

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

#### May 28, 2020 – 09:46 pm BST

PDB ID	:	6TWC
$\operatorname{Title}$	:	Crystal Structure of the Catalytic Domain of the Coagulation Factor XIa in
		Complex with Double Bridged Peptide F21
Authors	:	Kong, X.D.; Pojer, F.; Heinis, C.
Deposited on	:	2020-01-13
$\operatorname{Resolution}$	:	2.86  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 following versions of software and data (see references (1)) were used in the production of this report:

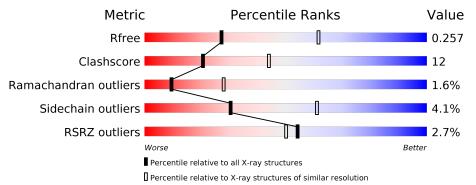
MolProbity		4.02b-467 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)		1.13
EDS	:	2.11
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
$\operatorname{Refmac}$	:	5.8.0158
$\operatorname{CCP4}$	:	$7.0.044 (\mathrm{Gargrove})$
Ideal 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 2.86 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	$egin{array}{c} { m Whole \ archive} \ (\#{ m Entries}) \end{array}$	${f Similar\ resolution}\ (\#{ m Entries},{ m resolution\ range}({ m \AA}))$
$R_{free}$	130704	3168 (2.90-2.82)
Clashscore	141614	3438 (2.90-2.82)
Ramachandran outliers	138981	3348 (2.90-2.82)
Sidechain outliers	138945	3351 (2.90-2.82)
RSRZ outliers	127900	3103 (2.90-2.82)

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% 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		$\mathbf{Q}$ ualit	y of chain		
1	А	244	<u>2%</u>	68%		27%	•••
2	Н	18	28%	17%	56%		
3	С	12	8% 	0%	33%		17%



#### 6TWC

# 2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 2053 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 Coagulation factor XI.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
1	А	238	Total 1888	C 1193	N 336	0 348	S 11	0	1	0

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
А	113	GLY	ASN	$\operatorname{conflict}$	UNP P03951
А	115	GLY	THR	$\operatorname{conflict}$	UNP P03951
А	246	HIS	-	expression tag	UNP P03951
А	247	HIS	-	expression tag	UNP P03951
A	248	HIS	-	expression tag	UNP P03951
А	249	HIS	-	expression tag	UNP P03951
А	250	HIS	-	expression tag	UNP P03951
А	251	HIS	_	expression tag	UNP P03951

• Molecule 2 is a protein called Coagulation factor XI.

Mol	Chain	Residues		At	oms			ZeroOcc	AltConf	Trace
2	Ц	8	Total	С	Ν	Ο	S	0	0	0
	11	0	62	39	10	12	1	0	0	0

There are 5 discrepancies between the modelled and reference sequences:

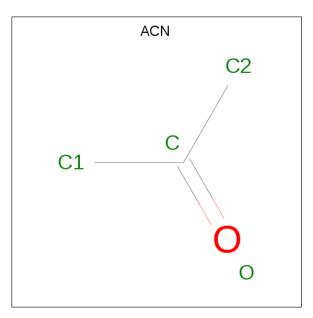
Chain	Residue	Modelled	Actual	Comment	Reference
Н	352	MET	-	initiating methionine	UNP P03951
Н	353	ASP	-	expression tag	UNP P03951
Н	354	ASP	-	expression tag	UNP P03951
Н	355	ASP	-	expression tag	UNP P03951
Н	356	ASP	-	expression tag	UNP P03951

• Molecule 3 is a protein called Double Bridged Peptide F21.



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
2	C	19	Total	С	Ν	Ο	S	0	0	0
J J	U	12	95	58	18	14	5	0	0	0

• Molecule 4 is ACETONE (three-letter code: ACN) (formula:  $C_3H_6O$ ) (labeled as "Ligand of Interest" by author).



