



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 17, 2026 – 11:52 PM UTC

PDB ID : 8PUY / pdb_00008puy
Title : TEAD2 with a covalent inhibitor
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Deposited on : 2023-07-17
Resolution : 2.20 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The 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 : **FAILED**
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

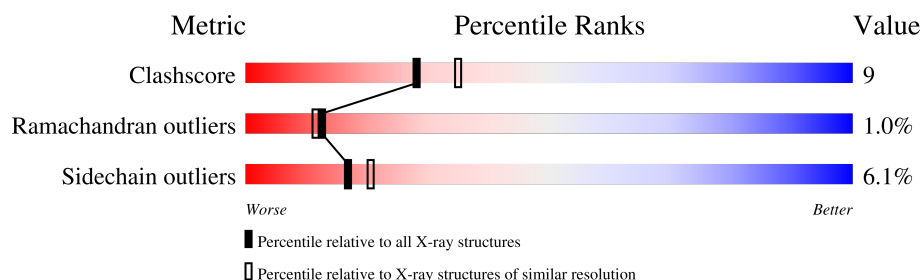
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)

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 $\leq 5\%$

Note EDS failed to run properly.

Mol	Chain	Length	Quality of chain
1	A	240	
1	B	240	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 3363 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.

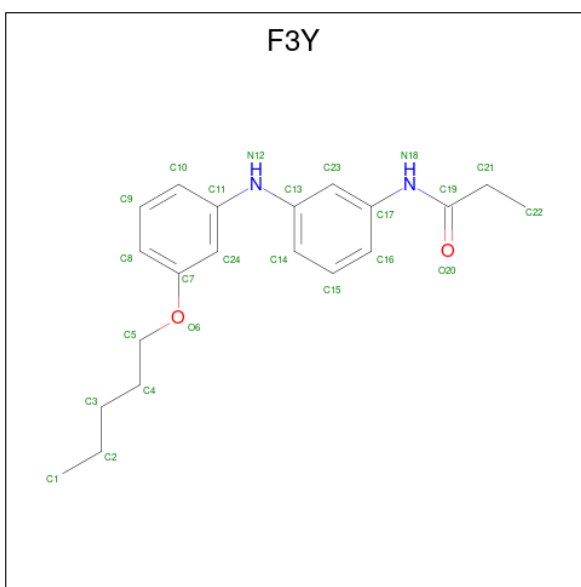
- Molecule 1 is a protein called Transcriptional enhancer factor TEF-4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	211	Total	C	N	O	S	0	1	0
			1738	1114	304	311	9			
1	A	195	Total	C	N	O	S	0	0	0
			1575	1013	268	286	8			

There are 18 discrepancies between the modelled and reference sequences:

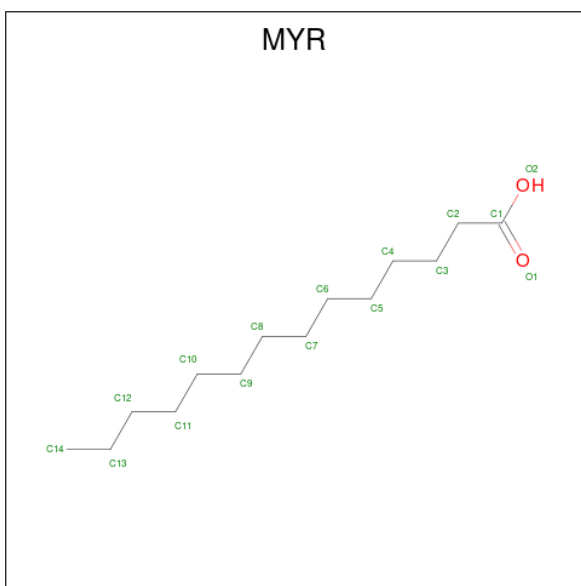
Chain	Residue	Modelled	Actual	Comment	Reference
B	216	MET	-	initiating methionine	UNP Q15562
B	448	VAL	-	expression tag	UNP Q15562
B	449	GLU	-	expression tag	UNP Q15562
B	450	HIS	-	expression tag	UNP Q15562
B	451	HIS	-	expression tag	UNP Q15562
B	452	HIS	-	expression tag	UNP Q15562
B	453	HIS	-	expression tag	UNP Q15562
B	454	HIS	-	expression tag	UNP Q15562
B	455	HIS	-	expression tag	UNP Q15562
A	216	MET	-	initiating methionine	UNP Q15562
A	448	VAL	-	expression tag	UNP Q15562
A	449	GLU	-	expression tag	UNP Q15562
A	450	HIS	-	expression tag	UNP Q15562
A	451	HIS	-	expression tag	UNP Q15562
A	452	HIS	-	expression tag	UNP Q15562
A	453	HIS	-	expression tag	UNP Q15562
A	454	HIS	-	expression tag	UNP Q15562
A	455	HIS	-	expression tag	UNP Q15562

- Molecule 2 is {N}-[3-[(3-pentoxypheyl)amino]phenyl]propanamide (CCD ID: F3Y) (formula: C₂₀H₂₆N₂O₂) (labeled as "Ligand of Interest" by depositor).



