

Full wwPDB NMR Structure Validation Report (i)

Jun 15, 2020 – 10:37 pm BST

PDB ID : 185D

Title : SEQUENCE SPECIFICITY OF QUINOXALINE ANTIBIOTICS. 1. SOLU-

TION STRUCTURE OF A 1:1 COMPLEX BETWEEN TRIOSTIN A AND [D(GACGTC)]2 AND COMPARISON WITH THE SOLUTION STRUCTURE OF THE [N-MECYS3, N-MECYS7]TANDEM-[D(GATATC)]2

COMPLEX

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Deposited on : 1994-08-10

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:

Cyrange: Kirchner and Güntert (2011)

NmrClust : Kelley et al. (1996)

MolProbity : 4.02b-467

Mogul : 1.8.5 (274361), CSD as541be (2020)

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 : 2.11

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

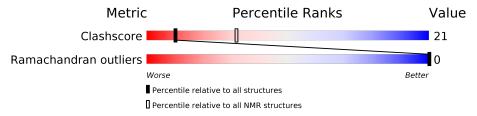
Validation Pipeline (wwPDB-VP) : 2.11

1 Overall quality at a glance (i)

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

The overall completeness of chemical shifts assignment was not calculated.

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	$egin{array}{c} ext{Whole archive} \ (\# ext{Entries}) \end{array}$	${ m NMR~archive} \ (\#{ m Entries})$
Clashscore	158937	12864
Ramachandran outliers	154571	11451

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	8	75%	25%						
2	С	6	83%	17%						
2	D	6	100%							

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA and RNA chains that are outliers for geometric criteria:

Mal	Chain	Compound	Dog	Total mo	dels with violations
IVIOI	Chain	Compound	nes	Chirality	Geometry
1	A	MVA	4	-	2
1	A	MVA	8	-	2



2 Ensemble composition and analysis (i)

This entry contains 5 models.

Cyrange was unable to find well-defined residues.

Error message: The number of core atoms (2) was below the domain threshold value (8).

NmrClust was unable to cluster the ensemble.

Error message: Wrapper check: not enough residues in core to run NmrClust



3 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 516 atoms, of which 200 are hydrogens and 0 are deuteriums.

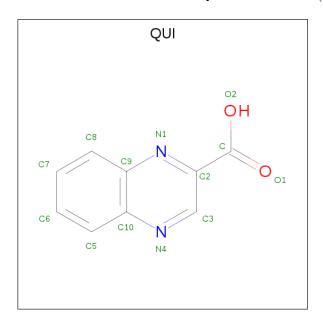
• Molecule 1 is a protein called TRIOSTIN A.

Mol	Chain	Residues		${f Atoms}$					
1	Λ	0	Total	С	Н	N	О	S	0
1	А	0	104	32	52	8	10	2	0

• Molecule 2 is a DNA chain called DNA (5'-D(*GP*AP*CP*GP*TP*C)-3').

Mol	Chain	Residues		Atoms					Trace
9	С	6	Total	С	Н	N	О	Р	0
	$\frac{2}{2}$	0	189	58	69	23	34	5	U
9	D	C	Total	С	Н	N	О	Р	0
	$\begin{array}{c c} 2 & D & \end{array}$	0	189	58	69	23	34	5	U

• Molecule 3 is 2-CARBOXYQUINOXALINE (three-letter code: QUI) (formula: C₉H₆N₂O₂).



Mol	Chain	Residues	Atoms				
3	Δ	1	Total	С	Н	N	О
)	Λ	1	17	9	5	2	1
2	Λ	1	Total	С	Н	N	О
3	A	1	17	9	5	2	1



4 Residue-property plots (i)

4.1 Average score per residue in the NMR ensemble

These plots are provided for all protein, RNA and DNA 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: TRIOSTIN A

Chain A: 75% 25%

• Molecule 2: DNA (5'-D(*GP*AP*CP*GP*TP*C)-3')

Chain C: 83% 17%

• Molecule 2: DNA (5'-D(*GP*AP*CP*GP*TP*C)-3')

Chain D: 100%

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: TRIOSTIN A

Chain A: 75% 25%



• Molecule 2: DNA (5'-D(*GP*AP*CP*GP*TP*C)-3')



