



# Full wwPDB X-ray Structure Validation Report ⓘ

Sep 7, 2020 – 02:53 PM BST

PDB ID : 6VZ9  
Title : Structure of proline utilization A with the FAD covalently modified by L-thiazolidine-2-carboxylate  
Authors : Campbell, A.C.; Tanner, J.J.  
Deposited on : 2020-02-28  
Resolution : 1.52 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](mailto: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.

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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 : **FAILED**  
buster-report : 1.1.7 (2018)  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.14.2

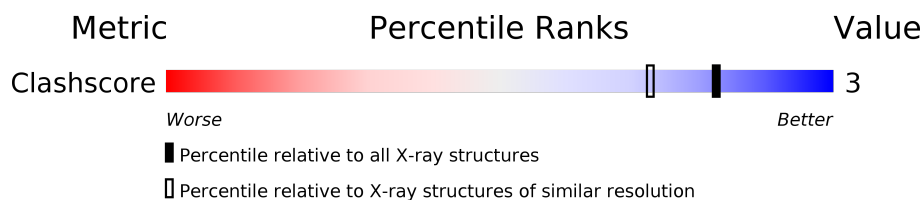
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.52 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))
Clashscore	141614	4249 (1.54-1.50)

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  $\geq 3$ , 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  $\leq 5\%$

Note EDS failed to run properly.

Mol	Chain	Length	Quality of chain
1	A	1235	 92% 7% •
1	B	1235	 93% 6% •

## 2 Entry composition [i](#)

There are 9 unique types of molecules in this entry. The entry contains 20839 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.

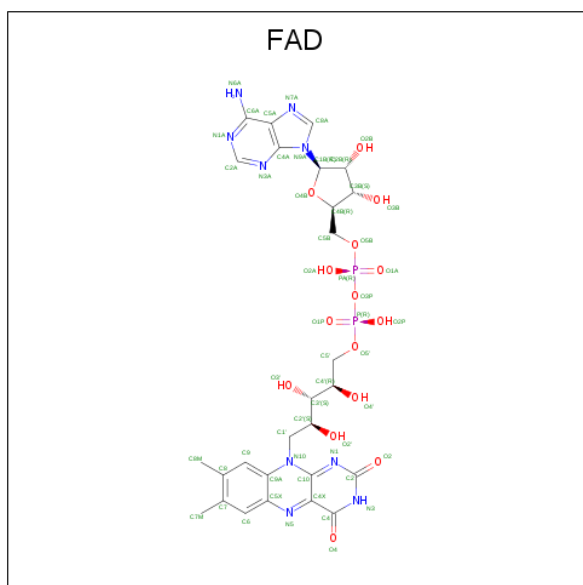
- Molecule 1 is a protein called Bifunctional protein PutA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	1217	Total	C	N	O	S	0	10	0
			9108	5727	1635	1713	33			
1	B	1217	Total	C	N	O	S	0	10	0
			9096	5721	1642	1700	33			

There are 4 discrepancies between the modelled and reference sequences:

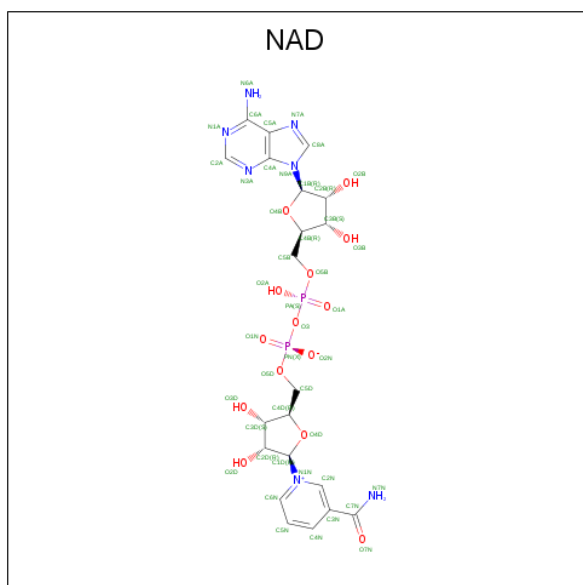
Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	SER	-	expression tag	UNP F7X6I3
A	0	MET	-	expression tag	UNP F7X6I3
B	-1	SER	-	expression tag	UNP F7X6I3
B	0	MET	-	expression tag	UNP F7X6I3

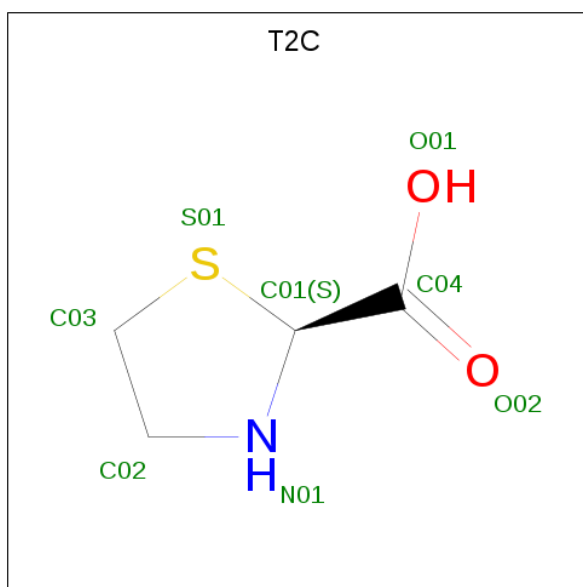
- Molecule 2 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula:  $C_{27}H_{33}N_9O_{15}P_2$ ) (labeled as "Ligand of Interest" by author).



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		

- Molecule 3 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (three-letter code: NAD) (formula: C<sub>21</sub>H<sub>27</sub>N<sub>7</sub>O<sub>14</sub>P<sub>2</sub>) (labeled as "Ligand of Interest" by author).





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

- Molecule 6 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula:  $C_4H_{10}O_3$ ).



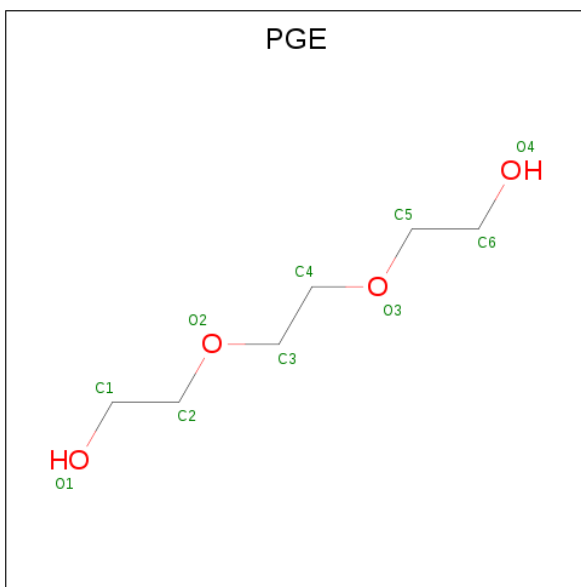
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			7	4	3		

- Molecule 7 is SULFATE ION (three-letter code: SO4) (formula:  $O_4S$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total	O	S	0	0
			5	4	1		
7	A	1	Total	O	S	0	0
			5	4	1		
7	A	1	Total	O	S	0	0
			5	4	1		
7	A	1	Total	O	S	0	0
			5	4	1		
7	B	1	Total	O	S	0	0
			5	4	1		
7	B	1	Total	O	S	0	0
			5	4	1		
7	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 8 is TRIETHYLENE GLYCOL (three-letter code: PGE) (formula: C<sub>6</sub>H<sub>14</sub>O<sub>4</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	B	1	Total	C	O	0	0
			10	6	4		

- Molecule 9 is water.