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

Chain A:	2% 	ô	27% ••
116 V17 W29 W29	T35 T35 T35 T35 T35 T35 T36 T49 T52 T52 T52 T63 T63 T63 T63	669 669 773 774 775 774 775 775 775 775 775 775 775	E108 E108 E108 T110 S117 S117 D120 D122 D122 D122 D122 D123 D122 D123 D123
<mark>141 0142 0142</mark> К145 К146 К146 L147	K150 1451 1452 1452 1450 1450 1450 1450 1450 1450 14170 14172 14172	117 H178 M180 M180 M180 G184 G184 G184 G186 G186 G186 G186 G186 G186 C191 C192 C192 C192	2150 1200 1200 1200 1200 1200 1216 1216 121
HIS HIS HIS HIS HIS HIS			
• Molecu	le 2: Coagulation factor	XI	
Chain H:	28% 1	7% 569	%
MET ASP ASP ASP ASP LYS MET	ASP ASU AS61 C362 C362 C365 P366 ARG		
• Molecu	le 3: Double Bridged Pe	eptide F21	
Chain C:	8% 50%	33%	17%
11 C2 M6 15 M6 C7 C7	명 10 11 712		

• Molecule 1: Coagulation factor XI



# 4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants	76.16Å 76.16Å 117.08Å	Depositor
a, b, c, $\alpha$ , $\beta$ , $\gamma$	$90.00^{\circ}$ $90.00^{\circ}$ $120.00^{\circ}$	-
Resolution (Å)	43.78 - 2.86	Depositor
	43.78 - 2.86	EDS
% Data completeness	99.6 (43.78 - 2.86)	Depositor
(in resolution range)	99.7(43.78-2.86)	EDS
R <sub>merge</sub>	0.12	Depositor
R <sub>sym</sub>	(Not available)	Depositor
$< I/\sigma(I) > 1$	$2.08 (at 2.86 \text{\AA})$	Xtriage
Refinement program	PHENIX dev-3374	Depositor
D D	0.213 , $0.259$	Depositor
$R, R_{free}$	0.213 , $0.257$	DCC
$R_{free}$ test set	511 reflections $(5.38\%)$	wwPDB-VP
Wilson B-factor $(Å^2)$	62.0	Xtriage
Anisotropy	0.786	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$ , $B_{sol}(Å^2)$	0.35 , $50.6$	EDS
L-test for twinning <sup>2</sup>	$<  L  > = 0.49, < L^2 > = 0.31$	Xtriage
Estimated twinning fraction	0.039 for -h,-k,l	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	2053	wwPDB-VP
Average B, all atoms $(Å^2)$	67.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 7.22% of the height of the origin peak. No significant pseudotranslation is detected.

<sup>&</sup>lt;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.



 $<sup>^1 {\</sup>rm Intensities}$  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: ACN, LTN

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	
Mol Chain	RMSZ	# Z  > 5	RMSZ	# Z  > 5	
1	А	0.44	0/1936	0.62	0/2623
2	Н	0.48	0/62	0.80	0/82
3	С	3.85	15/80~(18.8%)	1.99	2/107~(1.9%)
All	All	0.87	15/2078~(0.7%)	0.72	2/2812~(0.1%)

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	А	0	2

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	С	9	ARG	CZ-NH2	11.30	1.47	1.33
3	С	11	PRO	N-CD	10.66	1.62	1.47
3	С	9	ARG	NE-CZ	9.79	1.45	1.33
3	С	10	CYS	C-N	9.21	1.51	1.34
3	С	4	ASN	C-N	8.81	1.54	1.34
3	С	1	THR	C-N	8.37	1.53	1.34
3	С	8	CYS	C-N	7.84	1.52	1.34
3	С	7	CYS	C-N	7.20	1.50	1.34
3	С	6	MET	C-N	7.11	1.50	1.34
3	С	2	CYS	C-N	7.00	1.50	1.34
3	С	9	ARG	C-N	6.83	1.49	1.34
3	С	11	PRO	N-CA	-6.64	1.35	1.47
3	С	3	VAL	C-N	6.44	1.48	1.34
3	С	5	ILE	C-N	6.41	1.48	1.34
3	С	4	ASN	CG-ND2	6.07	1.48	1.32

All (15) bond length outliers are listed below:



Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
3	С	6	MET	CG-SD-CE	7.77	112.64	100.20
3	С	10	CYS	CA-CB-SG	-5.92	103.34	114.00

All (2) bond angle outliers are listed below:

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	А	76	ILE	Peptide
1	А	77	LYS	Peptide

### 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	1888	0	1855	46	0
2	Н	62	0	67	3	0
3	С	95	0	92	5	0
4	С	8	0	6	1	0
All	All	2053	0	2020	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 12.

