



# wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 1, 2016 – 08:19 AM GMT

PDB ID : 3E3I  
Title : H. influenzae beta-carbonic anhydrase, variant G41A with 100 mM bicarbonate  
Authors : Rowlett, R.S.; Failing, H.  
Deposited on : 2008-08-07  
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  
<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) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

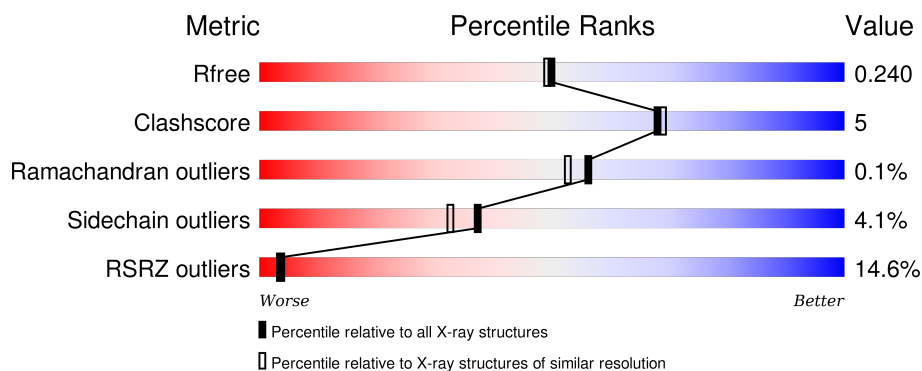
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.00 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))
$R_{free}$	91344	6249 (2.00-2.00)
Clashscore	102246	7340 (2.00-2.00)
Ramachandran outliers	100387	7248 (2.00-2.00)
Sidechain outliers	100360	7247 (2.00-2.00)
RSRZ outliers	91569	6262 (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	229	<div> <div>9%</div> <div> <div></div> <div>75%</div> <div>14%</div> <div>•</div> <div>9%</div> </div> </div>
1	B	229	<div> <div>12%</div> <div> <div></div> <div>80%</div> <div>11%</div> <div>•</div> <div>7%</div> </div> </div>
1	C	229	<div> <div>11%</div> <div> <div></div> <div>77%</div> <div>12%</div> <div>•</div> <div>9%</div> </div> </div>
1	D	229	<div> <div>15%</div> <div> <div></div> <div>83%</div> <div>10%</div> <div>•</div> <div>7%</div> </div> </div>
1	E	229	<div> <div>17%</div> <div> <div></div> <div>80%</div> <div>12%</div> <div>•</div> <div>7%</div> </div> </div>

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Mol	Chain	Length	Quality of chain
1	F	229	
1	G	229	
1	H	229	
1	I	229	
1	J	229	
1	K	229	
1	L	229	

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

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	SO4	A	231	-	-	-	X
3	SO4	H	231	-	-	-	X
3	SO4	I	231	-	-	-	X
4	BCT	B	233	-	-	X	-
4	BCT	C	232	-	-	X	X

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 21039 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 Carbonic anhydrase 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	208	Total	C	N	O	S	0	2	0
			1683	1069	300	305	9			
1	B	212	Total	C	N	O	S	0	1	0
			1711	1087	304	311	9			
1	C	208	Total	C	N	O	S	0	1	0
			1681	1067	300	305	9			
1	D	213	Total	C	N	O	S	0	1	0
			1722	1094	306	313	9			
1	E	212	Total	C	N	O	S	0	1	0
			1714	1088	305	312	9			
1	F	208	Total	C	N	O	S	0	1	0
			1680	1070	300	301	9			
1	G	214	Total	C	N	O	S	0	0	0
			1716	1090	305	312	9			
1	H	209	Total	C	N	O	S	0	1	0
			1693	1075	302	307	9			
1	I	210	Total	C	N	O	S	0	1	0
			1696	1078	302	307	9			
1	J	205	Total	C	N	O	S	0	2	0
			1669	1063	300	297	9			
1	K	209	Total	C	N	O	S	0	1	0
			1693	1075	302	307	9			
1	L	206	Total	C	N	O	S	0	0	0
			1656	1052	297	298	9			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	41	ALA	GLY	ENGINEERED	UNP P45148
B	41	ALA	GLY	ENGINEERED	UNP P45148
C	41	ALA	GLY	ENGINEERED	UNP P45148
D	41	ALA	GLY	ENGINEERED	UNP P45148
E	41	ALA	GLY	ENGINEERED	UNP P45148

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Chain	Residue	Modelled	Actual	Comment	Reference
F	41	ALA	GLY	ENGINEERED	UNP P45148
G	41	ALA	GLY	ENGINEERED	UNP P45148
H	41	ALA	GLY	ENGINEERED	UNP P45148
I	41	ALA	GLY	ENGINEERED	UNP P45148
J	41	ALA	GLY	ENGINEERED	UNP P45148
K	41	ALA	GLY	ENGINEERED	UNP P45148
L	41	ALA	GLY	ENGINEERED	UNP P45148