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

- Molecule 3 is MYRISTIC ACID (CCD ID: MYR) (formula: $C_{14}H_{28}O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			15	14	1		

- Molecule 4 is water.

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

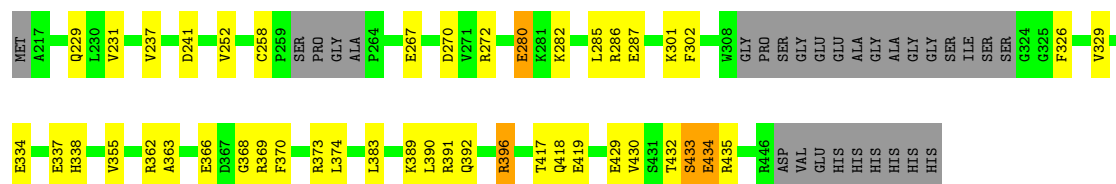
3 Residue-property plots [i](#)

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 failed to run properly.

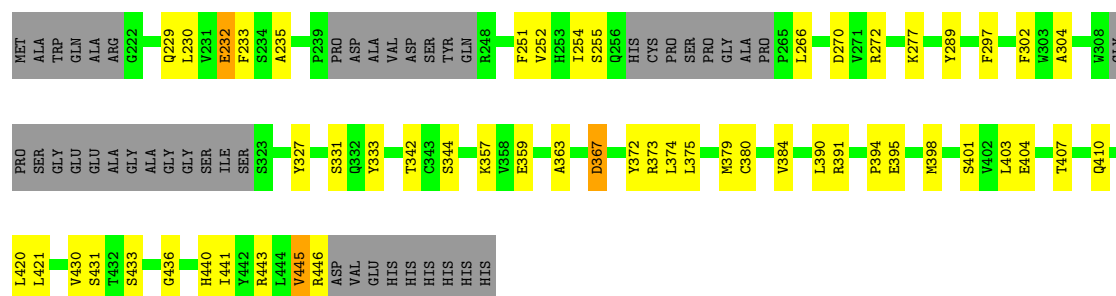
• Molecule 1: Transcriptional enhancer factor TEF-4

Chain B: 



• Molecule 1: Transcriptional enhancer factor TEF-4

Chain A: 



4 Data and refinement statistics

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

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	122.69Å 61.60Å 79.70Å 90.00° 117.51° 90.00°	Depositor
Resolution (Å)	48.38 – 2.20	Depositor
% Data completeness (in resolution range)	85.5 (48.38-2.20)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.33 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.221 , 0.252	Depositor
Wilson B-factor (Å ²)	63.4	Xtriage
Anisotropy	0.147	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3363	wwPDB-VP
Average B, all atoms (Å ²)	76.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.79% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: F3Y, MYR

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	0.96	0/1611	1.20	0/2177
1	B	0.91	0/1786	1.17	1/2416 (0.0%)
All	All	0.94	0/3397	1.19	1/4593 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	231	VAL	N-CA-C	-6.25	105.73	111.67

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	1575	0	1524	34	0
1	B	1738	0	1693	25	0
2	B	24	0	0	1	0
3	A	15	0	27	5	0
4	A	4	0	0	1	0
4	B	7	0	0	0	0
All	All	3363	0	3244	59	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 9.

