



# wwPDB X-ray Structure Validation Summary Report ⓘ

Nov 9, 2023 – 03:56 pm GMT

PDB ID : 8PC2  
Title : SelDeg51 in complex with FKBP51FK1 domain and pVHL:EloB:EloC  
Authors : Meyners, C.; Walz, M.; Geiger, T.M.; Hausch, F.  
Deposited on : 2023-06-09  
Resolution : 2.80 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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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 : 1.8.4, CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.36  
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) : 2.36

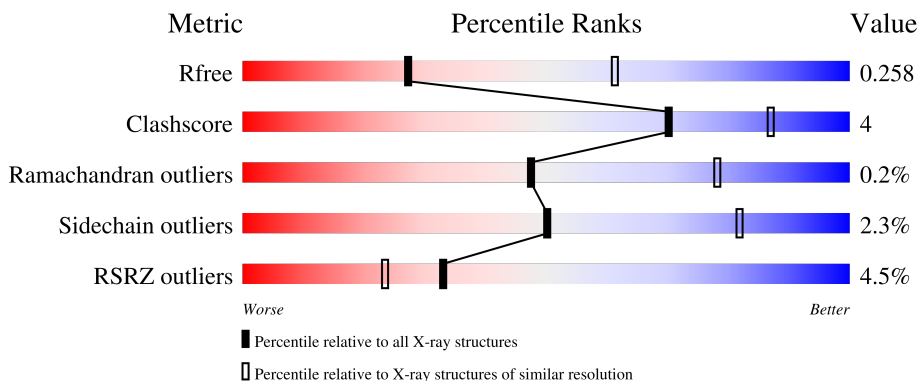
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.80 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.



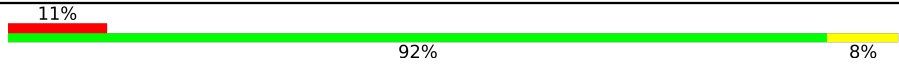


Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	130704	3140 (2.80-2.80)
Clashscore	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)
RSRZ outliers	127900	3078 (2.80-2.80)

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  $\geq 3$ , 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  $\leq 5\%$ . 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	A	162	80% 10% 8%
1	B	162	81% 10% 8%
2	C	97	77% 12% 10%
2	E	97	74% 14% 10%
3	D	104	91% 9%

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Mol	Chain	Length	Quality of chain
3	F	104	 11% 92% 8%
4	G	128	 7% 84% 9% 6%
4	H	128	 12% 83% 9% 8%

## 2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 14339 atoms, of which 7115 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 von Hippel-Lindau disease tumor suppressor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace	
			Total	C	H	N	O				S
1	B	149	2393	761	1198	220	212	2	38	0	0
1	A	149	2385	761	1188	217	217	2	37	0	0

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	52	GLY	-	expression tag	UNP P40337
B	53	SER	-	expression tag	UNP P40337
A	52	GLY	-	expression tag	UNP P40337
A	53	SER	-	expression tag	UNP P40337

- Molecule 2 is a protein called Elongin-C.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace	
			Total	C	H	N	O				S
2	C	87	1359	438	679	109	127	6	29	0	0
2	E	87	1368	441	685	108	129	5	28	0	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	16	MET	-	initiating methionine	UNP Q15369
E	16	MET	-	initiating methionine	UNP Q15369

- Molecule 3 is a protein called Elongin-B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace	
			Total	C	H	N	O				S
3	D	104	1590	506	794	132	153	5	26	0	0

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace	
			Total	C	H	N	O				S
3	F	104	1550	496	769	130	151	4	30	0	0

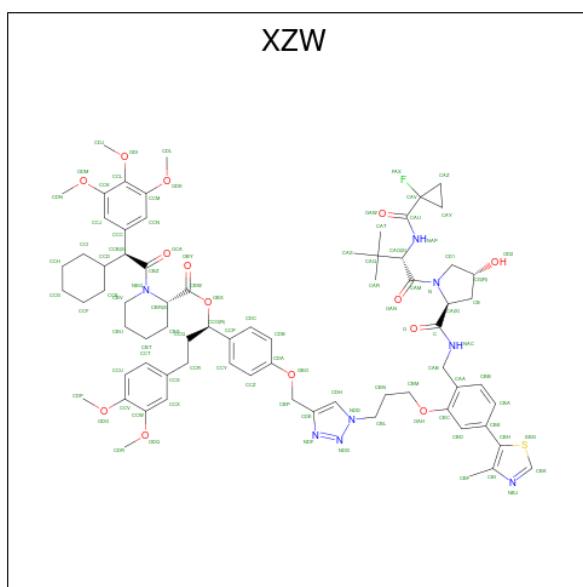
- Molecule 4 is a protein called Peptidyl-prolyl cis-trans isomerase FKBP5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace	
			Total	C	H	N	O				S
4	G	120	1721	561	850	143	165	2	38	0	0
4	H	118	1585	527	770	130	157	1	47	0	0

