Full wwPDB X-ray Structure Validation Report ⓘ

Oct 3, 2023 – 08:33 AM EDT

PDB ID	:	6TZ6
Title	:	Crystal Structure of Candida Albicans Calcineurin A, Calcineurin B, FKBP12 and FK506 (Tacrolimus)
Authors	:	Fox III, D.; Lukacs, C.M.
Deposited on	:	2019-08-10
Resolution	:	2.55 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	:	FAILED
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)	:	2.35.1

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 2.55 Å.

There are no overall percentile quality scores available for this entry.

MolProbity and EDS failed to run properly - the sequence quality summary graphics cannot be shown.

2 Entry composition (i)

There are 10 unique types of molecules in this entry. The entry contains 9858 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 Serine/threonine-protein phosphatase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	361	Total	C	N	O	S	0	5	0
			2928	1890	488	530	20			
1	D	352	Total	C	N	O	S	0	3	0
			2821	1822	467	514	18			

There are 38 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	35	MET	-	expression tag	UNP C4YFI3
A	36	GLY	-	expression tag	UNP C4YFI3
A	37	SER	-	expression tag	UNP C4YFI3
A	38	SER	-	expression tag	UNP C4YFI3
A	39	HIS	-	expression tag	UNP C4YFI3
A	40	HIS	-	expression tag	UNP C4YFI3
A	41	HIS	-	expression tag	UNP C4YFI3
A	42	HIS	-	expression tag	UNP C4YFI3
A	43	HIS	-	expression tag	UNP C4YFI3
A	44	HIS	-	expression tag	UNP C4YFI3
A	45	SER	-	expression tag	UNP C4YFI3
A	46	SER	-	expression tag	UNP C4YFI3
A	47	GLY	-	expression tag	UNP C4YFI3
A	48	LEU	-	expression tag	UNP C4YFI3
A	49	VAL	-	expression tag	UNP C4YFI3
A	50	PRO	-	expression tag	UNP C4YFI3
A	51	ARG	-	expression tag	UNP C4YFI3
A	52	GLY	-	expression tag	UNP C4YFI3
A	53	SER	-	expression tag	UNP C4YFI3
D	35	MET	-	expression tag	UNP C4YFI3
D	36	GLY	-	expression tag	UNP C4YFI3
D	37	SER	-	expression tag	UNP C4YFI3
D	38	SER	-	expression tag	UNP C4YFI3
D	39	HIS	-	expression tag	UNP C4YFI3
D	40	HIS	-	expression tag	UNP C4YFI3

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Chain	Residue	Modelled	Actual	Comment	Reference
D	41	HIS	-	expression tag	UNP C4YFI3
D	42	HIS	-	expression tag	UNP C4YFI3
D	43	HIS	-	expression tag	UNP C4YFI3
D	44	HIS	-	expression tag	UNP C4YFI3
D	45	SER	-	expression tag	UNP C4YFI3
D	46	SER	-	expression tag	UNP C4YFI3
D	47	GLY	-	expression tag	UNP C4YFI3
D	48	LEU	-	expression tag	UNP C4YFI3
D	49	VAL	-	expression tag	UNP C4YFI3
D	50	PRO	-	expression tag	UNP C4YFI3
D	51	ARG	-	expression tag	UNP C4YFI3
D	52	GLY	-	expression tag	UNP C4YFI3
D	53	SER	-	expression tag	UNP C4YFI3

- Molecule 2 is a protein called Calcineurin subunit B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	141	Total	C	N	O	S	0	0	0
			1012	641	164	202	5			
2	E	146	Total	C	N	O	S	0	0	0
			1034	653	166	209	6			