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

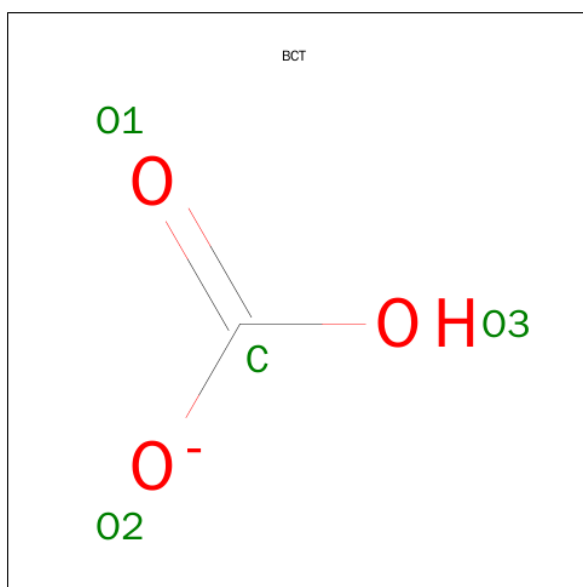
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	G	1	Total Zn 1 1	0	0
2	J	1	Total Zn 1 1	0	0
2	D	1	Total Zn 1 1	0	0
2	K	1	Total Zn 1 1	0	0
2	E	1	Total Zn 1 1	0	0
2	H	1	Total Zn 1 1	0	0
2	B	1	Total Zn 1 1	0	0
2	I	1	Total Zn 1 1	0	0
2	C	1	Total Zn 1 1	0	0
2	A	1	Total Zn 1 1	0	0
2	L	1	Total Zn 1 1	0	0
2	F	1	Total Zn 1 1	0	0

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		
3	E	1	Total	O	S	0	0
			5	4	1		
3	E	1	Total	O	S	0	0
			5	4	1		
3	H	1	Total	O	S	0	0
			5	4	1		
3	H	1	Total	O	S	0	0
			5	4	1		
3	I	1	Total	O	S	0	0
			5	4	1		
3	J	1	Total	O	S	0	0
			5	4	1		
3	K	1	Total	O	S	0	0
			5	4	1		
3	L	1	Total	O	S	0	0
			5	4	1		

- Molecule 4 is BICARBONATE ION (three-letter code: BCT) (formula:  $\text{CHO}_3$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	C	O	0	0
			4	1	3		
4	C	1	Total	C	O	0	0
			4	1	3		
4	D	1	Total	C	O	0	0
			4	1	3		
4	E	1	Total	C	O	0	0
			4	1	3		
4	F	1	Total	C	O	0	0
			4	1	3		
4	G	1	Total	C	O	0	0
			4	1	3		
4	H	1	Total	C	O	0	0
			4	1	3		
4	I	1	Total	C	O	0	0
			4	1	3		
4	J	1	Total	C	O	0	0
			4	1	3		
4	K	1	Total	C	O	0	0
			4	1	3		
4	L	1	Total	C	O	0	0
			4	1	3		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	49	Total	O	0	0
			49	49		

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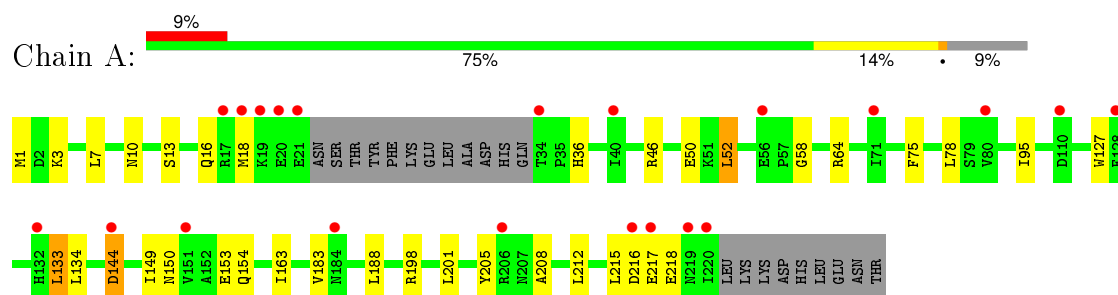
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	59	Total 59	O 59	0	0
5	C	43	Total 43	O 43	0	0
5	D	53	Total 53	O 53	0	0
5	E	65	Total 65	O 65	0	0
5	F	56	Total 56	O 56	0	0
5	G	54	Total 54	O 54	0	0
5	H	50	Total 50	O 50	0	0
5	I	46	Total 46	O 46	0	0
5	J	43	Total 43	O 43	0	0
5	K	44	Total 44	O 44	0	0
5	L	42	Total 42	O 42	0	0



