



Full wwPDB X-ray Structure Validation Report ⓘ

May 14, 2020 – 01:33 pm BST

PDB ID : 5JWB
Title : Structure of NDH2 from plasmodium falciparum in complex with NADH
Authors : Yu, Y.; Li, X.L.
Deposited on : 2016-05-12
Resolution : 2.70 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

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

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

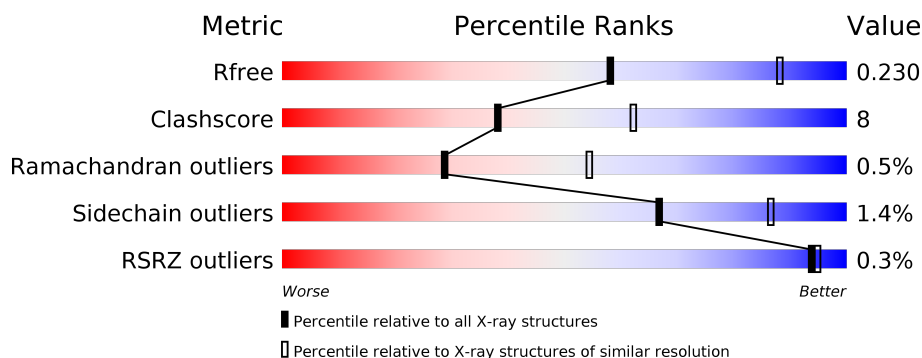
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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	2808 (2.70-2.70)
Clashscore	141614	3122 (2.70-2.70)
Ramachandran outliers	138981	3069 (2.70-2.70)
Sidechain outliers	138945	3069 (2.70-2.70)
RSRZ outliers	127900	2737 (2.70-2.70)

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

Mol	Chain	Length	Quality of chain
1	A	521	 81% 13% 5%
1	H	521	 81% 12% • 6%

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 8456 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.

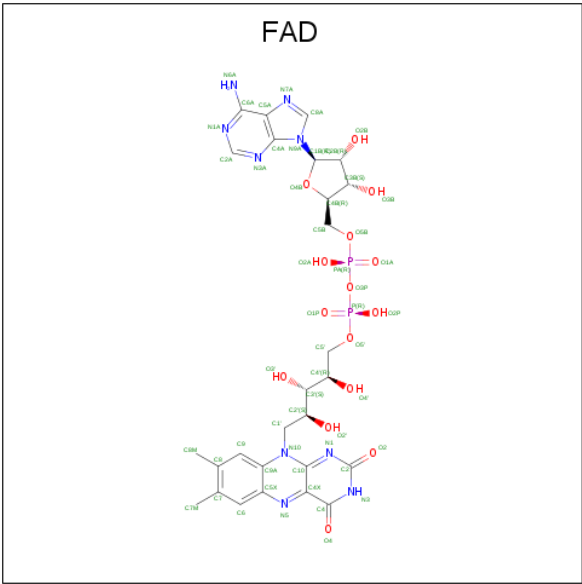
- Molecule 1 is a protein called Type II NADH:ubiquinone oxidoreductase.

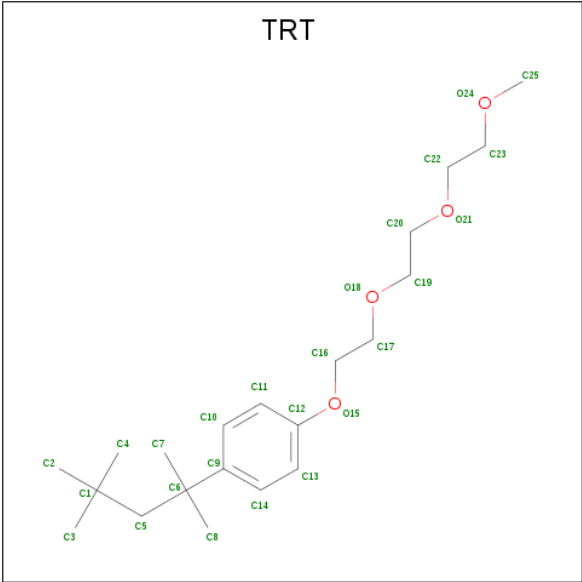
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	493	Total	C	N	O	S	0	0	0
			4030	2632	660	728	10			
1	H	492	Total	C	N	O	S	0	0	0
			4023	2628	659	726	10			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	13	MET	-	initiating methionine	UNP Q8I302
A	14	ARG	-	expression tag	UNP Q8I302
A	15	GLY	-	expression tag	UNP Q8I302
A	16	SER	-	expression tag	UNP Q8I302
A	17	HIS	-	expression tag	UNP Q8I302
A	18	HIS	-	expression tag	UNP Q8I302
A	19	HIS	-	expression tag	UNP Q8I302
A	20	HIS	-	expression tag	UNP Q8I302
A	21	HIS	-	expression tag	UNP Q8I302
A	22	HIS	-	expression tag	UNP Q8I302
A	23	GLY	-	expression tag	UNP Q8I302
A	24	SER	-	expression tag	UNP Q8I302
H	13	MET	-	initiating methionine	UNP Q8I302
H	14	ARG	-	expression tag	UNP Q8I302
H	15	GLY	-	expression tag	UNP Q8I302
H	16	SER	-	expression tag	UNP Q8I302
H	17	HIS	-	expression tag	UNP Q8I302
H	18	HIS	-	expression tag	UNP Q8I302
H	19	HIS	-	expression tag	UNP Q8I302
H	20	HIS	-	expression tag	UNP Q8I302
H	21	HIS	-	expression tag	UNP Q8I302
H	22	HIS	-	expression tag	UNP Q8I302
H	23	GLY	-	expression tag	UNP Q8I302
H	24	SER	-	expression tag	UNP Q8I302

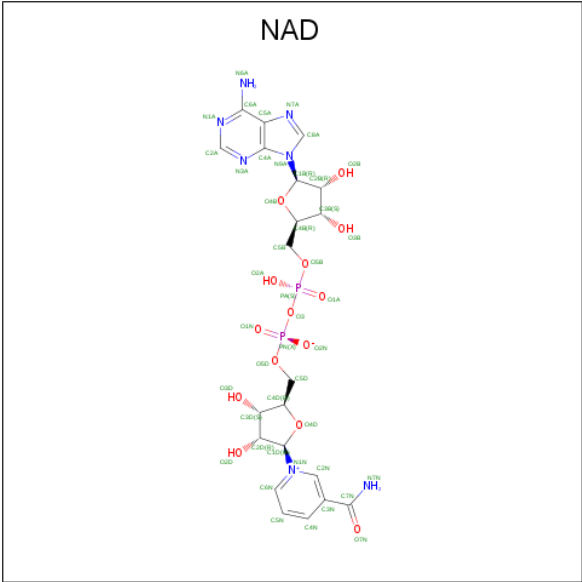
- Molecule 2 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula: C₂₇H₃₃N₉O₁₅P₂).





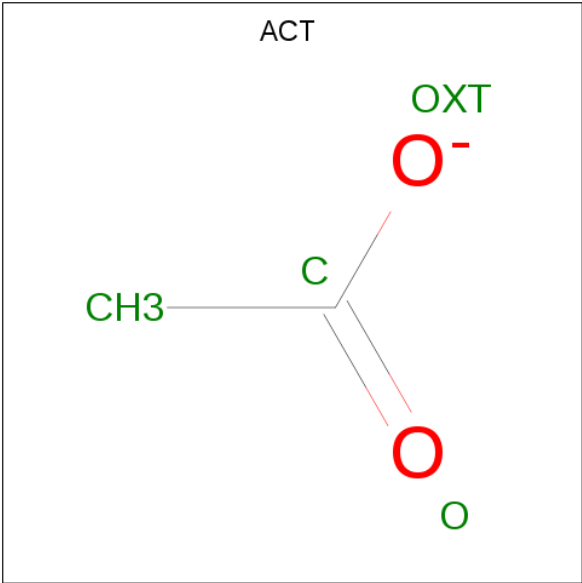
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			20	18	2		
4	A	1	Total	C	O	0	0
			20	18	2		
4	A	1	Total	C	O	0	0
			20	18	2		
4	H	1	Total	C	O	0	0
			20	18	2		
4	H	1	Total	C	O	0	0
			17	16	1		
4	H	1	Total	C	O	0	0
			20	18	2		

- Molecule 5 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (three-letter code: NAD) (formula: C₂₁H₂₇N₇O₁₄P₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	A	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
5	H	1	Total	C	N	O	P	0	0
			44	21	7	14	2		

- Molecule 6 is ACETATE ION (three-letter code: ACT) (formula: C₂H₃O₂).



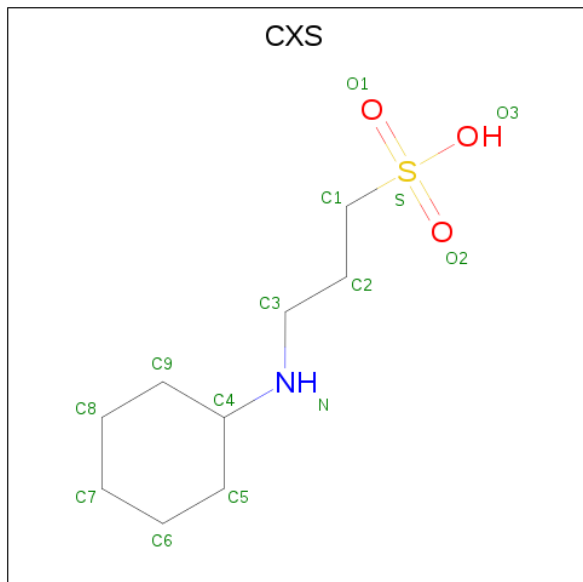
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			4	2	2		
6	A	1	Total	C	O	0	0
			4	2	2		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			4	2	2		
6	A	1	Total	C	O	0	0
			4	2	2		
6	H	1	Total	C	O	0	0
			4	2	2		
6	H	1	Total	C	O	0	0
			4	2	2		
6	H	1	Total	C	O	0	0
			4	2	2		

- Molecule 7 is 3-CYCLOHEXYL-1-PROPYLSULFONIC ACID (three-letter code: CXS) (formula: $C_9H_{19}NO_3S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
7	A	1	Total	C	N	O	S	0	0
			14	9	1	3	1		
7	H	1	Total	C	N	O	S	0	0
			14	9	1	3	1		