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
G	13	GLY	-	expression tag	UNP Q13451
G	14	ALA	-	expression tag	UNP Q13451
G	15	PRO	-	expression tag	UNP Q13451
G	19	THR	ALA	engineered mutation	UNP Q13451
G	103	ALA	CYS	engineered mutation	UNP Q13451
G	107	ILE	CYS	engineered mutation	UNP Q13451
H	13	GLY	-	expression tag	UNP Q13451
H	14	ALA	-	expression tag	UNP Q13451
H	15	PRO	-	expression tag	UNP Q13451
H	19	THR	ALA	engineered mutation	UNP Q13451
H	103	ALA	CYS	engineered mutation	UNP Q13451
H	107	ILE	CYS	engineered mutation	UNP Q13451

- Molecule 5 is [(1 {R})-3-(3,4-dimethoxyphenyl)-1-[4-[[1-[3-[2-[[[(2 {S}),4 {R})-1-[(2 {S})-2-[(1-fluoranyl)cyclopropyl]carbonylamino]-3,3-dimethyl-butanoyl]-4-oxidanyl-pyrrolidin-2-yl]carbonylamino]methyl]-5-(4-methyl-1,3-thiazol-5-yl)phenoxy]propyl]-1,2,3-triazol-4-yl]methoxy]phenyl]propyl] (2 {S})-1-[(2 {S})-2-cyclohexyl-2-(3,4,5-trimethoxyphenyl)ethanoyl]piperidine-2-carboxylate (three-letter code: XZW) (formula: C<sub>72</sub>H<sub>91</sub>FN<sub>8</sub>O<sub>14</sub>S) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
			Total	C	F	H	N	O		
5	B	1	187	72	1	91	8	14	1	0
5	A	1	187	72	1	91	8	14	1	0

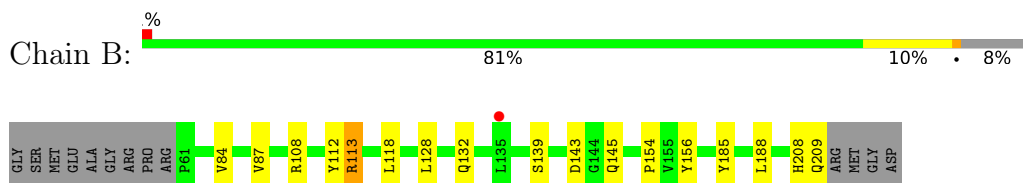
- Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	B	7	Total 7 7	0	0
6	G	3	Total 3 3	0	0
6	A	3	Total 3 3	0	0
6	E	1	Total 1 1	0	0

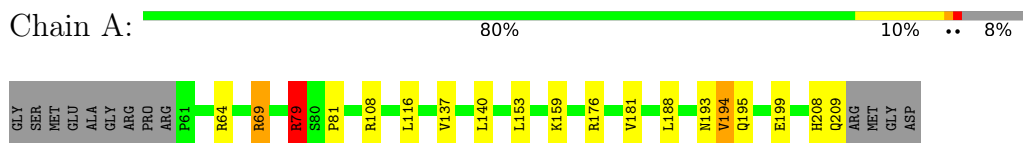
### 3 Residue-property plots [i](#)

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.

