

# wwPDB X-ray Structure Validation Summary Report (i)

#### May 19, 2020 - 11:34 am BST

PDB ID	:	1MBU
$\operatorname{Title}$	:	Crystal Structure Analysis of ClpSN heterodimer
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Deposited on		
Resolution	:	2.30  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 A user guide is available at https://www.wwpdb.org/validation/2017/XrayValidationReportHelp with specific help available everywhere you see the (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

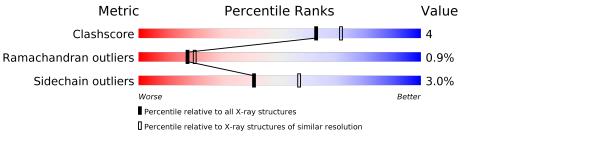
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	NOT EXECUTED
$\operatorname{EDS}$	:	NOT EXECUTED
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

# 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 2.30 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	Similar resolution
WEUTC	$(\# { m Entries})$	$(\# { m Entries}, { m resolution} { m range}({ m \AA}))$
Clashscore	141614	$5643 \ (2.30-2.30)$
Ramachandran outliers	138981	5575(2.30-2.30)
Sidechain outliers	138945	5575(2.30-2.30)

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 on 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%

Note EDS was not executed.

Mol	Chain	Length	Quality of chain				
1	А	142	90%			9%	•
1	В	142	86%			11%	••
2	С	106	73%	8%	•	19%	
2	D	106	75%	6%		19%	_



# 2 Entry composition (i)

There are 6 unique types of molecules in this entry. The entry contains 3851 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.

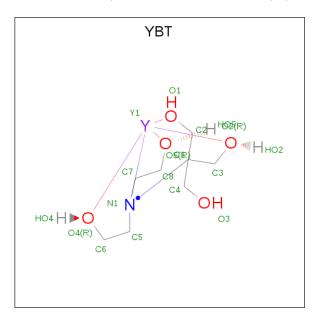
• Molecule 1 is a protein called ATP-Dependent clp Protease ATP-Binding Subunit clp A.

Mol	Chain	Residues		At	oms			ZeroOcc	AltConf	Trace
1	Δ	142	Total	С	Ν	Ο	S	0	0	0
	А	142	1127	702	203	218	4	0	0	0
1	р	142	Total	С	Ν	Ο	S	0	0	0
	D	142	1127	702	203	218	4	U	0	U

• Molecule 2 is a protein called Protein yljA.

Mol	Chain	Residues		At	oms			ZeroOcc	AltConf	Trace
0	C	86	Total	С	Ν	Ο	S	0	0	0
	U	80	686	444	109	127	6	0	0	0
0	л	86	Total	С	Ν	Ο	$\mathbf{S}$	0	0	0
			686	444	109	127	6			U

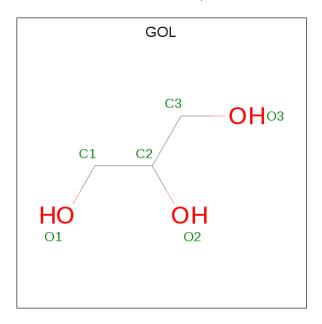
• Molecule 3 is BIS-(2-HYDROXYETHYL)AMINO-TRIS(HYDROXYMETHYL)METHAN E YTTRIUM (three-letter code: YBT) (formula: C<sub>8</sub>H<sub>19</sub>NO<sub>5</sub>Y).





Mol	Chain	Residues		Ate	oms			ZeroOcc	AltConf
3	Δ	1	Total	С	Ν	0	Υ	0	0
0	А	L	15	8	1	5	1	0	0
3	В	1	Total	С	Ν	0	Y	0	0
0	D	L	15	8	1	5	1	0	0
3	р	1	Total	С		0	Υ	0	0
0	D	L	15	8	1	5	1	0	0
3	Л	1	Total	С	Ν	0	Υ	0	0
J	D	L	15	8	1	5	1	0	0

• Molecule 4 is GLYCEROL (three-letter code: GOL) (formula:  $C_3H_8O_3$ ).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 6 & 3 & 3 \end{array}$	0	0
4	В	1	$\begin{array}{ccc} \text{Total}  \text{C}  \text{O} \\ 6  3  3 \end{array}$	0	0

• Molecule 5 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	В	2	Total Cl 2 2	0	0

• Molecule 6 is water.



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	А	42	$\begin{array}{cc} \text{Total} & \text{O} \\ 42 & 42 \end{array}$	0	0
6	В	40	Total         O           40         40	0	0
6	С	21	TotalO2121	0	0
6	D	48	Total         O           48         48	0	0

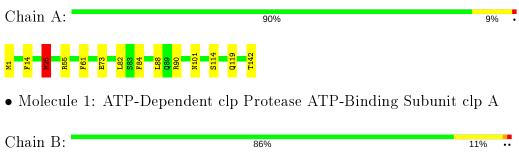


# 3 Residue-property plots (i)

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. Residues are colorcoded 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 was not executed.