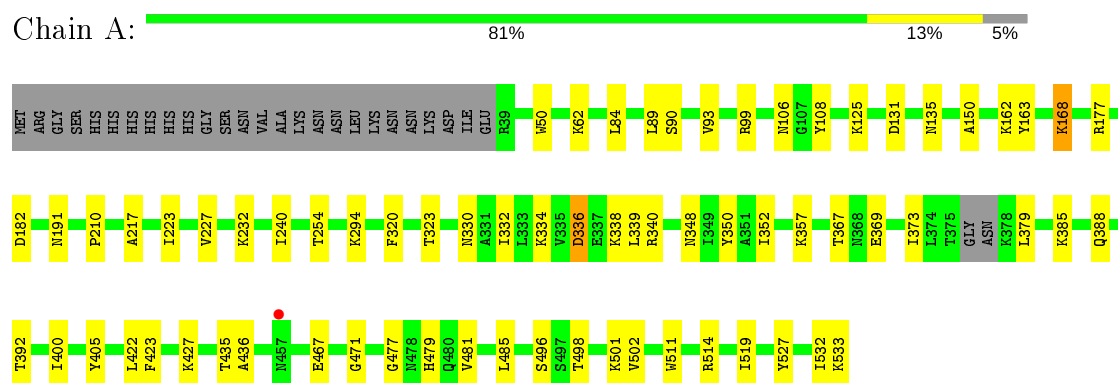
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	14	Total	O	0	0
			14	14		
8	H	14	Total	O	0	0
			14	14		

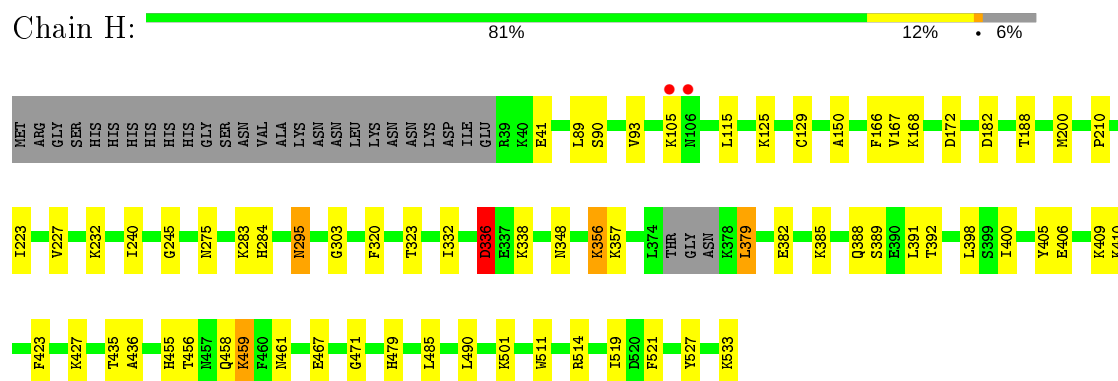
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

- Molecule 1: Type II NADH:ubiquinone oxidoreductase



- Molecule 1: Type II NADH:ubiquinone oxidoreductase



4 Data and refinement statistics

Property	Value	Source
Space group	I 41	Depositor
Cell constants a, b, c, α , β , γ	191.53 Å 191.53 Å 91.26 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	45.91 – 2.70 45.91 – 2.70	Depositor EDS
% Data completeness (in resolution range)	99.9 (45.91-2.70) 99.7 (45.91-2.70)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.65 (at 2.69 Å)	Xtriage
Refinement program	PHENIX PHENIX.REFINE:1.8.4_1496	Depositor
R, R_{free}	0.180 , 0.226 0.183 , 0.230	Depositor DCC
R_{free} test set	2295 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	50.9	Xtriage
Anisotropy	0.161	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 16.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.488 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	8456	wwPDB-VP
Average B, all atoms (Å ²)	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.40% 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: MG, CXS, NAD, ACT, TRT, 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	0.46	0/4127	0.55	0/5575
1	H	0.46	0/4120	0.56	0/5565
All	All	0.46	0/8247	0.56	0/11140

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	4030	0	4063	44	0
1	H	4023	0	4056	46	0
2	A	53	0	31	1	0
2	H	53	0	31	2	0
3	A	4	0	0	0	0
3	H	4	0	0	0	0
4	A	60	0	81	24	0
4	H	57	0	77	28	0
5	A	44	0	26	1	0
5	H	44	0	26	1	0
6	A	16	0	12	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	H	12	0	9	0	0
7	A	14	0	19	2	0
7	H	14	0	19	1	0
8	A	14	0	0	0	0
8	H	14	0	0	0	0
All	All	8456	0	8450	128	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 8.

All (128) 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)
4:H:609:TRT:C2	4:H:609:TRT:H7C3	1.62	1.29
4:A:606:TRT:C8	4:A:606:TRT:H4C3	1.61	1.29
4:A:609:TRT:C8	4:A:609:TRT:H2C2	1.71	1.20
4:H:608:TRT:C7	4:H:608:TRT:H4C2	1.74	1.17
4:H:608:TRT:H2C1	4:H:609:TRT:H3C3	1.20	1.15
4:H:608:TRT:H7C1	4:H:608:TRT:H4C2	1.18	1.13
4:H:609:TRT:H2C3	4:H:609:TRT:H7C3	1.32	1.11
4:H:609:TRT:C7	4:H:609:TRT:H2C2	1.82	1.10
4:H:608:TRT:C4	4:H:608:TRT:C7	2.30	1.10
4:H:608:TRT:H2C1	4:H:609:TRT:C3	1.81	1.09
4:A:606:TRT:C4	4:A:606:TRT:C8	2.30	1.09
4:H:609:TRT:C7	4:H:609:TRT:C2	2.30	1.08
4:A:609:TRT:H8C2	4:A:609:TRT:H2C2	1.12	1.08
4:H:609:TRT:H2C2	4:H:609:TRT:H7C3	1.35	1.05
4:A:606:TRT:H8C1	4:A:606:TRT:H4C3	1.05	1.03
4:H:608:TRT:C2	4:H:609:TRT:C3	2.38	1.01
4:A:609:TRT:C2	4:A:609:TRT:H8C2	1.94	0.96
4:H:608:TRT:C4	4:H:608:TRT:H7C3	1.96	0.95
4:H:608:TRT:C2	4:H:609:TRT:H3C3	1.97	0.94
4:H:608:TRT:C2	4:H:609:TRT:H3C2	2.01	0.90
4:H:608:TRT:H4C3	4:H:608:TRT:H7C3	1.52	0.89
4:A:609:TRT:H4C3	4:A:609:TRT:C14	2.03	0.88
4:A:609:TRT:C2	4:A:609:TRT:C8	2.49	0.87
4:A:606:TRT:C4	4:A:606:TRT:H8C2	2.03	0.87
4:A:609:TRT:H4C3	4:A:609:TRT:C9	2.04	0.86
4:H:608:TRT:H2C3	4:H:609:TRT:H3C2	1.55	0.85
1:A:150:ALA:O	1:A:168:LYS:NZ	2.08	0.84
1:H:332:ILE:HG12	1:H:357:LYS:HG2	1.60	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:A:609:TRT:C4	4:A:609:TRT:C9	2.60	0.80
1:H:458:GLN:OE1	1:H:461:ASN:ND2	2.13	0.79
4:H:609:TRT:H2C2	4:H:609:TRT:H7C1	1.67	0.77
1:H:295:ASN:N	1:H:295:ASN:HD22	1.85	0.73
4:H:601:TRT:H4C3	4:H:601:TRT:C10	2.19	0.73
1:H:223:ILE:HD12	1:H:240:ILE:HD11	1.72	0.72
1:A:336:ASP:HB3	1:A:338:LYS:H	1.56	0.70
1:A:177:ARG:NH2	1:A:532:ILE:O	2.25	0.69
4:A:609:TRT:H8C1	4:A:609:TRT:H2C2	1.71	0.68
1:H:501:LYS:HB3	4:H:608:TRT:H8C3	1.75	0.68
1:A:471:GLY:HA3	1:A:485:LEU:HD23	1.77	0.66
4:A:606:TRT:H4C2	4:A:606:TRT:H8C2	1.76	0.66
4:A:606:TRT:H2C1	4:A:607:TRT:H4C1	1.79	0.65
1:H:382:GLU:OE1	1:H:409:LYS:NZ	2.24	0.64
1:A:332:ILE:HD13	1:A:357:LYS:HD3	1.80	0.64
1:A:223:ILE:HD12	1:A:240:ILE:HD11	1.80	0.63
1:H:521:PHE:HB3	4:H:608:TRT:H161	1.82	0.62
1:H:41:GLU:OE1	1:H:455:HIS:NE2	2.20	0.61
1:H:295:ASN:H	1:H:295:ASN:HD22	1.49	0.60
1:H:336:ASP:HB3	1:H:338:LYS:H	1.66	0.60
4:H:601:TRT:H4C3	4:H:601:TRT:H10	1.82	0.60
1:A:514:ARG:NH2	4:A:607:TRT:H162	2.18	0.57
1:H:514:ARG:NH2	4:H:609:TRT:H162	2.19	0.57
1:A:210:PRO:HB2	5:A:608:NAD:C4N	2.35	0.57
1:A:514:ARG:HH22	4:A:607:TRT:H162	1.70	0.57
1:H:210:PRO:HB2	5:H:607:NAD:C4N	2.35	0.56
1:A:498:THR:O	1:A:502:VAL:HG23	2.07	0.55
1:H:105:LYS:N	1:H:105:LYS:HD3	2.22	0.54
1:A:511:TRP:HB3	4:H:601:TRT:H161	1.90	0.53
1:A:338:LYS:HD2	1:A:340:ARG:NH2	2.23	0.53
1:A:519:ILE:HG21	1:H:519:ILE:HG21	1.91	0.53
1:A:435:THR:HB	2:A:601:FAD:O2	2.09	0.53
1:H:388:GLN:O	1:H:392:THR:HB	2.09	0.52
1:A:89:LEU:HB3	1:A:93:VAL:HG11	1.92	0.52
1:H:356:LYS:O	1:H:356:LYS:HG3	2.10	0.52
1:H:89:LEU:HB3	1:H:93:VAL:HG11	1.91	0.52
1:A:367:THR:HG23	1:A:422:LEU:HD12	1.92	0.52
4:A:609:TRT:H161	1:H:511:TRP:HB3	1.91	0.52
4:A:609:TRT:H4C3	4:A:609:TRT:H8C2	1.91	0.51
1:H:182:ASP:OD2	1:H:533:LYS:NZ	2.41	0.51
1:A:423:PHE:CE2	1:A:427:LYS:HD2	2.46	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:338:LYS:HD2	1:A:340:ARG:HH21	1.75	0.50
1:A:479:HIS:NE2	7:A:614:CXS:H22	2.27	0.50
1:H:166:PHE:O	1:H:172:ASP:HB3	2.11	0.50
1:A:481:VAL:HB	1:A:496:SER:HB3	1.93	0.49
1:A:388:GLN:NE2	1:A:400:ILE:HG23	2.28	0.49
1:H:320:PHE:O	1:H:323:THR:HB	2.13	0.49
1:H:283:LYS:HE2	1:H:284:HIS:NE2	2.27	0.49
1:H:471:GLY:HA3	1:H:485:LEU:HD23	1.95	0.49
1:H:385:LYS:HB2	1:H:405:TYR:CD1	2.48	0.49
1:H:501:LYS:NZ	4:H:608:TRT:H162	2.28	0.48
1:A:162:LYS:HE3	1:A:163:TYR:OH	2.14	0.48
1:H:295:ASN:ND2	1:H:295:ASN:N	2.53	0.48
1:A:223:ILE:HA	1:A:227:VAL:HB	1.96	0.48
1:H:435:THR:HB	2:H:602:FAD:O2	2.14	0.48
1:A:125:LYS:HE3	1:A:348:ASN:ND2	2.28	0.47
1:A:62:LYS:HD2	1:A:108:TYR:CE2	2.49	0.47
1:A:340:ARG:HG3	1:A:350:TYR:CE1	2.49	0.47
1:A:182:ASP:OD2	1:A:533:LYS:NZ	2.33	0.46
1:A:501:LYS:NZ	4:A:606:TRT:O18	2.48	0.46
1:H:232:LYS:HE3	1:H:232:LYS:HB2	1.71	0.46
1:H:379:LEU:HD12	1:H:379:LEU:HA	1.77	0.46
1:A:50:TRP:CD2	1:A:436:ALA:HB1	2.51	0.46
1:A:369:GLU:O	1:A:373:ILE:HG13	2.16	0.45
1:H:389:SER:O	1:H:392:THR:HG22	2.16	0.45
1:H:223:ILE:HA	1:H:227:VAL:HB	1.99	0.45
1:H:388:GLN:NE2	1:H:400:ILE:HG23	2.30	0.45
1:A:232:LYS:HB2	1:A:232:LYS:HE2	1.62	0.45
1:H:391:LEU:HB3	1:H:398:LEU:CD1	2.46	0.45
1:H:423:PHE:CE2	1:H:427:LYS:HD3	2.51	0.45
1:H:527:TYR:OH	4:H:601:TRT:H162	2.17	0.45
1:H:436:ALA:HB2	2:H:602:FAD:H2'	1.99	0.44
1:H:90:SER:O	1:H:93:VAL:HG12	2.17	0.44
1:A:479:HIS:CD2	7:A:614:CXS:H22	2.52	0.44
1:H:456:THR:O	1:H:459:LYS:HG2	2.17	0.44
1:H:479:HIS:CE1	7:H:613:CXS:H12	2.52	0.44
1:A:339:LEU:HD13	1:A:352:ILE:HG22	2.00	0.44
1:A:217:ALA:HB1	1:A:477:GLY:N	2.32	0.44
1:H:410:LYS:HB3	1:H:410:LYS:HE3	1.82	0.43
1:A:320:PHE:O	1:A:323:THR:HB	2.18	0.43
1:A:106:ASN:OD1	1:A:108:TYR:N	2.39	0.43
1:A:385:LYS:HB2	1:A:405:TYR:CD1	2.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:H:608:TRT:C2	4:H:609:TRT:H5C2	2.49	0.43
1:H:150:ALA:O	1:H:168:LYS:HE2	2.19	0.42
4:H:601:TRT:C9	4:H:601:TRT:H4C3	2.48	0.42
1:H:245:GLY:O	1:H:275:ASN:HA	2.18	0.42
4:A:609:TRT:C4	4:A:609:TRT:H8C2	2.48	0.42
1:A:527:TYR:OH	4:A:609:TRT:C16	2.68	0.42
1:H:115:LEU:HB3	1:H:129:CYS:HB3	2.01	0.42
1:A:501:LYS:NZ	4:A:606:TRT:C20	2.83	0.42
1:H:200:MET:O	1:H:303:GLY:HA3	2.20	0.42
4:A:606:TRT:H13	4:A:606:TRT:H161	1.55	0.41
4:A:606:TRT:H8C2	4:A:606:TRT:H4C3	1.63	0.41
1:A:99:ARG:NH2	1:H:188:THR:HG21	2.35	0.41
1:A:90:SER:O	1:A:93:VAL:HG12	2.20	0.41
1:A:131:ASP:OD1	1:A:135:ASN:HB2	2.20	0.41
1:A:330:ASN:ND2	1:A:334:LYS:HD3	2.36	0.41
1:A:84:LEU:HD12	1:A:84:LEU:HA	1.91	0.41
4:H:608:TRT:H2C2	4:H:609:TRT:H5C2	2.03	0.41
1:H:125:LYS:HE2	1:H:348:ASN:ND2	2.36	0.40

