



# wwPDB X-ray Structure Validation Summary Report ⓘ

Jan 31, 2016 – 07:40 PM GMT

PDB ID : 1GN1  
Title : CRYSTAL STRUCTURE OF THE MOUSE CCT GAMMA APICAL DOMAIN (MONOCLINIC)  
Authors : Pappenberger, G.; Wilsher, J.A.; Roe, S.M.; Willison, K.R.; Pearl, L.H.  
Deposited on : 2001-10-01  
Resolution : 2.80 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

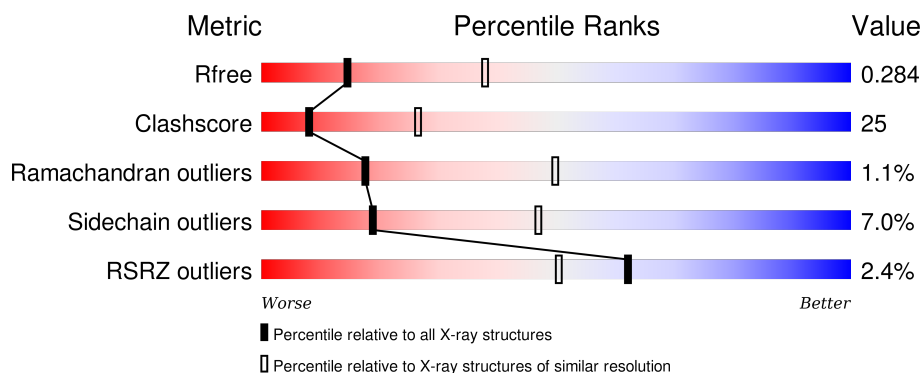
# 1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.80 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	2393 (2.80-2.80)
Clashscore	102246	2827 (2.80-2.80)
Ramachandran outliers	100387	2782 (2.80-2.80)
Sidechain outliers	100360	2784 (2.80-2.80)
RSRZ outliers	91569	2404 (2.80-2.80)

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	178	<div> <div>56%</div> <div>28%</div> <div>•</div> <div>15%</div> </div>
1	B	178	<div> <div>45%</div> <div>35%</div> <div>•</div> <div>18%</div> </div>
1	C	178	<div> <div>44%</div> <div>32%</div> <div>•</div> <div>21%</div> </div>
1	D	178	<div> <div>52%</div> <div>28%</div> <div>•</div> <div>18%</div> </div>
1	E	178	<div> <div>45%</div> <div>28%</div> <div>•</div> <div>24%</div> </div>

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Mol	Chain	Length	Quality of chain
1	F	178	
1	G	178	
1	H	178	

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
2	CA	F	1378	-	-	-	X

## 2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 9018 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 CCT-GAMMA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	151	Total	C	N	O	S	0	0	0
			1175	734	213	221	7			
1	B	146	Total	C	N	O	S	0	0	0
			1140	711	208	213	8			
1	C	140	Total	C	N	O	S	0	0	0
			1069	666	191	204	8			
1	D	146	Total	C	N	O	S	0	0	0
			1103	691	196	209	7			
1	E	135	Total	C	N	O	S	0	0	0
			997	621	173	196	7			
1	F	152	Total	C	N	O	S	0	0	0
			1191	748	215	220	8			
1	G	153	Total	C	N	O	S	0	0	0
			1166	731	206	220	9			
1	H	145	Total	C	N	O	S	0	0	0
			1130	708	202	213	7			

- Molecule 2 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	H	1	Total	Ca	0	0
			1	1		
2	F	1	Total	Ca	0	0
			1	1		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	7	Total	O	0	0
			7	7		
3	B	6	Total	O	0	0
			6	6		

