



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 11:04 AM GMT

PDB ID : 3NZU
Title : Structure-based Optimization of Pyrazolo -Pyrimidine and -Pyridine Inhibitors of PI3-Kinase
Authors : Murray, J.M.; Wiesmann, C.
Deposited on : 2010-07-16
Resolution : 2.60 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

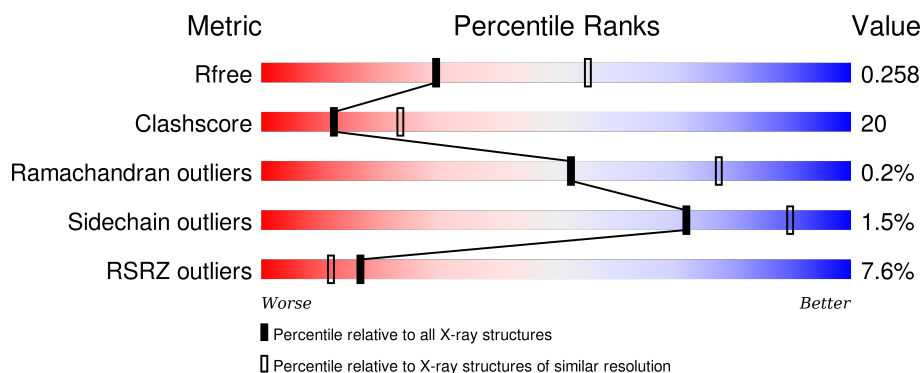
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	2328 (2.60-2.60)
Clashscore	102246	2679 (2.60-2.60)
Ramachandran outliers	100387	2635 (2.60-2.60)
Sidechain outliers	100360	2635 (2.60-2.60)
RSRZ outliers	91569	2334 (2.60-2.60)

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	954	<div> <div>6%</div> <div>55%</div> <div>30%</div> <div>•</div> <div>14%</div> </div>

2 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 6697 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.

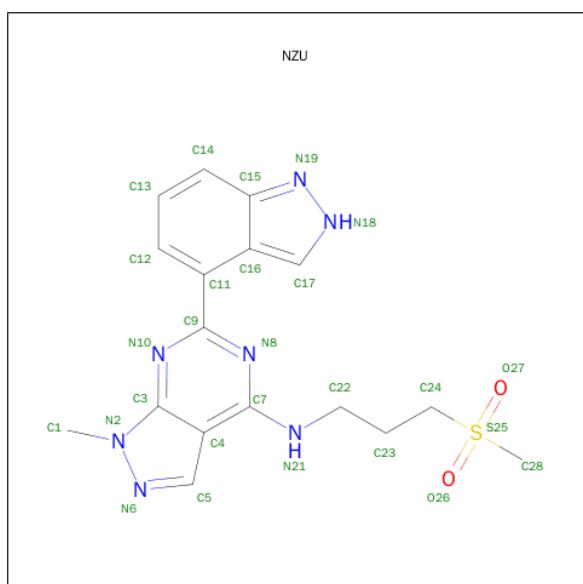
- Molecule 1 is a protein called Phosphatidylinositol-4,5-bisphosphate 3-kinase catalytic subunit gamma isoform.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	822	Total	C	N	O	S	37	0	0
			6670	4292	1137	1207	34			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1095	HIS	-	EXPRESSION TAG	UNP P48736
A	1096	HIS	-	EXPRESSION TAG	UNP P48736
A	1097	HIS	-	EXPRESSION TAG	UNP P48736
A	1098	HIS	-	EXPRESSION TAG	UNP P48736
A	1099	HIS	-	EXPRESSION TAG	UNP P48736
A	1100	HIS	-	EXPRESSION TAG	UNP P48736

- Molecule 2 is 6-(2H-INDAZOL-4-YL)-1-METHYL-N-[3-(METHYLSULFONYL)PROPYL]-1H-PYRAZOLO[3,4-D]PYRIMIDIN-4-AMINE (three-letter code: NZU) (formula: C₁₇H₁₉N₇O₂S).

