



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 15, 2017 – 08:53 PM EST

PDB ID : 5I6U  
Title : The crystal structure of PI3Kdelta with compound 32  
Authors : Somoza, J.R.; Villasenor, A.G.  
Deposited on : 2016-02-16  
Resolution : 2.84 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 : 1.7.1 (RC1), CSD as537be (2016)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20028442  
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) : rb-20028442

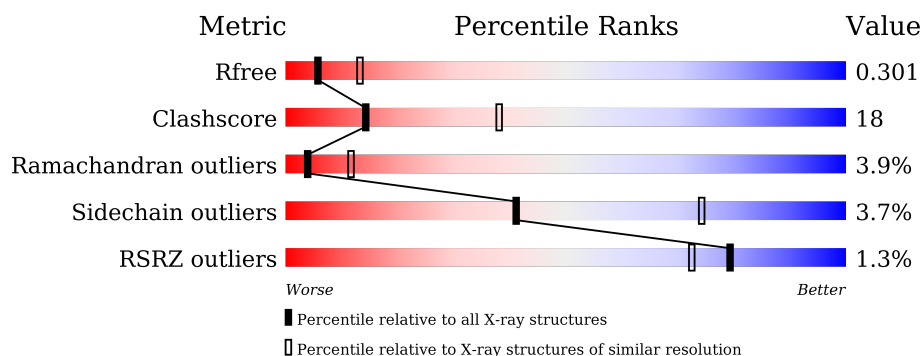
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.84 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	3170 (2.88-2.80)
Clashscore	102246	3658 (2.88-2.80)
Ramachandran outliers	100387	3591 (2.88-2.80)
Sidechain outliers	100360	3594 (2.88-2.80)
RSRZ outliers	91569	3184 (2.88-2.80)

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	939	<div> <div></div> <div>52%</div> <div>32%</div> <div>•</div> <div>13%</div> </div>

## 2 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 6622 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.

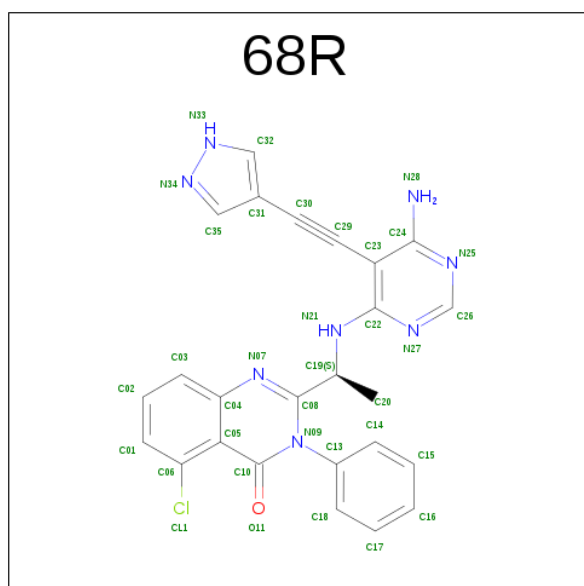
- Molecule 1 is a protein called Phosphatidylinositol 4,5-bisphosphate 3-kinase catalytic sub-unit delta isoform.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	816	6587	4224	1118	1191	54	0	0	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	508	GLN	-	insertion	UNP O35904

- Molecule 2 is 2-[(1S)-1-({6-amino-5-[(1H-pyrazol-4-yl)ethynyl]pyrimidin-4-yl}amino)ethyl]-5-chloro-3-phenylquinazolin-4(3H)-one (three-letter code: 68R) (formula: C<sub>25</sub>H<sub>19</sub>ClN<sub>8</sub>O).