All (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)
1:A:74:SER:O	1:A:76:ILE:N	2.10	0.83
1:A:164:THR:HG23	1:A:167:GLU:H	1.52	0.74
1:A:203:HIS:O	1:A:204:ASN:ND2	2.24	0.70
3:C:1:THR:N	3:C:2:CYS:HA	2.08	0.68
1:A:86:GLN:HB2	1:A:109:GLU:HG3	1.76	0.67
1:A:32:THR:HG23	1:A:67:TYR:HB2	1.77	0.65
1:A:76:ILE:O	1:A:76:ILE:HD12	1.99	0.63
3:C:1:THR:H3	3:C:2:CYS:HA	1.62	0.63
1:A:217:GLU:HG3	3:C:5:ILE:HD12	1.82	0.61

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Continued from previou		Interatomic	Clash	
Atom-1	Atom-2	distance (Å)	overlap (Å)	
1:A:150:LYS:HD2	1:A:151:ILE:H	1.66	0.60	
1:A:49:ASN:ND2	2:H:368:PRO:HB3	2.19	0.57	
1:A:114:TYR:CE1	1:A:120:PRO:HD3	2.40	0.57	
1:A:243:GLN:HE22	2:H:367:LYS:H	1.54	0.56	
1:A:204:ASN:O	1:A:205:GLU:HB2	2.06	0.54	
1:A:73:GLN:HG2	1:A:141:TRP:CD1	2.45	0.52	
1:A:85:VAL:HG11	1:A:106:LEU:HD22	1.90	0.52	
1:A:150:LYS:CD	1:A:151:ILE:H	2.23	0.51	
1:A:16:ILE:O	1:A:145:ARG:HA	2.11	0.51	
1:A:52:ILE:HB	1:A:106:LEU:HB2	1.92	0.51	
1:A:35:THR:O	1:A:38:GLN:HA	2.11	0.50	
3:C:11:PRO:O	3:C:12:LTN:N	2.44	0.50	
1:A:195:SER:HA	1:A:213:THR:HG22	1.94	0.49	
1:A:62:SER:OG	1:A:64:LYS:HG3	2.12	0.49	
1:A:17:VAL:O	1:A:188:LYS:HA	2.13	0.48	
1:A:184(A):TYR:HB3	1:A:185:GLU:OE1	2.14	0.48	
1:A:143:TYR:HB3	1:A:146:LYS:O	2.14	0.48	
1:A:170:ARG:HG2	1:A:223:GLU:HG3	1.95	0.48	
1:A:51:TRP:CD2	1:A:242:THR:HG22	2.49	0.47	
1:A:145:ARG:HE	1:A:152:GLN:NE2	2.13	0.47	
1:A:184(A):TYR:O	1:A:186:GLY:N	2.38	0.47	
1:A:77:LYS:O	1:A:79:ASP:N	2.48	0.47	
1:A:167:GLU:O	1:A:170:ARG:HD2	2.16	0.46	
1:A:160:ILE:HD11	1:A:188:LYS:HB2	1.97	0.46	
1:A:67:TYR:CE2	1:A:82:PHE:HB3	2.51	0.46	
1:A:189:ASP:OD1	1:A:190:ALA:N	2.47	0.46	
1:A:160:ILE:HG22	1:A:183:ALA:HB1	1.98	0.45	
1:A:214:SER:O	4:C:102:ACN:H21	2.16	0.45	
1:A:145:ARG:HE	1:A:152:GLN:HE22	1.65	0.45	
1:A:214:SER:OG	1:A:215:TRP:HD1	2.01	0.44	
1:A:177:THR:HG22	1:A:180:MET:SD	2.58	0.44	
1:A:168(A):GLN:O	1:A:168(A):GLN:NE2	2.48	0.43	
1:A:51:TRP:CZ3	1:A:107:LYS:HB2	2.53	0.43	
1:A:146:LYS:HG3	1:A:147:LEU:O	2.18	0.43	
1:A:69:GLY:HA2	1:A:117:SER:O	2.19	0.42	
1:A:51:TRP:CH2	1:A:107:LYS:HD3	2.55	0.42	
1:A:28:PRO:HB2	1:A:119[B]:ARG:H	1.84	0.41	
1:A:192:LYS:HE2	3:C:8:CYS:O	2.19	0.41	
1:A:83:PHE:HD1	1:A:110:THR:HB	1.87	0.40	
1:A:29:TRP:CE2	1:A:121:ILE:HD12	2.56	0.40	
2:H:367:LYS:HA	2:H:368:PRO:HD3	1.83	0.40	

