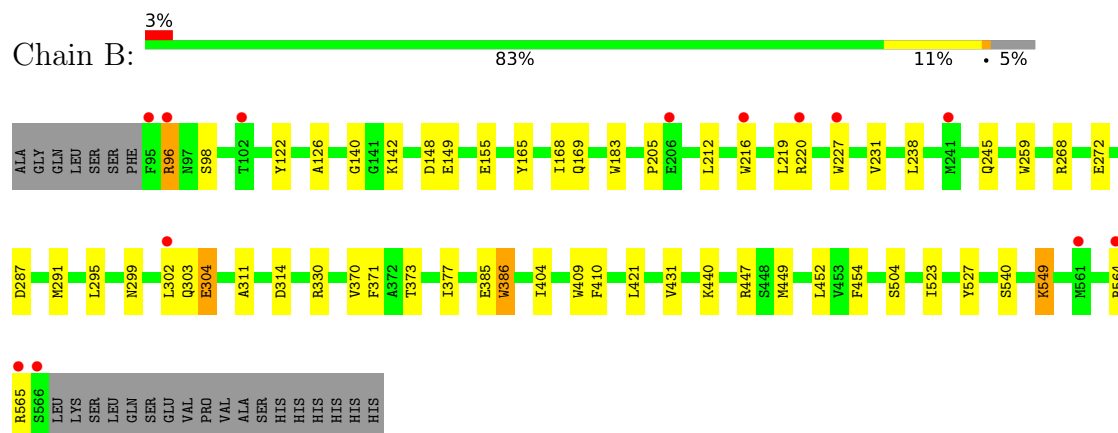
3 Residue-property plots [i](#)

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.

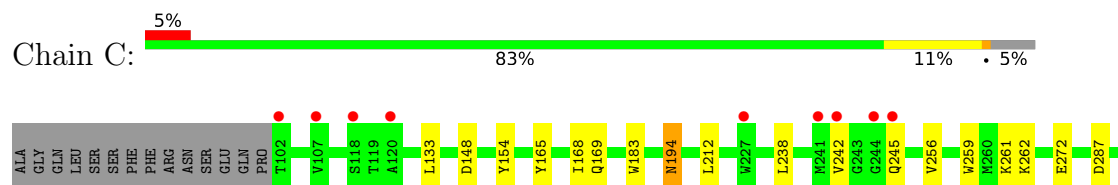
- Molecule 1: Phytoene dehydrogenase, chloroplastic/chromoplastic

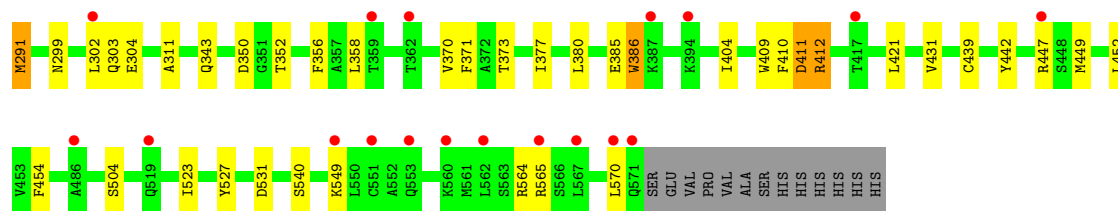


- Molecule 1: Phytoene dehydrogenase, chloroplastic/chromoplastic

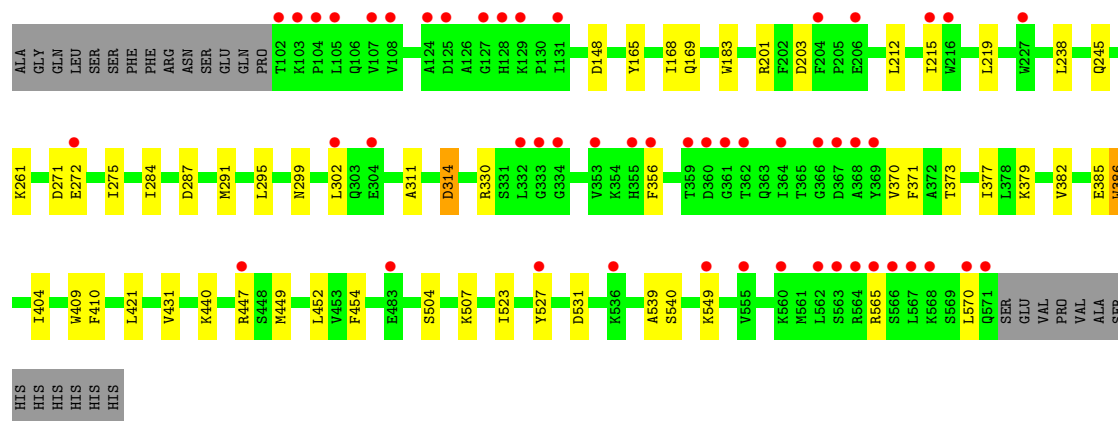
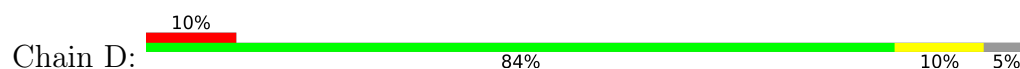


- Molecule 1: Phytoene dehydrogenase, chloroplastic/chromoplastic

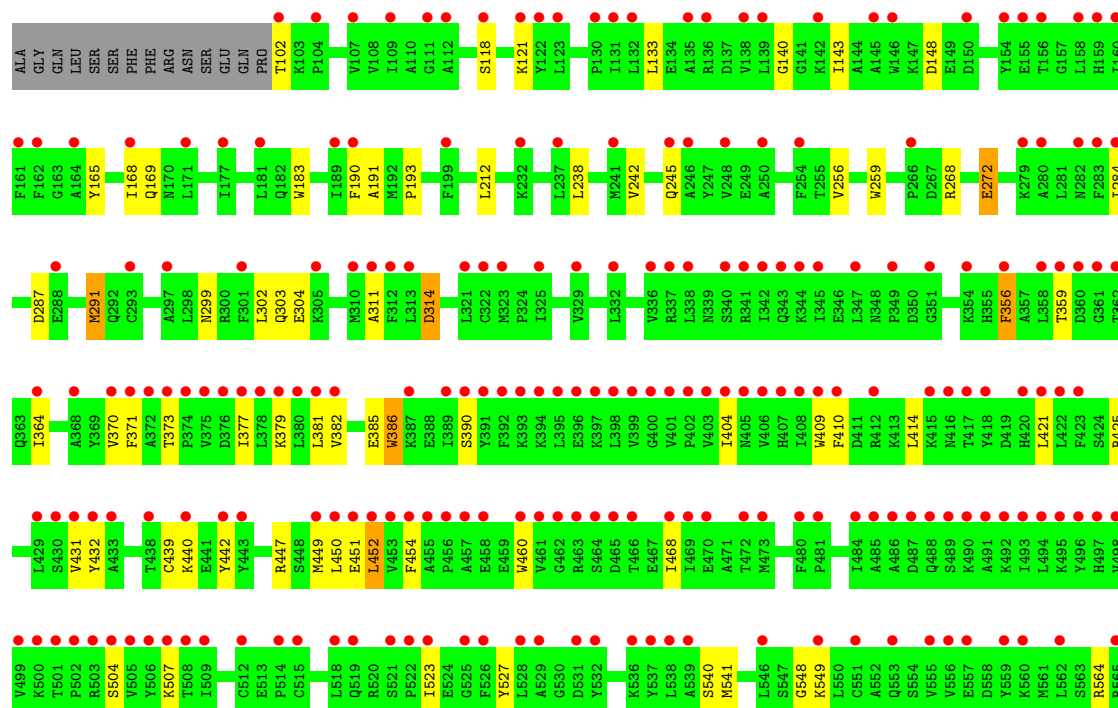
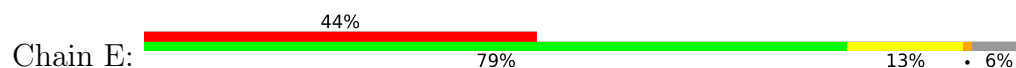




• Molecule 1: Phytoene dehydrogenase, chloroplatic/chromoplatic



• Molecule 1: Phytoene dehydrogenase, chloroplatic/chromoplatic



5566

L567

LYS	SER	LEU	GLN	SER	GLU	VAL	PRO	VAL	ALA	SER	HIS	HIS	HIS	HIS	HIS
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4 Data and refinement statistics