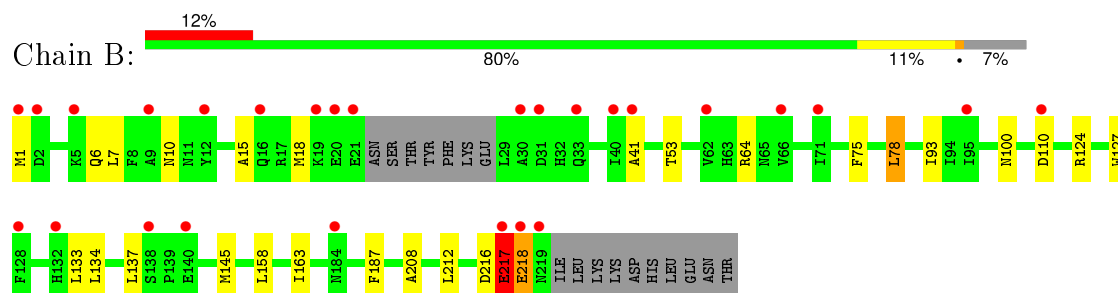
### 3 Residue-property plots [i](#)

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 ( $\text{RSRZ} > 2$ ). 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.

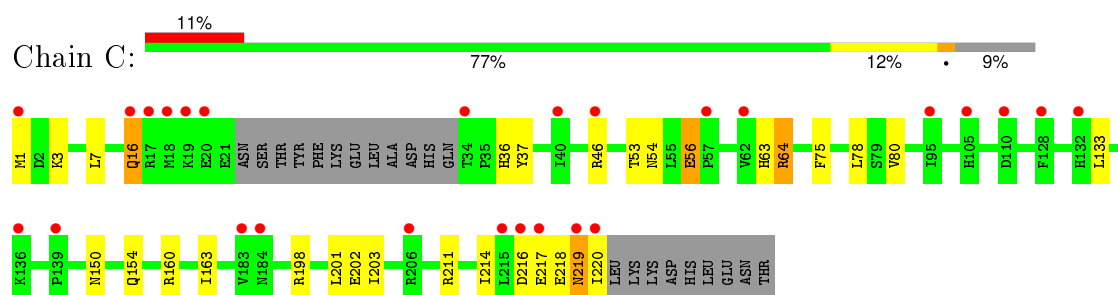
#### • Molecule 1: Carbonic anhydrase 2



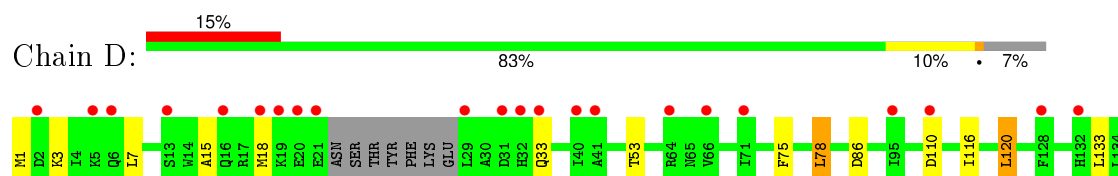
#### • Molecule 1: Carbonic anhydrase 2

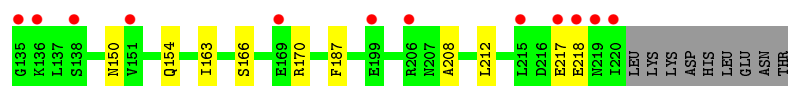


#### • Molecule 1: Carbonic anhydrase 2

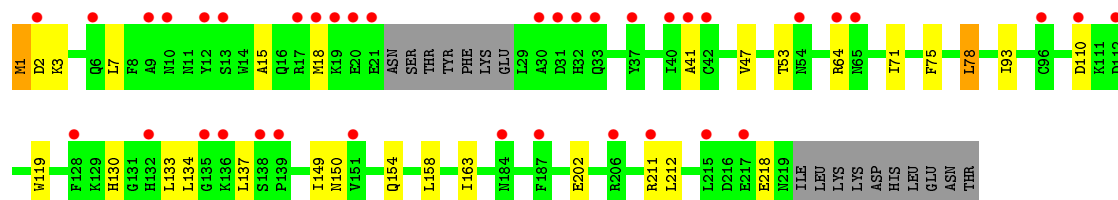
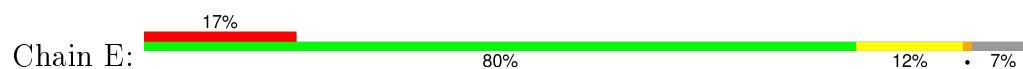


