



# wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 14, 2018 – 02:54 am GMT

PDB ID : 2V8H  
Title : Crystal structure of mutant E159A of beta-alanine synthase from *Saccharomyces kluyveri* in complex with its substrate N-carbamyl-beta- alanine  
Authors : Lundgren, S.; Andersen, B.; Piskur, J.; Dobritsch, D.  
Deposited on : 2007-08-08  
Resolution : 2.00 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](mailto: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 ⓘ 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 : 1.7.3 (157068), CSD as539be (2018)  
Xtriage (Phenix) : 1.13  
EDS : trunk31020  
Percentile statistics : 20171227.v01 (using entries in the PDB archive December 27th 2017)  
Refmac : 5.8.0158  
CCP4 : 7.0 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk31020

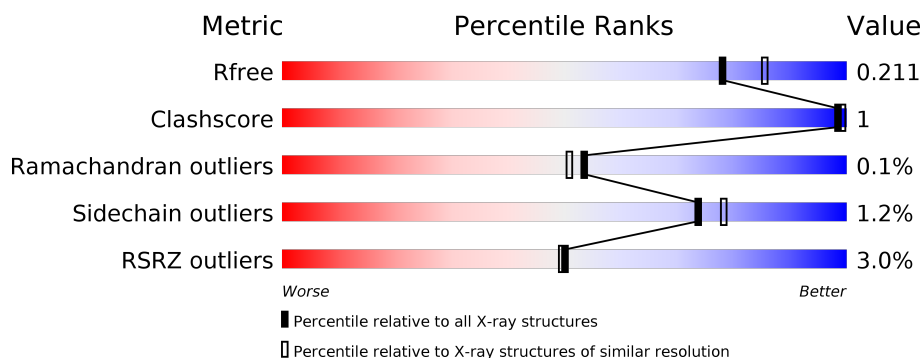
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.00 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))
$R_{free}$	111664	7193 (2.00-2.00)
Clashscore	122126	8267 (2.00-2.00)
Ramachandran outliers	120053	8166 (2.00-2.00)
Sidechain outliers	120020	8165 (2.00-2.00)
RSRZ outliers	108989	7011 (2.00-2.00)

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 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.

Mol	Chain	Length	Quality of chain
1	A	474	<div> <div>2%</div> <div> <div></div> <div>89%</div> <div>9%</div> </div> </div>
1	B	474	<div> <div>4%</div> <div> <div></div> <div>89%</div> <div>9%</div> </div> </div>
1	C	474	<div> <div>4%</div> <div> <div></div> <div>88%</div> <div>9%</div> </div> </div>
1	D	474	<div> <div>0%</div> <div> <div></div> <div>88%</div> <div>9%</div> </div> </div>

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 14629 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 BETA-ALANINE SYNTHASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	431	Total	C	N	O	S	0	9	0
			3390	2138	582	654	16			
1	B	431	Total	C	N	O	S	0	11	0
			3406	2151	586	652	17			
1	C	432	Total	C	N	O	S	0	4	0
			3370	2128	579	647	16			
1	D	431	Total	C	N	O	S	0	15	0
			3419	2163	588	652	16			

There are 84 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	159	ALA	GLU	engineered mutation	UNP Q96W94
A	456	GLN	-	expression tag	UNP Q96W94
A	457	PHE	-	expression tag	UNP Q96W94
A	458	PRO	-	expression tag	UNP Q96W94
A	459	GLY	-	expression tag	UNP Q96W94
A	460	ASP	-	expression tag	UNP Q96W94
A	461	ASP	-	expression tag	UNP Q96W94
A	462	ASP	-	expression tag	UNP Q96W94
A	463	ASP	-	expression tag	UNP Q96W94
A	464	LYS	-	expression tag	UNP Q96W94
A	465	HIS	-	expression tag	UNP Q96W94
A	466	HIS	-	expression tag	UNP Q96W94
A	467	HIS	-	expression tag	UNP Q96W94
A	468	HIS	-	expression tag	UNP Q96W94
A	469	HIS	-	expression tag	UNP Q96W94
A	470	HIS	-	expression tag	UNP Q96W94
A	471	HIS	-	expression tag	UNP Q96W94
A	472	HIS	-	expression tag	UNP Q96W94
A	473	SER	-	expression tag	UNP Q96W94
A	474	GLY	-	expression tag	UNP Q96W94
A	475	ASP	-	expression tag	UNP Q96W94

