



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 09:09 PM GMT

PDB ID : 1NSA  
Title : THREE-DIMENSIONAL STRUCTURE OF PORCINE PROCAR-  
BOXYPEPTIDASE B: A STRUCTURAL BASIS OF ITS INACTIVITY  
Authors : Huber, R.  
Deposited on : 1997-07-25  
Resolution : 2.30 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](mailto:validation@mail.wwpdb.org)  
A user guide is available at  
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>  
with specific help available everywhere you see the ⓘ symbol.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : **NOT EXECUTED**  
EDS : **NOT EXECUTED**  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

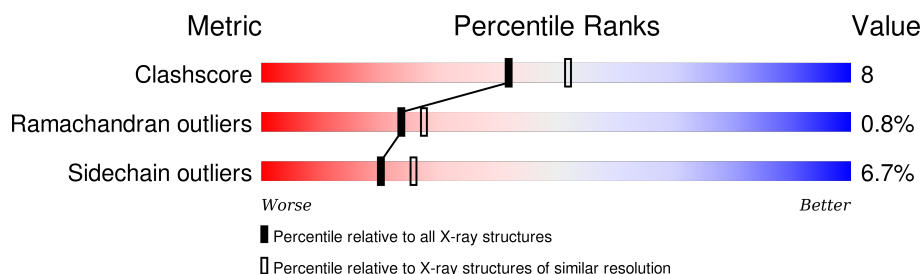
# 1 Overall quality at a glance

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 (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	102246	4452 (2.30-2.30)
Ramachandran outliers	100387	4410 (2.30-2.30)
Sidechain outliers	100360	4409 (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  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. 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  $\leq 5\%$ . 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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	395	 <div>69%27%. .</div>

## 2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 3406 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 PROCARBOXYPEPTIDASE B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	395	3174	2012	524	626	12	64	0	0

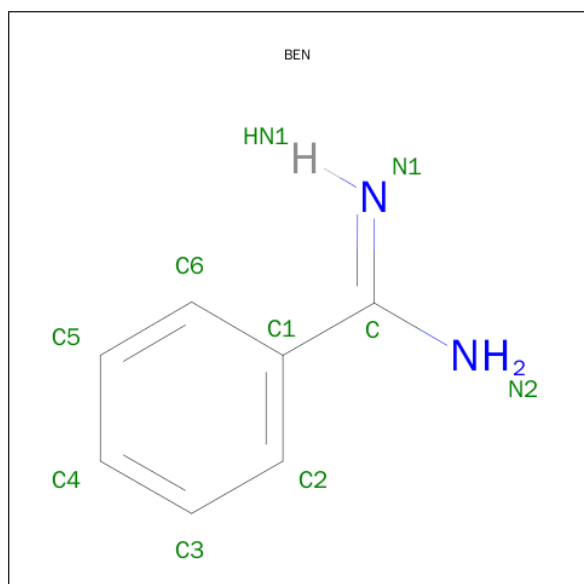
There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	93A	VAL	ARG	CONFLICT	UNP P09955
A	94A	SER	CYS	CONFLICT	UNP P09955

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Zn	0	0
			1	1		

- Molecule 3 is BENZAMIDINE (three-letter code: BEN) (formula: C<sub>7</sub>H<sub>8</sub>N<sub>2</sub>).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C N 9 7 2	0	0
3	A	1	Total C N 9 7 2	0	0
3	A	1	Total C N 9 7 2	0	0

- Molecule 4 is water.