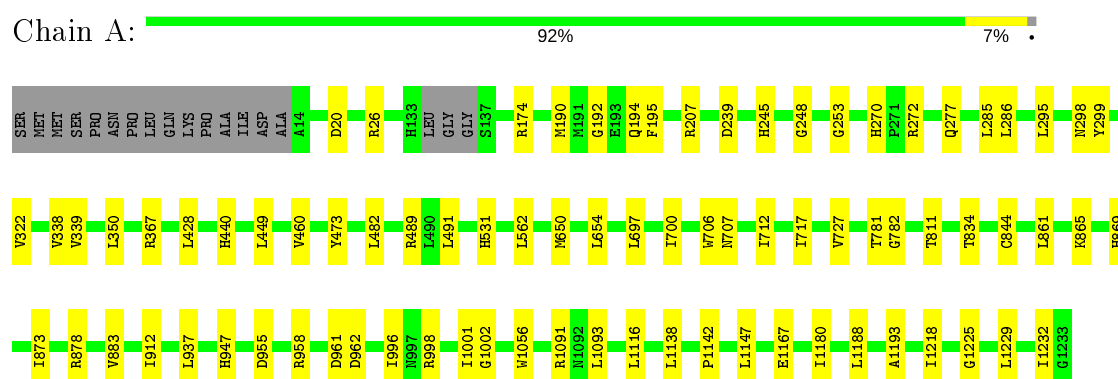
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	1129	Total	O	0	0
			1129	1129		
9	B	1237	Total	O	0	0
			1237	1237		

### 3 Residue-property plots

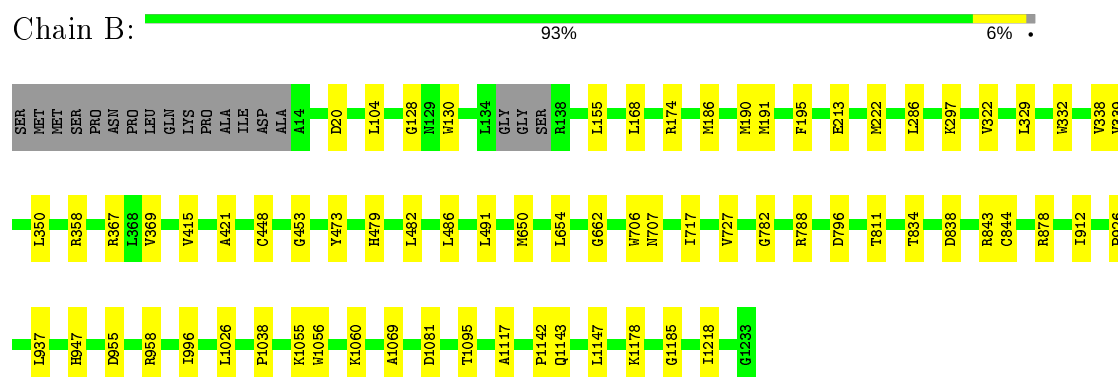
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded 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. 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.

Note EDS failed to run properly.

- Molecule 1: Bifunctional protein PutA



- Molecule 1: Bifunctional protein PutA





## 4 Data and refinement statistics

EDS failed to run properly - this section is therefore incomplete.

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	100.96Å 102.10Å 126.38Å 90.00° 106.35° 90.00°	Depositor
Resolution (Å)	48.44 – 1.52	Depositor
% Data completeness (in resolution range)	95.0 (48.44-1.52)	Depositor
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.36 (at 1.52Å)	Xtriage
Refinement program	PHENIX 1.14	Depositor
R, $R_{free}$	0.180 , 0.207	Depositor
Wilson B-factor (Å <sup>2</sup> )	17.2	Xtriage
Anisotropy	0.166	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	20839	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.40% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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, PGE, NAD, T2C, SO4, PEG, 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	0.33	0/9297	0.54	1/12644 (0.0%)
1	B	0.34	0/9285	0.56	1/12627 (0.0%)
All	All	0.33	0/18582	0.55	2/25271 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	174	ARG	NE-CZ-NH1	6.24	123.42	120.30
1	B	174	ARG	NE-CZ-NH1	5.12	122.86	120.30

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	9108	0	9180	49	0
1	B	9096	0	9177	45	0
2	A	53	0	31	1	0
2	B	53	0	31	1	0
3	A	44	0	26	1	0
3	B	44	0	26	2	0
4	A	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	1	0	0	0	0
5	A	8	0	0	0	0
5	B	8	0	0	0	0
6	A	7	0	10	0	0
7	A	20	0	0	0	0
7	B	20	0	0	0	0
8	B	10	0	14	1	0
9	A	1129	0	0	5	0
9	B	1237	0	0	4	0
All	All	20839	0	18495	93	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.

