

# Full wwPDB NMR Structure Validation Report (i)

Jun 17, 2021 – 12:09 PM BST

PDB ID : 6GZ7

Title : Polyamide - DNA complex NMR structure

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This is a Full wwPDB NMR 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/NMRValidationReportHelp 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)

buster-report : 1.1.7 (2018)

Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)

RCI : v 1n 11 5 13 A (Berjanski et al., 2005)

PANAV : Wang et al. (2010)

ShiftChecker : 2.20

Ideal geometry (proteins) : Engh & Huber (2001) Ideal geometry (DNA, RNA) : Parkinson et al. (1996)

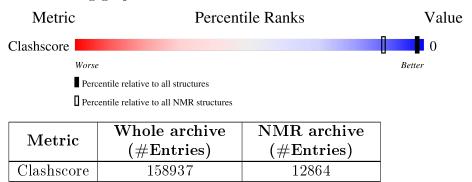
Validation Pipeline (wwPDB-VP) : 2.20

## 1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:  $SOLUTION\ NMR$ 

The overall completeness of chemical shifts assignment is 43%.

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.



The table below summarises the geometric issues observed across the polymeric chains and their fit to the experimental data. The red, orange, yellow and green segments indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria. A cyan segment indicates the fraction of residues that are not part of the well-defined cores, and 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%

Mol	Chain	Length		Quality of chain		
1	A	13	15%	54%	23%	8%
1	В	13		77%	15%	8%



# 2 Ensemble composition and analysis (i)

This entry contains 11 models. This entry does not contain polypeptide chains, therefore identification of well-defined residues and clustering analysis are not possible. All residues are included in the validation scores.



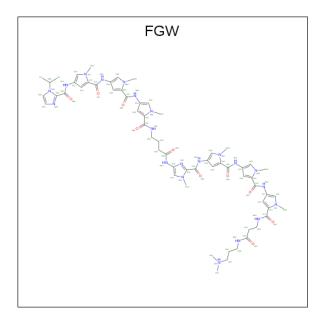
# 3 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 925 atoms, of which 348 are hydrogens and 0 are deuteriums.

• Molecule 1 is a DNA chain called DNA (5'-D(\*CP\*GP\*AP\*TP\*GP\*TP\*AP\*CP\*AP\*TP\*CP\*G)-3').

Mol	Chain	Residues		Atoms					Trace	
1	Λ	19	Total	С	Η	N	О	Р	0	
1   A	12	379	117	136	45	70	11			
1	D	19	Total	С	Н	N	О	Р	0	
1		Б 12		379	117	136	45	70	11	U

• Molecule 2 is dimethyl-[3-[3-[[1-methyl-4-[[1-methyl-4-[[1-methyl-4-[[1-methyl-4-[4-[[1-methyl-4-[4-[[1-methyl-4-[1-methyl-4



Mol	Chain	Residues	Atoms					
9	Λ	1	Total	С	Н	N	О	
2	А	1	167	60	76	21	10	

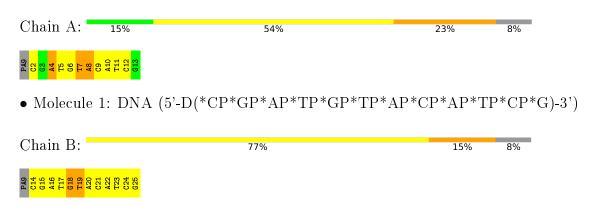


## 4 Residue-property plots (i)

### 4.1 Average score per residue in the NMR ensemble

These plots are provided for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic is the same as shown in the summary in section 1 of this report. The second graphic shows the sequence where residues are colour-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. Stretches of 2 or more consecutive residues without any outliers are shown as green connectors. Residues which are classified as ill-defined in the NMR ensemble, are shown in cyan with an underline colour-coded according to the previous scheme. Residues which were present in the experimental sample, but not modelled in the final structure are shown in grey.

