



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 11:57 AM GMT

PDB ID : 3QJZ
Title : Crystal structure of PI3K-gamma in complex with benzothiazole 1
Authors : Whittington, D.A.; Tang, J.; Yakowec, P.
Deposited on : 2011-01-31
Resolution : 2.90 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 (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

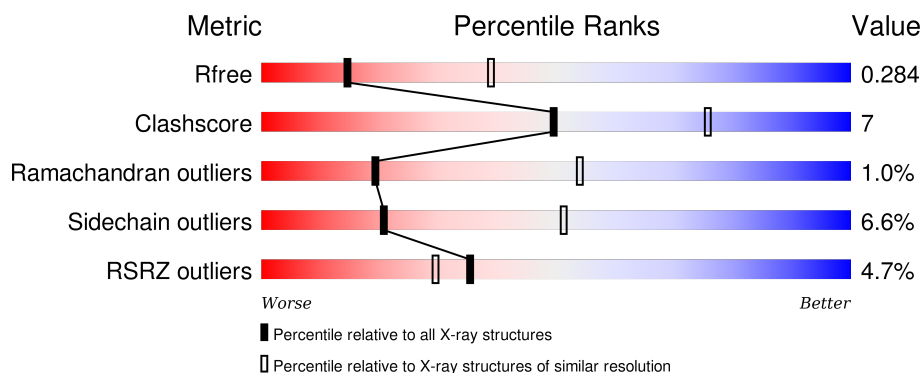
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION


The reported resolution of this entry is 2.90 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	1451 (2.90-2.90)
Clashscore	102246	1668 (2.90-2.90)
Ramachandran outliers	100387	1630 (2.90-2.90)
Sidechain outliers	100360	1632 (2.90-2.90)
RSRZ outliers	91569	1456 (2.90-2.90)

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	960	

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 6487 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.

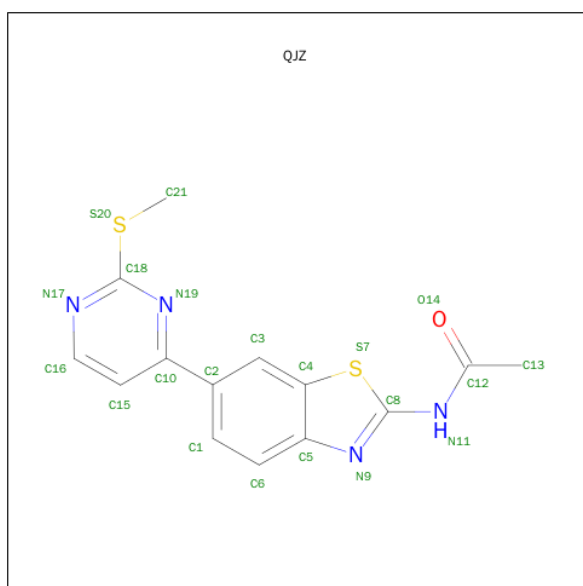
- Molecule 1 is a protein called Phosphatidylinositol-4,5-bisphosphate 3-kinase catalytic subunit gamma isoform.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	795	Total	C	N	O	S	0	0	0
			6458	4153	1100	1171	34			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	143	GLY	-	EXPRESSION TAG	UNP P48736

- Molecule 2 is N-{6-[2-(METHYLSULFANYL)PYRIMIDIN-4-YL]-1,3-BENZOTHIAZOL-2-YL}ACETAMIDE (three-letter code: QJZ) (formula: C₁₄H₁₂N₄OS₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	S	0	0
			21	14	4	1	2		

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



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

- Molecule 4 is water.

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

4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	141.51Å 66.87Å 105.72Å 90.00° 96.24° 90.00°	Depositor
Resolution (Å)	30.00 – 2.90 29.39 – 2.90	Depositor EDS
% Data completeness (in resolution range)	96.2 (30.00-2.90) 96.2 (29.39-2.90)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.18 (at 2.90Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.230 , 0.291 0.221 , 0.284	Depositor DCC
R_{free} test set	1248 reflections (6.27%)	DCC
Wilson B-factor (Å ²)	78.4	Xtriage
Anisotropy	0.083	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 47.0	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 21161 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	6487	wwPDB-VP
Average B, all atoms (Å ²)	81.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.04% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: QJZ, 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	0/6593	0.56	0/8909