• Molecule 1: ATP-Dependent clp Protease ATP-Binding Subunit clp A



Cham D.	0070		1170 ••
M1 14 14 15 15 15 15 16 16 16 16 16 16 16 16 16 16 16 16 16	E14 176 178 178 1125 1125 1125		
• Molecule 2: Protein yljA			
Chain C:	73%	8% •	19%
MET GLY THR ASN ASN ASP ASP ASP ASP ASP ASP ASP ASP ASP ASP	132 132 133 133 133 133 133 133 133 133		
• Molecule 2: Protein yljA			
Chain D:	75%	6%	19%
MET 0.11Y 1.1NS 1.1NS A.SN A.SP A.SP A.SP A.SP 0.1U 0.1U 0.1U 0.1U 0.1U 0.1U 0.1U 0.1U	E56 E56 L61 L103 A106		



## 4 Data and refinement statistics (i)

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source	
Space group	P 31 2 1	Depositor	
Cell constants	87.91Å 87.91Å 209.42Å	Depositor	
a, b, c, $\alpha$ , $\beta$ , $\gamma$	$90.00^{\circ}$ $90.00^{\circ}$ $120.00^{\circ}$	Depositor	
Resolution (Å)	19.85 - 2.30	Depositor	
% Data completeness	97.8 (19.85-2.30)	Depositor	
(in resolution range)	· · · · · ·	Depositor	
$R_{merge}$	(Not available)	Depositor	
R <sub>sym</sub>	0.07	Depositor	
Refinement program	CNS 1.1	Depositor	
$R, R_{free}$	0.217 , $0.251$	Depositor	
Estimated twinning fraction	No twinning to report.	Xtriage	
Total number of atoms	3851	wwPDB-VP	
Average B, all atoms $(Å^2)$	43.0	wwPDB-VP	



# 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: GOL, YBT, CL

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	А	0.75	0/1144	0.75	1/1548~(0.1%)
1	В	0.77	0/1144	0.83	2/1548~(0.1%)
2	С	0.74	1/700~(0.1%)	0.78	0/947
2	D	0.83	0/700	0.79	0/947
All	All	0.77	1/3688~(0.0%)	0.79	3/4990~(0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	$\operatorname{Ideal}(\operatorname{\AA})$
2	С	73	CYS	CB-SG	-5.51	1.72	1.81

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
1	А	25	MET	CG-SD-CE	-5.99	90.62	100.20
1	В	25	MET	CG-SD-CE	-5.86	90.83	100.20
1	В	8	LEU	CA-CB-CG	5.65	128.30	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	А	1127	0	1117	9	0
1	В	1127	0	1117	17	0
2	С	686	0	692	4	1
2	D	686	0	692	3	0
3	А	15	0	19	0	0
3	В	30	0	38	0	1
3	D	15	0	19	0	0
4	А	6	0	8	1	0
4	В	6	0	8	0	0
5	В	2	0	0	1	0
6	А	42	0	0	0	0
6	В	40	0	0	0	0
6	С	21	0	0	0	0
6	D	48	0	0	0	0
All	All	3851	0	3710	32	1

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 32 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:B:131:ARG:NH2	5:B:144:CL:CL	2.60	0.71
1:B:14:PHE:CE2	1:B:18:ARG:CZ	2.76	0.68
1:B:14:PHE:CE2	1:B:18:ARG:NH2	2.66	0.63
1:A:14:PHE:CD1	1:A:25:MET:HE2	2.36	0.59
1:A:14:PHE:HD1	1:A:25:MET:HE2	1.68	0.58

All (1) 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)
2:C:41:GLU:OE2	3:B:145:YBT:Y1[1_655]	2.16	0.04

### 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	Percen	ntiles
1	А	140/142~(99%)	136~(97%)	3(2%)	1 (1%)	22	26
1	В	140/142~(99%)	137~(98%)	0	3~(2%)	7	5
2	С	84/106~(79%)	82~(98%)	2(2%)	0	100	100
2	D	84/106~(79%)	82~(98%)	2(2%)	0	100	100
All	All	448/496~(90%)	437 (98%)	7 (2%)	4 (1%)	17	20

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	В	72	SER
1	В	73	GLU
1	А	73	GLU
1	В	71	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	А	124/124~(100%)	122~(98%)	2(2%)	62 78
1	В	124/124~(100%)	119~(96%)	5(4%)	31 44
2	С	75/93~(81%)	73~(97%)	2(3%)	44 61
2	D	75/93~(81%)	72~(96%)	3~(4%)	31 44
All	All	398/434~(92%)	386~(97%)	12 (3%)	41 57

5 of 12 residues with a non-rotameric side chain are listed below:

Mol	Chain	Res	Type
1	В	39	SER
1	В	76	ARG
2	D	32	LEU

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Mol	Chain	Res	Type
1	В	25	MET
2	С	103	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	А	64	GLN
1	А	94	HIS
2	С	68	GLN
2	D	68	GLN
2	D	97	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 carbohydrates in this entry.

