

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

Nov 30, 2021 – 03:09 pm GMT

PDB ID : 7B46

Title : Structural basis of reactivation of oncogenic p53 mutants by a small molecule:

methylene quinuclidinone (MQ). Human wild-type p53DBD bound to DNA

and MQ: wt-DNA-MQ (P1)

Authors : Degtjarik, O.; Rozenberg, H.; Shakked, Z.

Deposited on : 2020-12-02

Resolution : 2.02 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: 1.8.4 (270009), CSD as541be (2020)

Xtriage (Phenix) : 1.13

EDS : 2.23.2buster-report : 1.1.7 (2018)

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

 $Refmac \quad : \quad 5.8.0267$ 

CCP4 : 7.1.010 (Gargrove)

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

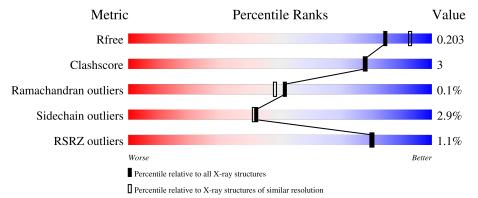
Validation Pipeline (wwPDB-VP) : 2.23.2

## 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 2.02 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	$\begin{array}{c} \textbf{Whole archive} \\ (\#\text{Entries}) \end{array}$	$\begin{array}{c} {\rm Similar\ resolution} \\ (\#{\rm Entries,\ resolution\ range(\mathring{A})}) \end{array}$
$R_{free}$	130704	10434 (2.04-2.00)
Clashscore	141614	11643 (2.04-2.00)
Ramachandran outliers	138981	11493 (2.04-2.00)
Sidechain outliers	138945	11492 (2.04-2.00)
RSRZ outliers	127900	10220 (2.04-2.00)

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 >=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	Q	uality of chain	
1	A	200		88%	8% •
1	В	200	2%	33%	12% •
1	С	200		89%	5% • 5%
1	D	200	2%	90%	6% •
2	Е	12	42%	33%	17% 8%

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Mol	Chain	Length	Quality of chain			
2	F	12	67%	17%	8%	8%
2	G	12	75%	8%	17	%
2	Н	12	75%	17	<b>'</b> %	8%



## 2 Entry composition (i)

There are 7 unique types of molecules in this entry. The entry contains 7337 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 Cellular tumor antigen p53.

Mol	Chain	Residues		Atoms				ZeroOcc	AltConf	Trace
1	Λ	192	Total	С	N	О	S	0	12	0
1	A	192	1573	971	290	290	22	0	12	0
1	B	191	Total	С	N	О	S	0	6	0
1	Ъ	191	1489	923	273	275	18	U		
1	С	190	Total	С	N	О	S	0	10	0
1		190	1519	942	278	278	21		10	0
1	1 D	D 192	Total	С	N	О	S	0	4	0
			1470	909	265	277	19	U	4	0

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

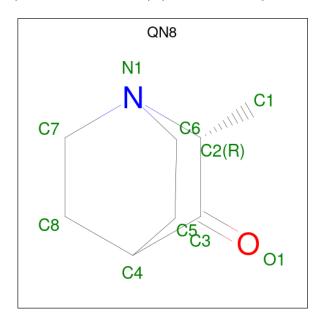
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
2	E	11	Total	С	N	О	Р	0	1	0
2	ت ا		221	105	42	64	10	Ŭ	1	
2	F	11	Total	С	N	О	Р	0	0	0
2	_ Z	11	227	106	44	66	11			
2	G	12	Total	С	N	О	Р	0	1	1
2	G	12	228	106	44	67	11	0		
2	2 H	12	Total	С	N	О	Р	0	1	0
		12	224	105	42	66	11	U		U

• Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total Zn 1 1	0	0
3	В	1	Total Zn 1 1	0	0
3	C	1	Total Zn 1 1	0	0
3	D	1	Total Zn 1 1	0	0



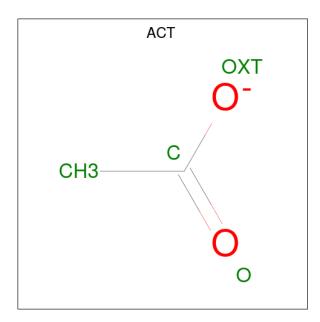
• Molecule 4 is  $(2 \{R\})$ -2-methyl-1-azabicyclo[2.2.2]octan-3-one (three-letter code: QN8) (formula:  $C_8H_{13}NO$ ) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C N O 10 8 1 1	0	1
4	A	1	Total C N O 10 8 1 1	0	1
4	A	1	Total C N O 10 8 1 1	0	1
4	В	1	Total C N O 10 8 1 1	0	1
4	С	1	Total C N O 10 8 1 1	0	1
4	С	1	Total C N O 10 8 1 1	0	1
4	D	1	Total C N O 10 8 1 1	0	1

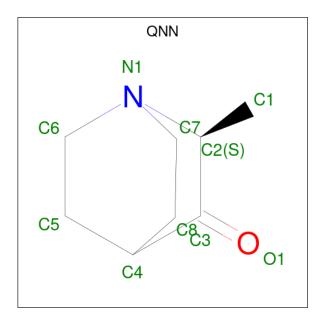
 $\bullet$  Molecule 5 is ACETATE ION (three-letter code: ACT) (formula:  $\mathrm{C_2H_3O_2}).$ 





Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	В	1	Total 4	C 2	O 2	0	0

• Molecule 6 is  $(2 \{S\})$ -2-methyl-1-azabicyclo[2.2.2]octan-3-one (three-letter code: QNN) (formula:  $C_8H_{13}NO$ ) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	D	1	Total 10	C 8	N 1	O 1	0	1

• Molecule 7 is water.



