



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 07:24 AM GMT

PDB ID : 3APC  
Title : Crystal structure of human PI3K-gamma in complex with CH5132799  
Authors : Nakamura, M.; Fukami, T.A.; Miyazaki, T.; Yoshida, M.  
Deposited on : 2010-10-14  
Resolution : 2.54 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  
<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

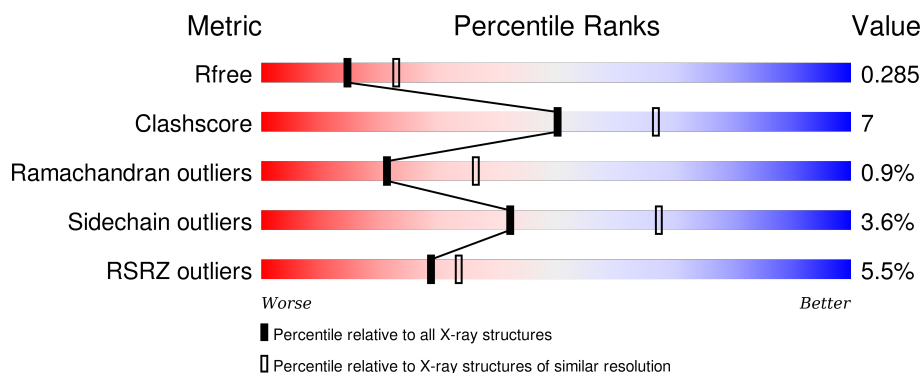
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.54 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	4549 (2.58-2.50)
Clashscore	102246	5292 (2.58-2.50)
Ramachandran outliers	100387	5194 (2.58-2.50)
Sidechain outliers	100360	5196 (2.58-2.50)
RSRZ outliers	91569	4561 (2.58-2.50)

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	966	

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6652 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.

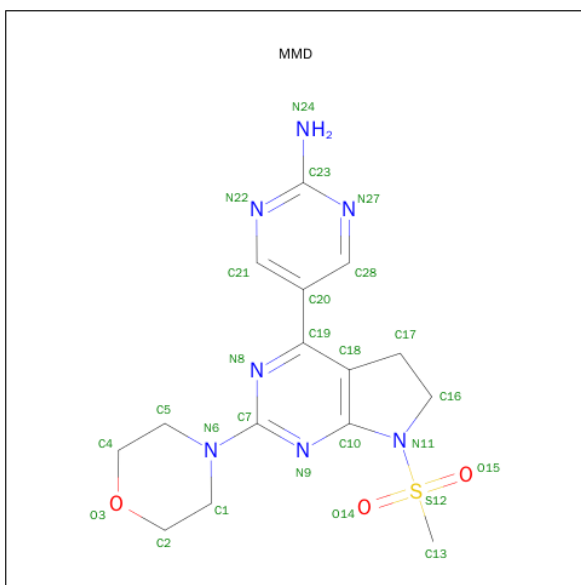
- Molecule 1 is a protein called Phosphatidylinositol-4,5-bisphosphate 3-kinase catalytic subunit gamma isoform.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	839	Total	C	N	O	S	0	0	0
			6594	4236	1114	1210	34			

There are 7 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	137	GLY	-	EXPRESSION TAG	UNP P48736
A	138	PRO	-	EXPRESSION TAG	UNP P48736
A	139	LEU	-	EXPRESSION TAG	UNP P48736
A	140	HIS	-	EXPRESSION TAG	UNP P48736
A	141	MET	-	EXPRESSION TAG	UNP P48736
A	142	GLY	-	EXPRESSION TAG	UNP P48736
A	143	SER	-	EXPRESSION TAG	UNP P48736

- Molecule 2 is 5-(7-METHANESULFONYL-2-MORPHOLIN-4-YL-6,7-DIHYDRO-5H-PYRROLO[2,3-D]PYRIMIDIN-4-YL)-PYRIMIDIN-2-YLAMINE (three-letter code: MMD) (formula: C<sub>15</sub>H<sub>19</sub>N<sub>7</sub>O<sub>3</sub>S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	S	0	0
			26	15	7	3	1		

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



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

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	22	Total	O	0	0
			22	22		



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	142.91Å 67.55Å 106.49Å 90.00° 95.74° 90.00°	Depositor
Resolution (Å)	51.92 – 2.54 51.92 – 2.54	Depositor EDS
% Data completeness (in resolution range)	100.0 (51.92-2.54) 99.7 (51.92-2.54)	Depositor EDS
$R_{merge}$	0.04	Depositor
$R_{sym}$	0.04	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.44 (at 2.55Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, $R_{free}$	0.236 , 0.290 0.229 , 0.285	Depositor DCC
$R_{free}$ test set	1688 reflections (5.33%)	DCC
Wilson B-factor (Å <sup>2</sup> )	63.9	Xtriage
Anisotropy	0.084	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 54.9	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 33352 reflections	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	6652	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	60.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.92% 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: MMD, 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	0.39	0/6736	0.53	0/9146

There are no bond length outliers.