#### • Molecule 1: Carbonic anhydrase 2

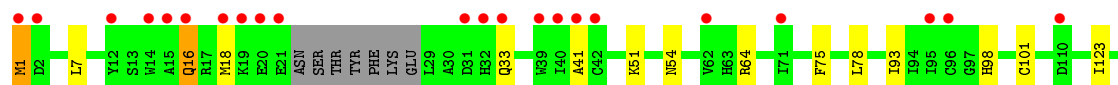
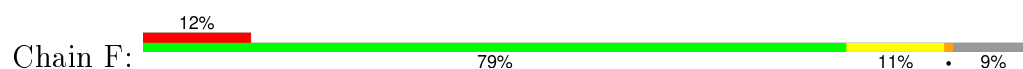




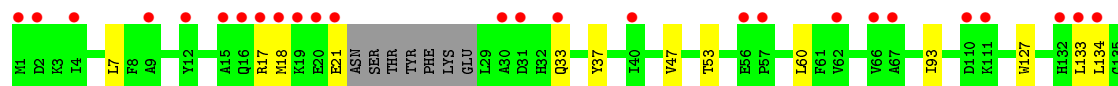
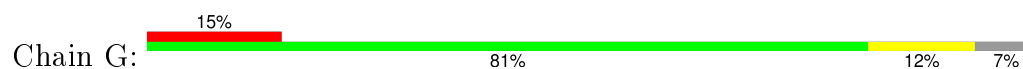
• Molecule 1: Carbonic anhydrase 2



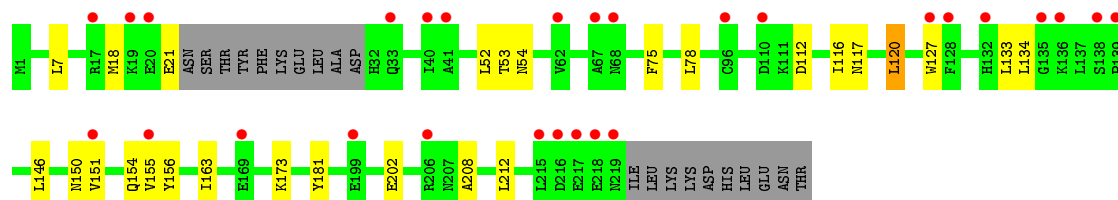
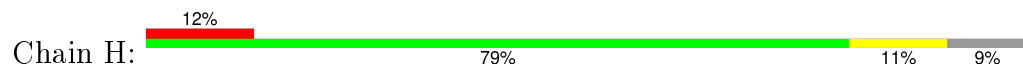
• Molecule 1: Carbonic anhydrase 2



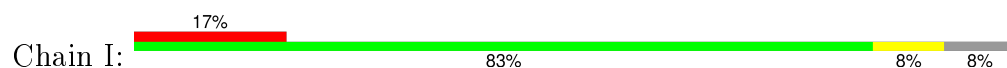
• Molecule 1: Carbonic anhydrase 2

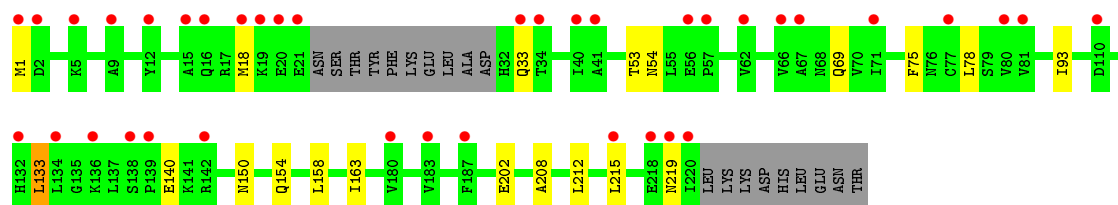


• Molecule 1: Carbonic anhydrase 2

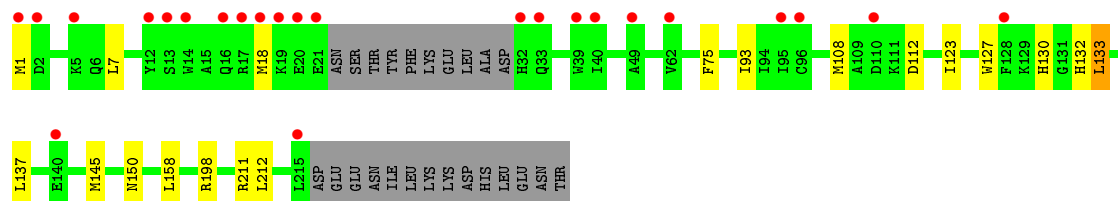
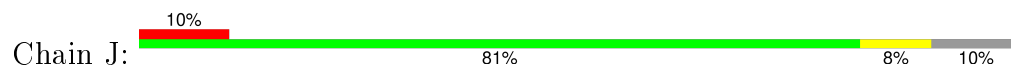