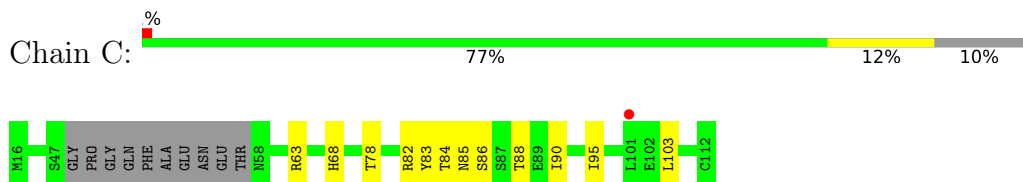
- Molecule 1: von Hippel-Lindau disease tumor suppressor



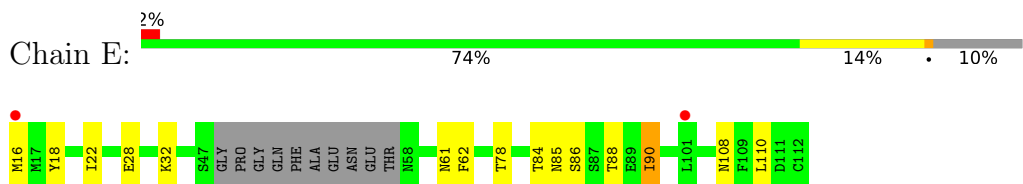
- Molecule 1: von Hippel-Lindau disease tumor suppressor



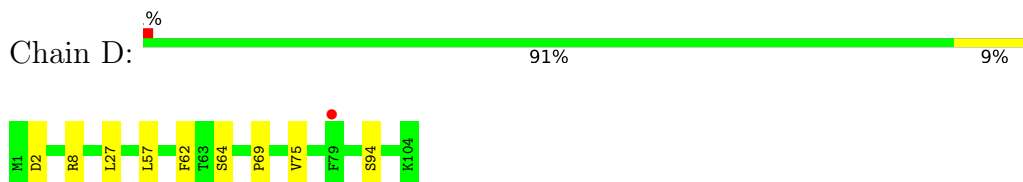
- Molecule 2: Elongin-C



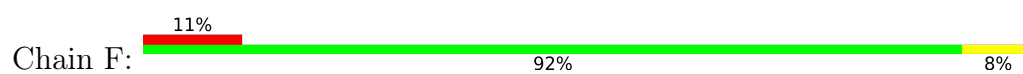
- Molecule 2: Elongin-C



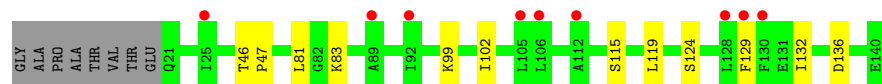
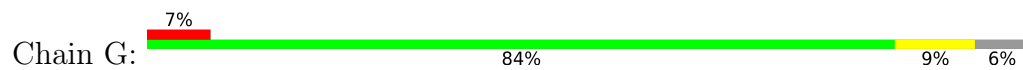
- Molecule 3: Elongin-B



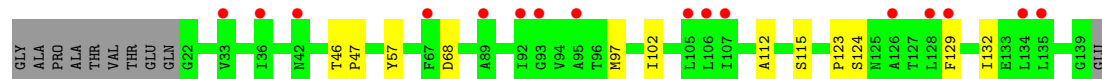
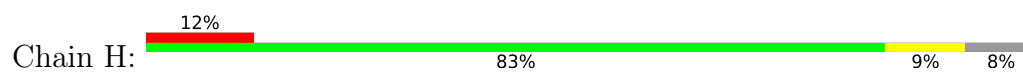
- Molecule 3: Elongin-B



- Molecule 4: Peptidyl-prolyl cis-trans isomerase FKBP5



- Molecule 4: Peptidyl-prolyl cis-trans isomerase FKBP5





## 4 Data and refinement statistics

Property	Value	Source
Space group	I 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	138.65Å 68.42Å 159.26Å 90.00° 114.45° 90.00°	Depositor
Resolution (Å)	48.74 – 2.80 48.69 – 2.80	Depositor EDS
% Data completeness (in resolution range)	98.9 (48.74-2.80) 98.4 (48.69-2.80)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.19 (at 2.81Å)	Xtrriage
Refinement program	REFMAC 5.8.0411	Depositor
R, $R_{free}$	0.210 , 0.261 0.212 , 0.258	Depositor DCC
$R_{free}$ test set	1664 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	47.6	Xtrriage
Anisotropy	0.536	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 44.0	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	14339	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	79.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.96% 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: XZW

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	0.41	0/1229	0.78	1/1682 (0.1%)
1	B	0.45	0/1227	0.78	0/1678
2	C	0.46	0/694	0.70	0/939
2	E	0.45	0/697	0.73	0/942
3	D	0.42	0/812	0.79	0/1100
3	F	0.39	0/797	0.75	0/1084
4	G	0.40	0/890	0.74	1/1206 (0.1%)
4	H	0.34	0/834	0.66	0/1141
All	All	0.41	0/7180	0.75	2/9772 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	4
1	B	0	2
2	C	0	1
2	E	0	1
3	D	0	1
3	F	0	1
All	All	0	10

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	G	83	LYS	CB-CA-C	5.64	121.67	110.40
1	A	79	ARG	CG-CD-NE	5.03	122.36	111.80

There are no chirality outliers.

