



# Full wwPDB X-ray Structure Validation Report ⓘ

May 22, 2020 – 09:23 pm BST

PDB ID : 6DHX  
Title : Structure of TipC2 from Streptococcus intermedius B196  
Authors : Bryant, D.; Klein, T.A.; Whitney, J.C.  
Deposited on : 2018-05-21  
Resolution : 1.75 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  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.11  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.11

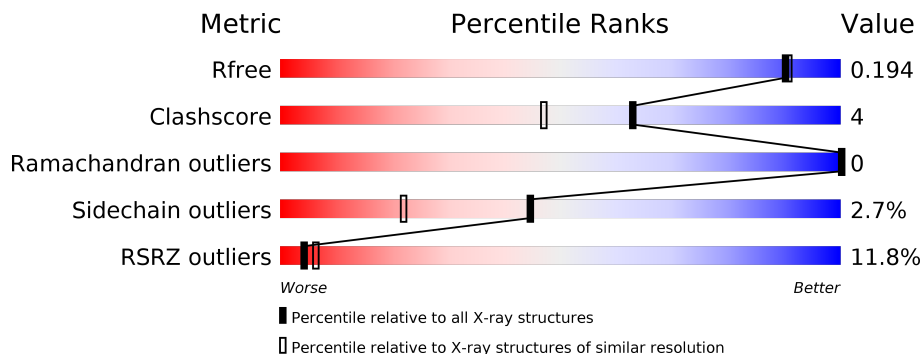
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.75 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))
$R_{free}$	130704	2340 (1.76-1.76)
Clashscore	141614	2466 (1.76-1.76)
Ramachandran outliers	138981	2437 (1.76-1.76)
Sidechain outliers	138945	2437 (1.76-1.76)
RSRZ outliers	127900	2298 (1.76-1.76)

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 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  $\leq 5\%$ . 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	195	 8% 86% 6% • 7%
1	B	195	 12% 78% 12% • 8%
1	C	195	 12% 83% 8% • 9%

## 2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 4978 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 TipC2.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	S	Se			
1	A	182	1535	988	252	292	1	2	0	0	0
1	B	180	1493	961	245	285	1	1	0	0	0
1	C	177	1456	941	238	275	1	1	0	0	0

There are 42 discrepancies between the modelled and reference sequences:

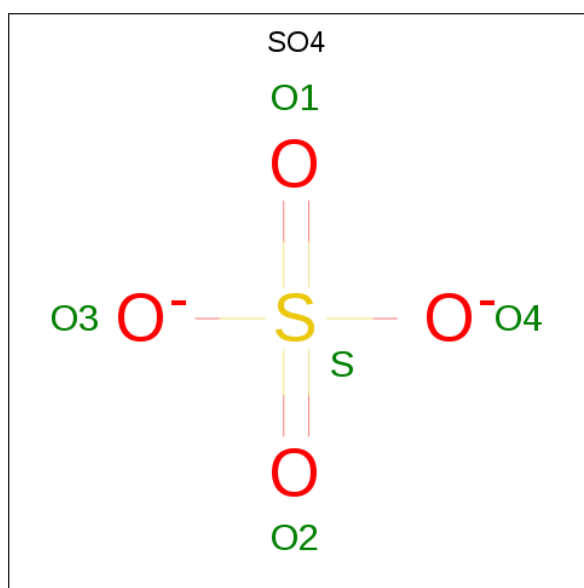
Chain	Residue	Modelled	Actual	Comment	Reference
A	9	MSE	-	initiating methionine	UNP T1ZH71
A	10	GLY	-	expression tag	UNP T1ZH71
A	11	SER	-	expression tag	UNP T1ZH71
A	12	SER	-	expression tag	UNP T1ZH71
A	13	HIS	-	expression tag	UNP T1ZH71
A	14	HIS	-	expression tag	UNP T1ZH71
A	15	HIS	-	expression tag	UNP T1ZH71
A	16	HIS	-	expression tag	UNP T1ZH71
A	17	HIS	-	expression tag	UNP T1ZH71
A	18	HIS	-	expression tag	UNP T1ZH71
A	19	SER	-	expression tag	UNP T1ZH71
A	20	GLN	-	expression tag	UNP T1ZH71
A	21	ASP	-	expression tag	UNP T1ZH71
A	22	PRO	-	expression tag	UNP T1ZH71
B	9	MSE	-	initiating methionine	UNP T1ZH71
B	10	GLY	-	expression tag	UNP T1ZH71
B	11	SER	-	expression tag	UNP T1ZH71
B	12	SER	-	expression tag	UNP T1ZH71
B	13	HIS	-	expression tag	UNP T1ZH71
B	14	HIS	-	expression tag	UNP T1ZH71
B	15	HIS	-	expression tag	UNP T1ZH71
B	16	HIS	-	expression tag	UNP T1ZH71
B	17	HIS	-	expression tag	UNP T1ZH71