• Molecule 1: DNA (5'-D(\*CP\*GP\*AP\*TP\*GP\*TP\*AP\*CP\*AP\*TP\*CP\*G)-3')

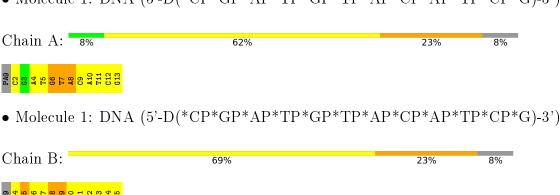


### 4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

#### 4.2.1 Score per residue for model 1

• Molecule 1: DNA (5'-D(\*CP\*GP\*AP\*TP\*GP\*TP\*AP\*CP\*AP\*TP\*CP\*G)-3')





#### 4.2.2 Score per residue for model 2

• Molecule 1: DNA (5'-D(\*CP\*GP\*AP\*TP\*GP\*TP\*AP\*CP\*AP\*TP\*CP\*G)-3')

Chain A: 15% 62% 15% 8%

• Molecule 1: DNA (5'-D(\*CP\*GP\*AP\*TP\*GP\*TP\*AP\*CP\*AP\*TP\*CP\*G)-3')

Chain B: 8% 69% 15% 8%

PA9 C14 A16 T17 G18 T19 A20 C21 A22 T23 C24 G26

#### 4.2.3 Score per residue for model 3

• Molecule 1: DNA (5'-D(\*CP\*GP\*AP\*TP\*GP\*TP\*AP\*CP\*AP\*TP\*CP\*G)-3')

Chain A: 15% 46% 31% 8%

PA9 C2 C3 A4 A4 T7 T7 A8 C9 A10 T111 C12 C12

Chain B: 8% 69% 15% 8%

PA9 C14 G15 A16 T17 T19 A20 C21 T23 T23 C24 G25

#### 4.2.4 Score per residue for model 4

Chain A: 15% 62% 15% 8%

PA9

C2

G3

A4

T5

T6

C9

C9

A10

T111

C112

• Molecule 1: DNA (5'-D(\*CP\*GP\*AP\*TP\*GP\*TP\*AP\*CP\*AP\*TP\*CP\*G)-3')

Chain B: 8% 69% 15% 8%



#### 4.2.5 Score per residue for model 5

• Molecule 1: DNA (5'-D(\*CP\*GP\*AP\*TP\*GP\*TP\*AP\*CP\*AP\*TP\*CP\*G)-3')

Chain A: 15% 46% 31% 8%

C2 C2 C3 A4 T5 C6 C9 C9 C12 C12 C12

• Molecule 1: DNA (5'-D(\*CP\*GP\*AP\*TP\*GP\*TP\*AP\*CP\*AP\*TP\*CP\*G)-3')

Chain B: 77% 15% 8%

PA9 C14 G15 A16 T17 T19 A20 C21 A22 T23 C24 C24

#### 4.2.6 Score per residue for model 6

• Molecule 1: DNA (5'-D(\*CP\*GP\*AP\*TP\*GP\*TP\*AP\*CP\*AP\*TP\*CP\*G)-3')

Chain A: 62% 31% 8%

PA9 G3 G3 A4 T5 G6 T7 A8 C9 A10 T111 C12 G13

Chain B: 46% 46% 8%

PA9 C14 G15 A16 T117 T119 A20 C21 A22 T23 C24 G25

#### 4.2.7 Score per residue for model 7

• Molecule 1: DNA (5'-D(\*CP\*GP\*AP\*TP\*GP\*TP\*AP\*CP\*AP\*TP\*CP\*G)-3')

Chain A: 8% 62% 23% 8%

• Molecule 1: DNA (5'-D(\*CP\*GP\*AP\*TP\*GP\*TP\*AP\*CP\*AP\*TP\*CP\*G)-3')

Chain B: 85% 8% 8%

PA9 C14 G15 A16 T17 T19 A20 C21 A22 T23 C24 G25



#### 4.2.8 Score per residue for model 8

• Molecule 1: DNA (5'-D(\*CP\*GP\*AP\*TP\*GP\*TP\*AP\*CP\*AP\*TP\*CP\*G)-3')