All (59) 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:327:TYR:HB3	1:A:379:MET:HE2	1.39	1.02
1:A:327:TYR:HB3	1:A:379:MET:CE	2.11	0.81
1:A:342:THR:HG21	4:A:603:HOH:O	1.85	0.77
1:B:301:LYS:HE3	1:B:429:GLU:OE2	1.87	0.74
1:A:394:PRO:HG2	1:A:398:MET:HE1	1.70	0.74
1:A:357:LYS:HD3	3:A:501:MYR:H21	1.74	0.70
1:A:390:LEU:HD21	1:A:403:LEU:HD21	1.77	0.67
1:B:280:GLU:OE2	1:B:286:ARG:NH2	2.30	0.65
1:B:301:LYS:CE	1:B:429:GLU:OE2	2.49	0.60
1:B:432:THR:O	1:B:434:GLU:N	2.34	0.60
1:A:401:SER:O	1:A:404:GLU:HG2	2.01	0.60
1:A:394:PRO:CG	1:A:398:MET:HE1	2.31	0.59
1:B:389:LYS:O	1:B:392:GLN:HG2	2.04	0.58
1:A:380:CYS:N	3:A:501:MYR:O1	2.31	0.56
1:A:420:LEU:HD12	1:A:421:LEU:N	2.23	0.53
1:B:270:ASP:OD1	1:B:272:ARG:HB2	2.08	0.53
1:B:338:HIS:ND1	1:B:363:ALA:HB3	2.22	0.53
1:A:270:ASP:HA	1:A:445:VAL:O	2.08	0.53
1:A:333:TYR:OH	1:A:410:GLN:NE2	2.42	0.52
1:A:270:ASP:OD1	1:A:272:ARG:N	2.36	0.51
1:B:338:HIS:HB2	1:B:370:PHE:CZ	2.45	0.51
1:A:233:PHE:HZ	3:A:501:MYR:H71	1.76	0.50
1:A:289:TYR:CE2	1:A:446:ARG:HG3	2.47	0.50
1:B:417:THR:O	1:B:418:GLN:HB2	2.11	0.50
1:A:363:ALA:HB2	1:A:372:TYR:CE2	2.47	0.49
1:A:266:LEU:HD22	1:A:441:ILE:O	2.13	0.49
1:A:235:ALA:HB3	3:A:501:MYR:H101	1.97	0.47
1:B:237:VAL:HA	1:B:326:PHE:O	2.15	0.47
1:B:280:GLU:OE1	1:B:280:GLU:HA	2.15	0.47
1:A:430:VAL:HG12	1:A:431:SER:O	2.15	0.47
1:B:334:GLU:OE1	1:B:369:ARG:HD3	2.15	0.46
1:B:366:GLU:O	1:B:368:GLY:N	2.48	0.46
1:A:297:PHE:CZ	1:A:420:LEU:HD21	2.51	0.45
1:A:331:SER:OG	1:A:374:LEU:HB2	2.16	0.45
1:A:373:ARG:HB3	1:A:375:LEU:HG	1.98	0.45
1:B:396:ARG:HG2	1:B:430:VAL:CG2	2.46	0.45
1:A:395:GLU:HB2	1:A:398:MET:HE2	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:252:VAL:HB	1:B:302:PHE:CE1	2.52	0.44
1:B:338:HIS:CE1	1:B:363:ALA:HB3	2.53	0.44
1:A:251:PHE:O	1:A:304:ALA:HA	2.18	0.44
1:A:357:LYS:HE2	1:A:359:GLU:OE2	2.19	0.43
1:A:344:SER:O	1:A:410:GLN:HA	2.19	0.43
1:B:383:LEU:HD23	1:B:383:LEU:HA	1.92	0.42
1:B:390:LEU:HD12	1:B:390:LEU:HA	1.85	0.42
1:A:390:LEU:HD12	1:A:390:LEU:HA	1.93	0.42
1:B:417:THR:OG1	1:B:419:GLU:HG2	2.19	0.42
1:A:367:ASP:OD1	1:A:367:ASP:N	2.52	0.42
1:A:394:PRO:CD	1:A:398:MET:HE1	2.50	0.42
1:A:252:VAL:HB	1:A:302:PHE:CE1	2.55	0.41
1:B:285:LEU:HD23	1:B:285:LEU:HA	1.91	0.41
1:B:329:VAL:HG11	2:B:501:F3Y:C14	2.49	0.41
1:B:374:LEU:HD12	1:B:374:LEU:N	2.35	0.41
1:A:380:CYS:O	1:A:384:VAL:HG23	2.21	0.41
1:A:235:ALA:CB	3:A:501:MYR:H101	2.50	0.41
1:A:277:LYS:HB3	1:A:407:THR:HG21	2.03	0.41
1:B:432:THR:O	1:B:435:ARG:N	2.41	0.41
1:B:366:GLU:CD	1:B:373:ARG:HH21	2.29	0.40
1:B:396:ARG:HG2	1:B:430:VAL:HG23	2.03	0.40
1:A:232:GLU:HA	1:A:254:ILE:O	2.21	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	187/240 (78%)	180 (96%)	5 (3%)	2 (1%)	11	10
1	B	206/240 (86%)	195 (95%)	9 (4%)	2 (1%)	12	11
All	All	393/480 (82%)	375 (95%)	14 (4%)	4 (1%)	12	11