Property	Value	Source
Space group	I 4	Depositor
Cell constants a, b, c, α , β , γ	230.21Å 230.21Å 179.23Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	60.25 – 2.77 60.25 – 2.77	Depositor EDS
% Data completeness (in resolution range)	100.0 (60.25-2.77) 100.0 (60.25-2.77)	Depositor EDS
R_{merge}	0.17	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.78 (at 2.77Å)	Xtriage
Refinement program	BUSTER	Depositor
R, R_{free}	0.209 , 0.224 0.217 , 0.237	Depositor DCC
R_{free} test set	5852 reflections (4.95%)	wwPDB-VP
Wilson B-factor (Å ²)	66.6	Xtriage
Anisotropy	0.147	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 37.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.017 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	19520	wwPDB-VP
Average B, all atoms (Å ²)	79.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.94% 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: NRF, TRS, EPE, IMD, 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	0.75	0/3794	1.23	8/5141 (0.2%)
1	B	0.76	0/3852	1.23	5/5218 (0.1%)
1	C	0.76	1/3829 (0.0%)	1.22	10/5186 (0.2%)
1	D	0.75	0/3818	1.23	6/5172 (0.1%)
1	E	0.77	1/3782 (0.0%)	1.23	11/5126 (0.2%)
All	All	0.76	2/19075 (0.0%)	1.23	40/25843 (0.2%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	291	MET	SD-CE	-8.54	1.58	1.79
1	C	291	MET	SD-CE	-7.99	1.59	1.79

The worst 5 of 40 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	194	ASN	CA-CB-CG	6.86	119.46	112.60
1	B	148	ASP	CA-CB-CG	6.14	118.74	112.60
1	D	148	ASP	CA-CB-CG	6.02	118.62	112.60
1	E	148	ASP	CA-CB-CG	6.00	118.60	112.60
1	C	148	ASP	CA-CB-CG	5.93	118.53	112.60

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	3703	0	3718	23	0
1	B	3757	0	3766	30	0
1	C	3736	0	3760	20	0
1	D	3728	0	3747	19	0
1	E	3692	0	3699	37	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
3	A	20	0	0	1	0
3	B	20	0	0	1	0
3	C	20	0	0	1	0
3	D	20	0	0	1	0
3	E	20	0	0	1	0
4	A	30	0	34	1	0
4	B	30	0	34	0	0
4	E	15	0	17	1	0
5	A	5	0	5	0	0
5	B	10	0	10	0	0
5	D	15	0	15	0	0
6	B	8	0	12	3	0
6	D	8	0	12	1	0
7	A	144	0	0	0	0
7	B	139	0	0	2	0
7	C	62	0	0	0	0
7	D	62	0	0	0	0
7	E	11	0	0	0	0
All	All	19520	0	18984	122	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 3.

The worst 5 of 122 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:432:TYR:HB2	1:E:452:LEU:HD22	1.46	0.96
1:E:256:VAL:HA	1:E:291:MET:HE2	1.49	0.94
1:C:256:VAL:HA	1:C:291:MET:HE2	1.49	0.91

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:259:TRP:HB3	1:C:291:MET:HE3	1.56	0.87
1:E:259:TRP:HB3	1:E:291:MET:HE3	1.57	0.86

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	465/497 (94%)	449 (97%)	15 (3%)	1 (0%)	43	70
1	B	471/497 (95%)	455 (97%)	15 (3%)	1 (0%)	43	70
1	C	469/497 (94%)	454 (97%)	14 (3%)	1 (0%)	43	70
1	D	468/497 (94%)	452 (97%)	15 (3%)	1 (0%)	43	70
1	E	464/497 (93%)	448 (97%)	15 (3%)	1 (0%)	43	70
All	All	2337/2485 (94%)	2258 (97%)	74 (3%)	5 (0%)	43	70

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	504	SER
1	B	504	SER
1	C	504	SER
1	E	504	SER
1	D	504	SER

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	400/427 (94%)	379 (95%)	21 (5%)	20	49
1	B	406/427 (95%)	385 (95%)	21 (5%)	21	50
1	C	404/427 (95%)	382 (95%)	22 (5%)	20	48
1	D	403/427 (94%)	387 (96%)	16 (4%)	28	60
1	E	398/427 (93%)	377 (95%)	21 (5%)	20	49
All	All	2011/2135 (94%)	1910 (95%)	101 (5%)	22	51

5 of 101 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	377	ILE
1	D	302	LEU
1	E	549	LYS
1	C	412	ARG
1	D	212	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 31 such sidechains are listed below:

Mol	Chain	Res	Type
1	C	245	GLN
1	E	285	ASN
1	C	355	HIS
1	E	348	ASN
1	D	355	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

23 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$
5	IMD	D	605	-	5,5,5	0.32	0	5,5,5	0.55	0
4	EPE	B	603	-	15,15,15	1.16	1 (6%)	19,20,20	0.34	0
2	FAD	D	600	-	58,58,58	0.41	0	85,89,89	0.60	1 (1%)
5	IMD	B	606	-	5,5,5	0.31	0	5,5,5	0.54	0
6	TRS	D	602	-	7,7,7	0.28	0	9,9,9	0.30	0
5	IMD	A	604	-	5,5,5	0.31	0	5,5,5	0.54	0
6	TRS	B	604	-	7,7,7	0.29	0	9,9,9	0.37	0
4	EPE	B	602	-	15,15,15	1.29	1 (6%)	19,20,20	0.45	0
3	NRF	E	601	-	19,21,21	1.19	2 (10%)	27,31,31	1.05	3 (11%)
3	NRF	B	601	-	19,21,21	1.18	2 (10%)	27,31,31	0.97	3 (11%)
4	EPE	A	602	-	15,15,15	1.08	1 (6%)	19,20,20	0.26	0
3	NRF	A	601	-	19,21,21	1.17	2 (10%)	27,31,31	1.02	3 (11%)
2	FAD	A	600	-	58,58,58	0.40	0	85,89,89	0.50	0
5	IMD	B	605	-	5,5,5	0.32	0	5,5,5	0.55	0
4	EPE	E	602	-	15,15,15	1.10	1 (6%)	19,20,20	0.47	0
2	FAD	B	600	-	58,58,58	0.42	0	85,89,89	0.56	0
2	FAD	C	600	-	58,58,58	0.43	0	85,89,89	0.59	0
4	EPE	A	603	-	15,15,15	0.66	1 (6%)	19,20,20	0.33	0
3	NRF	D	601	-	19,21,21	1.11	2 (10%)	27,31,31	1.03	4 (14%)
3	NRF	C	601	-	19,21,21	1.18	2 (10%)	27,31,31	0.98	3 (11%)
5	IMD	D	604	-	5,5,5	0.31	0	5,5,5	0.55	0
2	FAD	E	600	-	58,58,58	0.40	0	85,89,89	0.55	0
5	IMD	D	603	-	5,5,5	0.32	0	5,5,5	0.55	0

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
5	IMD	D	605	-	-	-	0/1/1/1
4	EPE	B	603	-	-	1/9/19/19	0/1/1/1
2	FAD	D	600	-	-	7/34/50/50	0/6/6/6
5	IMD	B	606	-	-	-	0/1/1/1
6	TRS	D	602	-	-	0/9/9/9	-
5	IMD	A	604	-	-	-	0/1/1/1
6	TRS	B	604	-	-	1/9/9/9	-
4	EPE	B	602	-	-	5/9/19/19	0/1/1/1
3	NRF	E	601	-	-	0/11/12/12	0/2/2/2
3	NRF	B	601	-	-	0/11/12/12	0/2/2/2
4	EPE	A	602	-	-	8/9/19/19	0/1/1/1
3	NRF	A	601	-	-	0/11/12/12	0/2/2/2
2	FAD	A	600	-	-	7/34/50/50	0/6/6/6
5	IMD	B	605	-	-	-	0/1/1/1
4	EPE	E	602	-	-	5/9/19/19	0/1/1/1
2	FAD	B	600	-	-	7/34/50/50	0/6/6/6
2	FAD	C	600	-	-	7/34/50/50	0/6/6/6
4	EPE	A	603	-	-	4/9/19/19	0/1/1/1
3	NRF	D	601	-	-	0/11/12/12	0/2/2/2
3	NRF	C	601	-	-	0/11/12/12	0/2/2/2
5	IMD	D	604	-	-	-	0/1/1/1
2	FAD	E	600	-	-	7/34/50/50	0/6/6/6
5	IMD	D	603	-	-	-	0/1/1/1

The worst 5 of 15 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	602	EPE	C10-S	-4.86	1.70	1.77
4	B	603	EPE	C10-S	-4.45	1.71	1.77
4	A	602	EPE	C10-S	-4.15	1.71	1.77
4	E	602	EPE	C10-S	-4.13	1.71	1.77
3	C	601	NRF	C4-C1	3.90	1.48	1.43

The worst 5 of 17 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	E	601	NRF	C2-C1-N	-3.20	122.60	129.89
3	A	601	NRF	C2-C1-N	-3.18	122.64	129.89
3	B	601	NRF	C2-C1-N	-3.17	122.67	129.89
3	D	601	NRF	C2-C1-N	-3.16	122.68	129.89
3	C	601	NRF	C2-C1-N	-3.16	122.69	129.89

There are no chirality outliers.