Chain C:	67%	33%
611 A12 C13 614 T15 C16		
• Molecule 2	: DNA (5'-D(*GP*AP*CP*GP*TP*C)-3')	
Chain D:	17% 83%	
617 A 18 C 19 C 20 T 21 C 22		
4.2.2 Scor	e per residue for model 2	
• Molecule 1	: TRIOSTIN A	
Chain A:	75%	25%
V4 V8		
• Molecule 2	: DNA (5'-D(*GP*AP*CP*GP*TP*C)-3')	
Chain C:	17% 67%	17%
G11 C13 C13 C16 C16		
• Molecule 2	: DNA (5'-D(*GP*AP*CP*GP*TP*C)-3')	
Chain D:	83%	17%
617 A18 C19 G20 T21 C22		
4.2.3 Scor	e per residue for model 3	
• Molecule 1	: TRIOSTIN A	
Chain A:	75%	25%
N8 V8		
• Molecule 2	: DNA (5'-D(*GP*AP*CP*GP*TP*C)-3')	
Chain C:	17% 33%	50%



2	
• Molecule 2: DNA (5'-D(*GP*AP*CP*GP*TP*C)-3')	
Chain D: 100%	
414 418 619 620 622 721	
4.2.4 Score per residue for model 4	
• Molecule 1: TRIOSTIN A	
Chain A: 13% 88%	
8 X X X X X X X X X X X X X X X X X X X	
• Molecule 2: DNA (5'-D(*GP*AP*CP*GP*TP*C)-3')	
Chain C: 67%	33%
10	
• Molecule 2: DNA (5'-D(*GP*AP*CP*GP*TP*C)-3')	
Chain D: 17% 83%	
Chain D: 17% 83%	
A117 C19 C20 C22	
5 និក្សិត្តិ និង Score per residue for model 5	
4.2.5 Score per residue for model 5 • Molecule 1: TRIOSTIN A	
4.2.5 Score per residue for model 5 • Molecule 1: TRIOSTIN A Chain A: 38% 63%	
4.2.5 Score per residue for model 5 • Molecule 1: TRIOSTIN A Chain A: 38% 63%	
4.2.5 Score per residue for model 5 • Molecule 1: TRIOSTIN A Chain A: 38% 63% Molecule 2: DNA (5'-D(*GP*AP*CP*GP*TP*C)-3')	



Chain D: 67% 33%

G17 A18 C19 G20 T21 C22



Refinement protocol and experimental data overview (i) 5



The models were refined using the following method: DISTANCE GEOMETRY, MOLECULAR DYNAMICS.

Of the? calculated structures, 5 were deposited, based on the following criterion:?.

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

Software name	Classification	Version
X-PLOR	${ m refinement}$	

No chemical shift data was provided. No validations of the models with respect to experimental NMR restraints is performed at this time.



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: DSN, NCY, QUI, MVA

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	В	Sond lengths	Bond angles			
	Chain	RMSZ	#Z>5	RMSZ	#Z>5	
1	A	1.38 ± 0.10	$0{\pm}0/8~(~0.0{\pm}~0.0\%)$	1.64 ± 0.25	$0\pm0/8~(~0.0\pm~0.0\%)$	
2	С	1.37 ± 0.04	$1\pm0/134$ ($0.7\pm$ 0.0%)	2.30 ± 0.07	$14\pm 2/205~(~6.8\pm~1.0\%)$	
2	D	1.32 ± 0.01	$1\pm0/134$ ($0.7\pm$ 0.0%)	2.25 ± 0.08	$11\pm 2/205~(~5.6\pm~0.9\%)$	
All	All	1.35	10/1380 (0.7%)	2.26	127/2090 (6.1%)	

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

Mol	Chain	Res	Tuna	Atoms	\mathbf{z}	Observed (Å)	Ideal(Å)	Mod	dels
MIOI	Chain	nes	Type	Atoms	L	$Observed(\AA) \mid Ideal(\AA) \mid$		Worst	Total
2	D	21	DT	C5-C7	7.12	1.54	1.50	1	5
2	С	15	DT	C5-C7	6.97	1.54	1.50	2	5

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	Z	$Observed(^o)$	Ideal(0)	Mod	dels
MIOI	Chain	nes	Type	Atoms		Observed()	$\operatorname{Ideal}({}^o)$	Worst	Total
2	С	16	DC	O4'-C1'-N1	8.58	114.00	108.00	4	2
2	D	20	DG	O4'-C1'-C2'	-8.49	99.11	105.90	2	5
2	С	15	DT	O4'-C1'-C2'	-8.44	99.15	105.90	3	2
2	С	13	DC	O4'-C1'-N1	7.78	113.45	108.00	4	3
2	D	17	DG	P-O3'-C3'	7.68	128.92	119.70	3	1
2	С	14	DG	O4'-C1'-C2'	-7.32	100.05	105.90	2	5
2	D	17	DG	N7-C8-N9	7.08	116.64	113.10	4	5
2	С	11	DG	O4'-C1'-N9	6.90	112.83	108.00	2	2
2	D	19	DC	C1'-O4'-C4'	-6.82	103.28	110.10	1	1
2	D	20	DG	N7-C8-N9	6.76	116.48	113.10	4	5
2	С	13	DC	C4'-C3'-C2'	-6.74	97.03	103.10	4	3
2	D	19	DC	C4'-C3'-C2'	-6.73	97.04	103.10	5	2
2	С	11	DG	O4'-C1'-C2'	-6.72	100.52	105.90	4	4