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	433	SER
1	B	258	CYS
1	A	232	GLU
1	A	436	GLY

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	171/210 (81%)	162 (95%)	9 (5%)	20	26
1	B	190/210 (90%)	177 (93%)	13 (7%)	14	17
All	All	361/420 (86%)	339 (94%)	22 (6%)	17	20

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	229	GLN
1	B	241	ASP
1	B	267	GLU
1	B	280	GLU
1	B	282	LYS
1	B	287	GLU
1	B	337	GLU
1	B	355	VAL
1	B	362	ARG
1	B	391	ARG
1	B	396	ARG
1	B	433	SER
1	B	434	GLU
1	A	229	GLN
1	A	230	LEU
1	A	255	SER
1	A	367	ASP
1	A	391	ARG
1	A	433	SER

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Mol	Chain	Res	Type
1	A	440	HIS
1	A	443	ARG
1	A	445	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	219	GLN
1	B	229	GLN
1	B	418	GLN
1	A	273	GLN
1	A	410	GLN
1	A	439	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

2 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	MYR	A	501	1	13,14,15	0.42	0	12,13,15	0.35	0
2	F3Y	B	501	1	25,25,25	0.28	0	31,31,31	0.59	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	MYR	A	501	1	-	5/12/12/13	-
2	F3Y	B	501	1	-	2/16/16/16	0/2/2/2

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	501	F3Y	C3-C4-C5-O6
3	A	501	MYR	C10-C11-C12-C13
3	A	501	MYR	C7-C8-C9-C10
3	A	501	MYR	C11-C10-C9-C8
3	A	501	MYR	C1-C2-C3-C4
2	B	501	F3Y	C2-C3-C4-C5
3	A	501	MYR	C6-C7-C8-C9

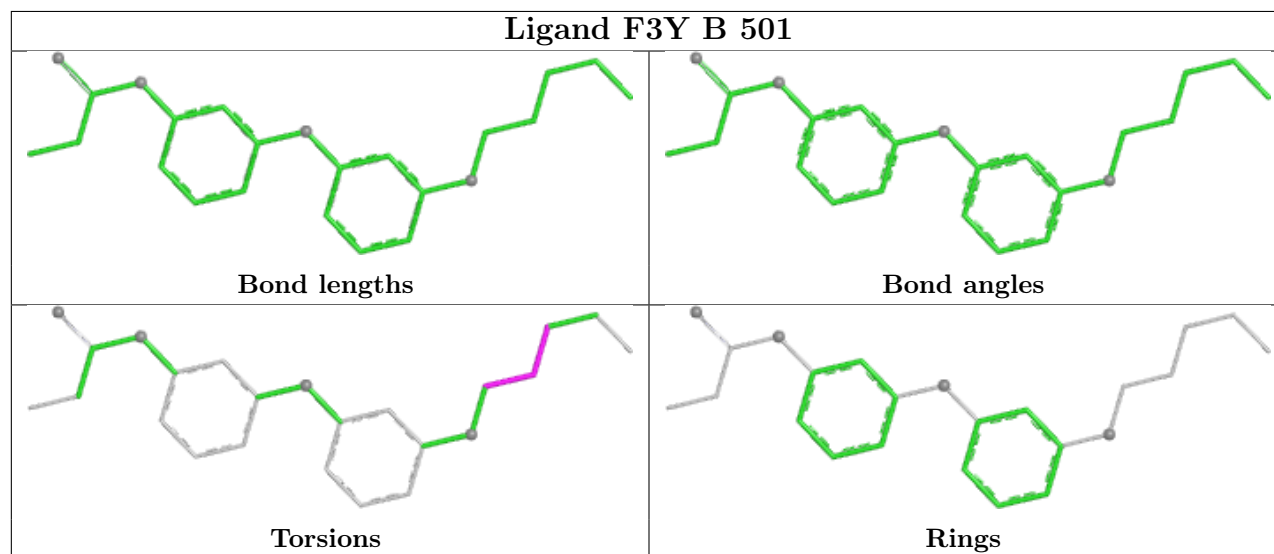
There are no ring outliers.

2 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	501	MYR	5	0
2	B	501	F3Y	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 [i](#)

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

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

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

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

6.3 Carbohydrates [i](#)

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

6.4 Ligands [i](#)

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

6.5 Other polymers [i](#)

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