Full wwPDB X-ray Structure Validation Report (i)

Jan 27, 2024 – 01:48 PM EST

PDB ID	:	1B5T
Title	:	ESCHERICHIA COLI METHYLENETETRAHYDROFOLATE REDUCTASE
Authors	:	Guenther, B.D.; Sheppard, C.A.; Tran, P.; Rozen, R.; Matthews, R.G.; Ludwig, M.L.
Deposited on	:	1999-01-07
Resolution	:	2.50 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 (i) 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 (1)) were used in the production of this report:

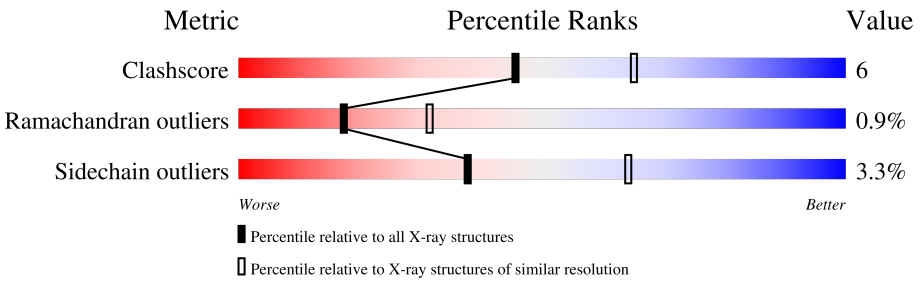
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

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

The reported resolution of this entry is 2.50 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	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)

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 of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=5%

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	275	86% 12% •
1	B	275	80% 19% •
1	C	275	84% 12% • •

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 6564 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 (METHYLENETETRAHYDROFOLATE REDUCTASE).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	275	2116	1342	366	397	11	0	0	0
1	B	274	2095	1328	364	392	11	0	0	0
1	C	267	2007	1280	340	377	10	0	0	0

There are 3 discrepancies between the modelled and reference sequences:

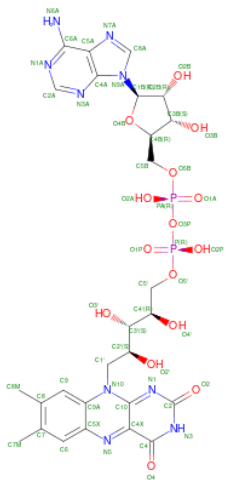
Chain	Residue	Modelled	Actual	Comment	Reference
A	295	ALA	GLY	conflict	UNP P00394
B	295	ALA	GLY	conflict	UNP P00394
C	295	ALA	GLY	conflict	UNP P00394

- Molecule 2 is MERCURY (II) ION (three-letter code: HG) (formula: Hg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Hg		
2	A	1	1	1	0	0
2	B	1	1	1	0	0
2	C	1	1	1	0	0

- Molecule 3 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula: $C_{27}H_{33}N_9O_{15}P_2$).

FAD



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

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
4	A	81	81	81	0	0
4	B	61	61	61	0	0
4	C	42	42	42	0	0

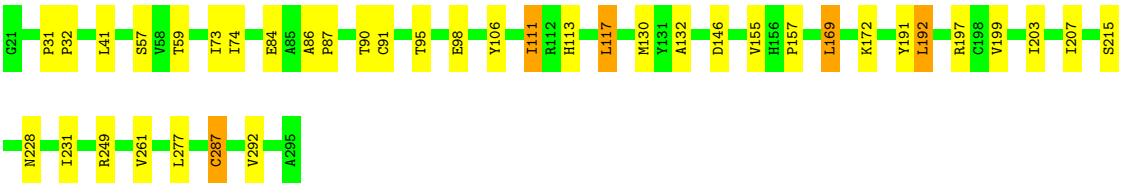
3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. 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 was not executed.

