



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 5, 2026 – 11:35 AM UTC

PDB ID : 1BVR / pdb_00001bvr
Title : M.TB. ENOYL-ACP REDUCTASE (INHA) IN COMPLEX WITH NAD+ AND C16-FATTY-ACYL-SUBSTRATE
Authors : Rozwarski, D.A.; Vilcheze, C.; Sugantino, M.; Bittman, R.; Jacobs, W.; TB Structural Genomics Consortium (TBSGC)
Deposited on : 1998-09-17
Resolution : 2.80 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

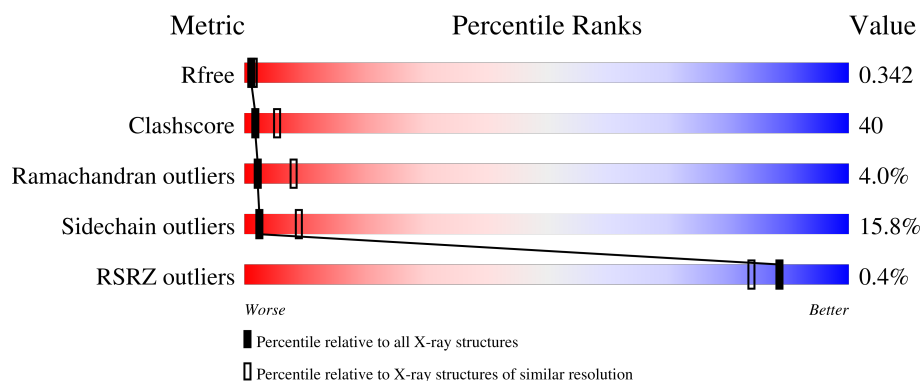
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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	268	46% 44% 9% .
1	B	268	49% 41% 9% .
1	C	268	45% 46% 8% .
1	D	268	44% 42% 13% .

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
1	E	268	 17% 44% 32% 7%
1	F	268	 43% 46% 10% •

2 Entry composition [i](#)

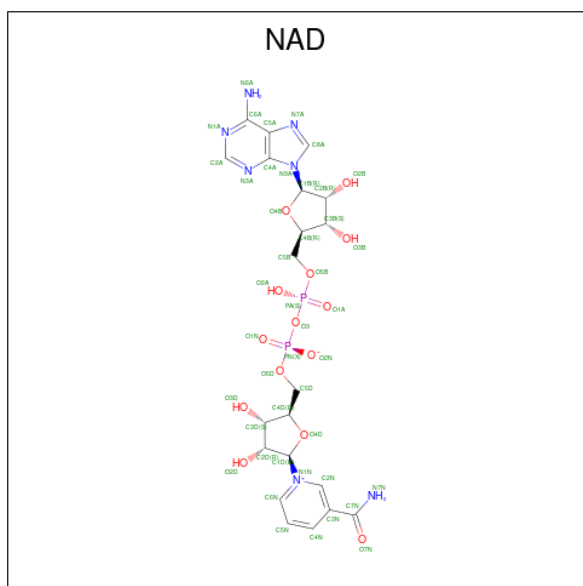
There are 4 unique types of molecules in this entry. The entry contains 12987 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 PROTEIN (ENOYL-ACYL CARRIER PROTEIN (ACP) REDUCTASE).

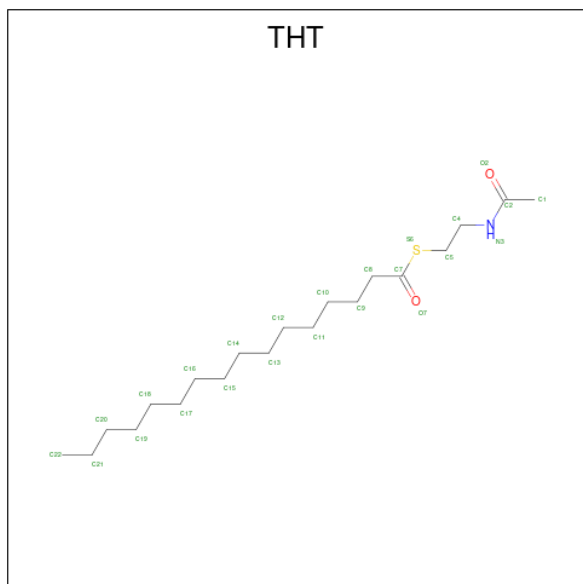
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	268	Total	C	N	O	S	0	0	0
			1994	1263	348	373	10			
1	B	268	Total	C	N	O	S	0	0	0
			1994	1263	348	373	10			
1	C	268	Total	C	N	O	S	0	0	0
			1994	1263	348	373	10			
1	D	268	Total	C	N	O	S	0	0	0
			1994	1263	348	373	10			
1	E	268	Total	C	N	O	S	0	0	0
			1994	1263	348	373	10			
1	F	268	Total	C	N	O	S	0	0	0
			1994	1263	348	373	10			

- Molecule 2 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (CCD ID: NAD) (formula: $C_{21}H_{27}N_7O_{14}P_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	B	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	C	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	D	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	E	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	F	1	Total	C	N	O	P	0	0
			44	21	7	14	2		

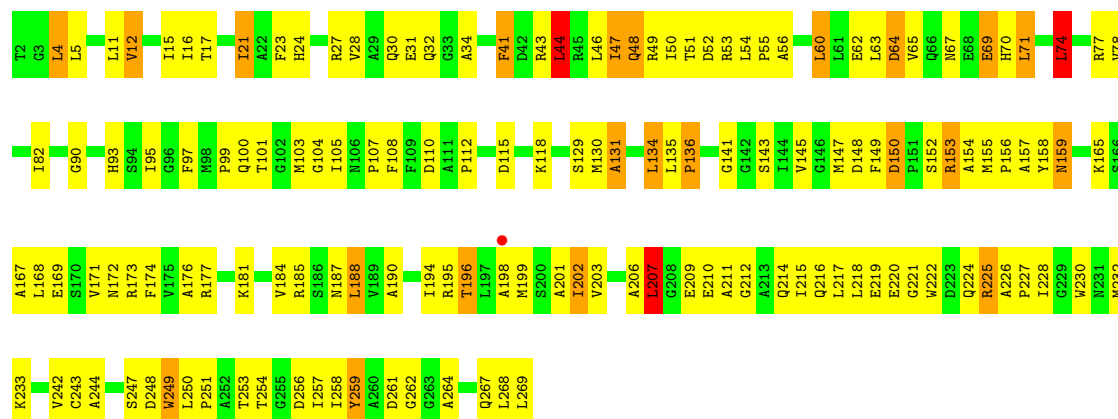
- Molecule 3 is TRANS-2-HEXADECENOYL-(N-ACETYL-CYSTEAMINE)-THIOESTER (CCD ID: THT) (formula: $C_{20}H_{39}NO_2S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	S	0	0
			24	20	1	2	1		
3	B	1	Total	C	N	O	S	0	0
			24	20	1	2	1		
3	C	1	Total	C	N	O	S	0	0
			24	20	1	2	1		
3	F	1	Total	C	N	O	S	0	0
			24	20	1	2	1		

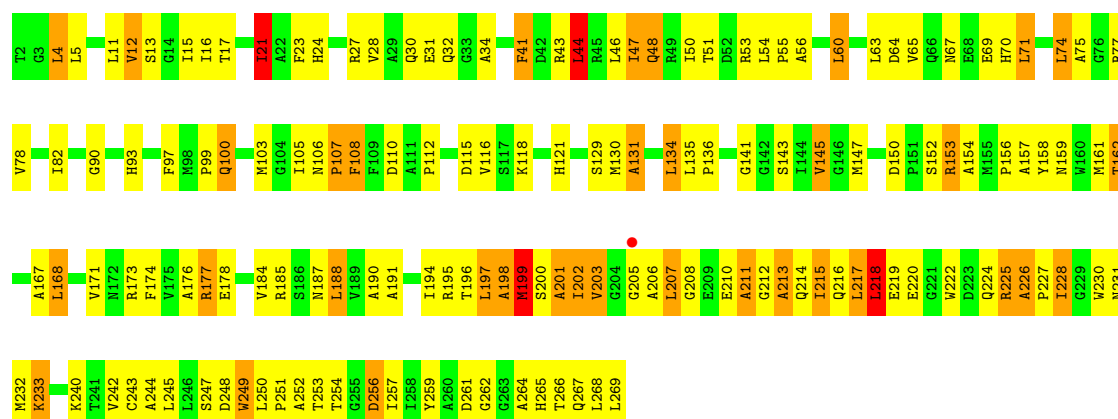
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	110	Total 110	O 110	0	0
4	B	114	Total 114	O 114	0	0
4	C	106	Total 106	O 106	0	0
4	D	104	Total 104	O 104	0	0
4	E	119	Total 119	O 119	0	0
4	F	110	Total 110	O 110	0	0



