



# Full wwPDB X-ray Structure Validation Report ⓘ

Jun 23, 2018 – 07:55 PM EDT

PDB ID : 6FE2  
Title : Three dimensional structure of human carbonic anhydrase IX  
Authors : Leitans, J.; Tars, K.  
Deposited on : 2017-12-28  
Resolution : 1.87 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

<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  
Xtriage (Phenix) : 1.13  
EDS : rb-20031172  
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) : rb-20031172

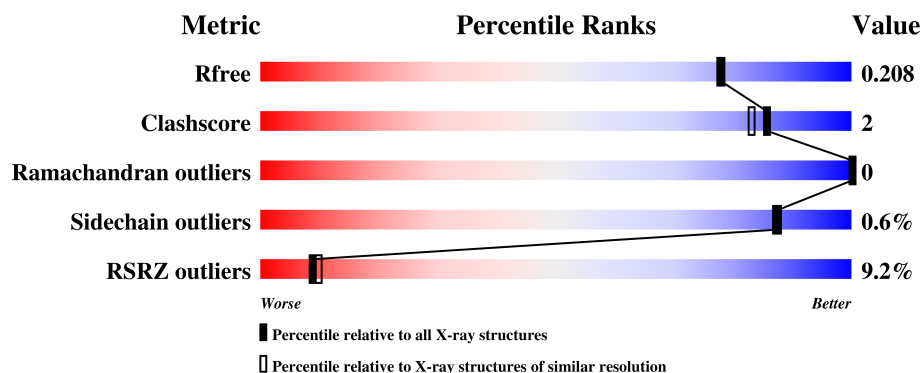
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.87 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	8255 (1.90-1.86)
Clashscore	122126	9028 (1.90-1.86)
Ramachandran outliers	120053	8930 (1.90-1.86)
Sidechain outliers	120020	8930 (1.90-1.86)
RSRZ outliers	108989	8087 (1.90-1.86)

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	257	<div> <div>7%</div> <div>91%</div> <div>5% ..</div> </div>
1	B	257	<div> <div>6%</div> <div>88%</div> <div>6% • 5%</div> </div>
1	C	257	<div> <div>13%</div> <div>92%</div> <div>5% ..</div> </div>
1	D	257	<div> <div>11%</div> <div>94%</div> <div>..</div> </div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 8481 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 9.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	251	Total	C	N	O	S	7	3	0
			1948	1243	342	359	4			
1	B	244	Total	C	N	O	S	0	3	0
			1890	1205	336	345	4			
1	C	251	Total	C	N	O	S	0	3	0
			1937	1240	339	354	4			
1	D	251	Total	C	N	O	S	0	1	0
			1939	1237	340	358	4			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	expression tag	UNP Q16790
A	0	PRO	-	expression tag	UNP Q16790
A	41	SER	CYS	engineered mutation	UNP Q16790
B	2	GLY	-	expression tag	UNP Q16790
B	3	PRO	-	expression tag	UNP Q16790
B	41	SER	CYS	engineered mutation	UNP Q16790
C	-1	GLY	-	expression tag	UNP Q16790
C	0	PRO	-	expression tag	UNP Q16790
C	41	SER	CYS	engineered mutation	UNP Q16790
D	-1	GLY	-	expression tag	UNP Q16790
D	0	PRO	-	expression tag	UNP Q16790
D	41	SER	CYS	engineered mutation	UNP Q16790

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

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

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	D	1	Total 1	Zn 1	0	0
2	C	1	Total 1	Zn 1	0	0

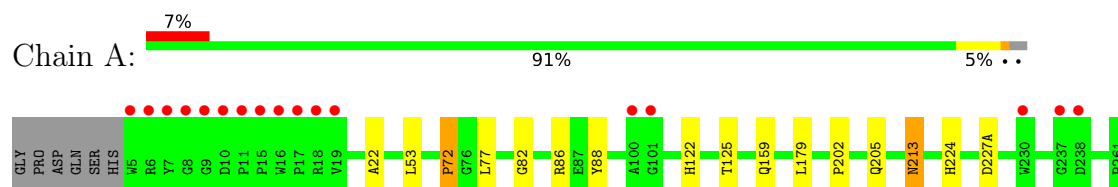
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	250	Total 250	O 250	0	0
3	B	176	Total 176	O 176	0	0
3	C	166	Total 166	O 166	0	0
3	D	171	Total 171	O 171	0	0