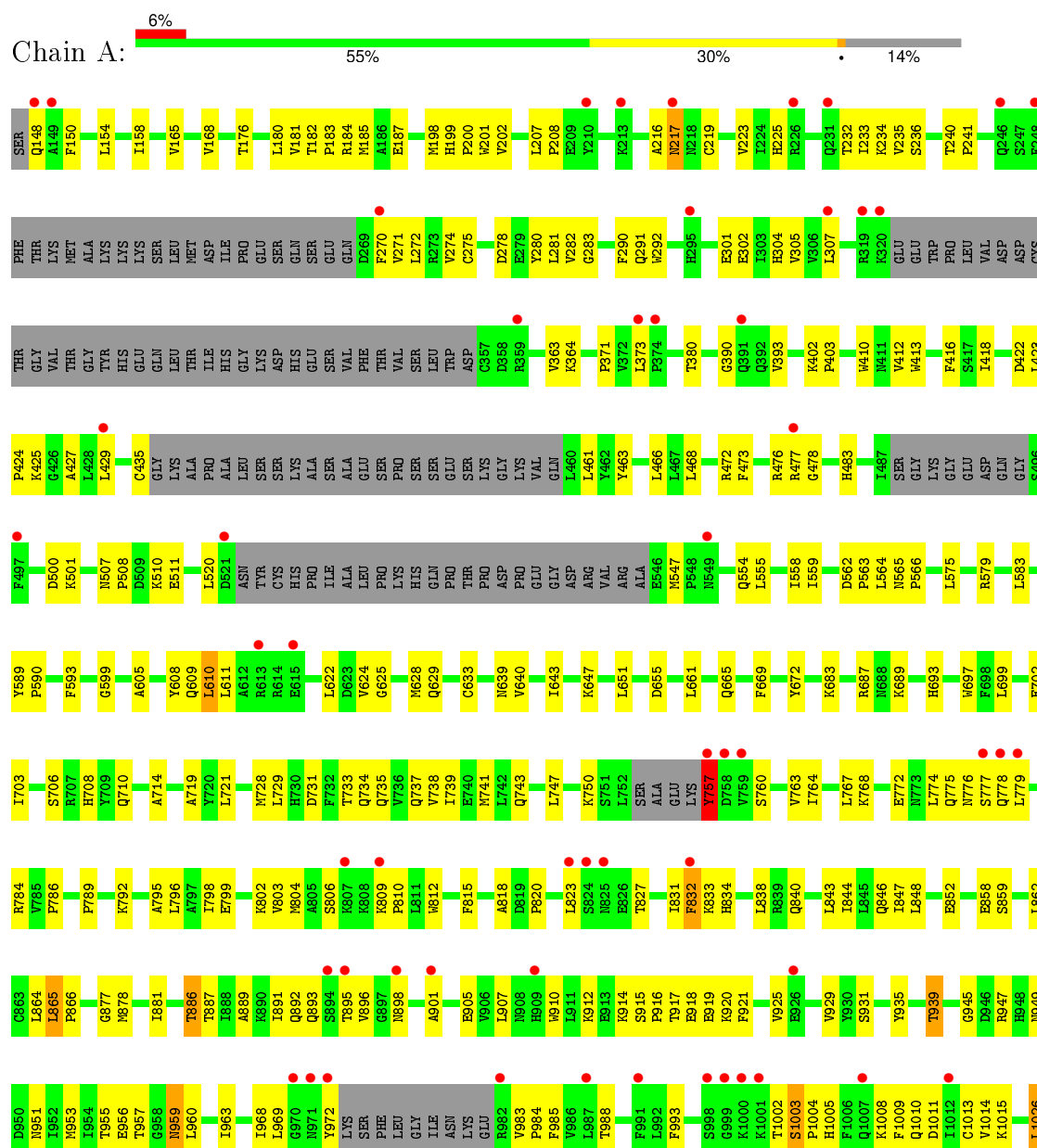


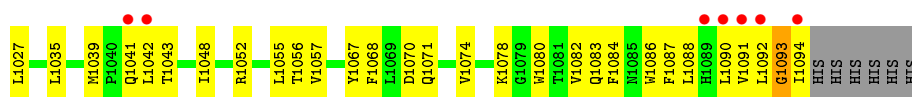
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	S	0	0
			27	17	7	2	1		

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 gamma isoform





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	144.00Å 67.68Å 106.36Å 90.00° 95.52° 90.00°	Depositor
Resolution (Å)	19.88 – 2.60 19.88 – 2.60	Depositor EDS
% Data completeness (in resolution range)	99.9 (19.88-2.60) 99.9 (19.88-2.60)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.78 (at 2.59Å)	Xtriage
Refinement program	PHENIX (phenix.refine: dev_392)	Depositor
R, R_{free}	0.213 , 0.240 0.239 , 0.258	Depositor DCC
R_{free} test set	1588 reflections (5.31%)	DCC
Wilson B-factor (Å ²)	70.7	Xtriage
Anisotropy	0.053	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 69.0	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 31497 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6697	wwPDB-VP
Average B, all atoms (Å ²)	97.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.65% 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 [i](#)

5.1 Standard geometry [i](#)

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

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	0.22	0/6814	0.41	3/9221 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	390	GLY	N-CA-C	-11.78	83.66	113.10
1	A	1003	SER	C-N-CD	-6.08	107.22	120.60
1	A	757	TYR	N-CA-C	5.76	126.56	111.00

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	6670	0	6728	262	0
2	A	27	0	19	3	0
All	All	6697	0	6747	264	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 20.