Chain A: 15% 54% 23% 8%

PA9
C2
C3
A4
T5
G6
C7
T7
T7
T7
T11
C12
C12

• Molecule 1: DNA (5'-D(\*CP\*GP\*AP\*TP\*GP\*TP\*AP\*CP\*AP\*TP\*CP\*G)-3')

Chain B: 8% 62% 23% 8%

PA9 C14 A16 T17 G18 T19 A20 C21 A22 C21 T23 C24 G24

#### 4.2.9 Score per residue for model 9

• Molecule 1: DNA (5'-D(\*CP\*GP\*AP\*TP\*GP\*TP\*AP\*CP\*AP\*TP\*CP\*G)-3')

Chain A: 15% 38% 38% 8%

• Molecule 1: DNA (5'-D(\*CP\*GP\*AP\*TP\*GP\*TP\*AP\*CP\*AP\*TP\*CP\*G)-3')

Chain B: 69% 23% 8%

PA9 C14 G15 A16 T17 G18 T19 A20 C21 T23 C24 G25

#### 4.2.10 Score per residue for model 10

• Molecule 1: DNA (5'-D(\*CP\*GP\*AP\*TP\*GP\*TP\*AP\*CP\*AP\*TP\*CP\*G)-3')

Chain A: 15% 46% 31% 8%

• Molecule 1: DNA (5'-D(\*CP\*GP\*AP\*TP\*GP\*TP\*AP\*CP\*AP\*TP\*CP\*G)-3')

Chain B: 8% 8% 8%

PA9 (14 (14 (14 (15) (18) (18) (19)



### 4.2.11 Score per residue for model 11

Chain A: 8% 54% 31% 8%

PA9

C2

G3

A4

T5

G6

G6

G7

T7

A8

C9

A10

T111

C112

• Molecule 1: DNA (5'-D(\*CP\*GP\*AP\*TP\*GP\*TP\*AP\*CP\*AP\*TP\*CP\*G)-3')

Chain B: 8% 62% 23% 8%

PA9 C144 C145 C154 C117 C21 C21 C21 C24 C24 C25



#### 5 Refinement protocol and experimental data overview (i)



The models were refined using the following method: molecular dynamics.

Of the 11 calculated structures, 11 were deposited, based on the following criterion: all calculated structures submitted.

The following table shows the software used for structure solution, optimisation and refinement.

Software name	Classification	Version
Amber	structure calculation	

The following table shows chemical shift validation statistics as aggregates over all chemical shift files. Detailed validation can be found in section 7 of this report.

Chemical shift file(s)	$working\_cs.cif$
Number of chemical shift lists	1
Total number of shifts	250
Number of shifts mapped to atoms	250
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	43%



## 6 Model quality (i)

### 6.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: FGW

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 (average) root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	E	ond lengths	Bond angles		
Wioi Chain		RMSZ	#Z>5	RMSZ	#Z>5	
1	A	$1.56 \pm 0.01$	$0\pm0/272~(~0.0\pm~0.0\%)$	$2.43 \pm 0.03$	$24\pm2/418~(~5.8\pm~0.4\%)$	
1	В	$1.59 \pm 0.01$	$0\pm0/272~(~0.0\pm~0.0\%)$	$2.47 \pm 0.04$	$26\pm2/418~(~6.1\pm~0.6\%)$	
All	All	1.57	0/5984~(~0.0%)	2.45	548/9196 ( 6.0%)	

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	Planarity
1	A	$0.0 \pm 0.0$	$4.2 \pm 0.9$
1	В	$0.0 \pm 0.0$	$2.7{\pm}1.1$
All	All	0	76

There are no bond-length outliers.

All unique angle outliers are listed below. They are sorted according to the Z-score of the worst occurrence in the ensemble.