5 of 10 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	69	ARG	Sidechain
1	B	108	ARG	Sidechain
1	B	113	ARG	Sidechain
2	C	82	ARG	Sidechain
3	D	8	ARG	Sidechain

## 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	1197	1188	1172	9	0
1	B	1195	1198	1180	8	0
2	C	680	679	664	8	0
2	E	683	685	672	7	0
3	D	796	794	779	5	0
3	F	781	769	746	5	0
4	G	871	850	813	7	0
4	H	815	770	715	6	0
5	A	96	91	0	0	0
5	B	96	91	0	2	0
6	A	3	0	0	0	0
6	B	7	0	0	0	0
6	E	1	0	0	0	0
6	G	3	0	0	0	0
All	All	7224	7115	6741	50	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 4.

The worst 5 of 50 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)
2:E:86:SER:HB2	2:E:90:ILE:HD11	1.69	0.73
2:C:78:THR:HG22	3:D:69:PRO:HB3	1.80	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:84:THR:O	2:E:85:ASN:HB2	2.00	0.62
1:B:87:VAL:HB	1:B:118:LEU:HD12	1.83	0.61
2:C:86:SER:HB2	2:C:90:ILE:HD11	1.82	0.61

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	147/162 (91%)	140 (95%)	6 (4%)	1 (1%)	22	53
1	B	147/162 (91%)	142 (97%)	5 (3%)	0	100	100
2	C	83/97 (86%)	79 (95%)	4 (5%)	0	100	100
2	E	83/97 (86%)	80 (96%)	3 (4%)	0	100	100
3	D	102/104 (98%)	97 (95%)	5 (5%)	0	100	100
3	F	102/104 (98%)	97 (95%)	5 (5%)	0	100	100
4	G	118/128 (92%)	114 (97%)	4 (3%)	0	100	100
4	H	116/128 (91%)	113 (97%)	2 (2%)	1 (1%)	17	46
All	All	898/982 (91%)	862 (96%)	34 (4%)	2 (0%)	47	78

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	194	VAL
4	H	112	ALA

### 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	134/148 (90%)	129 (96%)	5 (4%)	34	68
1	B	133/148 (90%)	128 (96%)	5 (4%)	33	67
2	C	75/86 (87%)	74 (99%)	1 (1%)	69	91
2	E	76/86 (88%)	73 (96%)	3 (4%)	32	66
3	D	85/92 (92%)	85 (100%)	0	100	100
3	F	81/92 (88%)	79 (98%)	2 (2%)	47	80
4	G	83/105 (79%)	83 (100%)	0	100	100
4	H	72/105 (69%)	71 (99%)	1 (1%)	67	90
All	All	739/862 (86%)	722 (98%)	17 (2%)	50	82

5 of 17 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	F	22	SER
4	H	46	THR
1	A	79	ARG
1	A	140	LEU
1	A	181	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	209	GLN
1	A	209	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 monosaccharides 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
5	XZW	B	301	-	97,105,105	2.54	16 (16%)	127,150,150	1.92	34 (26%)
5	XZW	A	301	-	97,105,105	2.18	14 (14%)	127,150,150	1.90	31 (24%)

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
5	XZW	B	301	-	-	10/88/127/127	0/10/10/10
5	XZW	A	301	-	-	8/88/127/127	0/10/10/10

The worst 5 of 30 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	301	XZW	CBE-CBH	-12.15	1.36	1.48
5	B	301	XZW	CBE-CBH	-10.95	1.37	1.48
5	B	301	XZW	NDG-NDD	-10.46	1.14	1.34
5	B	301	XZW	CCC-CCB	-8.95	1.39	1.52
5	A	301	XZW	CCP-CCO	-8.56	1.36	1.51

The worst 5 of 65 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	301	XZW	CBL-NDD-CDH	-8.49	109.86	129.82
5	A	301	XZW	OBX-CCO-CCQ	8.19	117.06	105.34
5	A	301	XZW	NDF-NDG-NDD	5.68	111.60	107.31
5	B	301	XZW	ODQ-CCW-CCV	5.68	123.33	115.41
5	B	301	XZW	CCE-CCD-CCB	-5.39	104.75	111.91

There are no chirality outliers.