### 5.6 Ligand geometry (i)

Of 8 ligands modelled in this entry, 2 are monoatomic - leaving 6 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	Tuno	Chain	Ros	Link	B	ond leng	gths	В	ond ang	gles
INIOI	туре	Unam	nes		Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
3	YBT	D	107	-	$6,\!18,\!18$	1.63	1 (16%)	$3,\!33,\!33$	2.37	1 (33%)



Mol	Tune	Chain	Res	Link	B	ond leng	gths	B	ond ang	gles
	Type	Chain	nes		Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
4	GOL	А	144	-	5, 5, 5	0.70	0	$5,\!5,\!5$	0.36	0
3	YBT	В	145	-	$6,\!18,\!18$	0.80	0	$3,\!33,\!33$	1.96	1 (33%)
3	YBT	В	146	-	6,18,18	1.46	1 (16%)	3,33,33	2.50	1(33%)
3	YBT	А	143	-	6,18,18	1.32	1 (16%)	3,33,33	2.38	1 (33%)
4	GOL	В	147	-	5, 5, 5	0.92	0	$5,\!5,\!5$	0.32	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	YBT	D	107	-	-	0/3/58/58	0/4/4/4
4	GOL	А	144	-	-	0/4/4/4	-
3	YBT	В	145	-	-	1/3/58/58	0/4/4/4
3	YBT	В	146	-	-	3/3/58/58	0/4/4/4
3	YBT	А	143	-	-	3/3/58/58	0/4/4/4
4	GOL	В	147	-	-	2/4/4/4	-

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$\operatorname{Observed}(\operatorname{\AA})$	Ideal(Å)
3	D	107	YBT	C3-C1	3.64	1.57	1.53
3	В	146	YBT	C3-C1	2.97	1.57	1.53
3	А	143	YBT	C3-C1	2.58	1.56	1.53

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$\mathbf{Observed}(^{o})$	$Ideal(^{o})$
3	В	146	YBT	C3-C1-C4	-4.28	104.08	110.46
3	D	107	YBT	C3-C1-C4	-4.09	104.36	110.46
3	А	143	YBT	C3-C1-C4	-4.00	104.50	110.46
3	В	145	YBT	C3-C1-C4	-3.34	105.49	110.46

There are no chirality outliers.

 $5~{\rm of}~9$  torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	В	146	YBT	N1-C1-C4-O3
3	А	143	YBT	N1-C1-C4-O3

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Mol	Chain	$\mathbf{Res}$	Type	Atoms				
3	А	143	YBT	C2-C1-C4-O3				
3	А	143	YBT	C3-C1-C4-O3				
4	В	147	GOL	C1-C2-C3-O3				

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There are no ring outliers.

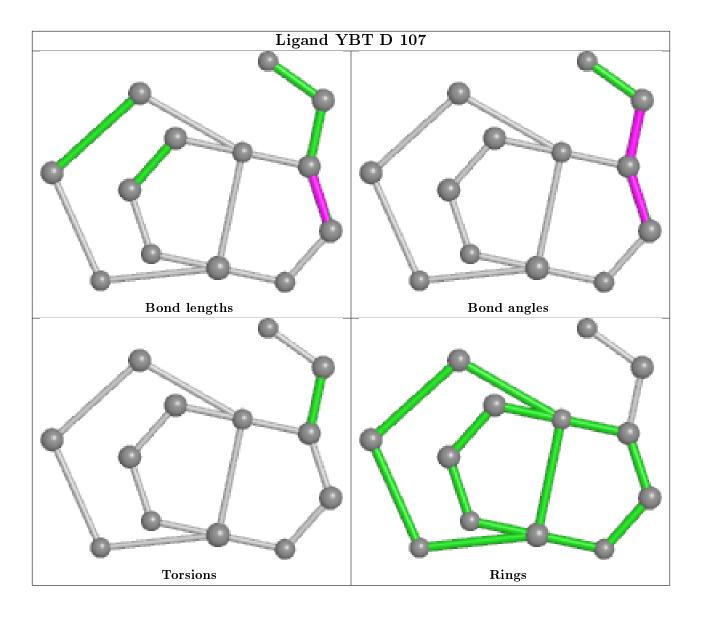
2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	А	144	GOL	1	0
3	В	145	YBT	0	1