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Chain	Residue	Modelled	Actual	Comment	Reference
B	159	ALA	GLU	engineered mutation	UNP Q96W94
B	456	GLN	-	expression tag	UNP Q96W94
B	457	PHE	-	expression tag	UNP Q96W94
B	458	PRO	-	expression tag	UNP Q96W94
B	459	GLY	-	expression tag	UNP Q96W94
B	460	ASP	-	expression tag	UNP Q96W94
B	461	ASP	-	expression tag	UNP Q96W94
B	462	ASP	-	expression tag	UNP Q96W94
B	463	ASP	-	expression tag	UNP Q96W94
B	464	LYS	-	expression tag	UNP Q96W94
B	465	HIS	-	expression tag	UNP Q96W94
B	466	HIS	-	expression tag	UNP Q96W94
B	467	HIS	-	expression tag	UNP Q96W94
B	468	HIS	-	expression tag	UNP Q96W94
B	469	HIS	-	expression tag	UNP Q96W94
B	470	HIS	-	expression tag	UNP Q96W94
B	471	HIS	-	expression tag	UNP Q96W94
B	472	HIS	-	expression tag	UNP Q96W94
B	473	SER	-	expression tag	UNP Q96W94
B	474	GLY	-	expression tag	UNP Q96W94
B	475	ASP	-	expression tag	UNP Q96W94
C	159	ALA	GLU	engineered mutation	UNP Q96W94
C	456	GLN	-	expression tag	UNP Q96W94
C	457	PHE	-	expression tag	UNP Q96W94
C	458	PRO	-	expression tag	UNP Q96W94
C	459	GLY	-	expression tag	UNP Q96W94
C	460	ASP	-	expression tag	UNP Q96W94
C	461	ASP	-	expression tag	UNP Q96W94
C	462	ASP	-	expression tag	UNP Q96W94
C	463	ASP	-	expression tag	UNP Q96W94
C	464	LYS	-	expression tag	UNP Q96W94
C	465	HIS	-	expression tag	UNP Q96W94
C	466	HIS	-	expression tag	UNP Q96W94
C	467	HIS	-	expression tag	UNP Q96W94
C	468	HIS	-	expression tag	UNP Q96W94
C	469	HIS	-	expression tag	UNP Q96W94
C	470	HIS	-	expression tag	UNP Q96W94
C	471	HIS	-	expression tag	UNP Q96W94
C	472	HIS	-	expression tag	UNP Q96W94
C	473	SER	-	expression tag	UNP Q96W94
C	474	GLY	-	expression tag	UNP Q96W94
C	475	ASP	-	expression tag	UNP Q96W94