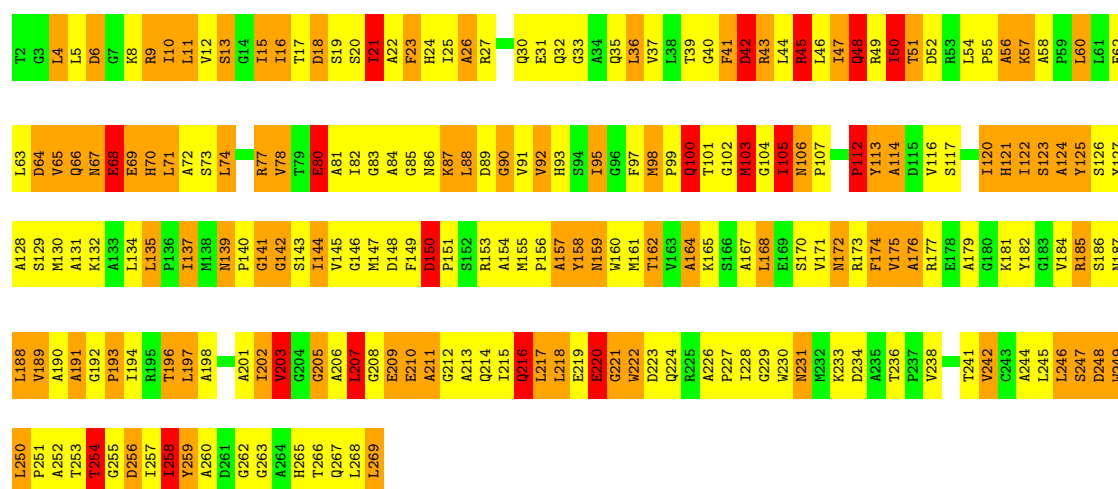
• Molecule 1: PROTEIN (ENOYL-ACYL CARRIER PROTEIN (ACP) REDUCTASE)

Chain D: 44% 42% 13%

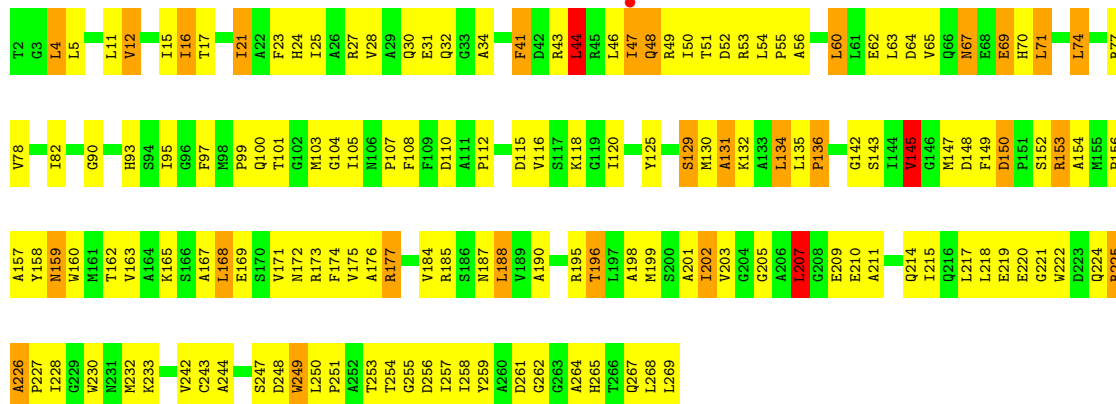


• Molecule 1: PROTEIN (ENOYL-ACYL CARRIER PROTEIN (ACP) REDUCTASE)

Chain E: 17% 44% 32% 7%



• Molecule 1: PROTEIN (ENOYL-ACYL CARRIER PROTEIN (ACP) REDUCTASE)

Chain F:  43% 46% 10%

4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	101.08Å 83.45Å 192.81Å 90.00° 95.24° 90.00°	Depositor
Resolution (Å)	10.00 – 2.80 10.00 – 2.80	Depositor EDS
% Data completeness (in resolution range)	90.2 (10.00-2.80) 88.0 (10.00-2.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.16	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.29 (at 2.80Å)	Xtriage
Refinement program	X-PLOR	Depositor
R, R_{free}	0.217 , 0.344 0.218 , 0.342	Depositor DCC
R_{free} test set	3460 reflections (9.93%)	wwPDB-VP
Wilson B-factor (Å ²)	51.7	Xtriage
Anisotropy	0.129	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.12 , 107.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.36$, $\langle L^2 \rangle = 0.18$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	12987	wwPDB-VP
Average B, all atoms (Å ²)	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.70% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

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

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	0.77	0/2032	1.16	16/2758 (0.6%)
1	B	0.78	0/2032	1.19	17/2758 (0.6%)
1	C	0.76	0/2032	1.19	16/2758 (0.6%)
1	D	0.83	2/2032 (0.1%)	1.25	22/2758 (0.8%)
1	E	1.12	4/2032 (0.2%)	1.69	61/2758 (2.2%)
1	F	0.73	1/2032 (0.0%)	1.18	18/2758 (0.7%)
All	All	0.84	7/12192 (0.1%)	1.29	150/16548 (0.9%)

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	E	0	1

The worst 5 of 7 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	15	ILE	CA-CB	8.13	1.63	1.54
1	D	202	ILE	CA-CB	6.76	1.62	1.54
1	E	103	MET	SD-CE	6.12	1.94	1.79
1	D	145	VAL	CA-CB	5.96	1.62	1.54
1	E	258	ILE	CA-CB	5.27	1.60	1.54

The worst 5 of 150 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	87	LYS	N-CA-C	11.03	124.56	110.33
1	E	135	LEU	N-CA-C	9.76	126.02	112.75

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	69	GLU	N-CA-C	-9.05	101.49	112.54
1	E	255	GLY	N-CA-C	-8.82	98.50	114.46
1	E	78	VAL	N-CA-C	-8.81	102.24	111.58

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	E	113	TYR	Sidechain

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	1994	0	2008	131	0
1	B	1994	0	2008	125	0
1	C	1994	0	2008	153	0
1	D	1994	0	2008	144	0
1	E	1994	0	2008	354	0
1	F	1994	0	2008	146	0
2	A	44	0	26	2	0
2	B	44	0	26	1	0
2	C	44	0	26	1	0
2	D	44	0	26	2	0
2	E	44	0	25	9	0
2	F	44	0	26	1	0
3	A	24	0	39	17	0
3	B	24	0	39	16	0
3	C	24	0	39	17	0
3	F	24	0	39	17	0
4	A	110	0	0	13	0
4	B	114	0	0	15	0
4	C	106	0	0	11	0
4	D	104	0	0	11	0
4	E	119	0	0	47	0
4	F	110	0	0	13	0
All	All	12987	0	12359	988	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 40.