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	489/521 (94%)	471 (96%)	15 (3%)	3 (1%)	25	50
1	H	488/521 (94%)	470 (96%)	16 (3%)	2 (0%)	34	60
All	All	977/1042 (94%)	941 (96%)	31 (3%)	5 (0%)	29	54

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	336	ASP
1	A	379	LEU
1	H	379	LEU
1	H	336	ASP
1	A	191	ASN

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	449/474 (95%)	444 (99%)	5 (1%)	73	90
1	H	448/474 (94%)	440 (98%)	8 (2%)	59	83
All	All	897/948 (95%)	884 (99%)	13 (1%)	67	86

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	168	LYS
1	A	254	THR
1	A	294	LYS
1	A	392	THR
1	A	467	GLU
1	H	167	VAL
1	H	295	ASN
1	H	336	ASP
1	H	356	LYS
1	H	406	GLU
1	H	459	LYS
1	H	467	GLU
1	H	490	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	H	295	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 carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 27 ligands modelled in this entry, 8 are monoatomic - leaving 19 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	$\# Z > 2$
6	ACT	A	611	-	1,3,3	2.65	1 (100%)	0,3,3	0.00	-
5	NAD	H	607	-	42,48,48	2.45	16 (38%)	50,73,73	1.65	8 (16%)
2	FAD	H	602	3	51,58,58	1.27	6 (11%)	60,89,89	1.85	9 (15%)
5	NAD	A	608	-	42,48,48	2.42	15 (35%)	50,73,73	1.55	8 (16%)
4	TRT	H	601	-	20,20,25	0.91	1 (5%)	28,28,33	4.62	7 (25%)
6	ACT	A	612	-	1,3,3	2.23	1 (100%)	0,3,3	0.00	-
6	ACT	H	611	-	1,3,3	2.15	1 (100%)	0,3,3	0.00	-
6	ACT	A	610	-	1,3,3	2.54	1 (100%)	0,3,3	0.00	-
4	TRT	H	609	-	20,20,25	1.04	1 (5%)	28,28,33	0.82	1 (3%)
4	TRT	A	606	-	20,20,25	0.79	1 (5%)	28,28,33	1.28	4 (14%)
6	ACT	H	612	-	1,3,3	2.09	1 (100%)	0,3,3	0.00	-
4	TRT	H	608	-	17,17,25	0.99	1 (5%)	25,25,33	0.91	1 (4%)
2	FAD	A	601	3	51,58,58	1.25	6 (11%)	60,89,89	1.84	11 (18%)
7	CXS	H	613	-	14,14,14	1.68	3 (21%)	18,18,18	1.40	3 (16%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	ACT	A	613	-	1,3,3	1.57	0	0,3,3	0.00	-
7	CXS	A	614	-	14,14,14	1.59	4 (28%)	18,18,18	1.64	4 (22%)
4	TRT	A	607	-	20,20,25	1.35	2 (10%)	28,28,33	1.18	1 (3%)
6	ACT	H	610	-	1,3,3	2.65	1 (100%)	0,3,3	0.00	-
4	TRT	A	609	-	20,20,25	0.88	1 (5%)	28,28,33	1.19	2 (7%)

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	NAD	H	607	-	-	1/26/62/62	0/5/5/5
2	FAD	H	602	3	-	1/30/50/50	0/6/6/6
4	TRT	H	601	-	-	7/18/18/23	0/1/1/1
5	NAD	A	608	-	-	2/26/62/62	0/5/5/5
4	TRT	H	609	-	-	7/18/18/23	0/1/1/1
4	TRT	A	606	-	-	4/18/18/23	0/1/1/1
4	TRT	H	608	-	-	13/15/15/23	0/1/1/1
2	FAD	A	601	3	-	1/30/50/50	0/6/6/6
7	CXS	H	613	-	-	1/8/16/16	0/1/1/1
7	CXS	A	614	-	-	4/8/16/16	0/1/1/1
4	TRT	A	607	-	-	6/18/18/23	0/1/1/1
4	TRT	A	609	-	-	2/18/18/23	0/1/1/1

All (63) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	H	607	NAD	O4B-C1B	6.38	1.50	1.41
5	A	608	NAD	O4D-C1D	6.27	1.49	1.41
5	H	607	NAD	O4D-C1D	6.16	1.49	1.41
5	A	608	NAD	O4B-C1B	6.13	1.49	1.41
5	H	607	NAD	C7N-N7N	5.27	1.43	1.33
5	A	608	NAD	C7N-N7N	4.70	1.41	1.33
5	A	608	NAD	C2N-N1N	4.56	1.40	1.35
5	A	608	NAD	C2B-C1B	-4.29	1.47	1.53
5	H	607	NAD	C2A-N3A	4.16	1.38	1.32
5	H	607	NAD	C2N-N1N	4.05	1.39	1.35
5	A	608	NAD	C2A-N3A	3.78	1.38	1.32

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	607	TRT	O15-C12	3.76	1.46	1.37
5	H	607	NAD	C2B-C1B	-3.59	1.48	1.53
2	H	602	FAD	C2A-N3A	3.57	1.37	1.32
5	H	607	NAD	C2D-C1D	-3.54	1.48	1.53
5	A	608	NAD	C4N-C3N	3.48	1.45	1.39
7	H	613	CXS	C1-S	3.48	1.82	1.77
2	H	602	FAD	C10-N1	3.47	1.37	1.33
5	H	607	NAD	C4N-C3N	3.41	1.45	1.39
5	A	608	NAD	C2D-C1D	-3.37	1.48	1.53
4	H	609	TRT	O15-C12	3.34	1.45	1.37
2	A	601	FAD	C2A-N3A	3.22	1.37	1.32
2	H	602	FAD	C4X-N5	3.15	1.37	1.33
7	A	614	CXS	C1-S	3.13	1.82	1.77
2	A	601	FAD	C4X-N5	3.08	1.37	1.33
7	H	613	CXS	O1-S	3.02	1.53	1.45
5	A	608	NAD	C8A-N7A	3.01	1.40	1.34
2	A	601	FAD	C10-N1	3.00	1.37	1.33
7	H	613	CXS	O2-S	2.99	1.53	1.45
5	H	607	NAD	C8A-N7A	2.93	1.39	1.34
2	A	601	FAD	C4-N3	2.88	1.38	1.33
2	H	602	FAD	C4-N3	2.88	1.38	1.33
5	A	608	NAD	C2N-C3N	2.84	1.43	1.39
7	A	614	CXS	O1-S	2.78	1.53	1.45
5	H	607	NAD	C6A-N6A	2.78	1.44	1.34
5	A	608	NAD	O4D-C4D	2.71	1.51	1.45
6	A	611	ACT	CH3-C	2.65	1.52	1.48
6	H	610	ACT	CH3-C	2.65	1.52	1.48
7	A	614	CXS	O2-S	2.64	1.52	1.45
4	H	608	TRT	O15-C12	2.61	1.43	1.37
5	H	607	NAD	O4D-C4D	2.58	1.50	1.45
6	A	610	ACT	CH3-C	2.54	1.52	1.48
2	H	602	FAD	C2A-N1A	2.47	1.38	1.33
5	H	607	NAD	C2N-C3N	2.47	1.42	1.39
4	H	601	TRT	O15-C12	2.41	1.43	1.37
5	H	607	NAD	C3D-C4D	-2.36	1.47	1.53
5	A	608	NAD	C6A-N6A	2.33	1.42	1.34
5	A	608	NAD	C3D-C4D	-2.31	1.47	1.53
4	A	607	TRT	C13-C12	2.26	1.43	1.38
6	A	612	ACT	CH3-C	2.23	1.51	1.48
2	A	601	FAD	C2A-N1A	2.21	1.38	1.33
5	H	607	NAD	C2D-C3D	-2.17	1.47	1.53
6	H	611	ACT	CH3-C	2.15	1.51	1.48