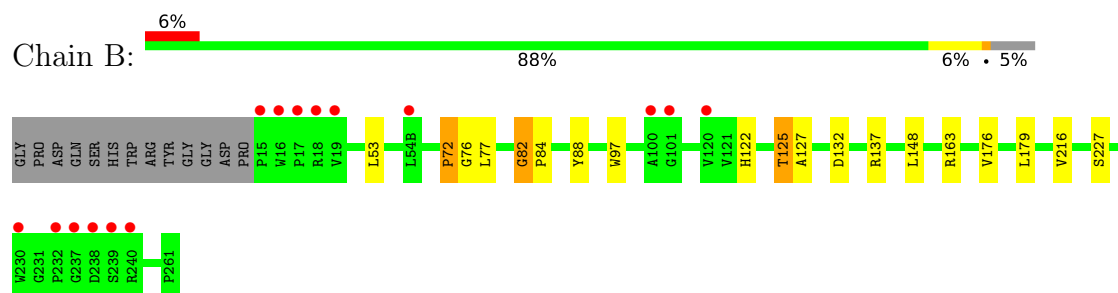
### 3 Residue-property plots [i](#)

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

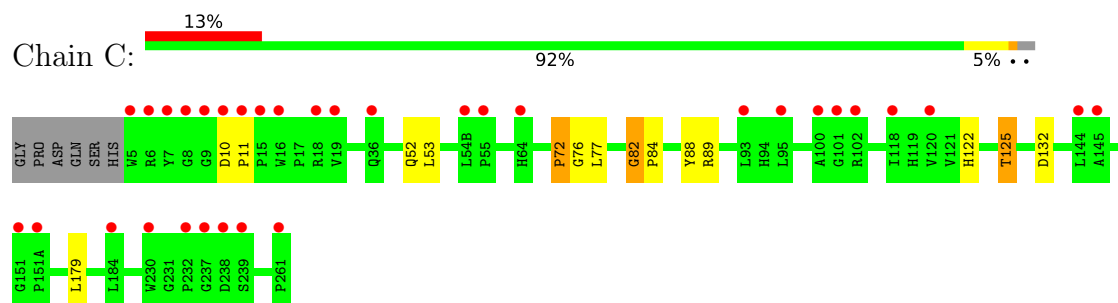
- Molecule 1: Carbonic anhydrase 9



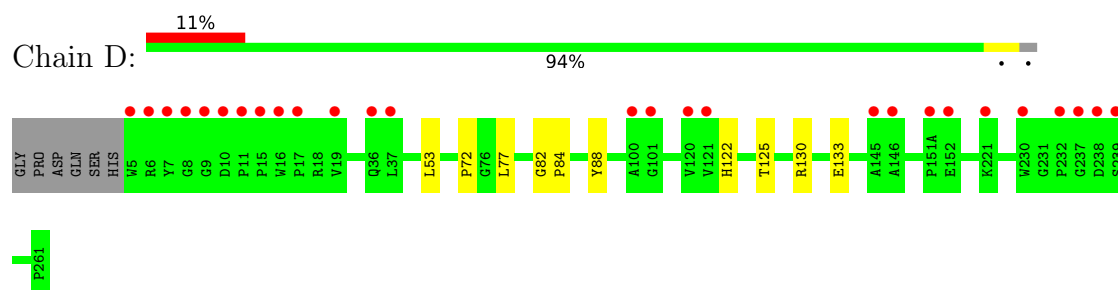
- Molecule 1: Carbonic anhydrase 9



- Molecule 1: Carbonic anhydrase 9



- Molecule 1: Carbonic anhydrase 9



## 4 Data and refinement statistics

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	152.67Å 152.67Å 170.70Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	31.70 – 1.87 31.70 – 1.87	Depositor EDS
% Data completeness (in resolution range)	99.3 (31.70-1.87) 99.3 (31.70-1.87)	Depositor EDS
$R_{merge}$	0.05	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.83 (at 1.87Å)	Xtriage
Refinement program	REFMAC 5.8.0158	Depositor
R, $R_{free}$	0.172 , 0.199 0.181 , 0.208	Depositor DCC
$R_{free}$ test set	6280 reflections (5.16%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	26.8	Xtriage
Anisotropy	0.025	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.39 , 54.9	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.021 for h,-h-k,-l	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	8481	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.48% 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