The worst 5 of 988 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:F:198:ALA:HB2	3:F:1310:THT:HC12	1.22	1.15
1:A:198:ALA:HB2	3:A:1302:THT:HC12	1.18	1.08
1:B:198:ALA:HB2	3:B:1304:THT:HC12	1.25	1.08
1:C:198:ALA:HB2	3:C:1306:THT:HC12	1.19	1.08
1:E:135:LEU:HD11	1:E:144:ILE:HD11	1.38	1.02

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	266/268 (99%)	238 (90%)	22 (8%)	6 (2%)	5	18
1	B	266/268 (99%)	238 (90%)	21 (8%)	7 (3%)	4	15
1	C	266/268 (99%)	241 (91%)	18 (7%)	7 (3%)	4	15
1	D	266/268 (99%)	231 (87%)	27 (10%)	8 (3%)	3	12
1	E	266/268 (99%)	190 (71%)	46 (17%)	30 (11%)	0	1
1	F	266/268 (99%)	239 (90%)	21 (8%)	6 (2%)	5	18
All	All	1596/1608 (99%)	1377 (86%)	155 (10%)	64 (4%)	2	8

5 of 64 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	209	GLU
1	D	103	MET
1	D	197	LEU

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	E	21	ILE
1	E	56	ALA

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	203/204 (100%)	176 (87%)	27 (13%)	4	14
1	B	203/204 (100%)	176 (87%)	27 (13%)	4	14
1	C	203/204 (100%)	177 (87%)	26 (13%)	4	15
1	D	203/204 (100%)	176 (87%)	27 (13%)	4	14
1	E	203/204 (100%)	143 (70%)	60 (30%)	0	1
1	F	203/204 (100%)	177 (87%)	26 (13%)	4	15
All	All	1218/1224 (100%)	1025 (84%)	193 (16%)	2	9

5 of 193 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	E	36	LEU
1	E	139	ASN
1	E	47	ILE
1	E	77	ARG
1	E	185	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 36 such sidechains are listed below:

Mol	Chain	Res	Type
1	E	121	HIS
1	F	214	GLN
1	E	172	ASN
1	E	231	ASN
1	C	70	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 ⓘ

10 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$
3	THT	C	1306	-	23,23,23	1.53	4 (17%)	23,24,24	2.01	4 (17%)
2	NAD	F	1309	-	46,48,48	1.52	6 (13%)	64,73,73	2.06	10 (15%)
2	NAD	A	1301	-	46,48,48	1.55	7 (15%)	64,73,73	2.06	11 (17%)
3	THT	A	1302	-	23,23,23	1.16	2 (8%)	23,24,24	1.92	4 (17%)
3	THT	B	1304	-	23,23,23	1.39	4 (17%)	23,24,24	2.00	5 (21%)
2	NAD	D	1307	-	46,48,48	1.88	7 (15%)	64,73,73	2.12	16 (25%)
2	NAD	E	1308	-	46,48,48	2.02	9 (19%)	64,73,73	2.65	26 (40%)
2	NAD	C	1305	-	46,48,48	1.57	5 (10%)	64,73,73	2.05	13 (20%)
2	NAD	B	1303	-	46,48,48	1.52	5 (10%)	64,73,73	2.15	15 (23%)
3	THT	F	1310	-	23,23,23	1.20	3 (13%)	23,24,24	1.82	4 (17%)

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
3	THT	C	1306	-	-	14/22/22/22	-
2	NAD	F	1309	-	-	13/30/62/62	0/5/5/5
2	NAD	A	1301	-	-	13/30/62/62	0/5/5/5
3	THT	A	1302	-	-	15/22/22/22	-
3	THT	B	1304	-	-	14/22/22/22	-
2	NAD	D	1307	-	-	13/30/62/62	0/5/5/5
2	NAD	E	1308	-	-	18/30/62/62	0/5/5/5
2	NAD	C	1305	-	-	13/30/62/62	0/5/5/5
2	NAD	B	1303	-	-	12/30/62/62	0/5/5/5
3	THT	F	1310	-	-	14/22/22/22	-

The worst 5 of 52 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	1308	NAD	PA-O3	7.69	1.67	1.59
2	D	1307	NAD	C2N-N1N	7.01	1.42	1.35
2	A	1301	NAD	C2N-N1N	6.37	1.42	1.35
2	B	1303	NAD	C2N-N1N	6.21	1.41	1.35
2	C	1305	NAD	C2N-N1N	6.05	1.41	1.35

The worst 5 of 108 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	1308	NAD	C4B-O4B-C1B	-9.80	87.84	109.47
2	E	1308	NAD	O4B-C1B-N9A	8.07	123.59	108.09
2	B	1303	NAD	C5A-C4A-N3A	-7.07	116.98	126.72
2	B	1303	NAD	O4B-C1B-N9A	6.83	121.21	108.09
2	C	1305	NAD	C5A-C4A-N3A	-6.75	117.42	126.72

There are no chirality outliers.

5 of 139 torsion outliers are listed below:

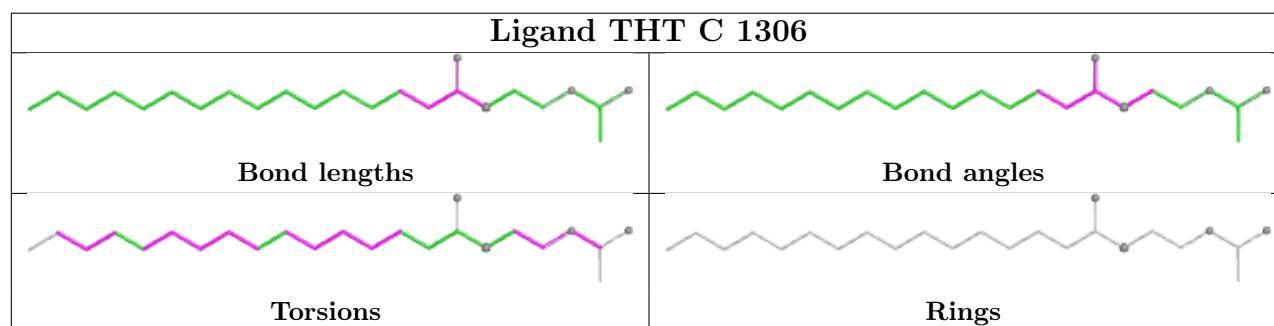
Mol	Chain	Res	Type	Atoms
2	A	1301	NAD	C5B-O5B-PA-O1A
2	A	1301	NAD	C5B-O5B-PA-O3
2	A	1301	NAD	O4D-C1D-N1N-C2N
2	A	1301	NAD	O4D-C1D-N1N-C6N
2	A	1301	NAD	C2D-C1D-N1N-C2N

