

Full wwPDB X-ray Structure Validation Report (i)

Oct 15, 2023 – 01:57 AM EDT

PDB ID	:	8ER6
Title	:	FKBP12-FRB in Complex with Compound 11
Authors	:	Tomlinson, A.C.A.; Yano, J.K.
Deposited on		
Resolution	:	2.81 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:

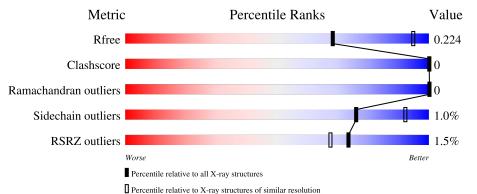
Xtriage (Phenix) EDS buster-report Percentile statistics Refmac CCP4 Ideal geometry (proteins) Ideal geometry (DNA, RNA)	:::::::::::::::::::::::::::::::::::::::	20191225.v01 (using entries in the PDB archive December 25th 2019) 5.8.0158 7.0.044 (Gargrove) Engh & Huber (2001) Parkinson et al. (1996)
Ideal geometry (DNA, RNA) Validation Pipeline (wwPDB-VP)		Parkinson et al. (1996) 2.36

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: $X\text{-}RAY \, DIFFRACTION$

The reported resolution of this entry is 2.81 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	$\begin{array}{c} \textbf{Whole archive} \\ (\#\textbf{Entries}) \end{array}$	${f Similar\ resolution}\ (\#{ m Entries,\ resolution\ range}({ m \AA}))$
R_{free}	130704	3617 (2.84 - 2.80)
Clashscore	141614	4060 (2.84-2.80)
Ramachandran outliers	138981	3978 (2.84-2.80)
Sidechain outliers	138945	3980 (2.84-2.80)
RSRZ outliers	127900	3552 (2.84-2.80)

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% The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	А	107	4% 99%
1	С	107	98% •
1	Е	107	<u>5%</u> 99%
2	В	95	99%
2	D	95	98%

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Mol	Chain	Length	Quality of chain
2	F	95	97% .



2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 5146 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.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
1	А	107	Total	С	Ν	0	\mathbf{S}	15	0	0
	Л	107	832	527	146	155	4	10		
1	C	107	Total	С	Ν	0	S	6	0	0
	U	107	832	527	146	155	4			
1	F	107	Total	С	Ν	0	S	15	0	0
	Ľ	107	832	527	146	155	4	10	0	U

• Molecule 1 is a protein called Peptidyl-prolyl cis-trans isomerase FKBP1A.

• Molecule 2 is a protein called non-specific serine/threenine protein kinase.

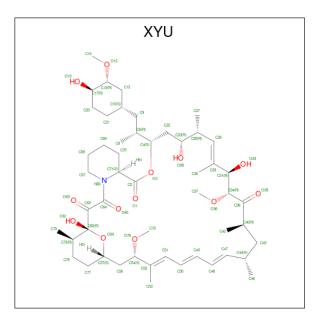
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
0	Р	95	Total	С	Ν	Ο	S	0	1	0
	D	90	809	515	142	145	$\overline{7}$	0	1	0
2	р	95	Total	С	Ν	0	S	8	0	0
	D	95	800	510	140	143	$\overline{7}$			
0	F	95	Total	С	Ν	0	S	0	9	0
	Г	95	817	520	143	146	8			0

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
В	2018	GLY	-	expression tag	UNP B1AKP8
D	2018	GLY	-	expression tag	UNP B1AKP8
F	2018	GLY	-	expression tag	UNP B1AKP8

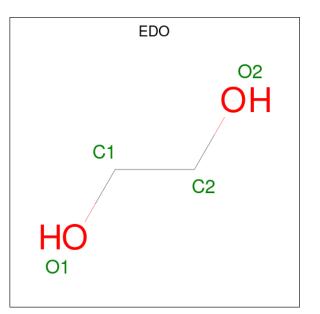
• Molecule 3 is (3S,5R,6R,7E,9R,10R,12R,14S,15E,17E,19E,21S,23S,26R,27R,30R,34aS)-5,9,27-trihydroxy-3-{ $(2R)-1-[(1S,3R,4R)-4-hydroxy-3-methoxycyclohexyl]propan-2-yl}-10,21-dim ethoxy-6,8,12,14,20,26-hexamethyl-5,6,9,10,12,13,14,21,22,23,24,25,26,27,32,33,34,34a-octa decahydro-3H-23,27-epoxypyrido[2,1-c][1,4]oxazacyclohentriacontine-1,11,28,29(4H,31H)-t etrone (three-letter code: XYU) (formula: <math>C_{51}H_{81}NO_{13}$) (labeled as "Ligand of Interest" by depositor).





Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	А	1	Total C N O 65 51 1 13	0	0
3	С	1	Total C N O 65 51 1 13	0	0
3	Е	1	Total C N O 65 51 1 13	0	0

• Molecule 4 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: $C_2H_6O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	D	1	Total 4	$\begin{array}{c} \mathrm{C} \\ \mathrm{2} \end{array}$	O 2	0	0



• Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	А	2	Total O 2 2	0	0
5	В	8	Total O 8 8	0	0
5	С	3	Total O 3 3	0	0
5	D	4	Total O 4 4	0	0
5	Ε	2	Total O 2 2	0	0
5	F	6	Total O 6 6	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 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: Peptidyl-prolyl cis-trans isomerase FKBP1A

Chain A:
T15 T22 E101 E103
\bullet Molecule 1: Peptidyl-prolyl cis-trans isomerase FKBP1A
Chain C: 98% ·
10 11 11 11 11 11 11 11 11 11 11 11 11 1
\bullet Molecule 1: Peptidyl-prolyl cis-trans isomerase FKBP1A
Chain E: 99%
C C C C C C C C C C C C C C C C C C C
\bullet Molecule 2: non-specific serine/three onine protein kinase
Chain B: 99%
C2 018 R2 036 S2 112 S2 112
\bullet Molecule 2: non-specific serine/three onine protein kinase
Chain D: 98% ·
R2018 82018 22112 22112

• Molecule 2: non-specific serine/threenine protein kinase



•

Chain F:

97%





4 Data and refinement statistics (i)

Property	Value	Source	
Space group	P 64 2 2	Depositor	
Cell constants	125.31Å 125.31Å 252.85Å	Depositor	
a, b, c, α , β , γ	90.00° 90.00° 120.00°	Depositor	
Resolution (Å)	45.84 - 2.81	Depositor	
Resolution (A)	45.84 - 2.81	EDS	
% Data completeness	96.5 (45.84-2.81)	Depositor	
(in resolution range)	96.5(45.84-2.81)	EDS	
R _{merge}	(Not available)	Depositor	
R _{sym}	0.07	Depositor	
$< I/\sigma(I) > 1$	$2.62 (at 2.81 \text{\AA})$	Xtriage	
Refinement program	PHENIX 1.20.1_4487	Depositor	
D D.	0.201 , 0.224	Depositor	
R, R_{free}	0.204 , 0.224	DCC	
R_{free} test set	669 reflections $(2.36%)$	wwPDB-VP	
Wilson B-factor $(Å^2)$	67.7	Xtriage	
Anisotropy	0.399	Xtriage	
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.33 , 36.5	EDS	
L-test for twinning ²	$ \langle L \rangle = 0.48, \langle L^2 \rangle = 0.30$	Xtriage	
Estimated twinning fraction	No twinning to report.	Xtriage	
F_o, F_c correlation	0.95	EDS	
Total number of atoms	5146	wwPDB-VP	
Average B, all atoms $(Å^2)$	74.0	wwPDB-VP	

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 3.90% of the height of the origin peak. No significant pseudotranslation is detected.

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.



¹Intensities estimated from amplitudes.

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: XYU, EDO

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	Bo	nd lengths	Bond angles		
IVIOI			$RMSZ \qquad \# Z > 5$		# Z > 5	
1	А	0.58	0/851	0.78	0/1146	
1	С	0.57	0/851	0.80	1/1146~(0.1%)	
1	Е	0.58	0/851	0.80	0/1146	
2	В	0.60	0/831	0.80	1/1116~(0.1%)	
2	D	0.69	1/822~(0.1%)	0.80	0/1104	
2	F	0.61	0/839	0.77	2/1126~(0.2%)	
All	All	0.60	1/5045~(0.0%)	0.79	4/6784~(0.1%)	

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	2076	ARG	NE-CZ	10.62	1.46	1.33

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Ζ	$\mathbf{Observed}(^{o})$	$Ideal(^{o})$
2	В	2036	ARG	NE-CZ-NH1	6.07	123.33	120.30
2	F	2036	ARG	NE-CZ-NH1	6.03	123.32	120.30
1	С	14	ARG	NE-CZ-NH1	5.06	122.83	120.30
2	F	2109	ARG	NE-CZ-NH1	5.05	122.82	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



Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	А	832	0	828	0	0
1	С	832	0	828	0	0
1	Е	832	0	828	0	0
2	В	809	0	770	0	0
2	D	800	0	763	0	0
2	F	817	0	778	0	0
3	А	65	0	0	0	0
3	С	65	0	0	0	0
3	Е	65	0	0	0	0
4	D	4	0	6	0	0
5	А	2	0	0	0	0
5	В	8	0	0	0	0
5	С	3	0	0	0	0
5	D	4	0	0	0	0
5	Е	2	0	0	0	0
5	F	6	0	0	0	0
All	All	5146	0	4801	0	0

the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

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 0.

There are no clashes within the asymmetric unit.

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	Perce	ntiles
1	А	105/107~(98%)	102 (97%)	3~(3%)	0	100	100
1	С	105/107~(98%)	101 (96%)	4 (4%)	0	100	100
1	Е	105/107~(98%)	102 (97%)	3~(3%)	0	100	100
2	В	94/95~(99%)	94 (100%)	0	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Perce	ntiles
2	D	93/95~(98%)	93 (100%)	0	0	100	100
2	F	95/95~(100%)	95 (100%)	0	0	100	100
All	All	597/606~(98%)	587~(98%)	10 (2%)	0	100	100

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There are no Ramachandran outliers to report.

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	Perce	ntiles
1	А	89/89~(100%)	88~(99%)	1 (1%)	73	91
1	С	89/89~(100%)	88~(99%)	1 (1%)	73	91
1	Е	89/89~(100%)	88 (99%)	1 (1%)	73	91
2	В	83/82~(101%)	83 (100%)	0	100	100
2	D	82/82 (100%)	81 (99%)	1 (1%)	71	91
2	F	84/82~(102%)	83~(99%)	1 (1%)	71	91
All	All	516/513~(101%)	511 (99%)	5 (1%)	76	92

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	А	107	LEU
1	С	19	ARG
2	D	2079	MET
1	Е	30	MET
2	F	2087	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.



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)

4 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	Mol Type		Chain Res	Res Link	Bo	Bond lengths			Bond angles		
	туре	Chain	nes		Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2	
3	XYU	С	201	-	64,68,68	1.12	6 (9%)	74,96,96	1.51	10 (13%)	
3	XYU	А	201	-	64,68,68	1.08	6 (9%)	74,96,96	1.54	12 (16%)	
3	XYU	Е	201	-	64,68,68	1.04	5 (7%)	74,96,96	1.66	12 (16%)	
4	EDO	D	2201	-	3,3,3	0.55	0	2,2,2	0.55	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	XYU	С	201	-	-	5/81/124/124	0/3/4/4
3	XYU	А	201	-	-	4/81/124/124	0/3/4/4
3	XYU	Е	201	-	-	4/81/124/124	0/3/4/4
4	EDO	D	2201	-	-	1/1/1/1	-



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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	С	201	XYU	C64-C62	-4.79	1.47	1.53
3	А	201	XYU	C64-C62	-4.50	1.47	1.53
3	Ε	201	XYU	C64-C62	-3.96	1.48	1.53
3	С	201	XYU	O61-C60	2.91	1.44	1.39
3	Ε	201	XYU	O61-C60	2.89	1.44	1.39
3	С	201	XYU	C6-C4	2.86	1.56	1.53
3	А	201	XYU	C71-N66	2.71	1.50	1.47
3	С	201	XYU	C67-N66	2.61	1.51	1.47
3	Ε	201	XYU	C71-N66	2.57	1.50	1.47
3	А	201	XYU	C67-N66	2.55	1.51	1.47
3	А	201	XYU	C6-C4	2.47	1.55	1.53
3	А	201	XYU	O61-C60	2.45	1.43	1.39
3	С	201	XYU	C71-N66	2.40	1.50	1.47
3	Ε	201	XYU	C67-N66	2.38	1.51	1.47
3	С	201	XYU	C51-C52	2.07	1.35	1.33
3	А	201	XYU	C51-C52	2.05	1.35	1.33
3	Е	201	XYU	C6-C4	2.02	1.55	1.53

All (17) bond length outliers are listed below:

All (34) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Ζ	$Observed(^{o})$	$Ideal(^{o})$
3	Е	201	XYU	C62-C64-N66	5.38	125.40	119.25
3	С	201	XYU	C62-C64-N66	4.93	124.89	119.25
3	А	201	XYU	C62-C64-N66	4.89	124.84	119.25
3	Ε	201	XYU	C4-O3-C2	4.72	124.98	117.89
3	Е	201	XYU	C56-C54-C52	-4.59	107.12	113.50
3	Е	201	XYU	O3-C2-C71	4.46	120.60	110.78
3	С	201	XYU	C4-O3-C2	4.35	124.42	117.89
3	А	201	XYU	O3-C2-C71	4.24	120.12	110.78
3	С	201	XYU	O3-C2-C71	4.09	119.79	110.78
3	Е	201	XYU	C8-C6-C9	-3.95	104.76	110.69
3	А	201	XYU	C8-C6-C9	-3.49	105.46	110.69
3	С	201	XYU	C8-C6-C9	-3.45	105.52	110.69
3	А	201	XYU	C4-O3-C2	3.29	122.83	117.89
3	А	201	XYU	O63-C62-C64	-3.16	115.53	119.08
3	С	201	XYU	O63-C62-C64	-3.13	115.56	119.08
3	С	201	XYU	O1-C2-C71	-3.08	117.55	124.49
3	А	201	XYU	O1-C2-C71	-3.08	117.55	124.49
3	Е	201	XYU	O1-C2-C71	-3.07	117.57	124.49
3	А	201	XYU	C56-C54-C52	-3.06	109.25	113.50
3	Е	201	XYU	O63-C62-C64	-2.87	115.85	119.08
3	А	201	XYU	C9-C10-C12	-2.63	106.02	111.75

PROTEIN DATA BANK

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Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
3	А	201	XYU	C48-C49-C50	-2.57	119.08	124.81
3	А	201	XYU	C53-C52-C51	-2.57	120.86	124.03
3	С	201	XYU	C48-C49-C50	-2.54	119.16	124.81
3	Ε	201	XYU	C9-C6-C4	2.52	117.04	111.51
3	Е	201	XYU	C48-C49-C50	-2.52	119.21	124.81
3	С	201	XYU	C42-C40-C38	2.45	115.33	109.44
3	С	201	XYU	C9-C10-C12	-2.43	106.45	111.75
3	А	201	XYU	C50-C51-C52	-2.39	123.80	127.32
3	С	201	XYU	C53-C52-C51	-2.39	121.07	124.03
3	Е	201	XYU	C9-C10-C12	-2.35	106.63	111.75
3	А	201	XYU	C2-C71-N66	2.19	116.13	112.05
3	Е	201	XYU	C50-C51-C52	-2.08	124.27	127.32
3	Е	201	XYU	C42-C40-C38	2.05	114.37	109.44

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There are no chirality outliers.

All (14) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	Е	201	XYU	C54-C56-C57-C77
3	Е	201	XYU	C54-C56-C57-O59
3	С	201	XYU	C47-C48-C49-C50
4	D	2201	EDO	O1-C1-C2-O2
3	С	201	XYU	C46-C44-C47-C48
3	С	201	XYU	C43-C44-C47-C48
3	А	201	XYU	C4-C22-C23-O80
3	С	201	XYU	C4-C22-C23-O80
3	Е	201	XYU	C4-C22-C23-O80
3	С	201	XYU	C4-C22-C23-C25
3	А	201	XYU	C29-C31-C34-C38
3	А	201	XYU	C4-C22-C23-C25
3	Е	201	XYU	C4-C22-C23-C25
3	А	201	XYU	O39-C38-C40-C43

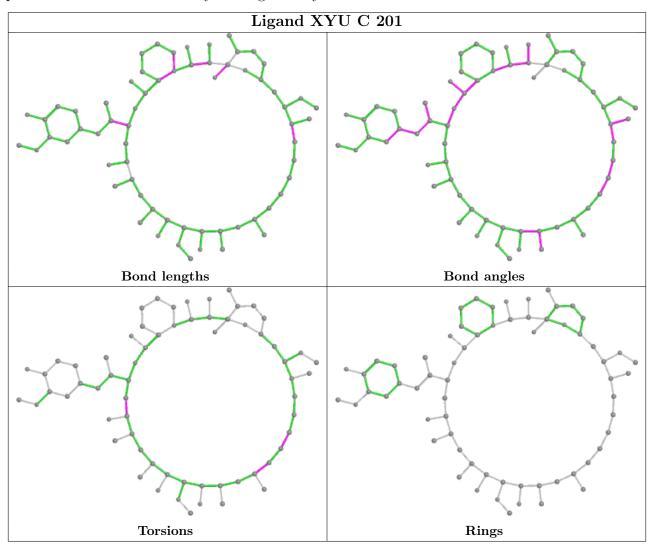
There are no ring outliers.