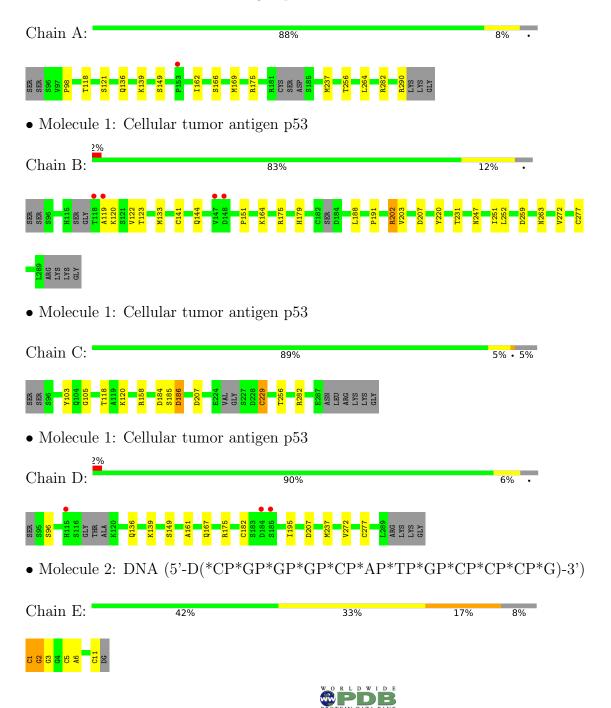
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	89	Total O 90 90	0	1
7	В	54	Total O 54 54	0	1
7	С	90	Total O 91 91	0	1
7	D	44	Total O 45 45	0	1
7	Е	4	Total O 4 4	0	0
7	F	4	Total O 4 4	0	0
7	G	4	Total O 4 4	0	0
7	Н	6	Total O 6 6	0	0

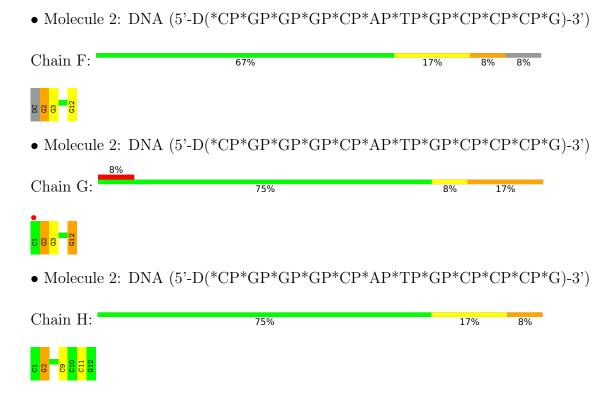


## 3 Residue-property plots (i)

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: Cellular tumor antigen p53







## 4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1	Depositor
Cell constants	54.87Å 58.31Å 78.12Å	Donositor
a, b, c, $\alpha$ , $\beta$ , $\gamma$	82.88° 87.47° 72.81°	Depositor
Resolution (Å)	47.73 - 2.02	Depositor
Resolution (A)	47.73 - 2.00	EDS
% Data completeness	93.8 (47.73-2.02)	Depositor
(in resolution range)	92.2 (47.73-2.00)	EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$< I/\sigma(I) > 1$	3.17 (at 2.00Å)	Xtriage
Refinement program	PHENIX 1.18.2-3874	Depositor
D D.	0.179 , 0.203	Depositor
$R, R_{free}$	0.179 , 0.203	DCC
$R_{free}$ test set	2000 reflections (3.41%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	32.9	Xtriage
Anisotropy	0.277	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$ , $B_{sol}(Å^2)$	(Not available), (Not available)	EDS
L-test for twinning <sup>2</sup>	$  <  L  > = 0.49, < L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	7337	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 6.16% of the height of the origin peak. No significant pseudotranslation is detected.

<sup>&</sup>lt;sup>2</sup>Theoretical values of <|L|>,  $<L^2>$  for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.