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	6458	0	6496	92	0
2	A	21	0	12	1	0
3	A	5	0	0	0	0
4	A	3	0	0	0	0
All	All	6487	0	6508	92	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 (92) 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:784:ARG:HG3	1:A:784:ARG:HH11	1.29	0.97
1:A:613:ARG:HH11	1:A:613:ARG:HG3	1.39	0.88
1:A:359:ARG:HG3	1:A:359:ARG:HH11	1.37	0.86
1:A:808:LYS:O	1:A:810:PRO:HD3	1.79	0.83
1:A:366:ARG:HD2	1:A:479:GLU:OE2	1.81	0.80
1:A:887:THR:HG22	1:A:889:ALA:H	1.53	0.73
1:A:167:ASN:HD22	1:A:711:GLN:HE22	1.39	0.70
1:A:651:LEU:HD22	1:A:655:ASP:HB3	1.73	0.70
1:A:792:LYS:HB3	1:A:818:ALA:HB3	1.76	0.68
1:A:629:GLN:HG2	1:A:1029:ILE:HG13	1.76	0.66
1:A:564:LEU:HD11	1:A:1048:ILE:HG22	1.77	0.66
1:A:939:THR:HG23	1:A:945:GLY:HA2	1.79	0.64
1:A:784:ARG:HG3	1:A:784:ARG:NH1	2.05	0.63
1:A:947:ARG:NH2	1:A:963:ILE:O	2.32	0.62
1:A:576:TRP:O	1:A:579:ARG:HB2	2.00	0.62
1:A:955:THR:OG1	1:A:957:THR:HG22	1.99	0.61
1:A:429:LEU:HB2	1:A:468:LEU:HD21	1.82	0.60
1:A:238:ASP:HA	1:A:286:PRO:HB3	1.84	0.60
1:A:178:ARG:HH11	1:A:178:ARG:HG2	1.68	0.59
1:A:359:ARG:CG	1:A:359:ARG:HH11	2.14	0.59
1:A:944:ILE:O	1:A:947:ARG:HD3	2.03	0.58
1:A:498:ASN:HD21	1:A:500:ASP:HB2	1.68	0.58
1:A:767:LEU:HD22	1:A:803:VAL:HG23	1.86	0.57
1:A:370:ILE:HG13	1:A:371:PRO:HD2	1.88	0.56
1:A:888:ILE:HD12	1:A:952:ILE:HG22	1.88	0.55
1:A:705:GLN:HG3	1:A:839:ARG:NH2	2.22	0.55
1:A:273:ARG:NH2	1:A:819:ASP:OD2	2.41	0.54
1:A:887:THR:HG22	1:A:889:ALA:N	2.22	0.53
1:A:925:VAL:O	1:A:929:VAL:HG23	2.09	0.53
1:A:804:MET:HB2	1:A:810:PRO:HD2	1.90	0.52
1:A:730:HIS:O	1:A:734:GLN:HG2	2.10	0.52
1:A:1044:SER:O	1:A:1048:ILE:HG12	2.10	0.51
1:A:1068:PHE:O	1:A:1072:ILE:HG12	2.10	0.51
1:A:1017:TYR:CZ	1:A:1021:ARG:HD2	2.46	0.51
1:A:882:VAL:O	2:A:1:QJZ:H13	2.10	0.50
1:A:381:VAL:HG23	1:A:435:CYS:SG	2.51	0.50
1:A:622:LEU:HA	1:A:647:LYS:HE2	1.92	0.50
1:A:547:MET:HG2	1:A:552:ARG:NH1	2.26	0.50
1:A:466:LEU:HD21	1:A:476:ARG:HG2	1.94	0.50
1:A:815:PHE:O	1:A:827:THR:HB	2.12	0.49
1:A:278:ASP:OD1	1:A:784:ARG:NH2	2.46	0.49
1:A:784:ARG:CG	1:A:784:ARG:NH1	2.72	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1014:VAL:O	1:A:1018:LEU:HG	2.13	0.48
1:A:641:ARG:O	1:A:645:VAL:HG23	2.13	0.48