• Molecule 1: Carbonic anhydrase 2

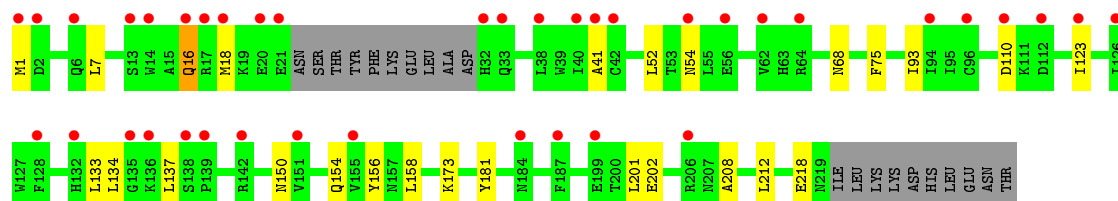
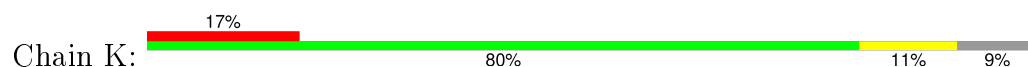




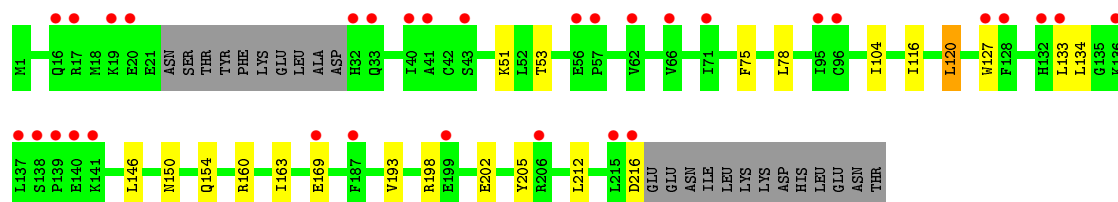
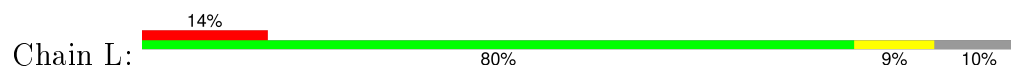
• Molecule 1: Carbonic anhydrase 2



• Molecule 1: Carbonic anhydrase 2



• Molecule 1: Carbonic anhydrase 2



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	229.59Å 144.44Å 104.89Å 90.00° 94.43° 90.00°	Depositor
Resolution (Å)	29.75 – 2.00 29.75 – 2.00	Depositor EDS
% Data completeness (in resolution range)	89.6 (29.75-2.00) 89.6 (29.75-2.00)	Depositor EDS
$R_{merge}$	0.06	Depositor
$R_{sym}$	0.06	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.68 (at 2.00Å)	Xtriage
Refinement program	REFMAC 5.2	Depositor
R, $R_{free}$	0.201 , 0.236 0.207 , 0.240	Depositor DCC
$R_{free}$ test set	10347 reflections (5.30%)	DCC
Wilson B-factor (Å <sup>2</sup> )	32.2	Xtriage
Anisotropy	0.043	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.39 , 58.3	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	5 of 205402 reflections (0.002%)	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	21039	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 87.84 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 5.9966e-08. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

<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, 0.375 respectively for untwinned datasets, and 0.333, 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: BCT, ZN, SO4

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	0.54	0/1725	0.66	2/2336 (0.1%)
1	B	0.51	0/1751	0.62	0/2371
1	C	0.52	0/1717	0.67	3/2325 (0.1%)
1	D	0.50	0/1759	0.59	0/2382
1	E	0.56	0/1751	0.64	0/2371
1	F	0.54	0/1717	0.63	1/2325 (0.0%)
1	G	0.57	0/1752	0.65	2/2373 (0.1%)
1	H	0.53	0/1730	0.62	0/2342
1	I	0.52	0/1733	0.58	0/2346
1	J	0.49	0/1707	0.59	0/2311
1	K	0.50	0/1730	0.59	0/2342
1	L	0.48	0/1692	0.62	1/2291 (0.0%)
All	All	0.52	0/20764	0.62	9/28115 (0.0%)

There are no bond length outliers.

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

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	52	LEU	CA-CB-CG	8.34	134.47	115.30
1	G	170	ARG	NE-CZ-NH2	-6.66	116.97	120.30
1	F	160	ARG	NE-CZ-NH2	-6.22	117.19	120.30
1	C	64	ARG	NE-CZ-NH1	6.20	123.40	120.30
1	C	64	ARG	NE-CZ-NH2	-6.16	117.22	120.30

There are no chirality outliers.