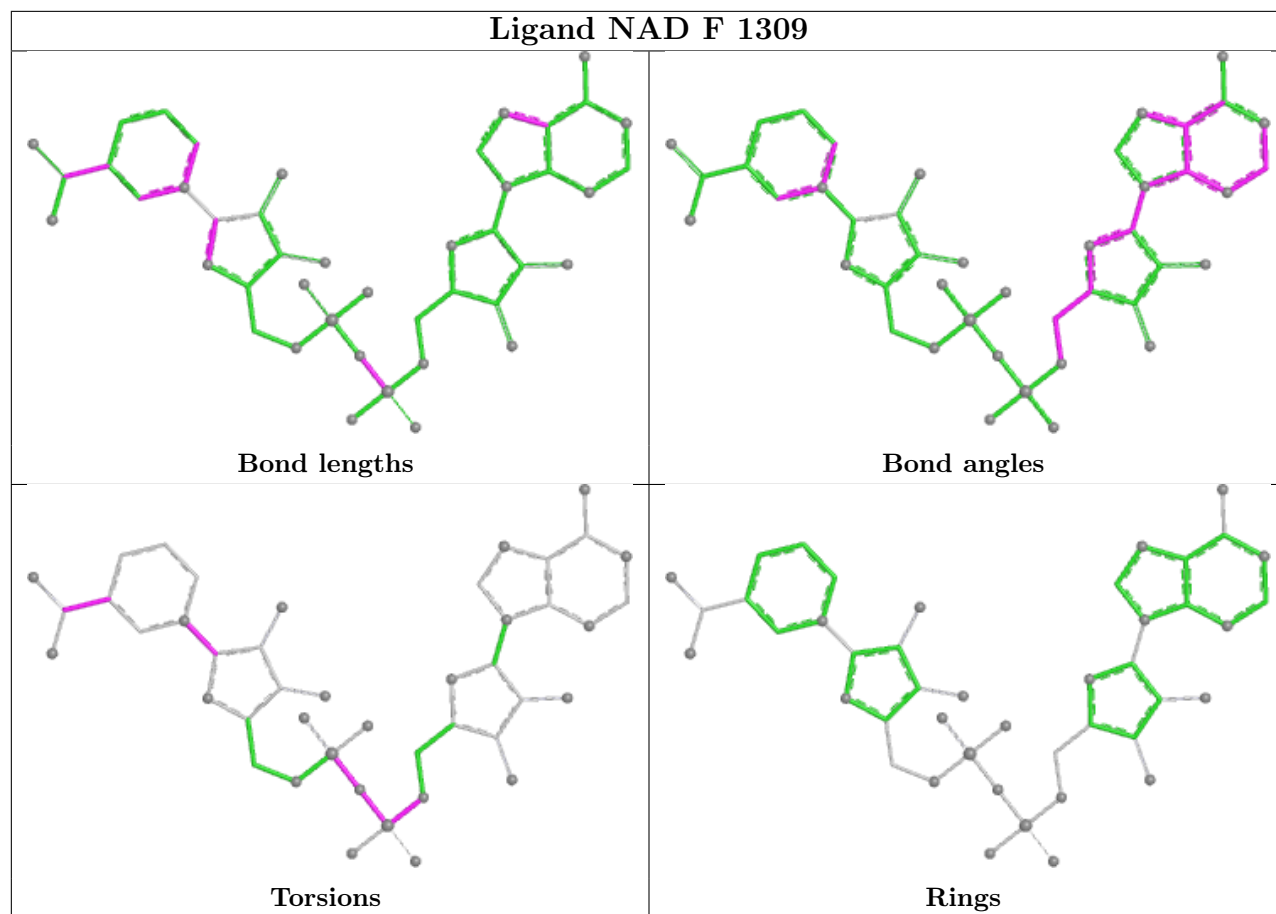
There are no ring outliers.

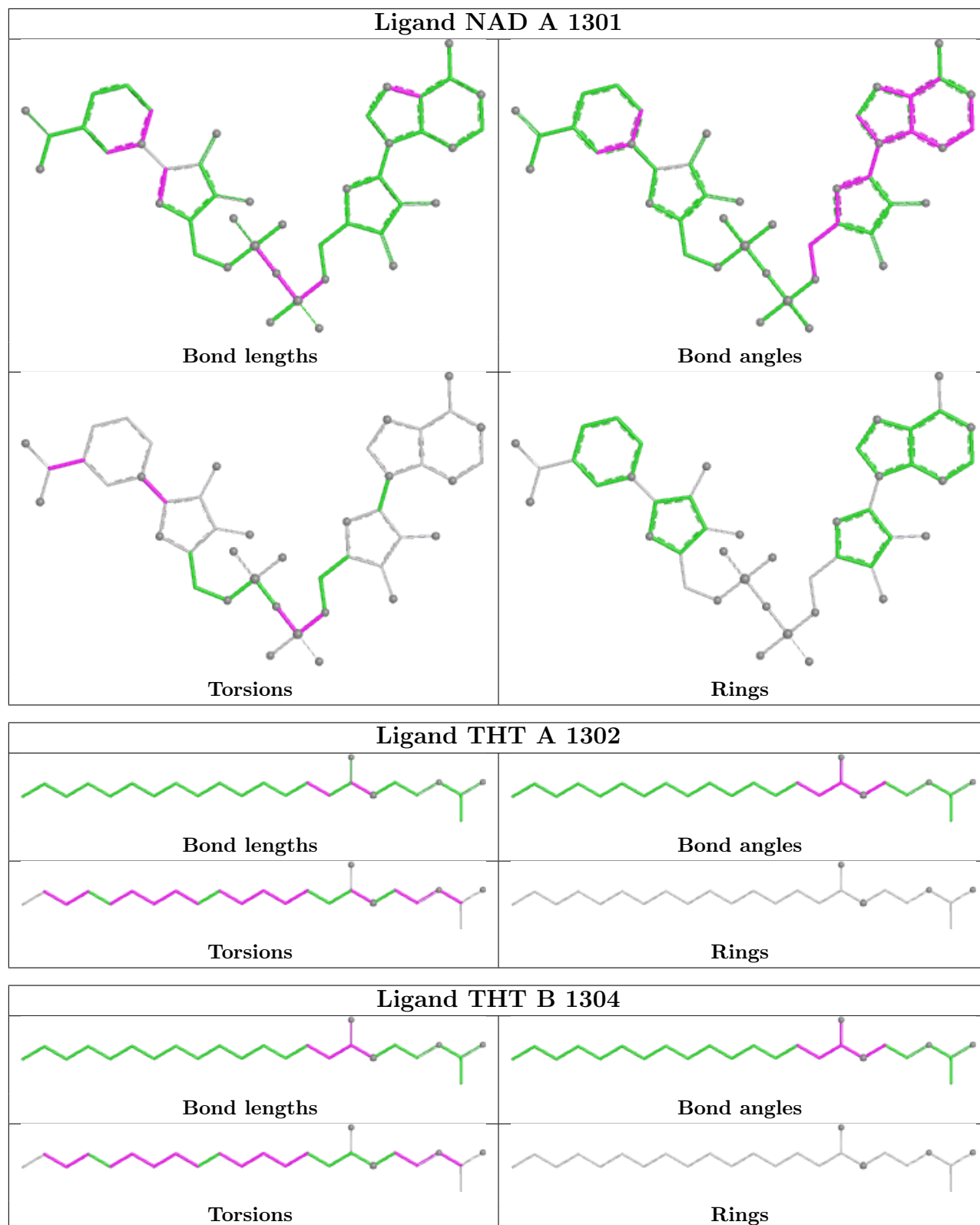
10 monomers are involved in 83 short contacts:

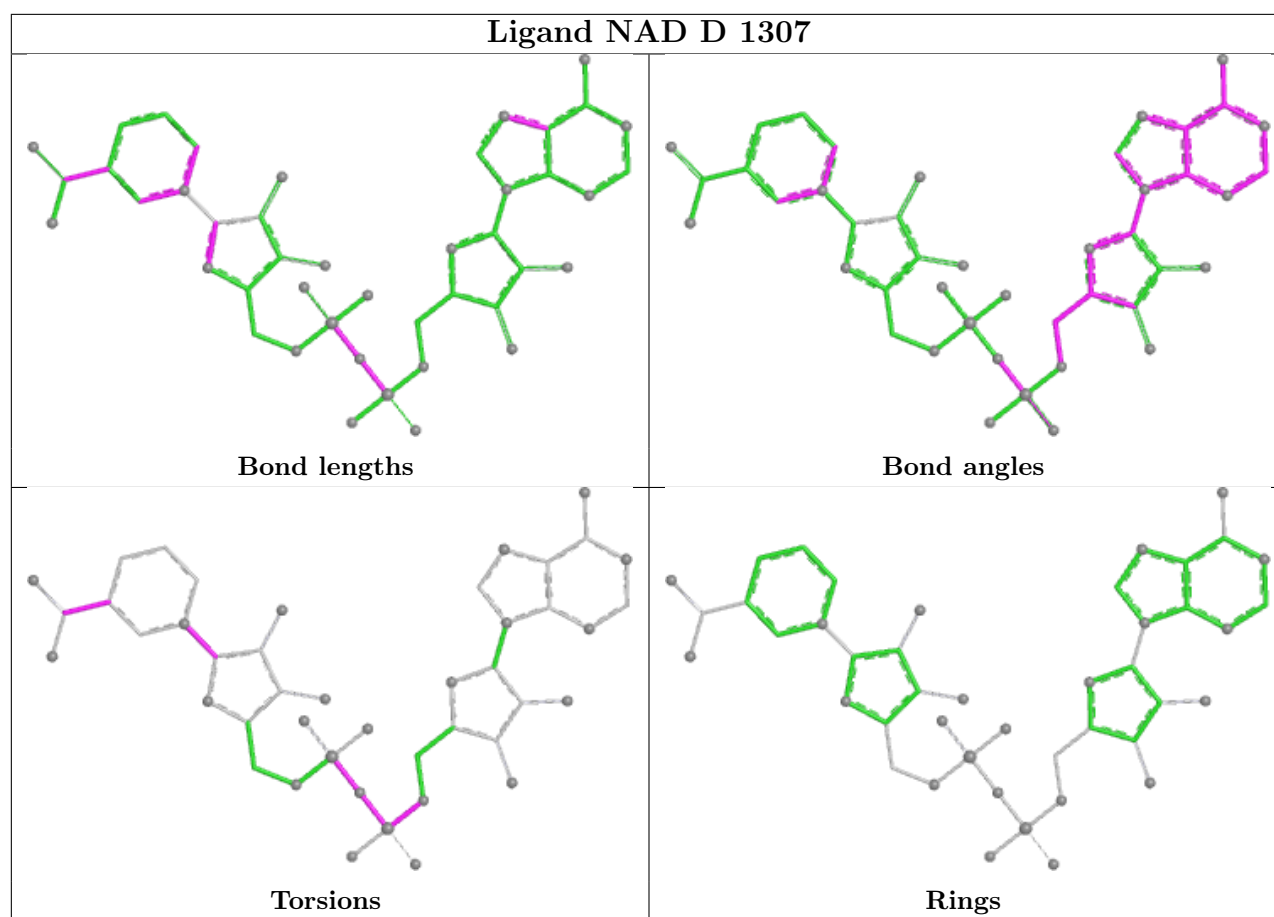
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	1306	THT	17	0
2	F	1309	NAD	1	0
2	A	1301	NAD	2	0
3	A	1302	THT	17	0
3	B	1304	THT	16	0
2	D	1307	NAD	2	0
2	E	1308	NAD	9	0
2	C	1305	NAD	1	0
2	B	1303	NAD	1	0
3	F	1310	THT	17	0

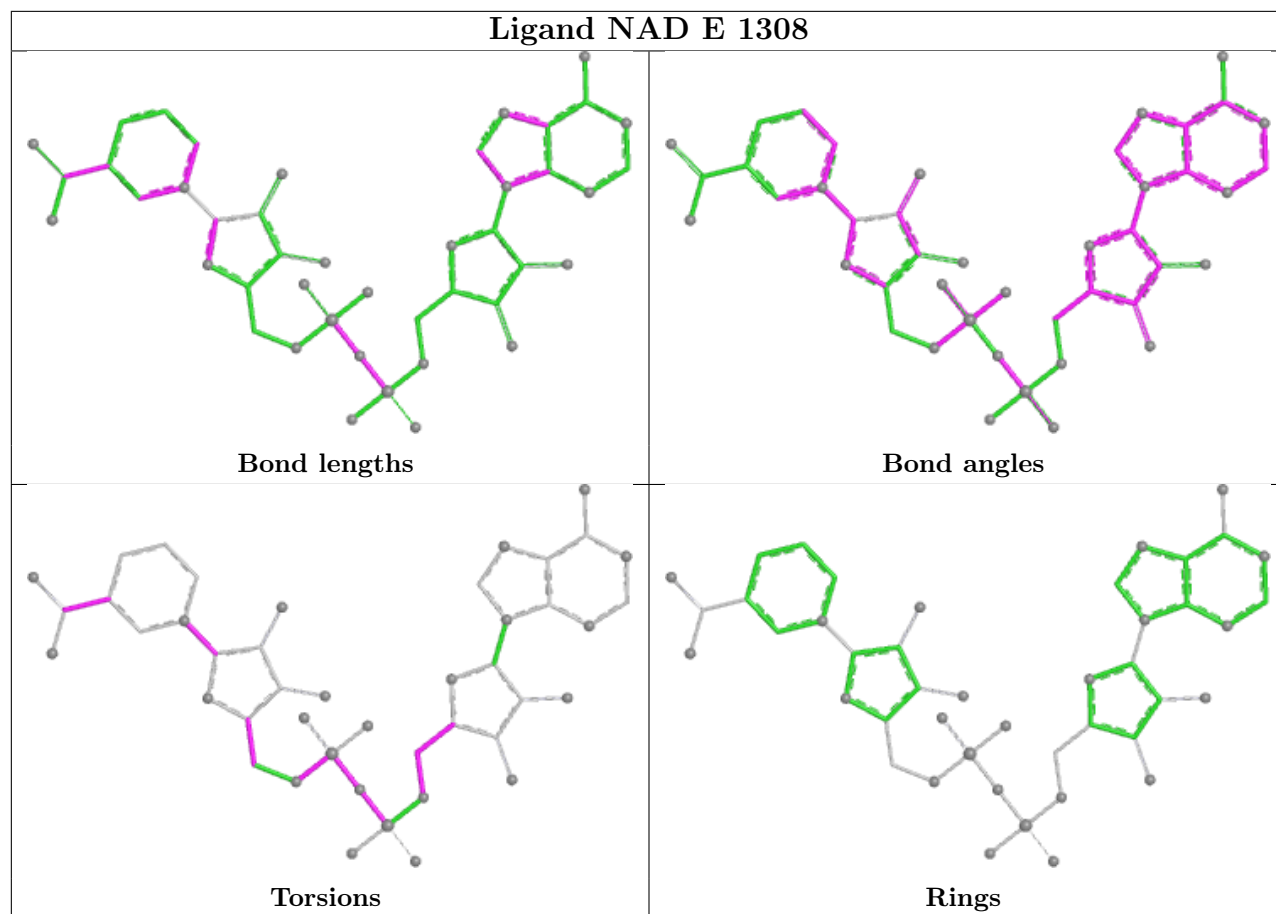
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