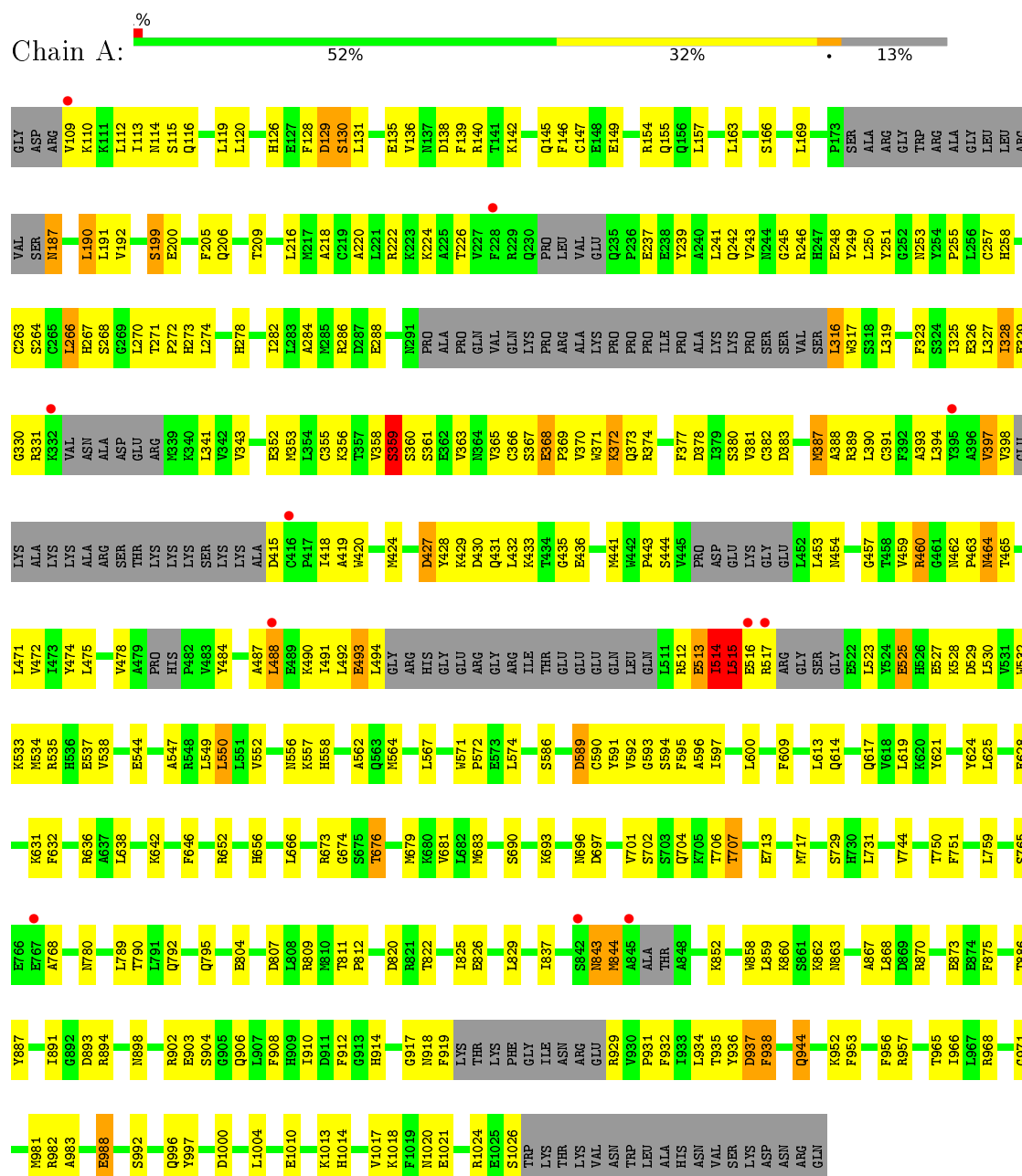


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	Cl	N	O		
2	A	1	35	25	1	8	1	0	0

### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

- Molecule 1: Phosphatidylinositol 4,5-bisphosphate 3-kinase catalytic subunit delta isoform



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	142.84Å 64.93Å 117.06Å 90.00° 103.17° 90.00°	Depositor
Resolution (Å)	29.66 – 2.84 29.66 – 2.84	Depositor EDS
% Data completeness (in resolution range)	96.9 (29.66-2.84) 90.0 (29.66-2.84)	Depositor EDS
$R_{merge}$	0.13	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.02 (at 2.85Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.9_1692)	Depositor
R, $R_{free}$	0.212 , 0.307 0.208 , 0.301	Depositor DCC
$R_{free}$ test set	1853 reflections (8.27%)	DCC
Wilson B-factor (Å <sup>2</sup> )	50.4	Xtriage
Anisotropy	0.172	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.27 , 37.5	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.46$ , $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.90	EDS
Total number of atoms	6622	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	53.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.00% 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: 68R