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: 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	0.78	4/2014 (0.2%)	0.73	1/2748 (0.0%)
1	B	0.90	3/1953 (0.2%)	0.73	1/2666 (0.0%)
1	C	0.97	3/2002 (0.1%)	0.72	3/2734 (0.1%)
1	D	0.70	2/1997 (0.1%)	0.68	0/2722
All	All	0.84	12/7966 (0.2%)	0.72	5/10870 (0.0%)

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	1

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	72	PRO	C-N	28.67	1.84	1.33
1	B	72	PRO	C-N	21.31	1.71	1.33
1	C	82	GLY	C-N	21.28	1.74	1.34
1	B	82	GLY	C-N	18.60	1.69	1.34
1	D	72	PRO	C-N	16.30	1.62	1.33
1	B	125	THR	C-N	15.05	1.68	1.34
1	D	125	THR	C-N	13.52	1.65	1.34
1	A	82	GLY	C-N	12.83	1.58	1.34
1	A	125	THR	C-N	12.81	1.63	1.34
1	A	72	PRO	C-N	12.74	1.55	1.33
1	C	125	THR	C-N	11.74	1.61	1.34
1	A	213	ASN	CG-OD1	5.24	1.35	1.24

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	72	PRO	O-C-N	-10.31	105.68	123.20
1	A	86	ARG	NE-CZ-NH1	6.62	123.61	120.30
1	B	132	ASP	CB-CG-OD1	6.03	123.72	118.30
1	C	132	ASP	CB-CG-OD1	5.48	123.23	118.30
1	C	72	PRO	CA-C-N	-5.39	105.41	116.20

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	72	PRO	Mainchain

## 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	1948	0	1890	6	0
1	B	1890	0	1848	13	0
1	C	1937	0	1882	12	0
1	D	1939	0	1880	6	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
3	A	250	0	0	1	0
3	B	176	0	0	0	0
3	C	166	0	0	2	0
3	D	171	0	0	0	0
All	All	8481	0	7500	36	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 2.