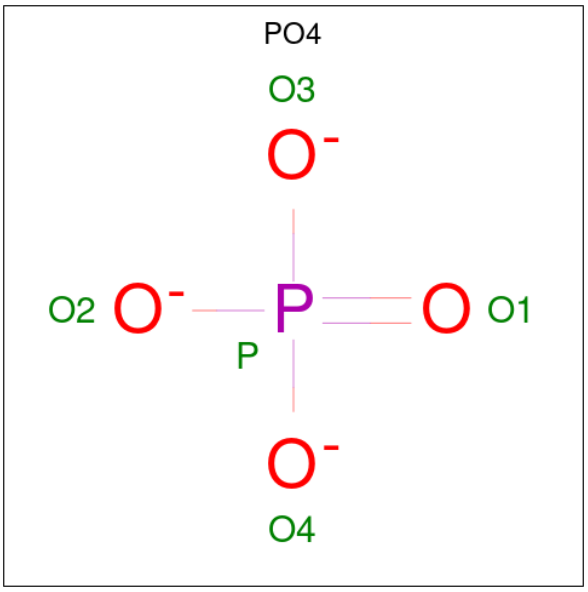
- Molecule 3 is a protein called FK506-binding protein 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	116	Total	C	N	O	S	0	1	0
			814	519	133	161	1			
3	F	114	Total	C	N	O	S	0	1	0
			820	522	137	160	1			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	-2	GLY	-	expression tag	UNP P28870
C	-1	SER	-	expression tag	UNP P28870
C	0	HIS	-	expression tag	UNP P28870
F	-2	GLY	-	expression tag	UNP P28870
F	-1	SER	-	expression tag	UNP P28870
F	0	HIS	-	expression tag	UNP P28870

- Molecule 4 is PHOSPHATE ION (three-letter code: PO4) (formula: O_4P).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total O P 5 4 1	0	0
4	D	1	Total O P 5 4 1	0	0
4	D	1	Total O P 5 4 1	0	0

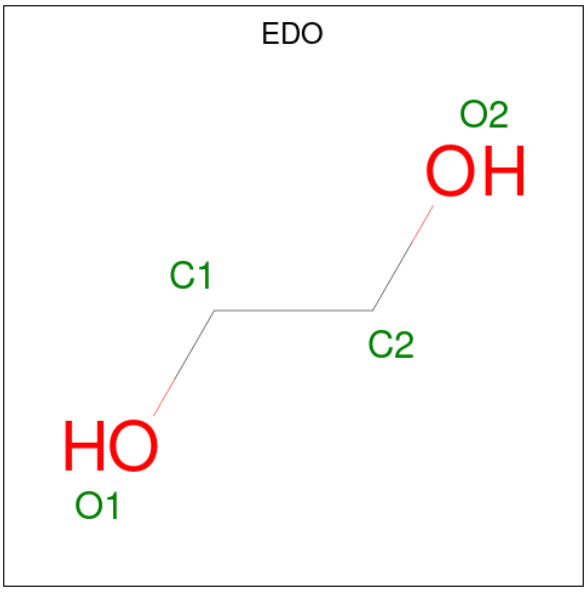
- Molecule 5 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total Zn 1 1	0	0
5	D	1	Total Zn 1 1	0	0

- Molecule 6 is FE (III) ION (three-letter code: FE) (formula: Fe).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	1	Total Fe 1 1	0	0
6	D	1	Total Fe 1 1	0	0

- Molecule 7 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: $C_2H_6O_2$).