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	Chain				\mathbf{z}	Observed(0)	Tdool(0)	Mod	dels
Mol	Chain	Res	Type	Atoms	L	$\operatorname{Observed}(^{o})$	$\operatorname{Ideal}({}^{o})$	Worst	Total
2	С	11	DG	N7-C8-N9	6.69	116.45	113.10	3	5
2	D	20	DG	O4'-C1'-N9	6.68	112.68	108.00	2	3
2	С	14	DG	N7-C8-N9	6.58	116.39	113.10	4	5
2	D	17	DG	C1'-O4'-C4'	-6.50	103.60	110.10	1	1
2	D	18	DA	O4'-C1'-N9	6.49	112.54	108.00	4	1
2	D	22	DC	O4'-C1'-N1	6.29	112.41	108.00	3	1
2	С	13	DC	N1-C2-O2	6.28	122.67	118.90	1	5
2	С	16	DC	C4'-C3'-C2'	-6.24	97.49	103.10	4	3
2	D	19	DC	N1-C2-O2	6.22	122.63	118.90	1	4
2	С	13	DC	C1'-O4'-C4'	-6.21	103.89	110.10	1	1
2	D	19	DC	O4'-C1'-N1	6.19	112.34	108.00	3	1
2	D	20	DG	C1'-O4'-C4'	-6.18	103.92	110.10	2	2
2	D	17	DG	O4'-C1'-C2'	-6.17	100.96	105.90	5	1
2	D	18	DA	O4'-C1'-C2'	-6.14	100.99	105.90	3	2
2	D	21	DT	O4'-C1'-C2'	-6.10	101.02	105.90	1	2
2	С	14	DG	O4'-C1'-N9	6.01	112.21	108.00	2	1
2	С	16	DC	N1-C2-O2	6.00	122.50	118.90	2	4
2	С	13	DC	O4'-C1'-C2'	-5.93	101.15	105.90	1	1
2	D	21	DT	C6-C5-C7	-5.89	119.36	122.90	1	4
2	D	22	DC	N1-C2-O2	5.81	122.39	118.90	2	3
2	D	20	DG	C8-N9-C4	-5.78	104.09	106.40	2	3
2	С	15	DT	C6-C5-C7	-5.75	119.45	122.90	1	5
2	С	11	DG	C1'-O4'-C4'	-5.68	104.42	110.10	4	3
2	С	14	DG	C8-N9-C4	-5.67	104.13	106.40	3	4
2	С	14	DG	C1'-O4'-C4'	-5.62	104.48	110.10	2	1
2	С	12	DA	N7-C8-N9	5.58	116.59	113.80	3	4
2	D	17	DG	C8-N9-C4	-5.47	104.21	106.40	2	5
2	D	18	DA	N7-C8-N9	5.43	116.52	113.80	5	4
2	С	11	DG	C8-N9-C4	-5.29	104.28	106.40	4	4
2	С	15	DT	P-O5'-C5'	5.16	129.15	120.90	5	1
2	С	12	DA	C4'-C3'-C2'	-5.11	98.50	103.10	4	1
2	D	17	DG	C4'-C3'-C2'	-5.11	98.50	103.10	4	1
2	С	11	DG	C5-N7-C8	-5.03	101.78	104.30	1	1

There are no chirality outliers.

There are no planarity outliers.

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	A	52	52	50	10±8
2	С	120	69	69	1±1
2	D	120	69	69	0±0
3	A	24	10	10	2±3
All	All	1580	1000	990	55

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 21.