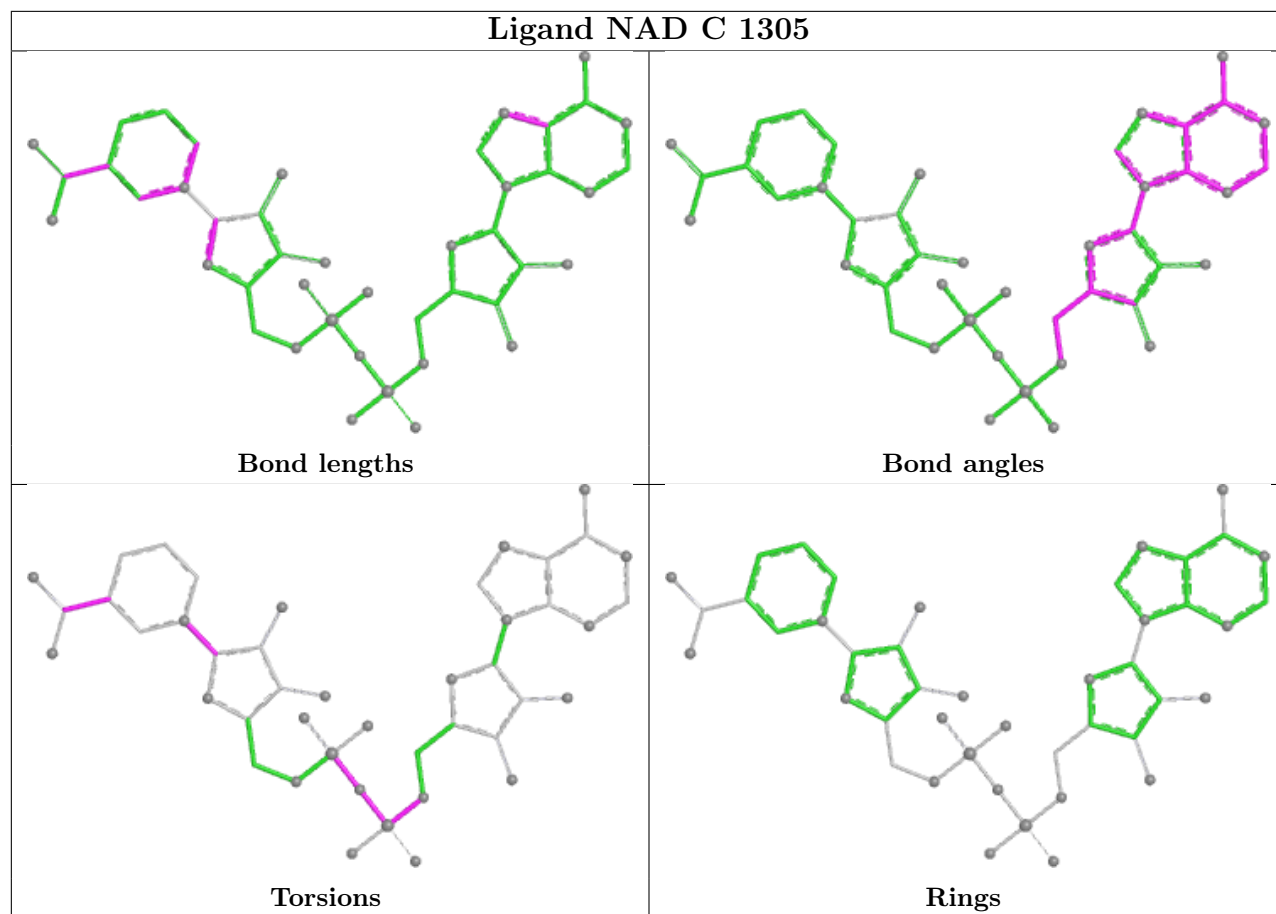


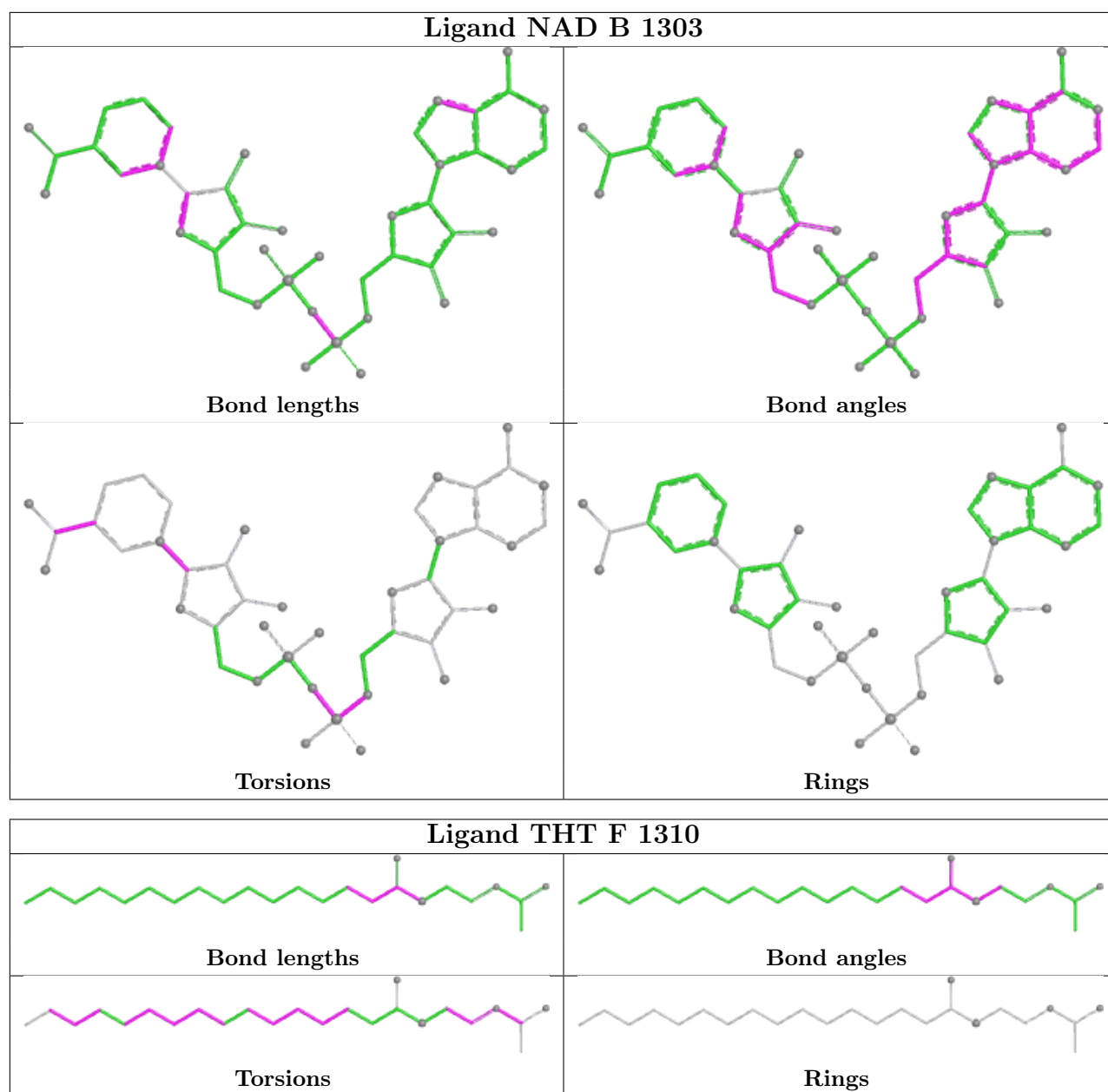












5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

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

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

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	268/268 (100%)	-0.31	2 (0%) 84 77	18, 43, 80, 104	0
1	B	268/268 (100%)	-0.30	1 (0%) 88 84	12, 43, 76, 104	0
1	C	268/268 (100%)	-0.19	1 (0%) 88 84	25, 47, 79, 109	0
1	D	268/268 (100%)	-0.21	1 (0%) 88 84	18, 50, 78, 102	0
1	E	268/268 (100%)	-0.58	0 100 100	17, 44, 62, 84	0
1	F	268/268 (100%)	-0.21	1 (0%) 88 84	19, 47, 78, 110	0
All	All	1608/1608 (100%)	-0.30	6 (0%) 88 84	12, 46, 76, 110	0

The worst 5 of 6 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	198	ALA	2.6
1	D	205	GLY	2.4
1	B	198	ALA	2.4
1	F	47	ILE	2.3
1	A	221	GLY	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 oligosaccharides in this entry.

6.4 Ligands

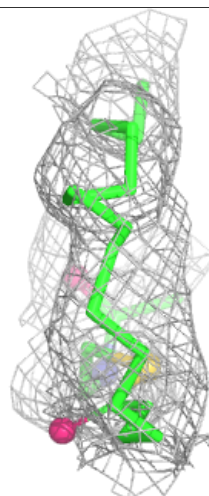
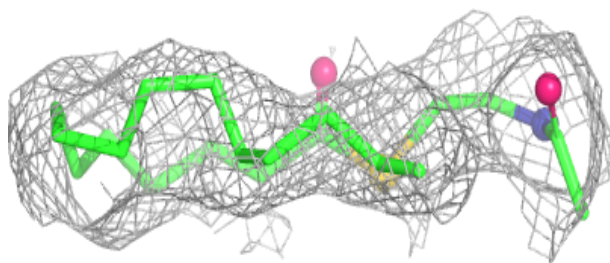
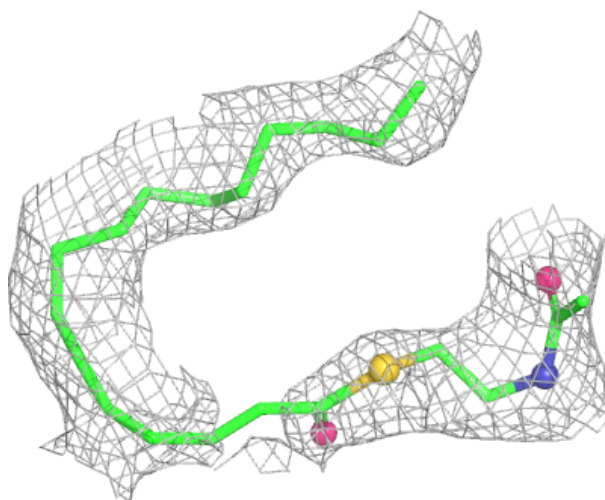
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	THT	A	1302	24/24	0.72	0.09	48,78,85,88	0
3	THT	F	1310	24/24	0.74	0.11	49,76,100,101	0
3	THT	B	1304	24/24	0.76	0.10	61,74,89,91	0
3	THT	C	1306	24/24	0.80	0.11	45,76,103,107	0
2	NAD	A	1301	44/44	0.86	0.10	55,67,75,77	0
2	NAD	D	1307	44/44	0.86	0.10	49,74,83,86	0
2	NAD	C	1305	44/44	0.87	0.09	57,66,76,80	0
2	NAD	F	1309	44/44	0.87	0.09	57,69,75,78	0
2	NAD	B	1303	44/44	0.89	0.10	47,65,74,78	0
2	NAD	E	1308	44/44	0.92	0.07	34,54,70,72	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.