<sup>&</sup>lt;sup>1</sup>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: QN8, QNN, ZN, ACT

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	Bo	nd lengths	Bond angles		
IVIOI	Chain	RMSZ	# Z  > 5	RMSZ	# Z  > 5	
1	A	0.30	0/1607	0.49	0/2176	
1	В	0.38	$2/1521 \ (0.1\%)$	0.48	0/2064	
1	С	0.35	0/1558	0.53	1/2115 (0.0%)	
1	D	0.34	2/1503 (0.1%)	0.45	0/2043	
2	Е	1.60	$4/247 \ (1.6\%)$	1.30	4/379 (1.1%)	
2	F	1.68	6/254~(2.4%)	1.21	3/390 (0.8%)	
2	G	1.11	2/255~(0.8%)	1.11	1/393 (0.3%)	
2	Н	1.10	4/250~(1.6%)	1.02	0/385	
All	All	0.61	20/7195~(0.3%)	0.64	9/9945 (0.1%)	

The worst 5 of 20 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$\operatorname{Observed}(\text{\AA})$	$\operatorname{Ideal}(\text{\AA})$
2	Ε	11	DC	N1-C2	-15.44	1.24	1.40
2	F	2	DG	C2-N3	-15.01	1.20	1.32
2	F	2	DG	C5-C6	-13.87	1.28	1.42
2	E	11	DC	C4-C5	-13.14	1.32	1.43
2	F	2	DG	N7-C5	-7.17	1.34	1.39

The worst 5 of 9 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	$\mathbf{Z}$	$Observed(^o)$	$\operatorname{Ideal}({}^{o})$
2	Ε	11	DC	N1-C2-O2	-8.68	113.69	118.90
2	E	11	DC	N3-C2-O2	8.50	127.85	121.90
2	F	2	DG	N1-C6-O6	8.25	124.85	119.90
2	F	2	DG	N3-C2-N2	-7.20	114.86	119.90
2	Ε	11	DC	N3-C4-N4	6.56	122.59	118.00

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	A	1573	0	1504	6	0
1	В	1489	0	1392	11	0
1	С	1519	0	1450	5	0
1	D	1470	0	1354	3	0
2	Ε	221	0	124	4	0
2	F	227	0	121	2	0
2	G	228	0	123	4	0
2	Н	224	0	120	1	0
3	A	1	0	0	0	0
3	В	1	0	0	0	0
3	С	1	0	0	0	0
3	D	1	0	0	0	0
4	A	30	0	0	0	0
4	В	10	0	0	0	0
4	С	20	0	0	0	0
4	D	10	0	0	0	0
5	В	4	0	3	0	0
6	D	10	0	0	0	0
7	A	90	0	0	1	0
7	В	54	0	0	1	0
7	С	91	0	0	1	0
7	D	45	0	0	0	0
7	Е	4	0	0	0	0
7	F	4	0	0	0	0
7	G	4	0	0	0	0
7	Н	6	0	0	0	0
All	All	7337	0	6191	33	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 3.

The worst 5 of 33 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	$egin{aligned} & & & & & & & & & & & & & & & & & & &$	Clash overlap (Å)	
1:B:144:GLN:HG2	1:B:231:THR:HG22	1.77	0.66	

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Atom-1	Atom-2	$\begin{array}{c} {\rm Interatomic} \\ {\rm distance} \ ({\rm \AA}) \end{array}$	$egin{array}{c} { m Clash} \\ { m overlap} \ ({ m \AA}) \end{array}$
1:B:202[A]:ARG:HG2	7:B:452[A]:HOH:O	1.94	0.66
1:B:119:ALA:O	1:B:122:VAL:HG22	2.02	0.58
2:F:2:DG:H2"	2:F:3:DG:H5'	1.86	0.58
2:E:1[A]:DC:C5	2:G:12[A]:DG:C6	2.93	0.56

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	Perce	entiles
1	A	$200/200\ (100\%)$	198 (99%)	2 (1%)	0	100	100
1	В	191/200~(96%)	188 (98%)	3 (2%)	0	100	100
1	С	194/200 (97%)	191 (98%)	2 (1%)	1 (0%)	29	22
1	D	191/200 (96%)	190 (100%)	1 (0%)	0	100	100
All	All	776/800 (97%)	767 (99%)	8 (1%)	1 (0%)	51	48

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	С	185	SER

#### 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	175/179~(98%)	172 (98%)	3 (2%)	60 63		
1	В	158/179 (88%)	151 (96%)	7 (4%)	28 24		
1	С	169/179~(94%)	164 (97%)	5 (3%)	41 40		
1	D	157/179 (88%)	151 (96%)	6 (4%)	33 30		
All	All	$659/716 \; (92\%)$	638 (97%)	21 (3%)	42 37		

5 of 21 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	С	229[B]	CYS
1	D	167	GLN
1	D	272	VAL
1	D	182	CYS
1	D	149	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

#### 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)

Of 13 ligands modelled in this entry, 4 are monoatomic - leaving 9 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

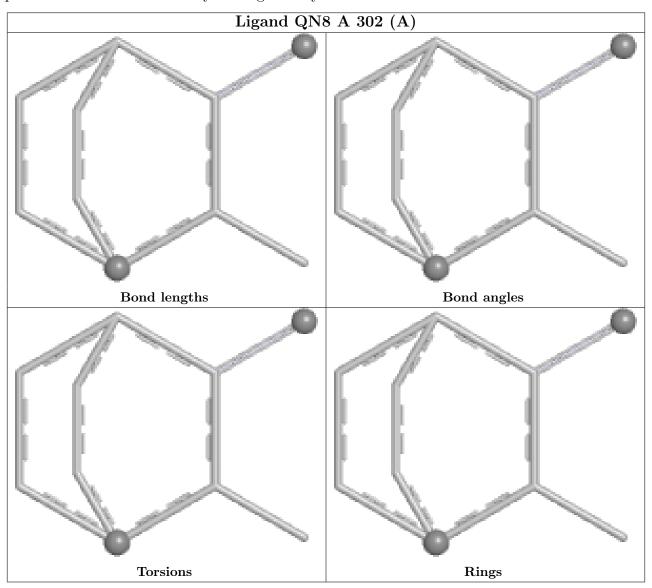
There are no torsion outliers.