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

A 4 0 mg 1	A + a ma - 2	Clash (Å)	Distance (8)	Mod	dels
Atom-1	Atom-2	$\operatorname{Clash}(ext{\AA})$	$oxed{ ext{Distance}(ext{Å}) }$	Worst	Total
1:A:4:MVA:HG12	1:A:4:MVA:HN3	0.84	1.47	4	2
1:A:8:MVA:HN3	1:A:8:MVA:HG12	0.83	1.48	4	2
1:A:8:MVA:CG2	3:A:0:QUI:H8	0.83	2.04	4	2
1:A:4:MVA:CG2	3:A:9:QUI:H8	0.81	2.06	4	2
1:A:7:NCY:HCN3	1:A:8:MVA:HN1	0.80	1.50	5	2
1:A:7:NCY:CN	1:A:8:MVA:HN1	0.76	2.11	5	2
1:A:3:NCY:HCN3	1:A:4:MVA:HN1	0.75	1.57	5	2
1:A:4:MVA:HG21	3:A:9:QUI:H8	0.74	1.58	5	2
1:A:3:NCY:CN	1:A:4:MVA:HN1	0.73	2.13	5	2
1:A:8:MVA:HG21	3:A:0:QUI:H8	0.71	1.59	5	2
1:A:7:NCY:HCN3	1:A:8:MVA:CN	0.70	2.16	5	2
1:A:3:NCY:HCN3	1:A:4:MVA:CN	0.68	2.18	5	2
2:C:13:DC:H2'	2:C:14:DG:O4'	0.67	1.89	4	2
1:A:8:MVA:CG1	1:A:8:MVA:HN3	0.66	2.21	5	2
1:A:4:MVA:HN3	1:A:4:MVA:CG1	0.62	2.19	4	2
2:C:15:DT:H6	2:C:15:DT:OP2	0.57	1.82	3	1
1:A:4:MVA:HN2	1:A:4:MVA:CG2	0.56	2.31	2	3
1:A:8:MVA:CG2	1:A:8:MVA:HN2	0.51	2.35	1	3
1:A:8:MVA:CG2	3:A:0:QUI:C8	0.50	2.87	4	1
1:A:8:MVA:CN	2:C:14:DG:N2	0.49	2.75	2	1
1:A:8:MVA:CB	3:A:0:QUI:H8	0.48	2.38	5	1
1:A:1:DSN:N	1:A:8:MVA:C	0.47	2.77	5	1
1:A:2:ALA:HB1	1:A:3:NCY:HCN1	0.46	1.88	4	1
1:A:8:MVA:CN	2:C:14:DG:H21	0.46	2.23	2	1
1:A:3:NCY:O	1:A:4:MVA:C	0.45	2.65	5	1
1:A:7:NCY:C	1:A:8:MVA:CG2	0.45	2.94	5	1
1:A:6:ALA:HB1	1:A:7:NCY:HCN1	0.44	1.90	4	1
1:A:4:MVA:C	1:A:5:DSN:N	0.43	2.81	4	1

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Atom-1	Atom-2	Clash(Å)	$\mathbf{Distance}(\mathbf{\mathring{A}})$	${f Models}$	
Atom-1	Atom-2	Clash(A)	Distance(A)	Worst	Total
1:A:8:MVA:HG21	3:A:0:QUI:C8	0.43	2.38	5	1
2:D:17:DG:C6	2:D:18:DA:C6	0.42	3.07	5	1
1:A:4:MVA:HN2	1:A:4:MVA:HG22	0.42	1.90	2	1
1:A:8:MVA:CG1	1:A:8:MVA:CN	0.42	2.97	4	1
1:A:4:MVA:CB	3:A:9:QUI:H8	0.41	2.46	5	1
1:A:3:NCY:C	1:A:4:MVA:CG2	0.41	2.99	5	1
1:A:4:MVA:CN	2:D:20:DG:N2	0.41	2.84	2	1
2:C:11:DG:C6	2:C:12:DA:C6	0.41	3.08	1	1

6.3 Torsion angles (i)

6.3.1 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 PDB entries followed by that with respect to all NMR entries. 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	2/8~(25%)	2±0 (100±0%)	0±0 (0±0%)	0±0 (0±0%)	100 100	
All	All	10/40~(25%)	10 (100%)	0 (0%)	0 (0%)	100 100	

There are no Ramachandran outliers.

6.3.2 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 PDB entries followed by that with respect to all NMR entries. 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	0	-	-	-
All	All	0	-	-	-

There are no protein residues with a non-rotameric sidechain to report.



6.3.3 RNA (i)

There are no RNA molecules in this entry.