All (264) 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:798:ILE:HD12	1:A:798:ILE:H	1.26	1.00
1:A:1092:LEU:HD11	1:A:1094:ILE:HG13	1.48	0.95
1:A:1041:GLN:HG2	1:A:1042:LEU:CD1	2.01	0.90
1:A:1041:GLN:HG2	1:A:1042:LEU:HD13	1.55	0.87
1:A:216:ALA:O	1:A:217:ASN:HB2	1.72	0.87
1:A:1092:LEU:CD1	1:A:1094:ILE:HG13	2.05	0.87
1:A:887:THR:HG22	1:A:889:ALA:H	1.40	0.86
1:A:955:THR:HG22	1:A:957:THR:H	1.40	0.84
1:A:905:GLU:HG2	1:A:993:PHE:CE1	2.15	0.81
1:A:182:THR:HB	1:A:183:PRO:HD3	1.63	0.79
1:A:776:ASN:O	1:A:777:SER:OG	2.01	0.79
1:A:429:LEU:HB2	1:A:468:LEU:HD21	1.65	0.79
1:A:278:ASP:HB2	1:A:792:LYS:HE3	1.64	0.78
1:A:969:LEU:HD13	1:A:1039:MET:HE1	1.67	0.76
1:A:774:LEU:O	1:A:778:GLN:HB2	1.88	0.73
1:A:809:LYS:N	1:A:810:PRO:HD3	2.04	0.73
1:A:905:GLU:HG2	1:A:993:PHE:CZ	2.24	0.72
1:A:579:ARG:HB2	1:A:610:LEU:HD11	1.72	0.72
1:A:798:ILE:CD1	1:A:798:ILE:H	2.02	0.72
1:A:969:LEU:HD13	1:A:1039:MET:CE	2.20	0.71
1:A:887:THR:HG22	1:A:889:ALA:N	2.06	0.71
1:A:697:TRP:CH2	1:A:739:ILE:HD13	2.26	0.71
1:A:947:ARG:HD2	1:A:968:ILE:HD12	1.71	0.70
1:A:625:GLY:O	1:A:629:GLN:HG3	1.92	0.70
1:A:983:VAL:HG23	1:A:984:PRO:HD2	1.74	0.70
1:A:775:GLN:HA	1:A:779:LEU:CD2	2.22	0.69
1:A:919:GLU:HG3	1:A:920:LYS:N	2.06	0.69
1:A:935:TYR:O	1:A:939:THR:HB	1.93	0.69
1:A:896:VAL:HG11	1:A:901:ALA:HB3	1.75	0.68
1:A:768:LYS:O	1:A:772:GLU:HG3	1.94	0.67
1:A:735:GLN:O	1:A:739:ILE:HD12	1.95	0.67
1:A:693:HIS:CD2	1:A:789:PRO:HG3	2.30	0.67
1:A:271:VAL:CG2	1:A:282:VAL:HG22	2.25	0.67
1:A:969:LEU:HB3	1:A:1039:MET:HE3	1.77	0.66
1:A:1092:LEU:HD11	1:A:1094:ILE:CG1	2.25	0.66
1:A:1041:GLN:HE21	1:A:1042:LEU:HD13	1.62	0.65
1:A:219:CYS:SG	1:A:234:LYS:HG3	2.37	0.65
1:A:1092:LEU:HD12	1:A:1093:GLY:N	2.11	0.64
1:A:969:LEU:HD22	1:A:1039:MET:CE	2.27	0.64
1:A:843:LEU:O	1:A:847:ILE:HG12	1.97	0.64
1:A:775:GLN:HA	1:A:779:LEU:HD21	1.80	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:738:VAL:HA	1:A:741:MET:HE2	1.82	0.62
1:A:232:THR:HG22	1:A:233:ILE:N	2.14	0.62
1:A:947:ARG:CD	1:A:968:ILE:HD12	2.29	0.62
1:A:425:LYS:HD2	1:A:672:TYR:OH	1.99	0.61
1:A:889:ALA:O	1:A:893:GLN:HG3	2.00	0.61
1:A:951:ASN:HA	1:A:963:ILE:HD11	1.83	0.61
1:A:1087:PHE:CD2	1:A:1091:VAL:HG21	2.36	0.61
1:A:181:VAL:O	1:A:185:MET:HG3	2.01	0.61
1:A:919:GLU:HG3	1:A:920:LYS:H	1.66	0.60
1:A:583:LEU:HD22	1:A:610:LEU:CD2	2.31	0.60
1:A:1010:GLN:O	1:A:1014:VAL:HG23	2.02	0.60