All (93) 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:A:491:LEU:HD13	1:A:1225:GLY:HA3	1.72	0.69
1:B:195:PHE:HB3	1:B:482:LEU:HD11	1.75	0.69
1:A:873:ILE:HG13	1:A:883:VAL:HB	1.75	0.68
1:B:844:CYS:SG	3:B:1302:NAD:C4N	2.81	0.68
1:A:958:ARG:NH1	1:A:962:ASP:OD2	2.30	0.65
1:B:339:VAL:HG21	1:B:350:LEU:HD21	1.79	0.64
1:B:358:ARG:HG2	1:B:415:VAL:HG11	1.81	0.62
1:A:195:PHE:HB3	1:A:482:LEU:HD11	1.80	0.62
1:A:286:LEU:HD21	1:A:322:VAL:HG11	1.81	0.62
1:A:996[B]:ILE:HD12	1:A:1218:ILE:HG12	1.84	0.60
1:A:1147:LEU:HD22	1:B:1147:LEU:HD13	1.83	0.60
1:B:955:ASP:HB3	1:B:958:ARG:HH21	1.67	0.60
1:B:650:MET:O	1:B:654:LEU:HG	2.02	0.59
1:B:297:LYS:HD2	1:B:329:LEU:HA	1.83	0.59
1:B:20:ASP:OD2	1:B:878:ARG:NH1	2.35	0.58
1:B:834:THR:O	1:B:838[B]:ASP:HB2	2.03	0.58
1:A:1116[B]:LEU:HD11	1:A:1138:LEU:HD11	1.84	0.58
1:B:297:LYS:HG3	1:B:332:TRP:HB2	1.84	0.58
1:A:298:ASN:ND2	9:A:1401:HOH:O	2.19	0.57
1:B:286:LEU:HD21	1:B:322:VAL:HG11	1.86	0.57
1:A:869:HIS:ND1	1:A:912:ILE:HG22	2.20	0.56
1:B:155:LEU:HD22	1:B:491:LEU:HD11	1.88	0.55
1:A:998:ARG:HH22	1:A:1002:GLY:HA3	1.72	0.55
1:A:861:LEU:HD22	1:A:865:LYS:HE3	1.88	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:844:CYS:SG	3:A:1302:NAD:C4N	2.95	0.54
1:B:1056:TRP:CD1	1:B:1142:PRO:HD3	2.41	0.54
1:A:473:TYR:HB2	2:A:1301:FAD:HM72	1.90	0.54
1:A:562:LEU:HD11	1:A:654:LEU:HD12	1.90	0.53
1:A:1056:TRP:CD1	1:A:1142:PRO:HD3	2.44	0.53
1:B:473:TYR:HB2	2:B:1301:FAD:HM72	1.90	0.52
1:A:26:ARG:HH22	1:A:322:VAL:HG23	1.74	0.52
1:A:270:HIS:HB2	1:A:285:LEU:HG	1.93	0.51
1:A:1229:LEU:HD23	1:A:1232:ILE:HD12	1.93	0.50
1:A:449:LEU:HD13	1:A:489:ARG:HD2	1.92	0.50
1:B:782:GLY:O	1:B:811:THR:HA	2.11	0.50
1:B:1026:LEU:HD23	1:B:1038:PRO:HG2	1.94	0.50
1:B:796:ASP:HA	1:B:1178:LYS:HE2	1.94	0.49
1:B:1143:GLN:O	1:B:1147:LEU:HG	2.12	0.49
1:A:961:ASP:OD2	1:B:1055:LYS:NZ	2.40	0.48
1:A:192:GLY:O	1:A:207:ARG:NH1	2.46	0.48
1:B:213:GLU:OE1	9:B:1401:HOH:O	2.20	0.47
1:A:190:MET:HA	1:A:194:GLN:HB3	1.97	0.46
1:B:1056:TRP:CZ2	1:B:1060:LYS:HD2	2.51	0.46
8:B:1305:PGE:H1	9:B:1602:HOH:O	2.16	0.45
1:A:428:LEU:HD11	1:A:460:VAL:HG21	1.99	0.45
1:A:245:HIS:CE1	1:A:295:LEU:HD11	2.52	0.45
1:A:697:LEU:HD12	1:A:700[B]:ILE:HD11	1.98	0.45
1:B:104:LEU:HD22	1:B:168:LEU:HD11	1.99	0.45
1:A:440:HIS:HE1	9:A:2359:HOH:O	1.99	0.44
1:B:128:GLY:O	1:B:130:TRP:N	2.46	0.44
1:B:448:CYS:HB2	1:B:453:GLY:HA3	1.99	0.44
1:B:706:TRP:CE3	1:B:707:ASN:HA	2.53	0.44
1:A:650:MET:O	1:A:654:LEU:HG	2.17	0.44
1:A:782:GLY:O	1:A:811:THR:HA	2.18	0.44
1:A:937:LEU:HD21	1:A:947:HIS:CD2	2.52	0.44
1:B:844:CYS:SG	3:B:1302:NAD:C3N	3.05	0.44
1:A:272:ARG:HB3	1:A:277:GLN:HG3	2.00	0.44
1:A:195:PHE:HB2	1:A:482:LEU:HD21	1.99	0.43
1:A:338:VAL:HG22	1:A:367:ARG:HB3	2.00	0.43
1:B:937:LEU:HD21	1:B:947:HIS:CD2	2.52	0.43
1:A:717:ILE:HG12	1:A:727:VAL:HG11	2.00	0.43
1:A:1180:ILE:HG23	1:A:1188:LEU:HD12	1.99	0.43
1:A:239:ASP:OD2	9:A:1402:HOH:O	2.21	0.43
1:A:706:TRP:CE3	1:A:707:ASN:HA	2.53	0.43
1:B:955:ASP:HB3	1:B:958:ARG:NH2	2.33	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:20:ASP:OD2	1:A:878:ARG:NH1	2.52	0.43
1:A:253:GLY:HA2	9:A:2239:HOH:O	2.18	0.43
1:B:834:THR:O	1:B:838[A]:ASP:HB3	2.18	0.43
1:A:834:THR:HG22	1:A:1001:ILE:HD11	2.01	0.43
1:B:369:VAL:HG12	1:B:421:ALA:HB3	2.01	0.43
1:A:1091:ARG:HG2	1:A:1093:LEU:HD21	2.00	0.42
1:B:838[A]:ASP:OD1	1:B:843:ARG:NH2	2.46	0.42
1:B:186:MET:O	1:B:190:MET:HG3	2.19	0.42
1:B:482:LEU:HD11	1:B:486:LEU:HD13	2.00	0.42
1:B:996[A]:ILE:HD12	1:B:1218:ILE:HG12	2.01	0.42
1:B:191:MET:HB3	1:B:479:HIS:CD2	2.55	0.41
1:B:662:GLY:HA2	9:B:1511:HOH:O	2.20	0.41
1:B:1069:ALA:HA	1:B:1117:ALA:HB1	2.02	0.41
1:A:955:ASP:HB3	1:A:958:ARG:HE	1.85	0.41
1:B:788:ARG:NH1	1:B:1185:GLY:O	2.54	0.41
1:B:195:PHE:HB2	1:B:482:LEU:HD21	2.02	0.41
1:A:1002:GLY:HA3	9:A:1531:HOH:O	2.20	0.41
1:B:1081:ASP:HA	1:B:1095[A]:THR:HG22	2.03	0.41
1:A:1167:GLU:HA	1:A:1193:ALA:O	2.21	0.41
1:A:248:GLY:HA3	1:A:299:TYR:CG	2.55	0.41
1:A:339[B]:VAL:HG21	1:A:350:LEU:HD21	2.03	0.41
1:A:958:ARG:O	1:A:961:ASP:HB2	2.21	0.41
1:A:531:HIS:CE1	1:A:1232:ILE:HG23	2.56	0.40
1:B:338:VAL:HG22	1:B:367:ARG:HB3	2.03	0.40
1:A:712:ILE:HD13	1:A:781:THR:HG21	2.03	0.40
1:B:222:MET:HG2	9:B:1612:HOH:O	2.22	0.40
1:B:717:ILE:HG12	1:B:727:VAL:HG11	2.04	0.40
1:B:912:ILE:HD11	1:B:926:PRO:HD2	2.03	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles [i](#)

### 5.3.1 Protein backbone [i](#)

There are no protein backbone outliers to report in this entry.

### 5.3.2 Protein sidechains [i](#)

There are no protein residues with a non-rotameric sidechain to report in this entry.

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates ⓘ

There are no monosaccharides in this entry.