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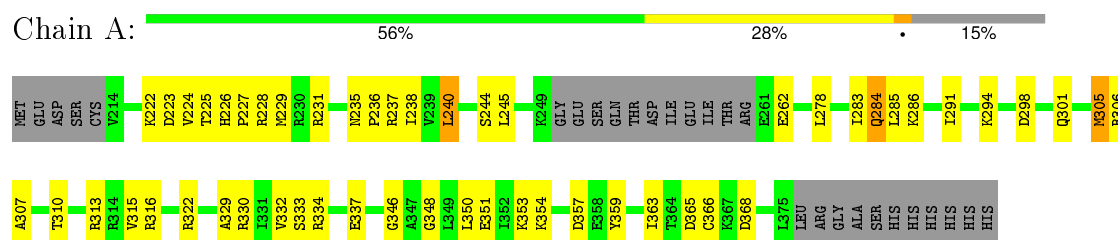
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	C	2	Total 2	O 2	0	0
3	D	7	Total 7	O 7	0	0
3	E	2	Total 2	O 2	0	0
3	F	11	Total 11	O 11	0	0
3	G	3	Total 3	O 3	0	0
3	H	7	Total 7	O 7	0	0

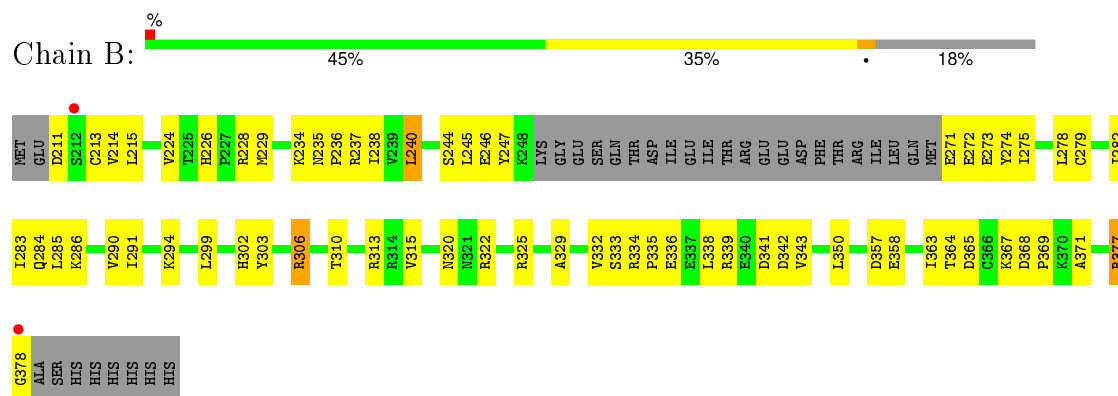
### 3 Residue-property plots

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.