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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	А	237/244~(97%)	223~(94%)	11 (5%)	3~(1%)	12 33
2	Н	6/18~(33%)	5 (83%)	0	1 (17%)	0 0
3	С	9/12~(75%)	9 (100%)	0	0	100 100
All	All	252/274~(92%)	237~(94%)	11 (4%)	4 (2%)	9 28

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	А	74	SER
1	А	75	GLU
1	А	205	GLU
2	Н	362	CYS

#### 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	203/208~(98%)	196~(97%)	7(3%)	37	67
2	Н	8/18 (44%)	8 (100%)	0	100	100
3	С	11/11~(100%)	8 (73%)	3(27%)	0	1
All	All	222/237~(94%)	212~(96%)	10 (4%)	30	57



Mol	Chain	Res	Type
1	А	36(A)	SER
1	А	119[A]	ARG
1	А	119[B]	ARG
1	А	170	ARG
1	А	172	ARG
1	А	178	HIS
1	А	184(B)	ARG
3	С	1	THR
3	С	3	VAL
3	С	5	ILE

All (10) residues with a non-rotameric sidechain are listed below:

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

Mol	Chain	$\mathbf{Res}$	Type
1	А	204	ASN
1	А	243	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 carbohydrates 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



Mol Type Chain		Type Chain Res Li		Link Bond lengths			Bond angles			
	Type	Chain	nes		Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
4	ACN	С	101	3	$^{3,3,3}$	1.23	0	$^{3,3,3}$	0.22	0
4	ACN	С	102	3	3, 3, 3	1.29	0	$^{3,3,3}$	0.31	0

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).

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

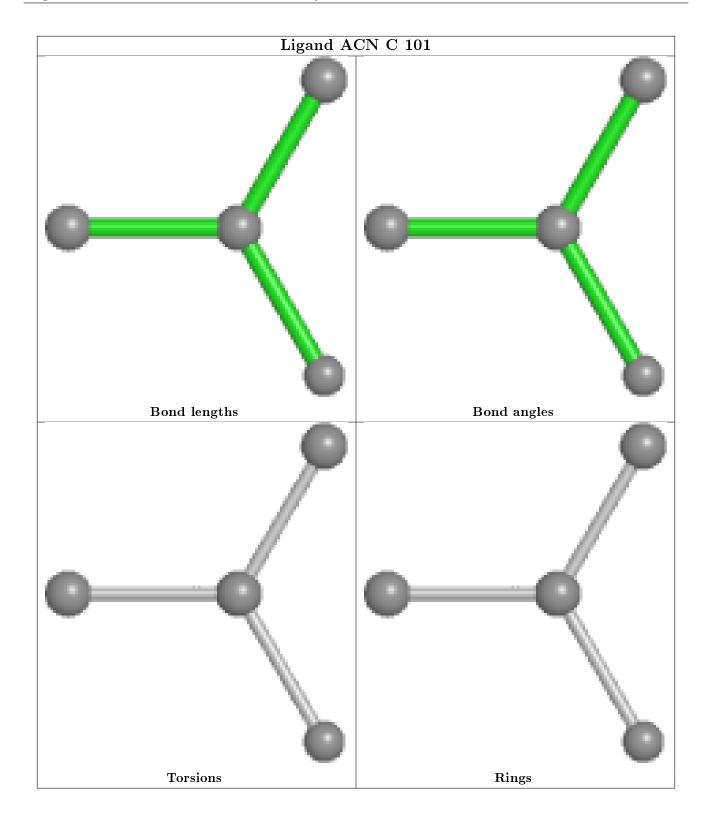
There are no ring outliers.

1 monomer is involved in 1 short contact:

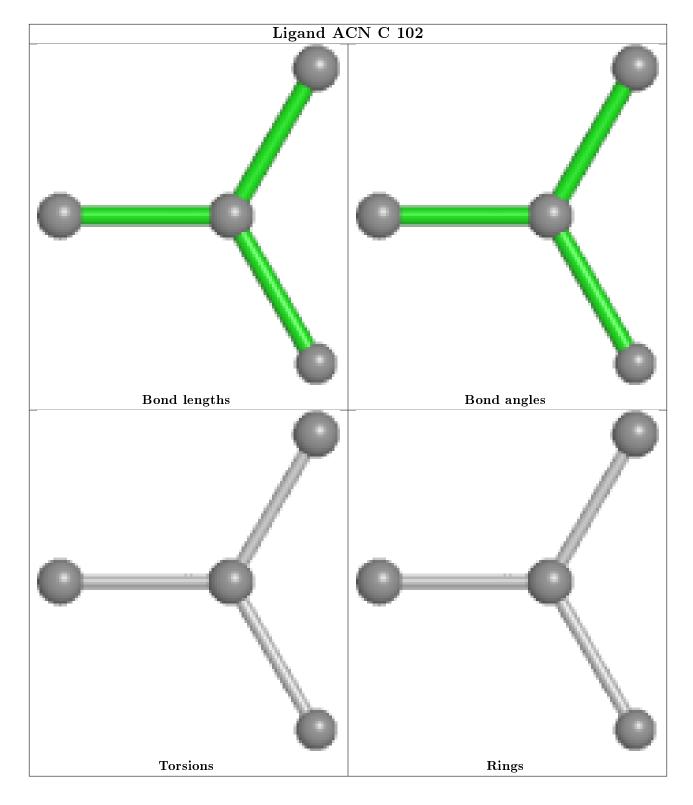
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	С	102	ACN	1	0

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 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	$\mathbf{OWAB}(\mathrm{\AA}^2)$	Q<0.9
1	А	238/244~(97%)	-0.04	6 (2%) 57 54	47, 63, 94, 126	0
2	Н	8/18 (44%)	0.27	0 100 100	83, 95, 97, 98	0
3	С	$11/12 \ (91\%)$	0.20	1 (9%) 9 6	57,61,85,85	0
All	All	257/274~(93%)	-0.02	7 (2%) 54 50	47, 63, 95, 126	0

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	А	130	ASN	6.1
1	А	131	VAL	3.5
1	А	132	ILE	3.3
1	А	127	GLY	3.0
3	С	1	THR	2.7
1	А	204	ASN	2.1
1	А	205	GLU	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 carbohydrates in this entry.

## 6.4 Ligands (i)

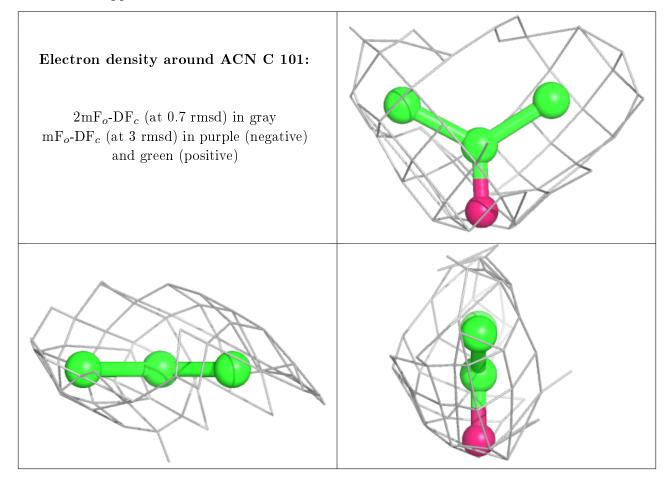
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum,



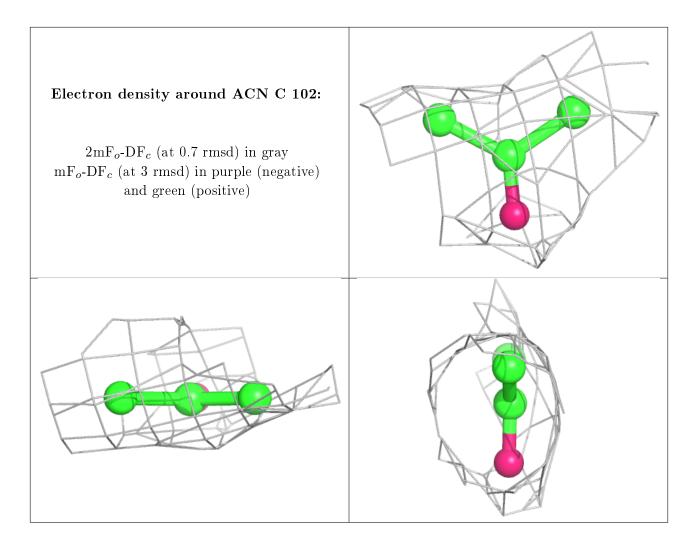
Mol	Type	Chain	Res	Atoms	RSCC	RSR	$B-factors(A^2)$	Q<0.9
4	ACN	С	101	4/4	0.92	0.17	64,73,75,80	0
4	ACN	С	102	4/4	0.98	0.14	$51,\!56,\!59,\!60$	0

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.

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.