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	0.68	0/6727	0.83	6/9071 (0.1%)

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	8

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	488	LEU	CB-CG-CD2	-7.26	98.66	111.00
1	A	488	LEU	CA-CB-CG	-7.04	99.11	115.30
1	A	190	LEU	CA-CB-CG	6.95	131.27	115.30
1	A	515	LEU	CA-CB-CG	6.66	130.61	115.30
1	A	266	LEU	CA-CB-CG	6.18	129.52	115.30
1	A	488	LEU	N-CA-C	5.18	124.98	111.00

There are no chirality outliers.

All (8) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	444	SER	Peptide
1	A	490	LYS	Peptide
1	A	514	ILE	Peptide
1	A	516	GLU	Peptide

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Mol	Chain	Res	Type	Group
1	A	523	LEU	Peptide
1	A	525	GLU	Peptide
1	A	537	GLU	Peptide
1	A	914	HIS	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	6587	0	6572	239	0
2	A	35	0	0	4	0
All	All	6622	0	6572	239	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 18.

All (239) 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:525:GLU:HB3	1:A:528:LYS:HB3	1.38	1.02
1:A:355:CYS:SG	1:A:356:LYS:N	2.47	0.84
1:A:241:LEU:HD13	1:A:274:LEU:HD13	1.58	0.83
1:A:957:ARG:NH1	1:A:1020:ASN:OD1	2.12	0.83
1:A:383:ASP:OD1	1:A:558:HIS:ND1	2.13	0.82
1:A:982:ARG:HH12	1:A:992:SER:HA	1.47	0.80
1:A:547:ALA:HA	1:A:550:LEU:HD12	1.65	0.78
1:A:245:GLY:HA3	1:A:768:ALA:HB2	1.65	0.78
1:A:887:TYR:OH	1:A:929:ARG:NH2	2.17	0.77
1:A:216:LEU:HD22	1:A:241:LEU:HD11	1.67	0.77
1:A:267:HIS:CD2	1:A:809:ARG:NH1	2.53	0.76
1:A:192:VAL:HG12	1:A:272:PRO:HB2	1.68	0.76
1:A:517:ARG:NH1	1:A:544:GLU:OE1	2.18	0.74
1:A:904:SER:OG	1:A:906:GLN:HG3	1.86	0.74
1:A:533:LYS:HG2	1:A:534:MET:HG3	1.70	0.72
1:A:843:ASN:OD1	1:A:844:MET:N	2.23	0.71
1:A:359:SER:OG	1:A:360:SER:N	2.23	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:154:ARG:NH2	1:A:674:GLY:O	2.25	0.69
1:A:356:LYS:HD3	1:A:378:ASP:OD2	1.92	0.69
1:A:826:GLU:O	2:A:1101:68R:N28	2.26	0.68
1:A:891:ILE:HG22	1:A:894:ARG:HD3	1.74	0.68
1:A:371:TRP:O	1:A:373:GLN:N	2.27	0.67
1:A:380:SER:HB3	1:A:557:LYS:NZ	2.09	0.67
1:A:696:ASN:HB2	1:A:759:LEU:HD11	1.76	0.67
1:A:155:GLN:NE2	1:A:286:ARG:O	2.26	0.66
1:A:894:ARG:NH2	1:A:910:ILE:O	2.29	0.65
1:A:389:ARG:HB3	1:A:424:MET:HE1	1.77	0.65
1:A:512:ARG:HD3	1:A:530:LEU:HD22	1.78	0.64
1:A:116:GLN:HA	1:A:119:LEU:HD12	1.78	0.64
1:A:765:SER:HB3	1:A:768:ALA:HB3	1.78	0.64
1:A:513:GLU:HA	1:A:515:LEU:HB2	1.80	0.64
1:A:191:LEU:O	1:A:272:PRO:HD2	1.98	0.64
1:A:512:ARG:HH11	1:A:530:LEU:CD2	2.11	0.64
1:A:199:SER:OG	1:A:200:GLU:N	2.28	0.63
1:A:341:LEU:HB2	1:A:363:VAL:HG23	1.79	0.63
1:A:435:GLY:HA2	1:A:475:LEU:O	1.98	0.63
1:A:205:PHE:HZ	1:A:220:ALA:HA	1.62	0.63
1:A:701:VAL:O	1:A:704:GLN:HB2	1.98	0.63
1:A:453:LEU:HD12	1:A:454:ASN:H	1.63	0.63
1:A:614:GLN:HG3	1:A:981:MET:HG2	1.80	0.63
1:A:110:LYS:O	1:A:114:ASN:ND2	2.32	0.63
1:A:355:CYS:HG	1:A:356:LYS:H	1.46	0.62
1:A:917:GLY:HA3	1:A:997:TYR:CE1	2.34	0.62