## 5.6 Ligand geometry ⓘ

Of 18 ligands modelled in this entry, 2 are monoatomic - leaving 16 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	$\# Z  > 2$
7	SO4	B	1308	-	4,4,4	0.13	0	6,6,6	0.09	0
7	SO4	A	1309	-	4,4,4	0.17	0	6,6,6	0.20	0
7	SO4	B	1307	-	4,4,4	0.14	0	6,6,6	0.10	0
7	SO4	B	1309	-	4,4,4	0.14	0	6,6,6	0.07	0
3	NAD	A	1302	4	42,48,48	3.57	17 (40%)	50,73,73	1.75	6 (12%)
7	SO4	A	1306	-	4,4,4	0.13	0	6,6,6	0.17	0
7	SO4	B	1306	-	4,4,4	0.25	0	6,6,6	0.33	0
3	NAD	B	1302	4	42,48,48	3.47	16 (38%)	50,73,73	1.64	6 (12%)
6	PEG	A	1305	-	6,6,6	0.51	0	5,5,5	0.25	0
7	SO4	A	1308	-	4,4,4	0.15	0	6,6,6	0.21	0
2	FAD	B	1301	5	51,58,58	2.82	17 (33%)	60,89,89	1.96	13 (21%)
7	SO4	A	1307	-	4,4,4	0.16	0	6,6,6	0.16	0
5	T2C	B	1304	2	4,8,8	0.70	0	3,10,10	2.35	2 (66%)
5	T2C	A	1304	2	4,8,8	0.66	0	3,10,10	2.41	2 (66%)
2	FAD	A	1301	5	51,58,58	2.74	18 (35%)	60,89,89	1.91	13 (21%)
8	PGE	B	1305	-	9,9,9	0.51	0	8,8,8	0.30	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
3	NAD	A	1302	4	-	4/26/62/62	0/5/5/5
3	NAD	B	1302	4	-	2/26/62/62	0/5/5/5
6	PEG	A	1305	-	-	2/4/4/4	-
8	PGE	B	1305	-	-	3/7/7/7	-
2	FAD	B	1301	5	-	2/30/50/50	0/6/6/6
5	T2C	B	1304	2	-	0/0/11/11	0/1/1/1
5	T2C	A	1304	2	-	0/0/11/11	0/1/1/1
2	FAD	A	1301	5	-	3/30/50/50	0/6/6/6

All (68) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1302	NAD	O4D-C1D	-9.02	1.28	1.41
3	B	1302	NAD	O4D-C1D	-8.90	1.28	1.41
2	B	1301	FAD	C10-N1	8.69	1.44	1.33
3	B	1302	NAD	C3B-C4B	-8.33	1.31	1.53
3	A	1302	NAD	C3B-C4B	-8.30	1.31	1.53
3	A	1302	NAD	C7N-N7N	8.24	1.48	1.33
2	B	1301	FAD	C4X-N5	8.18	1.45	1.33
2	A	1301	FAD	C4X-N5	8.12	1.44	1.33
2	A	1301	FAD	C10-N1	8.03	1.43	1.33
3	B	1302	NAD	C7N-N7N	8.03	1.48	1.33
3	A	1302	NAD	C3D-C4D	-7.66	1.33	1.53
3	B	1302	NAD	C3D-C4D	-7.25	1.34	1.53
3	A	1302	NAD	O4B-C4B	7.16	1.61	1.45
3	A	1302	NAD	O4D-C4D	7.15	1.61	1.45
3	B	1302	NAD	O4B-C4B	6.99	1.60	1.45
3	B	1302	NAD	O4D-C4D	6.88	1.60	1.45
2	A	1301	FAD	C5X-N5	6.61	1.46	1.35
2	B	1301	FAD	C5X-N5	6.16	1.45	1.35
3	A	1302	NAD	O4B-C1B	-5.33	1.33	1.41
2	B	1301	FAD	C4-C4X	4.97	1.49	1.41
2	B	1301	FAD	O4-C4	4.87	1.36	1.24
2	A	1301	FAD	C4-C4X	4.83	1.49	1.41
2	B	1301	FAD	C9A-N10	4.76	1.44	1.38
2	B	1301	FAD	C6-C5X	4.58	1.49	1.41
3	B	1302	NAD	O4B-C1B	-4.53	1.34	1.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1301	FAD	C6-C5X	4.50	1.48	1.41
2	A	1301	FAD	O4-C4	4.40	1.35	1.24
3	B	1302	NAD	C3N-C7N	4.39	1.57	1.50
3	B	1302	NAD	C6A-N6A	4.34	1.49	1.34
3	A	1302	NAD	C3N-C7N	4.19	1.56	1.50
3	A	1302	NAD	O3D-C3D	4.11	1.52	1.43
3	A	1302	NAD	C6A-N6A	4.02	1.48	1.34
3	B	1302	NAD	O3D-C3D	3.92	1.52	1.43
2	A	1301	FAD	C9A-N10	3.72	1.43	1.38
2	B	1301	FAD	C2B-C1B	-3.65	1.48	1.53
2	A	1301	FAD	C9-C9A	3.64	1.48	1.40
2	B	1301	FAD	C4-N3	3.61	1.39	1.33
2	A	1301	FAD	C4-N3	3.54	1.39	1.33
2	B	1301	FAD	C9-C9A	3.40	1.47	1.40
2	B	1301	FAD	C9A-C5X	-3.40	1.35	1.42
2	A	1301	FAD	C2B-C1B	-3.34	1.48	1.53
3	A	1302	NAD	O3B-C3B	3.28	1.50	1.43
3	B	1302	NAD	C2N-N1N	3.24	1.38	1.35
3	A	1302	NAD	C2N-N1N	3.14	1.38	1.35
2	A	1301	FAD	C9A-C5X	-3.05	1.36	1.42
2	B	1301	FAD	C2A-N3A	3.05	1.37	1.32
3	B	1302	NAD	O3B-C3B	2.91	1.49	1.43
2	B	1301	FAD	C2-N1	2.84	1.43	1.38
3	B	1302	NAD	C2A-N1A	2.83	1.39	1.33
2	B	1301	FAD	C6A-N6A	2.81	1.44	1.34
2	A	1301	FAD	C6A-N6A	2.74	1.44	1.34
3	A	1302	NAD	C2A-N1A	2.69	1.38	1.33
2	A	1301	FAD	PA-O5B	-2.66	1.48	1.59
2	A	1301	FAD	C2A-N3A	2.61	1.36	1.32
2	A	1301	FAD	O2'-C2'	-2.58	1.37	1.43
3	A	1302	NAD	C5A-C4A	-2.58	1.34	1.40
2	A	1301	FAD	C2-N1	2.52	1.43	1.38
2	B	1301	FAD	PA-O2A	-2.42	1.43	1.55
2	B	1301	FAD	O2'-C2'	-2.38	1.38	1.43
3	A	1302	NAD	O7N-C7N	-2.35	1.19	1.24
3	A	1302	NAD	O2B-C2B	-2.29	1.37	1.43
2	A	1301	FAD	O3B-C3B	-2.21	1.37	1.43
3	A	1302	NAD	C2A-N3A	2.20	1.35	1.32
2	B	1301	FAD	O3B-C3B	-2.18	1.37	1.43
3	B	1302	NAD	C5A-C4A	-2.18	1.35	1.40
3	B	1302	NAD	O7N-C7N	-2.16	1.20	1.24
3	B	1302	NAD	C2A-N3A	2.13	1.35	1.32