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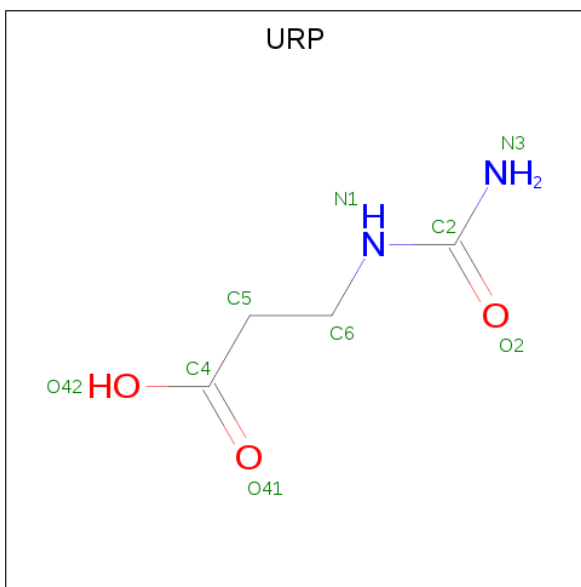
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Chain	Residue	Modelled	Actual	Comment	Reference
D	159	ALA	GLU	engineered mutation	UNP Q96W94
D	456	GLN	-	expression tag	UNP Q96W94
D	457	PHE	-	expression tag	UNP Q96W94
D	458	PRO	-	expression tag	UNP Q96W94
D	459	GLY	-	expression tag	UNP Q96W94
D	460	ASP	-	expression tag	UNP Q96W94
D	461	ASP	-	expression tag	UNP Q96W94
D	462	ASP	-	expression tag	UNP Q96W94
D	463	ASP	-	expression tag	UNP Q96W94
D	464	LYS	-	expression tag	UNP Q96W94
D	465	HIS	-	expression tag	UNP Q96W94
D	466	HIS	-	expression tag	UNP Q96W94
D	467	HIS	-	expression tag	UNP Q96W94
D	468	HIS	-	expression tag	UNP Q96W94
D	469	HIS	-	expression tag	UNP Q96W94
D	470	HIS	-	expression tag	UNP Q96W94
D	471	HIS	-	expression tag	UNP Q96W94
D	472	HIS	-	expression tag	UNP Q96W94
D	473	SER	-	expression tag	UNP Q96W94
D	474	GLY	-	expression tag	UNP Q96W94
D	475	ASP	-	expression tag	UNP Q96W94

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

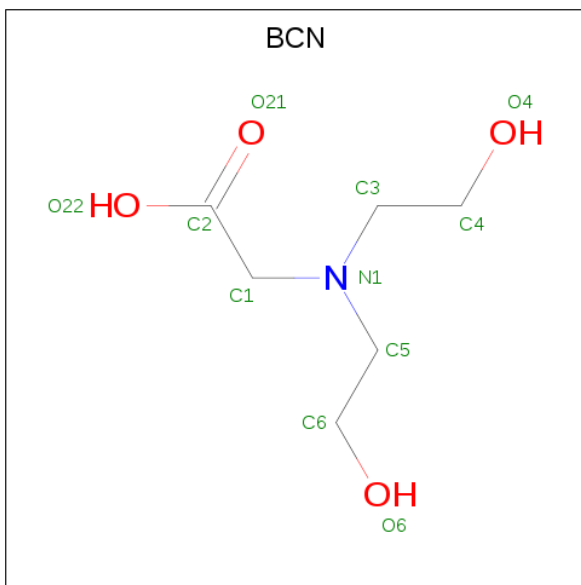
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	2	Total 2	Zn 2	0	0
2	A	2	Total 2	Zn 2	0	0
2	D	2	Total 2	Zn 2	0	0
2	C	2	Total 2	Zn 2	0	0

- Molecule 3 is N-(AMINOCARBONYL)-BETA-ALANINE (three-letter code: URP) (formula: C<sub>4</sub>H<sub>8</sub>N<sub>2</sub>O<sub>3</sub>).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			9	4	2	3		
3	B	1	Total	C	N	O	0	0
			9	4	2	3		
3	C	1	Total	C	N	O	0	0
			9	4	2	3		
3	D	1	Total	C	N	O	0	0
			9	4	2	3		

- Molecule 4 is BICINE (three-letter code: BCN) (formula:  $C_6H_{13}NO_4$ ).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			11	6	1	4		
4	B	1	Total	C	N	O	0	0
			11	6	1	4		
4	C	1	Total	C	N	O	0	0
			11	6	1	4		
4	D	1	Total	C	N	O	0	0
			11	6	1	4		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	303	Total	O	0	0
			303	303		
5	B	204	Total	O	0	0
			204	204		
5	C	180	Total	O	0	0
			180	180		
5	D	269	Total	O	0	0
			269	269		