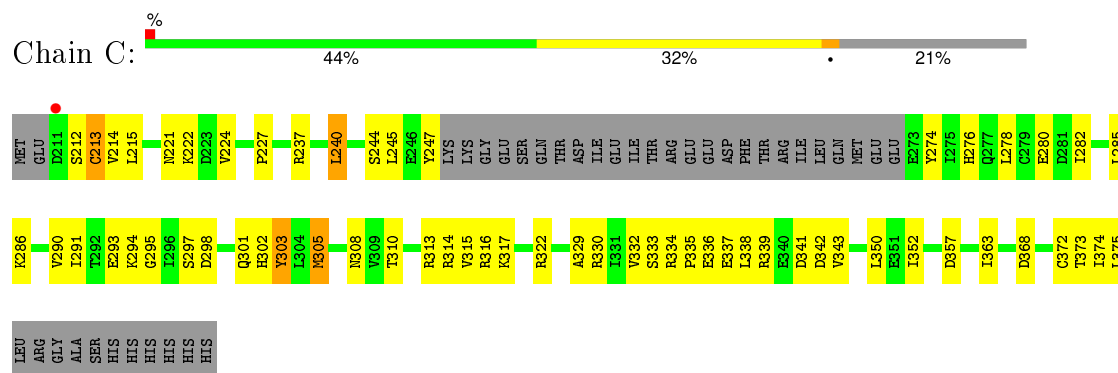
#### • Molecule 1: CCT-GAMMA



#### • Molecule 1: CCT-GAMMA



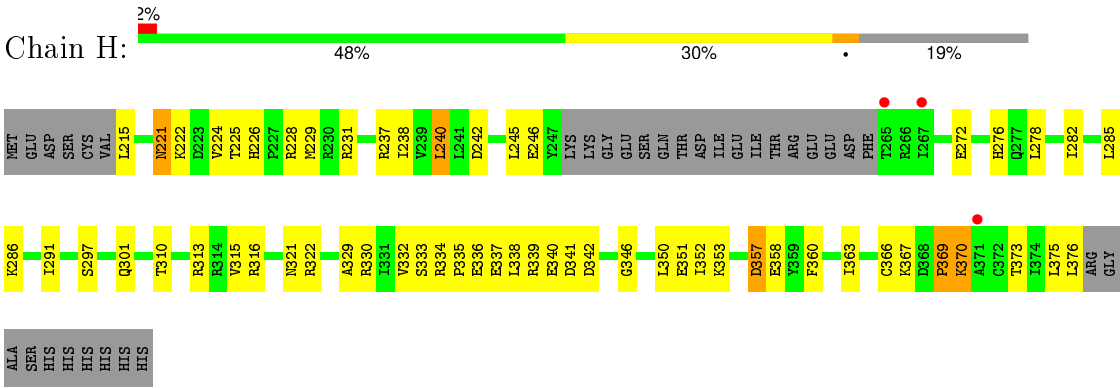
#### • Molecule 1: CCT-GAMMA



#### • Molecule 1: CCT-GAMMA









## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	60.24Å 234.23Å 62.70Å 90.00° 114.70° 90.00°	Depositor
Resolution (Å)	29.11 – 2.80 29.11 – 2.80	Depositor EDS
% Data completeness (in resolution range)	84.8 (29.11-2.80) 84.8 (29.11-2.80)	Depositor EDS
$R_{merge}$	0.05	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.89 (at 2.80Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, $R_{free}$	0.238 , 0.285 0.233 , 0.284	Depositor DCC
$R_{free}$ test set	1632 reflections (4.97%)	DCC
Wilson B-factor (Å <sup>2</sup> )	78.1	Xtriage
Anisotropy	0.467	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.30 , 68.9	EDS
Estimated twinning fraction	0.033 for l,-k,h	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 32815 reflections	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	9018	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	83.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.75% 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, 0.375 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: CA

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	0.49	0/1189	0.72	1/1607 (0.1%)
1	B	0.46	0/1153	0.69	0/1556
1	C	0.47	0/1081	0.74	1/1464 (0.1%)
1	D	0.44	0/1115	0.65	0/1511
1	E	0.40	0/1009	0.61	0/1372
1	F	0.51	0/1204	0.73	1/1624 (0.1%)
1	G	0.48	0/1180	0.69	0/1599
1	H	0.49	0/1143	0.69	0/1545
All	All	0.47	0/9074	0.69	3/12278 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	314	ARG	NE-CZ-NH2	-6.73	116.94	120.30
1	A	305	MET	CA-CB-CG	5.58	122.78	113.30
1	C	372	CYS	CA-CB-SG	-5.43	104.22	114.00

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	1175	0	1154	50	0
1	B	1140	0	1145	57	0
1	C	1069	0	1050	71	0
1	D	1103	0	1074	45	0
1	E	997	0	934	37	0
1	F	1191	0	1206	67	0
1	G	1166	0	1130	65	0
1	H	1130	0	1129	82	0
2	F	1	0	0	0	0
2	H	1	0	0	0	0
3	A	7	0	0	4	0
3	B	6	0	0	0	0
3	C	2	0	0	0	0
3	D	7	0	0	0	0
3	E	2	0	0	0	0
3	F	11	0	0	0	0
3	G	3	0	0	0	0
3	H	7	0	0	3	0
All	All	9018	0	8822	437	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 25.