1:A:982:ARG:NH1	1:A:992:SER:HA	2.14	0.62
1:A:488:LEU:HD21	1:A:491:ILE:HB	1.81	0.62
1:A:936:TYR:HD1	1:A:1026:SER:HB3	1.64	0.61
1:A:532:TRP:HZ3	1:A:564:MET:HE2	1.66	0.61
1:A:138:ASP:OD1	1:A:142:LYS:HB2	2.00	0.61
1:A:652:ARG:O	1:A:652:ARG:NH1	2.31	0.60
1:A:693:LYS:NZ	1:A:697:ASP:OD2	2.34	0.60
1:A:120:LEU:HD13	1:A:683:MET:HG3	1.82	0.60
1:A:435:GLY:O	1:A:475:LEU:N	2.34	0.60
1:A:706:THR:OG1	1:A:707:THR:N	2.34	0.60
1:A:263:CYS:HA	1:A:266:LEU:HG	1.83	0.60
1:A:656:HIS:CD2	1:A:820:ASP:OD2	2.55	0.60
1:A:329:GLU:HG3	1:A:370:VAL:HG22	1.84	0.59
1:A:136:VAL:HG13	1:A:666:LEU:HD11	1.84	0.59
1:A:982:ARG:HH12	1:A:992:SER:CA	2.14	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:192:VAL:HG23	1:A:205:PHE:HB2	1.84	0.59
1:A:488:LEU:CD2	1:A:491:ILE:HB	2.32	0.59
1:A:355:CYS:C	1:A:356:LYS:HD2	2.24	0.58
1:A:360:SER:OG	1:A:373:GLN:OE1	2.22	0.58
1:A:387:MET:HE3	1:A:590:CYS:H	1.66	0.58
1:A:971:GLY:HA3	1:A:1004:LEU:HD11	1.83	0.58
1:A:875:PHE:CE2	1:A:938:PHE:HB3	2.39	0.58
1:A:356:LYS:N	1:A:356:LYS:HD2	2.19	0.57
1:A:1010:GLU:O	1:A:1013:LYS:HB2	2.04	0.57
1:A:789:LEU:HD22	1:A:981:MET:HE3	1.86	0.57
1:A:154:ARG:HH11	1:A:154:ARG:HG2	1.70	0.57
1:A:267:HIS:CD2	1:A:809:ARG:HH11	2.23	0.57
1:A:157:LEU:O	1:A:286:ARG:NH1	2.38	0.56
1:A:389:ARG:HB3	1:A:424:MET:CE	2.35	0.56
1:A:621:TYR:CZ	1:A:983:ALA:HB2	2.40	0.56
1:A:255:PRO:HG2	1:A:258:HIS:HB2	1.87	0.56
1:A:187:ASN:OD1	1:A:187:ASN:N	2.39	0.55
1:A:431:GLN:HA	1:A:484:TYR:HA	1.88	0.55
1:A:128:PHE:HB2	1:A:140:ARG:NH1	2.22	0.55
1:A:549:LEU:HG	1:A:564:MET:HE3	1.89	0.54
1:A:352:GLU:OE2	1:A:353:MET:HG2	2.08	0.54
1:A:493:GLU:HG2	1:A:494:LEU:N	2.23	0.54
1:A:512:ARG:HG3	1:A:513:GLU:N	2.23	0.53
1:A:693:LYS:HE2	1:A:780:ASN:HD21	1.73	0.53
1:A:1017:VAL:O	1:A:1021:GLU:HG3	2.08	0.53
1:A:383:ASP:CG	1:A:557:LYS:HD2	2.28	0.53
1:A:380:SER:HB3	1:A:557:LYS:HZ2	1.73	0.53
1:A:693:LYS:NZ	1:A:697:ASP:OD1	2.41	0.53
1:A:935:THR:HG22	1:A:937:ASP:H	1.72	0.53
1:A:205:PHE:CZ	1:A:220:ALA:HA	2.43	0.53
1:A:326:GLU:HG2	1:A:327:LEU:O	2.09	0.53
1:A:390:LEU:HD12	1:A:391:CYS:H	1.74	0.53
1:A:525:GLU:C	1:A:527:GLU:H	2.13	0.52
1:A:571:TRP:CD1	1:A:572:PRO:HD2	2.44	0.52
1:A:424:MET:HG3	1:A:624:TYR:CE1	2.44	0.52
1:A:271:THR:O	1:A:273:HIS:ND1	2.42	0.52
1:A:367:SER:OG	1:A:369:PRO:HD3	2.09	0.52
1:A:397:VAL:HG22	1:A:398:VAL:H	1.73	0.52
1:A:693:LYS:NZ	1:A:697:ASP:CG	2.62	0.52
1:A:493:GLU:HG2	1:A:494:LEU:HB3	1.92	0.52
1:A:129:ASP:O	1:A:131:LEU:N	2.42	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:371:TRP:C	1:A:373:GLN:H	2.12	0.52
1:A:270:LEU:HD12	1:A:271:THR:H	1.75	0.52
1:A:209:THR:HB	1:A:257:CYS:HB3	1.93	0.51
1:A:529:ASP:OD1	1:A:530:LEU:N	2.40	0.51
1:A:609:PHE:CD2	1:A:642:LYS:NZ	2.76	0.51
1:A:387:MET:CE	1:A:590:CYS:H	2.23	0.51
1:A:129:ASP:C	1:A:131:LEU:H	2.14	0.51
1:A:329:GLU:HG2	1:A:330:GLY:H	1.74	0.51
1:A:343:VAL:HG22	1:A:394:LEU:HD12	1.93	0.51
1:A:996:GLN:O	1:A:996:GLN:HG3	2.09	0.51