There are no planarity outliers.

## 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	1683	0	1672	29	0
1	B	1711	0	1692	16	0
1	C	1681	0	1665	36	0
1	D	1722	0	1702	15	0
1	E	1714	0	1691	18	0
1	F	1680	0	1669	17	0
1	G	1716	0	1701	18	0
1	H	1693	0	1671	18	0
1	I	1696	0	1671	17	0
1	J	1669	0	1655	11	0
1	K	1693	0	1671	14	0
1	L	1656	0	1645	10	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
2	E	1	0	0	0	0
2	F	1	0	0	0	0
2	G	1	0	0	0	0
2	H	1	0	0	0	0
2	I	1	0	0	0	0
2	J	1	0	0	0	0
2	K	1	0	0	0	0
2	L	1	0	0	0	0
3	A	5	0	0	1	0
3	B	10	0	0	1	0
3	C	5	0	0	0	0
3	D	5	0	0	0	0
3	E	10	0	0	0	0
3	H	10	0	0	0	0
3	I	5	0	0	0	0
3	J	5	0	0	0	0
3	K	5	0	0	0	0
3	L	5	0	0	0	0
4	B	4	0	0	2	0
4	C	4	0	0	3	0
4	D	4	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	E	4	0	0	1	0
4	F	4	0	0	0	0
4	G	4	0	0	1	0
4	H	4	0	0	0	0
4	I	4	0	0	0	0
4	J	4	0	0	0	0
4	K	4	0	0	1	0
4	L	4	0	0	0	0
5	A	49	0	0	1	0
5	B	59	0	0	0	0
5	C	43	0	0	3	0
5	D	53	0	0	0	0
5	E	65	0	0	0	0
5	F	56	0	0	0	0
5	G	54	0	0	0	0
5	H	50	0	0	0	0
5	I	46	0	0	1	0
5	J	43	0	0	1	0
5	K	44	0	0	0	0
5	L	42	0	0	0	0
All	All	21039	0	20105	189	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 5.

The worst 5 of 189 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:I:133:LEU:HD21	1:I:215:LEU:HD21	1.34	1.09
1:C:203:ILE:HD12	1:F:214:ILE:HD11	1.37	1.04
1:C:214:ILE:HD11	1:F:203:ILE:HD12	1.35	1.03
1:H:78:LEU:HD12	1:H:163:ILE:HD12	1.43	1.00
1:B:53:THR:HG22	1:D:7:LEU:HD21	1.46	0.97

There are no symmetry-related clashes.

## 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	206/229 (90%)	200 (97%)	5 (2%)	1 (0%)	34	26
1	B	209/229 (91%)	203 (97%)	5 (2%)	1 (0%)	34	26
1	C	205/229 (90%)	200 (98%)	5 (2%)	0	100	100
1	D	210/229 (92%)	205 (98%)	5 (2%)	0	100	100
1	E	209/229 (91%)	204 (98%)	5 (2%)	0	100	100
1	F	205/229 (90%)	201 (98%)	4 (2%)	0	100	100
1	G	210/229 (92%)	204 (97%)	6 (3%)	0	100	100
1	H	206/229 (90%)	201 (98%)	5 (2%)	0	100	100
1	I	207/229 (90%)	199 (96%)	7 (3%)	1 (0%)	34	26
1	J	203/229 (89%)	196 (97%)	7 (3%)	0	100	100
1	K	206/229 (90%)	202 (98%)	4 (2%)	0	100	100
1	L	202/229 (88%)	199 (98%)	3 (2%)	0	100	100
All	All	2478/2748 (90%)	2414 (97%)	61 (2%)	3 (0%)	56	53

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	217	GLU
1	B	217	GLU
1	I	219	ASN

### 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	183/201 (91%)	177 (97%)	6 (3%)	45	43
1	B	185/201 (92%)	176 (95%)	9 (5%)	31	25
1	C	182/201 (90%)	175 (96%)	7 (4%)	40	36
1	D	186/201 (92%)	178 (96%)	8 (4%)	35	30
1	E	185/201 (92%)	176 (95%)	9 (5%)	31	25
1	F	181/201 (90%)	170 (94%)	11 (6%)	23	17
1	G	185/201 (92%)	180 (97%)	5 (3%)	52	52
1	H	183/201 (91%)	178 (97%)	5 (3%)	52	52
1	I	182/201 (90%)	176 (97%)	6 (3%)	45	43
1	J	180/201 (90%)	172 (96%)	8 (4%)	35	30
1	K	183/201 (91%)	174 (95%)	9 (5%)	31	25
1	L	179/201 (89%)	173 (97%)	6 (3%)	44	41
All	All	2194/2412 (91%)	2105 (96%)	89 (4%)	37	32