1:A:410:TRP:HB3	1:A:412:VAL:HG22	1.96	0.48
1:A:768:LYS:O	1:A:772:GLU:HG2	2.13	0.48
1:A:630:LEU:HB2	1:A:644:ALA:HB2	1.96	0.48
1:A:240:THR:HB	1:A:243:ALA:H	1.79	0.47
1:A:476:ARG:HB3	1:A:520:LEU:HD23	1.94	0.47
1:A:498:ASN:ND2	1:A:500:ASP:HB2	2.28	0.47
1:A:359:ARG:HG3	1:A:359:ARG:NH1	2.17	0.47
1:A:614:ARG:HG2	1:A:614:ARG:O	2.15	0.47
1:A:878:MET:O	1:A:879:ILE:HG13	2.15	0.47
1:A:990:ASP:O	1:A:994:VAL:HG23	2.14	0.47
1:A:613:ARG:HH11	1:A:613:ARG:CG	2.18	0.47
1:A:787:TYR:HB3	1:A:870:ILE:HD12	1.97	0.47
1:A:945:GLY:HA3	1:A:984:PRO:O	2.15	0.46
1:A:796:LEU:HG	1:A:815:PHE:CE2	2.50	0.46
1:A:549:ASN:HB3	1:A:550:GLN:HE21	1.80	0.46
1:A:210:TYR:OH	1:A:856:GLU:HG3	2.16	0.45
1:A:202:VAL:CG1	1:A:203:THR:N	2.80	0.45
1:A:833:LYS:O	1:A:876:ILE:HA	2.17	0.45
1:A:661:LEU:O	1:A:665:GLN:HG2	2.17	0.45
1:A:557:ALA:O	1:A:561:THR:HG23	2.18	0.44
1:A:287:ILE:HG13	1:A:287:ILE:H	1.62	0.44
1:A:741:MET:HB3	1:A:774:LEU:HD13	2.00	0.44
1:A:613:ARG:NH1	1:A:613:ARG:HG3	2.17	0.44
1:A:930:TYR:HD1	1:A:1012:ILE:HD13	1.81	0.44
1:A:693:HIS:O	1:A:696:PHE:HB3	2.18	0.43
1:A:808:LYS:HG2	1:A:835:GLY:HA3	1.99	0.43
1:A:963:ILE:HD12	1:A:964:ASP:HB2	2.00	0.43
1:A:365:ILE:HD13	1:A:518:ILE:HG22	2.00	0.43
1:A:472:ARG:O	1:A:473:PHE:HB2	2.18	0.43
1:A:935:TYR:CE1	1:A:961:PHE:HA	2.54	0.43
1:A:872:THR:OG1	1:A:877:GLY:HA2	2.17	0.43
1:A:1082:VAL:O	1:A:1085:ASN:HB2	2.18	0.43
1:A:400:SER:O	1:A:410:TRP:NE1	2.43	0.42
1:A:731:ASP:O	1:A:735:GLN:HG3	2.20	0.42
1:A:202:VAL:HG13	1:A:203:THR:N	2.34	0.41
1:A:645:VAL:HA	1:A:648:LEU:HD12	2.02	0.41
1:A:364:LYS:HB3	1:A:519:LEU:HB3	2.01	0.41
1:A:641:ARG:HE	1:A:670:GLU:CD	2.23	0.41
1:A:930:TYR:CD1	1:A:1012:ILE:HD13	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:624:VAL:HG11	1:A:659:TYR:CE1	2.55	0.41
1:A:943:GLY:CA	1:A:1050:TYR:CE2	3.03	0.41
1:A:737:GLN:O	1:A:741:MET:HG3	2.21	0.41
1:A:1009:PHE:HA	1:A:1012:ILE:HD12	2.02	0.41
1:A:1008:LYS:O	1:A:1012:ILE:HG13	2.21	0.41
1:A:172:GLU:HG3	1:A:471:HIS:CG	2.56	0.41
1:A:915:SER:HA	1:A:916:PRO:HD3	1.92	0.40
1:A:624:VAL:O	1:A:628:MET:HG2	2.20	0.40
1:A:1006:PHE:O	1:A:1009:PHE:HB3	2.22	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	769/960 (80%)	712 (93%)	49 (6%)	8 (1%)	19	54