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	609	TRT	C6-C9	-2.12	1.50	1.53
4	A	606	TRT	O15-C12	2.11	1.42	1.37
5	A	608	NAD	C5A-N7A	2.11	1.47	1.39
5	H	607	NAD	C5N-C4N	2.10	1.43	1.38
5	H	607	NAD	C5A-N7A	2.09	1.47	1.39
6	H	612	ACT	CH3-C	2.09	1.51	1.48
7	A	614	CXS	C4-N	-2.05	1.43	1.48
2	H	602	FAD	C1'-N10	2.05	1.50	1.48
5	A	608	NAD	C5N-C4N	2.04	1.43	1.38
2	A	601	FAD	C1'-N10	2.04	1.50	1.48

All (59) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	H	601	TRT	C5-C6-C9	-15.00	78.80	111.93
4	H	601	TRT	C8-C6-C9	-11.81	80.59	110.20
4	H	601	TRT	C7-C6-C8	-9.38	81.61	107.28
2	H	602	FAD	C4-N3-C2	7.89	121.80	115.14
2	A	601	FAD	C4-N3-C2	7.88	121.79	115.14
4	H	601	TRT	C7-C6-C5	-7.15	75.61	109.08
4	H	601	TRT	C7-C6-C9	6.64	126.85	110.20
2	H	602	FAD	N3A-C2A-N1A	-5.52	120.05	128.68
2	A	601	FAD	N3A-C2A-N1A	-5.28	120.43	128.68
5	H	607	NAD	C6N-N1N-C2N	-5.12	117.31	121.97
4	A	607	TRT	C5-C6-C9	-4.74	101.47	111.93
4	H	601	TRT	C8-C6-C5	4.49	130.10	109.08
5	A	608	NAD	C6N-N1N-C2N	-4.48	117.89	121.97
4	A	609	TRT	C1-C5-C6	-4.23	108.68	123.91
4	H	601	TRT	C1-C5-C6	-4.18	108.86	123.91
5	A	608	NAD	C5N-C4N-C3N	-4.15	115.43	120.34
5	H	607	NAD	C3B-C2B-C1B	3.97	106.96	100.98
2	H	602	FAD	C1'-N10-C9A	3.87	121.34	118.29
2	A	601	FAD	C5X-C9A-N10	3.82	120.49	117.72
2	A	601	FAD	C4'-C3'-C2'	-3.71	105.65	113.36
5	H	607	NAD	C5N-C4N-C3N	-3.70	115.97	120.34
2	A	601	FAD	C4X-C4-N3	-3.63	118.47	123.43
2	H	602	FAD	C4X-C4-N3	-3.59	118.52	123.43
2	H	602	FAD	C4X-N5-C5X	3.58	120.35	116.77
2	H	602	FAD	C1B-N9A-C4A	-3.54	120.42	126.64
5	H	607	NAD	C5N-C6N-N1N	3.51	125.43	120.40
7	A	614	CXS	C2-C1-S	-3.33	108.14	113.25
7	A	614	CXS	C9-C4-C5	3.32	116.58	110.82

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	602	FAD	C4'-C3'-C2'	-3.32	106.45	113.36
4	A	606	TRT	C16-O15-C12	-3.23	109.48	117.93
5	A	608	NAD	C3D-C2D-C1D	3.19	105.78	100.98
4	A	606	TRT	C14-C9-C6	-3.15	115.75	121.53
2	A	601	FAD	C1B-N9A-C4A	-3.13	121.14	126.64
5	A	608	NAD	C3B-C2B-C1B	3.10	105.64	100.98
5	H	607	NAD	C3D-C2D-C1D	3.06	105.59	100.98
5	A	608	NAD	C1B-N9A-C4A	-3.01	121.35	126.64
5	H	607	NAD	C1B-N9A-C4A	-2.98	121.40	126.64
2	H	602	FAD	C5X-C9A-N10	2.84	119.77	117.72
7	H	613	CXS	C2-C1-S	-2.81	108.95	113.25
5	A	608	NAD	C5N-C6N-N1N	2.76	124.36	120.40
2	A	601	FAD	C1'-N10-C9A	2.75	120.46	118.29
7	H	613	CXS	C3-N-C4	-2.63	108.97	114.14
7	H	613	CXS	O1-S-C1	2.57	110.01	106.92
2	A	601	FAD	P-O3P-PA	-2.51	124.21	132.83
5	H	607	NAD	C2D-C3D-C4D	2.47	107.45	102.64
2	A	601	FAD	C4X-N5-C5X	2.46	119.23	116.77
5	A	608	NAD	C2D-C3D-C4D	2.44	107.39	102.64
4	A	606	TRT	C5-C6-C9	-2.43	106.55	111.93
2	H	602	FAD	P-O3P-PA	-2.42	124.53	132.83
5	A	608	NAD	C2B-C3B-C4B	2.33	107.17	102.64
2	A	601	FAD	C1'-N10-C10	2.24	120.42	118.41
7	A	614	CXS	O3-S-C1	2.20	109.32	105.77
4	H	609	TRT	C1-C5-C6	-2.14	116.21	123.91
2	A	601	FAD	C9A-N10-C10	-2.13	119.12	121.91
7	A	614	CXS	O2-S-C1	2.10	109.44	106.92
4	A	606	TRT	C1-C5-C6	-2.09	116.39	123.91
5	H	607	NAD	C2B-C3B-C4B	2.09	106.70	102.64
4	H	608	TRT	C10-C9-C6	-2.08	117.72	121.53
4	A	609	TRT	C10-C9-C14	2.04	121.02	117.97

There are no chirality outliers.

All (49) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	H	609	TRT	C1-C5-C6-C9
4	H	608	TRT	C1-C5-C6-C9
4	H	608	TRT	C1-C5-C6-C8
4	H	608	TRT	C1-C5-C6-C7
7	A	614	CXS	C2-C1-S-O2
7	A	614	CXS	C1-C2-C3-N

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Mol	Chain	Res	Type	Atoms
4	A	606	TRT	C1-C5-C6-C9
7	H	613	CXS	C1-C2-C3-N
4	H	601	TRT	C1-C5-C6-C9
4	H	601	TRT	C1-C5-C6-C7
4	H	609	TRT	C1-C5-C6-C7
4	A	606	TRT	C1-C5-C6-C8
4	H	609	TRT	C2-C1-C5-C6
4	H	609	TRT	C4-C1-C5-C6
4	H	609	TRT	C3-C1-C5-C6
4	H	601	TRT	C4-C1-C5-C6
4	A	607	TRT	O15-C16-C17-O18
4	H	608	TRT	C4-C1-C5-C6
4	H	608	TRT	C3-C1-C5-C6
4	H	601	TRT	C2-C1-C5-C6
7	A	614	CXS	C2-C1-S-O3
4	H	608	TRT	C13-C12-O15-C16
4	H	609	TRT	O15-C16-C17-O18
4	H	608	TRT	C11-C12-O15-C16
4	H	608	TRT	C17-C16-O15-C12
4	H	608	TRT	C2-C1-C5-C6
4	H	601	TRT	C3-C1-C5-C6
4	H	601	TRT	O15-C16-C17-O18
4	H	609	TRT	C1-C5-C6-C8
4	A	606	TRT	C1-C5-C6-C7
7	A	614	CXS	C2-C1-S-O1
5	A	608	NAD	O4B-C4B-C5B-O5B
4	H	601	TRT	C16-C17-O18-C19
4	A	607	TRT	C16-C17-O18-C19
4	A	607	TRT	C17-C16-O15-C12
4	A	609	TRT	C20-C19-O18-C17
2	A	601	FAD	O4B-C4B-C5B-O5B
4	A	607	TRT	C1-C5-C6-C9
4	A	609	TRT	O15-C16-C17-O18
2	H	602	FAD	O4B-C4B-C5B-O5B
5	H	607	NAD	O4B-C4B-C5B-O5B
4	H	608	TRT	C7-C6-C9-C14
4	H	608	TRT	C8-C6-C9-C10
5	A	608	NAD	C3B-C4B-C5B-O5B
4	H	608	TRT	C7-C6-C9-C10
4	A	607	TRT	C3-C1-C5-C6
4	A	606	TRT	C20-C19-O18-C17
4	A	607	TRT	C4-C1-C5-C6