All (36) 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:B:82:GLY:C	1:B:84:PRO:N	1.69	1.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:125:THR:C	1:B:127:ALA:N	1.68	1.44
1:B:72:PRO:C	1:B:76:GLY:N	1.71	1.44
1:C:82:GLY:C	1:C:84:PRO:N	1.74	1.40
1:C:72:PRO:C	1:C:76:GLY:N	1.84	1.30
1:D:82:GLY:C	1:D:84:PRO:N	2.04	1.11
1:A:224[B]:HIS:CE1	1:A:227(A)[B]:ASP:OD2	2.46	0.69
1:B:82:GLY:C	1:B:84:PRO:CA	2.67	0.60
1:C:82:GLY:C	1:C:84:PRO:CA	2.67	0.59
1:B:82:GLY:CA	1:B:84:PRO:N	2.64	0.59
1:C:53:LEU:HD21	1:C:77:LEU:HD11	1.85	0.58
1:A:205:GLN:HE22	1:C:52:GLN:HE22	1.53	0.55
1:B:163:ARG:HD2	1:B:176:VAL:HG12	1.88	0.55
1:D:82:GLY:CA	1:D:84:PRO:N	2.69	0.55
1:D:82:GLY:C	1:D:84:PRO:CA	2.76	0.54
1:B:148:LEU:HD22	1:B:216:VAL:CG2	2.37	0.53
1:B:125:THR:C	1:B:127:ALA:CA	2.71	0.53
1:D:130:ARG:HG3	1:D:133[A]:GLU:HG2	1.89	0.52
1:C:72:PRO:CA	1:C:76:GLY:N	2.69	0.51
1:B:53:LEU:HD21	1:B:77:LEU:HD11	1.93	0.50
1:A:159:GLN:NE2	3:A:402:HOH:O	2.46	0.48
1:B:72:PRO:C	1:B:76:GLY:CA	2.72	0.48
1:A:53:LEU:HD21	1:A:77:LEU:HD11	1.96	0.48
1:C:10:ASP:HB2	1:C:11:PRO:HA	1.97	0.47
1:C:72:PRO:C	1:C:76:GLY:CA	2.80	0.46
1:A:88:TYR:HB3	1:A:122:HIS:HB3	1.96	0.46
1:B:72:PRO:CA	1:B:76:GLY:N	2.71	0.46
1:D:88:TYR:HB3	1:D:122:HIS:HB3	1.98	0.45
1:C:82:GLY:CA	1:C:84:PRO:N	2.71	0.43
1:C:88:TYR:HB3	1:C:122:HIS:HB3	2.00	0.43
1:D:53:LEU:HD21	1:D:77:LEU:HD11	2.00	0.43
1:C:125:THR:HG21	3:C:401:HOH:O	2.18	0.42
1:A:22:ALA:HB3	1:A:202:PRO:O	2.21	0.41
1:B:88:TYR:HB3	1:B:122:HIS:HB3	2.03	0.41
1:B:97:TRP:CD1	1:B:227:SER:HA	2.57	0.40
1:C:89[A]:ARG:NH2	3:C:401:HOH:O	2.54	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	248/257 (96%)	243 (98%)	5 (2%)	0	100	100
1	B	243/257 (95%)	236 (97%)	7 (3%)	0	100	100
1	C	248/257 (96%)	241 (97%)	7 (3%)	0	100	100
1	D	244/257 (95%)	240 (98%)	4 (2%)	0	100	100
All	All	983/1028 (96%)	960 (98%)	23 (2%)	0	100	100

There are no Ramachandran outliers to report.

### 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	205/210 (98%)	203 (99%)	2 (1%)	78	76
1	B	199/210 (95%)	197 (99%)	2 (1%)	78	76
1	C	202/210 (96%)	201 (100%)	1 (0%)	90	89
1	D	203/210 (97%)	203 (100%)	0	100	100
All	All	809/840 (96%)	804 (99%)	5 (1%)	87	87

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	179	LEU
1	A	213	ASN

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Mol	Chain	Res	Type
1	B	137	ARG
1	B	179	LEU
1	C	179	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	36	GLN
1	A	159	GLN
1	A	213	ASN
1	B	36	GLN
1	C	52	GLN
1	C	213	ASN
1	D	213	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 4 ligands modelled in this entry, 4 are monoatomic - leaving 0 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.

## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	D	5
1	C	5
1	B	4
1	A	3

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	C	11:PRO	C	15:PRO	N	3.14
1	B	232:PRO	C	237:GLY	N	3.07
1	C	232:PRO	C	237:GLY	N	2.99
1	D	11:PRO	C	15:PRO	N	2.85
1	A	11:PRO	C	15:PRO	N	2.84
1	D	232:PRO	C	237:GLY	N	2.80
1	A	232:PRO	C	237:GLY	N	2.77
1	D	82:GLY	C	84:PRO	N	2.04
1	C	72:PRO	C	76:GLY	N	1.84
1	C	82:GLY	C	84:PRO	N	1.74
1	B	72:PRO	C	76:GLY	N	1.71
1	B	82:GLY	C	84:PRO	N	1.69
1	B	125:THR	C	127:ALA	N	1.68
1	D	125:THR	C	127:ALA	N	1.65
1	A	125:THR	C	127:ALA	N	1.63
1	D	72:PRO	C	76:GLY	N	1.62
1	C	125:THR	C	127:ALA	N	1.61

