



# Full wwPDB X-ray Structure Validation Report ⓘ

Oct 10, 2019 – 09:54 AM EDT

PDB ID : 1V1F  
Title : Structure of the Arabidopsis thaliana SOS3 complexed with Calcium(II) and Manganese(II) ions  
Authors : Sanchez-Barrena, M.J.; Martinez-Ripoll, M.; Zhu, J.K.; Albert, A.  
Deposited on : 2004-04-15  
Resolution : 3.00 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 : 1.8.0 (224370), CSD as540be (2019)  
Xtriage (Phenix) : 1.13  
EDS : 2.4  
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) : 2.4

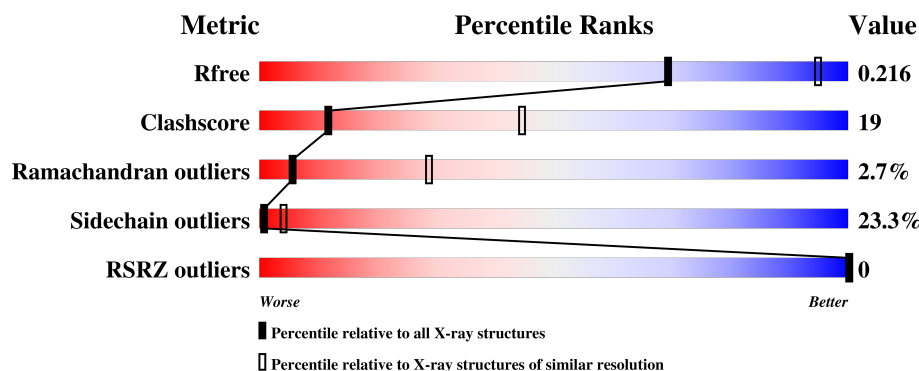
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 3.00 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	1851 (3.00-3.00)
Clashscore	122126	2167 (3.00-3.00)
Ramachandran outliers	120053	2101 (3.00-3.00)
Sidechain outliers	120020	2104 (3.00-3.00)
RSRZ outliers	108989	1751 (3.00-3.00)

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	222	

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	IOD	A	1206	-	-	X	-

## 2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 1614 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 CALCINEURIN B-LIKE PROTEIN 4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	188	Total	C	N	O	S	0	0	0
			1547	997	258	288	4			

- Molecule 2 is IODIDE ION (three-letter code: IOD) (formula: I).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	7	Total	I	0	0
			7	7		

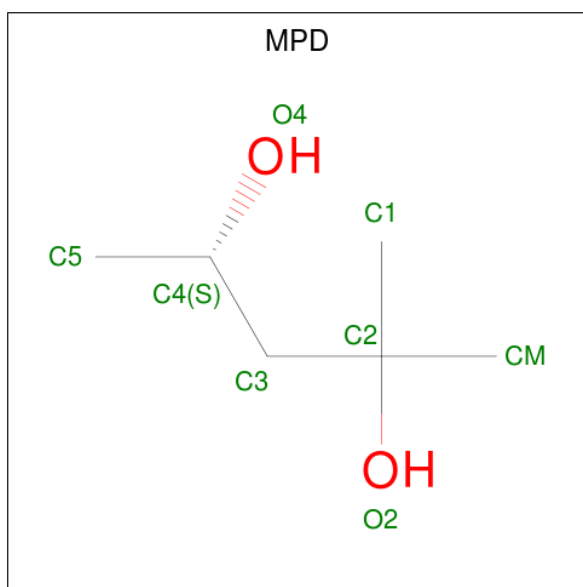
- Molecule 3 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	3	Total	Ca	0	0
			3	3		

- Molecule 4 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Mn	0	0
			1	1		

- Molecule 5 is (4S)-2-METHYL-2,4-PENTANEDIOL (three-letter code: MPD) (formula: C<sub>6</sub>H<sub>14</sub>O<sub>2</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			8	6	2		
5	A	1	Total	C	O	0	0
			8	6	2		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	40	Total	O	0	0
			40	40		



## 4 Data and refinement statistics

Property	Value	Source
Space group	I 41	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	91.90Å 91.90Å 85.60Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 3.00 19.40 – 3.00	Depositor EDS
% Data completeness (in resolution range)	97.2 (20.00-3.00) 97.2 (19.40-3.00)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.21 (at 2.98Å)	Xtriage
Refinement program	REFMAC 5.1.24	Depositor
R, $R_{free}$	0.237 , 0.262 0.233 , 0.216	Depositor DCC
$R_{free}$ test set	326 reflections (4.67%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	79.6	Xtriage
Anisotropy	0.056	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 36.4	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.42$ , $\langle L^2 \rangle = 0.25$	Xtriage
Estimated twinning fraction	0.179 for -k,-h,-l	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	1614	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	61.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.24% 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: CA, IOD, MPD, MN

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	0.79	0/1579	1.05	9/2129 (0.4%)

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

There are no bond length outliers.

