



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 05:37 PM UTC

PDB ID : 5RWY / pdb_00005rwy
Title : INPP5D PanDDA analysis group deposition – Crystal Structure of the phosphatase and C2 domains of SHIP1 in complex with Z1259335913
Authors : Bradshaw, W.J.; Newman, J.A.; von Delft, F.; Arrowsmith, C.H.; Edwards, A.M.; Bountra, C.; Gileadi, O.
Deposited on : 2020-10-30
Resolution : 1.35 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

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

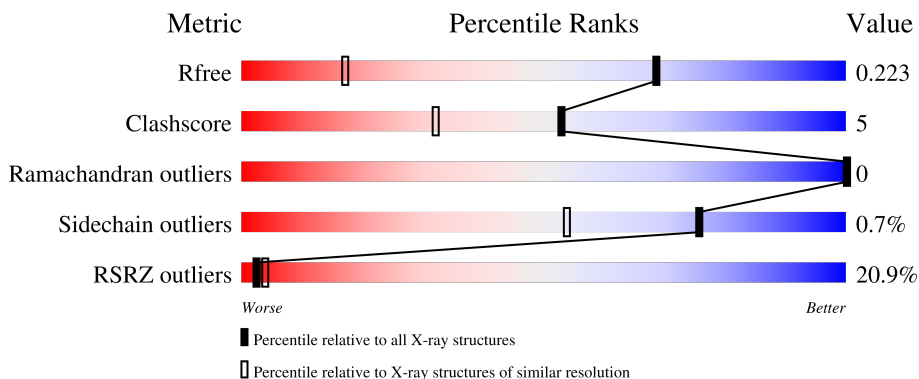
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.35 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}	180053	1216 (1.36-1.36)
Clashscore	190562	1232 (1.36-1.36)
Ramachandran outliers	187476	1220 (1.36-1.36)
Sidechain outliers	187428	1220 (1.36-1.36)
RSRZ outliers	180081	1214 (1.36-1.36)

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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 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	463	21% 90% 8%

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 4730 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.

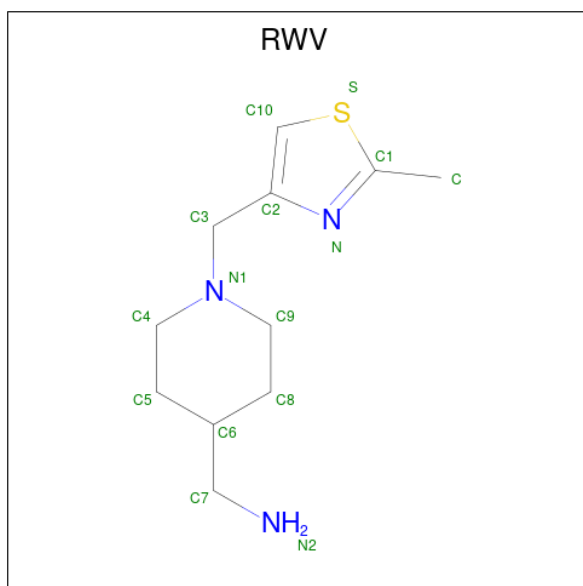
- Molecule 1 is a protein called Phosphatidylinositol 3,4,5-trisphosphate 5-phosphatase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	460	4256	2691	725	820	20	0	33	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	394	SER	-	expression tag	UNP Q92835
A	395	MET	-	expression tag	UNP Q92835

- Molecule 2 is 1-{1-[(2-methyl-1,3-thiazol-4-yl)methyl]piperidin-4-yl}methanamine (CCD ID: RWV) (formula: C₁₁H₁₉N₃S).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	S		
2	A	1	15	11	3	1	0	0

- Molecule 3 is DIMETHYL SULFOXIDE (CCD ID: DMS) (formula: C₂H₆OS).



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

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	444	Total	O	0	7
			447	447		

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	62.58Å 79.42Å 89.32Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	44.70 – 1.35 44.70 – 1.35	Depositor EDS
% Data completeness (in resolution range)	100.0 (44.70-1.35) 100.0 (44.70-1.35)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.40 (at 1.35Å)	Xtriage
Refinement program	REFMAC 5.8.0238	Depositor
R, R_{free}	0.175 , 0.201 0.195 , 0.223	Depositor DCC
R_{free} test set	4893 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	14.5	Xtriage
Anisotropy	0.028	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 31.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	4730	wwPDB-VP
Average B, all atoms (Å ²)	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.39% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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: RWV, DMS

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	1.00	3/4352 (0.1%)	1.08	2/5877 (0.0%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	811	ASP	C-N	-5.78	1.25	1.33
1	A	851	LYS	C-N	-5.19	1.26	1.33
1	A	812	GLU	C-N	5.01	1.40	1.33

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	416	PRO	N-CA-CB	5.68	106.37	103.19
1	A	732	PHE	CA-C-O	-5.06	115.28	120.54

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	4256	0	4154	43	0
2	A	15	0	0	0	0
3	A	12	0	18	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	447	0	0	13	0
All	All	4730	0	4172	43	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.