Mol	Chain	Res	Trno	Atoms	$\mathbf{z}$	$Observed(^o)$	$\operatorname{Ideal}({}^{o})$	Models	
IVIOI	Chain	nes	Type	Atoms		Observed(')	Ideal(*)	Worst	Total
1	В	22	DA	N1-C6-N6	-9.63	112.82	118.60	6	11
1	A	4	DA	N1-C6-N6	-9.53	112.88	118.60	6	11
1	A	10	DA	N1-C6-N6	-9.48	112.91	118.60	6	11
1	В	16	DA	O4'-C1'-N9	9.18	114.42	108.00	6	5
1	В	22	DA	O4'-C1'-N9	9.17	114.42	108.00	7	9
1	В	16	DA	N1-C6-N6	-8.77	113.34	118.60	9	11
1	A	11	DT	O4'-C1'-N1	8.62	114.03	108.00	9	9
1	A	8	DA	C5-C6-N1	8.46	121.93	117.70	6	11
1	В	20	DA	C5-C6-N1	8.43	121.92	117.70	10	11
1	В	21	DC	N3-C2-O2	-8.21	116.15	121.90	7	11

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			T		7	Observed(0)	T.J. (0)	Mod	dels
Mol	Chain	Res	Type	Atoms	$\mathbf{Z}$	$\operatorname{Observed}(^{o})$	$\operatorname{Ideal}({}^{o})$	Worst	Total
1	A	2	DC	N3-C2-O2	-7.91	116.36	121.90	5	11
1	A	9	DC	N3-C2-O2	-7.87	116.39	121.90	8	11
1	В	25	DG	O4'-C1'-N9	7.74	113.42	108.00	2	9
1	A	12	DC	N3-C2-O2	-7.65	116.55	121.90	2	11
1	В	16	DA	C5-C6-N1	7.58	121.49	117.70	2	11
1	В	14	DC	N3-C2-O2	-7.57	116.60	121.90	3	11
1	В	24	DC	N3-C2-O2	-7.49	116.66	121.90	9	11
1	A	5	DT	O4'-C1'-N1	7.46	113.22	108.00	10	10
1	В	22	DA	C5-C6-N1	7.39	121.39	117.70	5	11
1	A	4	DA	C5-C6-N1	7.35	121.38	117.70	1	11
1	В	22	DA	C4-C5-C6	-7.31	113.35	117.00	8	11
1	A	10	DA	C5-C6-N1	7.13	121.27	117.70	9	11
1	A	4	DA	C4-C5-C6	-7.00	113.50	117.00	6	11
1	В	21	DC	N1-C2-O2	6.94	123.07	118.90	7	11
1	В	20	DA	N1-C6-N6	-6.82	114.51	118.60	9	11
1	В	16	DA	C4-C5-C6	-6.81	113.59	117.00	11	11
1	A	10	DA	C4-C5-C6	-6.78	113.61	117.00	2	11
1	A	13	DG	O4'-C1'-N9	6.72	112.70	108.00	6	3
1	A	11	DT	C6-C5-C7	-6.71	118.88	122.90	1	11
1	В	23	DT	O4'-C1'-N1	6.66	112.66	108.00	9	3
1	В	14	DC	O4'-C1'-N1	6.62	112.64	108.00	6	3
1	A	12	DC	O4'-C1'-N1	6.61	112.62	108.00	10	5
1	A	8	DA	N1-C6-N6	-6.60	114.64	118.60	9	11
1	A	5	DT	N3-C2-O2	-6.59	118.35	122.30	3	11
1	В	17	DT	O4'-C1'-N1	6.58	112.61	108.00	7	1
1	A	5	DT	C6-C5-C7	-6.47	119.02	122.90	4	11
1	A	2	DC	N1-C2-O2	6.47	122.78	118.90	3	11
1	В	17	DT	C6-C5-C7	-6.45	119.03	122.90	8	11
1	A	11	DT	O4'-C1'-C2'	-6.40	100.78	105.90	7	1
1	В	15	DG	O4'-C1'-N9	6.39	112.47	108.00	9	2
1	В	24	DC	O4'-C1'-N1	6.35	112.44	108.00	9	3
1	В	23	DT	C6-C5-C7	-6.33	119.10	122.90	11	11
1	A	8	DA	C4-C5-C6	-6.32	113.84	117.00	3	11
1	A	9	DC	N1-C2-O2	6.26	122.66	118.90	10	11
1	A	2	DC	O4'-C1'-N1	6.19	112.34	108.00	2	3
1	A	7	DT	C6-C5-C7	-6.15	119.21	122.90	10	11
1	В	24	DC	N1-C2-O2	6.14	122.58	118.90	9	11
1	В	23	DT	N3-C2-O2	-6.13	118.62	122.30	3	10
1	В	14	DC	N1-C2-O2	6.12	122.57	118.90	3	5
1	A	12	DC	N1-C2-O2	6.11	122.57	118.90	2	9
1	В	19	DT	C6-C5-C7	-6.03	119.28	122.90	7	10