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	84	ASP	CB-CG-OD2	7.74	125.27	118.30
1	A	158	ASP	CB-CG-OD2	6.03	123.73	118.30
1	A	77	ASP	CB-CG-OD2	5.87	123.58	118.30
1	A	169	ASP	CB-CG-OD2	5.76	123.48	118.30
1	A	52	ASP	CB-CG-OD2	5.71	123.44	118.30
1	A	198	ASP	CB-CG-OD2	5.60	123.34	118.30
1	A	21	ASP	CB-CG-OD1	5.44	123.19	118.30
1	A	183	LEU	CA-CB-CG	5.32	127.54	115.30
1	A	121	ASP	CB-CG-OD1	5.21	122.99	118.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	15	ARG	Peptide

## 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	1547	0	1546	59	0
2	A	7	0	0	4	0
3	A	3	0	0	0	0
4	A	1	0	0	0	0
5	A	16	0	28	0	0
6	A	40	0	0	4	0
All	All	1614	0	1574	59	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 19.

All (59) 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:166:ARG:HG3	1:A:166:ARG:HH11	1.13	1.12
1:A:87:ARG:HG2	1:A:87:ARG:HH11	0.94	1.06
1:A:167:LYS:HE3	1:A:173:ASP:OD2	1.66	0.95
1:A:104:HIS:CD2	1:A:106:SER:H	1.86	0.94
1:A:87:ARG:HG2	1:A:87:ARG:NH1	1.73	0.88
1:A:104:HIS:HD2	1:A:106:SER:H	1.19	0.86
1:A:194:PRO:HB3	2:A:1205:IOD:I	2.55	0.75
1:A:42:LEU:HD12	1:A:46:LEU:HD21	1.68	0.75
1:A:166:ARG:HG3	1:A:166:ARG:NH1	1.92	0.74
1:A:166:ARG:HH11	1:A:166:ARG:CG	1.98	0.74
1:A:112:LYS:HG2	1:A:191:MET:HB3	1.72	0.72
1:A:195:TYR:HD2	6:A:2037:HOH:O	1.75	0.69
1:A:178:LYS:HD3	2:A:1214:IOD:I	2.63	0.69
1:A:104:HIS:HD2	1:A:106:SER:N	1.91	0.68
1:A:49:SER:OG	1:A:50:ILE:HG13	1.96	0.65
1:A:112:LYS:NZ	1:A:192:THR:HB	2.11	0.65
1:A:55:LEU:HD13	1:A:90:VAL:HG21	1.79	0.65
1:A:133:LEU:HD12	1:A:136:MET:HE3	1.82	0.62
1:A:133:LEU:HD12	1:A:136:MET:CE	2.31	0.60
1:A:112:LYS:HZ1	1:A:192:THR:HB	1.67	0.59
1:A:140:LEU:O	1:A:143:GLU:HB3	2.03	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:39:LEU:HD23	1:A:42:LEU:HD23	1.87	0.57
1:A:114:LYS:HG2	2:A:1206:IOD:I	2.74	0.57
1:A:71:ARG:HE	1:A:79:ILE:HD11	1.70	0.56
1:A:114:LYS:HD3	6:A:2021:HOH:O	2.04	0.56
1:A:80:PHE:CE1	1:A:91:ILE:HD12	2.43	0.54
1:A:87:ARG:HH11	1:A:87:ARG:CG	1.87	0.54
1:A:130:ARG:NH2	1:A:158:ASP:OD1	2.39	0.53
1:A:100:LEU:HD22	1:A:196:LEU:HD11	1.90	0.53
1:A:123:ARG:O	1:A:125:THR:HG23	2.10	0.51
1:A:42:LEU:CD1	1:A:46:LEU:HD21	2.41	0.49
1:A:133:LEU:O	1:A:137:VAL:HG23	2.12	0.49
1:A:120:TYR:CE1	1:A:136:MET:SD	3.06	0.48
1:A:82:VAL:CG2	1:A:115:PHE:CZ	2.97	0.48
1:A:127:PHE:CD2	1:A:171:LYS:HD3	2.49	0.47
1:A:75:PHE:O	1:A:79:ILE:HG12	2.14	0.47
1:A:66:PHE:CD2	1:A:66:PHE:C	2.89	0.46
1:A:78:ARG:HA	1:A:78:ARG:HE	1.80	0.46
1:A:77:ASP:O	1:A:81:ASP:HB2	2.14	0.46
1:A:15:ARG:N	1:A:15:ARG:HD2	2.32	0.45
1:A:87:ARG:NH1	1:A:87:ARG:CG	2.57	0.45
1:A:167:LYS:HG2	1:A:176:GLU:OE2	2.17	0.44
1:A:114:LYS:CG	2:A:1206:IOD:I	3.37	0.43
1:A:71:ARG:HH12	1:A:190:ASN:HD22	1.67	0.43
1:A:34:GLU:H	1:A:34:GLU:CD	2.21	0.43
1:A:81:ASP:OD1	1:A:87:ARG:NH2	2.51	0.43
1:A:142:HIS:HB2	6:A:2027:HOH:O	2.19	0.42
1:A:82:VAL:HG23	1:A:115:PHE:CZ	2.54	0.42
1:A:197:LYS:HZ3	1:A:197:LYS:HG3	1.73	0.42
1:A:78:ARG:NH1	1:A:121:ASP:O	2.53	0.42
1:A:142:HIS:CE1	6:A:2029:HOH:O	2.72	0.42
1:A:69:ARG:HG3	1:A:69:ARG:O	2.19	0.41
1:A:43:PHE:CE1	1:A:56:ILE:HG13	2.55	0.41
1:A:58:LYS:NZ	1:A:77:ASP:OD1	2.51	0.41
1:A:92:GLU:HG3	1:A:95:GLU:OE2	2.21	0.41
1:A:86:LYS:HG2	1:A:95:GLU:OE1	2.19	0.41
1:A:83:PHE:HB3	1:A:91:ILE:HG12	2.02	0.41
1:A:129:GLU:N	1:A:132:GLU:OE2	2.43	0.40
1:A:174:ILE:HG12	1:A:174:ILE:O	2.20	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	186/222 (84%)	157 (84%)	24 (13%)	5 (3%)	5	29