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Chain	Residue	Modelled	Actual	Comment	Reference
B	18	HIS	-	expression tag	UNP T1ZH71
B	19	SER	-	expression tag	UNP T1ZH71
B	20	GLN	-	expression tag	UNP T1ZH71
B	21	ASP	-	expression tag	UNP T1ZH71
B	22	PRO	-	expression tag	UNP T1ZH71
C	9	MSE	-	initiating methionine	UNP T1ZH71
C	10	GLY	-	expression tag	UNP T1ZH71
C	11	SER	-	expression tag	UNP T1ZH71
C	12	SER	-	expression tag	UNP T1ZH71
C	13	HIS	-	expression tag	UNP T1ZH71
C	14	HIS	-	expression tag	UNP T1ZH71
C	15	HIS	-	expression tag	UNP T1ZH71
C	16	HIS	-	expression tag	UNP T1ZH71
C	17	HIS	-	expression tag	UNP T1ZH71
C	18	HIS	-	expression tag	UNP T1ZH71
C	19	SER	-	expression tag	UNP T1ZH71
C	20	GLN	-	expression tag	UNP T1ZH71
C	21	ASP	-	expression tag	UNP T1ZH71
C	22	PRO	-	expression tag	UNP T1ZH71

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		


- Molecule 3 is water.

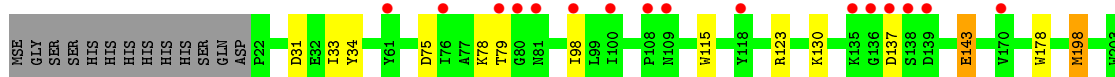
<b>Mol</b>	<b>Chain</b>	<b>Residues</b>	<b>Atoms</b>		<b>ZeroOcc</b>	<b>AltConf</b>
3	A	192	Total 192	O 192	0	0
3	B	167	Total 167	O 167	0	0
3	C	130	Total 130	O 130	0	0

### 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.

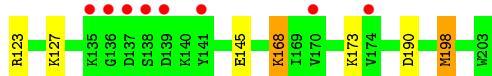
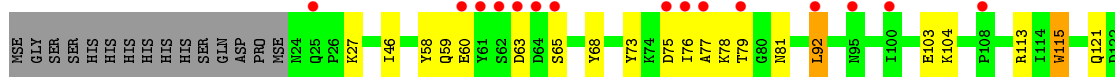
- Molecule 1: TipC2

Chain A: 




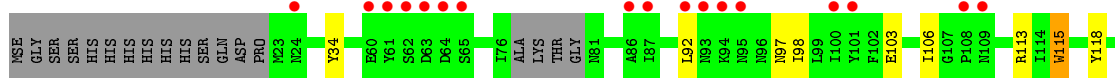
- Molecule 1: TipC2

Chain B: 



- Molecule 1: TipC2

Chain C: 



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	159.74Å 54.52Å 104.43Å 90.00° 108.05° 90.00°	Depositor
Resolution (Å)	33.60 – 1.75 33.60 – 1.75	Depositor EDS
% Data completeness (in resolution range)	96.9 (33.60-1.75) 96.5 (33.60-1.75)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.69 (at 1.75Å)	Xtrriage
Refinement program	PHENIX 1.11.1_2575	Depositor
R, $R_{free}$	0.170 , 0.194 0.170 , 0.194	Depositor DCC
$R_{free}$ test set	2000 reflections (2.33%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	27.7	Xtrriage
Anisotropy	0.208	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.37 , 55.1	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.97	EDS
Total number of atoms	4978	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	38.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.25% 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: 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	0.70	1/1569 (0.1%)	0.74	0/2114
1	B	0.73	1/1526 (0.1%)	0.74	0/2061
1	C	0.58	0/1488	0.67	1/2012 (0.0%)
All	All	0.67	2/4583 (0.0%)	0.72	1/6187 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	115	TRP	CB-CG	-5.63	1.40	1.50
1	B	115	TRP	CB-CG	-5.32	1.40	1.50

