



# Full wwPDB X-ray Structure Validation Report ⓘ

May 16, 2020 – 07:20 am BST

PDB ID : 5OEP  
Title : Mycobacterium tuberculosis DprE1 in complex with inhibitor TCA481  
Authors : Futterer, K.; Batt, S.M.; Besra, G.S.  
Deposited on : 2017-07-09  
Resolution : 2.35 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 : 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



## 2 Entry composition i

There are 5 unique types of molecules in this entry. The entry contains 6553 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 Decaprenylphosphoryl-beta-D-ribose oxidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	426	Total 3209	C 2037	N 564	O 598	S 10	0	0	0
1	B	416	Total 3121	C 1982	N 548	O 581	S 10	0	0	0

There are 32 discrepancies between the modelled and reference sequences:

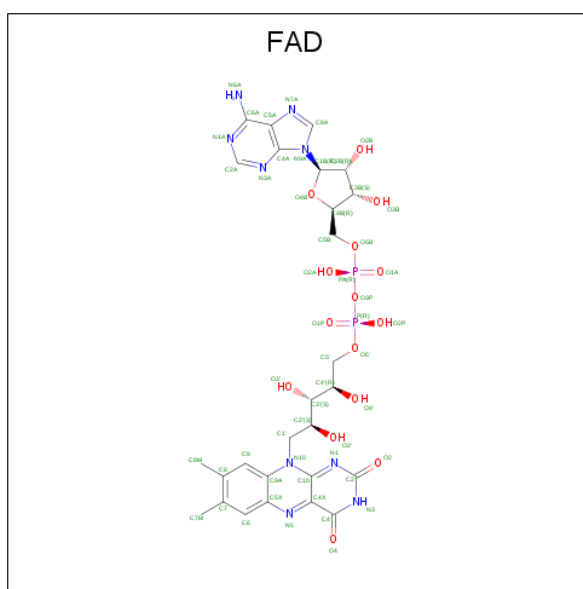
Chain	Residue	Modelled	Actual	Comment	Reference
A	-15	MET	-	initiating methionine	UNP P9WJF1
A	-14	GLY	-	expression tag	UNP P9WJF1
A	-13	SER	-	expression tag	UNP P9WJF1
A	-12	SER	-	expression tag	UNP P9WJF1
A	-11	HIS	-	expression tag	UNP P9WJF1
A	-10	HIS	-	expression tag	UNP P9WJF1
A	-9	HIS	-	expression tag	UNP P9WJF1
A	-8	HIS	-	expression tag	UNP P9WJF1
A	-7	HIS	-	expression tag	UNP P9WJF1
A	-6	HIS	-	expression tag	UNP P9WJF1
A	-5	SER	-	expression tag	UNP P9WJF1
A	-4	GLN	-	expression tag	UNP P9WJF1
A	-3	ASP	-	expression tag	UNP P9WJF1
A	-2	PRO	-	expression tag	UNP P9WJF1
A	-1	ASN	-	expression tag	UNP P9WJF1
A	0	SER	-	expression tag	UNP P9WJF1
B	-15	MET	-	initiating methionine	UNP P9WJF1
B	-14	GLY	-	expression tag	UNP P9WJF1
B	-13	SER	-	expression tag	UNP P9WJF1
B	-12	SER	-	expression tag	UNP P9WJF1
B	-11	HIS	-	expression tag	UNP P9WJF1
B	-10	HIS	-	expression tag	UNP P9WJF1
B	-9	HIS	-	expression tag	UNP P9WJF1
B	-8	HIS	-	expression tag	UNP P9WJF1
B	-7	HIS	-	expression tag	UNP P9WJF1

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-6	HIS	-	expression tag	UNP P9WJF1
B	-5	SER	-	expression tag	UNP P9WJF1
B	-4	GLN	-	expression tag	UNP P9WJF1
B	-3	ASP	-	expression tag	UNP P9WJF1
B	-2	PRO	-	expression tag	UNP P9WJF1
B	-1	ASN	-	expression tag	UNP P9WJF1
B	0	SER	-	expression tag	UNP P9WJF1

- Molecule 2 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula:  $C_{27}H_{33}N_9O_{15}P_2$ ).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
			Total	C	N	O			P
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 ethyl {N}-[2-(1,3-benzothiazol-2-ylcarbonylamino)-5-fluoranyl-thiophen-3-yl]carbonylcarbamate (three-letter code: 9T2) (formula:  $C_{16}H_{12}FN_3O_4S_2$ ).

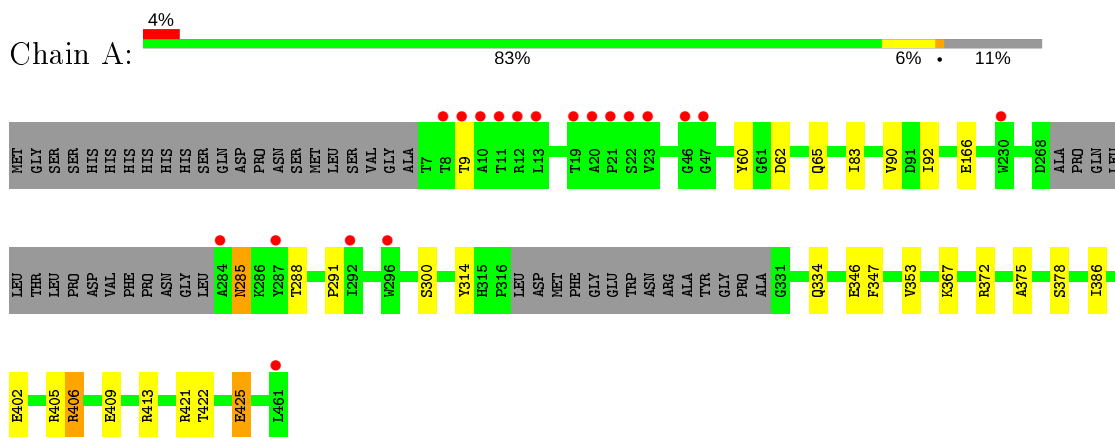


<b>Mol</b>	<b>Chain</b>	<b>Residues</b>	<b>Atoms</b>		<b>ZeroOcc</b>	<b>AltConf</b>
5	A	36	Total 36	O 36	0	0
5	B	24	Total 24	O 24	0	0