1:A:583:LEU:HD22	1:A:610:LEU:HD22	1.84	0.60
1:A:862:LEU:HD22	1:A:862:LEU:N	2.17	0.59
1:A:410:TRP:HB3	1:A:412:VAL:HG23	1.84	0.59
1:A:1087:PHE:O	1:A:1091:VAL:HB	2.01	0.59
1:A:799:GLU:CD	1:A:799:GLU:H	2.05	0.59
1:A:925:VAL:O	1:A:929:VAL:HG23	2.03	0.59
1:A:905:GLU:HA	1:A:993:PHE:CD2	2.38	0.59
1:A:838:LEU:HD12	1:A:877:GLY:HA3	1.85	0.59
1:A:640:VAL:O	1:A:643:ILE:HG12	2.02	0.58
1:A:784:ARG:HH12	1:A:792:LYS:HE3	1.68	0.58
1:A:953:MET:O	1:A:960:LEU:HD12	2.04	0.58
1:A:806:SER:HB3	1:A:810:PRO:CD	2.33	0.58
1:A:466:LEU:HD11	1:A:476:ARG:HD3	1.85	0.58
1:A:733:THR:O	1:A:737:GLN:HG3	2.02	0.58
1:A:1041:GLN:HG2	1:A:1042:LEU:HD12	1.84	0.58
1:A:622:LEU:HD13	1:A:647:LYS:HB3	1.85	0.58
1:A:154:LEU:O	1:A:158:ILE:HG13	2.04	0.57
1:A:564:LEU:HD11	1:A:1048:ILE:HG22	1.85	0.57
1:A:969:LEU:HD22	1:A:1039:MET:HE1	1.85	0.57
1:A:939:THR:HG23	1:A:945:GLY:HA3	1.87	0.57
1:A:838:LEU:HD12	1:A:877:GLY:CA	2.34	0.57
1:A:775:GLN:NE2	1:A:795:ALA:HB1	2.19	0.57
1:A:207:LEU:HD12	1:A:208:PRO:HD2	1.87	0.57
1:A:798:ILE:HD12	1:A:798:ILE:N	2.07	0.57
1:A:232:THR:HG22	1:A:233:ILE:H	1.69	0.56
1:A:1041:GLN:CG	1:A:1042:LEU:HD13	2.33	0.56
1:A:697:TRP:HH2	1:A:739:ILE:HD13	1.67	0.56
1:A:555:LEU:O	1:A:559:ILE:HG13	2.05	0.56
1:A:625:GLY:HA2	1:A:1026:LEU:HD23	1.86	0.56
1:A:1092:LEU:HD12	1:A:1093:GLY:H	1.70	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1088:LEU:HB3	1:A:1094:ILE:HB	1.87	0.56
1:A:402:LYS:HG3	1:A:403:PRO:HD2	1.87	0.56
1:A:988:THR:HG21	1:A:1083:GLN:HG3	1.88	0.56
1:A:914:LYS:O	1:A:916:PRO:HD3	2.06	0.56
1:A:274:VAL:HG21	1:A:292:TRP:CD1	2.41	0.56
1:A:483:HIS:CD2	1:A:510:LYS:HG2	2.41	0.56
1:A:271:VAL:HG21	1:A:282:VAL:CG2	2.36	0.55
1:A:760:SER:OG	1:A:763:VAL:HG12	2.06	0.55
1:A:796:LEU:HD21	1:A:815:PHE:CZ	2.41	0.55
1:A:767:LEU:HD22	1:A:803:VAL:HG23	1.88	0.55
1:A:461:LEU:HD22	1:A:461:LEU:N	2.22	0.55
1:A:779:LEU:O	1:A:779:LEU:HD12	2.07	0.55
1:A:757:TYR:CD2	1:A:757:TYR:N	2.72	0.55
1:A:1074:VAL:O	1:A:1078:LYS:HG2	2.06	0.55
1:A:743:GLN:O	1:A:747:LEU:HG	2.06	0.54
1:A:702:GLU:O	1:A:706:SER:HB3	2.07	0.54
1:A:864:LEU:C	1:A:866:PRO:HD3	2.28	0.54
1:A:731:ASP:O	1:A:735:GLN:HG3	2.08	0.54
1:A:477:ARG:HG2	1:A:478:GLY:N	2.22	0.54
1:A:983:VAL:CG2	1:A:984:PRO:HD2	2.38	0.54
1:A:955:THR:HG22	1:A:956:GLU:N	2.22	0.54
1:A:775:GLN:HA	1:A:779:LEU:HG	1.89	0.54
1:A:1082:VAL:HG12	1:A:1086:TRP:CD1	2.43	0.54