5 of 59 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	600	FAD	C5'-O5'-P-O1P
2	A	600	FAD	C5'-O5'-P-O3P
2	B	600	FAD	C5'-O5'-P-O1P
2	B	600	FAD	C5'-O5'-P-O3P
2	C	600	FAD	C5'-O5'-P-O2P

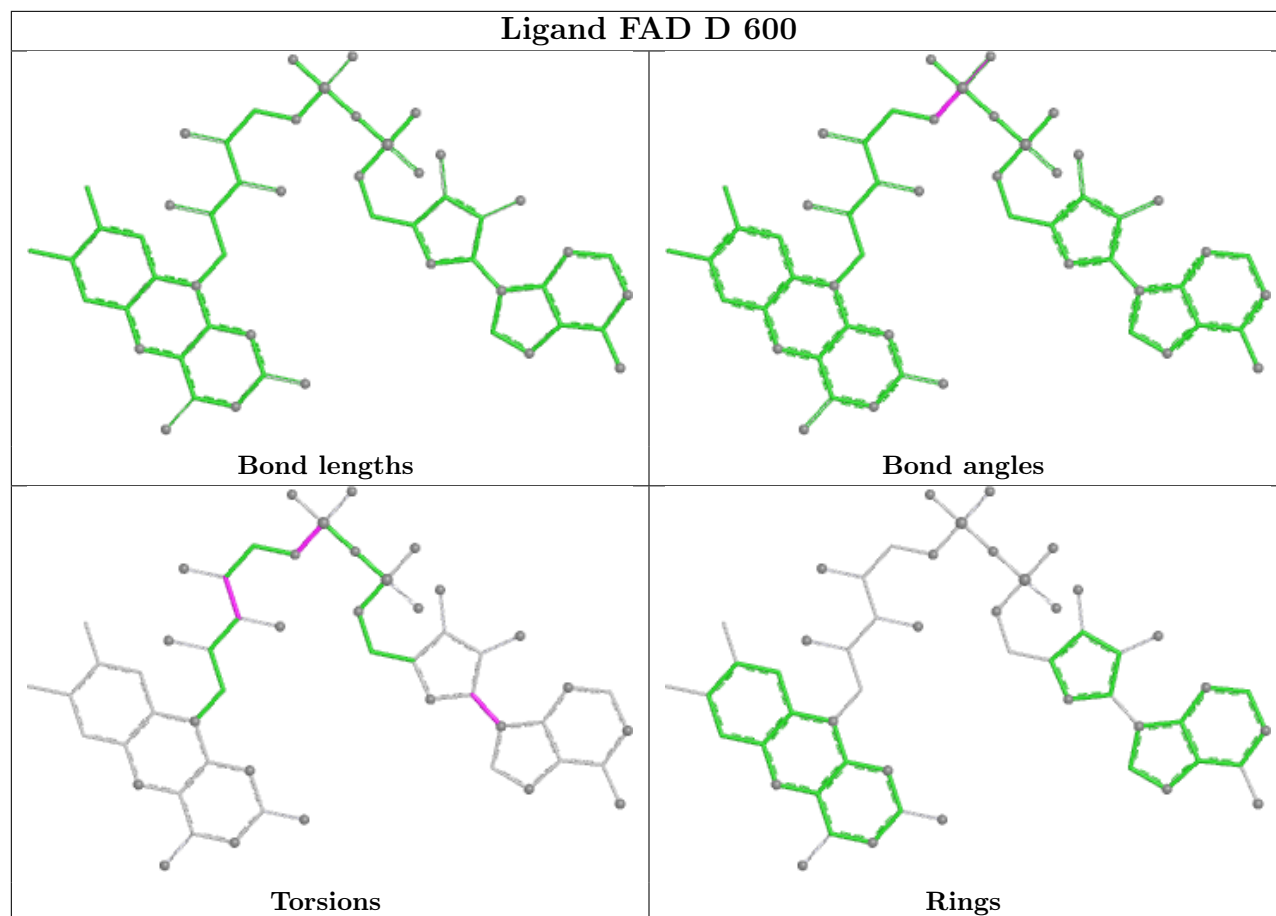
There are no ring outliers.

9 monomers are involved in 11 short contacts:

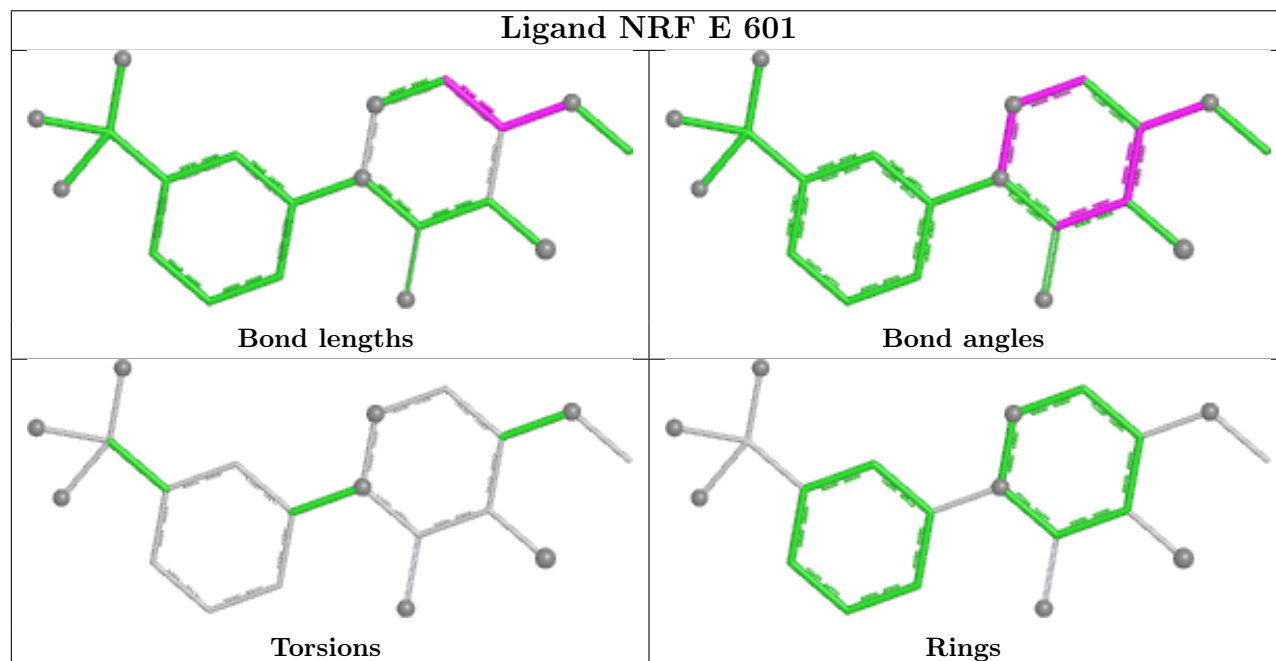
Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	D	602	TRS	1	0
6	B	604	TRS	3	0
3	E	601	NRF	1	0
3	B	601	NRF	1	0
3	A	601	NRF	1	0
4	E	602	EPE	1	0
4	A	603	EPE	1	0
3	D	601	NRF	1	0
3	C	601	NRF	1	0

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 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.