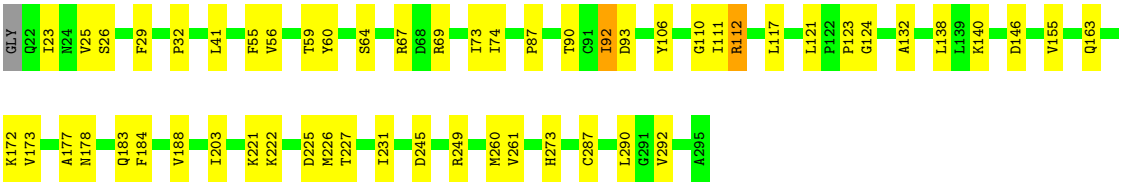
• Molecule 1: PROTEIN (METHYLENETETRAHYDROFOLATE REDUCTASE)

Chain A:



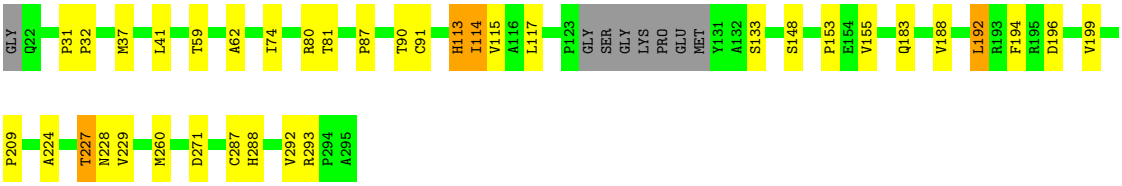
• Molecule 1: PROTEIN (METHYLENETETRAHYDROFOLATE REDUCTASE)

Chain B:



• Molecule 1: PROTEIN (METHYLENETETRAHYDROFOLATE REDUCTASE)

Chain C:



4 Data and refinement statistics (i)

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	103.00Å 128.20Å 98.20Å 90.00° 121.60° 90.00°	Depositor
Resolution (Å)	20.00 – 2.50	Depositor
% Data completeness (in resolution range)	97.0 (20.00-2.50)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
Refinement program	X-PLOR 3.851	Depositor
R, R_{free}	0.212 , 0.263	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	6564	wwPDB-VP
Average B, all atoms (Å^2)	27.0	wwPDB-VP

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: FAD, HG

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	0.38	0/2162	0.60	1/2937 (0.0%)
1	B	0.37	0/2141	0.58	0/2914
1	C	0.35	0/2051	0.56	0/2799
All	All	0.37	0/6354	0.58	1/8650 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	146	ASP	N-CA-C	-5.33	96.60	111.00

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	2116	0	2052	22	0
1	B	2095	0	2006	34	0
1	C	2007	0	1885	23	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
3	A	53	0	31	1	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	53	0	31	1	0
3	C	53	0	31	1	0
4	A	81	0	0	0	0
4	B	61	0	0	0	0
4	C	42	0	0	0	0
All	All	6564	0	6036	78	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 6.