5 of 18 torsion outliers are listed below:

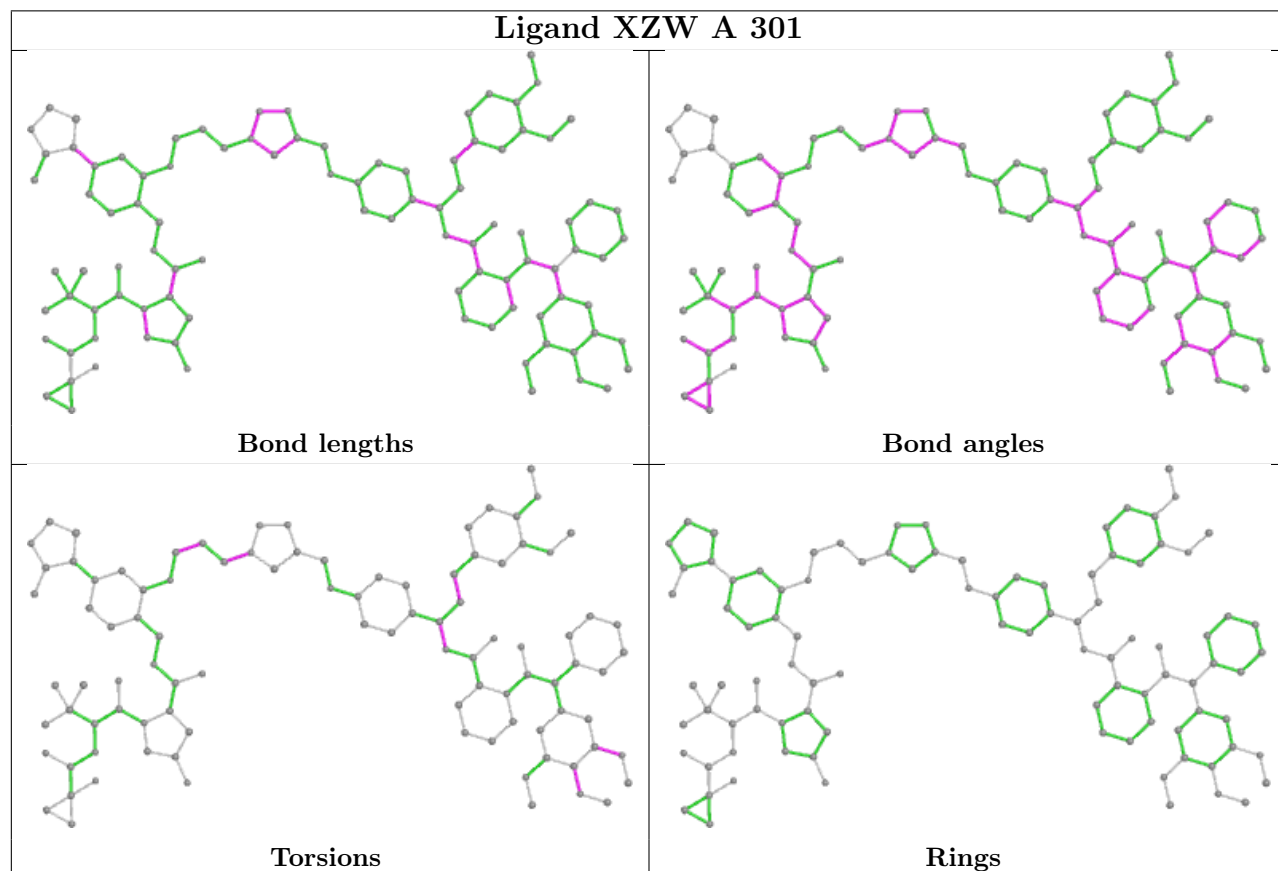
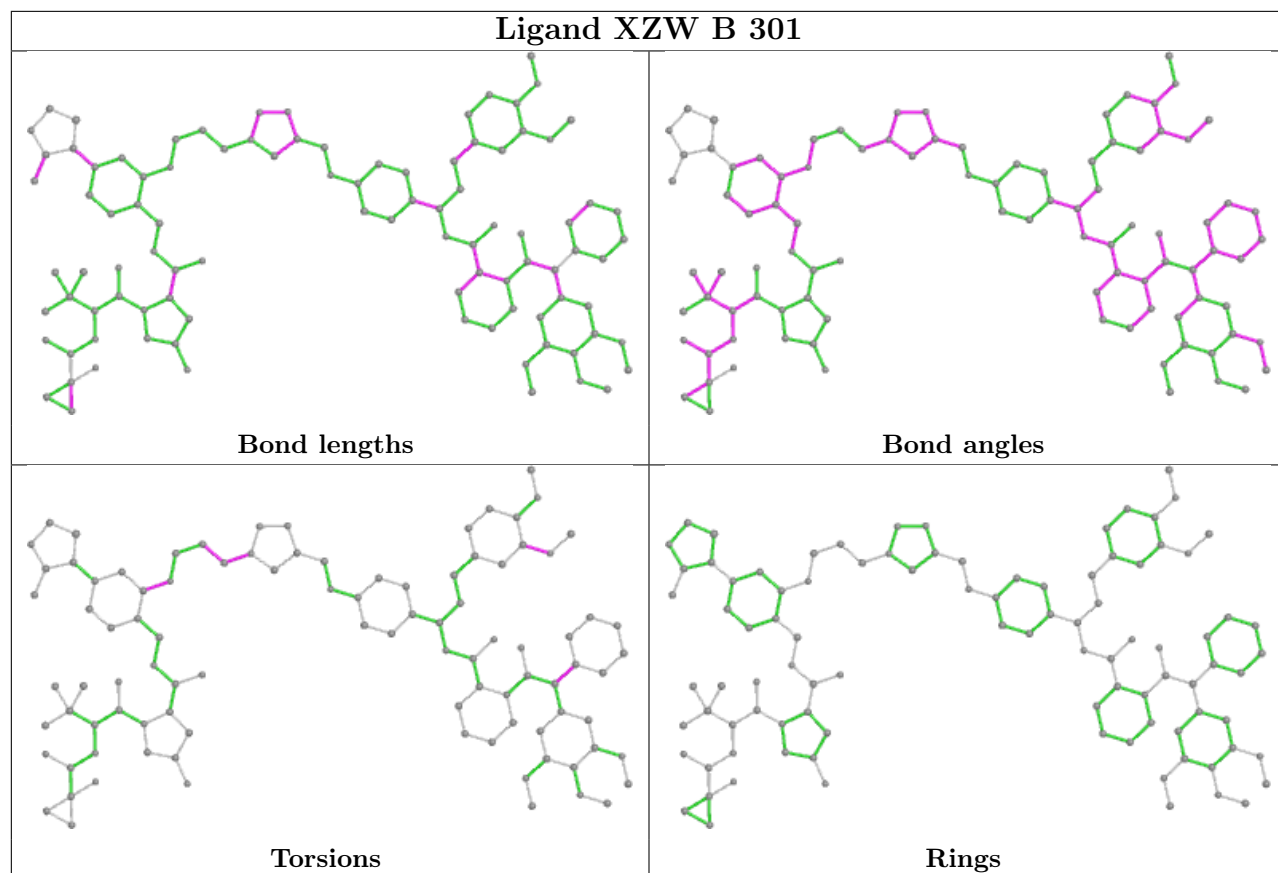
Mol	Chain	Res	Type	Atoms
5	B	301	XZW	CBN-CBL-NDD-CDH
5	B	301	XZW	CBZ-CCB-CCD-CCE
5	B	301	XZW	CCC-CCB-CCD-CCE
5	B	301	XZW	CCC-CCB-CCD-CCI
5	B	301	XZW	CCX-CCW-ODQ-CDR