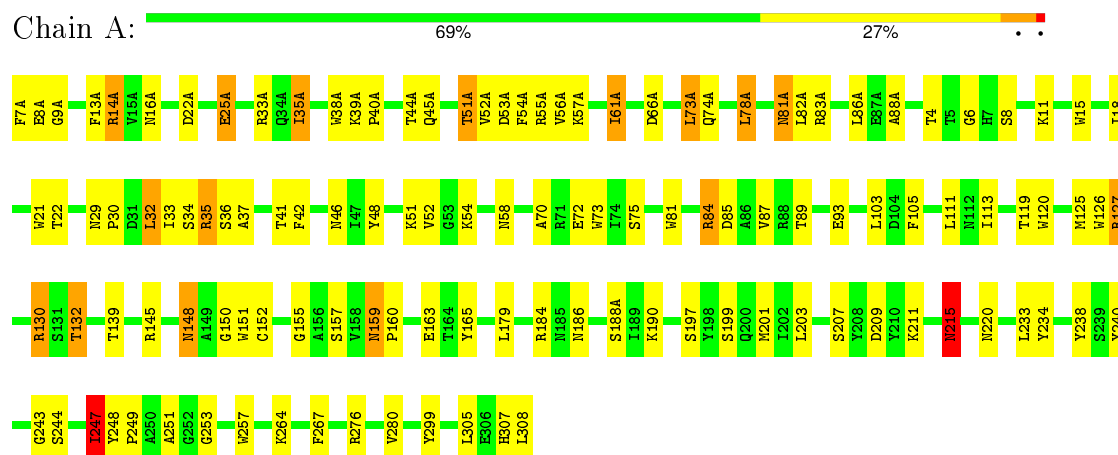
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	204	Total O 204 204	0	0

### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

Note EDS was not executed.

#### • Molecule 1: PROCARBOXYPEPTIDASE B



## 4 Data and refinement statistics

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

Property	Value	Source
Space group	P 43	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	103.10 Å   103.10 Å   46.60 Å 90.00°   90.00°   90.00°	Depositor
Resolution (Å)	8.00 – 2.30	Depositor
% Data completeness (in resolution range)	(Not available) (8.00-2.30)	Depositor
$R_{merge}$	0.09	Depositor
$R_{sym}$	(Not available)	Depositor
Refinement program	EREF	Depositor
R, $R_{free}$	0.169 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3406	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	22.0	wwPDB-VP

## 5 Model quality

### 5.1 Standard geometry

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

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	A	1.01	10/3254 (0.3%)	1.53	27/4421 (0.6%)

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	A	0	16

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	151	TRP	NE1-CE2	-8.58	1.26	1.37
1	A	120	TRP	NE1-CE2	-8.05	1.27	1.37
1	A	15	TRP	NE1-CE2	-7.53	1.27	1.37
1	A	38(A)	TRP	NE1-CE2	-7.49	1.27	1.37
1	A	21	TRP	NE1-CE2	-7.19	1.28	1.37
1	A	73	TRP	NE1-CE2	-6.11	1.29	1.37
1	A	257	TRP	NE1-CE2	-5.99	1.29	1.37
1	A	126	TRP	NE1-CE2	-5.86	1.29	1.37
1	A	81	TRP	NE1-CE2	-5.52	1.30	1.37
1	A	307	HIS	CE1-NE2	5.21	1.44	1.32

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	127	ARG	NE-CZ-NH1	16.32	128.46	120.30
1	A	14(A)	ARG	NE-CZ-NH1	15.51	128.06	120.30
1	A	14(A)	ARG	NE-CZ-NH2	-15.16	112.72	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	55(A)	ARG	NE-CZ-NH2	-12.92	113.84	120.30
1	A	130	ARG	NE-CZ-NH2	-11.49	114.56	120.30
1	A	184	ARG	NE-CZ-NH2	-10.94	114.83	120.30
1	A	127	ARG	NE-CZ-NH2	-9.84	115.38	120.30
1	A	125	MET	N-CA-CB	-8.46	95.37	110.60
1	A	55(A)	ARG	CG-CD-NE	-7.55	95.94	111.80
1	A	211	LYS	N-CA-CB	-7.13	97.77	110.60
1	A	25(A)	GLU	CB-CG-CD	-7.05	95.16	114.20
1	A	186	ASN	CA-CB-CG	-6.92	98.18	113.40
1	A	25(A)	GLU	N-CA-CB	-6.71	98.51	110.60
1	A	215	ASN	CA-CB-CG	-6.68	98.70	113.40
1	A	215	ASN	N-CA-CB	-6.11	99.61	110.60
1	A	130	ARG	NE-CZ-NH1	6.02	123.31	120.30
1	A	44(A)	THR	CA-CB-CG2	5.97	120.75	112.40
1	A	220	ASN	CA-CB-CG	-5.97	100.27	113.40
1	A	84	ARG	CD-NE-CZ	-5.90	115.34	123.60
1	A	132	THR	N-CA-CB	-5.72	99.44	110.30
1	A	66(A)	ASP	CB-CG-OD1	5.70	123.43	118.30
1	A	145	ARG	NE-CZ-NH2	5.70	123.15	120.30
1	A	35	ARG	NE-CZ-NH2	-5.59	117.51	120.30
1	A	238	TYR	CB-CG-CD2	-5.54	117.67	121.00
1	A	209	ASP	CB-CG-OD1	5.53	123.28	118.30
1	A	127	ARG	NH1-CZ-NH2	-5.12	113.77	119.40
1	A	14(A)	ARG	CD-NE-CZ	5.01	130.62	123.60

There are no chirality outliers.