All (78) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:188:VAL:HG22	1:B:260:MET:HG3	1.64	0.79
1:A:197:ARG:HH11	1:A:197:ARG:HG2	1.48	0.79
1:A:111:ILE:O	1:A:111:ILE:HD13	1.97	0.65
1:B:287:CYS:HB3	1:B:292:VAL:HG13	1.78	0.64
1:C:188:VAL:HG22	1:C:260:MET:SD	2.39	0.61
1:B:227:THR:O	1:B:227:THR:HG22	2.02	0.59
1:B:183:GLN:HE21	1:B:184:PHE:H	1.51	0.59
1:C:188:VAL:HG12	1:C:192:LEU:HD22	1.85	0.58
1:C:183:GLN:HG2	1:C:227:THR:HG21	1.87	0.56
1:A:90:THR:HA	1:A:117:LEU:O	2.06	0.56
1:B:64:SER:HA	1:B:67:ARG:HB2	1.89	0.55
1:C:183:GLN:NE2	1:C:227:THR:HG21	2.22	0.55
1:B:117:LEU:HD13	3:B:396:FAD:C4X	2.37	0.55
1:B:74:ILE:HG13	1:B:87:PRO:HB3	1.88	0.53
1:B:23:ILE:HG21	1:B:290:LEU:HG	1.91	0.53
1:A:157:PRO:HG3	1:A:228:ASN:HB2	1.92	0.52
1:C:32:PRO:HD3	1:C:41:LEU:HD22	1.92	0.51
1:A:132:ALA:HB3	1:A:172:LYS:HD3	1.92	0.51
1:B:245:ASP:O	1:B:249:ARG:HG3	2.10	0.51
1:C:87:PRO:HD2	1:C:113:HIS:O	2.11	0.51
1:A:287:CYS:HB3	1:A:292:VAL:HG23	1.94	0.50
1:B:146:ASP:HA	1:B:178:ASN:OD1	2.11	0.50
1:C:113:HIS:HE1	1:C:148:SER:OG	1.94	0.50
1:A:117:LEU:HD13	3:A:395:FAD:C4X	2.42	0.50
1:C:74:ILE:HG13	1:C:87:PRO:HB3	1.92	0.49
1:C:114:ILE:HG12	1:C:115:VAL:N	2.25	0.49
1:B:110:GLY:HA2	1:B:112:ARG:HH21	1.77	0.48
1:C:224:ALA:HB1	1:C:229:VAL:O	2.13	0.48

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:74:ILE:HG13	1:A:87:PRO:HB3	1.95	0.48
1:C:287:CYS:HB3	1:C:292:VAL:HB	1.94	0.48
1:B:26:SER:OG	1:B:273:HIS:HD2	1.96	0.48
1:B:32:PRO:HD3	1:B:41:LEU:HD22	1.94	0.48
1:C:227:THR:O	1:C:228:ASN:HB3	2.14	0.48
1:B:222:LYS:O	1:B:226:MET:HG2	2.14	0.48
1:C:183:GLN:HE21	1:C:227:THR:HG21	1.79	0.48
1:A:197:ARG:HG2	1:A:197:ARG:NH1	2.24	0.47
1:B:261:VAL:HG13	1:B:290:LEU:HD13	1.96	0.47
1:C:117:LEU:HD13	3:C:397:FAD:C4X	2.44	0.47
1:A:215:SER:O	1:A:249:ARG:HD2	2.15	0.47
1:A:57:SER:HA	1:A:86:ALA:O	2.14	0.47
1:A:32:PRO:HD3	1:A:41:LEU:HD22	1.97	0.47
1:A:95:THR:OG1	1:A:98:GLU:HG3	2.14	0.47
1:C:90:THR:HA	1:C:117:LEU:O	2.15	0.46
1:B:92:ILE:HG22	1:B:93:ASP:N	2.30	0.46
1:B:106:TYR:HB3	1:B:111:ILE:HB	1.98	0.46
1:C:288:HIS:CD2	1:C:293:ARG:HG2	2.51	0.46
1:C:80:ARG:HG3	1:C:81:THR:HG23	1.97	0.46
1:B:173:VAL:HG21	1:B:203:ILE:HG23	1.97	0.46
1:C:227:THR:HB	1:C:229:VAL:HG23	1.98	0.46
1:A:91:CYS:SG	1:A:130:MET:CE	3.03	0.46
1:A:169:LEU:HD13	1:A:203:ILE:HD13	1.99	0.45
1:A:231:ILE:HG13	1:A:231:ILE:O	2.16	0.45
1:A:32:PRO:HD2	1:A:73:ILE:HD11	1.99	0.44
1:B:69:ARG:O	1:B:73:ILE:HD13	2.17	0.44
1:A:207:ILE:O	1:A:207:ILE:HG13	2.18	0.43
1:A:292:VAL:HG23	1:A:292:VAL:O	2.17	0.43
1:B:132:ALA:HB3	1:B:172:LYS:HD3	1.99	0.43
1:C:31:PRO:HA	1:C:32:PRO:HD2	1.92	0.43
1:B:25:VAL:HG12	1:B:26:SER:N	2.33	0.43
1:B:25:VAL:CG1	1:B:26:SER:N	2.82	0.43
1:B:26:SER:HA	1:B:55:PHE:O	2.19	0.43
1:B:60:TYR:HD2	1:B:90:THR:HB	1.83	0.43
1:B:261:VAL:CG1	1:B:290:LEU:HD13	2.49	0.43
1:A:31:PRO:HA	1:A:32:PRO:HD2	1.93	0.42
1:C:153:PRO:HA	1:C:194:PHE:CG	2.54	0.42
1:B:221:LYS:O	1:B:225:ASP:HB2	2.19	0.42
1:C:37:MET:HB3	1:C:37:MET:HE2	1.96	0.42
1:A:199:VAL:HG23	1:B:163:GLN:NE2	2.34	0.42
1:B:173:VAL:HA	1:B:177:ALA:HB3	2.01	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:138:LEU:C	1:B:138:LEU:HD23	2.39	0.41
1:B:188:VAL:CG2	1:B:260:MET:HG3	2.43	0.41
1:A:191:TYR:HD2	1:A:192:LEU:HD13	1.85	0.41
1:B:29:PHE:CE2	1:B:56:VAL:HG13	2.56	0.41
1:C:196:ASP:O	1:C:199:VAL:HG12	2.20	0.41
1:C:209:PRO:HD2	1:C:271:ASP:O	2.21	0.41
1:B:60:TYR:CD2	1:B:90:THR:HB	2.57	0.40
1:B:140:LYS:HD3	1:B:140:LYS:HA	1.93	0.40
1:B:221:LYS:HA	1:B:231:ILE:HD11	2.04	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	273/275 (99%)	269 (98%)	3 (1%)	1 (0%)	34 54
1	B	272/275 (99%)	260 (96%)	8 (3%)	4 (2%)	10 18
1	C	263/275 (96%)	254 (97%)	7 (3%)	2 (1%)	19 35
All	All	808/825 (98%)	783 (97%)	18 (2%)	7 (1%)	17 31