## 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	251/257 (97%)	-0.09	17 (6%) 17 19	16, 24, 56, 94	0
1	B	244/257 (94%)	0.09	15 (6%) 21 23	18, 28, 60, 98	0
1	C	251/257 (97%)	0.48	33 (13%) 3 3	22, 37, 69, 95	0
1	D	251/257 (97%)	0.38	27 (10%) 6 6	22, 35, 62, 110	0
All	All	997/1028 (96%)	0.22	92 (9%) 9 10	16, 31, 64, 110	0

All (92) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	15	PRO	8.7
1	D	9	GLY	8.3
1	C	237	GLY	8.2
1	D	11	PRO	8.0
1	D	19	VAL	7.9
1	C	8	GLY	7.8
1	D	8	GLY	7.7
1	C	15	PRO	7.2
1	C	9	GLY	7.0
1	A	8	GLY	6.6
1	B	19	VAL	6.6
1	D	237	GLY	6.5
1	B	237	GLY	6.1
1	D	10	ASP	6.0
1	C	10	ASP	5.7
1	C	11	PRO	5.6
1	A	10	ASP	5.5
1	D	101	GLY	5.5
1	B	16	TRP	5.4
1	A	19	VAL	5.2
1	C	238	ASP	5.2

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Mol	Chain	Res	Type	RSRZ
1	A	9	GLY	5.2
1	C	101	GLY	5.0
1	A	11	PRO	5.0
1	B	100	ALA	4.8
1	B	18	ARG	4.7
1	A	237	GLY	4.7
1	B	101	GLY	4.5
1	A	5	TRP	4.5
1	C	19	VAL	4.4
1	D	238	ASP	4.2
1	C	230	TRP	4.2
1	B	232	PRO	4.1
1	C	118	ILE	4.0
1	D	6	ARG	4.0
1	D	15	PRO	3.9
1	B	17	PRO	3.9
1	D	5	TRP	3.9
1	A	15	PRO	3.9
1	B	238	ASP	3.7
1	D	100	ALA	3.5
1	C	7	TYR	3.4
1	B	230	TRP	3.3
1	C	54(B)	LEU	3.3
1	C	144	LEU	3.2
1	C	5	TRP	3.2
1	A	101	GLY	3.2
1	C	36	GLN	3.0
1	D	7	TYR	3.0
1	C	16	TRP	2.9
1	D	17	PRO	2.9
1	A	238	ASP	2.8
1	C	18	ARG	2.8
1	C	64	HIS	2.8
1	D	36	GLN	2.8
1	A	18	ARG	2.8
1	D	16	TRP	2.7
1	C	232	PRO	2.7
1	C	102	ARG	2.7
1	D	232	PRO	2.7
1	C	120	VAL	2.7
1	D	230	TRP	2.6
1	C	95[A]	LEU	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	239	SER	2.6
1	C	261	PRO	2.5
1	D	151(A)	PRO	2.5
1	D	121	VAL	2.4
1	C	151(A)	PRO	2.4
1	C	151	GLY	2.4
1	C	6	ARG	2.4
1	D	221	LYS	2.4
1	D	239	SER	2.4
1	A	17	PRO	2.3
1	B	54(B)	LEU	2.3
1	C	100	ALA	2.3
1	A	6	ARG	2.3
1	D	120	VAL	2.3
1	C	145	ALA	2.2
1	A	16	TRP	2.2
1	A	7	TYR	2.2
1	C	93	LEU	2.2
1	B	240	ARG	2.2
1	C	55	PRO	2.1
1	C	184	LEU	2.1
1	C	239	SER	2.1
1	A	100	ALA	2.1
1	D	37	LEU	2.1
1	D	146	ALA	2.1
1	D	145	ALA	2.1
1	B	120	VAL	2.0
1	A	230	TRP	2.0
1	D	152	GLU	2.0

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

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 [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(Å <sup>2</sup> )	Q<0.9
2	ZN	C	301	1/1	0.99	0.09	33,33,33,33	0
2	ZN	D	301	1/1	0.99	0.06	27,27,27,27	0
2	ZN	A	301	1/1	1.00	0.07	23,23,23,23	0
2	ZN	B	301	1/1	1.00	0.07	27,27,27,27	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.