



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 10:33 am GMT

PDB ID : 4FA6
Title : Design and Synthesis of a Novel Pyrrolidinyl Pyrido Pyrimidinone Derivative
as a Potent Inhibitor of PI3Ka and mTOR
Authors : Pannifer, A.; Greasley, S.E.
Deposited on : 2012-05-21
Resolution : 2.70 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

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : trunk28620
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
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) : recalc28949

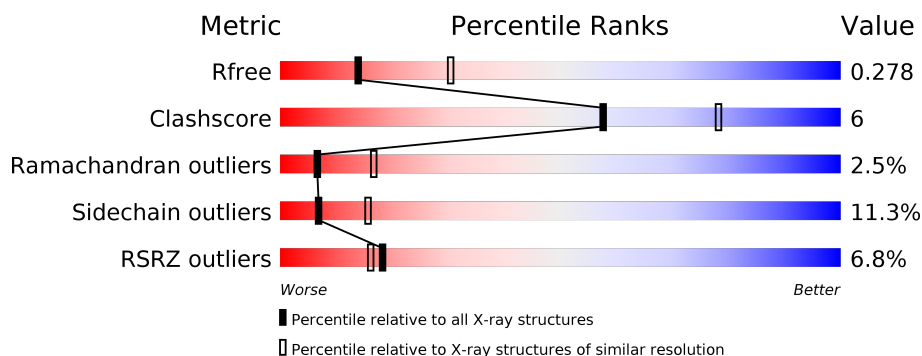
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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}	100719	2259 (2.70-2.70)
Clashscore	112137	2590 (2.70-2.70)
Ramachandran outliers	110173	2550 (2.70-2.70)
Sidechain outliers	110143	2550 (2.70-2.70)
RSRZ outliers	101464	2275 (2.70-2.70)

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 ≥ 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	<div> <div>6%</div> <div>68%</div> <div>17%</div> <div>•</div> <div>13%</div> </div>

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 6863 atoms, of which 18 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.

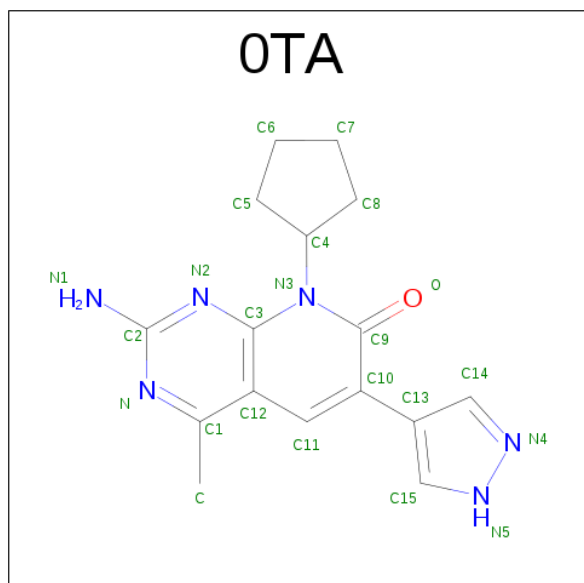
- Molecule 1 is a protein called Phosphatidylinositol 4,5-bisphosphate 3-kinase catalytic sub-unit gamma isoform.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	842	Total	C	N	O	S	0	0	0
			6818	4375	1165	1242	36			

There are 7 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	143	MET	-	EXPRESSION TAG	UNP P48736
A	1103	HIS	-	EXPRESSION TAG	UNP P48736
A	1104	HIS	-	EXPRESSION TAG	UNP P48736
A	1105	HIS	-	EXPRESSION TAG	UNP P48736
A	1106	HIS	-	EXPRESSION TAG	UNP P48736
A	1107	HIS	-	EXPRESSION TAG	UNP P48736
A	1108	HIS	-	EXPRESSION TAG	UNP P48736

- Molecule 2 is 2-AMINO-8-CYCLOPENTYL-4-METHYL-6-(1H-PYRAZOL-4-YL)PYRIDO [2,3-D]PYRIMIDIN-7(8H)-ONE (three-letter code: OTA) (formula: C₁₆H₁₈N₆O).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	H	N	O	0	0
			41	16	18	6	1		

- Molecule 3 is water.