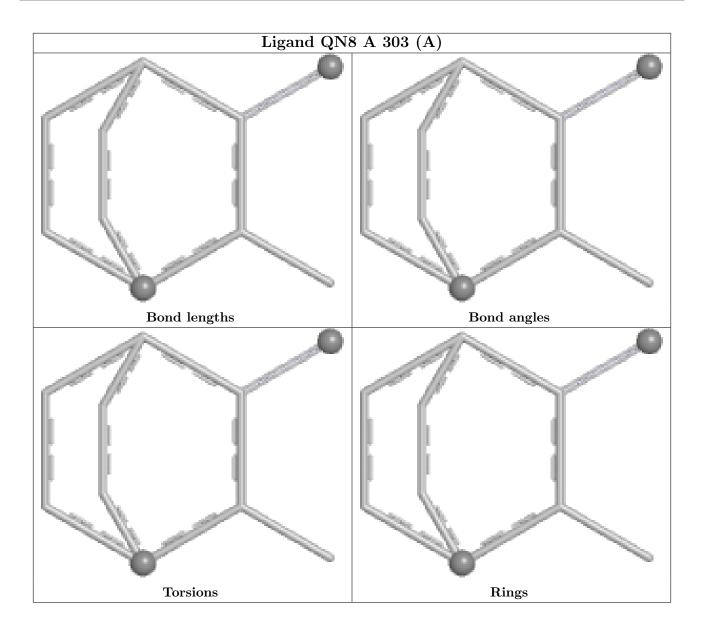
There are no ring outliers.

No monomer is involved in short contacts.

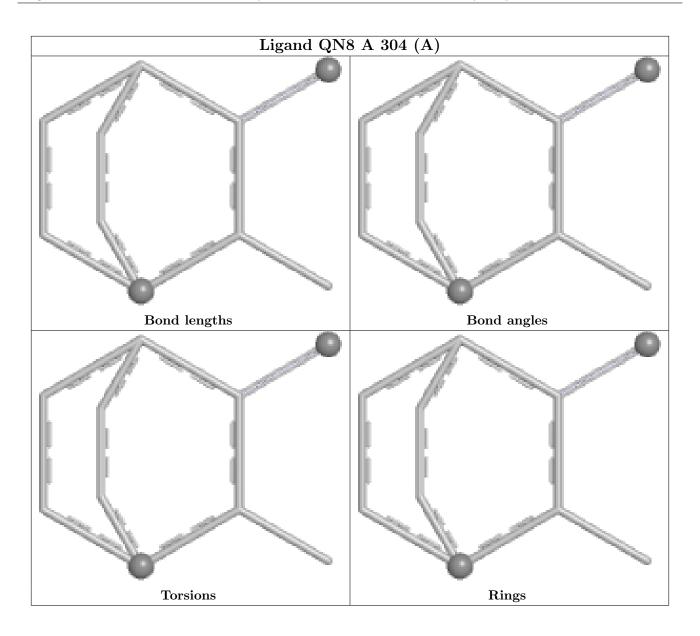
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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