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	549	ASN
1	A	164	ASP
1	A	579	ARG
1	A	411	ASN
1	A	1079	GLY
1	A	390	GLY
1	A	999	GLY
1	A	916	PRO

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	716/857 (84%)	669 (93%)	47 (7%)	21 51

All (47) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	165	VAL
1	A	178	ARG
1	A	240	THR
1	A	273	ARG
1	A	282	VAL
1	A	285	THR
1	A	287	ILE
1	A	359	ARG
1	A	372	VAL
1	A	381	VAL
1	A	395	CYS
1	A	398	ARG
1	A	409	LEU
1	A	430	ASN
1	A	464	VAL
1	A	479	GLU
1	A	512	ASN
1	A	520	LEU
1	A	591	LYS
1	A	600	GLN
1	A	613	ARG
1	A	614	ARG
1	A	624	VAL
1	A	626	LEU
1	A	632	ASP
1	A	729	LEU
1	A	734	GLN
1	A	758	ASP
1	A	784	ARG
1	A	845	LEU

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Mol	Chain	Res	Type
1	A	853	SER
1	A	862	LEU
1	A	865	LEU
1	A	870	ILE
1	A	886	THR
1	A	905	GLU
1	A	911	LEU
1	A	939	THR
1	A	959	ASN
1	A	964	ASP
1	A	983	VAL
1	A	1002	THR
1	A	1003	SER
1	A	1018	LEU
1	A	1039	MET
1	A	1059	LYS
1	A	1078	LYS

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	218	ASN
1	A	299	ASN
1	A	498	ASN
1	A	512	ASN
1	A	550	GLN
1	A	609	GLN
1	A	711	GLN
1	A	730	HIS
1	A	743	GLN
1	A	951	ASN
1	A	967	HIS
1	A	1085	ASN

5.3.3 RNA ⓘ

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](#)

2 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	QJZ	A	1	-	20,23,23	1.54	1 (5%)	21,32,32	3.06	8 (38%)
3	SO4	A	1103	-	4,4,4	0.21	0	6,6,6	0.17	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	QJZ	A	1	-	-	0/8/10/10	0/3/3/3
3	SO4	A	1103	-	-	0/0/0/0	0/0/0/0

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1	QJZ	C2-C10	-5.64	1.40	1.48

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1	QJZ	N17-C18-N19	-5.30	120.56	127.66

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1	QJZ	C15-C16-N17	-3.49	119.92	123.90
2	A	1	QJZ	C15-C10-N19	-2.51	118.55	122.01
2	A	1	QJZ	C4-C5-N9	2.39	113.70	108.16
2	A	1	QJZ	C2-C10-N19	2.49	119.44	116.13
2	A	1	QJZ	C16-N17-C18	3.81	121.01	114.93
2	A	1	QJZ	C18-N19-C10	4.91	121.22	116.22
2	A	1	QJZ	C21-S20-C18	9.48	108.34	102.26

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	1	QJZ	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	795/960 (82%)	0.23	37 (4%) 35 29	44, 77, 119, 134	0

All (37) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	217	ASN	5.6
1	A	216	ALA	4.9
1	A	1000	LYS	4.6
1	A	998	SER	4.3
1	A	232	THR	4.0
1	A	999	GLY	3.9
1	A	617	TRP	3.7
1	A	619	GLN	3.6
1	A	936	CYS	3.4
1	A	1089	HIS	2.9
1	A	919	GLU	2.8
1	A	981	GLU	2.8
1	A	989	PRO	2.7
1	A	753	SER	2.7
1	A	307	LEU	2.7
1	A	909	HIS	2.6
1	A	371	PRO	2.5
1	A	935	TYR	2.5
1	A	458	VAL	2.4
1	A	932	CYS	2.4
1	A	310	PRO	2.4
1	A	906	VAL	2.3
1	A	358	ASP	2.3
1	A	166	SER	2.3
1	A	231	GLN	2.2
1	A	1044	SER	2.2
1	A	213	LYS	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	911	LEU	2.2
1	A	314	ALA	2.2
1	A	937	VAL	2.1
1	A	245	LEU	2.1
1	A	1087	PHE	2.1
1	A	752	LEU	2.1
1	A	312	ASP	2.1
1	A	807	LYS	2.1
1	A	769	GLN	2.1
1	A	230	SER	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, 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	QJZ	A	1	21/21	0.93	0.22	0.59	73,75,76,77	0
3	SO4	A	1103	5/5	0.85	0.27	0.30	129,129,129,129	0

6.5 Other polymers [i](#)

There are no such residues in this entry.