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

4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	143.44Å 68.24Å 106.17Å 90.00° 94.82° 90.00°	Depositor
Resolution (Å)	29.85 – 2.70 29.86 – 2.70	Depositor EDS
% Data completeness (in resolution range)	96.7 (29.85-2.70) 96.7 (29.86-2.70)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.87 (at 2.68Å)	Xtriage
Refinement program	BUSTER 2.11.1	Depositor
R, R_{free}	0.205 , 0.257 0.217 , 0.278	Depositor DCC
R_{free} test set	1384 reflections (5.32%)	DCC
Wilson B-factor (Å ²)	63.7	Xtriage
Anisotropy	0.343	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 53.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6863	wwPDB-VP
Average B, all atoms (Å ²)	74.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.06% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: 0TA

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	0.51	0/6965	0.74	0/9423

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

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	6818	0	6850	77	0
2	A	23	18	18	1	0
3	A	4	0	0	0	0
All	All	6845	18	6868	77	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 6.

All (77) 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:210:TYR:O	1:A:213:LYS:HE3	1.64	0.95

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:213:LYS:NZ	1:A:214:LYS:HG3	1.87	0.88
1:A:213:LYS:HZ3	1:A:214:LYS:HG3	1.42	0.81
1:A:939:THR:HG23	1:A:945:GLY:HA2	1.69	0.72
1:A:1045:LYS:HG2	1:A:1049:GLU:CD	2.14	0.68
1:A:1045:LYS:HZ2	1:A:1045:LYS:HB2	1.57	0.68
1:A:808:LYS:O	1:A:810:PRO:HD3	1.94	0.68
1:A:916:PRO:HD2	1:A:920:LYS:HD3	1.77	0.66
1:A:214:LYS:O	1:A:215:ILE:HG13	1.99	0.63
1:A:1045:LYS:NZ	1:A:1045:LYS:HB2	2.14	0.62
1:A:523:TYR:O	1:A:524:CYS:HB2	1.99	0.62
1:A:1045:LYS:NZ	1:A:1045:LYS:CB	2.65	0.59
1:A:905:GLU:HG3	1:A:993:PHE:CE2	2.37	0.59
1:A:743:GLN:HG2	1:A:876:ILE:HD12	1.84	0.59
1:A:949:ASN:H	1:A:1083:GLN:HE22	1.51	0.59
1:A:213:LYS:HZ2	1:A:214:LYS:HG3	1.66	0.58
1:A:272:LEU:HB3	1:A:305:VAL:HG11	1.86	0.58
1:A:214:LYS:C	1:A:215:ILE:HG13	2.24	0.58
1:A:214:LYS:HZ1	1:A:300:GLY:HA2	1.69	0.57
1:A:935:TYR:O	1:A:939:THR:HB	2.04	0.57
1:A:724:CYS:HB2	1:A:728:MET:HE3	1.87	0.56
1:A:579:ARG:HD2	1:A:610:LEU:HD11	1.87	0.56
1:A:887:THR:HG22	1:A:890:LYS:H	1.72	0.54
1:A:807:LYS:H	1:A:807:LYS:HE3	1.73	0.54
1:A:205:LYS:O	1:A:206:PRO:O	2.26	0.54
1:A:222:ILE:HD12	1:A:235:VAL:HG21	1.90	0.54
1:A:206:PRO:O	1:A:294:ARG:NH2	2.40	0.53
1:A:838:LEU:HD23	1:A:877:GLY:HA3	1.90	0.53
1:A:947:ARG:HH21	1:A:951:ASN:HB3	1.74	0.52
1:A:207:LEU:CD2	1:A:211:LEU:HB3	2.40	0.52
1:A:925:VAL:O	1:A:929:VAL:HG23	2.09	0.52
1:A:1045:LYS:O	1:A:1049:GLU:HG3	2.09	0.52
1:A:1035:LEU:HA	1:A:1039:MET:HG2	1.93	0.49
1:A:882:VAL:HG23	2:A:1201:OTA:H2	1.92	0.49
1:A:889:ALA:HB2	1:A:949:ASN:HB3	1.94	0.49
1:A:558:ILE:O	1:A:561:THR:HG22	2.12	0.49
1:A:905:GLU:HG2	1:A:909:HIS:CE1	2.48	0.49
1:A:207:LEU:HD21	1:A:211:LEU:HB3	1.95	0.49
1:A:860:LEU:HD11	1:A:1015:LYS:HB3	1.93	0.49
1:A:212:TRP:C	1:A:214:LYS:H	2.16	0.48
1:A:215:ILE:HA	1:A:217:ASN:OD1	2.12	0.48
1:A:831:ILE:HB	1:A:879:ILE:HB	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1043:THR:C	1:A:1045:LYS:H	2.18	0.47
1:A:214:LYS:NZ	1:A:300:GLY:HA2	2.28	0.47
1:A:370:ILE:HD13	1:A:371:PRO:HD2	1.96	0.46
1:A:948:HIS:N	1:A:948:HIS:ND1	2.62	0.46
1:A:706:SER:O	1:A:710:GLN:HB3	2.16	0.46
1:A:834:HIS:HB2	1:A:876:ILE:HG12	1.97	0.45
1:A:1067:TYR:O	1:A:1071:GLN:HG2	2.17	0.44
1:A:274:VAL:HG21	1:A:292:TRP:CD1	2.53	0.44
1:A:422:ASP:HB3	1:A:599:GLY:O	2.18	0.44
1:A:629:GLN:HG2	1:A:1029:ILE:HG13	1.99	0.43
1:A:368:ILE:HG21	1:A:433:ILE:HD11	2.00	0.43
1:A:172:GLU:HG3	1:A:471:HIS:CG	2.53	0.43
1:A:391:GLN:HE21	1:A:502:LEU:HD21	1.83	0.43
1:A:841:ASP:O	1:A:845:LEU:HD22	2.19	0.43
1:A:410:TRP:HB3	1:A:412:VAL:HG22	2.01	0.43
1:A:651:LEU:HD22	1:A:655:ASP:HB3	2.01	0.43
1:A:211:LEU:HD23	1:A:211:LEU:HA	1.86	0.43
1:A:471:HIS:H	1:A:471:HIS:HD1	1.66	0.43
1:A:912:LYS:HE2	1:A:918:GLU:HG2	2.01	0.43
1:A:287:ILE:HG13	1:A:287:ILE:H	1.74	0.42
1:A:365:ILE:HD13	1:A:518:ILE:HG22	2.02	0.42
1:A:406:GLU:H	1:A:406:GLU:HG2	1.68	0.42
1:A:848:LEU:HD12	1:A:851:MET:HE3	2.02	0.42
1:A:363:VAL:HG12	1:A:416:PHE:HE1	1.85	0.41
1:A:370:ILE:HD12	1:A:372:VAL:O	2.20	0.41
1:A:808:LYS:H	1:A:808:LYS:HG2	1.74	0.41
1:A:731:ASP:O	1:A:735:GLN:HG3	2.20	0.41
1:A:629:GLN:CG	1:A:1029:ILE:HG13	2.51	0.41
1:A:819:ASP:HA	1:A:820:PRO:HD3	2.00	0.41
1:A:480:TYR:HB2	1:A:518:ILE:HG13	2.01	0.41
1:A:565:ASN:HA	1:A:566:PRO:HD3	1.89	0.41
1:A:500:ASP:O	1:A:503:THR:HG22	2.21	0.41
1:A:614:ARG:HB3	1:A:618:ASP:OD2	2.21	0.41
1:A:180:LEU:C	1:A:183:PRO:HD2	2.42	0.40
1:A:176:THR:HG23	1:A:674:ASP:HB2	2.04	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	828/966 (86%)	762 (92%)	45 (5%)	21 (2%)	6	17