There are no bond angle outliers.

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	6594	0	6435	89	0
2	A	26	0	19	3	0
3	A	10	0	0	0	0
4	A	22	0	0	0	0
All	All	6652	0	6454	91	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 7.

All (91) 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:729:LEU:O	1:A:733:THR:HG23	1.61	0.98
1:A:953:MET:HE2	1:A:963:ILE:HD13	1.58	0.86
1:A:227:SER:O	1:A:228:THR:HG22	1.79	0.83
1:A:862:LEU:HD11	1:A:1016:ALA:HB2	1.61	0.81
1:A:564:LEU:HD11	1:A:1048:ILE:HG22	1.63	0.81
1:A:1009:PHE:O	1:A:1012:ILE:HG22	1.81	0.80
1:A:761:SER:O	1:A:764:ILE:HG22	1.84	0.77
1:A:1090:LEU:O	1:A:1091:VAL:HG23	1.86	0.75
1:A:163:THR:O	1:A:165:VAL:HG23	1.86	0.75
1:A:843:LEU:HD13	1:A:1034:MET:HG3	1.72	0.69
1:A:165:VAL:HG12	1:A:165:VAL:O	1.91	0.69
1:A:168:VAL:CG1	1:A:170:ASP:O	2.43	0.67
1:A:983:VAL:HG22	1:A:984:PRO:HD2	1.77	0.65
1:A:168:VAL:HG13	1:A:170:ASP:O	1.99	0.63
1:A:1048:ILE:O	1:A:1051:ILE:HG22	2.00	0.62
1:A:583:LEU:HD22	1:A:610:LEU:HD12	1.80	0.62
1:A:899:THR:HG23	1:A:1087:PHE:CZ	2.34	0.62
1:A:363:VAL:HG23	1:A:520:LEU:CD2	2.30	0.62
2:A:1:MMD:C17	2:A:1:MMD:H28	2.31	0.61
1:A:899:THR:HG23	1:A:1087:PHE:CE1	2.36	0.60
2:A:1:MMD:H17A	2:A:1:MMD:H28	1.83	0.60
1:A:732:PHE:O	1:A:736:VAL:HG23	2.01	0.60
1:A:1059:LYS:O	1:A:1060:ASN:O	2.20	0.59
1:A:622:LEU:HD13	1:A:647:LYS:HB3	1.88	0.56
1:A:1010:GLN:O	1:A:1014:VAL:HG23	2.06	0.56
1:A:862:LEU:HD11	1:A:1016:ALA:CB	2.36	0.55
1:A:742:LEU:HD22	1:A:813:LEU:HD11	1.89	0.55
1:A:583:LEU:HD22	1:A:610:LEU:CD1	2.36	0.55
1:A:168:VAL:HG13	1:A:170:ASP:H	1.73	0.53
1:A:749:ILE:HG21	1:A:803:VAL:HG22	1.90	0.53
1:A:1040:PRO:O	1:A:1041:GLN:CB	2.55	0.53
1:A:858:GLU:OE2	1:A:1022:HIS:ND1	2.40	0.52
1:A:224:ILE:HD12	1:A:248:PHE:CD2	2.44	0.52
1:A:953:MET:CE	1:A:963:ILE:HD13	2.36	0.51
1:A:564:LEU:CD1	1:A:1048:ILE:HG22	2.39	0.51
1:A:176:THR:HG23	1:A:674:ASP:HB2	1.93	0.51
1:A:467:LEU:HD13	1:A:672:TYR:CE1	2.45	0.51
1:A:519:LEU:HD12	1:A:520:LEU:N	2.25	0.50
1:A:662:GLN:NE2	1:A:1030:LEU:HD22	2.27	0.50
1:A:739:ILE:HD12	1:A:739:ILE:C	2.32	0.50
1:A:519:LEU:HD12	1:A:520:LEU:H	1.76	0.50
1:A:749:ILE:HG21	1:A:803:VAL:CG2	2.41	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:363:VAL:HG23	1:A:520:LEU:HD23	1.93	0.49