5 of 89 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	F	16	GLN
1	F	211	ARG
1	K	218	GLU
1	F	18	MET
1	F	75	PHE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 25 such sidechains are listed below:

Mol	Chain	Res	Type
1	E	219	ASN
1	F	33	GLN
1	K	54	ASN
1	F	16	GLN
1	F	54	ASN

### 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 ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

Of 36 ligands modelled in this entry, 12 are monoatomic - leaving 24 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	SO4	A	231	-	4,4,4	0.38	0	6,6,6	0.20	0
3	SO4	B	231	-	4,4,4	0.37	0	6,6,6	0.35	0
3	SO4	B	232	-	4,4,4	0.45	0	6,6,6	0.37	0
4	BCT	B	233	-	0,3,3	0.00	-	0,3,3	0.00	-
3	SO4	C	231	-	4,4,4	0.34	0	6,6,6	0.25	0
4	BCT	C	232	-	0,3,3	0.00	-	0,3,3	0.00	-
3	SO4	D	231	-	4,4,4	0.39	0	6,6,6	0.46	0
4	BCT	D	232	-	0,3,3	0.00	-	0,3,3	0.00	-
3	SO4	E	231	-	4,4,4	0.27	0	6,6,6	0.61	0
3	SO4	E	232	-	4,4,4	0.22	0	6,6,6	0.39	0
4	BCT	E	233	-	0,3,3	0.00	-	0,3,3	0.00	-
4	BCT	F	231	-	0,3,3	0.00	-	0,3,3	0.00	-
4	BCT	G	231	-	0,3,3	0.00	-	0,3,3	0.00	-
3	SO4	H	231	-	4,4,4	0.19	0	6,6,6	0.35	0
3	SO4	H	232	-	4,4,4	0.18	0	6,6,6	0.23	0
4	BCT	H	233	-	0,3,3	0.00	-	0,3,3	0.00	-
3	SO4	I	231	-	4,4,4	0.31	0	6,6,6	0.59	0
4	BCT	I	232	-	0,3,3	0.00	-	0,3,3	0.00	-
3	SO4	J	231	-	4,4,4	0.27	0	6,6,6	0.59	0
4	BCT	J	232	-	0,3,3	0.00	-	0,3,3	0.00	-
3	SO4	K	231	-	4,4,4	0.21	0	6,6,6	0.23	0
4	BCT	K	232	-	0,3,3	0.00	-	0,3,3	0.00	-

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
3	SO4	L	231	-	4,4,4	0.20	0	6,6,6	0.28	0
4	BCT	L	232	-	0,3,3	0.00	-	0,3,3	0.00	-

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	SO4	A	231	-	-	0/0/0/0	0/0/0/0
3	SO4	B	231	-	-	0/0/0/0	0/0/0/0
3	SO4	B	232	-	-	0/0/0/0	0/0/0/0
4	BCT	B	233	-	-	0/0/0/0	0/0/0/0
3	SO4	C	231	-	-	0/0/0/0	0/0/0/0
4	BCT	C	232	-	-	0/0/0/0	0/0/0/0
3	SO4	D	231	-	-	0/0/0/0	0/0/0/0
4	BCT	D	232	-	-	0/0/0/0	0/0/0/0
3	SO4	E	231	-	-	0/0/0/0	0/0/0/0
3	SO4	E	232	-	-	0/0/0/0	0/0/0/0
4	BCT	E	233	-	-	0/0/0/0	0/0/0/0
4	BCT	F	231	-	-	0/0/0/0	0/0/0/0
4	BCT	G	231	-	-	0/0/0/0	0/0/0/0
3	SO4	H	231	-	-	0/0/0/0	0/0/0/0
3	SO4	H	232	-	-	0/0/0/0	0/0/0/0
4	BCT	H	233	-	-	0/0/0/0	0/0/0/0
3	SO4	I	231	-	-	0/0/0/0	0/0/0/0
4	BCT	I	232	-	-	0/0/0/0	0/0/0/0
3	SO4	J	231	-	-	0/0/0/0	0/0/0/0
4	BCT	J	232	-	-	0/0/0/0	0/0/0/0
3	SO4	K	231	-	-	0/0/0/0	0/0/0/0
4	BCT	K	232	-	-	0/0/0/0	0/0/0/0
3	SO4	L	231	-	-	0/0/0/0	0/0/0/0
4	BCT	L	232	-	-	0/0/0/0	0/0/0/0

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.