All (16) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	103	LEU	Mainchain,Peptide
1	A	157	SER	Mainchain
1	A	188(A)	SER	Mainchain
1	A	215	ASN	Mainchain
1	A	240	TYR	Sidechain
1	A	247	ILE	Mainchain
1	A	248	TYR	Sidechain
1	A	264	LYS	Mainchain
1	A	267	PHE	Mainchain
1	A	299	TYR	Sidechain
1	A	37	ALA	Mainchain
1	A	51	LYS	Mainchain
1	A	74(A)	GLN	Mainchain

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Mol	Chain	Res	Type	Group
1	A	84	ARG	Sidechain
1	A	88(A)	ALA	Mainchain

## 5.2 Too-close contacts

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	3174	0	3000	50	0
2	A	1	0	0	0	0
3	A	27	0	22	3	0
4	A	204	0	0	3	0
All	All	3406	0	3022	51	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 8.

All (51) 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:8:SER:HB3	1:A:11:LYS:HG2	1.46	0.93
1:A:201:MET:HE1	1:A:247:ILE:HD11	1.68	0.76
1:A:9(A):GLY:H	1:A:81(A):ASN:HD21	1.36	0.73
1:A:14(A):ARG:HD3	1:A:53(A):ASP:OD1	1.93	0.68
1:A:40(A):PRO:HD2	1:A:52(A):VAL:HG13	1.81	0.62
1:A:39(A):LYS:HB2	4:A:552:HOH:O	2.00	0.61
1:A:56(A):VAL:HG12	1:A:61(A):ILE:CD1	2.32	0.60
1:A:72:GLU:O	1:A:75:SER:HB3	2.03	0.58
1:A:253:GLY:HA3	3:A:1:BEN:HN21	1.71	0.56
1:A:25(A):GLU:HG2	1:A:73(A):LEU:HD21	1.86	0.56
3:A:3:BEN:H6	4:A:697:HOH:O	2.07	0.55
1:A:70:ALA:HB1	1:A:119:THR:HG23	1.89	0.54
1:A:8(A):GLU:OE1	1:A:83(A):ARG:HD2	2.08	0.53
1:A:233:LEU:HD13	1:A:234:TYR:CE2	2.44	0.53
1:A:18:ILE:O	1:A:22:THR:HG23	2.09	0.51
1:A:14(A):ARG:HD2	1:A:51(A):THR:HG23	1.93	0.51
1:A:45(A):GLN:HE22	1:A:155:GLY:HA2	1.77	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:190:LYS:HG3	1:A:308:LEU:CD2	2.43	0.49
1:A:56(A):VAL:HG12	1:A:61(A):ILE:HD13	1.94	0.48
1:A:160:PRO:HA	1:A:165:TYR:CG	2.48	0.48
1:A:9(A):GLY:H	1:A:81(A):ASN:ND2	2.08	0.48
1:A:41:THR:HG21	1:A:113:ILE:HB	1.95	0.48
1:A:89:THR:HB	1:A:93:GLU:HG3	1.95	0.48
1:A:233:LEU:HB3	1:A:234:TYR:CD2	2.49	0.47
1:A:150:GLY:O	1:A:251:ALA:HB1	2.14	0.47
1:A:253:GLY:HA3	3:A:1:BEN:N2	2.29	0.47
1:A:6:GLY:O	1:A:11:LYS:HE3	2.15	0.46
1:A:14(A):ARG:HD2	1:A:51(A):THR:CG2	2.45	0.46
1:A:36:SER:O	1:A:48:TYR:HA	2.16	0.46
1:A:35(A):ILE:HG23	1:A:54(A):PHE:HB2	1.99	0.45
1:A:130:ARG:CZ	1:A:139:THR:HG21	2.47	0.45
1:A:57(A):LYS:O	1:A:61(A):ILE:HD13	2.17	0.45
1:A:207:SER:O	1:A:249:PRO:HA	2.18	0.44
1:A:32:LEU:HA	1:A:32:LEU:HD12	1.86	0.44
1:A:33:ILE:HD11	1:A:87:VAL:CG1	2.47	0.44
1:A:14(A):ARG:HB2	1:A:78(A):LEU:HD13	2.01	0.43
1:A:82(A):LEU:HA	1:A:82(A):LEU:HD12	1.78	0.43
1:A:22(A):ASP:O	1:A:25(A):GLU:HB3	2.19	0.43
1:A:33(A):ARG:HB2	1:A:35(A):ILE:HD13	2.01	0.43
1:A:29:ASN:HB2	1:A:33:ILE:HD12	2.01	0.43
1:A:148:ASN:HD22	1:A:148:ASN:HA	1.62	0.42
1:A:29:ASN:N	1:A:30:PRO:HD3	2.35	0.42
1:A:111:LEU:HD13	1:A:179:LEU:HD12	2.02	0.41
1:A:52:VAL:HB	1:A:105:PHE:HB2	2.02	0.41
1:A:127:ARG:NH2	1:A:163:GLU:O	2.53	0.41
1:A:13(A):PHE:O	1:A:53(A):ASP:HA	2.21	0.41
1:A:203:LEU:HD22	1:A:243:GLY:HA2	2.03	0.41
1:A:42:PHE:HB2	4:A:584:HOH:O	2.20	0.40
1:A:159:ASN:HD22	1:A:160:PRO:HD2	1.87	0.40
1:A:54:LYS:HD3	1:A:54:LYS:HA	1.74	0.40
1:A:7(A):PHE:HD2	1:A:86(A):LEU:HD13	1.87	0.40