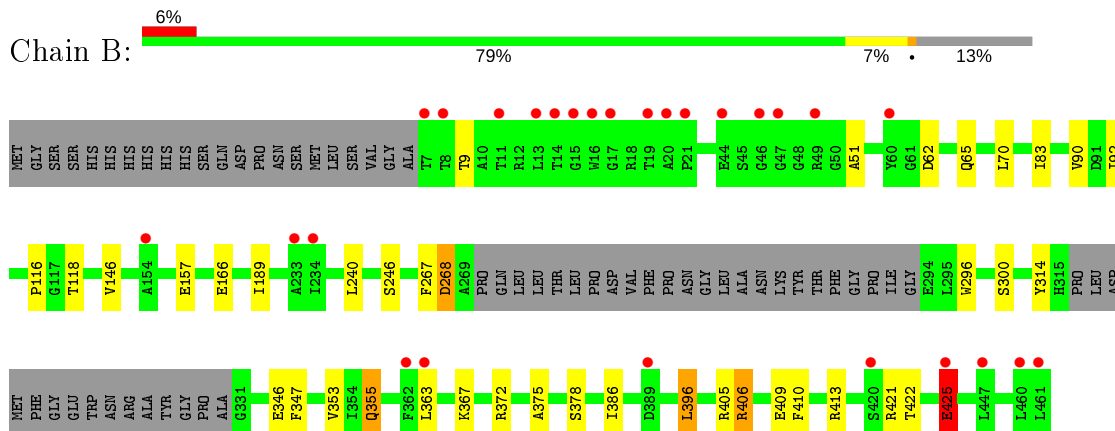
### 3 Residue-property plots [i](#)

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: Decaprenylphosphoryl-beta-D-ribose oxidase



- Molecule 1: Decaprenylphosphoryl-beta-D-ribose oxidase



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	78.30Å 84.54Å 80.97Å 90.00° 103.52° 90.00°	Depositor
Resolution (Å)	57.60 – 2.35 57.61 – 2.35	Depositor EDS
% Data completeness (in resolution range)	98.0 (57.60-2.35) 98.1 (57.61-2.35)	Depositor EDS
$R_{merge}$	0.04	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.08 (at 2.34Å)	Xtrriage
Refinement program	REFMAC 5.8.0135	Depositor
R, $R_{free}$	0.196 , 0.233 0.197 , 0.233	Depositor DCC
$R_{free}$ test set	2132 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	52.4	Xtrriage
Anisotropy	0.424	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.30 , 33.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	0.023 for l,-k,h	Xtrriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	6553	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	64.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.22% 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: 9T2, IMD, FAD

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 5$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# $ Z  > 5$	RMSZ	# $ Z  > 5$
1	A	0.70	1/3282 (0.0%)	0.79	1/4462 (0.0%)
1	B	0.65	0/3190	0.79	4/4339 (0.1%)
All	All	0.68	1/6472 (0.0%)	0.79	5/8801 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	291	PRO	C-O	6.34	1.35	1.23

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	425	GLU	CB-CA-C	-6.13	98.14	110.40
1	B	405	ARG	NE-CZ-NH2	-6.09	117.26	120.30
1	A	372	ARG	NE-CZ-NH2	-5.68	117.46	120.30
1	B	372	ARG	NE-CZ-NH2	-5.26	117.67	120.30
1	B	396	LEU	CA-CB-CG	5.13	127.10	115.30

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	3209	0	3126	16	3
1	B	3121	0	3027	17	3
2	A	53	0	31	0	0
2	B	53	0	31	0	0
3	A	26	0	0	0	0
3	B	26	0	0	0	0
4	A	5	0	5	0	0
5	A	36	0	0	0	0
5	B	24	0	0	0	0
All	All	6553	0	6220	33	3

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 (33) 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:347:PHE:HE1	1:A:386:ILE:CD1	1.64	1.10
1:B:347:PHE:HE1	1:B:386:ILE:CD1	1.65	1.10
1:A:347:PHE:CE1	1:A:386:ILE:HD11	1.90	1.05
1:B:347:PHE:CE1	1:B:386:ILE:HD11	1.92	1.04
1:A:353:VAL:HG21	1:A:406:ARG:NH1	1.75	1.00
1:A:347:PHE:CE1	1:A:386:ILE:CD1	2.50	0.91
1:B:347:PHE:CE1	1:B:386:ILE:CD1	2.53	0.90
1:B:347:PHE:HE1	1:B:386:ILE:HD11	1.28	0.89
1:B:353:VAL:HG21	1:B:406:ARG:NH1	1.90	0.85
1:A:353:VAL:HG21	1:A:406:ARG:HH11	1.47	0.79
1:A:347:PHE:CZ	1:A:386:ILE:HD11	2.20	0.75
1:A:402:GLU:OE2	1:A:405:ARG:NH2	2.19	0.74
1:A:62:ASP:HA	1:A:65:GLN:HE21	1.54	0.72
1:B:347:PHE:CZ	1:B:386:ILE:HD11	2.26	0.70
1:B:267:PHE:HA	1:B:268:ASP:CB	2.25	0.66
1:B:62:ASP:HA	1:B:65:GLN:HE21	1.60	0.65
1:B:240:LEU:HD22	1:B:355:GLN:HG2	1.78	0.64
1:A:353:VAL:HG21	1:A:406:ARG:HH12	1.63	0.60
1:A:83:ILE:HG12	1:A:90:VAL:HG12	1.91	0.53
1:A:402:GLU:CD	1:A:405:ARG:HH22	2.09	0.51
1:A:375:ALA:HB3	1:A:378:SER:HB2	1.94	0.49
1:B:83:ILE:HG12	1:B:90:VAL:HG12	1.95	0.49
1:B:375:ALA:HB3	1:B:378:SER:HB2	1.95	0.47
1:A:60:TYR:CE1	1:A:334:GLN:HG3	2.51	0.46
1:B:353:VAL:HG21	1:B:406:ARG:HH11	1.76	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:285:ASN:O	1:A:288:THR:HG22	2.17	0.44
1:B:146:VAL:HG22	1:B:189:ILE:HD12	2.00	0.43
1:A:346:GLU:N	1:A:346:GLU:OE1	2.51	0.42
1:B:146:VAL:HG22	1:B:189:ILE:CD1	2.50	0.42
1:A:347:PHE:HE1	1:A:386:ILE:HD12	1.69	0.42
1:B:116:PRO:HG2	1:B:118:THR:O	2.20	0.41
1:B:51:ALA:HA	1:B:70:LEU:O	2.20	0.41
1:B:346:GLU:HB3	1:B:410:PHE:CD2	2.56	0.41