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Mol	Chain	Res	Trno	Atoms	$\mathbf{z}$	$Observed(^o)$	$\mathrm{Ideal}(^{o})$	Mod	dels
IVIOI	Chain	nes	Type	Atoms		Observed()	Ideal( )	Worst	Total
1	В	19	DT	N3-C2-O2	-6.03	118.68	122.30	10	7
1	A	11	DT	N3-C2-O2	-5.94	118.74	122.30	2	11
1	В	17	DT	N3-C2-O2	-5.91	118.75	122.30	1	11
1	В	18	DG	N1-C6-O6	-5.84	116.40	119.90	8	8
1	В	20	DA	C4-C5-C6	-5.83	114.08	117.00	10	11
1	A	5	DT	O4'-C1'-C2'	-5.50	101.50	105.90	4	2
1	В	25	DG	C1'-O4'-C4'	-5.48	104.62	110.10	11	1
1	В	16	DA	O4'-C1'-C2'	-5.47	101.52	105.90	6	1
1	A	7	DT	N3-C2-O2	-5.41	119.05	122.30	6	6
1	В	25	DG	O4'-C1'-C2'	-5.38	101.60	105.90	1	3
1	В	15	DG	N1-C6-O6	-5.35	116.69	119.90	1	3
1	В	25	DG	N1-C6-O6	-5.28	116.73	119.90	11	5
1	В	18	DG	C8-N9-C4	-5.22	104.31	106.40	1	4
1	A	6	DG	N1-C6-O6	-5.20	116.78	119.90	1	4
1	A	13	DG	N1-C6-O6	-5.17	116.80	119.90	1	3
1	A	12	DC	N3-C4-C5	5.17	123.97	121.90	5	2
1	В	18	DG	N3-C4-C5	-5.11	126.05	128.60	11	1
1	A	6	DG	C5-C6-N1	5.06	114.03	111.50	5	1
1	В	21	DC	O4'-C1'-C2'	-5.04	101.86	105.90	1	1

There are no chirality outliers.

All unique planar outliers are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type	Group	Models (Total)
1	A	6	DG	Sidechain	11
1	A	7	DT	Sidechain	11
1	A	8	DA	Sidechain	11
1	В	18	DG	Sidechain	11
1	В	19	DT	Sidechain	11
1	A	4	DA	Sidechain	7
1	В	15	DG	Sidechain	4
1	A	11	DT	Sidechain	4
1	В	14	DC	Sidechain	2
1	A	3	DG	Sidechain	1
1	В	23	DT	Sidechain	1
1	В	24	DC	Sidechain	1
1	A	12	DC	Sidechain	1



### 6.2 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 each 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 averaged over the ensemble.

Mol	Chain	Non-H	H(model)	$\mathbf{H}(\mathbf{added})$	Clashes
1	В	243	136	137	0±0
2	A	91	76	0	0±0
All	All	6347	3828	3014	3

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.

All unique clashes are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Clack(Å)	${f Distance(\AA)}$	Mod	
	Atom-2	Clash(A)	Distance(A)	Worst	Total
2:A:101:FGW:C13	1:B:22:DA:H1'	0.43	2.43	8	3

### 6.3 Torsion angles (i)

### 6.3.1 Protein backbone (i)

There are no protein molecules in this entry.

### 6.3.2 Protein sidechains (i)

There are no protein molecules in this entry.

### 6.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

### 6.5 Carbohydrates (i)

There are no monosaccharides in this entry.