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1301	FAD	C2B-C3B	-2.06	1.47	1.53

All (42) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1302	NAD	C1B-N9A-C4A	-8.37	111.93	126.64
2	B	1301	FAD	C1'-N10-C9A	6.75	123.61	118.29
2	A	1301	FAD	C1'-N10-C9A	6.50	123.41	118.29
3	B	1302	NAD	C1B-N9A-C4A	-6.28	115.61	126.64
2	B	1301	FAD	C4-N3-C2	6.01	120.22	115.14
3	A	1302	NAD	N3A-C2A-N1A	-5.75	119.69	128.68
2	A	1301	FAD	N3A-C2A-N1A	-5.73	119.72	128.68
3	B	1302	NAD	N3A-C2A-N1A	-5.69	119.78	128.68
2	A	1301	FAD	C4-N3-C2	5.30	119.62	115.14
2	B	1301	FAD	N3A-C2A-N1A	-5.21	120.54	128.68
2	B	1301	FAD	C9A-N10-C10	-4.60	115.88	121.91
2	A	1301	FAD	C9A-N10-C10	-4.02	116.65	121.91
2	B	1301	FAD	C10-C4X-N5	-3.50	118.84	121.26
2	A	1301	FAD	C5X-C9A-N10	3.33	120.13	117.72
2	B	1301	FAD	C5X-C9A-N10	3.20	120.03	117.72
2	B	1301	FAD	C4X-C4-N3	-3.15	119.12	123.43
2	A	1301	FAD	C4-C4X-N5	3.01	122.03	118.60
3	B	1302	NAD	O4D-C1D-C2D	-2.99	102.55	106.93
3	A	1302	NAD	C3N-C2N-N1N	-2.96	117.54	120.43
3	B	1302	NAD	C3N-C2N-N1N	-2.94	117.55	120.43
5	A	1304	T2C	C02-N01-C01	2.94	112.21	106.19
5	B	1304	T2C	C02-N01-C01	2.88	112.10	106.19
5	B	1304	T2C	C03-C02-N01	2.81	111.73	107.23
2	A	1301	FAD	O5'-P-O1P	2.75	119.80	109.07
2	B	1301	FAD	C1B-N9A-C4A	-2.71	121.88	126.64
2	A	1301	FAD	C1B-N9A-C4A	-2.68	121.93	126.64
2	A	1301	FAD	C10-C4X-N5	-2.64	119.43	121.26
3	B	1302	NAD	C3N-C7N-N7N	2.63	120.91	117.75
2	B	1301	FAD	C4X-N5-C5X	-2.56	114.22	116.77
2	B	1301	FAD	C4-C4X-N5	2.49	121.44	118.60
5	A	1304	T2C	C03-C02-N01	2.45	111.16	107.23
2	A	1301	FAD	C2A-N1A-C6A	2.45	122.94	118.75
3	A	1302	NAD	O4B-C1B-C2B	-2.44	103.36	106.93
3	A	1302	NAD	O4D-C1D-C2D	-2.44	103.36	106.93
2	A	1301	FAD	C4-C4X-C10	-2.41	118.36	119.95
2	B	1301	FAD	C7-C6-C5X	-2.26	118.03	121.22
2	A	1301	FAD	C4X-C4-N3	-2.24	120.37	123.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	1302	NAD	O4B-C1B-C2B	-2.18	103.73	106.93
2	A	1301	FAD	C4X-N5-C5X	-2.13	114.64	116.77
3	A	1302	NAD	C2N-C3N-C4N	2.12	120.66	118.26
2	B	1301	FAD	C1'-N10-C10	2.12	120.30	118.41
2	B	1301	FAD	O5'-P-O1P	2.10	117.26	109.07

There are no chirality outliers.

All (16) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	1302	NAD	C5B-O5B-PA-O1A
8	B	1305	PGE	O3-C5-C6-O4
2	B	1301	FAD	PA-O3P-P-O1P
8	B	1305	PGE	C3-C4-O3-C5
3	A	1302	NAD	C5B-O5B-PA-O3
2	A	1301	FAD	PA-O3P-P-O1P
3	A	1302	NAD	C4D-C5D-O5D-PN
3	B	1302	NAD	C4D-C5D-O5D-PN
2	A	1301	FAD	C2'-C3'-C4'-O4'
2	A	1301	FAD	PA-O3P-P-O2P
6	A	1305	PEG	O1-C1-C2-O2
8	B	1305	PGE	O1-C1-C2-O2
6	A	1305	PEG	C4-C3-O2-C2
3	A	1302	NAD	C3D-C4D-C5D-O5D
2	B	1301	FAD	PA-O3P-P-O2P
3	B	1302	NAD	C5B-O5B-PA-O1A

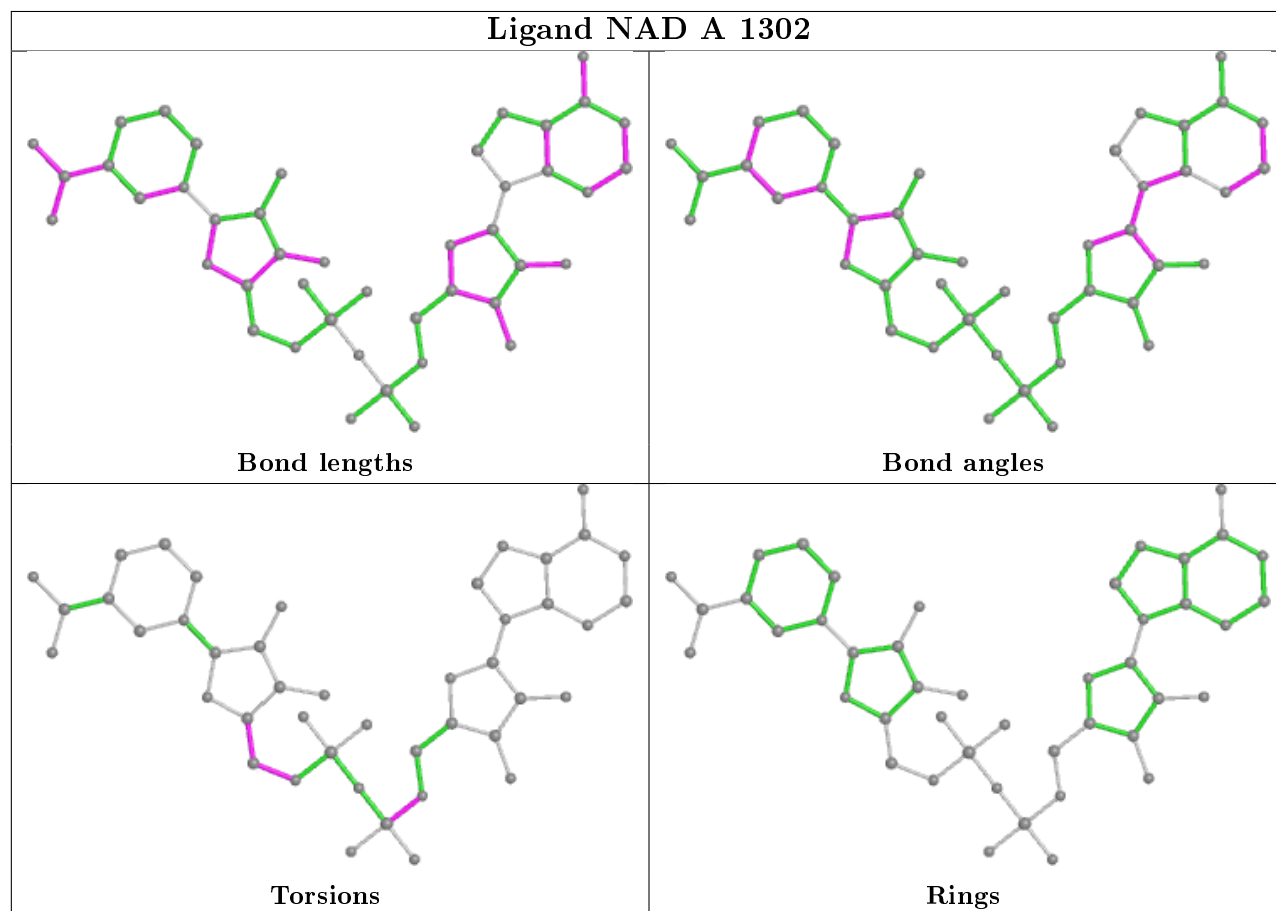
There are no ring outliers.