All (3) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:425:GLU:OE2	1:B:296:TRP:CG[2_556]	2.08	0.12
1:A:353:VAL:CG1	1:B:425:GLU:OE1[1_656]	2.10	0.10
1:A:425:GLU:OE2	1:B:296:TRP:CD1[2_556]	2.12	0.08

## 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	420/477 (88%)	412 (98%)	8 (2%)	0	100	100
1	B	410/477 (86%)	401 (98%)	8 (2%)	1 (0%)	47	56
All	All	830/954 (87%)	813 (98%)	16 (2%)	1 (0%)	51	63

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	268	ASP

### 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	324/383 (85%)	311 (96%)	13 (4%)	31	39
1	B	312/383 (82%)	295 (95%)	17 (5%)	22	25
All	All	636/766 (83%)	606 (95%)	30 (5%)	26	31

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	THR
1	A	92	ILE
1	A	166	GLU
1	A	285	ASN
1	A	300	SER
1	A	314	TYR
1	A	367	LYS
1	A	406	ARG
1	A	409	GLU
1	A	413	ARG
1	A	421	ARG
1	A	422	THR
1	A	425	GLU
1	B	9	THR
1	B	92	ILE
1	B	157	GLU
1	B	166	GLU
1	B	246	SER
1	B	300	SER
1	B	314	TYR
1	B	355	GLN
1	B	363	LEU
1	B	367	LYS
1	B	396	LEU
1	B	406	ARG
1	B	409	GLU
1	B	413	ARG

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Mol	Chain	Res	Type
1	B	421	ARG
1	B	422	THR
1	B	425	GLU

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

Mol	Chain	Res	Type
1	A	65	GLN

### 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 carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

5 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
2	FAD	A	501	-	51,58,58	1.86	9 (17%)	60,89,89	2.01	10 (16%)
2	FAD	B	501	-	51,58,58	1.99	6 (11%)	60,89,89	1.96	10 (16%)
3	9T2	B	502	-	25,28,28	1.96	3 (12%)	21,39,39	2.62	8 (38%)
3	9T2	A	502	-	25,28,28	1.97	3 (12%)	21,39,39	2.69	8 (38%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
4	IMD	A	503	-	3,5,5	0.32	0	4,5,5	0.51	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
2	FAD	A	501	-	-	2/30/50/50	0/6/6/6
2	FAD	B	501	-	-	2/30/50/50	0/6/6/6
3	9T2	B	502	-	-	2/11/19/19	0/3/3/3
3	9T2	A	502	-	-	2/11/19/19	0/3/3/3
4	IMD	A	503	-	-	-	0/1/1/1

All (21) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	501	FAD	C4X-C10	9.46	1.48	1.38
2	A	501	FAD	C4X-C10	8.96	1.47	1.38
3	B	502	9T2	O21-C21	6.02	1.46	1.35
3	A	502	9T2	O21-C21	5.96	1.46	1.35
3	A	502	9T2	C18-C14	4.91	1.51	1.41
3	B	502	9T2	C18-C14	4.85	1.51	1.41
3	B	502	9T2	C26-N30	4.40	1.37	1.31
2	B	501	FAD	C9A-C5X	4.37	1.51	1.42
3	A	502	9T2	C26-N30	4.32	1.37	1.31
2	B	501	FAD	C8-C7	4.06	1.51	1.40
2	B	501	FAD	C9A-N10	3.94	1.43	1.38
2	A	501	FAD	C4-C4X	3.90	1.48	1.41
2	B	501	FAD	C4-C4X	3.75	1.47	1.41
2	A	501	FAD	C8-C7	3.71	1.50	1.40
2	A	501	FAD	C9A-C5X	3.62	1.49	1.42
2	B	501	FAD	C5A-C4A	2.85	1.48	1.40
2	A	501	FAD	C5A-C4A	2.56	1.47	1.40
2	A	501	FAD	C2A-N3A	2.46	1.36	1.32
2	A	501	FAD	C10-N1	2.39	1.36	1.33
2	A	501	FAD	C2-N1	-2.02	1.34	1.38
2	A	501	FAD	C5'-C4'	2.01	1.54	1.51