### 6.6 Ligand geometry (i)

1 ligand is modelled in this entry.

In the following table, the Counts columns list the number of bonds for which Mogul statistics could be retrieved, the number of bonds that are observed in the model and the number of bonds 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 is the number of standard deviations the observed value is removed from the expected value. A bond length with |Z| > 2 is considered an outlier worth inspection. RMSZ is the average root-mean-square of all Z scores of the bond lengths.

Mol	Type	Chain	Pos	Link		Bond len	$_{ m igths}$
	туре	Chain	res	LIIIK	Counts	RMSZ	#Z>2
2	FGW	A	101	-	87,98,98	$1.31 \pm 0.02$	11±1 (12±1%)

In the following table, the Counts columns list the number of angles for which Mogul statistics could be retrieved, the number of angles that are observed in the model and the number of 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 angle is the number of standard deviations the observed value is removed from the expected value. A bond angle with |Z| > 2 is considered an outlier worth inspection. RMSZ is the average root-mean-square of all Z scores of the bond angles.

Mol	Tuno	Chain	Dog	Tinle		Bond ang	gles
	Type	Chain	res	LIIIK	Counts	RMSZ	#Z>2
2	FGW	A	101	-	83,140,140	$1.69 \pm 0.02$	12±0 (14±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
2	FGW	A	101	_	_	$0\pm0,44,84,84$	$0\pm0,8,8,8$

All unique bond outliers are listed below. They are sorted according to the Z-score of the worst occurrence in the ensemble.

Mal	Mol Chain	Pos	Trme	Res Type	Atoms	Z Observed(Å		Ideal(Å)	Mod	dels
IVIOI	Chain	nes	туре	Atoms	L	Observed(A)	Ideal(A)	Worst	Total	
2	A	101	FGW	C73-C74	3.65	1.42	1.39	3	11	
2	A	101	FGW	C33-C34	3.49	1.42	1.39	9	11	
2	A	101	FGW	C82-C81	3.40	1.50	1.53	3	11	
2	A	101	FGW	C23-C24	3.36	1.42	1.39	9	11	

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Mol	Chain	Res	Tuno	Atoma	Z	$Observed(\AA)$	$Ideal(\mathring{A})$	Mod	dels
IVIOI	Chain	nes	Type	Atoms	L	Observed(A)	ideai(A)	Worst	Total
2	A	101	FGW	C63-C64	3.04	1.42	1.39	3	11
2	A	101	FGW	C42-C41	2.82	1.51	1.53	4	11
2	A	101	FGW	C13-C14	2.63	1.41	1.39	9	11
2	A	101	FGW	C53-C54	2.57	1.41	1.39	11	11
2	A	101	FGW	C01-N01	2.45	1.53	1.46	7	11
2	A	101	FGW	C02-N01	2.42	1.53	1.46	10	11
2	A	101	FGW	C22-N21	2.08	1.40	1.36	11	1
2	A	101	FGW	C06-N02	2.08	1.38	1.33	11	2
2	A	101	FGW	C32-N31	2.04	1.40	1.36	7	3
2	A	101	FGW	C11-N03	2.03	1.38	1.33	4	2
2	A	101	FGW	C52-N51	2.03	1.40	1.36	3	1

All unique angle outliers are listed below. They are sorted according to the Z-score of the worst occurrence in the ensemble.

Mol	Chain	Res	Type	Atoms	Z	$Observed(^o)$	$Ideal(^{o})$	Mod	dels
WIOI	Chain	ites	Type	Atoms		Observed()	ideai()	Worst	Total
2	A	101	FGW	C75-C74-C73	6.01	108.99	106.05	7	11
2	A	101	FGW	C35-C34-C33	5.52	108.75	106.05	2	11
2	A	101	FGW	C65-C64-C63	5.35	108.67	106.05	10	11
2	A	101	FGW	C55-C54-C53	5.17	108.58	106.05	2	11
2	A	101	FGW	C15-C14-C13	5.10	108.55	106.05	4	11
2	A	101	FGW	C25-C24-C23	4.52	108.26	106.05	5	11
2	A	101	FGW	C83-N81-C88	3.96	121.49	125.40	9	11
2	A	101	FGW	C9-C88-N81	3.89	116.66	110.30	5	11
2	A	101	FGW	C83-N81-C82	3.28	104.93	108.58	5	11
2	A	101	FGW	C8-C88-N81	2.93	115.09	110.30	6	11
2	A	101	FGW	C44-N41-C42	2.53	105.87	108.60	3	11
2	A	101	FGW	CM4-N41-C42	2.52	129.97	126.39	4	9
2	A	101	FGW	CM2-N21-C25	2.24	120.69	124.90	11	1