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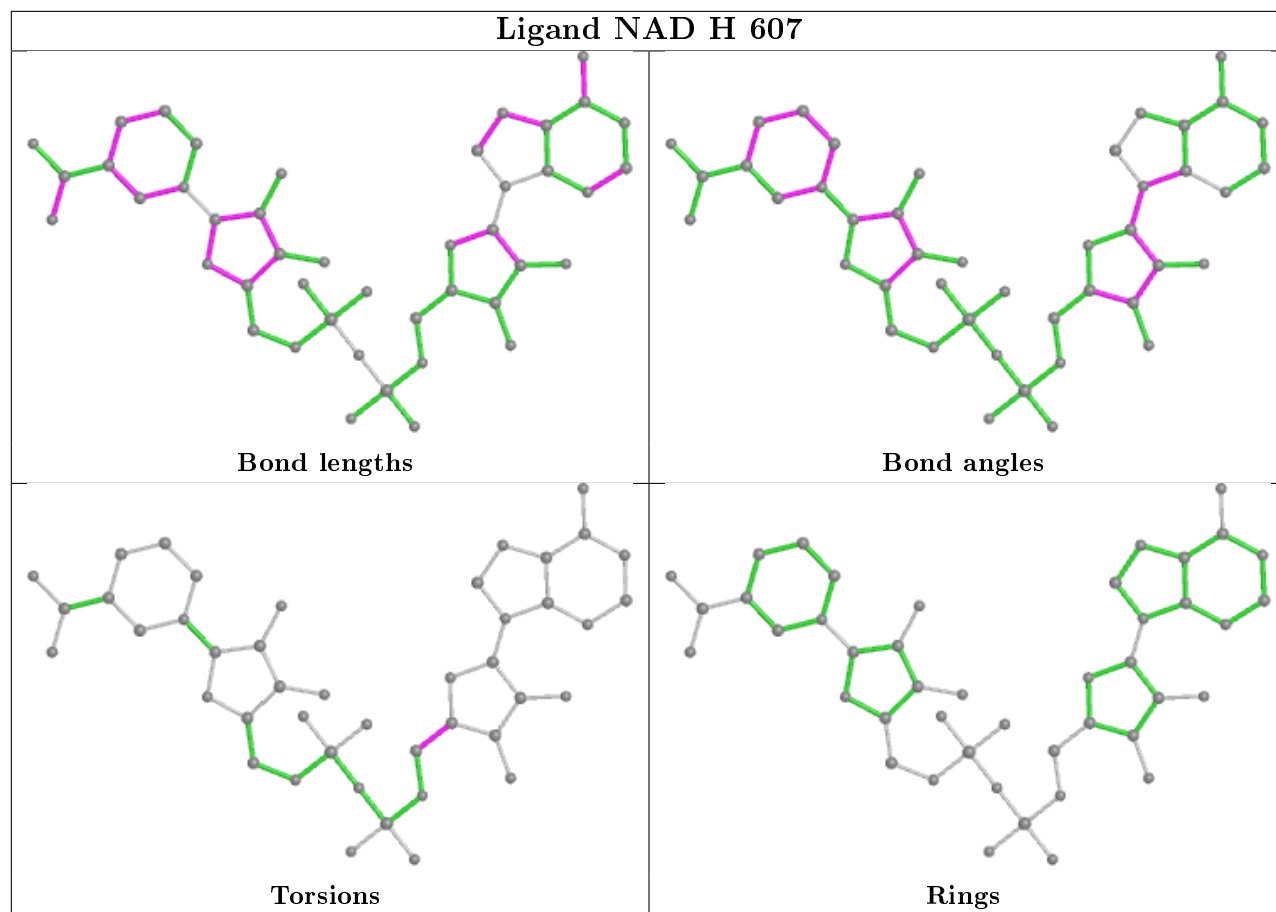
Mol	Chain	Res	Type	Atoms
4	H	608	TRT	C8-C6-C9-C14

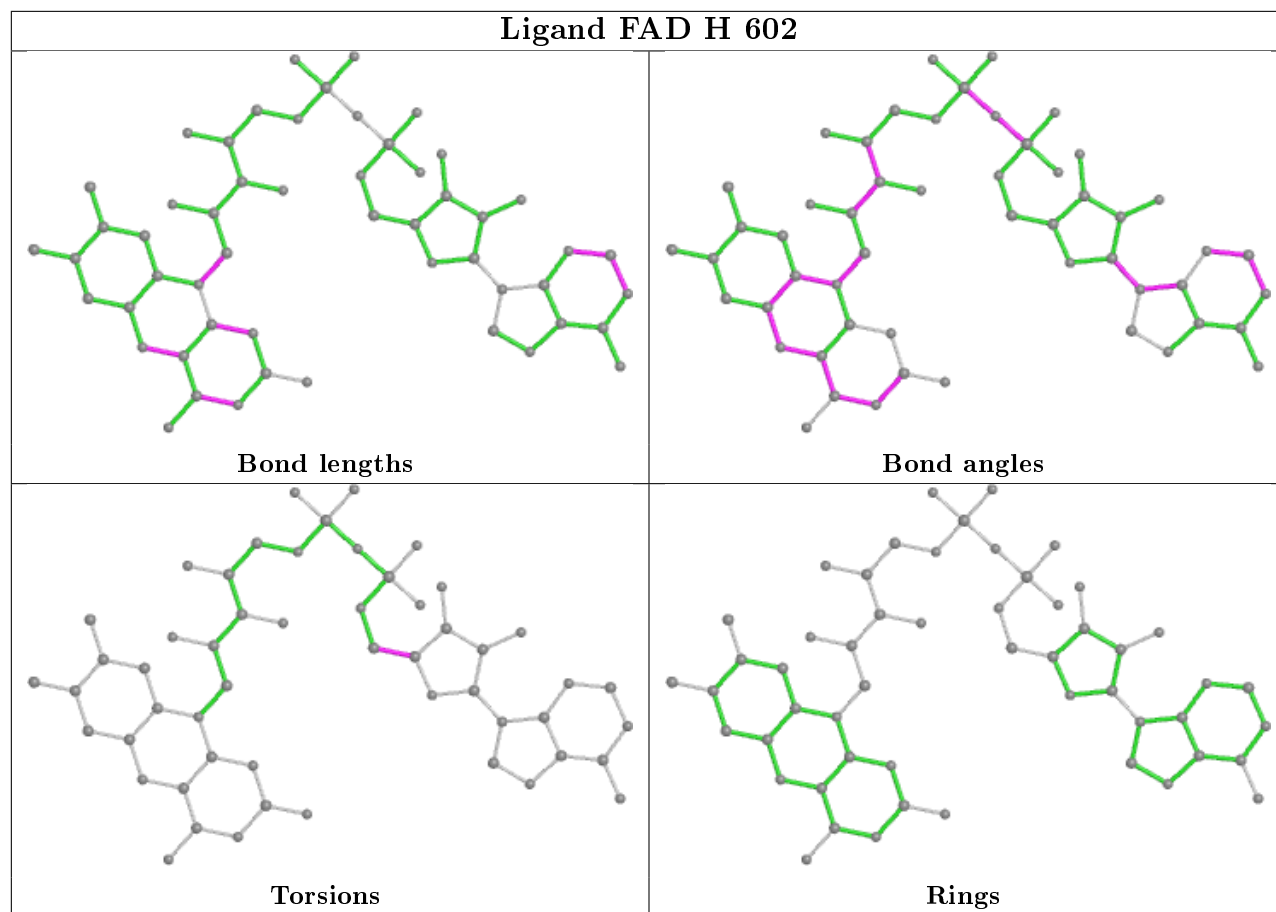
There are no ring outliers.

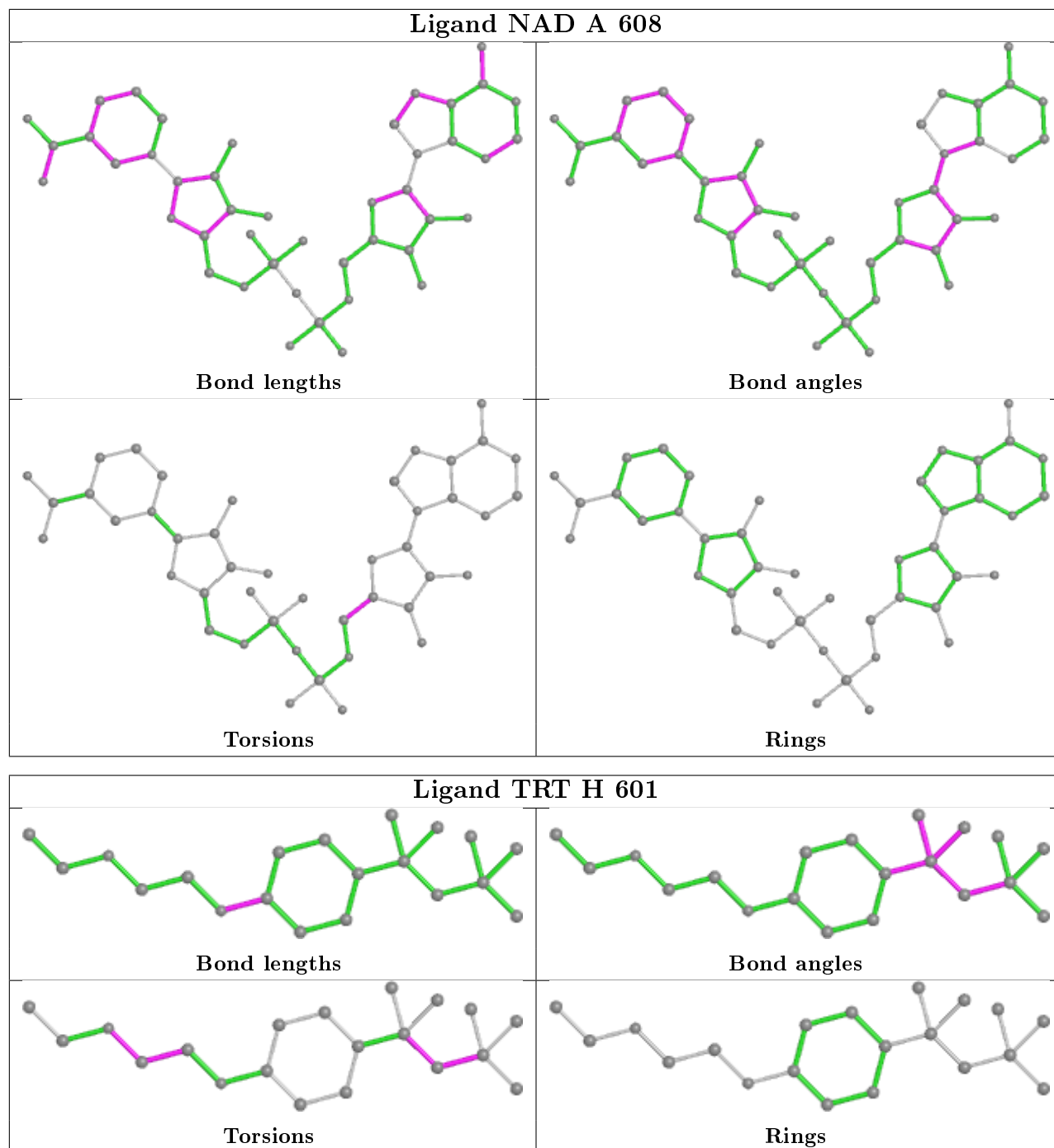
12 monomers are involved in 60 short contacts:

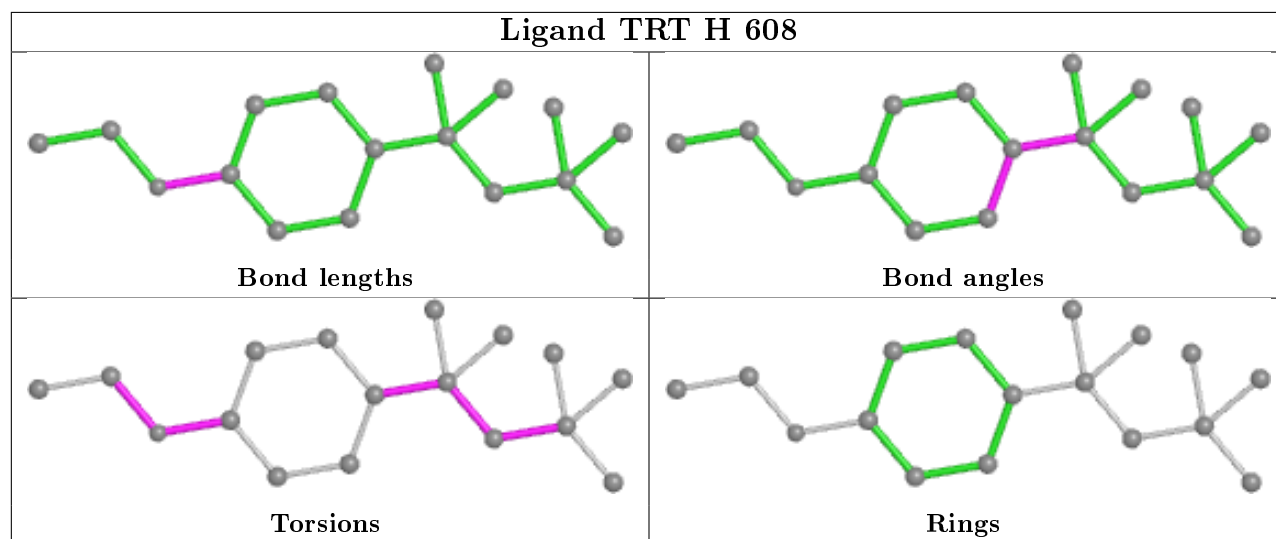
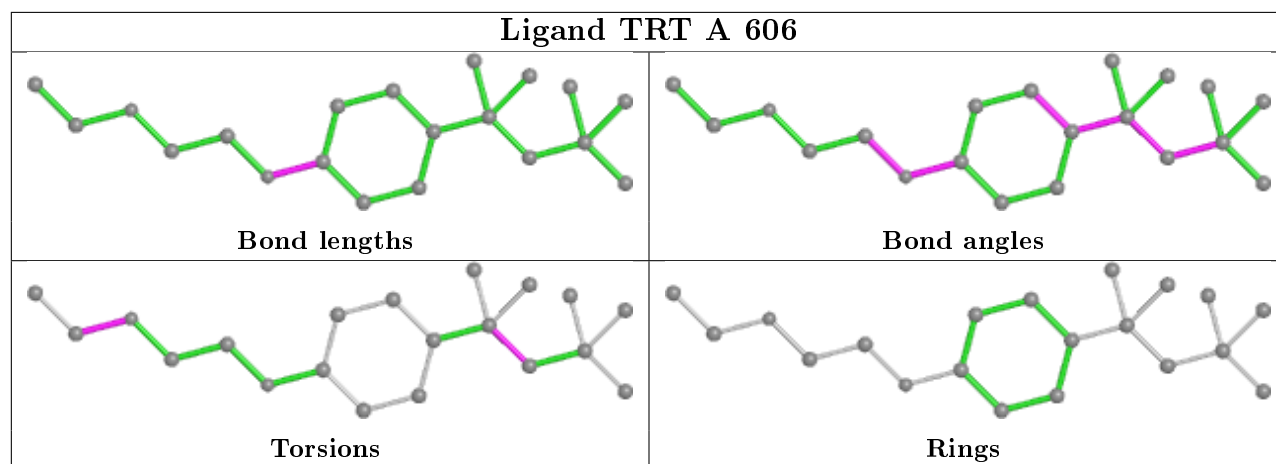
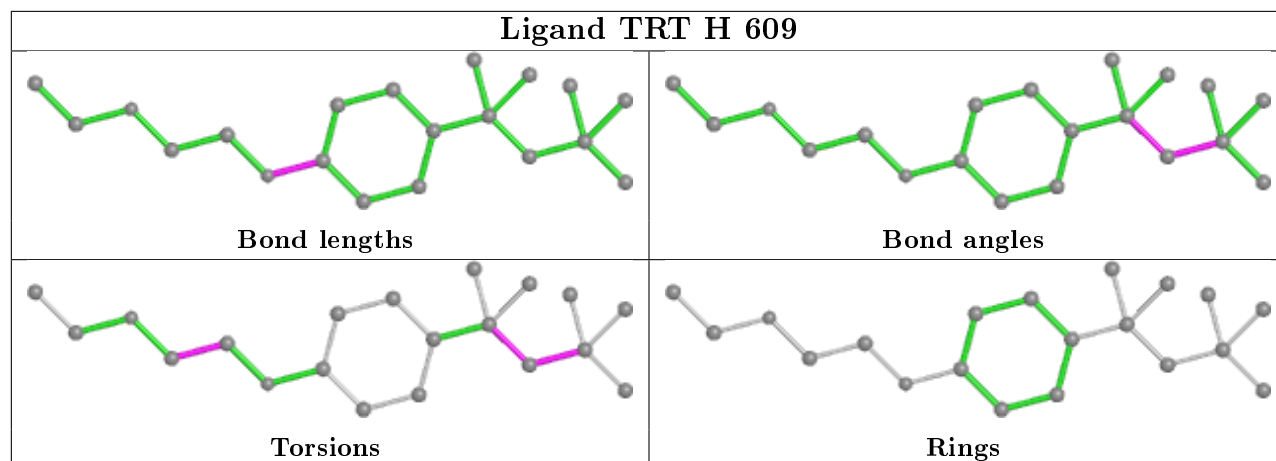
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	H	607	NAD	1	0
2	H	602	FAD	2	0
5	A	608	NAD	1	0
4	H	601	TRT	5	0
4	H	609	TRT	15	0
4	A	606	TRT	10	0
4	H	608	TRT	16	0
2	A	601	FAD	1	0
7	H	613	CXS	1	0
7	A	614	CXS	2	0
4	A	607	TRT	3	0
4	A	609	TRT	12	0

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.