1:A:860:LEU:HD11	1:A:1015:LYS:HB3	1.95	0.49
1:A:596:VAL:HG13	1:A:603:ILE:HG22	1.95	0.49
1:A:775:GLN:HE21	1:A:775:GLN:HA	1.78	0.49
1:A:620:SER:O	1:A:647:LYS:NZ	2.45	0.49
1:A:363:VAL:HG23	1:A:520:LEU:HD21	1.95	0.48
1:A:559:ILE:HG23	1:A:588:ALA:HB2	1.94	0.48
1:A:844:ILE:HD13	1:A:965:PHE:HB3	1.94	0.48
1:A:274:VAL:HG21	1:A:292:TRP:CD1	2.48	0.48
1:A:727:ALA:O	1:A:730:HIS:HB3	2.14	0.48
1:A:862:LEU:N	1:A:862:LEU:HD23	2.29	0.47
1:A:231:GLN:O	1:A:233:ILE:HD12	2.14	0.47
1:A:198:MET:SD	1:A:282:VAL:HG21	2.54	0.47
1:A:930:TYR:CE1	1:A:1012:ILE:HD13	2.50	0.47
1:A:1070:ASP:O	1:A:1074:VAL:HG23	2.15	0.47
1:A:863:CYS:SG	1:A:927:ARG:NH1	2.89	0.46
1:A:854:ILE:HG23	1:A:1023:HIS:CD2	2.50	0.46
1:A:471:HIS:H	1:A:471:HIS:CD2	2.34	0.46
1:A:428:LEU:HD22	1:A:465:ASN:HB3	1.98	0.45
1:A:933:ALA:CB	1:A:1012:ILE:HG23	2.47	0.45
1:A:953:MET:CE	1:A:963:ILE:HG21	2.47	0.45
1:A:738:VAL:HG12	1:A:742:LEU:HD12	2.00	0.44
1:A:742:LEU:HD22	1:A:813:LEU:CD1	2.48	0.44
1:A:228:THR:HG23	1:A:228:THR:O	2.18	0.44
1:A:168:VAL:HG11	1:A:170:ASP:O	2.18	0.44
1:A:199:HIS:O	1:A:199:HIS:ND1	2.50	0.44
1:A:157:LEU:HD21	1:A:733:THR:HG22	2.00	0.43
1:A:933:ALA:HB1	1:A:1012:ILE:HG23	1.99	0.43
1:A:596:VAL:HG13	1:A:603:ILE:CG2	2.49	0.43
1:A:500:ASP:HB3	1:A:708:HIS:CD2	2.54	0.43
1:A:740:GLU:HA	1:A:743:GLN:HG2	2.00	0.43
1:A:838:LEU:HD12	1:A:877:GLY:CA	2.49	0.43
1:A:617:TRP:O	1:A:647:LYS:NZ	2.46	0.42
1:A:843:LEU:CD1	1:A:1034:MET:HG3	2.47	0.42
1:A:831:ILE:HB	1:A:879:ILE:HB	2.02	0.42
1:A:838:LEU:HD12	1:A:877:GLY:HA3	2.00	0.42
1:A:997:THR:CG2	1:A:998:SER:N	2.82	0.42
1:A:879:ILE:HD13	2:A:1:MMD:C21	2.50	0.42
1:A:184:ARG:HD3	1:A:719:ALA:O	2.19	0.42
1:A:793:ALA:HA	1:A:828:ILE:HD12	2.01	0.41
1:A:853:SER:O	1:A:857:THR:HG23	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:799:GLU:HG2	1:A:800:LYS:N	2.36	0.41
1:A:463:TYR:CE2	1:A:501:LYS:HA	2.56	0.41
1:A:366:ARG:NH1	1:A:519:LEU:HD22	2.35	0.41
1:A:227:SER:O	1:A:228:THR:CG2	2.60	0.41
1:A:956:GLU:O	1:A:957:THR:HG23	2.21	0.41
1:A:891:ILE:HG22	1:A:906:VAL:HG12	2.03	0.40
1:A:997:THR:HG22	1:A:998:SER:N	2.36	0.40
1:A:984:PRO:HG3	1:A:1071:GLN:HE21	1.87	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	823/966 (85%)	780 (95%)	36 (4%)	7 (1%)	21	36