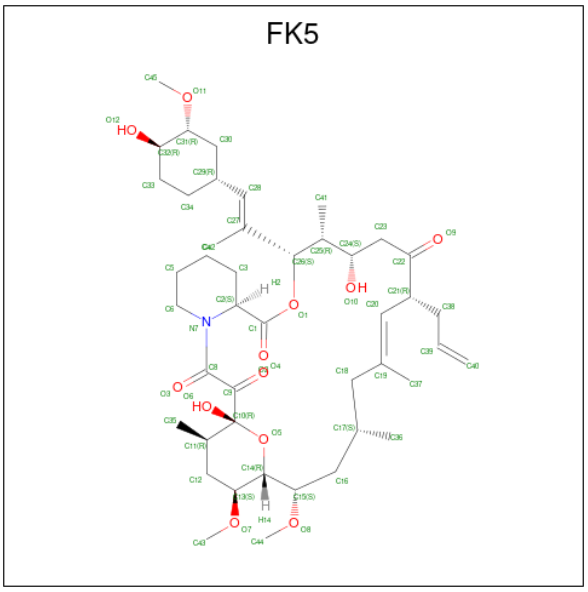


Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	1	Total C O 4 2 2	0	0
7	D	1	Total C O 4 2 2	0	0

- Molecule 8 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	B	3	Total Ca 3 3	0	0
8	E	4	Total Ca 4 4	0	0

- Molecule 9 is 8-DEETHYL-8-[BUT-3-ENYL]-ASCOMYCIN (three-letter code: FK5) (formula: $C_{44}H_{69}NO_{12}$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
9	C	1	Total C N O 57 44 1 12	0	0
9	F	1	Total C N O 57 44 1 12	0	0

- Molecule 10 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
10	A	165	Total O 165 165	0	0
10	B	7	Total O 7 7	0	0
10	C	5	Total O 5 5	0	0
10	D	92	Total O 92 92	0	0
10	E	4	Total O 4 4	0	0
10	F	8	Total O 8 8	0	0

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

3 Data and refinement statistics (i)

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

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	62.47Å 142.85Å 175.61Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	35.00 – 2.55	Depositor
% Data completeness (in resolution range)	98.5 (35.00-2.55)	Depositor
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	3.41 (at 2.54Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.182 , 0.229	Depositor
Wilson B-factor (Å^2)	36.0	Xtriage
Anisotropy	0.479	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	9858	wwPDB-VP
Average B, all atoms (Å^2)	52.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.79% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

4 Model quality

4.1 Standard geometry

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

4.2 Too-close contacts

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

4.3 Torsion angles

4.3.1 Protein backbone

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

4.3.2 Protein sidechains

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

4.3.3 RNA

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

4.4 Non-standard residues in protein, DNA, RNA chains

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

4.5 Carbohydrates

There are no monosaccharides in this entry.

4.6 Ligand geometry

Of 18 ligands modelled in this entry, 11 are monoatomic - leaving 7 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	EDO	D	505	-	3,3,3	0.60	0	2,2,2	0.18	0
9	FK5	F	200	-	54,60,60	3.44	10 (18%)	64,86,86	1.64	12 (18%)
4	PO4	A	501	6,5	4,4,4	0.76	0	6,6,6	0.72	0
4	PO4	D	504	-	4,4,4	0.85	0	6,6,6	0.51	0
9	FK5	C	200	-	54,60,60	3.45	11 (20%)	64,86,86	1.70	11 (17%)
4	PO4	D	501	6,5	4,4,4	0.70	0	6,6,6	1.15	0
7	EDO	A	504	-	3,3,3	0.43	0	2,2,2	0.47	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
7	EDO	D	505	-	-	1/1/1/1	-
9	FK5	F	200	-	-	5/68/114/114	0/3/4/4
7	EDO	A	504	-	-	0/1/1/1	-
9	FK5	C	200	-	-	4/68/114/114	0/3/4/4