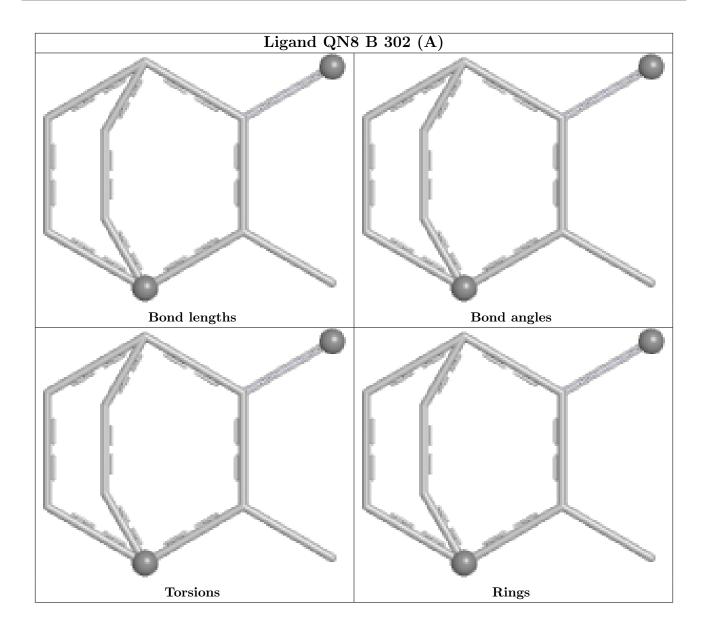




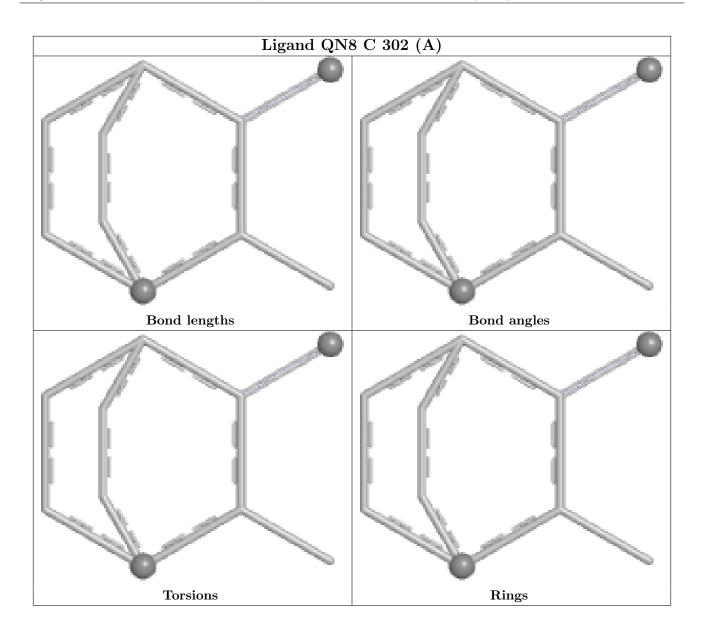




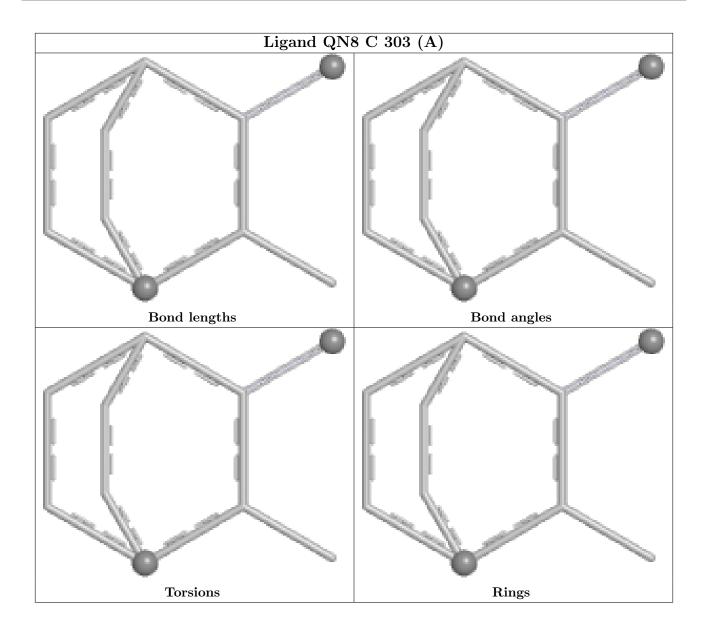




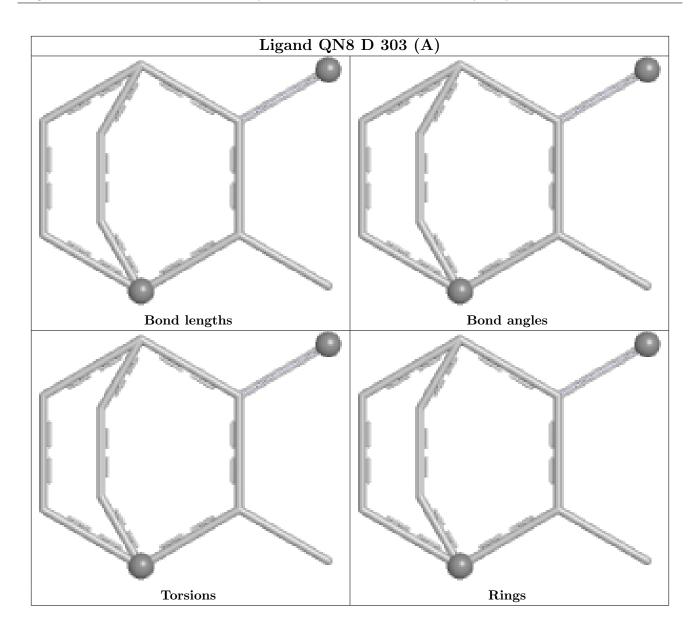




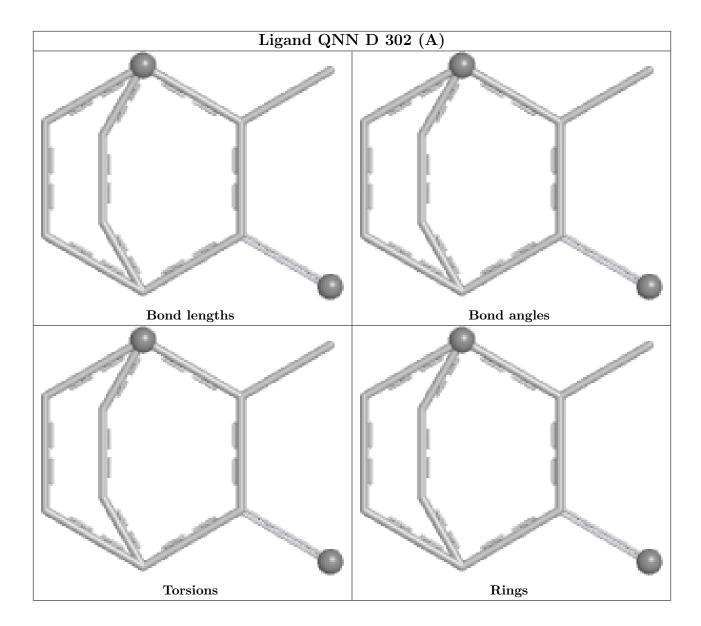












## 5.7 Other polymers (i)

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>	$\#\mathrm{RSRZ}{>}2$	$\mathbf{OWAB}(\mathbf{\mathring{A}}^2)$	Q<0.9
1	A	192/200~(96%)	-0.05	1 (0%) 91 91	23, 36, 68, 93	2 (1%)
1	В	191/200~(95%)	0.07	4 (2%) 63 63	30, 48, 79, 90	5 (2%)
1	С	190/200 (95%)	0.00	0 100 100	21, 38, 63, 98	1 (0%)
1	D	192/200 (96%)	0.02	3 (1%) 72 71	30, 48, 84, 98	1 (0%)
2	E	11/12 (91%)	0.33	0 100 100	45, 58, 90, 105	0
2	F	11/12 (91%)	0.20	0 100 100	43, 61, 105, 140	0
2	G	12/12 (100%)	0.66	1 (8%) 11 10	46, 71, 92, 100	3 (25%)
2	Н	12/12 (100%)	0.21	0 100 100	50, 70, 98, 114	0
All	All	811/848 (95%)	0.03	9 (1%) 80 80	21, 44, 79, 140	12 (1%)

The worst 5 of 9 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	В	118	THR	4.4
2	G	1	DC	3.9