All (36) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	501	FAD	C4-N3-C2	8.10	121.98	115.14
2	A	501	FAD	C4-C4X-C10	-7.05	115.28	119.95
2	A	501	FAD	C4-N3-C2	7.02	121.06	115.14
2	A	501	FAD	C1'-N10-C9A	5.99	123.00	118.29
2	B	501	FAD	C1'-N10-C9A	5.87	122.91	118.29
3	B	502	9T2	O21-C21-N20	5.83	117.83	108.92
3	A	502	9T2	O21-C21-N20	5.81	117.79	108.92
3	A	502	9T2	C26-C12-N13	5.73	119.87	113.74
3	B	502	9T2	C26-N30-C29	5.41	114.52	103.78
3	A	502	9T2	C26-N30-C29	5.38	114.45	103.78
2	B	501	FAD	C4-C4X-C10	-5.32	116.43	119.95
3	B	502	9T2	C26-C12-N13	4.77	118.84	113.74
2	A	501	FAD	C4X-N5-C5X	4.02	120.78	116.77
2	B	501	FAD	C4X-N5-C5X	3.97	120.74	116.77
2	B	501	FAD	N3A-C2A-N1A	-3.78	122.78	128.68
3	B	502	9T2	C18-C19-N20	3.76	122.70	115.31
3	A	502	9T2	O12-C12-C26	-3.57	115.36	120.70
2	A	501	FAD	C4-C4X-N5	3.52	122.62	118.60
3	B	502	9T2	O19-C19-C18	-3.43	114.75	121.01
3	A	502	9T2	O21-C21-O22	-3.42	117.69	124.25
2	B	501	FAD	C4X-C4-N3	-3.36	118.84	123.43
2	A	501	FAD	N3A-C2A-N1A	-3.34	123.46	128.68
3	B	502	9T2	O21-C21-O22	-3.25	118.01	124.25
3	A	502	9T2	C37-O21-C21	3.25	122.19	116.04
3	A	502	9T2	C18-C19-N20	3.14	121.47	115.31
3	B	502	9T2	C37-O21-C21	2.97	121.68	116.04
2	A	501	FAD	C4A-C5A-N7A	-2.92	106.36	109.40
2	B	501	FAD	C4A-C5A-N7A	-2.72	106.56	109.40
3	B	502	9T2	O12-C12-C26	-2.62	116.77	120.70
3	A	502	9T2	O19-C19-C18	-2.62	116.22	121.01
2	B	501	FAD	C4-C4X-N5	2.53	121.49	118.60
2	A	501	FAD	O4'-C4'-C3'	-2.47	103.08	109.10
2	B	501	FAD	C9A-N10-C10	-2.41	118.75	121.91
2	A	501	FAD	C4X-C4-N3	-2.39	120.16	123.43
2	B	501	FAD	P-O3P-PA	-2.23	125.18	132.83
2	A	501	FAD	O3B-C3B-C4B	-2.18	104.74	111.05

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	502	9T2	N20-C21-O21-C37
3	A	502	9T2	O22-C21-O21-C37

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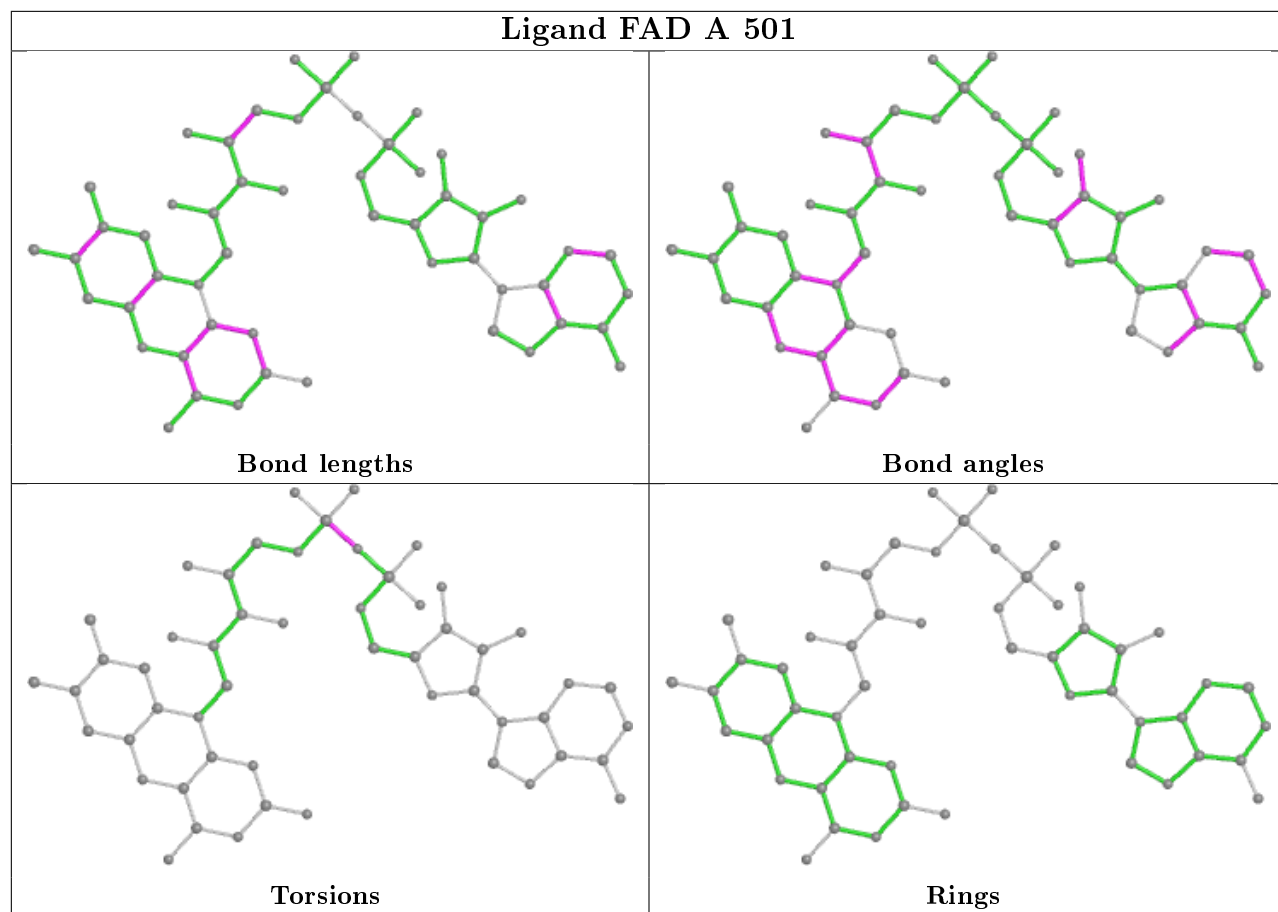
<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>Atoms</b>
3	B	502	9T2	O22-C21-O21-C37
3	B	502	9T2	N20-C21-O21-C37
2	A	501	FAD	PA-O3P-P-O2P
2	B	501	FAD	PA-O3P-P-O2P
2	A	501	FAD	PA-O3P-P-O1P
2	B	501	FAD	PA-O3P-P-O1P