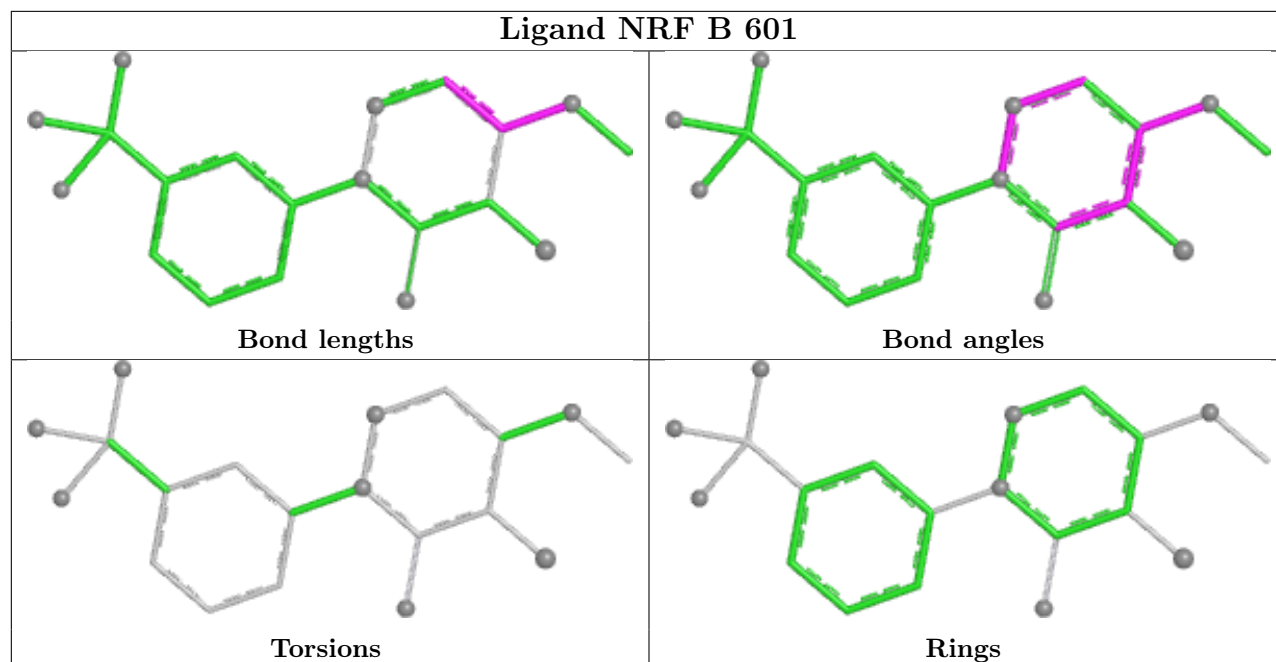
Ligand FAD D 600



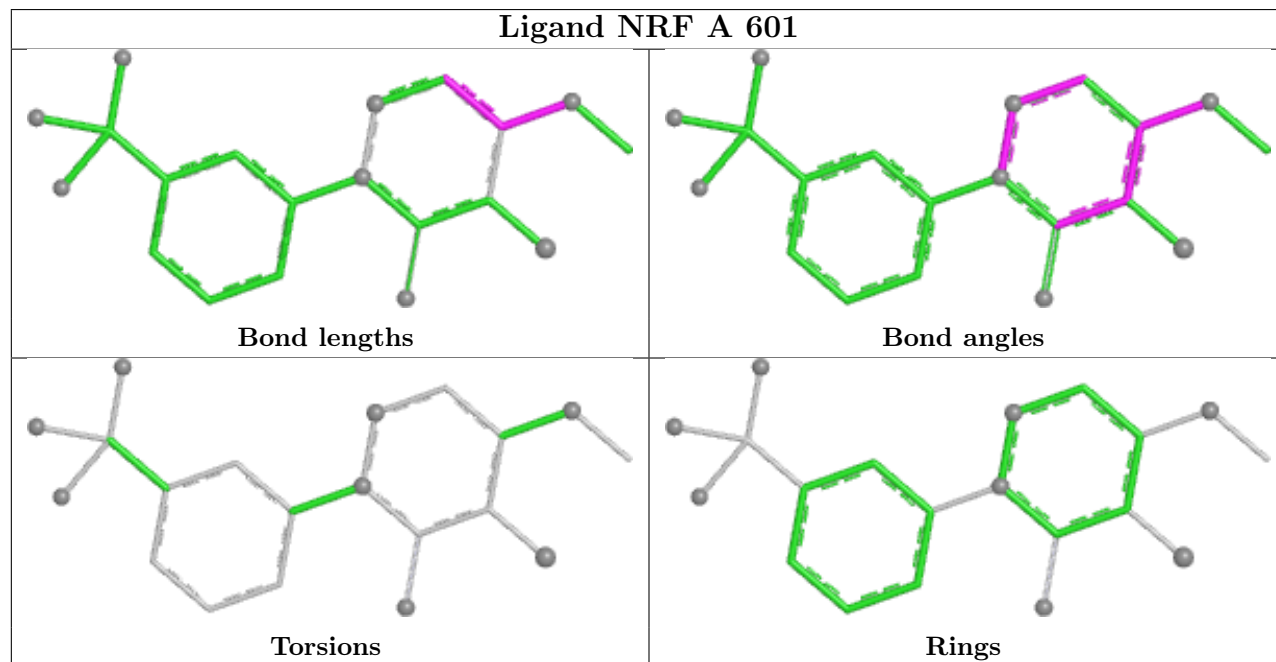
Ligand NRF E 601

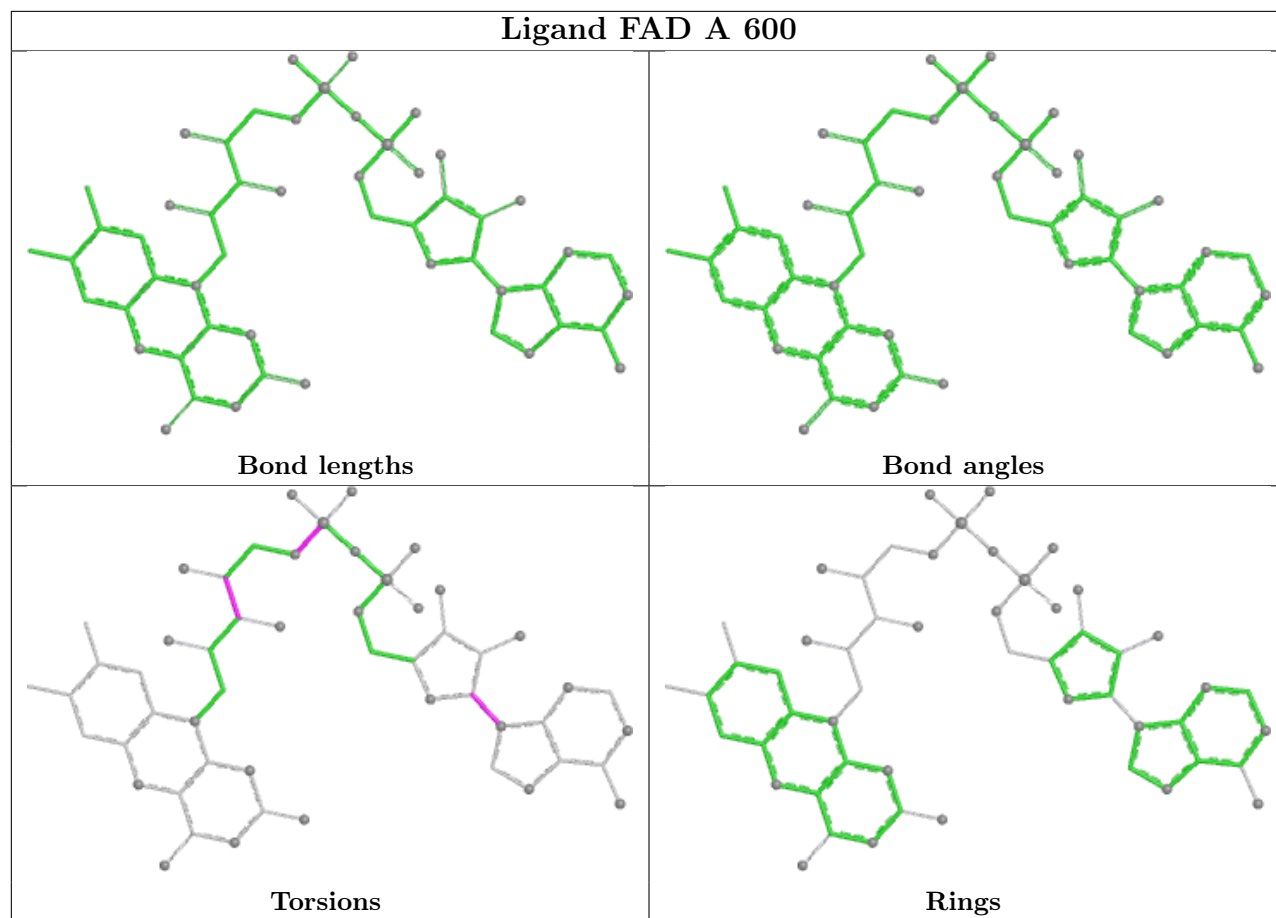


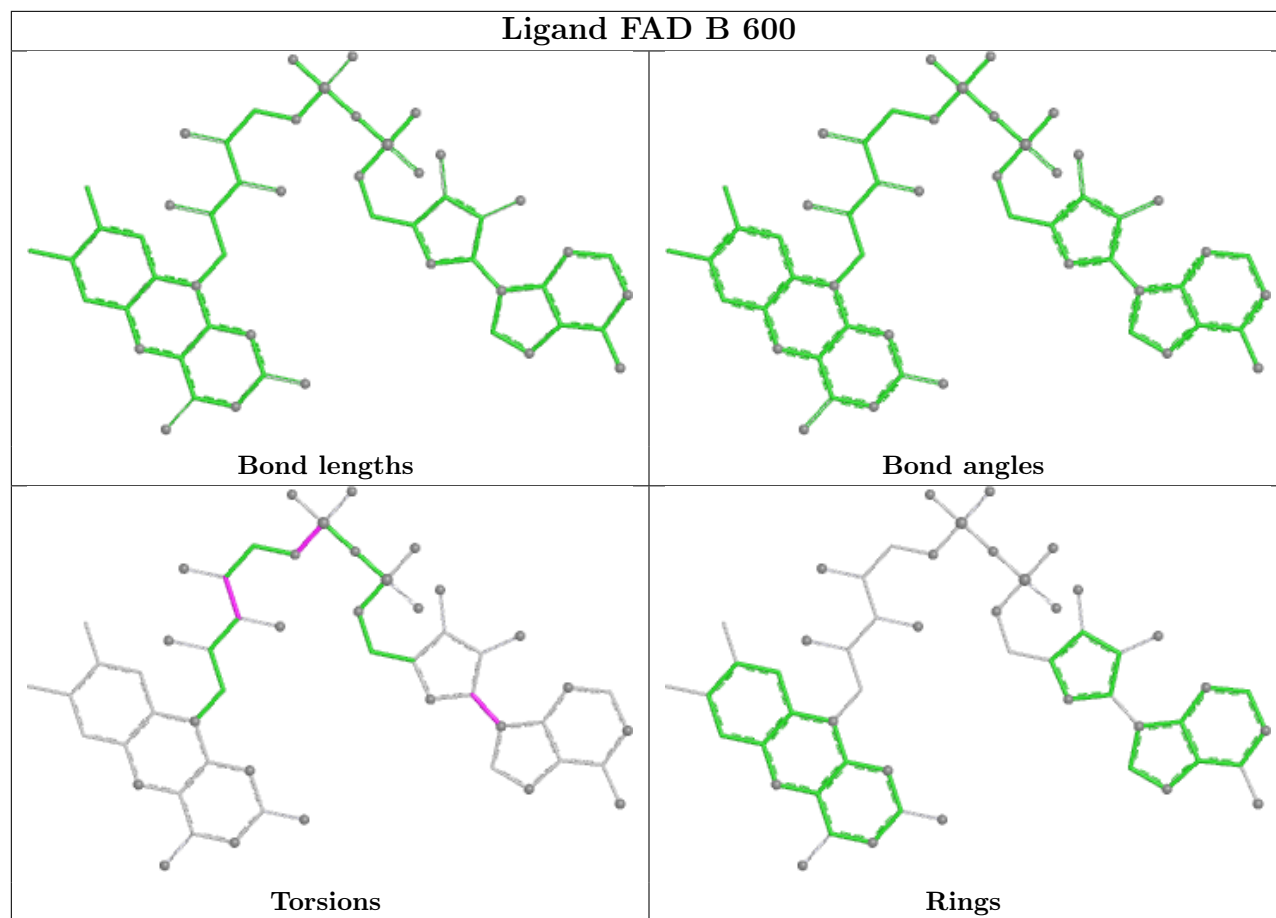
Ligand NRF B 601



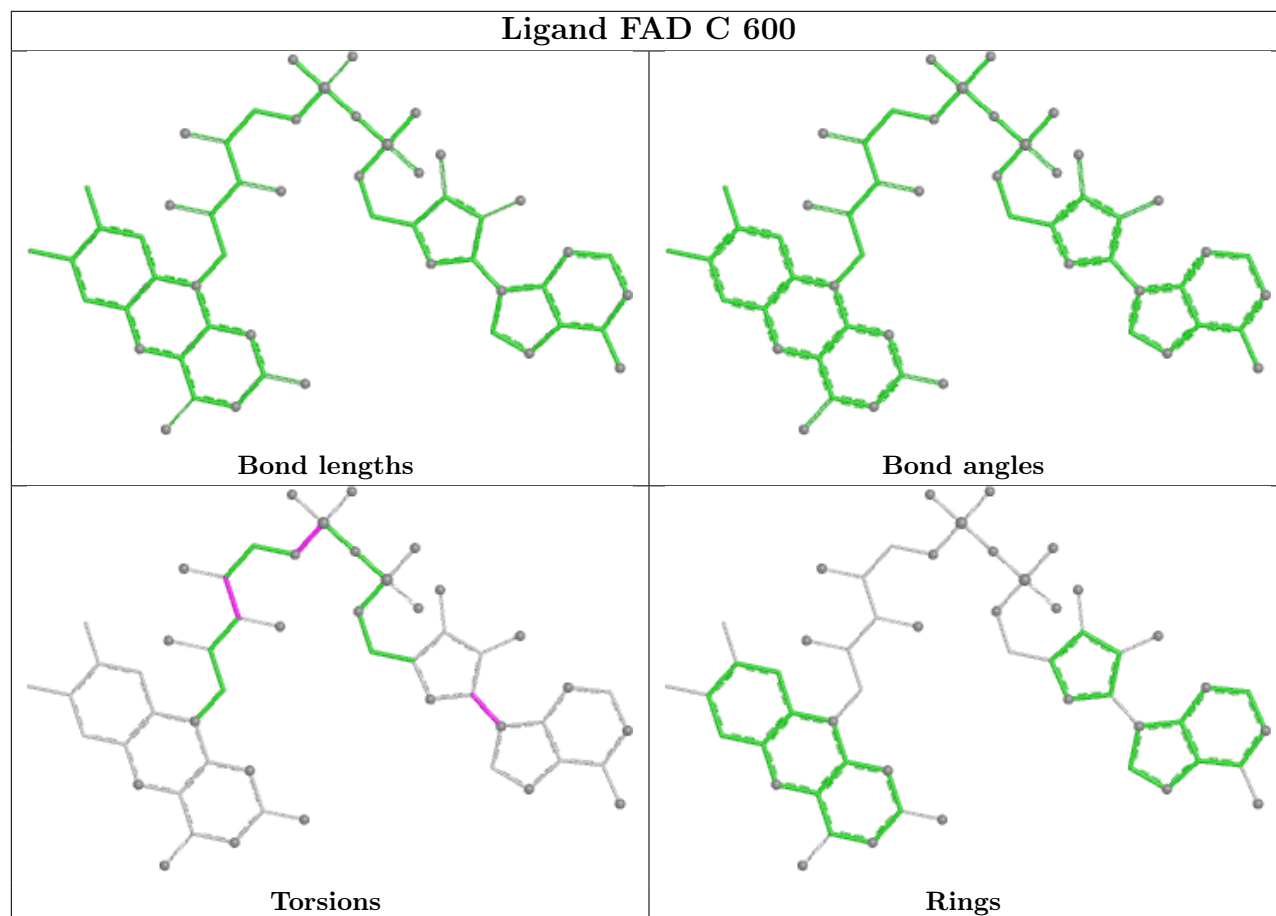
Ligand NRF A 601



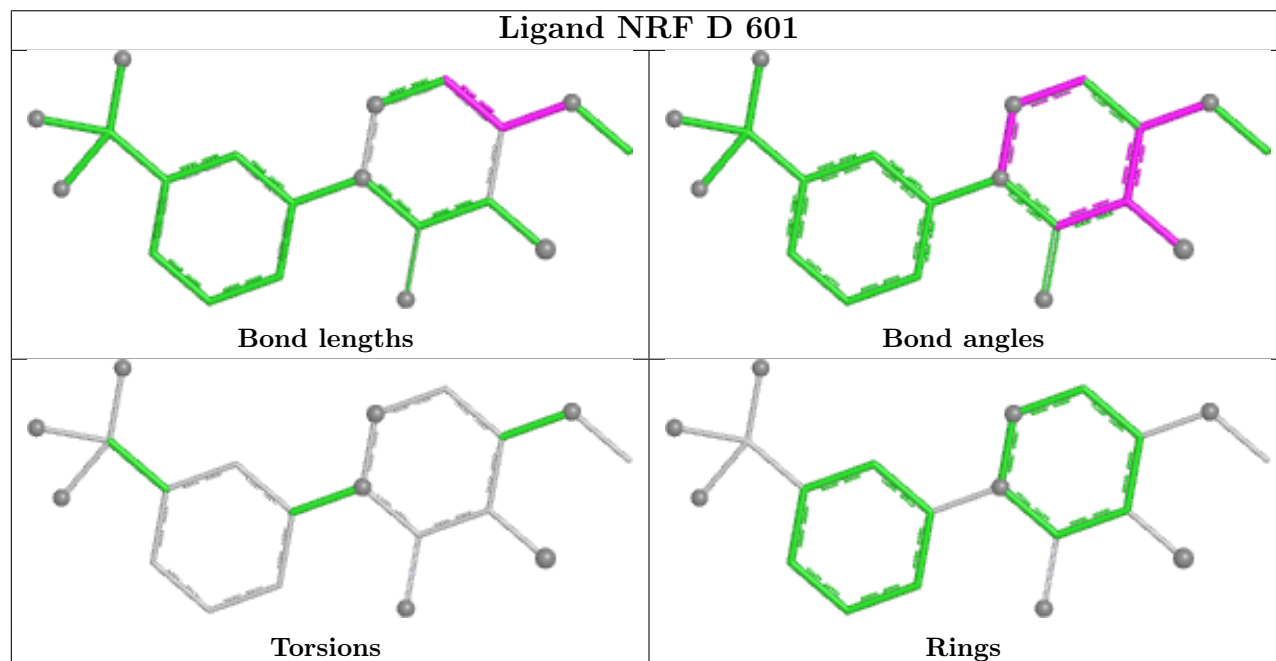


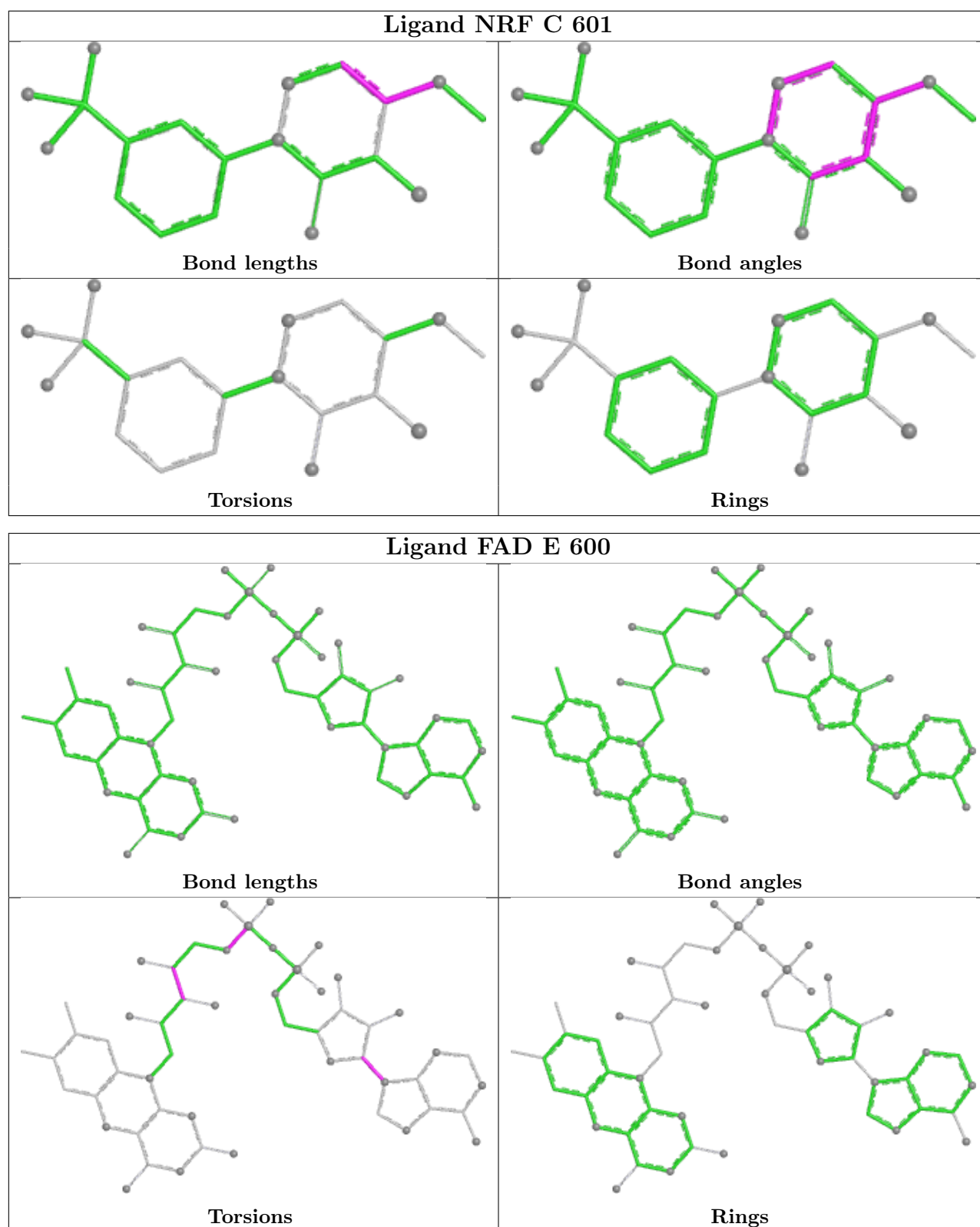


Ligand FAD C 600



Ligand NRF D 601





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 [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	467/497 (93%)	0.21	14 (2%)	52 45	38, 58, 83, 139	0
1	B	472/497 (94%)	0.17	13 (2%)	55 47	33, 57, 89, 135	1 (0%)
1	C	470/497 (94%)	0.75	27 (5%)	29 24	47, 75, 102, 180	1 (0%)
1	D	470/497 (94%)	0.78	51 (10%)	10 8	47, 73, 105, 211	0
1	E	466/497 (93%)	2.09	221 (47%)	0 0	68, 125, 162, 215	0
All	All	2345/2485 (94%)	0.80	326 (13%)	6 5	33, 71, 142, 215	2 (0%)

The worst 5 of 326 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	567	LEU	11.8
1	E	567	LEU	5.7
1	E	461	VAL	5.6
1	E	499	VAL	5.6
1	E	454	PHE	5.4

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 oligosaccharides 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,