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	1060	ASN
1	A	1091	VAL
1	A	509	ASP
1	A	1041	GLN
1	A	964	ASP
1	A	164	ASP
1	A	1001	LYS

### 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	702/862 (81%)	677 (96%)	25 (4%)	42 67

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	178	ARG
1	A	219	CYS
1	A	226	ARG
1	A	228	THR
1	A	373	LEU
1	A	404	PHE
1	A	501	LYS
1	A	643	ILE
1	A	707	ARG
1	A	739	ILE
1	A	743	GLN
1	A	775	GLN
1	A	776	ASN
1	A	799	GLU
1	A	841	ASP
1	A	853	SER
1	A	876	ILE
1	A	893	GLN
1	A	908	ASN
1	A	915	SER
1	A	922	GLN
1	A	957	THR
1	A	983	VAL
1	A	1003	SER
1	A	1061	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	396	GLN
1	A	471	HIS
1	A	662	GLN
1	A	743	GLN

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Mol	Chain	Res	Type
1	A	775	GLN
1	A	776	ASN
1	A	893	GLN
1	A	898	ASN
1	A	922	GLN
1	A	959	ASN
1	A	1023	HIS
1	A	1071	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](#)

3 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	$\# Z  > 2$
2	MMD	A	1	-	28,29,29	2.27	4 (14%)	34,43,43	2.85	17 (50%)
3	SO4	A	1103	-	4,4,4	0.16	0	6,6,6	0.21	0
3	SO4	A	2	-	4,4,4	0.21	0	6,6,6	0.14	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
2	MMD	A	1	-	-	0/14/31/31	0/4/4/4
3	SO4	A	1103	-	-	0/0/0/0	0/0/0/0
3	SO4	A	2	-	-	0/0/0/0	0/0/0/0

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1	MMD	C20-C19	-7.12	1.41	1.49
2	A	1	MMD	C18-C10	-3.29	1.39	1.46
2	A	1	MMD	O15-S12	5.81	1.49	1.43
2	A	1	MMD	O14-S12	5.97	1.50	1.43

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1	MMD	O15-S12-O14	-6.91	109.05	118.66
2	A	1	MMD	N27-C23-N22	-4.47	120.08	124.58
2	A	1	MMD	C18-C19-N8	-4.04	117.27	122.73
2	A	1	MMD	N9-C7-N8	-3.15	121.05	126.35
2	A	1	MMD	C20-C28-N27	-3.01	119.46	124.34
2	A	1	MMD	C28-C20-C21	2.17	117.22	114.25
2	A	1	MMD	C20-C19-N8	2.21	118.20	115.17
2	A	1	MMD	N9-C7-N6	2.35	119.72	117.08
2	A	1	MMD	C18-C10-N11	2.39	110.17	107.65
2	A	1	MMD	N24-C23-N27	2.83	119.94	117.39
2	A	1	MMD	N24-C23-N22	2.87	119.98	117.39
2	A	1	MMD	C21-N22-C23	3.02	120.22	116.05
2	A	1	MMD	C7-N9-C10	3.32	120.39	115.37
2	A	1	MMD	C28-N27-C23	4.20	121.84	116.05
2	A	1	MMD	C5-N6-C1	4.43	120.90	111.59
2	A	1	MMD	O14-S12-N11	4.91	112.63	107.66
2	A	1	MMD	N9-C10-N11	5.53	130.61	126.27

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1	MMD	3	0

## 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	839/966 (86%)	0.36	46 (5%) 29 33	25, 58, 94, 114	0

All (46) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	374	PRO	7.0
1	A	216	ALA	6.2
1	A	221	PHE	5.7
1	A	215	ILE	5.6
1	A	376	ASN	5.0
1	A	220	ILE	4.9
1	A	212	TRP	4.7
1	A	234	LYS	4.3
1	A	1064	ALA	4.2
1	A	375	ARG	4.1
1	A	968	ILE	3.9
1	A	377	THR	3.9
1	A	149	ALA	3.8
1	A	373	LEU	3.8
1	A	211	LEU	3.7
1	A	248	PHE	3.7
1	A	270	PHE	3.6
1	A	378	ASP	3.4
1	A	300	GLY	3.4
1	A	895	THR	3.3
1	A	217	ASN	3.3
1	A	379	LEU	3.3
1	A	207	LEU	3.2
1	A	219	CYS	3.2
1	A	250	THR	3.2
1	A	383	VAL	3.2
1	A	233	ILE	3.1

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Mol	Chain	Res	Type	RSRZ
1	A	404	PHE	3.1
1	A	901	ALA	2.9
1	A	996	GLY	2.8
1	A	896	VAL	2.6
1	A	287	ILE	2.5
1	A	235	VAL	2.5
1	A	232	THR	2.4
1	A	249	PHE	2.4
1	A	1040	PRO	2.3
1	A	222	ILE	2.3
1	A	823	LEU	2.2
1	A	902	PHE	2.2
1	A	825	ASN	2.2
1	A	253	ALA	2.1
1	A	869	CYS	2.1
1	A	370	ILE	2.1
1	A	892	GLN	2.0
1	A	321	GLU	2.0
1	A	911	LEU	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. 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( $\text{\AA}^2$ )	Q<0.9
3	SO4	A	2	5/5	0.95	0.15	-0.28	81,85,88,91	0
2	MMD	A	1	26/26	0.96	0.15	-0.93	38,50,53,54	0
3	SO4	A	1103	5/5	0.96	0.07	-1.25	68,71,74,75	0

## 6.5 Other polymers

There are no such residues in this entry.