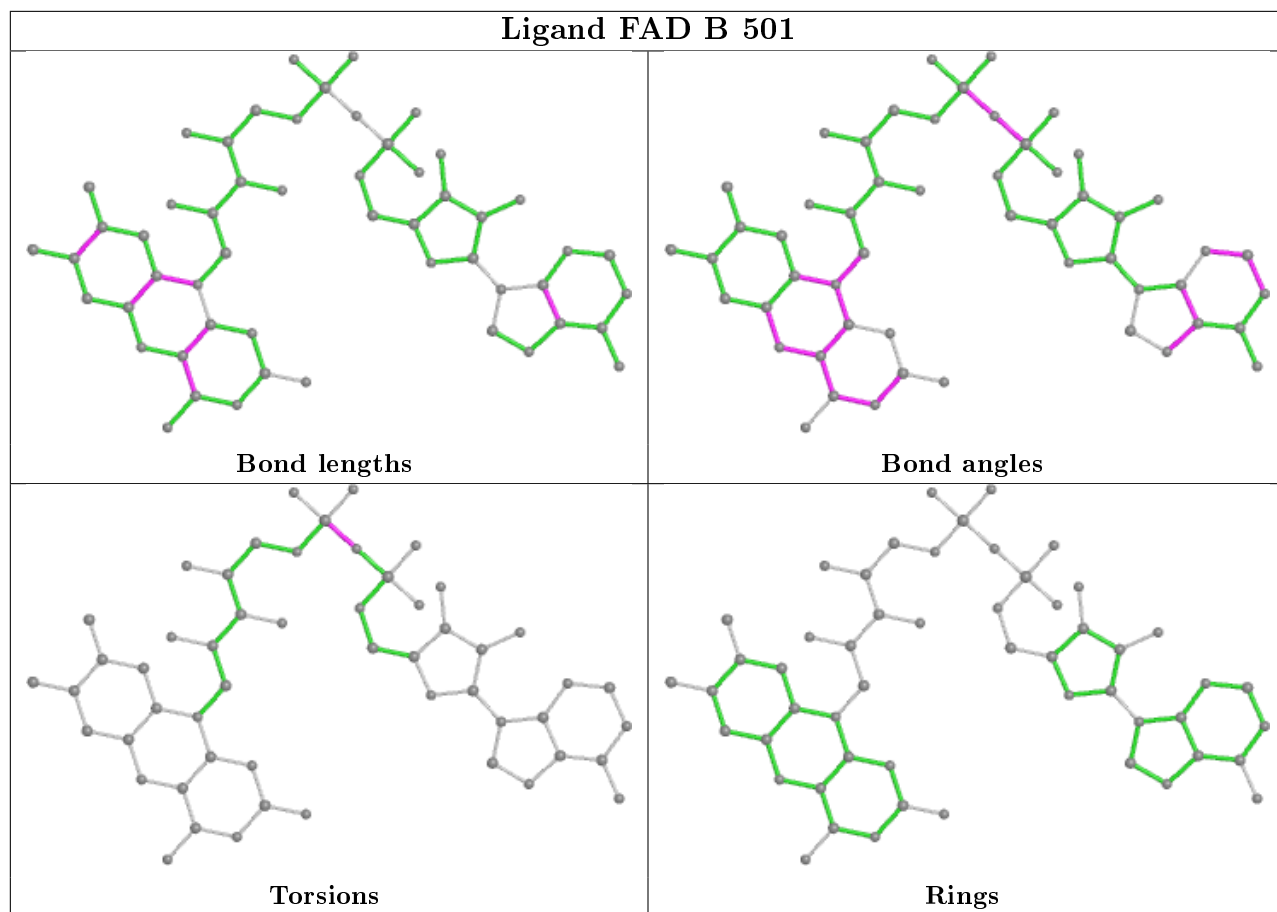
There are no ring outliers.

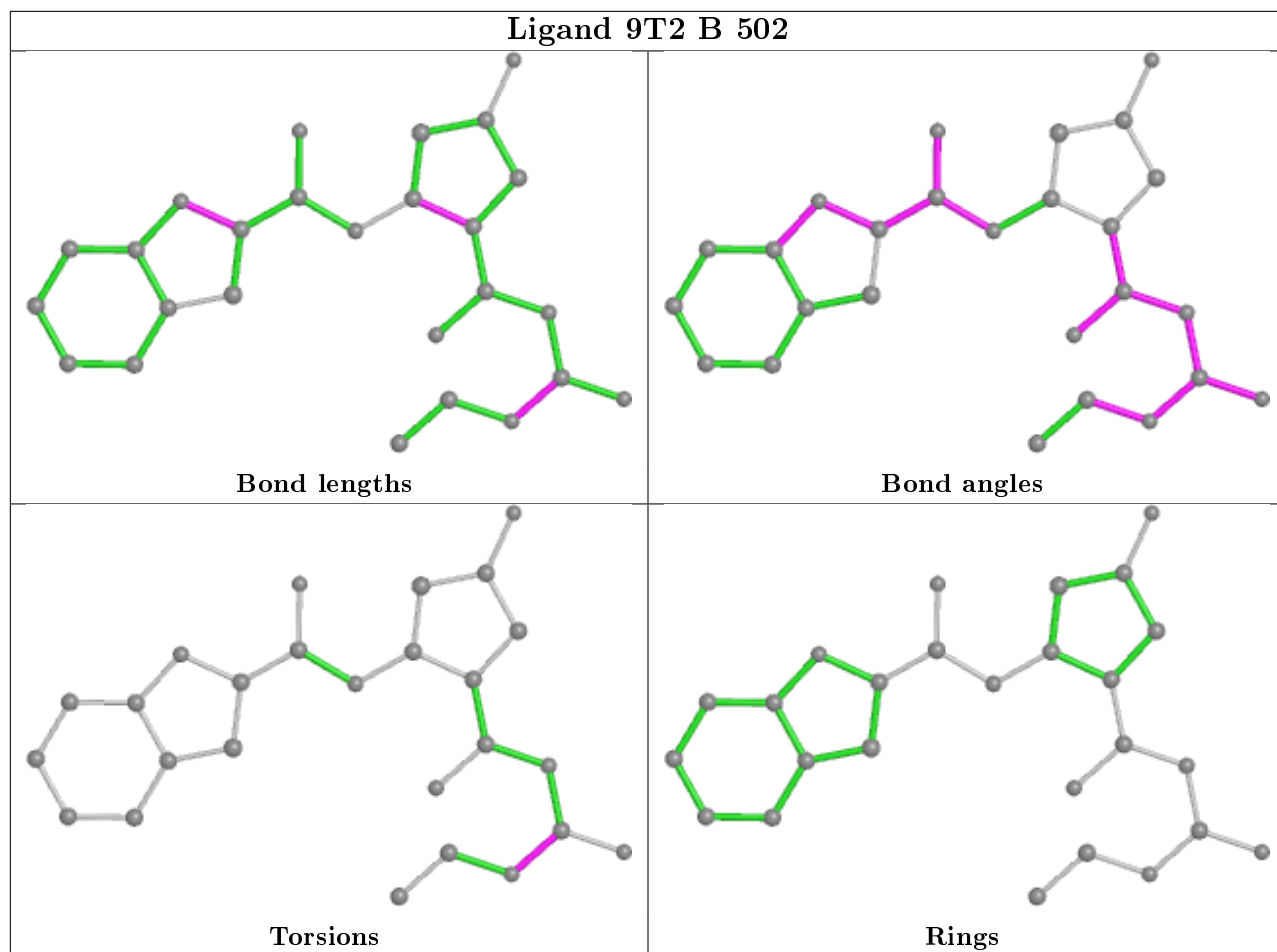
No monomer is involved in short contacts.