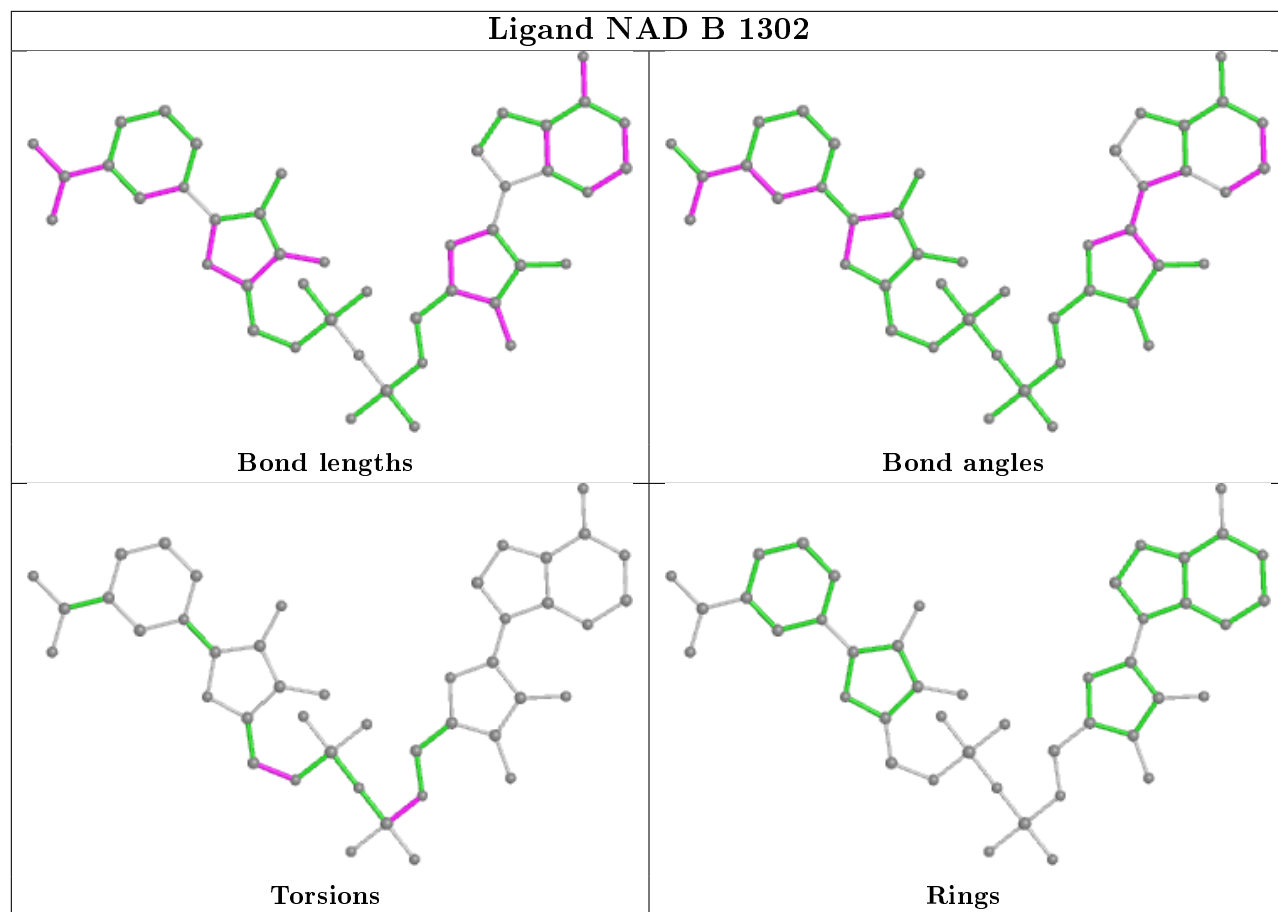
5 monomers are involved in 6 short contacts:

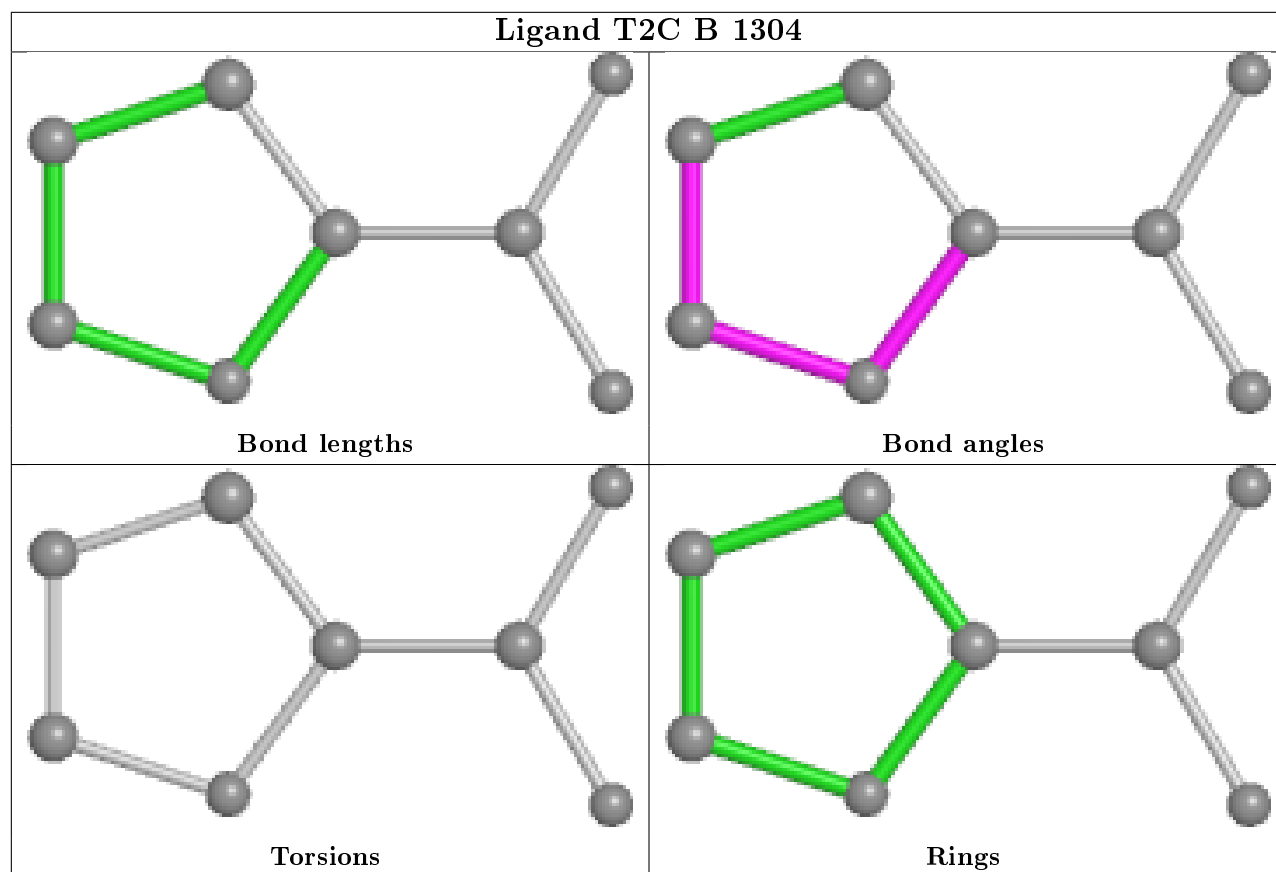
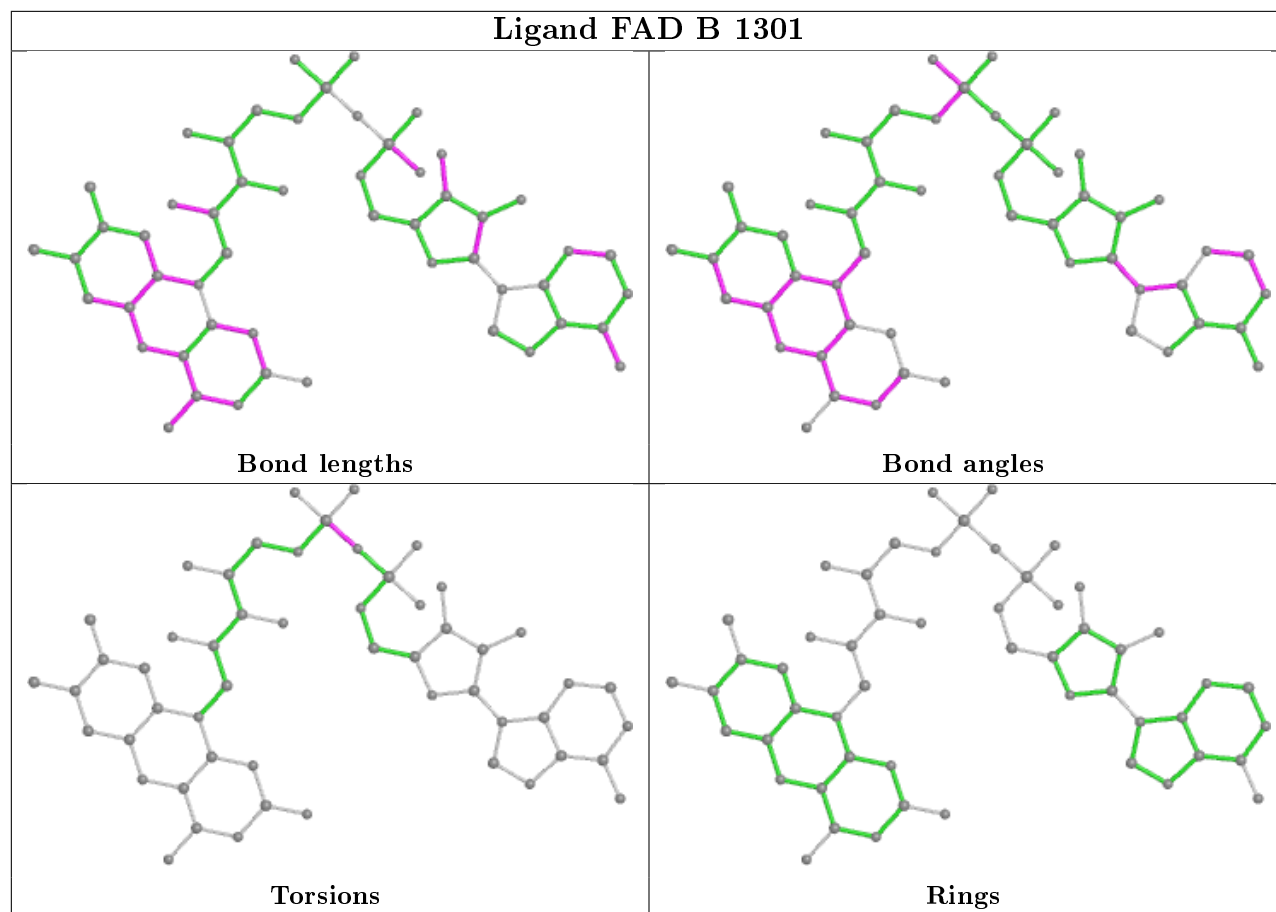
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1302	NAD	1	0
3	B	1302	NAD	2	0
2	B	1301	FAD	1	0
2	A	1301	FAD	1	0
8	B	1305	PGE	1	0