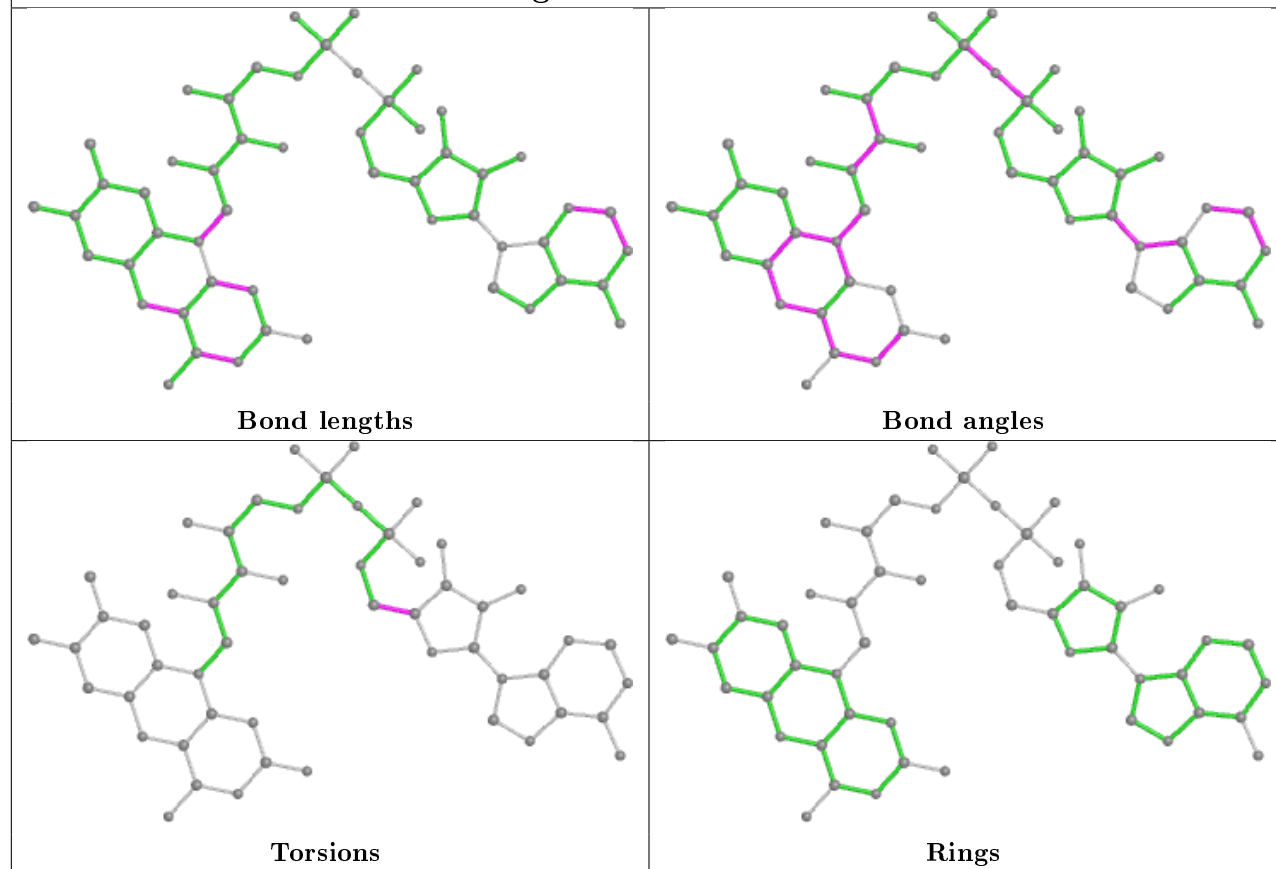




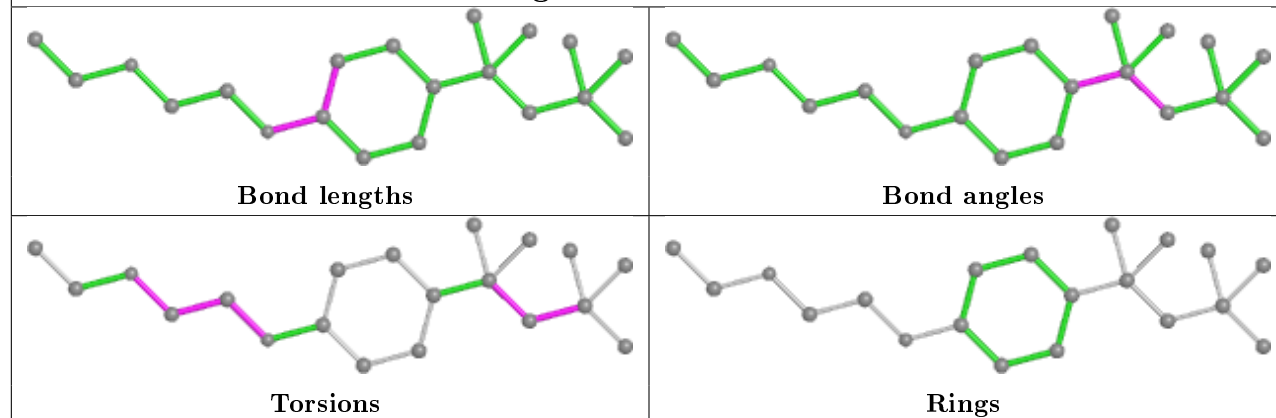


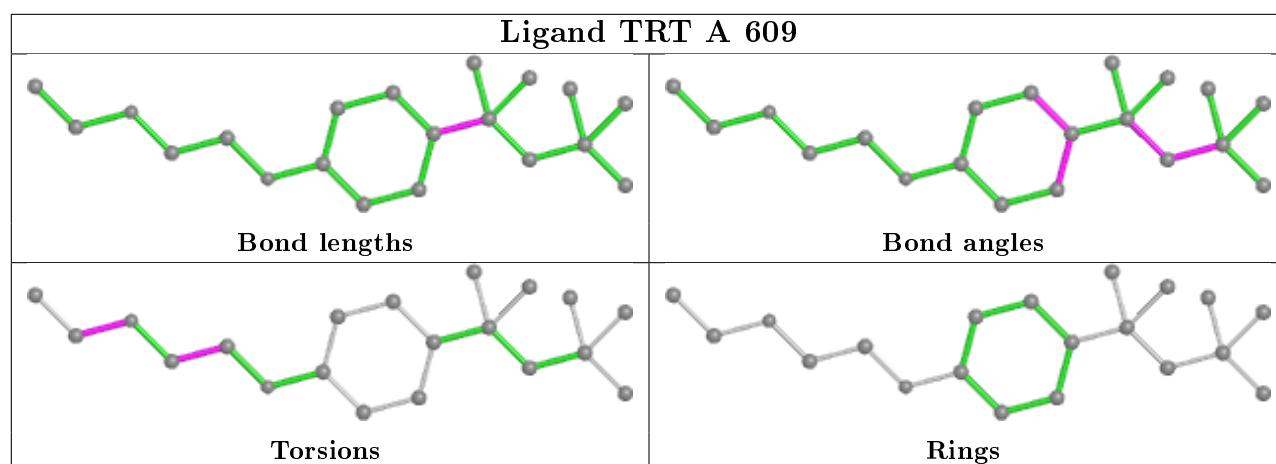


Ligand FAD A 601



Ligand TRT A 607





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	493/521 (94%)	0.04	1 (0%) 95 96	29, 44, 66, 89	0
1	H	492/521 (94%)	0.05	2 (0%) 92 93	29, 43, 66, 88	0
All	All	985/1042 (94%)	0.05	3 (0%) 94 95	29, 43, 66, 89	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	457	ASN	3.9
1	H	106	ASN	3.7
1	H	105	LYS	2.3

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

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
6	ACT	H	610	4/4	0.74	0.17	45,56,60,68	0
6	ACT	A	611	4/4	0.81	0.28	56,65,66,70	0

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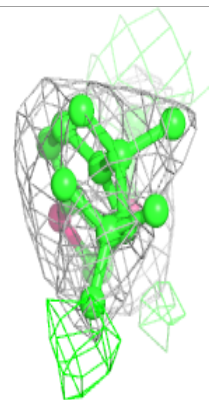
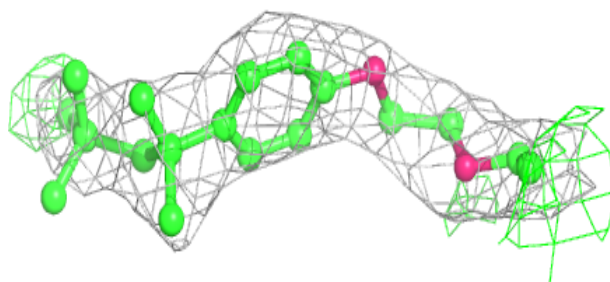
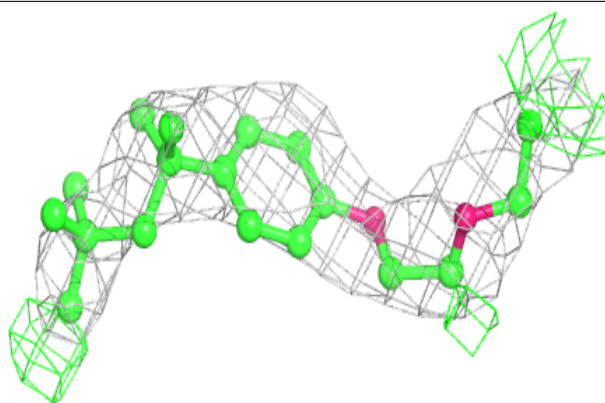
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
6	ACT	H	612	4/4	0.82	0.39	69,74,88,89	0
4	TRT	A	606	20/25	0.86	0.29	51,65,75,78	0
4	TRT	A	607	20/25	0.87	0.24	52,67,75,78	0
4	TRT	H	608	17/25	0.88	0.37	48,61,75,77	0
4	TRT	H	609	20/25	0.89	0.38	53,66,74,77	0
6	ACT	A	610	4/4	0.90	0.14	49,50,62,65	0
6	ACT	A	612	4/4	0.90	0.19	52,66,70,70	0
7	CXS	H	613	14/14	0.92	0.17	46,54,76,78	0
4	TRT	H	601	20/25	0.92	0.28	55,62,66,66	0
7	CXS	A	614	14/14	0.93	0.21	47,57,80,80	0
3	MG	A	604	1/1	0.93	0.21	45,45,45,45	0
6	ACT	A	613	4/4	0.94	0.11	51,54,63,66	0
4	TRT	A	609	20/25	0.94	0.30	55,62,65,65	0
3	MG	A	605	1/1	0.94	0.47	47,47,47,47	0
5	NAD	H	607	44/44	0.96	0.13	37,45,55,63	0
3	MG	H	606	1/1	0.96	0.31	50,50,50,50	0
6	ACT	H	611	4/4	0.96	0.20	47,51,58,59	0
3	MG	H	605	1/1	0.96	0.26	45,45,45,45	0
5	NAD	A	608	44/44	0.97	0.13	37,45,52,54	0
3	MG	H	603	1/1	0.98	0.20	43,43,43,43	0
3	MG	A	602	1/1	0.98	0.28	45,45,45,45	0
2	FAD	H	602	53/53	0.98	0.17	26,32,37,40	0
3	MG	A	603	1/1	0.98	0.15	37,37,37,37	0
3	MG	H	604	1/1	0.98	0.14	39,39,39,39	0
2	FAD	A	601	53/53	0.99	0.17	25,33,36,38	0

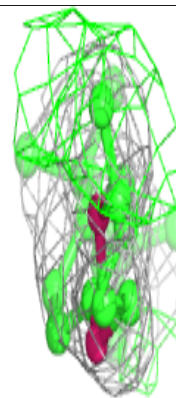
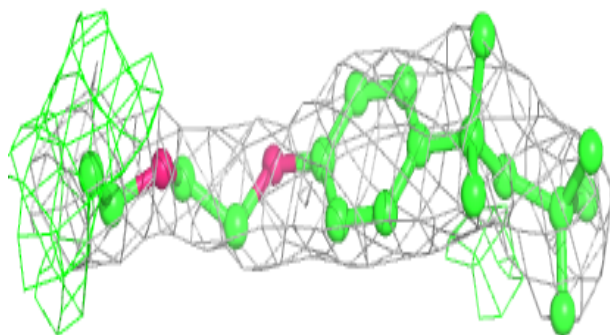
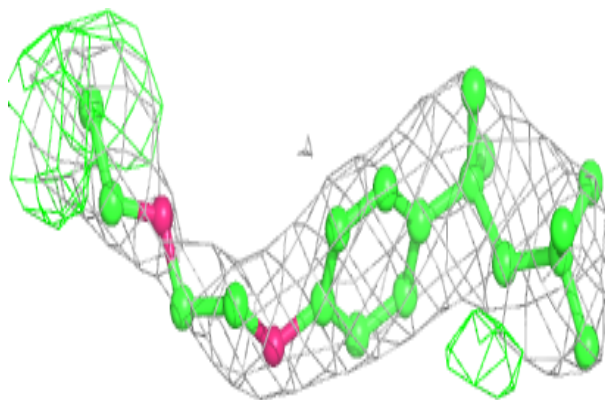
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around TRT A 606:

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

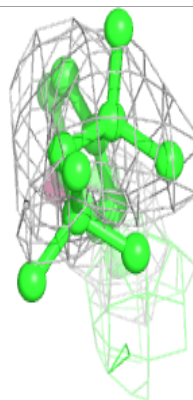
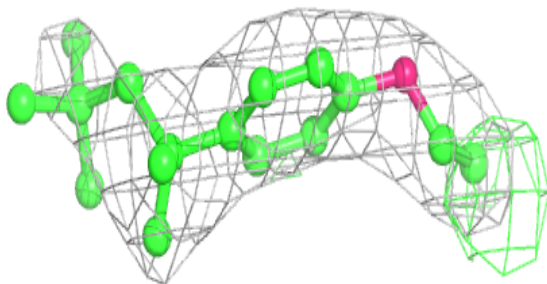
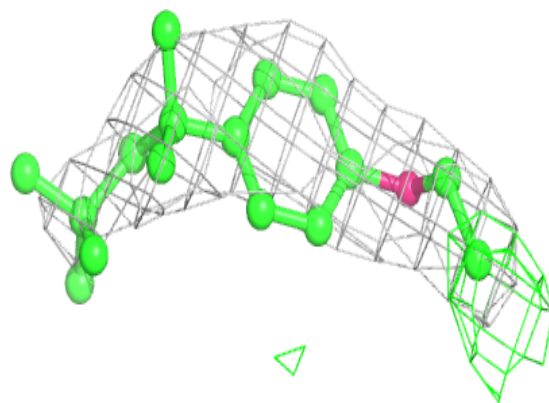
**Electron density around TRT A 607:**

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

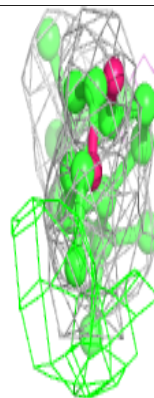
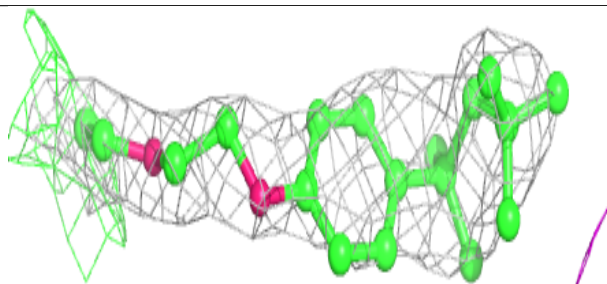
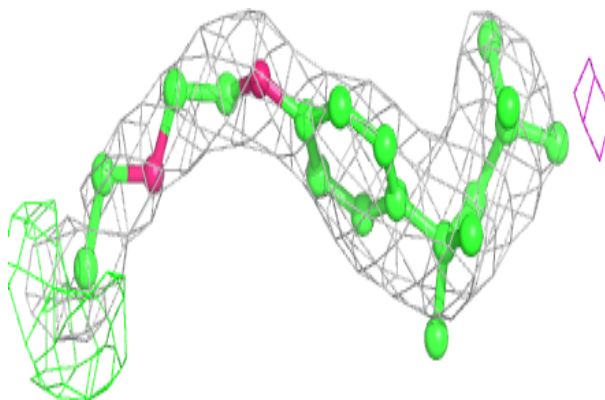