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

6 non-standard protein/DNA/RNA residues are 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	Chain Res Link Bond len				gths
WIOI	Type	Chain	rtes	LIIIK	Counts	RMSZ	#Z>2
1	MVA	A	8	1	6,7,8	2.38 ± 0.57	$0\pm0 (3\pm6\%)$
1	NCY	A	3	1	5,6,7	2.56 ± 0.32	0±0 (4±8%)
1	MVA	A	4	1	6,7,8	2.40 ± 0.53	0±0 (3±6%)
1	NCY	A	7	1	5,6,7	2.43 ± 0.27	0±0 (4±8%)

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	Res	Link	Bond angles			
MIOI	Type	Chain	nes	LIIIK	Counts	RMSZ	#Z>2	
1	MVA	A	8	1	7,8,10	1.88 ± 0.51	0±0 (0±0%)	
1	NCY	A	3	1	5,6,8	3.49 ± 0.67	1±1 (24±14%)	
1	MVA	A	4	1	7,8,10	1.87 ± 0.54	0±0 (2±5%)	
1	NCY	A	7	1	5,6,8	3.49 ± 0.78	1±1 (24±14%)	

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
1	MVA	A	8	1	-	$0\pm0,6,8,10$	-
1	MVA	A	4	1	-	$0\pm0,6,8,10$	-
1	NCY	A	7	1	-	$0\pm0,3,6,8$	-
1	NCY	A	3	1	-	$0\pm0,3,6,8$	-

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

Mol	Chain	Pos	Tuno	Atoms	\mathbf{z}	Observed(Å)	Ideal(Å)	Models	
WIOI	Chain	res	Type	Atoms		Observed(A)	Ideal(A)	Worst	Total
1	A	8	MVA	CB-CA	7.27	1.66	1.54	4	1
1	A	4	MVA	CB-CA	7.13	1.66	1.54	4	1
1	A	3	NCY	CB-CA	5.89	1.59	1.53	4	1
1	A	7	NCY	CB-CA	5.16	1.58	1.53	5	1

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

Mol	Iol Chain Res Type		Atoms Z		$Observed(^o)$	$\operatorname{Ideal}({}^{o})$	Models		
MIOI	Chain	nes	туре	Atoms	L	Observed()	ideai()	Worst	Total
1	A	3	NCY	CA-CB-SG	8.48	123.72	114.19	4	3
1	A	7	NCY	CA-CB-SG	7.95	123.14	114.19	5	3
1	A	7	NCY	CB-CA-N	6.28	102.02	111.19	1	3
1	A	3	NCY	CB-CA-N	5.66	102.93	111.19	1	3
1	A	4	MVA	C-CA-N	5.14	128.09	110.88	4	1

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

6.5 Carbohydrates (i)

There are no carbohydrates in this entry.

6.6 Ligand geometry (i)

2 ligands are 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 Counts		Bond lengths		
			nes		Counts	RMSZ	#Z>2	
3	QUI	A	9	1	13,13,14	1.63 ± 0.01	0±0 (0±0%)	
3	QUI	A	0	1	13,13,14	1.64 ± 0.01	0±0 (0±0%)	

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.

Mal	Т	Chain	Res	Tiple	Bond angles		
MIOI	туре			LIIIK	Counts	RMSZ	#Z>2
3	QUI	A	9	1	17,17,19	2.01 ± 0.12	$0\pm 1 \ (2\pm 4\%)$
3	QUI	A	0	1	17,17,19	2.02 ± 0.09	$0\pm0 \ (2\pm2\%)$

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	QUI	A	0	1	-	$0\pm0,2,2,4$	$0\pm0,2,2,2$
Ī	3	QUI	A	9	1	-	$0\pm0,2,2,4$	$0\pm0,2,2,2$

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	Tuno	Atoms	Z	$\mathbf{Observed}(^o)$	$\mathrm{Ideal}(^o)$	Models	
MIOI			Type					Worst	Total
3	A	9	QUI	C-C2-N1	5.53	120.08	114.66	2	1
3	A	0	QUI	C-C2-N1	5.14	119.69	114.66	2	1
3	A	0	QUI	O1-C-C2	5.12	119.37	124.22	5	1
3	A	9	QUI	O1-C-C2	5.05	119.44	124.22	2	1

There are no chirality outliers.

There are no torsion outliers.



There are no ring outliers.

6.7 Other polymers (i)

There are no such molecules in this entry.

6.8 Polymer linkage issues (i)

The following chains have linkage breaks:

Mol	Chain	Number of breaks			
1	4-A	1			
1 5-A		1			

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	$oxed{ ext{ Distance (Å) } }$
5	A	4:MVA	С	5:DSN	N	2.88
4	A	4:MVA	С	5:DSN	N	2.81



7 Chemical shift validation (i)

No chemical shift data were provided