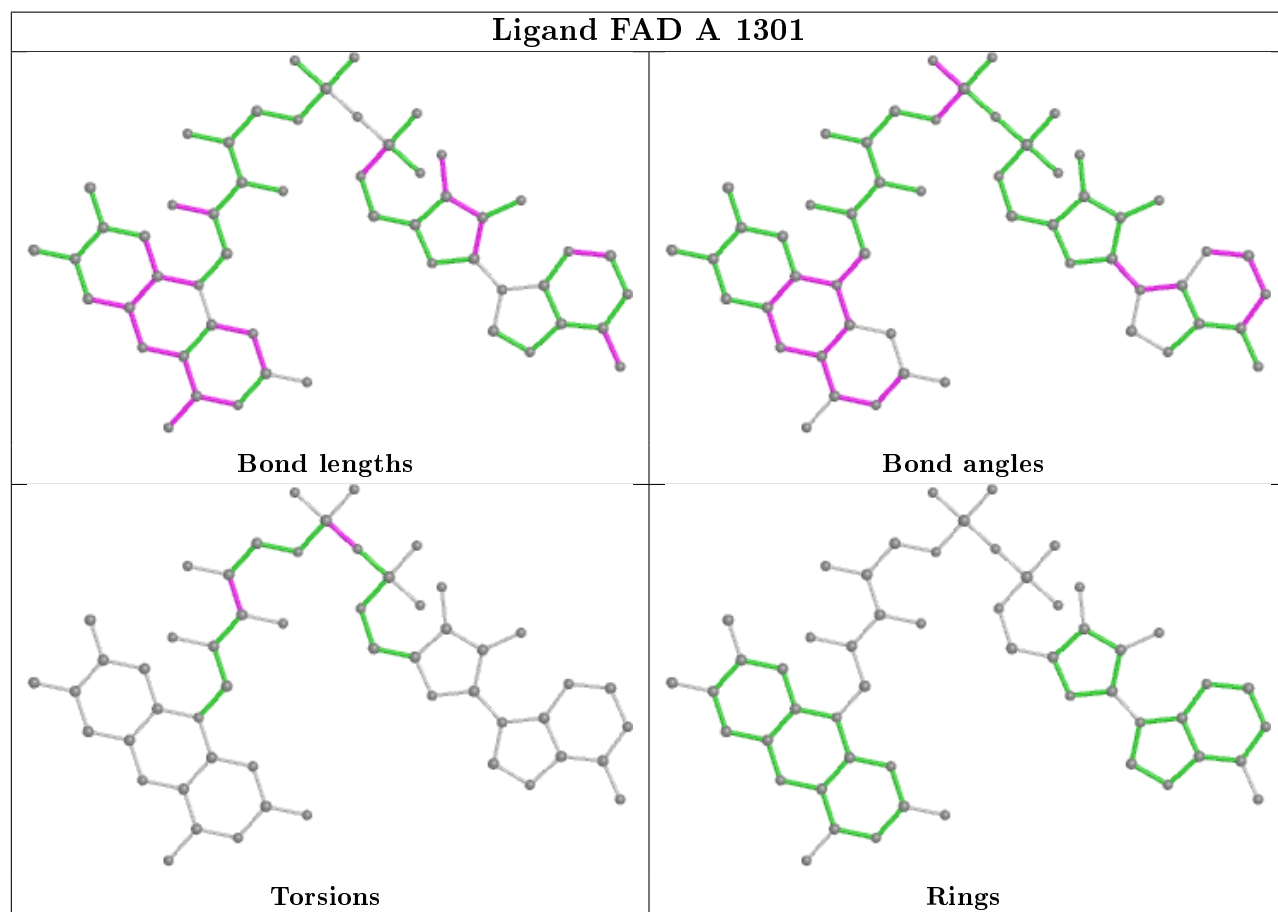
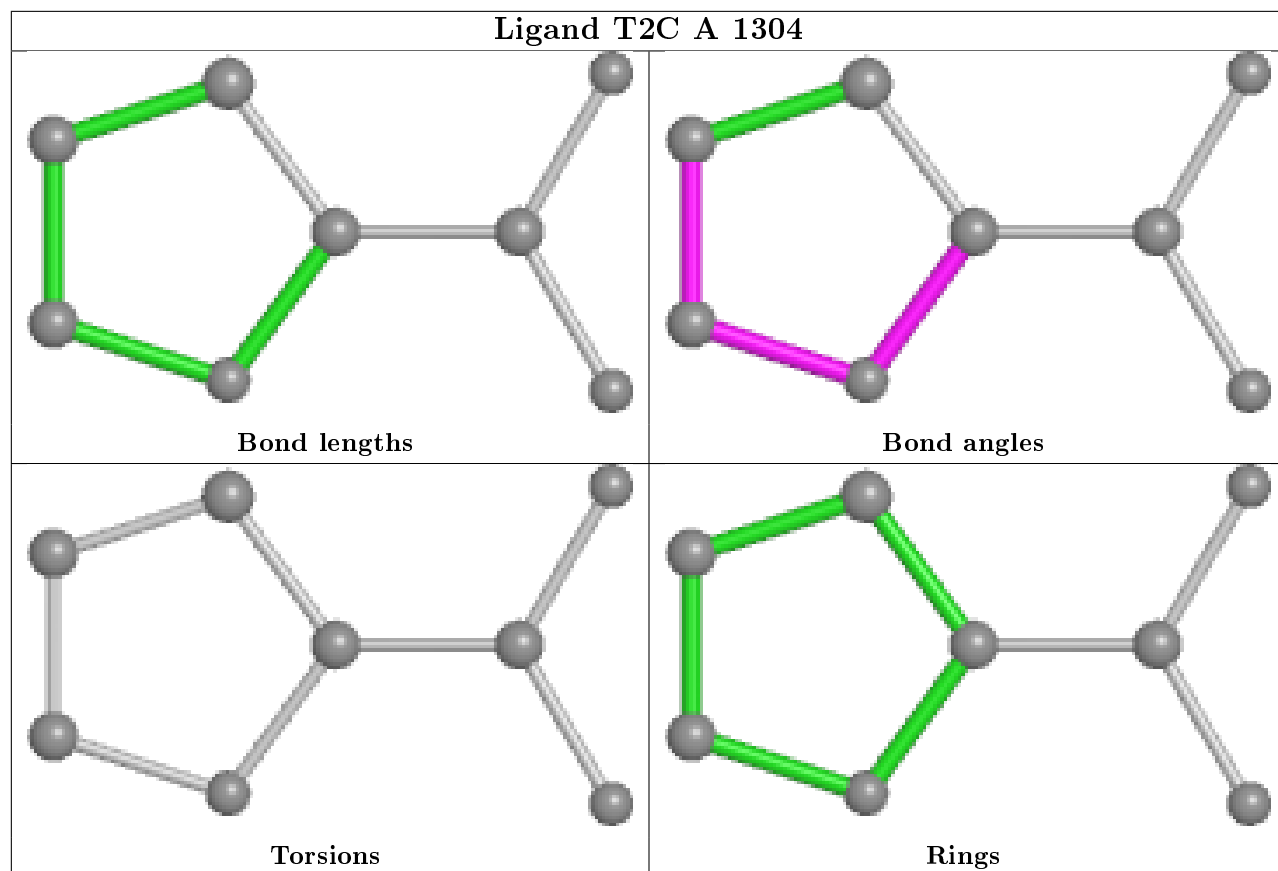
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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.









## 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 ⓘ

### 6.1 Protein, DNA and RNA chains ⓘ

EDS failed to run properly - this section is therefore empty.

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

EDS failed to run properly - this section is therefore empty.

### 6.3 Carbohydrates ⓘ

EDS failed to run properly - this section is therefore empty.

### 6.4 Ligands ⓘ

EDS failed to run properly - this section is therefore empty.

### 6.5 Other polymers ⓘ

EDS failed to run properly - this section is therefore empty.