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	62	ALA
1	A	155	VAL
1	B	92	ILE
1	B	155	VAL
1	C	155	VAL
1	B	124	GLY
1	B	123	PRO

5.3.2 Protein sidechains (i)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	220/232 (95%)	209 (95%)	11 (5%)	24	46
1	B	215/232 (93%)	212 (99%)	3 (1%)	67	86
1	C	199/232 (86%)	192 (96%)	7 (4%)	36	62
All	All	634/696 (91%)	613 (97%)	21 (3%)	38	64

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	59	THR
1	A	84	GLU
1	A	106	TYR
1	A	111	ILE
1	A	113	HIS
1	A	117	LEU
1	A	169	LEU
1	A	192	LEU
1	A	261	VAL
1	A	277	LEU
1	A	287	CYS
1	B	59	THR
1	B	112	ARG
1	B	121	LEU
1	C	59	THR
1	C	91	CYS
1	C	113	HIS
1	C	114	ILE
1	C	133	SER
1	C	192	LEU
1	C	227	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	113	HIS
1	B	163	GLN
1	B	183	GLN
1	B	273	HIS
1	C	113	HIS
1	C	178	ASN
1	C	288	HIS

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 6 ligands modelled in this entry, 3 are monoatomic - leaving 3 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	$\#\|Z\| > 2$
3	FAD	C	397	-	53,58,58	1.19	4 (7%)	68,89,89	1.45	11 (16%)
3	FAD	B	396	-	53,58,58	1.25	6 (11%)	68,89,89	1.30	7 (10%)
3	FAD	A	395	-	53,58,58	1.32	6 (11%)	68,89,89	1.36	7 (10%)