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	301	XZW	2	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](#)

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	149/162 (91%)	0.20	0 <span style="border: 1px solid blue;">100</span> <span style="border: 1px solid blue;">100</span>	51, 71, 101, 131	0
1	B	149/162 (91%)	0.20	1 (0%) <span style="border: 1px solid blue;">87</span> <span style="border: 1px solid blue;">84</span>	47, 63, 106, 138	0
2	C	87/97 (89%)	0.25	1 (1%) <span style="border: 1px solid blue;">80</span> <span style="border: 1px solid blue;">75</span>	50, 68, 102, 110	0
2	E	87/97 (89%)	0.21	2 (2%) <span style="border: 1px solid blue;">60</span> <span style="border: 1px solid blue;">51</span>	58, 77, 102, 117	0
3	D	104/104 (100%)	0.10	1 (0%) <span style="border: 1px solid blue;">82</span> <span style="border: 1px solid blue;">77</span>	54, 73, 101, 123	0
3	F	104/104 (100%)	0.47	11 (10%) <span style="border: 1px solid red;">6</span> <span style="border: 1px solid red;">3</span>	60, 87, 112, 151	0
4	G	120/128 (93%)	0.22	9 (7%) <span style="border: 1px solid red;">14</span> <span style="border: 1px solid red;">8</span>	55, 76, 102, 115	0
4	H	118/128 (92%)	0.47	16 (13%) <span style="border: 1px solid red;">3</span> <span style="border: 1px solid red;">1</span>	79, 120, 153, 172	0
All	All	918/982 (93%)	0.26	41 (4%) <span style="border: 1px solid red;">33</span> <span style="border: 1px solid red;">23</span>	47, 76, 130, 172	0

The worst 5 of 41 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
4	H	128	LEU	3.9
4	H	135	LEU	3.9
3	F	88	LEU	3.8
4	H	107	ILE	3.6
4	H	105	LEU	3.5

### 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](#)

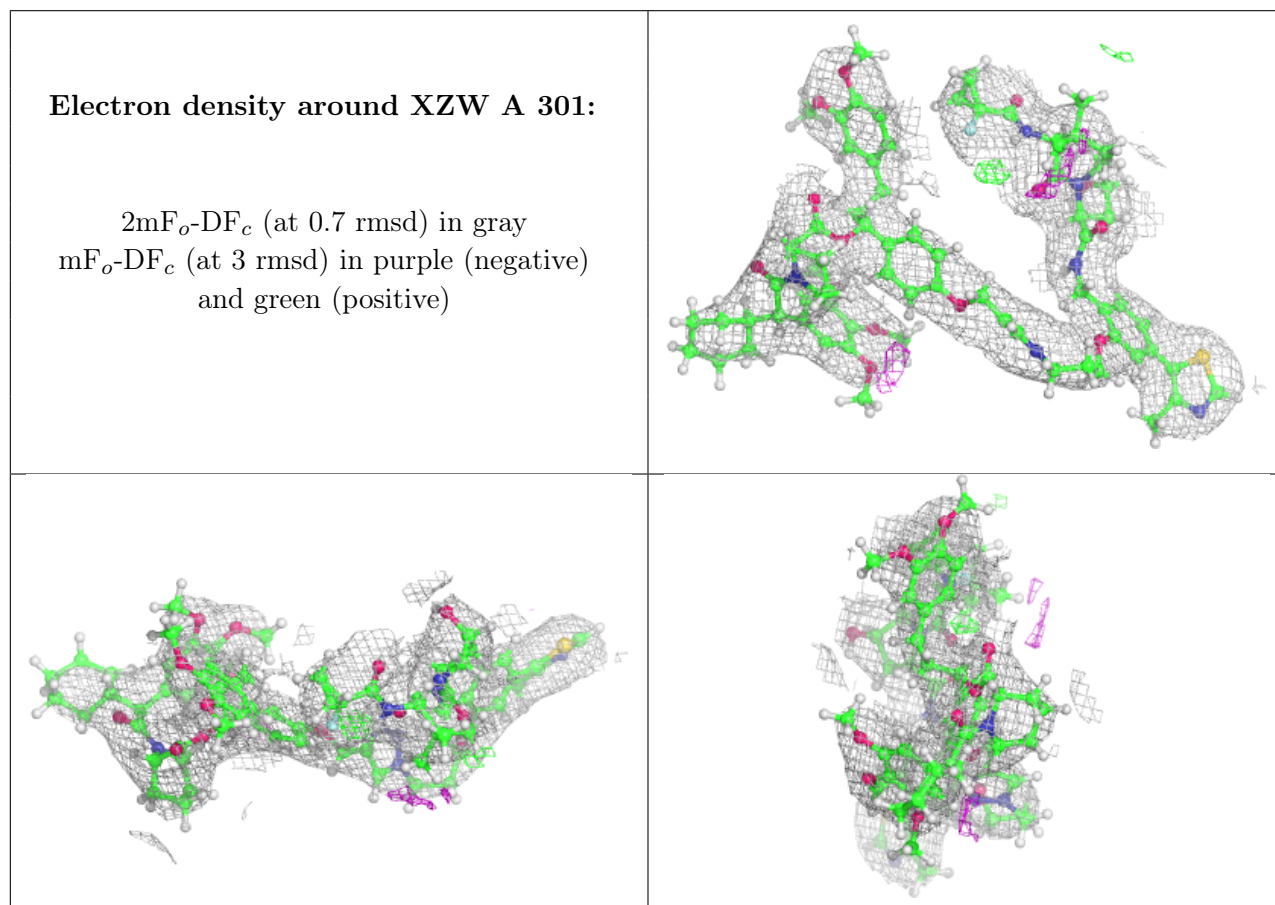
There are no monosaccharides in this entry.