All (21) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	206	PRO
1	A	377	THR
1	A	615	GLU
1	A	896	VAL
1	A	898	ASN
1	A	899	THR
1	A	213	LYS
1	A	215	ILE
1	A	524	CYS
1	A	1040	PRO
1	A	210	TYR
1	A	375	ARG
1	A	758	ASP
1	A	1000	LYS
1	A	1045	LYS
1	A	1046	GLU
1	A	374	PRO
1	A	874	ASP
1	A	1044	SER
1	A	964	ASP
1	A	759	VAL

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	755/864 (87%)	670 (89%)	85 (11%)	7 16

All (85) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	168	VAL
1	A	202	VAL
1	A	205	LYS
1	A	207	LEU
1	A	213	LYS
1	A	214	LYS
1	A	215	ILE
1	A	226	ARG
1	A	229	THR
1	A	230	SER
1	A	232	THR
1	A	236	SER
1	A	285	THR
1	A	302	GLU
1	A	307	LEU
1	A	309	THR
1	A	320	LYS
1	A	358	ASP
1	A	366	ARG
1	A	369	ASP
1	A	370	ILE
1	A	373	LEU
1	A	378	ASP
1	A	379	LEU
1	A	381	VAL
1	A	387	ILE
1	A	391	GLN
1	A	395	CYS
1	A	406	GLU
1	A	470	ASP
1	A	498	ASN
1	A	521	ASP
1	A	524	CYS
1	A	561	THR
1	A	570	GLU