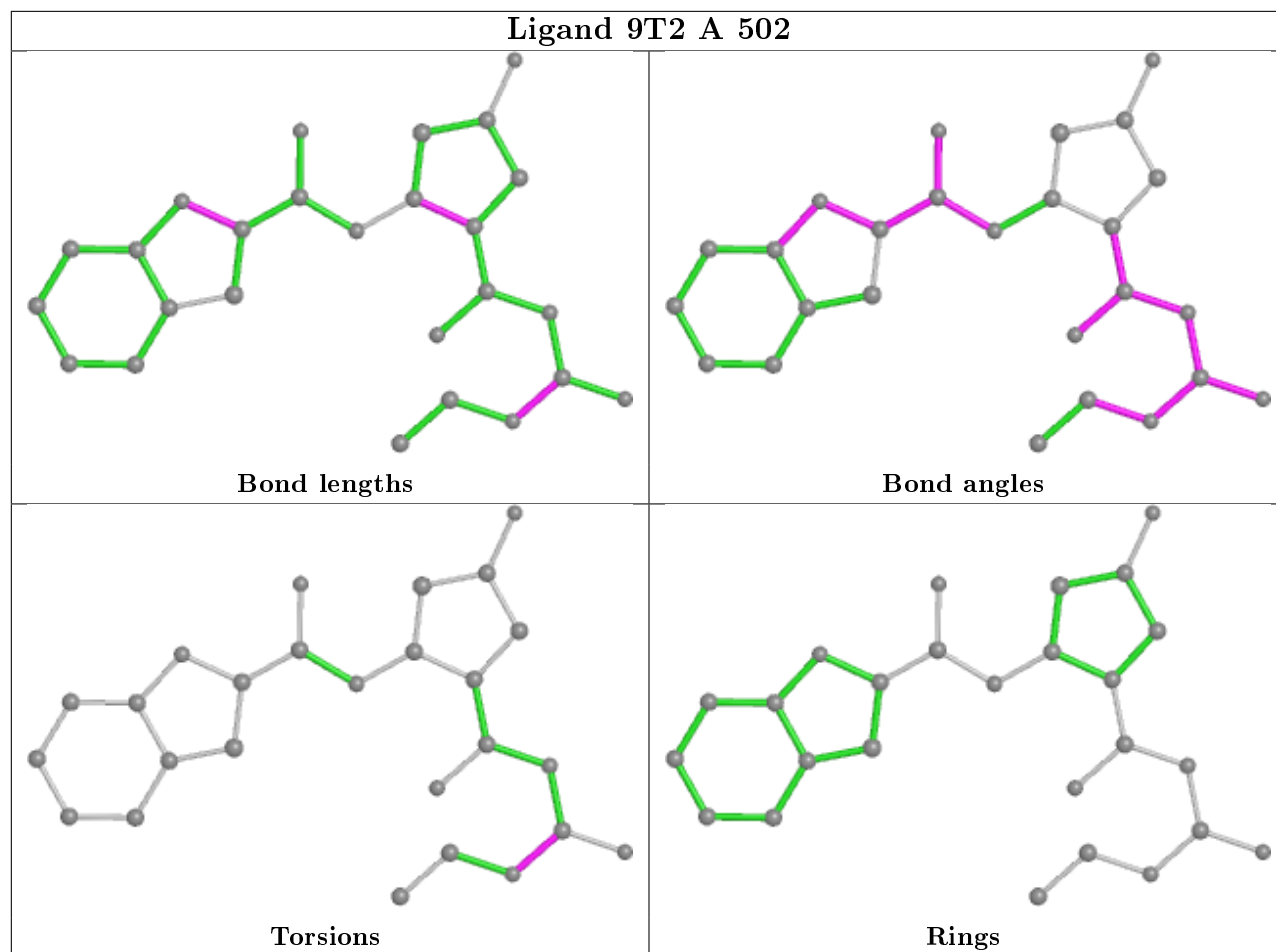
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	426/477 (89%)	0.48	19 (4%) 33 46	33, 57, 103, 135	0
1	B	416/477 (87%)	0.55	27 (6%) 18 27	37, 65, 113, 146	0
All	All	842/954 (88%)	0.52	46 (5%) 25 36	33, 61, 108, 146	0

All (46) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	7	THR	5.5
1	B	46	GLY	5.4
1	B	20	ALA	5.3
1	A	46	GLY	5.2
1	A	23	VAL	5.1
1	B	14	THR	5.1
1	A	47	GLY	5.0
1	B	47	GLY	4.9
1	B	8	THR	4.9
1	B	16	TRP	4.7
1	B	363	LEU	4.6
1	B	60	TYR	4.4
1	A	22	SER	4.4
1	A	19	THR	4.3
1	B	19	THR	4.0
1	B	17	GLY	4.0
1	B	461	LEU	4.0
1	B	460	LEU	3.8
1	B	420	SER	3.5
1	A	20	ALA	3.4
1	B	233	ALA	3.4
1	B	13	LEU	3.4
1	A	8	THR	3.3
1	A	292	ILE	3.1

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Mol	Chain	Res	Type	RSRZ
1	A	9	THR	3.1
1	B	11	THR	3.0
1	A	13	LEU	3.0
1	B	49	ARG	3.0
1	B	362	PHE	2.8
1	A	230	TRP	2.7
1	A	10	ALA	2.6
1	B	15	GLY	2.6
1	B	389	ASP	2.6
1	B	234	ILE	2.6
1	B	44	GLU	2.5
1	A	21	PRO	2.5
1	A	287	TYR	2.4
1	A	461	LEU	2.4
1	B	21	PRO	2.3
1	A	284	ALA	2.3
1	A	296	TRP	2.3
1	B	154	ALA	2.1
1	B	425	GLU	2.1
1	B	447	LEU	2.0
1	A	11	THR	2.0
1	A	12	ARG	2.0

## 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, 95<sup>th</sup> 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.