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

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.

### 6.7 Other polymers (i)

There are no such molecules in this entry.

### 6.8 Polymer linkage issues (i)

There are no chain breaks in this entry.



## 7 Chemical shift validation (i)

The completeness of assignment taking into account all chemical shift lists is 43% for the well-defined parts and 43% for the entire structure.

#### 7.1 Chemical shift list 1

File name: working cs.cif

Chemical shift list name: NMR\_STAR\_FILE\_3.txt

### 7.1.1 Bookkeeping (i)

The following table shows the results of parsing the chemical shift list and reports the number of nuclei with statistically unusual chemical shifts.

Total number of shifts	250
Number of shifts mapped to atoms	250
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	7

### 7.1.2 Chemical shift referencing (i)

No chemical shift referencing corrections were calculated (not enough data).

### 7.1.3 Completeness of resonance assignments (i)

The following table shows the completeness of the chemical shift assignments for the well-defined regions of the structure. The overall completeness is 43%, i.e. 206 atoms were assigned a chemical shift out of a possible 474. 0 out of 0 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	$^{1}\mathrm{H}$	$^{13}\mathbf{C}$	$^{15}{ m N}$
Backbone	0/0 (%)	0/0 (%)	0/0 (%)	0/0 (%)
Sidechain	0/0 (%)	0/0 (%)	0/0 (%)	0/0 (%)
Aromatic	0/0 (%)	0/0 (%)	0/0 (%)	0/0 (%)
Overall	206/474~(43%)	$206/282 \ (73\%)$	0/162~(0%)	0/30 (0%)

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 43%, i.e. 206 atoms were assigned a chemical shift out of a possible 474. 0 out of 0 assigned methyl groups (LEU and VAL) were assigned stereospecifically.



	Total	$^{1}\mathrm{H}$	$^{13}\mathbf{C}$	$^{15}{ m N}$
Backbone	0/0 (—%)	0/0 (—%)	0/0 (%)	0/0 (%)
Sidechain	0/0 (%)	0/0 (%)	0/0 (%)	0/0 (%)
Aromatic	0/0 (%)	0/0 (%)	0/0 (%)	0/0 (%)
Overall	206/474~(43%)	$206/282 \ (73\%)$	0/162~(0%)	0/30 (0%)

#### 7.1.4 Statistically unusual chemical shifts (i)

The following table lists the statistically unusual chemical shifts. These are statistical measures, and large deviations from the mean do not necessarily imply incorrect assignments. Molecules containing paramagnetic centres or hemes are expected to give rise to anomalous chemical shifts.

Mol	Chain	Res	Type	Atom	Shift, ppm	Expected range, ppm	Z-score
1	A	7	DT	H4'	2.00	5.65 - 2.55	-6.8
1	В	19	DT	H4'	2.06	5.65 - 2.55	-6.6
1	В	20	DA	H4'	2.65	5.86 - 2.76	-5.4
1	A	10	DA	H4'	2.67	5.86 - 2.76	-5.3
1	A	9	DC	H4'	2.22	5.92 - 2.32	-5.3
1	В	22	DA	H4'	2.68	5.86 - 2.76	-5.3
1	В	21	DC	H4'	2.30	5.92 - 2.32	-5.1

### 7.1.5 Random Coil Index (RCI) plots (i)

No random coil index (RCI) plot could be generated from the current chemical shift list (NMR\_STAR\_FILE\_RCI is only applicable to proteins.