No monomer is involved in short contacts.

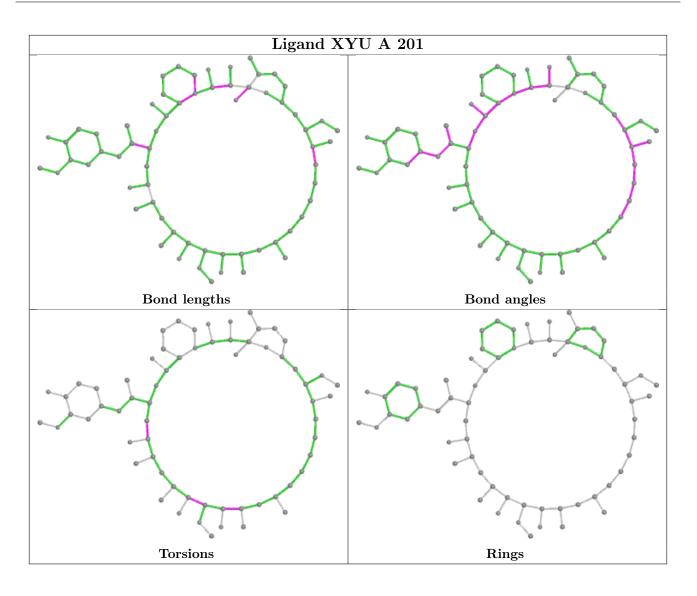
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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



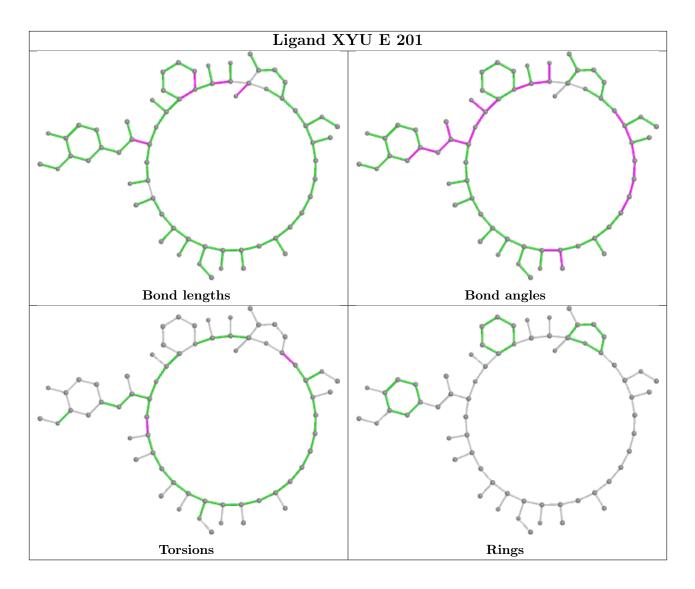
highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified 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.











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)

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^{th} 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 $>$	$\# RSRZ {>}2$	$\mathbf{OWAB}(\mathbf{A}^2)$	Q < 0.9
1	А	107/107~(100%)	0.33	4 (3%) 41 31	59, 93, 120, 137	5(4%)
1	С	107/107~(100%)	0.01	0 100 100	56, 67, 83, 114	3 (2%)
1	Ε	107/107~(100%)	0.44	5 (4%) 31 22	59, 86, 117, 132	5 (4%)
2	В	95/95~(100%)	0.03	0 100 100	46, 58, 70, 83	0
2	D	95/95~(100%)	0.12	0 100 100	59, 69, 85, 93	3~(3%)
2	F	95/95~(100%)	0.11	0 100 100	51, 68, 89, 98	0
All	All	606/606~(100%)	0.18	9 (1%) 73 67	46, 70, 112, 137	16 (2%)

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	А	16	PHE	5.3
1	Е	2	GLY	3.3
1	Е	33	ASP	2.7
1	А	15	THR	2.4
1	А	51	LEU	2.3
1	Е	99	VAL	2.2
1	Е	31	LEU	2.1
1	Е	3	VAL	2.1
1	А	22	THR	2.1

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 monosaccharides in this entry.