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	115	TRP	CA-CB-CG	5.46	124.08	113.70

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	1535	0	1506	6	0
1	B	1493	0	1447	24	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	1456	0	1389	8	0
2	A	5	0	0	0	0
3	A	192	0	0	0	0
3	B	167	0	0	8	0
3	C	130	0	0	1	0
All	All	4978	0	4342	38	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 4.

All (38) 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:34:TYR:HB2	1:A:98:ILE:HD11	1.55	0.87
1:B:190:ASP:OD1	3:B:301:HOH:O	1.96	0.84
1:B:75:ASP:HA	1:B:78:LYS:HE3	1.77	0.67
1:C:34:TYR:HB2	1:C:98:ILE:HD11	1.76	0.66
1:B:60:GLU:OE2	3:B:302:HOH:O	2.13	0.65
1:B:103:GLU:OE2	1:B:113:ARG:NE	2.31	0.64
1:B:145:GLU:OE1	3:B:303:HOH:O	2.15	0.63
1:C:113:ARG:HE	1:C:115:TRP:HE3	1.48	0.62
1:B:113:ARG:HH21	1:B:115:TRP:HZ3	1.47	0.62
1:B:81:ASN:O	3:B:304:HOH:O	2.17	0.58
1:B:190:ASP:OD2	3:B:305:HOH:O	2.19	0.52
1:B:77:ALA:O	1:B:79:THR:HG23	2.08	0.52
1:B:113:ARG:NH2	1:B:115:TRP:HZ3	2.08	0.52
1:A:31:ASP:OD2	1:A:123:ARG:NH2	2.43	0.51
1:B:123:ARG:NH1	3:B:308:HOH:O	2.44	0.49
1:B:81:ASN:HB3	1:B:104:LYS:NZ	2.28	0.49
1:B:113:ARG:HD3	1:B:115:TRP:HE3	1.78	0.48
1:A:198:MSE:HE2	1:A:198:MSE:HB2	1.70	0.48
1:B:121:GLN:HG2	3:B:364:HOH:O	2.13	0.48
1:C:106:ILE:HB	1:C:157:ILE:HD11	1.96	0.47
1:B:168:LYS:NZ	3:B:311:HOH:O	2.47	0.46
1:B:113:ARG:HD3	1:B:115:TRP:CE3	2.51	0.46
1:B:73:TYR:CD2	1:B:173:LYS:HE3	2.52	0.45
1:B:58:TYR:HB2	1:B:68:TYR:CZ	2.52	0.44
1:B:27:LYS:HG3	1:B:123:ARG:NH1	2.32	0.44
1:B:63:ASP:OD2	1:B:65:SER:HB3	2.19	0.42
1:B:198:MSE:HE2	1:B:198:MSE:HB2	1.63	0.42
1:B:127:LYS:HZ3	1:B:198:MSE:HE2	1.85	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:75:ASP:HA	1:A:78:LYS:HE2	2.02	0.41
1:A:33:ILE:HG12	1:A:178:TRP:CD2	2.56	0.41
1:C:103:GLU:OE2	1:C:113:ARG:NH1	2.54	0.41
1:C:167:GLU:HG3	1:C:172:GLN:NE2	2.36	0.41
1:C:97:ASN:HB2	1:C:118:TYR:O	2.21	0.40
1:B:103:GLU:OE2	1:B:113:ARG:NH1	2.54	0.40
1:A:130:LYS:HD3	1:A:143:GLU:OE2	2.22	0.40
1:B:92:LEU:HA	1:B:92:LEU:HD12	1.87	0.40
1:C:153:LYS:HD3	3:C:403:HOH:O	2.21	0.40
1:C:198:MSE:HB2	1:C:198:MSE:HE3	1.73	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	180/195 (92%)	174 (97%)	6 (3%)	0	100	100
1	B	178/195 (91%)	172 (97%)	6 (3%)	0	100	100
1	C	173/195 (89%)	166 (96%)	7 (4%)	0	100	100
All	All	531/585 (91%)	512 (96%)	19 (4%)	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	168/180 (93%)	164 (98%)	4 (2%)	49	26
1	B	160/180 (89%)	154 (96%)	6 (4%)	33	11
1	C	152/180 (84%)	149 (98%)	3 (2%)	55	34
All	All	480/540 (89%)	467 (97%)	13 (3%)	44	22