Electron density around TRT H 608:

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

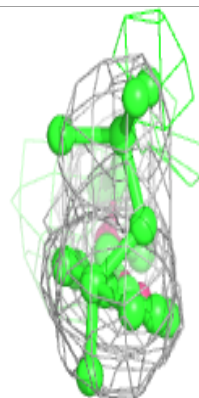
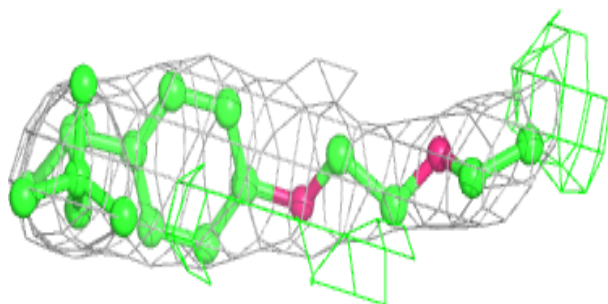
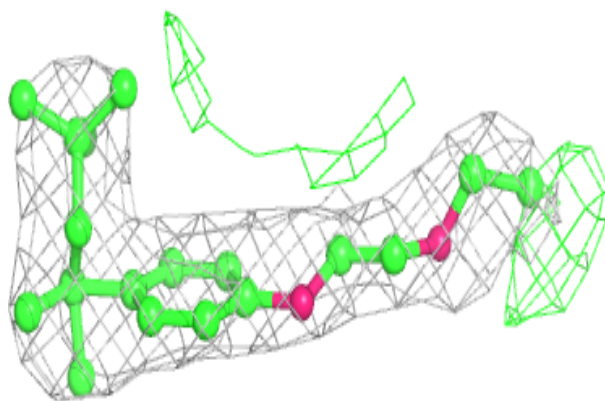
**Electron density around TRT H 609:**

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

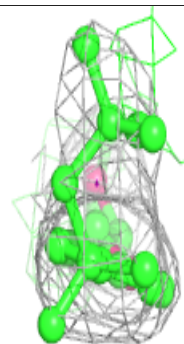
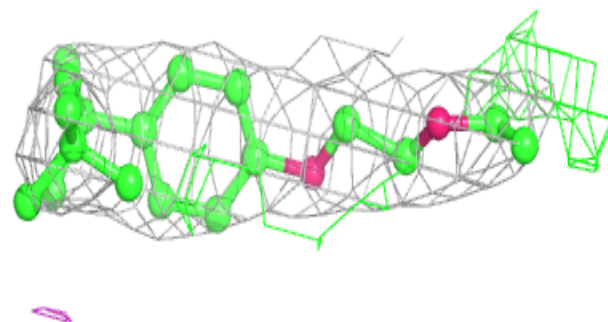
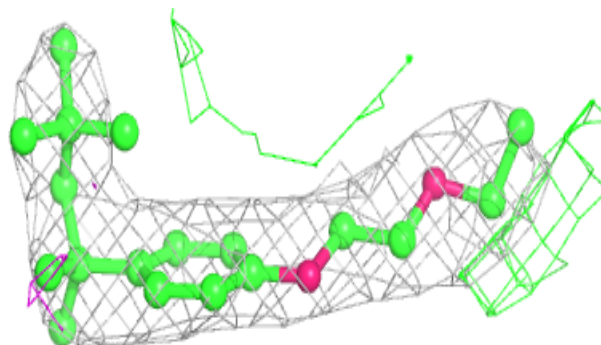


Electron density around TRT H 601:

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

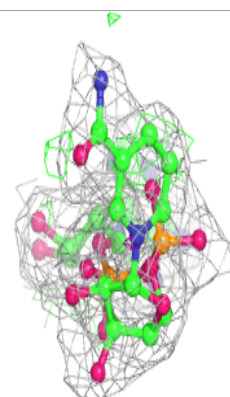
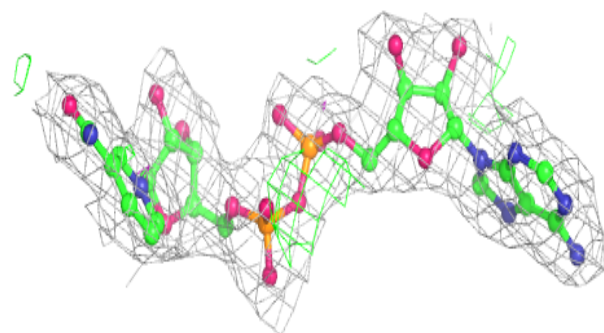
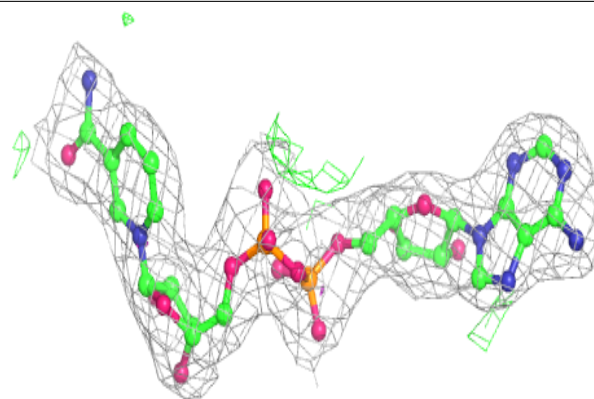
**Electron density around TRT A 609:**

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

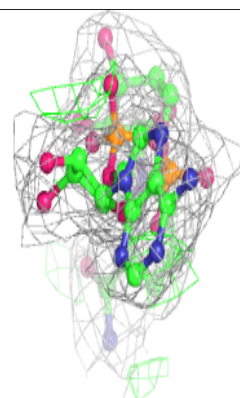
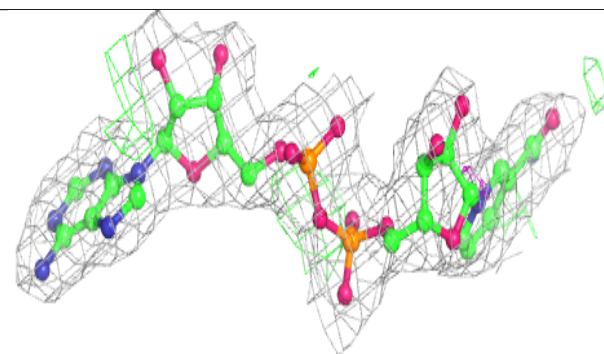
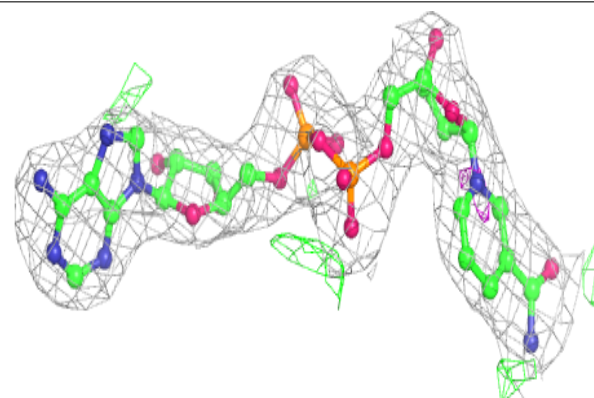


Electron density around NAD H 607:

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

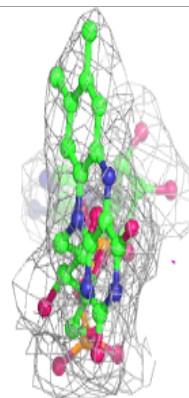
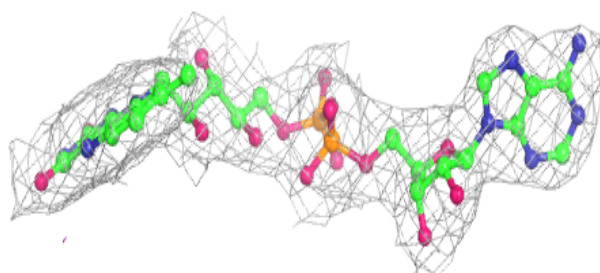
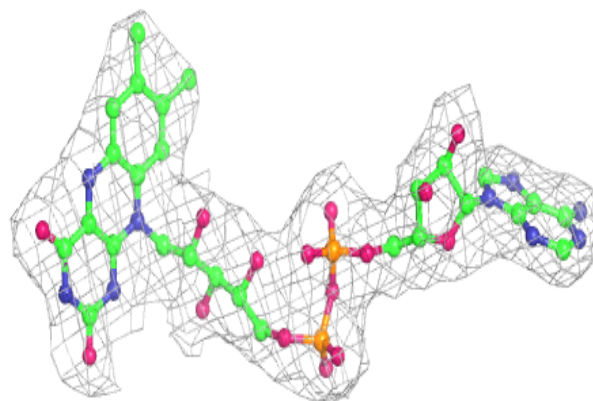
**Electron density around NAD A 608:**

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

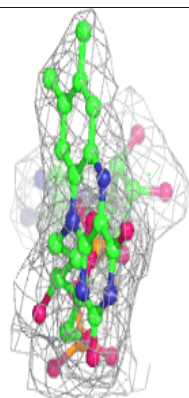
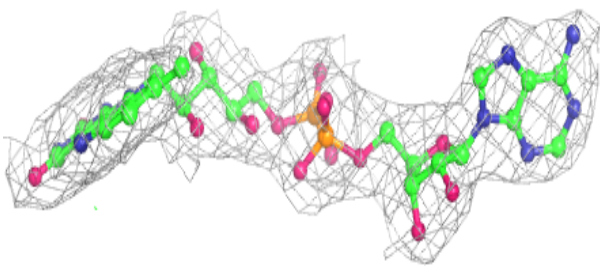
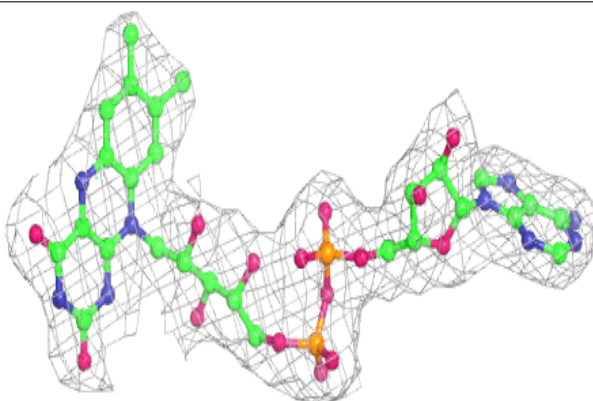


Electron density around FAD H 602:

$2mF_o-DF_c$ (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

There are no such residues in this entry.