- Molecule 1: BETA-ALANINE SYNTHASE





Q441	
	P458
	GLY
	ASP
	ASP
	ASP
	ASP
	LYS
	HIS
	HIS
	HIS
	HIS
	HIS
	HIS
	SER
	GLY
	ASP

## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	49.78Å 218.30Å 81.58Å 90.00° 92.19° 90.00°	Depositor
Resolution (Å)	19.76 – 2.00 19.76 – 2.00	Depositor EDS
% Data completeness (in resolution range)	99.6 (19.76-2.00) 99.6 (19.76-2.00)	Depositor EDS
$R_{merge}$	0.10	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.28 (at 2.01Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, $R_{free}$	0.178 , 0.208 0.181 , 0.211	Depositor DCC
$R_{free}$ test set	5792 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	21.5	Xtriage
Anisotropy	0.889	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.38 , 49.7	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.053 for h,-k,-l	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	14629	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.55% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<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.

## 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: BCN, URP, ZN

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	0.46	0/3486	0.55	0/4723
1	B	0.41	0/3517	0.52	0/4762
1	C	0.40	0/3461	0.52	0/4690
1	D	0.44	0/3542	0.54	0/4797
All	All	0.43	0/14006	0.53	0/18972

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	C	0	1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	120	GLU	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	A	3390	0	3283	4	0
1	B	3406	0	3322	3	0
1	C	3370	0	3266	6	0
1	D	3419	0	3352	5	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
2	C	2	0	0	0	0
2	D	2	0	0	0	0
3	A	9	0	7	0	0
3	B	9	0	7	0	0
3	C	9	0	7	0	0
3	D	9	0	7	0	0
4	A	11	0	12	0	0
4	B	11	0	12	0	0
4	C	11	0	12	0	0
4	D	11	0	12	0	0
5	A	303	0	0	0	0
5	B	204	0	0	0	0
5	C	180	0	0	0	1
5	D	269	0	0	1	1
All	All	14629	0	13299	17	1

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 1.

The worst 5 of 17 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:C:118:GLN:HB3	1:C:119:PRO:HD2	1.55	0.88
1:D:252:GLN:HG3	5:D:2147:HOH:O	2.01	0.59
1:C:118:GLN:HB3	1:C:119:PRO:CD	2.33	0.56
1:A:264:GLY:HA2	1:B:322:ARG:HD2	1.91	0.52
1:C:175:SER:HB2	1:C:405:HIS:CD2	2.47	0.49

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:C:2137:HOH:O	5:D:2221:HOH:O[2_656]	2.11	0.09

## 5.3 Torsion angles

### 5.3.1 Protein backbone

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	438/474 (92%)	429 (98%)	9 (2%)	0	100	100
1	B	441/474 (93%)	434 (98%)	7 (2%)	0	100	100
1	C	434/474 (92%)	427 (98%)	6 (1%)	1 (0%)	49	46
1	D	444/474 (94%)	436 (98%)	8 (2%)	0	100	100
All	All	1757/1896 (93%)	1726 (98%)	30 (2%)	1 (0%)	53	51

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	113	SER

### 5.3.2 Protein sidechains

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	365/392 (93%)	361 (99%)	4 (1%)	76	80
1	B	368/392 (94%)	362 (98%)	6 (2%)	65	70
1	C	361/392 (92%)	358 (99%)	3 (1%)	83	87
1	D	371/392 (95%)	365 (98%)	6 (2%)	65	70
All	All	1465/1568 (93%)	1446 (99%)	19 (1%)	74	76

5 of 19 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	250	ASN
1	C	134	LEU
1	D	202	LYS
1	B	134	LEU
1	D	250	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 10 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	250	ASN
1	B	387	GLN
1	C	401	GLN
1	A	441	GLN
1	C	250	ASN