1:A:270:PHE:CG	1:A:307:LEU:HD21	2.44	0.53
1:A:1088:LEU:O	1:A:1092:LEU:HG	2.08	0.53
1:A:187:GLU:OE2	1:A:683:LYS:HG3	2.08	0.53
1:A:463:TYR:CE1	1:A:501:LYS:HA	2.44	0.53
1:A:939:THR:HG23	1:A:945:GLY:CA	2.39	0.53
1:A:364:LYS:HE3	1:A:413:TRP:CE2	2.44	0.53
1:A:202:VAL:HG12	1:A:290:PHE:CD2	2.44	0.52
1:A:918:GLU:O	1:A:921:PHE:HB3	2.09	0.52
1:A:775:GLN:OE1	1:A:798:ILE:HD11	2.09	0.52
1:A:806:SER:HB3	1:A:810:PRO:HD2	1.91	0.52
1:A:957:THR:HG22	1:A:957:THR:O	2.10	0.52
1:A:775:GLN:HA	1:A:779:LEU:CG	2.40	0.52
1:A:466:LEU:HD11	1:A:476:ARG:CD	2.40	0.52
1:A:216:ALA:O	1:A:217:ASN:CB	2.48	0.51
1:A:271:VAL:HG22	1:A:272:LEU:N	2.25	0.51
1:A:917:THR:OG1	1:A:920:LYS:HG3	2.10	0.51
1:A:429:LEU:HB2	1:A:468:LEU:CD2	2.40	0.50
1:A:422:ASP:HB3	1:A:599:GLY:O	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:182:THR:CB	1:A:183:PRO:HD3	2.39	0.50
1:A:1041:GLN:NE2	1:A:1042:LEU:HD13	2.26	0.50
1:A:500:ASP:HB3	1:A:708:HIS:CD2	2.47	0.50
1:A:895:THR:HG22	1:A:896:VAL:HG23	1.94	0.50
1:A:852:GLU:OE2	1:A:864:LEU:HB2	2.11	0.50
1:A:472:ARG:O	1:A:473:PHE:HB2	2.12	0.49
1:A:831:ILE:HG13	1:A:881:ILE:CG1	2.42	0.49
1:A:852:GLU:HG3	1:A:864:LEU:HD12	1.94	0.49
1:A:844:ILE:O	1:A:848:LEU:HG	2.12	0.49
2:A:1:NZU:H17	2:A:1:NZU:N10	2.28	0.49
2:A:1:NZU:C17	2:A:1:NZU:N10	2.76	0.49
1:A:301:GLU:HG2	1:A:302:GLU:N	2.28	0.49
1:A:271:VAL:HG21	1:A:282:VAL:HG21	1.95	0.48
1:A:786:PRO:HG2	1:A:878:MET:SD	2.53	0.48
1:A:554:GLN:O	1:A:558:ILE:HG13	2.12	0.48
1:A:608:TYR:CZ	1:A:639:ASN:ND2	2.80	0.48
1:A:898:ASN:O	1:A:1087:PHE:HZ	1.97	0.48
1:A:886:THR:CG2	1:A:910:TRP:CZ2	2.96	0.48
1:A:886:THR:HG21	1:A:910:TRP:CZ2	2.48	0.48
1:A:779:LEU:C	1:A:779:LEU:HD12	2.34	0.48
1:A:234:LYS:NZ	1:A:234:LYS:HB2	2.29	0.47
1:A:734:GLN:O	1:A:738:VAL:HG23	2.13	0.47
1:A:887:THR:O	1:A:891:ILE:HG13	2.14	0.47
1:A:589:TYR:HB2	1:A:590:PRO:HD3	1.95	0.47
1:A:605:ALA:O	1:A:609:GLN:HG3	2.14	0.47
1:A:1005:HIS:O	1:A:1008:LYS:HB3	2.15	0.47
1:A:750:LYS:NZ	1:A:834:HIS:HD2	2.12	0.47
1:A:840:GLN:O	1:A:844:ILE:HG12	2.14	0.47
1:A:373:LEU:O	1:A:373:LEU:HD12	2.15	0.47
1:A:760:SER:O	1:A:763:VAL:HG13	2.13	0.47
1:A:963:ILE:C	1:A:963:ILE:HD12	2.35	0.47
1:A:706:SER:O	1:A:710:GLN:HB3	2.15	0.47
1:A:1086:TRP:HE3	1:A:1090:LEU:HD13	1.80	0.46
1:A:181:VAL:HG22	1:A:184:ARG:HH22	1.81	0.46
1:A:969:LEU:CD1	1:A:1039:MET:HE1	2.41	0.46
1:A:702:GLU:HA	1:A:702:GLU:OE1	2.15	0.46