All (43) 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:567:LYS:HD2	4:A:1203:HOH:O	1.53	1.09
1:A:810:SER:OG	1:A:812:GLU:HG2	1.65	0.96
1:A:781:GLU:OE1	4:A:1001:HOH:O	1.88	0.89
1:A:499[B]:ASN:ND2	4:A:1002:HOH:O	2.16	0.79
1:A:499[C]:ASN:ND2	4:A:1002:HOH:O	2.16	0.79
1:A:432[B]:LYS:HG2	4:A:1104:HOH:O	1.91	0.70
1:A:546[A]:GLU:H	1:A:546[A]:GLU:CD	2.00	0.69
1:A:546[C]:GLU:H	1:A:546[C]:GLU:CD	2.00	0.69
1:A:567:LYS:CD	4:A:1203:HOH:O	2.26	0.66
1:A:556[C]:MET:SD	1:A:839:HIS:HB3	2.40	0.61
1:A:519:ASN:HD21	1:A:520[C]:LYS:NZ	2.01	0.58
1:A:402[A]:MET:HE3	1:A:708[A]:GLU:CG	2.34	0.58
1:A:402[C]:MET:HE3	1:A:708[C]:GLU:CG	2.34	0.58
1:A:519:ASN:HD21	1:A:520[A]:LYS:NZ	2.01	0.58
1:A:402[A]:MET:CE	1:A:708[A]:GLU:CD	2.78	0.56
1:A:402[C]:MET:CE	1:A:708[C]:GLU:CD	2.78	0.56
1:A:810:SER:OG	1:A:812:GLU:CG	2.48	0.55
1:A:402[C]:MET:HE3	1:A:708[C]:GLU:HG2	1.87	0.55
1:A:402[A]:MET:HE3	1:A:708[A]:GLU:HG2	1.87	0.55
1:A:646:GLU:OE2	1:A:652:LYS:HD2	2.09	0.53
1:A:559[B]:LEU:C	1:A:559[B]:LEU:HD12	2.38	0.49
1:A:603[B]:GLN:NE2	4:A:1018:HOH:O	2.44	0.48
1:A:750:GLU:OE1	1:A:752:HIS:HE1	1.97	0.48
1:A:681:LEU:O	1:A:822[B]:ARG:NH2	2.47	0.48
1:A:512:GLY:O	4:A:1005:HOH:O	2.20	0.47
1:A:402[A]:MET:HE1	1:A:708[A]:GLU:CD	2.39	0.47
1:A:402[C]:MET:HE1	1:A:708[C]:GLU:CD	2.40	0.47
1:A:595[B]:THR:HG21	1:A:665:TYR:CE2	2.50	0.47
1:A:546[A]:GLU:OE2	4:A:1004:HOH:O	2.20	0.47
1:A:546[C]:GLU:OE2	4:A:1004:HOH:O	2.20	0.47
1:A:519:ASN:HD21	1:A:520[C]:LYS:HZ3	1.62	0.46
1:A:774:GLU:HB3	4:A:1358:HOH:O	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:432[B]:LYS:HG3	4:A:1045:HOH:O	2.17	0.44
1:A:442:PRO:HG2	1:A:494[C]:LYS:HE2	2.00	0.44
1:A:399:GLU:HG3	4:A:1341:HOH:O	2.17	0.43
1:A:779:PHE:O	1:A:782:THR:OG1	2.28	0.43
1:A:442:PRO:HG2	1:A:494[A]:LYS:HE2	2.00	0.43
1:A:596[B]:TRP:CZ3	1:A:597:GLU:CD	2.97	0.42
1:A:402[C]:MET:HE3	1:A:708[C]:GLU:CD	2.45	0.42
1:A:402[A]:MET:HE3	1:A:708[A]:GLU:CD	2.45	0.42
1:A:552:ASN:O	1:A:556[B]:MET:HG2	2.21	0.41
1:A:519:ASN:HD21	1:A:520[B]:LYS:HE3	1.85	0.41
1:A:725:ASP:OD1	1:A:727:GLN:NE2	2.54	0.41

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	524/463 (113%)	514 (98%)	10 (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	479/417 (115%)	476 (99%)	3 (1%)	78	58

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	743	SER
1	A	769	GLU
1	A	781	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	469	GLN
1	A	486	ASN
1	A	519	ASN
1	A	552	ASN
1	A	744	GLN
1	A	752	HIS
1	A	839	HIS
1	A	856	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

4 ligands are 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
3	DMS	A	904	-	3,3,3	0.36	0	3,3,3	0.29	0
3	DMS	A	903	-	3,3,3	0.38	0	3,3,3	0.38	0
3	DMS	A	902	-	3,3,3	0.28	0	3,3,3	0.41	0
2	RWV	A	901	-	15,16,16	3.63	9 (60%)	18,21,21	1.45	3 (16%)

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
2	RWV	A	901	-	-	0/6/16/16	0/2/2/2