All (21) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	F	200	FK5	O9-C22	11.33	1.41	1.21
9	C	200	FK5	O9-C22	11.10	1.40	1.21
9	C	200	FK5	C8-C9	-10.85	1.39	1.53
9	F	200	FK5	O4-C9	10.56	1.41	1.22
9	C	200	FK5	O4-C9	10.40	1.40	1.22
9	F	200	FK5	C8-C9	-9.73	1.41	1.53
9	C	200	FK5	C20-C19	9.09	1.47	1.33
9	F	200	FK5	C20-C19	8.82	1.46	1.33
9	F	200	FK5	O3-C8	8.77	1.41	1.23
9	C	200	FK5	O3-C8	7.86	1.39	1.23
9	F	200	FK5	O2-C1	7.75	1.40	1.21
9	C	200	FK5	O2-C1	7.44	1.39	1.21
9	C	200	FK5	C28-C27	6.33	1.46	1.33
9	F	200	FK5	C28-C27	5.98	1.46	1.33
9	F	200	FK5	O1-C1	-3.50	1.26	1.34
9	F	200	FK5	O6-C10	3.50	1.45	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	C	200	FK5	O1-C1	-3.42	1.26	1.34
9	C	200	FK5	O6-C10	3.30	1.45	1.39
9	F	200	FK5	C40-C39	2.79	1.47	1.28
9	C	200	FK5	C40-C39	2.79	1.47	1.28
9	C	200	FK5	O5-C14	-2.79	1.39	1.44

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
9	F	200	FK5	C21-C20-C19	-5.72	121.78	127.64
9	C	200	FK5	O1-C1-C2	5.13	122.08	110.78
9	F	200	FK5	O1-C1-C2	4.77	121.29	110.78
9	C	200	FK5	C21-C20-C19	-4.62	122.90	127.64
9	C	200	FK5	O1-C26-C25	4.17	110.69	105.91
9	C	200	FK5	O9-C22-C21	-3.81	117.20	121.13
9	C	200	FK5	C9-C8-N7	3.79	123.58	119.25
9	F	200	FK5	C42-C27-C26	3.15	121.12	115.68
9	C	200	FK5	O2-C1-C2	-2.91	117.94	124.49
9	C	200	FK5	C42-C27-C26	2.86	120.63	115.68
9	F	200	FK5	O9-C22-C21	-2.74	118.31	121.13
9	F	200	FK5	O2-C1-C2	-2.68	118.45	124.49
9	C	200	FK5	O3-C8-N7	-2.65	116.54	122.15
9	F	200	FK5	C23-C22-C21	2.55	121.02	117.75
9	F	200	FK5	C12-C11-C10	2.42	112.87	110.47
9	C	200	FK5	C5-C6-N7	-2.35	106.98	110.67
9	F	200	FK5	O3-C8-C9	2.27	120.02	116.28
9	F	200	FK5	C9-C8-N7	2.27	121.84	119.25
9	F	200	FK5	O1-C26-C25	2.22	108.45	105.91
9	F	200	FK5	O8-C15-C16	2.21	114.66	109.29
9	C	200	FK5	O3-C8-C9	2.19	119.87	116.28
9	C	200	FK5	O1-C1-O2	-2.12	119.97	123.94
9	F	200	FK5	O5-C14-C13	-2.06	106.39	109.42

There are no chirality outliers.

All (10) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
9	F	200	FK5	C20-C21-C38-C39
9	F	200	FK5	C22-C21-C38-C39
9	F	200	FK5	C21-C38-C39-C40
7	D	505	EDO	O1-C1-C2-O2
9	C	200	FK5	C36-C17-C18-C19

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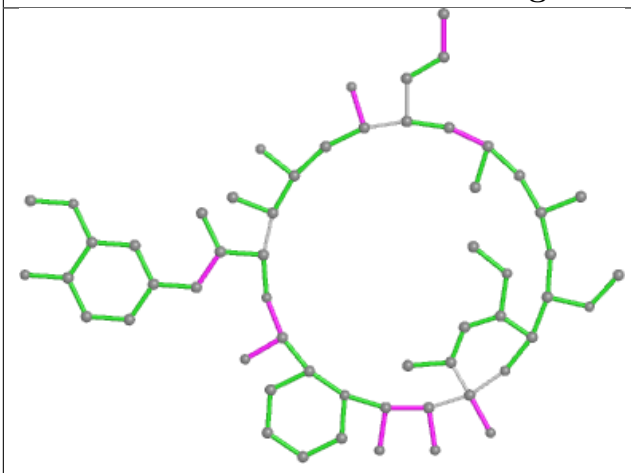
Mol	Chain	Res	Type	Atoms
9	C	200	FK5	C19-C20-C21-C22
9	F	200	FK5	C19-C20-C21-C22
9	C	200	FK5	C21-C38-C39-C40
9	C	200	FK5	C15-C16-C17-C36
9	F	200	FK5	C13-C14-C15-O8

There are no ring outliers.