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	67	ARG
1	A	121	ASP
1	A	63	LEU
1	A	71	ARG
1	A	49	SER

### 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	172/204 (84%)	132 (77%)	40 (23%)	1	4

All (40) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	15	ARG
1	A	27	SER
1	A	42	LEU
1	A	45	LYS
1	A	46	LEU
1	A	48	SER

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Mol	Chain	Res	Type
1	A	49	SER
1	A	51	ILE
1	A	58	LYS
1	A	59	GLU
1	A	61	PHE
1	A	66	PHE
1	A	69	ARG
1	A	73	ASN
1	A	78	ARG
1	A	81	ASP
1	A	86	LYS
1	A	87	ARG
1	A	92	GLU
1	A	100	LEU
1	A	106	SER
1	A	108	PRO
1	A	111	GLU
1	A	114	LYS
1	A	123	ARG
1	A	124	GLN
1	A	143	GLU
1	A	144	SER
1	A	147	VAL
1	A	151	ASP
1	A	152	MET
1	A	163	GLN
1	A	166	ARG
1	A	171	LYS
1	A	174	ILE
1	A	178	LYS
1	A	183	LEU
1	A	192	THR
1	A	198	ASP
1	A	202	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	62	GLN
1	A	104	HIS
1	A	163	GLN
1	A	190	ASN

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Mol	Chain	Res	Type
1	A	200	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 13 ligands modelled in this entry, 11 are monoatomic - leaving 2 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	$\# Z  > 2$
5	MPD	A	1211	-	7,7,7	0.29	0	9,10,10	0.62	0
5	MPD	A	1212	-	7,7,7	0.51	0	9,10,10	0.44	0

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
5	MPD	A	1211	-	-	2/5/5/5	-
5	MPD	A	1212	-	-	3/5/5/5	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	1212	MPD	O2-C2-C3-C4
5	A	1212	MPD	C2-C3-C4-O4
5	A	1212	MPD	C2-C3-C4-C5
5	A	1211	MPD	C2-C3-C4-C5
5	A	1211	MPD	C2-C3-C4-O4

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 [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	188/222 (84%)	-0.35	0 100 100	40, 56, 95, 104	0

There are no RSRZ outliers to report.

### 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(Å <sup>2</sup> )	Q<0.9
5	MPD	A	1212	8/8	0.84	0.25	77,79,81,81	0
5	MPD	A	1211	8/8	0.90	0.18	72,74,75,75	0
3	CA	A	1209	1/1	0.91	0.11	74,74,74,74	0
2	IOD	A	1205	1/1	0.96	0.05	110,110,110,110	0
2	IOD	A	1204	1/1	0.97	0.10	84,84,84,84	0
2	IOD	A	1203	1/1	0.98	0.18	84,84,84,84	0
4	MN	A	1210	1/1	0.98	0.19	59,59,59,59	0
2	IOD	A	1214	1/1	0.98	0.09	83,83,83,83	0
2	IOD	A	1215	1/1	0.98	0.10	97,97,97,97	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
3	CA	A	1207	1/1	0.98	0.11	79,79,79,79	0
2	IOD	A	1206	1/1	0.99	0.08	115,115,115,115	0
2	IOD	A	1213	1/1	0.99	0.09	86,86,86,86	0
3	CA	A	1208	1/1	0.99	0.10	71,71,71,71	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.