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	FAD	C	397	-	-	3/30/50/50	0/6/6/6
3	FAD	B	396	-	-	0/30/50/50	0/6/6/6
3	FAD	A	395	-	-	6/30/50/50	0/6/6/6

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	397	FAD	C4X-N5	3.98	1.38	1.30
3	A	395	FAD	C4X-N5	3.59	1.37	1.30
3	A	395	FAD	C10-N10	3.48	1.44	1.37
3	B	396	FAD	C10-N10	3.37	1.44	1.37
3	C	397	FAD	C1'-C2'	3.19	1.57	1.52
3	B	396	FAD	C9-C8	-3.12	1.35	1.39
3	A	395	FAD	C8A-N7A	-3.08	1.29	1.34
3	B	396	FAD	C4X-N5	3.02	1.36	1.30
3	B	396	FAD	C6-C7	-2.98	1.35	1.39
3	C	397	FAD	C10-N10	2.87	1.43	1.37
3	A	395	FAD	C7M-C7	2.55	1.56	1.51
3	A	395	FAD	C9-C8	-2.24	1.36	1.39
3	B	396	FAD	C8A-N7A	-2.18	1.30	1.34
3	A	395	FAD	C6-C7	-2.15	1.36	1.39
3	B	396	FAD	C4X-C10	-2.07	1.38	1.44
3	C	397	FAD	C8A-N7A	-2.05	1.31	1.34

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	A	395	FAD	C9A-C5X-N5	5.40	128.30	122.43
3	C	397	FAD	C9A-C5X-N5	5.20	128.09	122.43
3	B	396	FAD	C9A-C5X-N5	4.65	127.49	122.43
3	C	397	FAD	C5X-N5-C4X	-3.77	111.81	118.07
3	A	395	FAD	C5X-N5-C4X	-3.58	112.12	118.07
3	B	396	FAD	C5X-N5-C4X	-3.45	112.34	118.07
3	B	396	FAD	C6-C5X-N5	-2.60	113.98	118.51
3	C	397	FAD	C6-C5X-N5	-2.58	114.01	118.51
3	A	395	FAD	O5'-C5'-C4'	2.50	116.03	109.36
3	A	395	FAD	C6-C5X-N5	-2.50	114.15	118.51
3	C	397	FAD	O3B-C3B-C4B	-2.46	103.93	111.05
3	B	396	FAD	O3B-C3B-C4B	-2.44	104.00	111.05
3	A	395	FAD	O3'-C3'-C4'	2.43	114.67	108.81
3	C	397	FAD	O2-C2-N3	2.36	123.24	118.65

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	C	397	FAD	C1B-N9A-C4A	-2.35	122.51	126.64
3	A	395	FAD	P-O3P-PA	-2.34	124.80	132.83
3	B	396	FAD	C4-C4X-N5	-2.22	115.08	118.23
3	C	397	FAD	O2P-P-O1P	2.18	123.00	112.24
3	B	396	FAD	O2'-C2'-C1'	2.17	115.05	109.80
3	B	396	FAD	C4'-C3'-C2'	2.17	117.87	113.36
3	C	397	FAD	C4A-C5A-N7A	2.10	111.59	109.40
3	A	395	FAD	C5'-C4'-C3'	2.03	116.14	112.20
3	C	397	FAD	C9-C9A-N10	2.03	124.58	121.84
3	C	397	FAD	C10-N1-C2	2.01	120.92	116.90
3	C	397	FAD	O4'-C4'-C5'	-2.01	105.40	109.92

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	395	FAD	C5B-O5B-PA-O3P
3	C	397	FAD	C5B-O5B-PA-O3P
3	A	395	FAD	P-O3P-PA-O1A
3	A	395	FAD	O2'-C2'-C3'-C4'
3	C	397	FAD	P-O3P-PA-O1A
3	A	395	FAD	P-O3P-PA-O2A
3	C	397	FAD	P-O3P-PA-O2A
3	A	395	FAD	C5B-O5B-PA-O1A
3	A	395	FAD	C5B-O5B-PA-O2A

There are no ring outliers.