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Mol	Chain	Res	Type
1	A	574	LEU
1	A	575	LEU
1	A	587	LYS
1	A	601	GLN
1	A	609	GLN
1	A	626	LEU
1	A	646	GLN
1	A	647	LYS
1	A	649	GLU
1	A	682	LEU
1	A	701	SER
1	A	707	ARG
1	A	717	LEU
1	A	728	MET
1	A	767	LEU
1	A	775	GLN
1	A	798	ILE
1	A	807	LYS
1	A	808	LYS
1	A	823	LEU
1	A	825	ASN
1	A	838	LEU
1	A	843	LEU
1	A	845	LEU
1	A	848	LEU
1	A	876	ILE
1	A	878	MET
1	A	886	THR
1	A	887	THR
1	A	895	THR
1	A	896	VAL
1	A	907	LEU
1	A	918	GLU
1	A	927	ARG
1	A	939	THR
1	A	948	HIS
1	A	981	GLU
1	A	982	ARG
1	A	997	THR
1	A	1002	THR
1	A	1007	GLN
1	A	1026	LEU

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Mol	Chain	Res	Type
1	A	1029	ILE
1	A	1042	LEU
1	A	1048	ILE
1	A	1059	LYS
1	A	1078	LYS
1	A	1085	ASN
1	A	1087	PHE
1	A	1088	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	304	HIS
1	A	376	ASN
1	A	391	GLN
1	A	549	ASN
1	A	554	GLN
1	A	600	GLN
1	A	705	GLN
1	A	834	HIS
1	A	951	ASN
1	A	1041	GLN
1	A	1083	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](#)

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	0TA	A	1201	-	21,26,26	1.69	2 (9%)	21,38,38	3.15	7 (33%)

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	0TA	A	1201	-	-	0/2/15/15	0/4/4/4

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1201	0TA	C1-N	2.06	1.34	1.32
2	A	1201	0TA	C9-N3	6.77	1.48	1.38

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1201	0TA	C9-C10-C13	-2.84	119.07	122.34
2	A	1201	0TA	C2-N2-C3	2.16	117.68	115.16
2	A	1201	0TA	C11-C12-C3	2.32	119.76	117.01
2	A	1201	0TA	C11-C10-C9	2.55	120.17	117.67
2	A	1201	0TA	C10-C9-N3	3.97	118.87	116.11
2	A	1201	0TA	C5-C4-C8	4.04	108.86	104.11
2	A	1201	0TA	C15-C13-C10	11.83	131.55	127.37

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1201	OTA	1	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, 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	842/966 (87%)	0.23	57 (6%) 18 16	38, 71, 117, 147	0

All (57) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	895	THR	7.8
1	A	253	ALA	7.0
1	A	1044	SER	6.0
1	A	216	ALA	5.8
1	A	754	ALA	5.7
1	A	254	LYS	5.6
1	A	374	PRO	5.5
1	A	1092	LEU	5.3
1	A	757	TYR	5.2
1	A	376	ASN	5.0
1	A	896	VAL	4.9
1	A	458	VAL	4.2
1	A	1043	THR	4.1
1	A	525	HIS	3.9
1	A	375	ARG	3.8
1	A	377	THR	3.7
1	A	1086	TRP	3.6
1	A	758	ASP	3.5
1	A	986	VAL	3.5
1	A	1091	VAL	3.5
1	A	755	GLU	3.4
1	A	613	ARG	3.4
1	A	523	TYR	3.4
1	A	985	PHE	3.4
1	A	776	ASN	3.3
1	A	1000	LYS	3.2
1	A	221	PHE	3.1

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Mol	Chain	Res	Type	RSRZ
1	A	544	ARG	3.1
1	A	234	LYS	3.0
1	A	522	ASN	3.0
1	A	373	LEU	2.9
1	A	378	ASP	2.7
1	A	756	LYS	2.6
1	A	250	THR	2.6
1	A	248	PHE	2.5
1	A	147	SER	2.5
1	A	936	CYS	2.5
1	A	1041	GLN	2.5
1	A	779	LEU	2.5
1	A	967	HIS	2.4
1	A	1087	PHE	2.4
1	A	1077	ASP	2.4
1	A	226	ARG	2.4
1	A	320	LYS	2.4
1	A	778	GLN	2.3
1	A	526	PRO	2.2
1	A	898	ASN	2.2
1	A	984	PRO	2.2
1	A	252	MET	2.2
1	A	993	PHE	2.2
1	A	913	GLU	2.2
1	A	406	GLU	2.1
1	A	270	PHE	2.1
1	A	1089	HIS	2.1
1	A	524	CYS	2.1
1	A	916	PRO	2.0
1	A	1062	GLU	2.0

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, 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	LLDF	B-factors(\AA^2)	Q<0.9
2	0TA	A	1201	23/23	0.96	0.17	0.07	59,68,72,72	0

6.5 Other polymers [i](#)

There are no such residues in this entry.