1:A:242:GLN:HG2	1:A:243:VAL:O	2.11	0.51
1:A:147:CYS:SG	1:A:638:LEU:HD11	2.50	0.50
1:A:530:LEU:HA	1:A:533:LYS:HB3	1.94	0.50
1:A:592:VAL:O	1:A:596:ALA:N	2.36	0.50
1:A:593:GLY:O	1:A:597:ILE:HG12	2.11	0.50
1:A:367:SER:HA	1:A:368:GLU:C	2.32	0.49
1:A:424:MET:O	1:A:433:LYS:HD2	2.11	0.49
1:A:146:PHE:O	1:A:149:GLU:HG2	2.12	0.49
1:A:491:ILE:HG23	1:A:562:ALA:HB1	1.94	0.49
1:A:431:GLN:HG3	1:A:432:LEU:O	2.12	0.49
1:A:394:LEU:O	1:A:418:ILE:HB	2.13	0.49
1:A:239:TYR:CZ	1:A:278:HIS:HD2	2.31	0.49
1:A:114:ASN:HD22	1:A:114:ASN:H	1.60	0.49
1:A:825:ILE:HD13	2:A:1101:68R:C35	2.43	0.49
1:A:243:VAL:HA	1:A:274:LEU:HD23	1.95	0.49
1:A:549:LEU:O	1:A:552:VAL:HG22	2.13	0.49
1:A:944:GLN:OE1	1:A:952:LYS:HE3	2.13	0.49
1:A:420:TRP:CH2	1:A:457:GLY:HA3	2.48	0.48
1:A:453:LEU:HD12	1:A:454:ASN:N	2.29	0.48
1:A:617:GLN:HG3	1:A:981:MET:SD	2.53	0.48
1:A:218:ALA:O	1:A:222:ARG:HG3	2.14	0.48
1:A:936:TYR:CD1	1:A:1026:SER:HB3	2.48	0.48
1:A:617:GLN:HB3	1:A:983:ALA:HB3	1.96	0.48
1:A:419:ALA:HA	1:A:443:PRO:HA	1.95	0.48
1:A:862:LYS:HD2	1:A:903:GLU:HG2	1.95	0.48
1:A:397:VAL:HG22	1:A:398:VAL:HG13	1.95	0.48
1:A:358:VAL:HG21	1:A:377:PHE:CD1	2.49	0.48
1:A:380:SER:HB3	1:A:557:LYS:HZ1	1.78	0.48
1:A:246:ARG:NH1	1:A:248:GLU:OE1	2.47	0.47
1:A:512:ARG:O	1:A:514:ILE:N	2.47	0.47
1:A:804:GLU:OE2	1:A:966:ILE:HG23	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:968:ARG:HB3	1:A:968:ARG:NH1	2.30	0.47
1:A:427:ASP:HB3	1:A:431:GLN:O	2.14	0.47
1:A:690:SER:O	1:A:693:LYS:HB3	2.14	0.47
1:A:112:LEU:O	1:A:115:SER:HB3	2.14	0.47
1:A:114:ASN:CG	1:A:673:ARG:HH22	2.18	0.46
1:A:381:VAL:C	1:A:383:ASP:H	2.19	0.46
1:A:237:GLU:N	1:A:237:GLU:OE1	2.43	0.46
1:A:284:ALA:O	1:A:288:GLU:HG3	2.16	0.46
1:A:110:LYS:HE2	1:A:110:LYS:HB3	1.51	0.46
1:A:424:MET:CE	1:A:459:VAL:HG11	2.46	0.46
1:A:574:LEU:HG	1:A:595:PHE:HE2	1.81	0.46
1:A:807:ASP:OD1	1:A:807:ASP:C	2.54	0.46
1:A:875:PHE:CZ	1:A:938:PHE:HB3	2.51	0.46
1:A:852:LYS:HD2	1:A:936:TYR:CE2	2.51	0.45
1:A:953:PHE:O	1:A:956:PHE:HB3	2.16	0.45
1:A:613:LEU:O	1:A:617:GLN:HG2	2.16	0.45
1:A:860:LYS:HG2	1:A:868:LEU:HD22	1.97	0.45
1:A:628:GLU:OE2	1:A:631:LYS:HE3	2.16	0.45
1:A:652:ARG:HA	1:A:652:ARG:HD2	1.68	0.45
1:A:870:ARG:O	1:A:873:GLU:HB3	2.16	0.45
1:A:535:ARG:O	1:A:538:VAL:HB	2.16	0.45
1:A:163:LEU:HG	1:A:282:ILE:HG21	1.98	0.45
1:A:609:PHE:HE1	1:A:646:PHE:CD2	2.34	0.45
1:A:811:THR:HG23	1:A:902:ARG:HH21	1.81	0.45
1:A:415:ASP:OD1	1:A:415:ASP:N	2.48	0.45
1:A:428:TYR:CE2	1:A:429:LYS:HG2	2.51	0.45
1:A:135:GLU:HG3	1:A:428:TYR:CG	2.52	0.44
1:A:383:ASP:HB3	1:A:556:ASN:O	2.17	0.44
1:A:126:HIS:O	1:A:130:SER:HB3	2.18	0.44
1:A:136:VAL:O	1:A:140:ARG:HG3	2.17	0.44
1:A:330:GLY:C	1:A:331:ARG:HG3	2.38	0.44
1:A:327:LEU:O	1:A:372:LYS:HA	2.18	0.44
1:A:436:GLU:HB2	1:A:474:TYR:HD1	1.83	0.44
1:A:886:THR:HG22	1:A:891:ILE:HD12	1.99	0.44
1:A:932:PHE:HE2	1:A:934:LEU:CD2	2.31	0.44
1:A:729:SER:OG	1:A:744:VAL:HG12	2.18	0.44