1	A	153	PRO	3.7
1	D	185	SER	3.0
1	В	119	ALA	2.7

### 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.



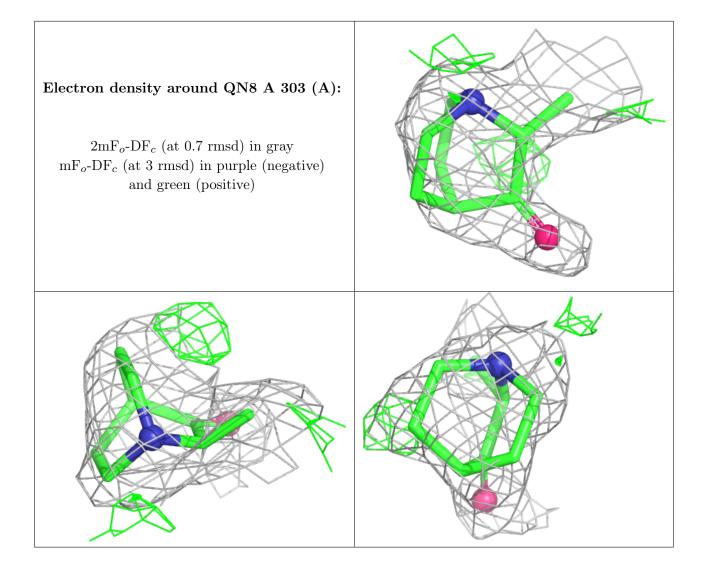
## 6.4 Ligands (i)

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,  $95^{th}$  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.

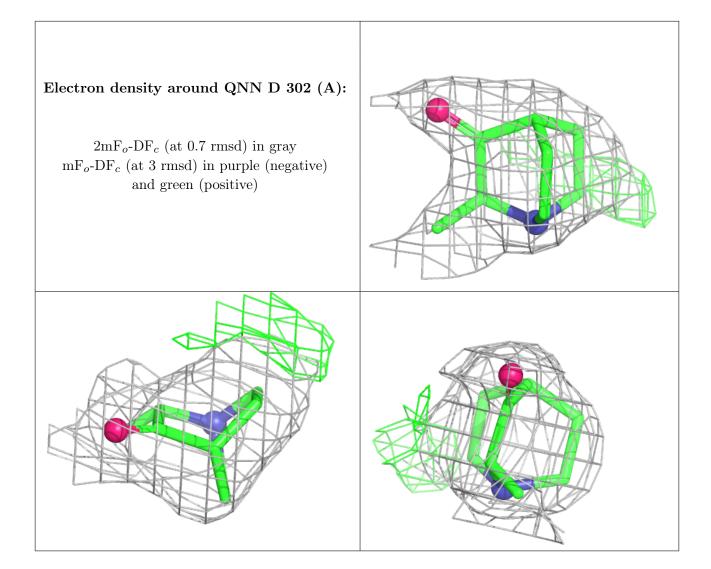
Mol	Type	Chain	Res	Atoms	RSCC	RSR	$\operatorname{B-factors}(\mathring{\mathbf{A}}^2)$	Q<0.9
4	QN8	A	303[A]	10/10	0.81	0.17	55,61,65,66	10
6	QNN	D	302[A]	10/10	0.83	0.16	60,63,70,70	10
4	QN8	С	302[A]	10/10	0.84	0.21	42,47,52,55	10
4	QN8	A	302[A]	10/10	0.90	0.14	48,51,53,53	10
4	QN8	С	303[A]	10/10	0.91	0.21	47,52,56,58	10
4	QN8	В	302[A]	10/10	0.91	0.16	56,67,72,76	10
5	ACT	В	303	4/4	0.92	0.15	44,45,52,54	4
4	QN8	A	304[A]	10/10	0.92	0.16	43,53,58,61	10
4	QN8	D	303[A]	10/10	0.93	0.15	61,67,70,70	10
3	ZN	A	301	1/1	0.98	0.10	32,32,32,32	0
3	ZN	С	301	1/1	0.98	0.11	38,38,38,38	0
3	ZN	D	301	1/1	0.98	0.09	36,36,36,36	0
3	ZN	В	301	1/1	0.99	0.08	33,33,33,33	0