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	79	THR
1	A	137	ASP
1	A	143	GLU
1	A	198	MSE
1	B	46	ILE
1	B	59	GLN
1	B	76	ILE
1	B	92	LEU
1	B	168	LYS
1	B	198	MSE
1	C	92	LEU
1	C	129	VAL
1	C	187	SER

Some 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 carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

1 ligand is modelled in this entry.

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	$\# Z  > 2$
2	SO4	A	301	-	4,4,4	0.39	0	6,6,6	0.52	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.

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](#)

There are no chain breaks in this entry.

## 6 Fit of model and data

### 6.1 Protein, DNA and RNA chains

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	180/195 (92%)	0.35	16 (8%) <b>9</b> <b>12</b>	19, 32, 59, 109	0
1	B	179/195 (91%)	0.49	23 (12%) <b>3</b> <b>5</b>	18, 29, 89, 107	0
1	C	175/195 (89%)	0.45	24 (13%) <b>3</b> <b>4</b>	23, 37, 81, 104	0
All	All	534/585 (91%)	0.43	63 (11%) <b>4</b> <b>6</b>	18, 34, 82, 109	0

All (63) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	61	TYR	9.9
1	B	76	ILE	9.5
1	A	137	ASP	8.6
1	B	62	SER	8.4
1	B	138	SER	8.0
1	C	137	ASP	7.6
1	A	61	TYR	7.6
1	B	137	ASP	7.5
1	A	136	GLY	6.7
1	B	136	GLY	6.3
1	C	136	GLY	5.9
1	A	76	ILE	5.3
1	B	77	ALA	5.2
1	C	93	ASN	5.2
1	C	138	SER	5.1
1	A	79	THR	5.0
1	C	92	LEU	4.9
1	B	64	ASP	4.9
1	B	63	ASP	4.9
1	A	138	SER	4.3
1	A	135	LYS	4.2
1	C	135	LYS	4.2
1	B	79	THR	4.0

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
1	C	62	SER	3.8
1	C	100	ILE	3.7
1	C	94	LYS	3.6
1	B	135	LYS	3.5
1	C	174	VAL	3.5
1	C	108	PRO	3.2
1	C	87	ILE	3.1
1	B	65	SER	3.1
1	A	100	ILE	3.1
1	A	170	VAL	3.1
1	C	24	ASN	3.0
1	C	64	ASP	3.0
1	B	92	LEU	3.0
1	A	108	PRO	3.0
1	A	80	GLY	3.0
1	A	98	ILE	2.9
1	A	139	ASP	2.9
1	B	141	TYR	2.8
1	C	65	SER	2.8
1	B	100	ILE	2.8
1	C	175	LEU	2.7
1	A	109	ASN	2.7
1	B	174	VAL	2.7
1	C	95	ASN	2.7
1	A	118	TYR	2.6
1	C	170	VAL	2.6
1	A	81	ASN	2.6
1	B	139	ASP	2.4
1	B	60	GLU	2.4
1	B	75	ASP	2.4
1	C	61	TYR	2.3
1	B	108	PRO	2.3
1	C	60	GLU	2.3
1	C	86	ALA	2.2
1	C	63	ASP	2.2
1	B	25	GLN	2.1
1	C	101	TYR	2.1
1	B	95	ASN	2.0
1	B	170	VAL	2.0
1	C	109	ASN	2.0

## 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
2	SO4	A	301	5/5	0.99	0.09	29,31,33,35	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.