1:A:325:ILE:HG22	1:A:475:LEU:HD23	2.00	0.44
1:A:731:LEU:H	1:A:731:LEU:HD23	1.82	0.43
1:A:491:ILE:HG23	1:A:562:ALA:CB	2.48	0.43
1:A:462:ASN:HA	1:A:463:PRO:HD2	1.81	0.43
1:A:514:ILE:HG13	1:A:514:ILE:O	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:589:ASP:OD2	1:A:591:TYR:HD2	2.01	0.43
1:A:371:TRP:C	1:A:373:GLN:N	2.71	0.43
1:A:525:GLU:C	1:A:527:GLU:N	2.72	0.43
1:A:239:TYR:CD1	1:A:278:HIS:HA	2.53	0.43
1:A:837:ILE:HD13	1:A:858:TRP:CD2	2.53	0.43
1:A:266:LEU:HD12	1:A:267:HIS:ND1	2.34	0.43
1:A:525:GLU:HA	1:A:528:LYS:H	1.83	0.43
1:A:825:ILE:HD13	2:A:1101:68R:C31	2.49	0.42
1:A:330:GLY:C	1:A:369:PRO:HD2	2.40	0.42
1:A:982:ARG:HH12	1:A:992:SER:N	2.17	0.42
1:A:393:ALA:HB2	1:A:453:LEU:HD13	2.01	0.42
1:A:656:HIS:CG	1:A:820:ASP:HB2	2.55	0.42
1:A:702:SER:O	1:A:706:THR:HG22	2.19	0.42
1:A:169:LEU:HD21	1:A:251:TYR:CE2	2.55	0.42
1:A:139:PHE:CD1	1:A:625:LEU:HB3	2.55	0.42
1:A:460:ARG:HE	1:A:460:ARG:HB2	1.50	0.42
1:A:326:GLU:O	1:A:474:TYR:N	2.47	0.42
1:A:383:ASP:OD1	1:A:558:HIS:N	2.29	0.42
1:A:829:LEU:HD23	1:A:829:LEU:HA	1.83	0.42
1:A:918:ASN:HB2	1:A:988:GLU:OE2	2.20	0.42
1:A:1024:ARG:HG3	1:A:1024:ARG:NH1	2.35	0.41
1:A:224:LYS:C	1:A:226:THR:H	2.23	0.41
1:A:441:MET:HE3	1:A:441:MET:HB2	1.86	0.41
1:A:513:GLU:HA	1:A:515:LEU:HD12	2.01	0.41
1:A:619:LEU:HD23	1:A:619:LEU:HA	1.84	0.41
1:A:902:ARG:HG3	1:A:908:PHE:HE2	1.85	0.41
1:A:328:ILE:HD12	1:A:472:VAL:HG12	2.02	0.41
1:A:241:LEU:O	1:A:250:LEU:N	2.48	0.41
1:A:323:PHE:O	1:A:377:PHE:HB2	2.20	0.41
1:A:858:TRP:CZ2	1:A:862:LYS:HE3	2.55	0.41
1:A:676:THR:O	1:A:679:MET:HB3	2.20	0.41
1:A:428:TYR:CZ	1:A:429:LYS:HG2	2.55	0.41
1:A:525:GLU:O	1:A:528:LYS:HG2	2.20	0.41
1:A:750:THR:HG23	2:A:1101:68R:CL1	2.57	0.41
1:A:492:LEU:O	1:A:492:LEU:HD23	2.20	0.41
1:A:600:LEU:HA	1:A:600:LEU:HD23	1.85	0.41
1:A:632:PHE:O	1:A:636:ARG:HG2	2.20	0.41
1:A:790:THR:HG21	1:A:912:PHE:CD1	2.56	0.41
1:A:898:ASN:HB2	1:A:910:ILE:O	2.20	0.41
1:A:316:LEU:HB2	1:A:317:TRP:H	1.69	0.41
1:A:360:SER:OG	1:A:361:SER:N	2.54	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:319:LEU:O	1:A:382:CYS:HB3	2.21	0.41
1:A:713:GLU:O	1:A:717:MET:HG3	2.21	0.41
1:A:557:LYS:HD2	1:A:557:LYS:HA	1.69	0.41
1:A:491:ILE:O	1:A:491:ILE:CG2	2.69	0.41
1:A:191:LEU:HD12	1:A:206:GLN:HG2	2.03	0.40
1:A:935:THR:HB	1:A:938:PHE:CE2	2.57	0.40
1:A:365:VAL:O	1:A:367:SER:N	2.54	0.40
1:A:567:LEU:O	1:A:571:TRP:HE3	2.05	0.40
1:A:129:ASP:C	1:A:131:LEU:N	2.74	0.40
1:A:145:GLN:O	1:A:149:GLU:OE2	2.39	0.40
1:A:750:THR:OG1	1:A:751:PHE:N	2.53	0.40
1:A:795:GLN:HE21	1:A:812:PRO:HB2	1.87	0.40
1:A:1014:HIS:NE2	1:A:1018:LYS:HE2	2.37	0.40
1:A:525:GLU:HA	1:A:527:GLU:H	1.87	0.40
1:A:242:GLN:HB2	1:A:249:TYR:CZ	2.57	0.40
1:A:271:THR:O	1:A:273:HIS:CE1	2.75	0.40
1:A:424:MET:HE2	1:A:459:VAL:HG11	2.03	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	792/939 (84%)	682 (86%)	79 (10%)	31 (4%)	<b>4</b> <b>12</b>