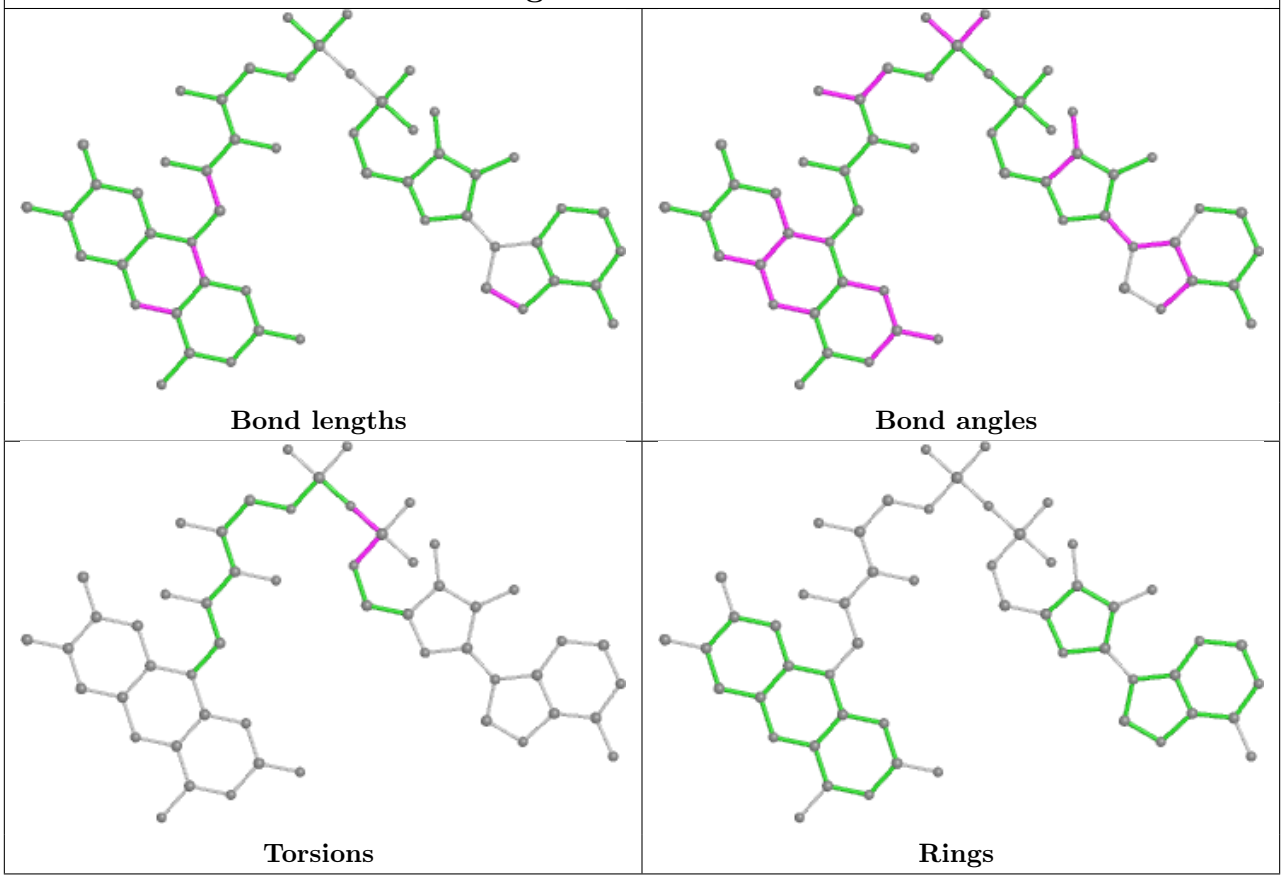
3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	397	FAD	1	0
3	B	396	FAD	1	0
3	A	395	FAD	1	0