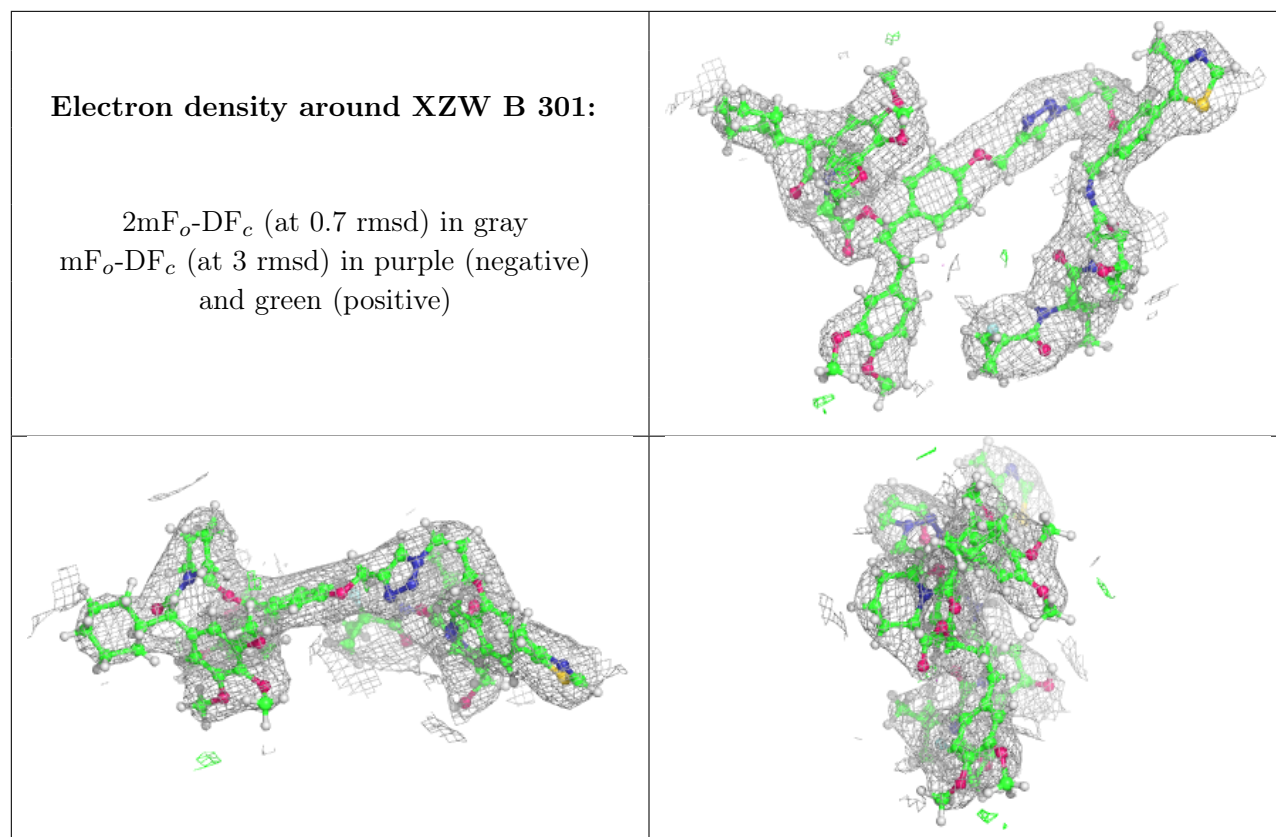
## 6.4 Ligands [i](#)

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<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
5	XZW	A	301	96/96	0.94	0.23	30,77,102,108	1
5	XZW	B	301	96/96	0.96	0.23	30,66,75,90	1

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.