All (31) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	130	SER
1	A	253	ASN
1	A	264	SER
1	A	366	CYS

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Mol	Chain	Res	Type
1	A	372	LYS
1	A	465	THR
1	A	843	ASN
1	A	844	MET
1	A	944	GLN
1	A	359	SER
1	A	388	ALA
1	A	397	VAL
1	A	478	VAL
1	A	487	ALA
1	A	893	ASP
1	A	387	MET
1	A	427	ASP
1	A	430	ASP
1	A	493	GLU
1	A	268	SER
1	A	513	GLU
1	A	589	ASP
1	A	129	ASP
1	A	368	GLU
1	A	464	ASN
1	A	586	SER
1	A	867	ALA
1	A	514	ILE
1	A	328	ILE
1	A	113	ILE
1	A	931	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	724/827 (88%)	697 (96%)	27 (4%)	41 74

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	109	VAL
1	A	166	SER
1	A	187	ASN
1	A	190	LEU
1	A	199	SER
1	A	316	LEU
1	A	359	SER
1	A	374	ARG
1	A	460	ARG
1	A	464	ASN
1	A	471	LEU
1	A	515	LEU
1	A	550	LEU
1	A	594	SER
1	A	676	THR
1	A	681	VAL
1	A	707	THR
1	A	792	GLN
1	A	822	THR
1	A	859	LEU
1	A	863	ASN
1	A	919	PHE
1	A	937	ASP
1	A	938	PHE
1	A	965	THR
1	A	988	GLU
1	A	1000	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	114	ASN
1	A	278	HIS
1	A	656	HIS
1	A	795	GLN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