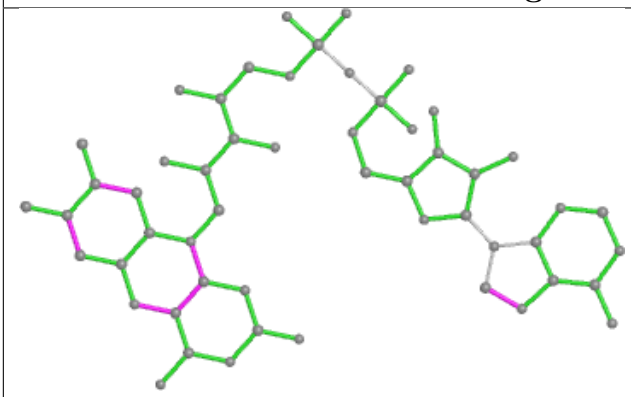
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less then 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the

average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

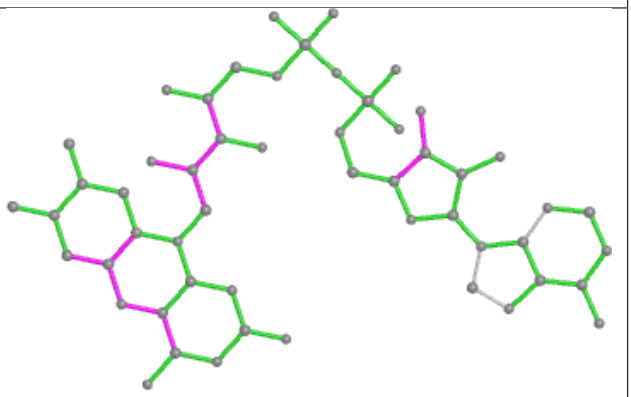
Ligand FAD C 397



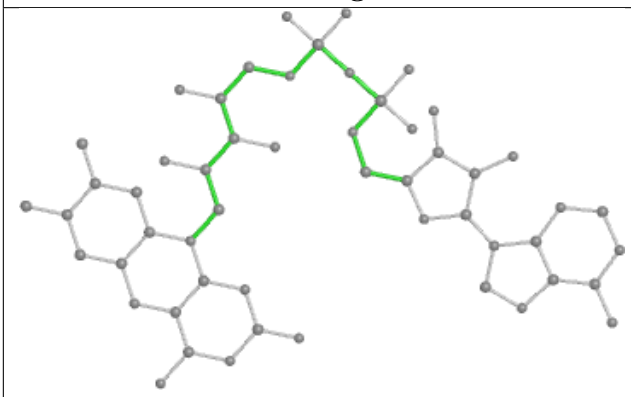
Ligand FAD B 396



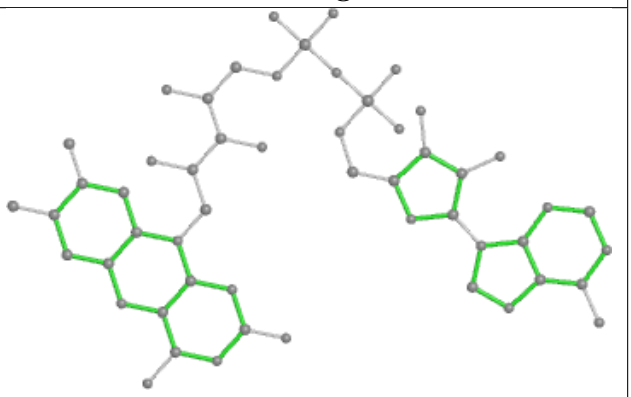
Bond lengths



Bond angles