7 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	231	SO4	1	0
3	B	232	SO4	1	0
4	B	233	BCT	2	0
4	C	232	BCT	3	0
4	E	233	BCT	1	0
4	G	231	BCT	1	0
4	K	232	BCT	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 [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	208/229 (90%)	0.66	21 (10%) 9 10	28, 34, 48, 57	0
1	B	212/229 (92%)	0.74	27 (12%) 5 5	27, 33, 46, 51	0
1	C	208/229 (90%)	0.70	26 (12%) 5 5	27, 33, 48, 54	0
1	D	213/229 (93%)	0.71	34 (15%) 3 3	27, 33, 45, 52	0
1	E	212/229 (92%)	0.89	38 (17%) 2 2	27, 32, 45, 48	0
1	F	208/229 (90%)	0.75	27 (12%) 5 5	27, 32, 44, 51	0
1	G	214/229 (93%)	0.99	34 (15%) 3 3	27, 34, 46, 49	0
1	H	209/229 (91%)	0.77	28 (13%) 4 5	28, 34, 46, 60	0
1	I	210/229 (91%)	1.06	38 (18%) 2 2	30, 37, 50, 54	0
1	J	205/229 (89%)	0.72	24 (11%) 6 7	28, 35, 47, 53	0
1	K	209/229 (91%)	1.20	39 (18%) 2 2	32, 38, 49, 55	0
1	L	206/229 (89%)	0.81	32 (15%) 3 3	30, 36, 47, 50	0
All	All	2514/2748 (91%)	0.83	368 (14%) 3 4	27, 34, 47, 60	0

The worst 5 of 368 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	20	GLU	8.8
1	K	135	GLY	7.2
1	J	20	GLU	6.5
1	I	20	GLU	6.4
1	C	219	ASN	6.3

### 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 ⓘ

There are no carbohydrates in this entry.

## 6.4 Ligands ⓘ

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
4	BCT	C	232	4/4	0.92	0.38	5.46	18,19,19,20	4
3	SO4	H	231	5/5	0.81	0.28	3.38	71,73,74,75	0
3	SO4	I	231	5/5	0.83	0.23	2.82	73,74,75,76	0
3	SO4	A	231	5/5	0.95	0.25	2.57	38,38,39,39	5
3	SO4	K	231	5/5	0.61	0.28	1.41	89,91,91,91	0
3	SO4	E	232	5/5	0.76	0.26	0.94	73,74,75,76	0
3	SO4	H	232	5/5	0.82	0.21	0.48	75,78,78,79	0
3	SO4	L	231	5/5	0.80	0.18	0.10	75,77,77,78	0
3	SO4	D	231	5/5	0.93	0.18	-0.03	67,69,70,70	0
3	SO4	J	231	5/5	0.92	0.15	-0.13	72,72,74,74	0
4	BCT	G	231	4/4	0.89	0.14	-0.14	40,40,40,40	0
3	SO4	E	231	5/5	0.91	0.16	-0.19	64,67,68,68	0
3	SO4	B	231	5/5	0.96	0.13	-0.55	57,59,60,60	0
4	BCT	B	233	4/4	0.96	0.13	-0.77	45,46,46,46	0
3	SO4	C	231	5/5	0.97	0.10	-1.13	60,62,62,62	0
4	BCT	I	232	4/4	0.96	0.10	-1.31	43,44,44,45	0
2	ZN	G	230	1/1	0.98	0.05	-1.65	31,31,31,31	0
4	BCT	D	232	4/4	0.95	0.10	-1.66	41,41,41,42	0
4	BCT	L	232	4/4	0.96	0.08	-1.71	41,41,41,41	0
4	BCT	E	233	4/4	0.97	0.08	-1.75	38,38,38,39	0
3	SO4	B	232	5/5	0.98	0.07	-1.77	55,55,55,56	0
4	BCT	J	232	4/4	0.97	0.07	-1.86	41,41,42,42	0
4	BCT	F	231	4/4	0.98	0.06	-2.02	38,38,38,38	0
4	BCT	H	233	4/4	0.98	0.06	-2.16	38,38,38,39	0
2	ZN	A	230	1/1	0.97	0.04	-2.22	31,31,31,31	0
2	ZN	K	230	1/1	0.98	0.05	-2.63	36,36,36,36	0
4	BCT	K	232	4/4	0.95	0.09	-2.79	37,38,38,38	0
2	ZN	H	230	1/1	0.99	0.03	-2.96	32,32,32,32	0
2	ZN	E	230	1/1	0.99	0.04	-3.05	28,28,28,28	0
2	ZN	I	230	1/1	0.95	0.04	-3.12	36,36,36,36	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
2	ZN	F	230	1/1	0.99	0.03	-3.59	31,31,31,31	0
2	ZN	J	230	1/1	0.99	0.03	-3.72	33,33,33,33	0
2	ZN	C	230	1/1	1.00	0.02	-3.88	32,32,32,32	0
2	ZN	L	230	1/1	0.99	0.02	-4.72	32,32,32,32	0
2	ZN	D	230	1/1	0.99	0.03	-6.44	30,30,30,30	0
2	ZN	B	230	1/1	0.99	0.02	-7.26	31,31,31,31	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.