## 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

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	68R	A	1101	-	37,39,39	1.55	5 (13%)	31,55,55	1.67	7 (22%)

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	68R	A	1101	-	-	0/11/17/17	0/5/5/5

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1101	68R	C23-C22	-3.60	1.37	1.42
2	A	1101	68R	C14-C13	2.06	1.42	1.38
2	A	1101	68R	C10-N09	2.66	1.42	1.37
2	A	1101	68R	C22-N21	3.04	1.40	1.35
2	A	1101	68R	C10-C05	5.61	1.50	1.41

All (7) bond angle outliers are listed below:



Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1101	68R	C26-N25-C24	-2.97	113.44	118.77
2	A	1101	68R	N28-C24-N25	-2.24	113.20	118.28
2	A	1101	68R	C14-C13-C18	-2.01	117.99	121.20
2	A	1101	68R	C15-C14-C13	2.85	122.52	118.71
2	A	1101	68R	C24-C23-C29	2.87	122.86	118.09
2	A	1101	68R	C20-C19-N21	3.38	114.24	109.08
2	A	1101	68R	C18-C13-N09	4.10	124.36	119.25

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1101	68R	4	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 [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	816/939 (86%)	-0.26	11 (1%) 79 72	23, 52, 88, 141	0

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	488	LEU	3.9
1	A	842	SER	3.4
1	A	109	VAL	3.2
1	A	416	CYS	3.0
1	A	845	ALA	2.7
1	A	332	LYS	2.7
1	A	516	GLU	2.4
1	A	517	ARG	2.4
1	A	395	TYR	2.4
1	A	767	GLU	2.3
1	A	228	PHE	2.2

### 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
2	68R	A	1101	35/35	0.96	0.14	0.09	18,31,37,73	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.