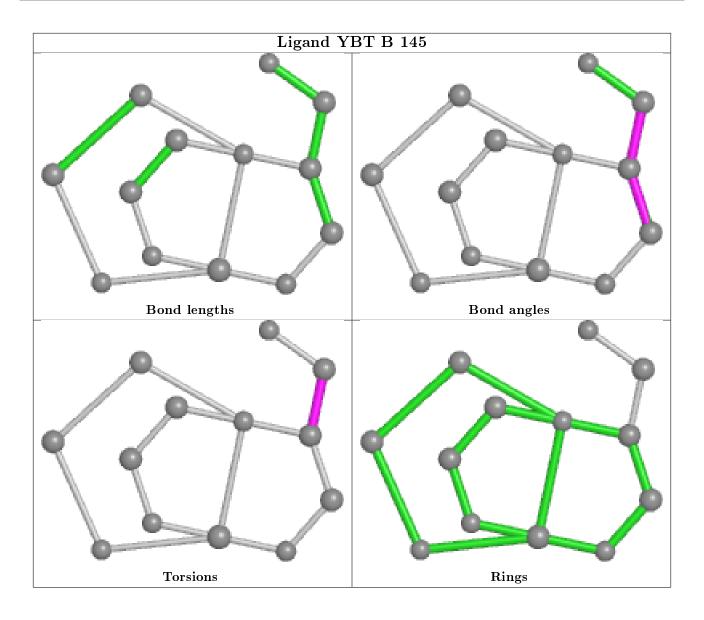
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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



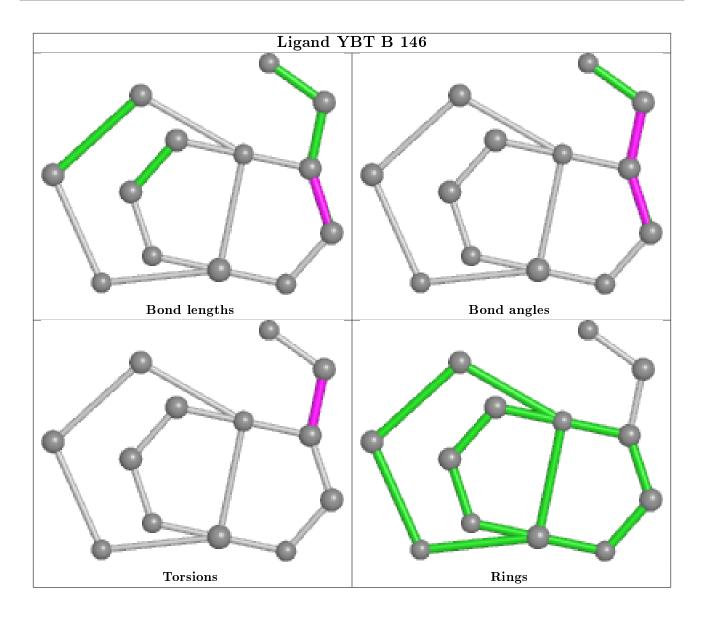






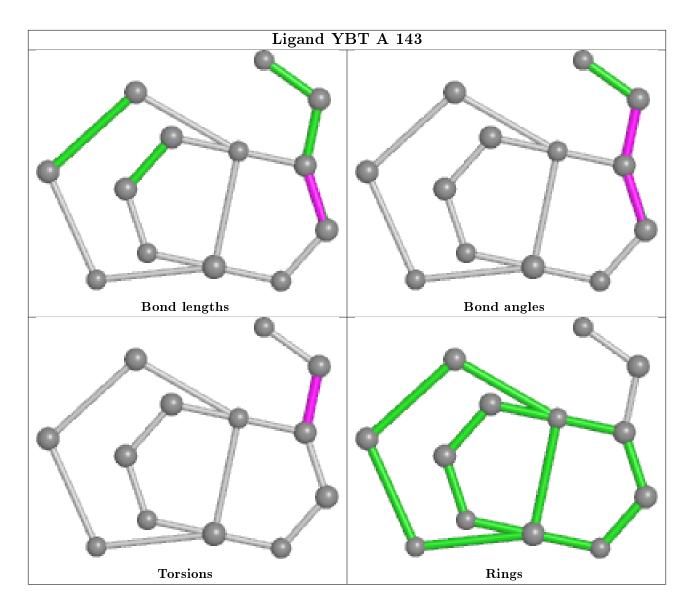












### 5.7 Other polymers (i)

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)

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

### 6.3 Carbohydrates (i)

EDS was not executed - this section is therefore empty.

### 6.4 Ligands (i)

EDS was not executed - this section is therefore empty.

#### 6.5 Other polymers (i)

EDS was not executed - this section is therefore empty.