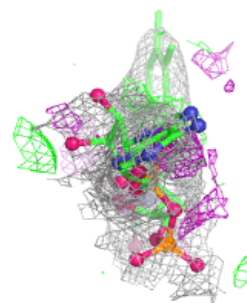
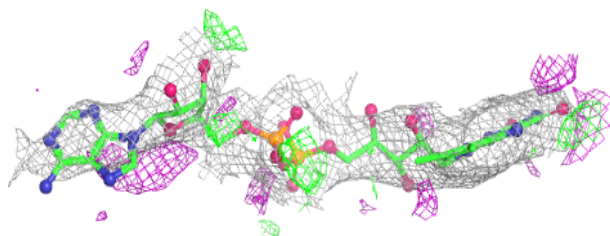
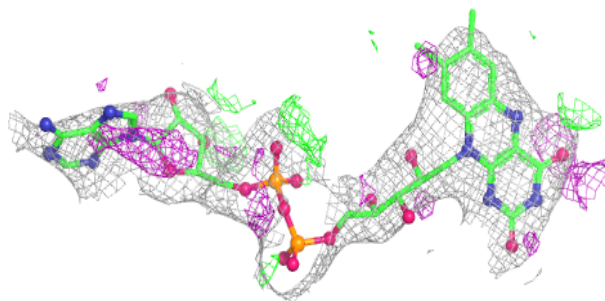
median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
5	IMD	D	603	5/5	0.52	0.41	122,122,123,123	0
4	EPE	E	602	15/15	0.53	0.26	143,148,161,161	0
4	EPE	A	603	15/15	0.54	0.34	140,147,160,160	0
4	EPE	B	603	15/15	0.69	0.22	100,110,131,132	0
5	IMD	B	606	5/5	0.78	0.26	94,94,95,95	0
6	TRS	B	604	8/8	0.78	0.22	102,105,107,108	0
5	IMD	D	604	5/5	0.80	0.31	114,114,114,114	0
2	FAD	E	600	53/53	0.80	0.15	93,117,145,145	0
5	IMD	D	605	5/5	0.85	0.23	98,98,98,98	0
5	IMD	B	605	5/5	0.86	0.33	112,112,113,113	0
6	TRS	D	602	8/8	0.86	0.17	100,103,104,105	0
3	NRF	E	601	20/20	0.87	0.14	95,100,102,102	0
5	IMD	A	604	5/5	0.88	0.24	108,108,108,108	0
4	EPE	A	602	15/15	0.90	0.17	80,84,91,92	0
4	EPE	B	602	15/15	0.90	0.16	78,85,91,91	0
2	FAD	D	600	53/53	0.95	0.08	48,56,62,63	0
2	FAD	C	600	53/53	0.95	0.08	44,57,62,63	0
3	NRF	C	601	20/20	0.96	0.08	59,66,68,68	0
3	NRF	D	601	20/20	0.96	0.09	57,63,68,68	0
2	FAD	B	600	53/53	0.97	0.07	38,45,48,49	0
2	FAD	A	600	53/53	0.97	0.06	31,39,45,47	0
3	NRF	A	601	20/20	0.97	0.07	42,48,51,52	0
3	NRF	B	601	20/20	0.97	0.07	39,48,54,58	0

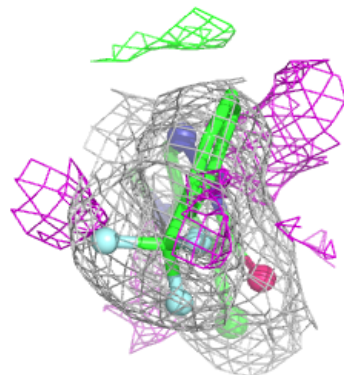
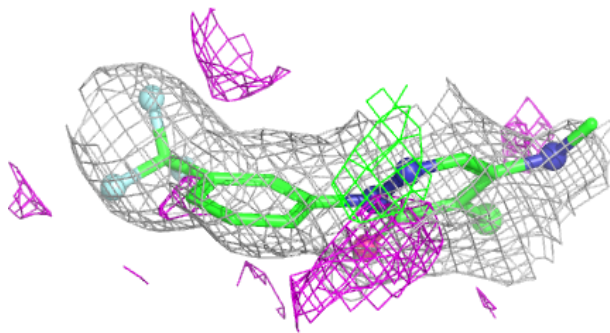
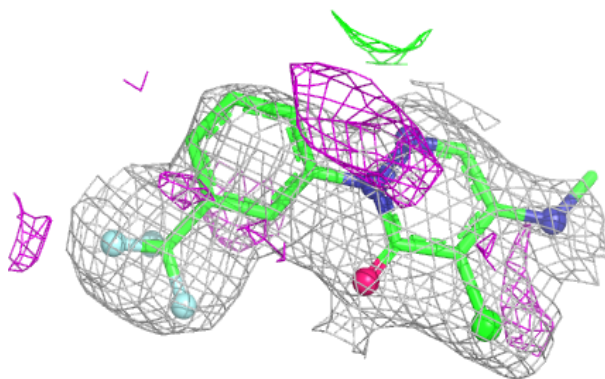
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around FAD E 600:

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

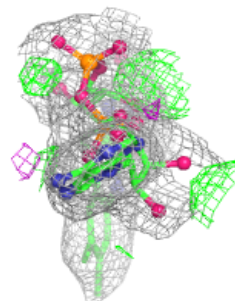
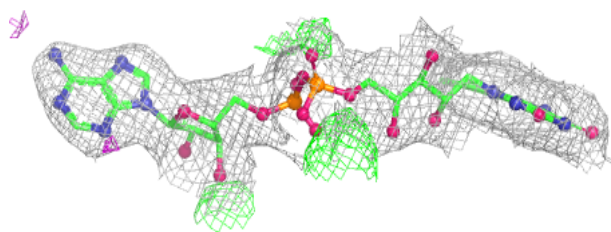
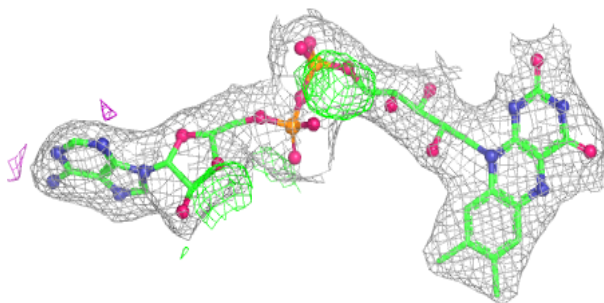
**Electron density around NRF E 601:**

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

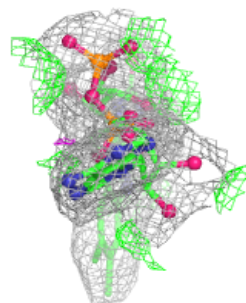
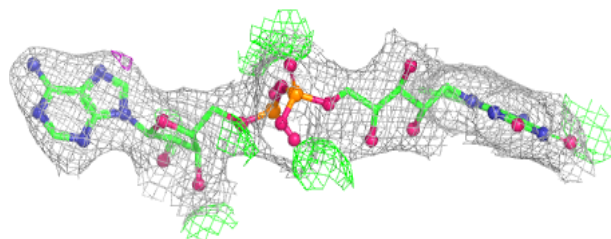
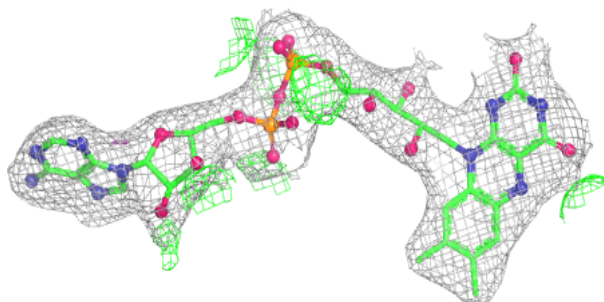


Electron density around FAD D 600:

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

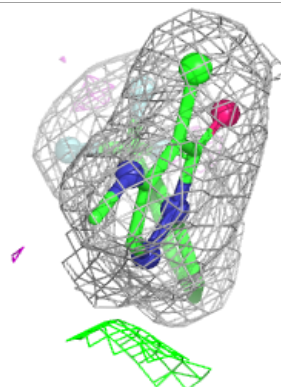
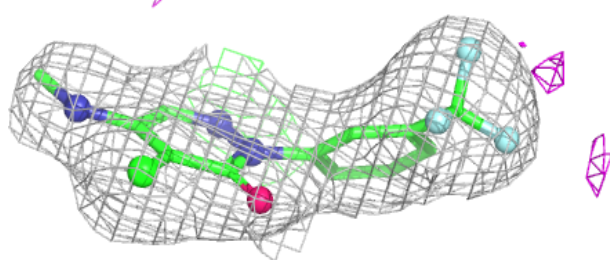
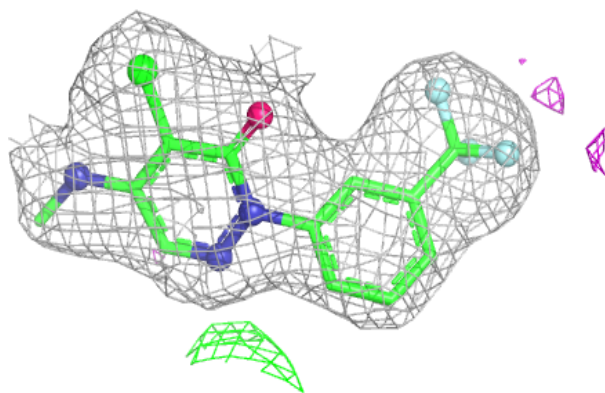
**Electron density around FAD C 600:**