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	901	RWV	C3-N1	-10.90	1.23	1.47
2	A	901	RWV	C2-N	4.21	1.46	1.38
2	A	901	RWV	C10-S	-3.74	1.62	1.71
2	A	901	RWV	C10-C2	3.55	1.42	1.35
2	A	901	RWV	C8-C6	-2.79	1.44	1.52
2	A	901	RWV	C3-C2	2.69	1.53	1.50
2	A	901	RWV	C5-C6	-2.55	1.45	1.52
2	A	901	RWV	C1-S	-2.40	1.64	1.72
2	A	901	RWV	C1-N	2.17	1.38	1.31

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	901	RWV	C10-C2-N	-3.32	109.96	115.12
2	A	901	RWV	C10-S-C1	3.13	95.56	89.73
2	A	901	RWV	C9-N1-C4	2.21	113.60	108.84

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, 95th 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 ²)	Q<0.9
1	A	460/463 (99%)	1.19	96 (20%) 2 4	2, 14, 50, 84	40 (8%)

All (96) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	596[A]	TRP	14.3
1	A	448[A]	ILE	11.2
1	A	721	PRO	10.8
1	A	724	VAL	10.5
1	A	559[A]	LEU	10.3
1	A	656[A]	THR	10.0
1	A	505[A]	CYS	9.6
1	A	722	GLY	9.5
1	A	513	ILE	9.4
1	A	467[A]	SER	8.6
1	A	461	LEU	8.5
1	A	723	THR	8.0
1	A	556[A]	MET	7.8
1	A	595[A]	THR	7.8
1	A	716	VAL	7.8
1	A	499[A]	ASN	7.6
1	A	761[A]	LYS	7.4
1	A	603[A]	GLN	7.3
1	A	550[A]	ARG	7.2
1	A	754[A]	SER	7.1
1	A	713[A]	SER	7.1
1	A	494[A]	LYS	6.9
1	A	592[A]	ASP	6.9
1	A	546[A]	GLU	6.9
1	A	516	THR	6.7
1	A	664[A]	LYS	6.7
1	A	853[A]	GLN	6.6

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Mol	Chain	Res	Type	RSRZ
1	A	696[A]	MET	6.5
1	A	715	PHE	6.4
1	A	773	GLY	6.3
1	A	469	GLN	6.3
1	A	466	HIS	6.2
1	A	520[A]	LYS	6.2
1	A	414[A]	ASN	6.2
1	A	822[A]	ARG	6.2
1	A	771	SER	5.9
1	A	720	GLY	5.9
1	A	514	ALA	5.8
1	A	512	GLY	5.8
1	A	733	LEU	5.4
1	A	517	LEU	5.3
1	A	732	PHE	5.2
1	A	515	ASN	4.9
1	A	719	ASN	4.8
1	A	770	GLY	4.8
1	A	402[A]	MET	4.6
1	A	718	LYS	4.6
1	A	470	GLU	4.5
1	A	462[A]	GLU	4.5
1	A	725	ASP	4.3
1	A	782	THR	4.3
1	A	856	GLN	4.2
1	A	518	GLY	4.0
1	A	812	GLU	3.9
1	A	717	SER	3.8
1	A	744	GLN	3.7
1	A	397	GLN	3.5
1	A	742	LYS	3.4
1	A	772	GLU	3.2
1	A	714	GLN	3.2
1	A	726	SER	3.1
1	A	828	THR	3.1
1	A	784	PRO	3.1
1	A	398	PRO	3.0
1	A	769	GLU	2.9
1	A	727	GLN	2.9
1	A	734	ARG	2.8
1	A	757	GLU	2.8
1	A	423[A]	SER	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	778	LYS	2.8
1	A	419	LYS	2.7
1	A	624	ARG	2.7
1	A	743	SER	2.7
1	A	736	TYR	2.7
1	A	567	LYS	2.6
1	A	660	ALA	2.5
1	A	687[A]	GLN	2.4
1	A	774	GLU	2.4
1	A	785	LYS	2.4
1	A	623	ARG	2.4
1	A	432[A]	LYS	2.4
1	A	783	LEU	2.3
1	A	430[A]	GLN	2.3
1	A	428[A]	LYS	2.3
1	A	635[A]	GLU	2.3
1	A	781	GLU	2.3
1	A	730	ILE	2.3
1	A	633	GLU	2.2
1	A	665	TYR	2.2
1	A	763	GLN	2.2
1	A	775	LEU	2.1
1	A	519	ASN	2.1
1	A	768	GLU	2.1
1	A	612	ASP	2.1
1	A	787	LYS	2.1
1	A	401[A]	ASP	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 oligosaccharides 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, 95th 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 ²)	Q<0.9
2	RWV	A	901	15/15	0.90	0.15	20,24,32,34	15
3	DMS	A	904	4/4	0.98	0.07	13,14,14,15	0
3	DMS	A	903	4/4	0.99	0.06	16,16,19,19	0
3	DMS	A	902	4/4	0.99	0.07	16,16,16,16	0

6.5 Other polymers [i](#)

There are no such residues in this entry.