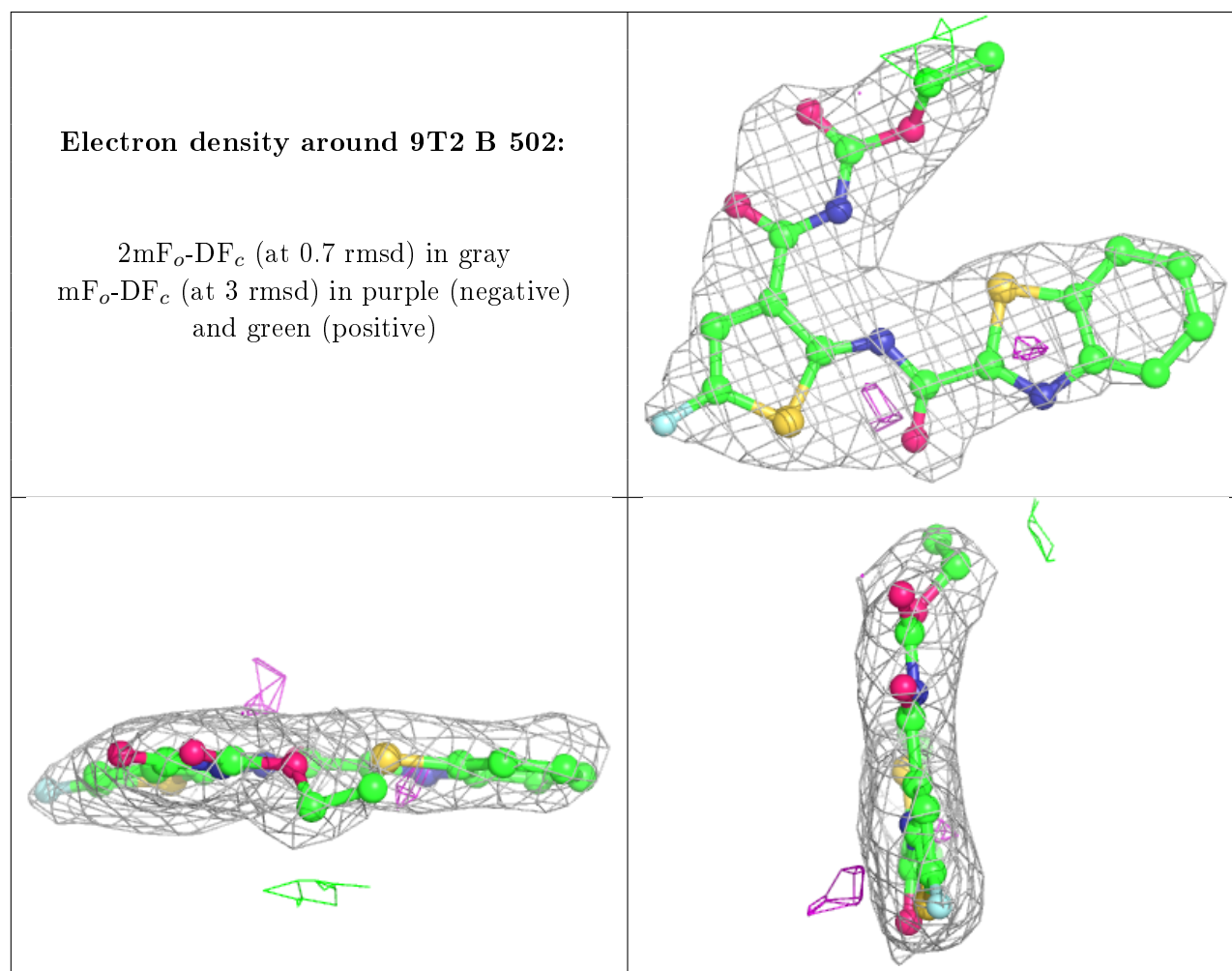
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
3	9T2	B	502	26/26	0.88	0.25	76,85,104,106	0

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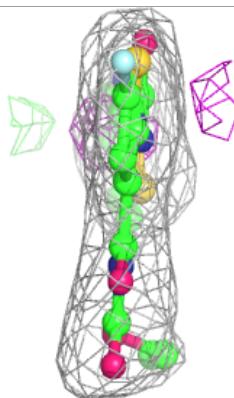
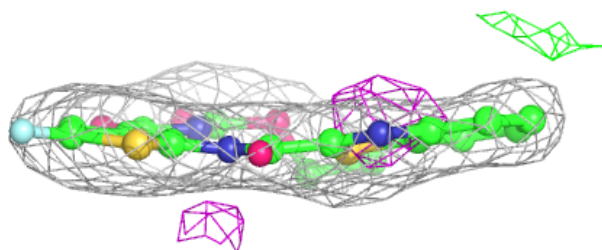
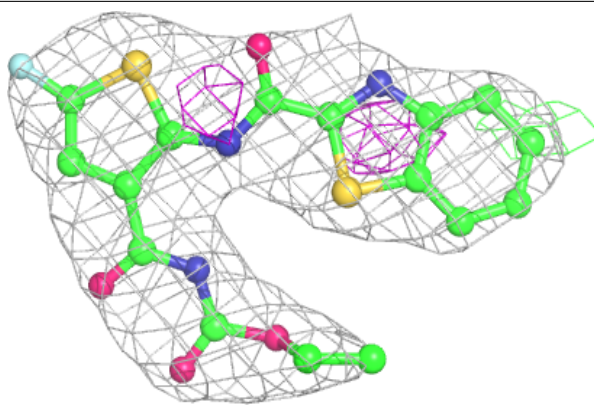
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
3	9T2	A	502	26/26	0.91	0.19	70,78,86,86	0
4	IMD	A	503	5/5	0.95	0.21	70,71,74,75	0
2	FAD	B	501	53/53	0.97	0.14	42,47,51,53	0
2	FAD	A	501	53/53	0.98	0.14	33,40,44,45	0

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.