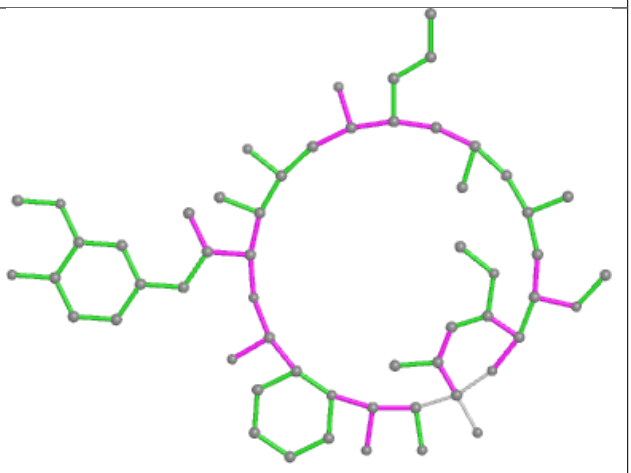
No monomer is involved in short contacts.

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 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.

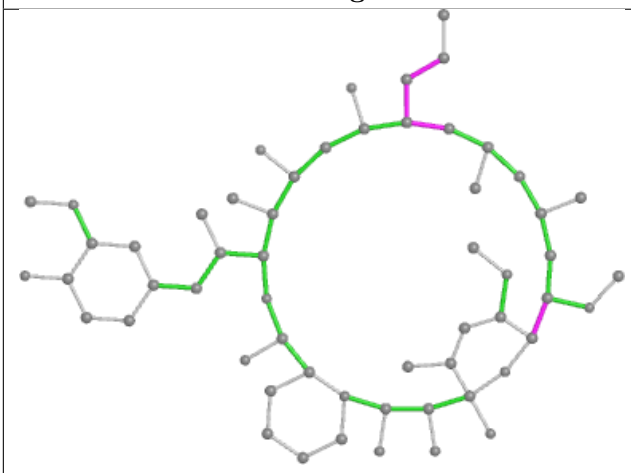
Ligand FK5 F 200



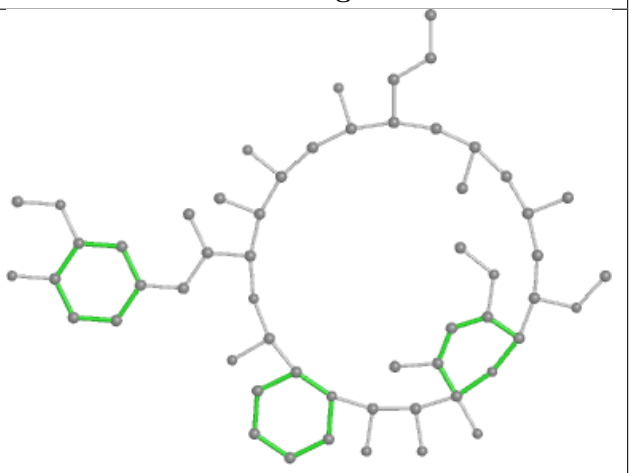
Bond lengths



Bond angles

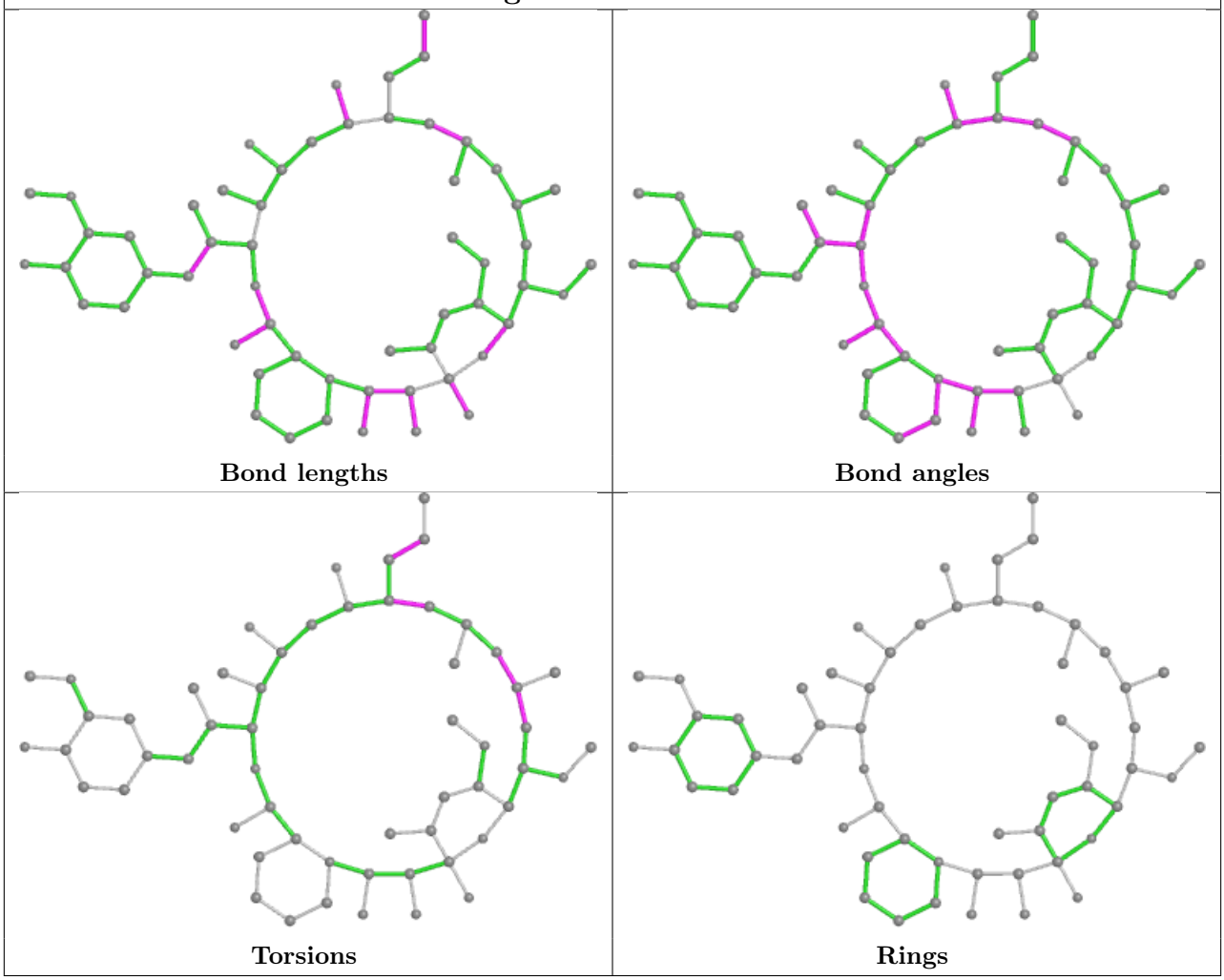


Torsions



Rings

Ligand FK5 C 200



4.7 Other polymers (i)

There are no such residues in this entry.

4.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

5 Fit of model and data

5.1 Protein, DNA and RNA chains

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

5.2 Non-standard residues in protein, DNA, RNA chains

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

5.3 Carbohydrates

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

5.4 Ligands

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

5.5 Other polymers

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