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

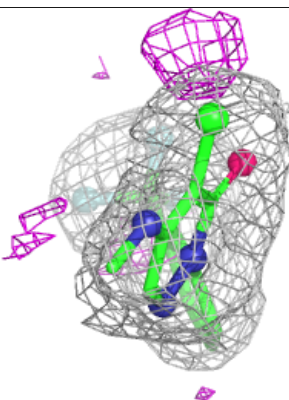
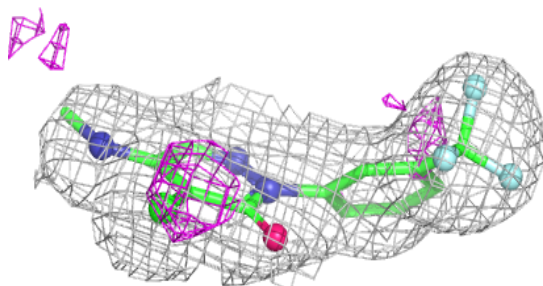
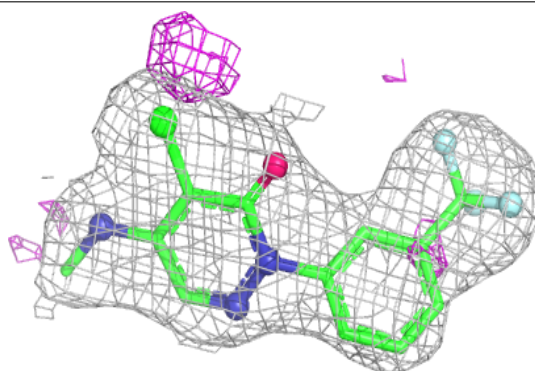


Electron density around NRF C 601:

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

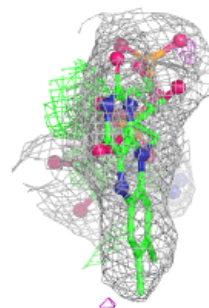
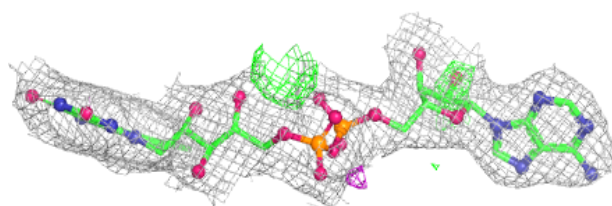
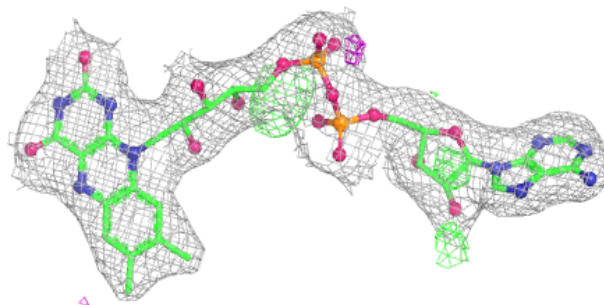
**Electron density around NRF D 601:**

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and green (positive)

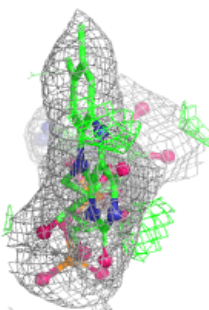
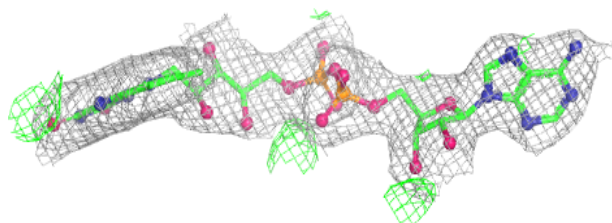
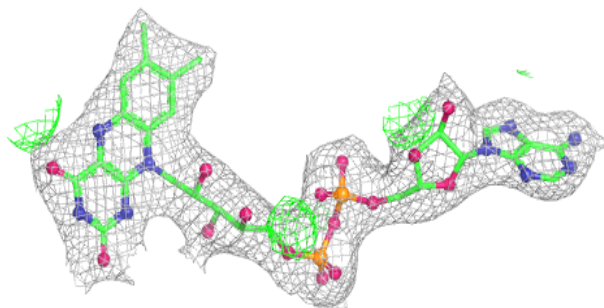


Electron density around FAD B 600:

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

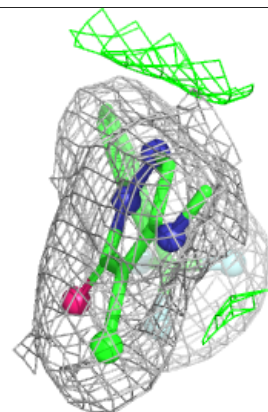
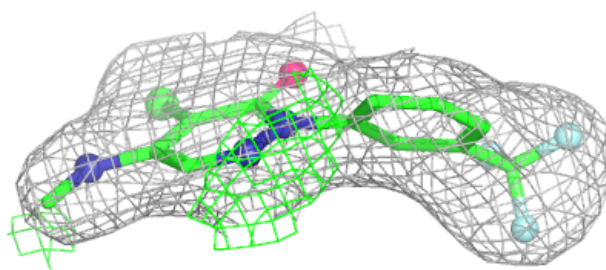
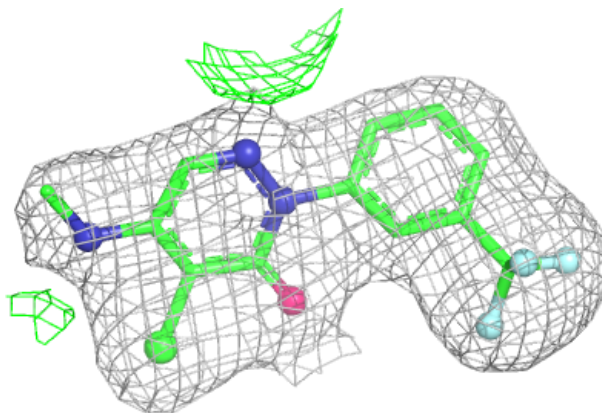
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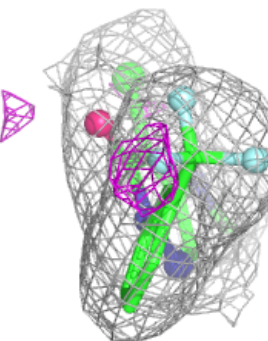
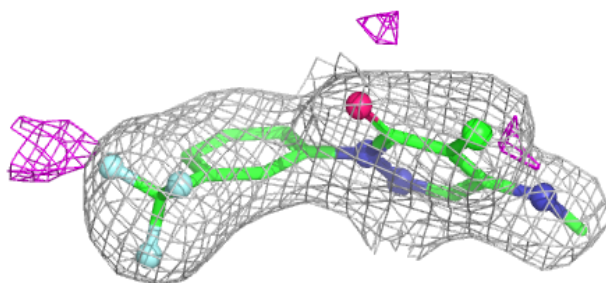
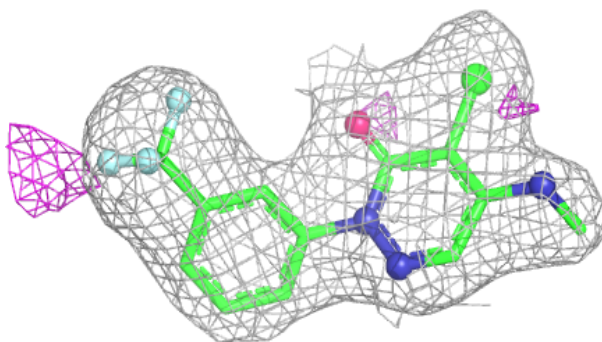


Electron density around NRF A 601:

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

**Electron density around NRF B 601:**

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



6.5 Other polymers [i](#)

There are no such residues in this entry.