1:A:815:PHE:O	1:A:827:THR:HB	2.16	0.46
1:A:917:THR:HB	1:A:919:GLU:HG2	1.97	0.46
1:A:280:TYR:HB3	1:A:282:VAL:HG23	1.97	0.46
1:A:1070:ASP:O	1:A:1074:VAL:HG23	2.16	0.46
1:A:661:LEU:O	1:A:665:GLN:HG2	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:912:LYS:HG3	1:A:921:PHE:CE1	2.51	0.46
1:A:235:VAL:HG12	1:A:236:SER:N	2.31	0.46
1:A:1052:ARG:O	1:A:1057:VAL:HG23	2.17	0.46
1:A:955:THR:CG2	1:A:956:GLU:N	2.78	0.45
1:A:373:LEU:HD12	1:A:373:LEU:C	2.37	0.45
1:A:651:LEU:HD22	1:A:655:ASP:HB3	1.97	0.45
1:A:589:TYR:CD2	1:A:593:PHE:HE1	2.34	0.45
1:A:363:VAL:HG12	1:A:416:PHE:HE1	1.81	0.45
1:A:274:VAL:HG11	1:A:292:TRP:CE2	2.52	0.45
1:A:760:SER:O	1:A:763:VAL:CG1	2.64	0.45
1:A:959:ASN:C	1:A:959:ASN:HD22	2.20	0.45
1:A:689:LYS:HG2	1:A:728:MET:SD	2.56	0.45
1:A:818:ALA:O	1:A:820:PRO:HD3	2.16	0.45
1:A:555:LEU:HD11	1:A:575:LEU:HD23	1.98	0.45
1:A:1035:LEU:HD12	1:A:1039:MET:HG3	1.99	0.44
1:A:302:GLU:HG3	1:A:304:HIS:NE2	2.32	0.44
1:A:858:GLU:O	1:A:859:SER:HB2	2.17	0.44
1:A:721:LEU:O	1:A:729:LEU:HD21	2.17	0.44
1:A:187:GLU:OE1	1:A:687:ARG:HG3	2.17	0.44
1:A:201:TRP:CE2	1:A:291:GLN:HG3	2.53	0.44
1:A:1084:PHE:O	1:A:1087:PHE:HB3	2.18	0.44
1:A:832:PHE:CE2	1:A:878:MET:HE2	2.53	0.44
1:A:687:ARG:HH11	1:A:687:ARG:HG3	1.82	0.44
1:A:275:CYS:HB3	1:A:823:LEU:HD13	1.99	0.44
1:A:775:GLN:NE2	1:A:796:LEU:H	2.16	0.43
1:A:796:LEU:CD2	1:A:815:PHE:CE2	3.01	0.43
1:A:184:ARG:HD3	1:A:719:ALA:O	2.17	0.43
1:A:562:ASP:HB2	1:A:563:PRO:HD2	1.99	0.43
1:A:889:ALA:HB2	1:A:949:ASN:OD1	2.18	0.43
1:A:804:MET:CE	2:A:1:NZU:H22	2.47	0.43
1:A:423:LEU:HA	1:A:424:PRO:HD3	1.88	0.43
1:A:802:LYS:HG3	1:A:812:TRP:HB3	1.99	0.43
1:A:565:ASN:HA	1:A:566:PRO:HD3	1.82	0.43
1:A:809:LYS:N	1:A:810:PRO:CD	2.76	0.43
1:A:862:LEU:HD22	1:A:862:LEU:H	1.83	0.43
1:A:862:LEU:CD2	1:A:862:LEU:N	2.81	0.43
1:A:435:CYS:SG	1:A:461:LEU:HD21	2.58	0.43
1:A:418:ILE:HD13	1:A:423:LEU:HD23	1.99	0.43
1:A:199:HIS:N	1:A:200:PRO:HD3	2.34	0.43
1:A:307:LEU:C	1:A:307:LEU:HD23	2.39	0.43
1:A:271:VAL:HG22	1:A:272:LEU:H	1.82	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:234:LYS:CB	1:A:234:LYS:NZ	2.82	0.43
1:A:232:THR:CG2	1:A:233:ILE:N	2.82	0.43
1:A:393:VAL:O	1:A:393:VAL:HG23	2.18	0.43
1:A:775:GLN:NE2	1:A:796:LEU:N	2.67	0.43
1:A:1041:GLN:CD	1:A:1041:GLN:H	2.22	0.43
1:A:282:VAL:HG12	1:A:283:GLY:N	2.34	0.42
1:A:235:VAL:CG1	1:A:236:SER:N	2.82	0.42