The worst 5 of 437 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:F:374:ILE:H	1:F:374:ILE:HD12	1.12	1.14
1:A:225:THR:HB	1:A:301:GLN:HE21	1.08	1.14
1:F:246:GLU:HG2	1:F:297:SER:HB3	1.35	1.07
1:F:370:LYS:HE3	1:F:370:LYS:HA	1.39	1.04
1:A:225:THR:CB	1:A:301:GLN:HE21	1.72	1.00

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	147/178 (83%)	135 (92%)	12 (8%)	0	100	100
1	B	142/178 (80%)	127 (89%)	14 (10%)	1 (1%)	26	62
1	C	136/178 (76%)	125 (92%)	9 (7%)	2 (2%)	13	40
1	D	142/178 (80%)	127 (89%)	13 (9%)	2 (1%)	14	42
1	E	131/178 (74%)	120 (92%)	9 (7%)	2 (2%)	13	40
1	F	148/178 (83%)	130 (88%)	16 (11%)	2 (1%)	14	42
1	G	149/178 (84%)	137 (92%)	11 (7%)	1 (1%)	26	62
1	H	141/178 (79%)	129 (92%)	9 (6%)	3 (2%)	9	29
All	All	1136/1424 (80%)	1030 (91%)	93 (8%)	13 (1%)	17	50

5 of 13 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	H	336	GLU
1	H	370	LYS
1	B	272	GLU
1	D	229	MET
1	C	213	CYS

### 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	123/160 (77%)	118 (96%)	5 (4%)	37	72
1	B	123/160 (77%)	116 (94%)	7 (6%)	25	58
1	C	114/160 (71%)	103 (90%)	11 (10%)	10	29
1	D	113/160 (71%)	107 (95%)	6 (5%)	28	61
1	E	101/160 (63%)	90 (89%)	11 (11%)	8	23
1	F	128/160 (80%)	119 (93%)	9 (7%)	19	47

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	G	121/160 (76%)	109 (90%)	12 (10%)	10	28
1	H	121/160 (76%)	116 (96%)	5 (4%)	37	72
All	All	944/1280 (74%)	878 (93%)	66 (7%)	19	47

5 of 66 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	E	234	LYS
1	E	305	MET
1	G	377	ARG
1	E	240	LEU
1	E	278	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 20 such sidechains are listed below:

Mol	Chain	Res	Type
1	C	301	GLN
1	C	308	ASN
1	F	284	GLN
1	C	235	ASN
1	C	277	GLN

### 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 2 ligands modelled in this entry, 2 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](#)

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	151/178 (84%)	-0.35	0 100 100	56, 72, 97, 124	0
1	B	146/178 (82%)	-0.37	2 (1%) 78 69	54, 75, 101, 122	0
1	C	140/178 (78%)	-0.27	1 (0%) 89 84	52, 72, 107, 135	0
1	D	146/178 (82%)	-0.10	3 (2%) 67 56	67, 90, 128, 141	0
1	E	135/178 (75%)	0.13	6 (4%) 38 26	75, 103, 140, 152	0
1	F	152/178 (85%)	-0.00	8 (5%) 30 20	52, 76, 114, 146	0
1	G	153/178 (85%)	-0.23	5 (3%) 50 38	57, 79, 112, 127	0
1	H	145/178 (81%)	-0.03	3 (2%) 67 56	52, 75, 116, 134	0
All	All	1168/1424 (82%)	-0.16	28 (2%) 62 50	52, 80, 121, 152	0

The worst 5 of 28 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	264	PHE	4.6
1	C	211	ASP	4.3
1	F	263	ASP	4.3
1	G	213	CYS	3.6
1	E	247	TYR	3.6

### 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. 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
2	CA	F	1378	1/1	0.97	0.29	4.98	73,73,73,73	0
2	CA	H	1377	1/1	0.81	0.26	-	89,89,89,89	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.