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	A	393/395 (100%)	376 (96%)	14 (4%)	3 (1%)	24 27

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	199	SER
1	A	280	VAL
1	A	247	ILE

### 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	344/344 (100%)	321 (93%)	23 (7%)	20 26

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	16(A)	ASN
1	A	35(A)	ILE
1	A	51(A)	THR
1	A	61(A)	ILE
1	A	73(A)	LEU
1	A	78(A)	LEU
1	A	81(A)	ASN
1	A	4	THR

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Mol	Chain	Res	Type
1	A	32	LEU
1	A	34	SER
1	A	35	ARG
1	A	46	ASN
1	A	58	ASN
1	A	85	ASP
1	A	132	THR
1	A	148	ASN
1	A	152	CYS
1	A	159	ASN
1	A	197	SER
1	A	215	ASN
1	A	244	SER
1	A	276	ARG
1	A	305	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	45(A)	GLN
1	A	81(A)	ASN
1	A	46	ASN
1	A	102	ASN
1	A	112	ASN
1	A	148	ASN
1	A	159	ASN
1	A	185	ASN
1	A	215	ASN
1	A	220	ASN
1	A	287	GLN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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 4 ligands modelled in this entry, 1 is monoatomic - leaving 3 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	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
3	BEN	A	1	-	9,9,9	2.30	1 (11%)	9,11,11	0.70	0
3	BEN	A	2	-	9,9,9	1.54	1 (11%)	9,11,11	0.87	1 (11%)
3	BEN	A	3	-	9,9,9	1.37	3 (33%)	9,11,11	0.73	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	BEN	A	1	-	-	0/4/4/4	0/1/1/1
3	BEN	A	2	-	-	0/4/4/4	0/1/1/1
3	BEN	A	3	-	-	0/4/4/4	0/1/1/1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1	BEN	C1-C	-5.91	1.37	1.47
3	A	2	BEN	C1-C	-3.87	1.41	1.47
3	A	3	BEN	C1-C	-2.17	1.43	1.47
3	A	3	BEN	C-N2	2.18	1.38	1.33
3	A	3	BEN	C-N1	2.46	1.37	1.28

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
3	A	2	BEN	C1-C-N2	2.35	121.86	118.11

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1	BEN	2	0
3	A	3	BEN	1	0

## 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 ⓘ

### 6.1 Protein, DNA and RNA chains ⓘ

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

### 6.3 Carbohydrates ⓘ

EDS was not executed - this section will therefore be empty.

### 6.4 Ligands ⓘ

EDS was not executed - this section will therefore be empty.

### 6.5 Other polymers ⓘ

EDS was not executed - this section will therefore be empty.