### 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](#)

Of 16 ligands modelled in this entry, 8 are monoatomic - leaving 8 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	URP	A	600	2	5,8,8	0.65	0	6,9,9	0.97	0
4	BCN	A	601	-	7,10,10	0.42	0	8,11,11	0.88	0
3	URP	B	600	2	5,8,8	0.50	0	6,9,9	0.66	0
4	BCN	B	601	-	7,10,10	0.32	0	8,11,11	1.16	1 (12%)
3	URP	C	600	2	5,8,8	0.54	0	6,9,9	1.03	0
4	BCN	C	601	-	7,10,10	0.39	0	8,11,11	0.72	0
3	URP	D	600	2	5,8,8	0.48	0	6,9,9	0.69	0
4	BCN	D	601	-	7,10,10	0.30	0	8,11,11	1.51	1 (12%)

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	URP	A	600	2	-	0/4/6/6	0/0/0/0
4	BCN	A	601	-	-	0/8/10/10	0/0/0/0
3	URP	B	600	2	-	0/4/6/6	0/0/0/0
4	BCN	B	601	-	-	0/8/10/10	0/0/0/0
3	URP	C	600	2	-	0/4/6/6	0/0/0/0
4	BCN	C	601	-	-	0/8/10/10	0/0/0/0
3	URP	D	600	2	-	0/4/6/6	0/0/0/0
4	BCN	D	601	-	-	0/8/10/10	0/0/0/0

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	601	BCN	C2-C1-N1	-3.47	108.54	113.48
4	B	601	BCN	C2-C1-N1	-2.12	110.46	113.48

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

## 5.7 Other polymers ⓘ

There are no such residues in this entry.

## 5.8 Polymer linkage issues ⓘ

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<sup>th</sup> 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	OWAB(Å <sup>2</sup> )	Q<0.9
1	A	431/474 (90%)	0.19	9 (2%) 63 62	23, 27, 31, 37	0
1	B	431/474 (90%)	0.38	18 (4%) 36 35	23, 27, 30, 35	0
1	C	432/474 (91%)	0.45	19 (4%) 34 33	23, 27, 30, 37	0
1	D	431/474 (90%)	0.20	6 (1%) 75 74	24, 27, 30, 40	0
All	All	1725/1896 (90%)	0.31	52 (3%) 50 49	23, 27, 30, 40	0

The worst 5 of 52 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	192	ASP	6.5
1	D	458	PRO	5.2
1	A	458	PRO	5.1
1	C	105	GLY	5.0
1	C	193	LYS	5.0

### 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, median, 95<sup>th</sup> 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	B-factors( $\text{\AA}^2$ )	Q<0.9
4	BCN	C	601	11/11	0.66	0.18	46,47,47,48	0
4	BCN	B	601	11/11	0.68	0.20	51,52,53,53	0
4	BCN	A	601	11/11	0.69	0.22	58,61,61,61	0
4	BCN	D	601	11/11	0.78	0.20	51,51,52,52	0
3	URP	B	600	9/9	0.88	0.18	33,35,36,37	0
2	ZN	B	501	1/1	0.89	0.10	39,39,39,39	1
3	URP	C	600	9/9	0.93	0.13	33,34,35,36	0
3	URP	A	600	9/9	0.93	0.14	33,33,36,37	0
2	ZN	D	501	1/1	0.93	0.10	30,30,30,30	1
3	URP	D	600	9/9	0.95	0.16	32,33,35,36	0
2	ZN	C	500	1/1	0.97	0.07	46,46,46,46	0
2	ZN	C	501	1/1	0.98	0.16	35,35,35,35	1
2	ZN	A	500	1/1	0.98	0.09	46,46,46,46	0
2	ZN	A	501	1/1	0.98	0.11	32,32,32,32	1
2	ZN	B	500	1/1	0.99	0.17	49,49,49,49	0
2	ZN	D	500	1/1	0.99	0.08	47,47,47,47	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.