1:A:931:SER:O	1:A:935:TYR:HD1	2.02	0.42
1:A:182:THR:HB	1:A:183:PRO:CD	2.43	0.42
1:A:1039:MET:O	1:A:1043:THR:HG23	2.19	0.42
1:A:1011:ASP:O	1:A:1015:LYS:HG3	2.20	0.42
1:A:891:ILE:HG22	1:A:891:ILE:O	2.19	0.42
1:A:148:GLN:C	1:A:150:PHE:H	2.23	0.42
1:A:624:VAL:O	1:A:628:MET:HG2	2.19	0.42
1:A:939:THR:CG2	1:A:945:GLY:CA	2.98	0.42
1:A:234:LYS:HZ2	1:A:234:LYS:HB2	1.84	0.42
1:A:1013:CYS:HB3	1:A:1068:PHE:CE2	2.54	0.42
1:A:972:TYR:HB3	1:A:1042:LEU:HD11	2.02	0.42
1:A:968:ILE:O	1:A:968:ILE:HG12	2.19	0.42
1:A:564:LEU:CD1	1:A:1048:ILE:HG22	2.47	0.42
1:A:281:LEU:HD12	1:A:305:VAL:HG11	2.02	0.41
1:A:507:ASN:HA	1:A:508:PRO:HD3	1.85	0.41
1:A:165:VAL:O	1:A:168:VAL:HG22	2.20	0.41
1:A:1067:TYR:O	1:A:1071:GLN:HG2	2.19	0.41
1:A:1092:LEU:CD1	1:A:1093:GLY:H	2.33	0.41
1:A:476:ARG:HG2	1:A:520:LEU:HD22	2.02	0.41
1:A:180:LEU:C	1:A:183:PRO:HD2	2.41	0.41
1:A:301:GLU:HG2	1:A:302:GLU:H	1.85	0.41
1:A:865:LEU:HD12	1:A:865:LEU:HA	1.80	0.41
1:A:547:MET:HB3	1:A:547:MET:HE2	1.99	0.41
1:A:1002:THR:O	1:A:1003:SER:HB3	2.21	0.41
1:A:240:THR:HB	1:A:241:PRO:HD2	2.02	0.41
1:A:892:GLN:HE22	1:A:1080:TRP:HH2	1.68	0.41
1:A:371:PRO:HG2	1:A:511:GLU:O	2.21	0.41
1:A:810:PRO:HB3	1:A:833:LYS:HB2	2.02	0.41
1:A:945:GLY:O	1:A:985:PHE:HA	2.20	0.41
1:A:201:TRP:NE1	1:A:291:GLN:HG3	2.36	0.41
1:A:223:VAL:HG12	1:A:225:HIS:NE2	2.36	0.41
1:A:775:GLN:CA	1:A:779:LEU:HD21	2.49	0.41
1:A:919:GLU:CG	1:A:920:LYS:N	2.81	0.41
1:A:915:SER:OG	1:A:921:PHE:HB2	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:831:ILE:HG13	1:A:881:ILE:HG12	2.02	0.41
1:A:424:PRO:HD2	1:A:427:ALA:HB2	2.02	0.41
1:A:699:LEU:O	1:A:703:ILE:HG13	2.21	0.41
1:A:703:ILE:CD1	1:A:714:ALA:HA	2.51	0.41
1:A:776:ASN:C	1:A:777:SER:HG	2.10	0.40
1:A:461:LEU:CD2	1:A:461:LEU:N	2.85	0.40
1:A:176:THR:O	1:A:180:LEU:HG	2.22	0.40
1:A:929:VAL:HG13	1:A:1009:PHE:HB2	2.02	0.40
1:A:846:GLN:HA	1:A:846:GLN:OE1	2.22	0.40
1:A:198:MET:SD	1:A:282:VAL:HG21	2.61	0.40
1:A:393:VAL:O	1:A:393:VAL:CG2	2.69	0.40
1:A:1003:SER:HA	1:A:1004:PRO:HD3	1.87	0.40
1:A:1055:LEU:O	1:A:1056:THR:C	2.59	0.40
1:A:278:ASP:HB2	1:A:792:LYS:CE	2.44	0.40
1:A:380:THR:O	1:A:435:CYS:HA	2.22	0.40
1:A:364:LYS:HE3	1:A:413:TRP:NE1	2.37	0.40
1:A:593:PHE:CE2	1:A:611:LEU:HD21	2.56	0.40
1:A:633:CYS:HB3	1:A:669:PHE:CE2	2.57	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	806/954 (84%)	759 (94%)	45 (6%)	2 (0%)	52 77