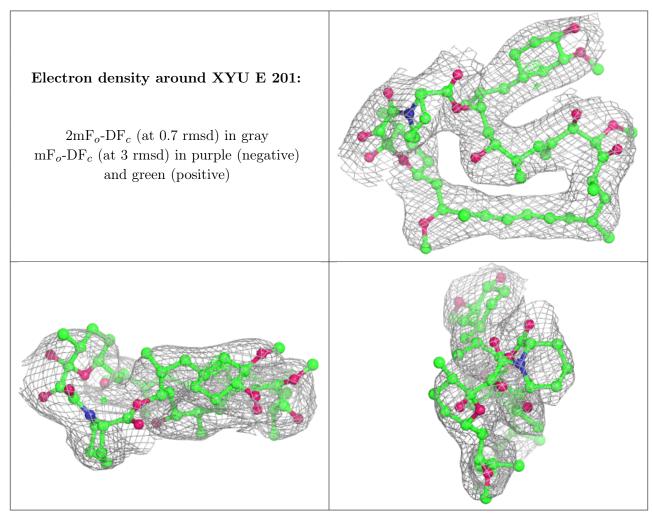


6.4 Ligands (i)

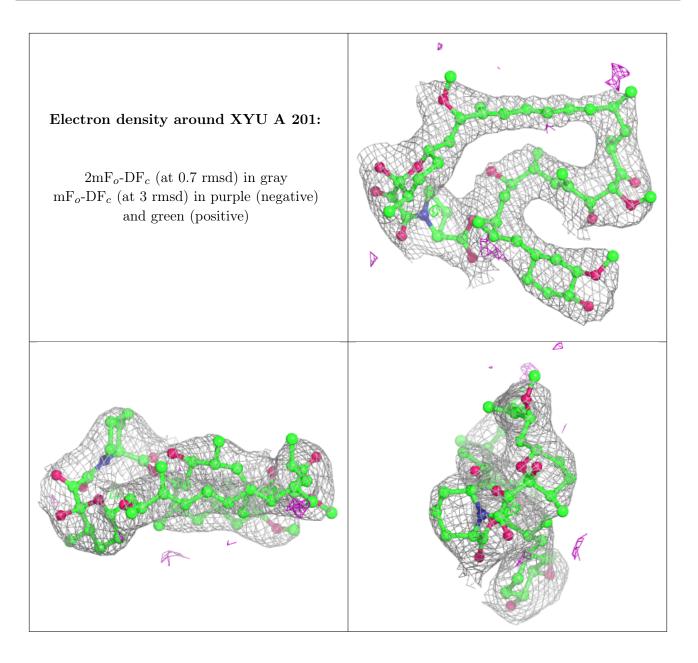
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	$\mathbf{B} ext{-factors}(\mathrm{\AA}^2)$	Q<0.9
4	EDO	D	2201	4/4	0.89	0.26	$57,\!59,\!65,\!73$	0
3	XYU	Е	201	65/65	0.96	0.25	58,72,93,111	0
3	XYU	А	201	65/65	0.97	0.26	52,67,77,78	0
3	XYU	С	201	65/65	0.97	0.23	45,57,63,72	0

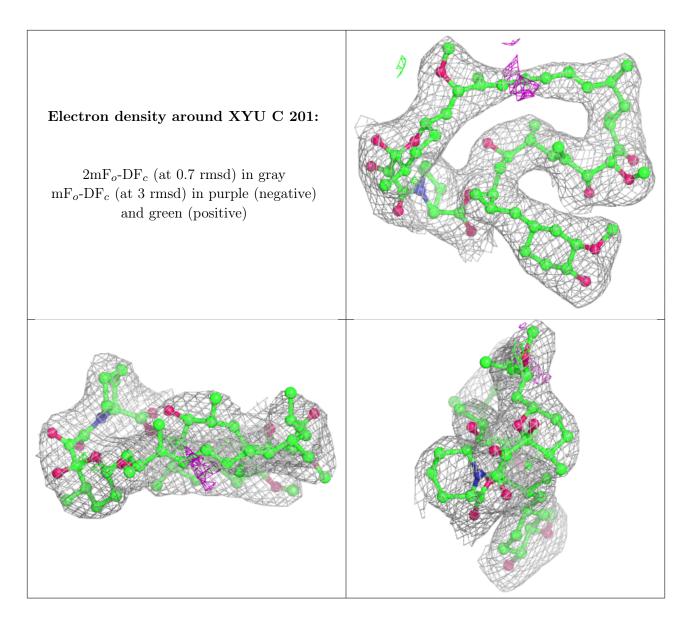
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.











6.5 Other polymers (i)

There are no such residues in this entry.