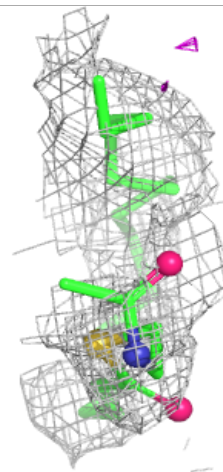
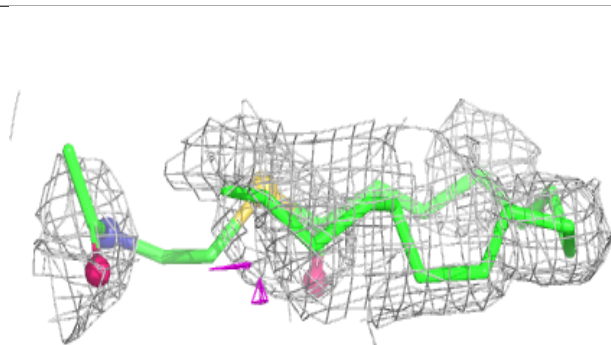
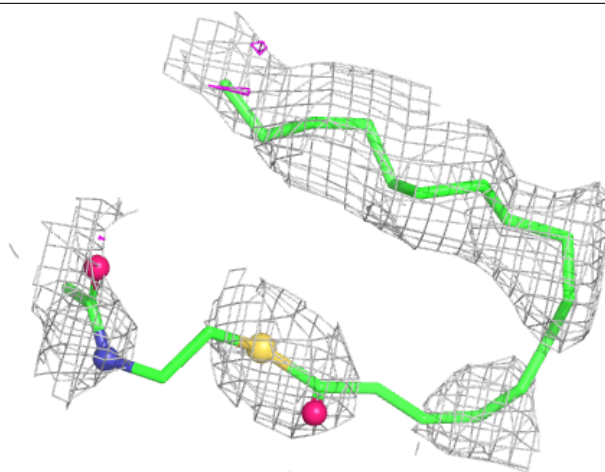
Electron density around THT A 1302:

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



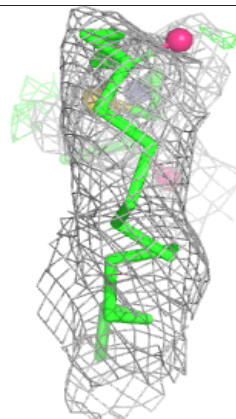
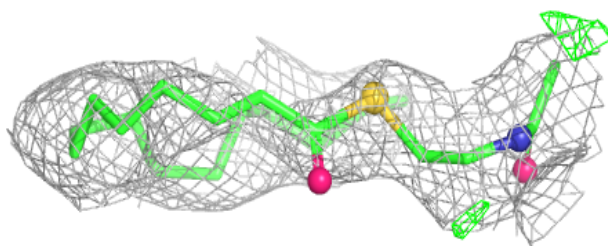
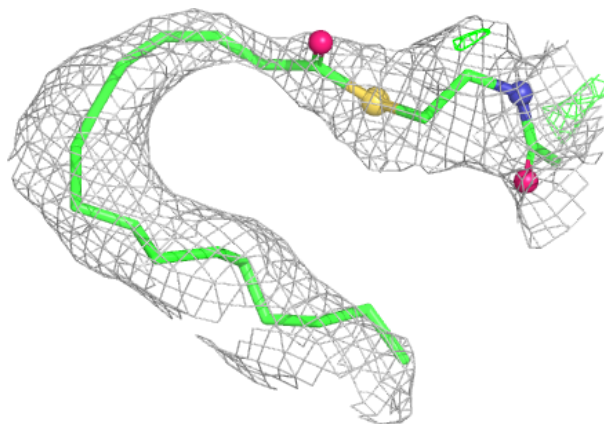
Electron density around THT F 1310:

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



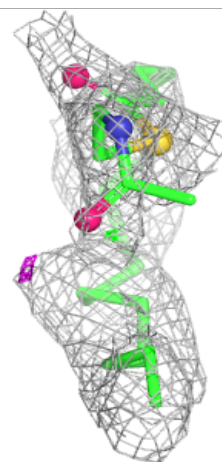
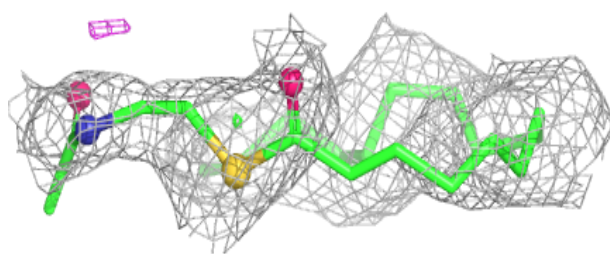
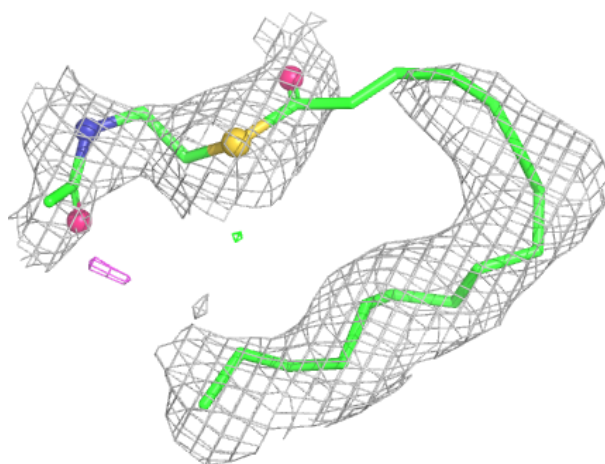
Electron density around THT B 1304:

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



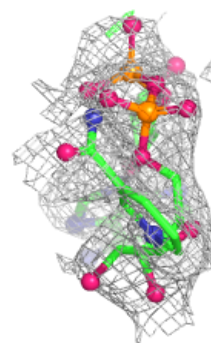
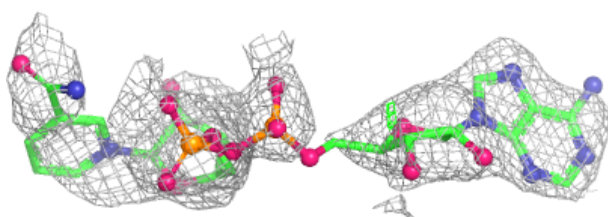
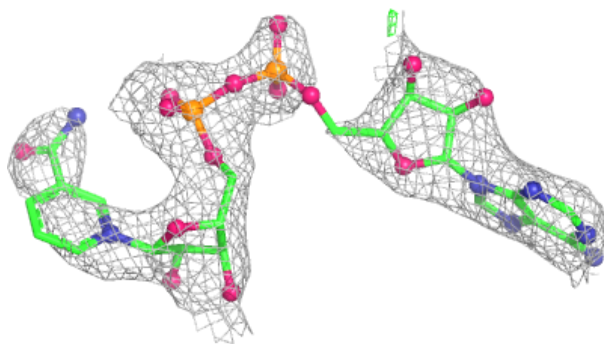
Electron density around THT C 1306:

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

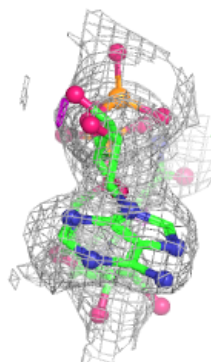
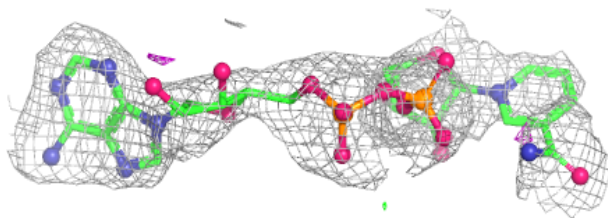
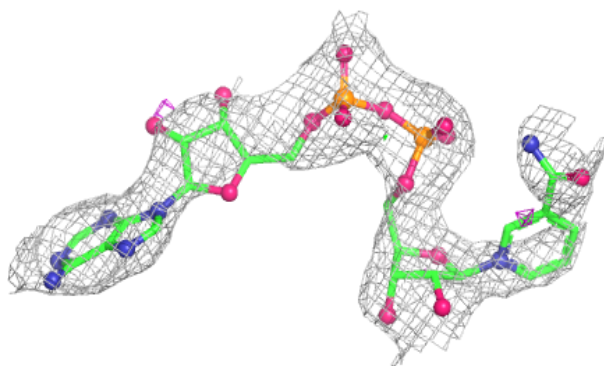


Electron density around NAD A 1301:

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

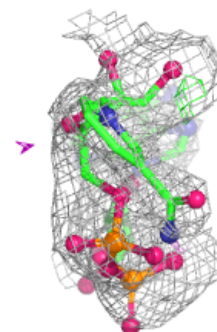
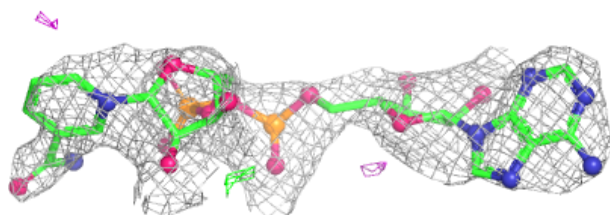
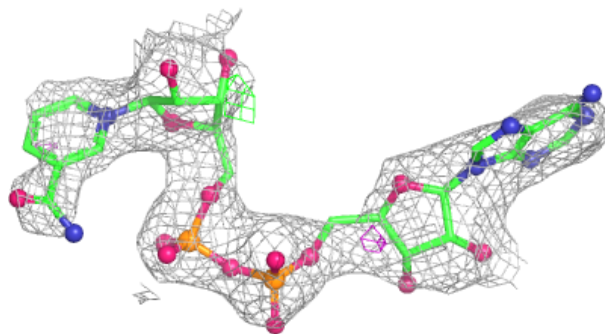
**Electron density around NAD D 1307:**

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

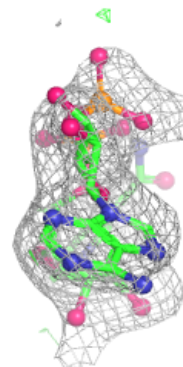
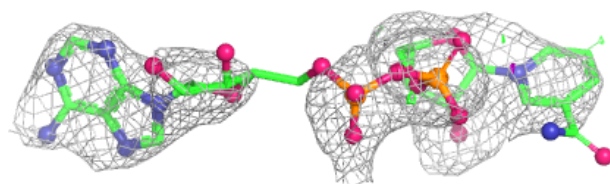
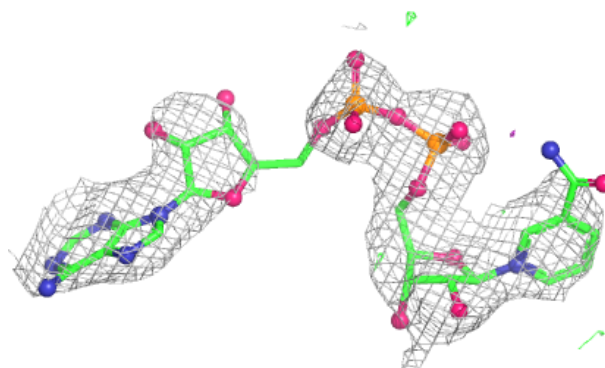


Electron density around NAD C 1305:

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

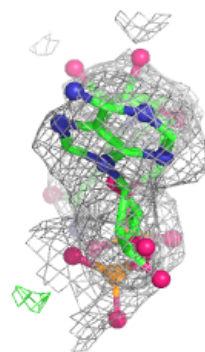
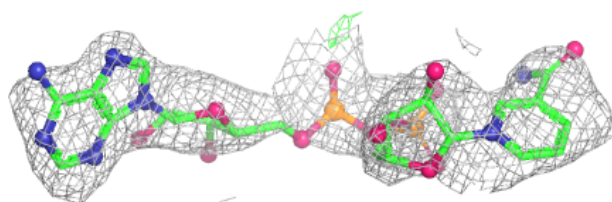
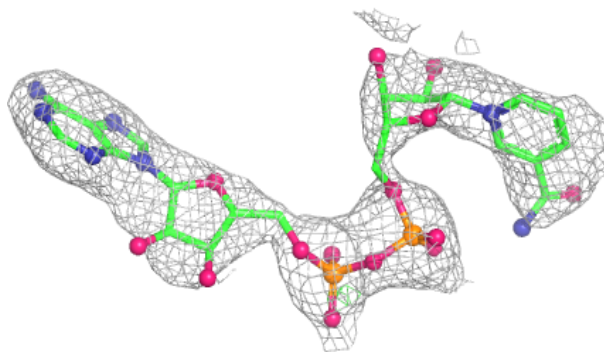
**Electron density around NAD F 1309:**

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

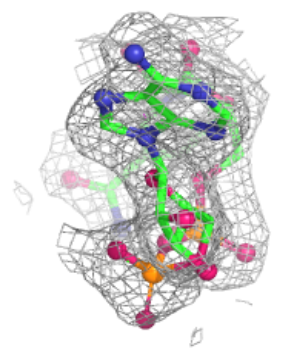
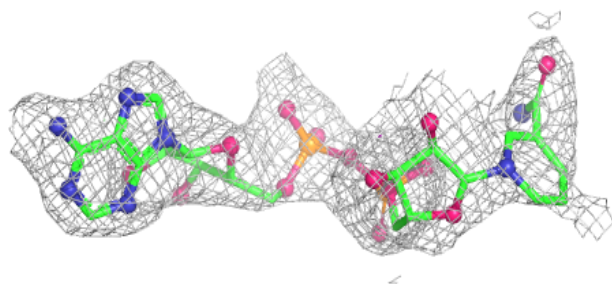
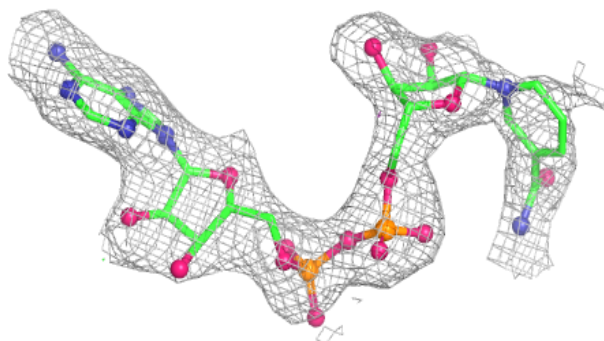


Electron density around NAD B 1303:

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

**Electron density around NAD E 1308:**

$2mF_o-DF_c$ (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.