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.

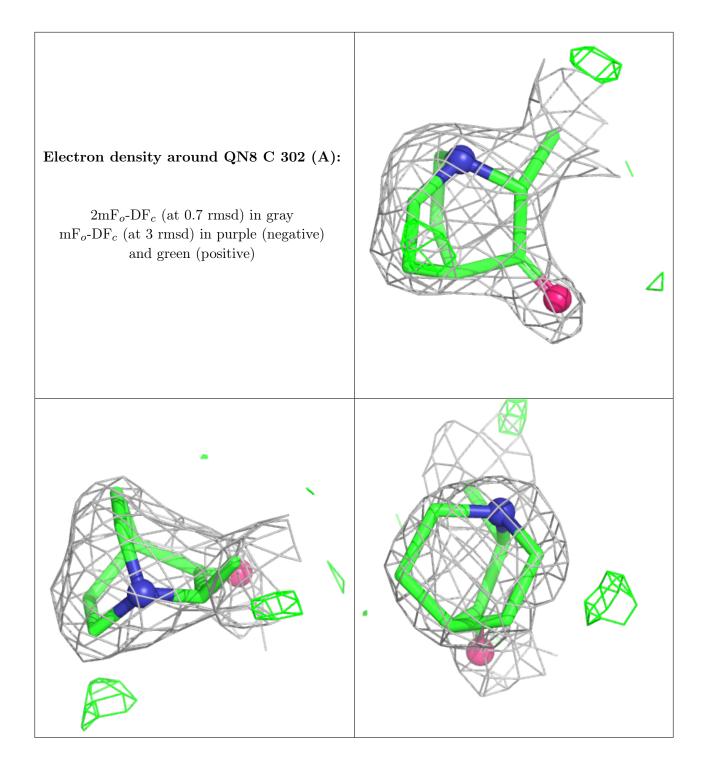




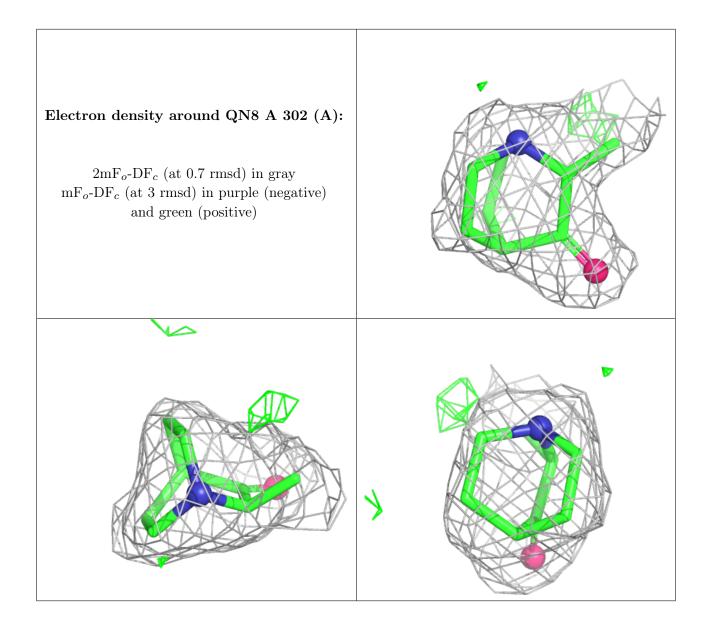




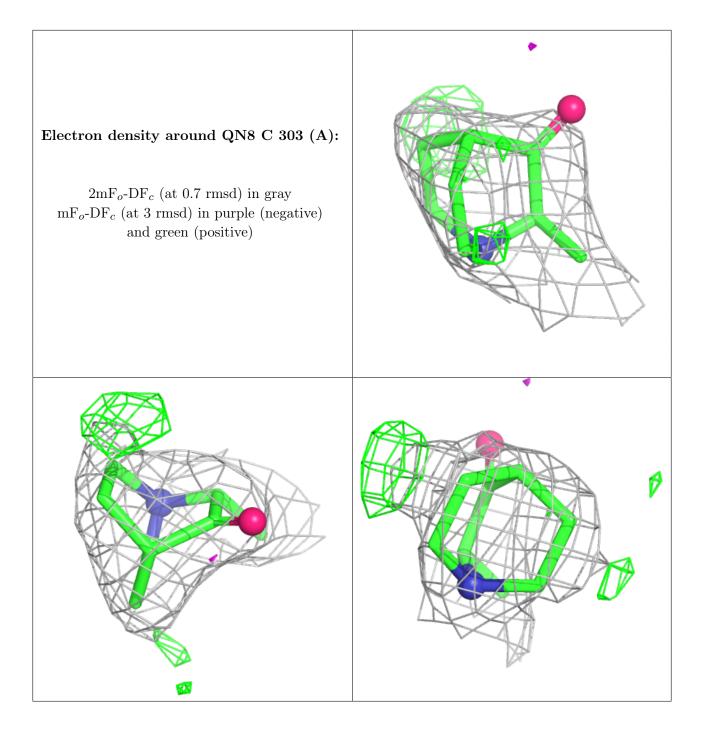




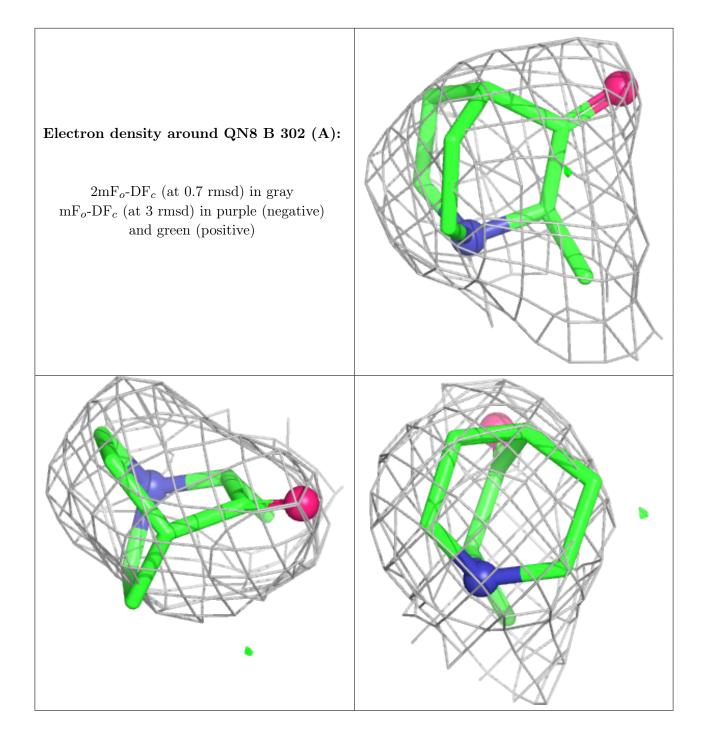








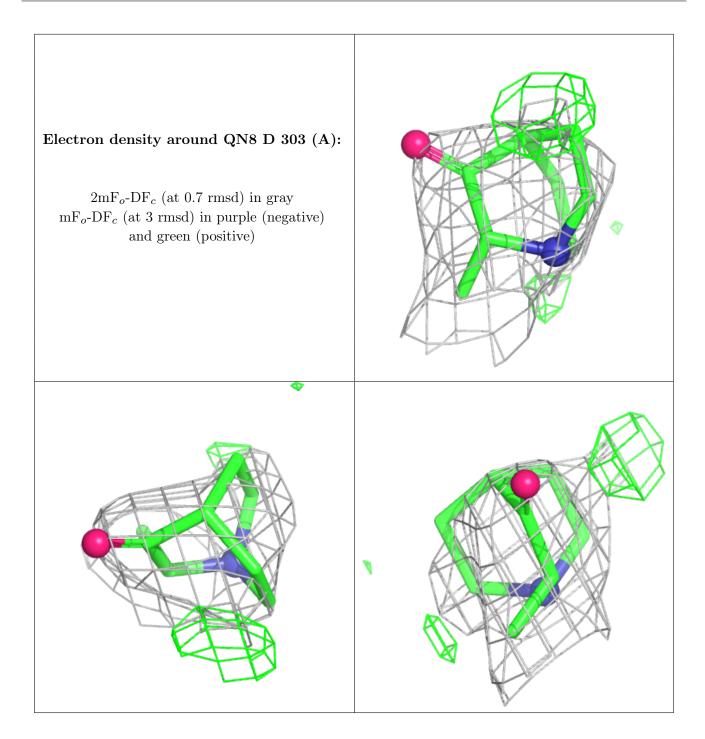






# Electron density around QN8 A 304 (A): $2 \mathrm{mF}_o\text{-}\mathrm{DF}_c$ (at 0.7 rmsd) in gray ${ m mF}_o{ m -DF}_c$ (at 3 rmsd) in purple (negative) and green (positive)





## 6.5 Other polymers (i)

There are no such residues in this entry.