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	217	ASN
1	A	1093	GLY

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	740/854 (87%)	729 (98%)	11 (2%)	72	90

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	610	LEU
1	A	757	TYR
1	A	764	ILE
1	A	832	PHE
1	A	865	LEU
1	A	886	THR
1	A	907	LEU
1	A	939	THR
1	A	959	ASN
1	A	1026	LEU
1	A	1027	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	769	GLN
1	A	773	ASN
1	A	775	GLN
1	A	834	HIS
1	A	840	GLN
1	A	959	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](#)

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	NZU	A	1	-	26,30,30	2.08	11 (42%)	31,44,44	1.92	6 (19%)

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	NZU	A	1	-	-	0/12/12/12	0/4/4/4

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1	NZU	C7-C4	-4.59	1.39	1.44
2	A	1	NZU	C3-N10	-3.86	1.29	1.35
2	A	1	NZU	C5-N6	-3.04	1.27	1.33
2	A	1	NZU	C17-N18	-2.75	1.28	1.33
2	A	1	NZU	C7-N8	-2.53	1.29	1.33
2	A	1	NZU	C1-N2	-2.52	1.43	1.46
2	A	1	NZU	C9-N10	-2.46	1.28	1.33
2	A	1	NZU	C14-C15	-2.34	1.37	1.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1	NZU	O26-S25	-2.31	1.38	1.44
2	A	1	NZU	C4-C3	-2.30	1.37	1.43
2	A	1	NZU	C24-S25	2.38	1.81	1.78

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1	NZU	N10-C9-N8	-4.43	123.34	126.20
2	A	1	NZU	O27-S25-C24	2.43	109.84	108.28
2	A	1	NZU	C7-N8-C9	3.09	120.61	116.81
2	A	1	NZU	C9-N10-C3	3.76	120.54	115.00
2	A	1	NZU	C5-N6-N2	4.50	108.90	104.24
2	A	1	NZU	C1-N2-N6	4.73	125.91	118.05

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	NZU	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, 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	818/954 (85%)	0.29	62 (7%) 17 12	43, 91, 157, 208	1 (0%)

All (62) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1000	LYS	7.8
1	A	895	THR	6.7
1	A	1090	LEU	6.3
1	A	374	PRO	5.8
1	A	823	LEU	5.5
1	A	824	SER	4.8
1	A	982	ARG	4.6
1	A	1091	VAL	4.5
1	A	779	LEU	4.4
1	A	1089	HIS	4.4
1	A	1041	GLN	4.4
1	A	1042	LEU	4.2
1	A	758	ASP	4.2
1	A	971	ASN	4.0
1	A	894	SER	3.9
1	A	778	GLN	3.9
1	A	231	GLN	3.9
1	A	1094	ILE	3.9
1	A	757	TYR	3.9
1	A	210	TYR	3.7
1	A	1001	LYS	3.7
1	A	972	TYR	3.7
1	A	1007	GLN	3.6
1	A	777	SER	3.6
1	A	898	ASN	3.6
1	A	148	GLN	3.5
1	A	999	GLY	3.5

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Mol	Chain	Res	Type	RSRZ
1	A	319	ARG	3.5
1	A	248	PHE	3.3
1	A	613	ARG	3.3
1	A	1092	LEU	3.3
1	A	825	ASN	3.3
1	A	373	LEU	3.2
1	A	149	ALA	3.0
1	A	213	LYS	3.0
1	A	901	ALA	3.0
1	A	320	LYS	2.9
1	A	759	VAL	2.7
1	A	998	SER	2.6
1	A	1012	ILE	2.5
1	A	359	ARG	2.5
1	A	549	ASN	2.5
1	A	615	GLU	2.5
1	A	809	LYS	2.4
1	A	226	ARG	2.4
1	A	477	ARG	2.4
1	A	926	GLU	2.4
1	A	270	PHE	2.4
1	A	307	LEU	2.4
1	A	295	HIS	2.3
1	A	497	PHE	2.3
1	A	970	GLY	2.2
1	A	246	GLN	2.2
1	A	807	LYS	2.2
1	A	987	LEU	2.2
1	A	521	ASP	2.2
1	A	909	HIS	2.1
1	A	429	LEU	2.1
1	A	832	PHE	2.1
1	A	991	PHE	2.1
1	A	391	GLN	2.0
1	A	217	ASN	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 [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	NZU	A	1	27/27	0.90	0.17	-0.10	70,77,118,124	0

6.5 Other polymers [i](#)

There are no such residues in this entry.