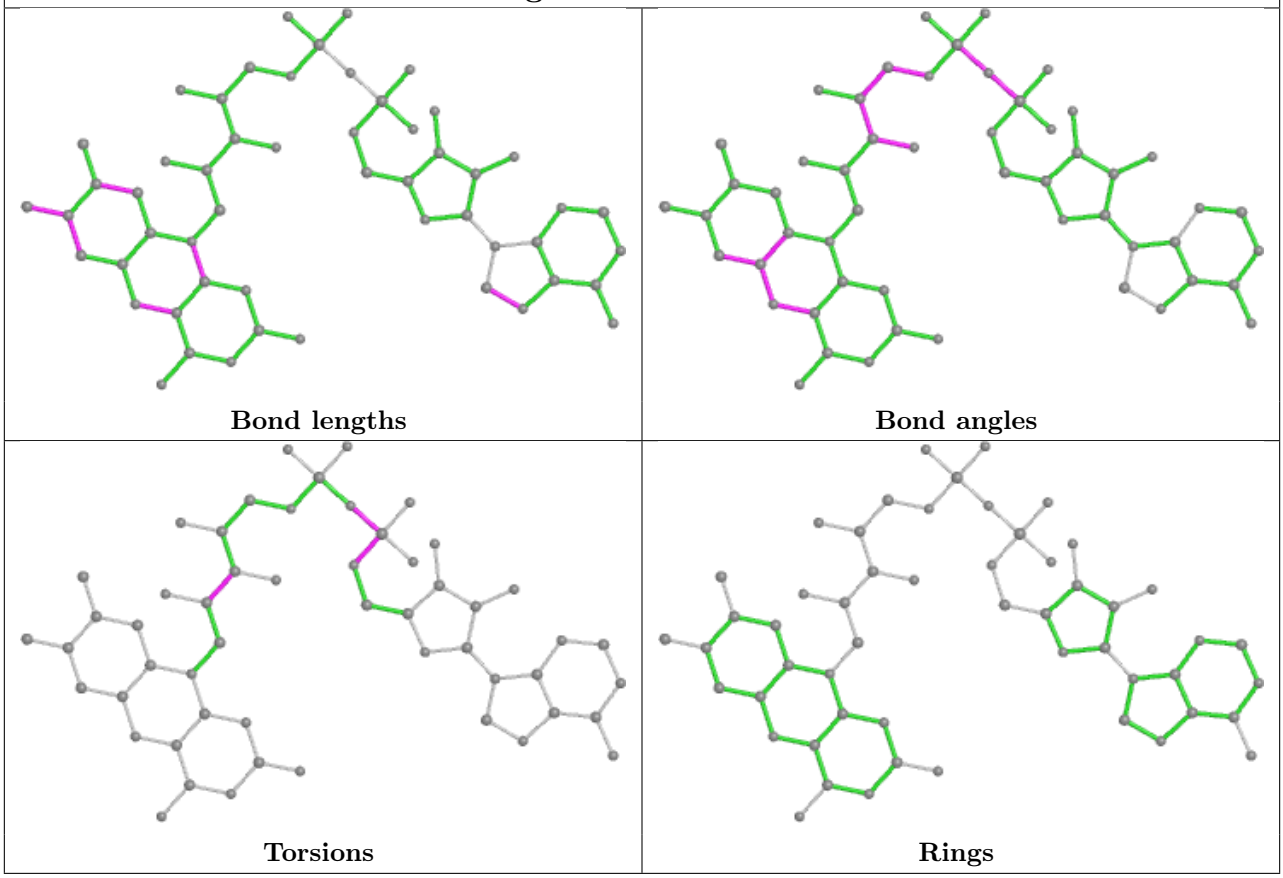


Torsions



Rings

Ligand FAD A 395



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)

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates (i)

EDS was not executed - this section is therefore empty.

6.4 Ligands (i)

EDS was not executed - this section is therefore empty.

6.5 Other polymers (i)

EDS was not executed - this section is therefore empty.