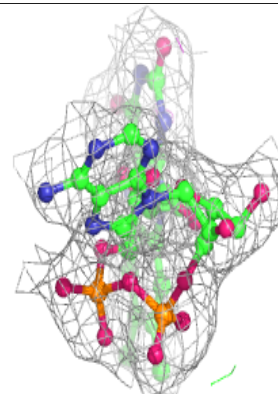
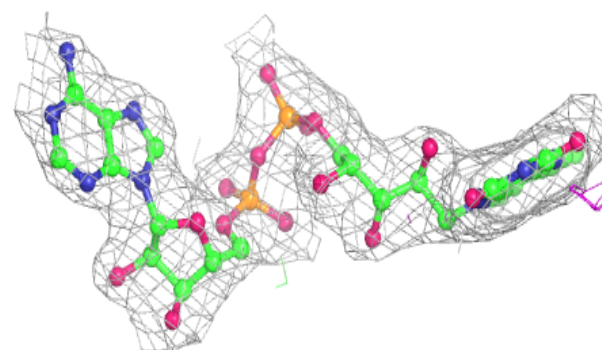
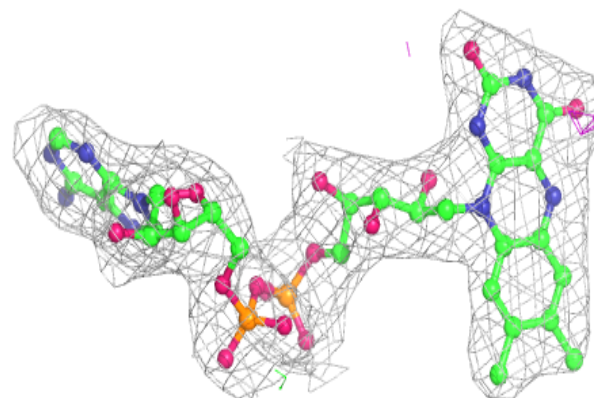


**Electron density around 9T2 A 502:**

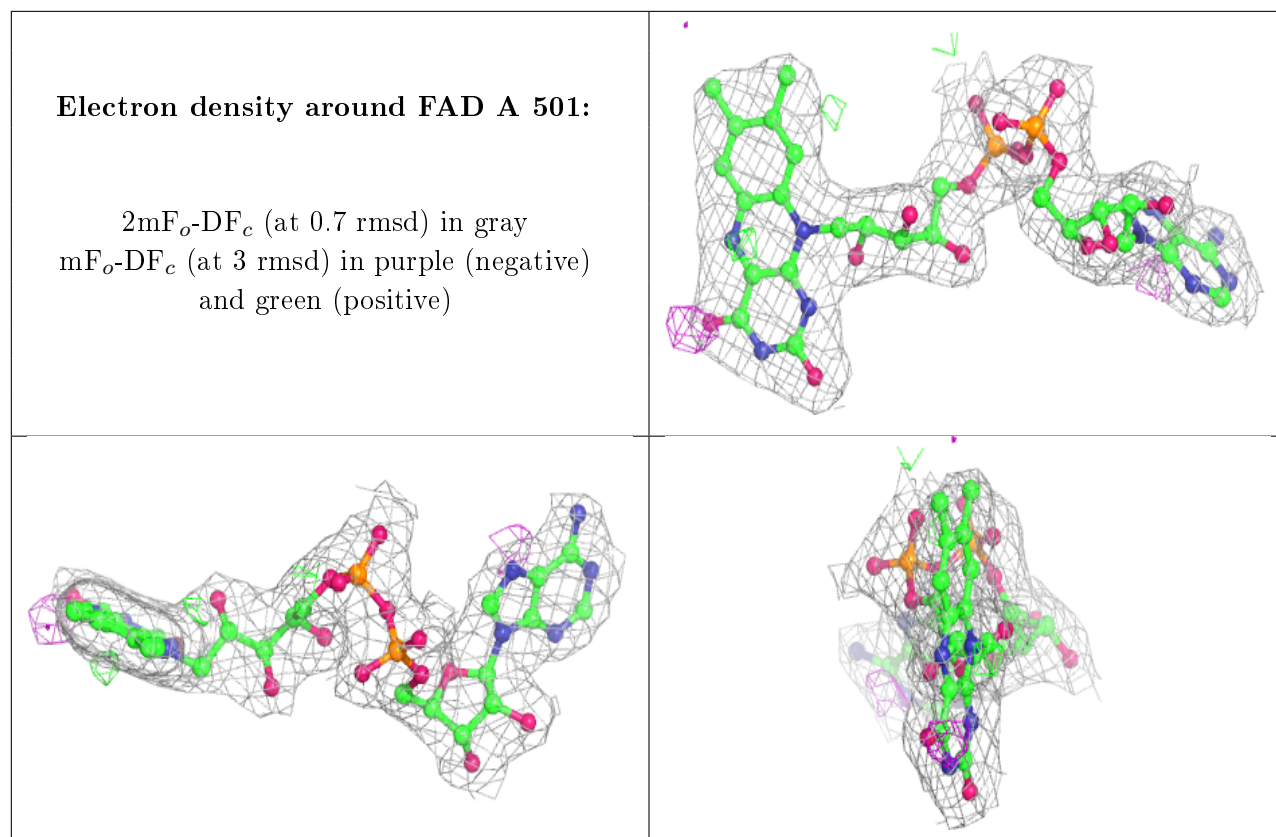
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

**Electron density around FAD B 501:**

$2mF_o-DF_c$  (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.