



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 27, 2014 – 11:38 AM GMT

PDB ID : 1E90
Title : STRUCTURE DETERMINANTS OF PHOSPHOINOSITIDE 3-KINASE INHIBITION BY WORTMANNIN, LY294002, QUERCETIN, MYRICETIN AND STAUROSPORINE
Authors : Walker, E.H.; Pacold, M.E.; Perisic, O.; Stephens, L.; Hawkins, P.T.; Wymann, M.P.; Williams, R.L.
Deposited on : 2000-10-03
Resolution : 2.70 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

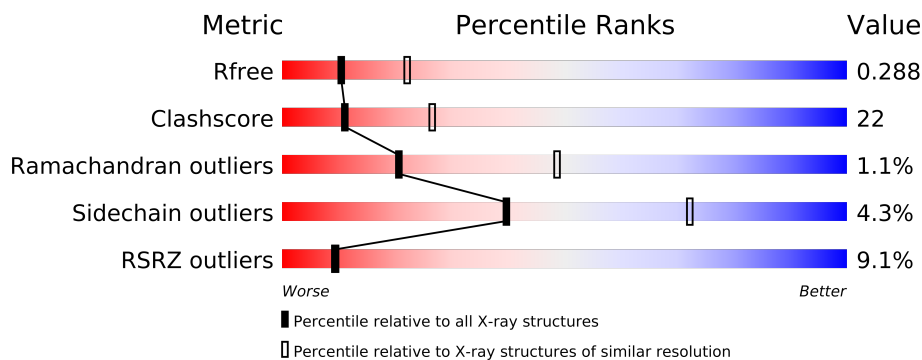
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

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}	66092	1557 (2.70-2.70)
Clashscore	79885	1939 (2.70-2.70)
Ramachandran outliers	78287	1905 (2.70-2.70)
Sidechain outliers	78261	1905 (2.70-2.70)
RSRZ outliers	66119	1559 (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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	961	

2 Entry composition i

There are 2 unique types of molecules in this entry. The entry contains 6862 atoms, of which 0 are hydrogen and 0 are deuterium.

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.

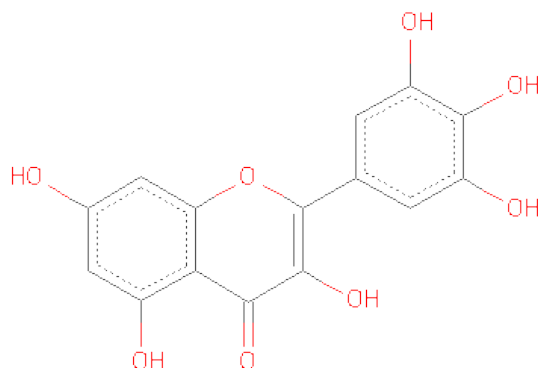
- Molecule 1 is a protein called PHOSPHATIDYLINOSITOL 3-KINASE CATALYTIC SUB-UNIT.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	844	6839	4399	1160	1244	36	0	0	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	143	ALA	PRO	EXPRESSION TAG	UNP O02697
A	505	ALA	ARG	CONFLICT	UNP O02697

- Molecule 2 is 3,5,7-TRIHYDROXY-2-(3,4,5-TRIHYDROXYPHENYL)-4H-CHROMEN-4-ONE (three-letter code: MYC) (formula: C₁₅H₁₀O₈).

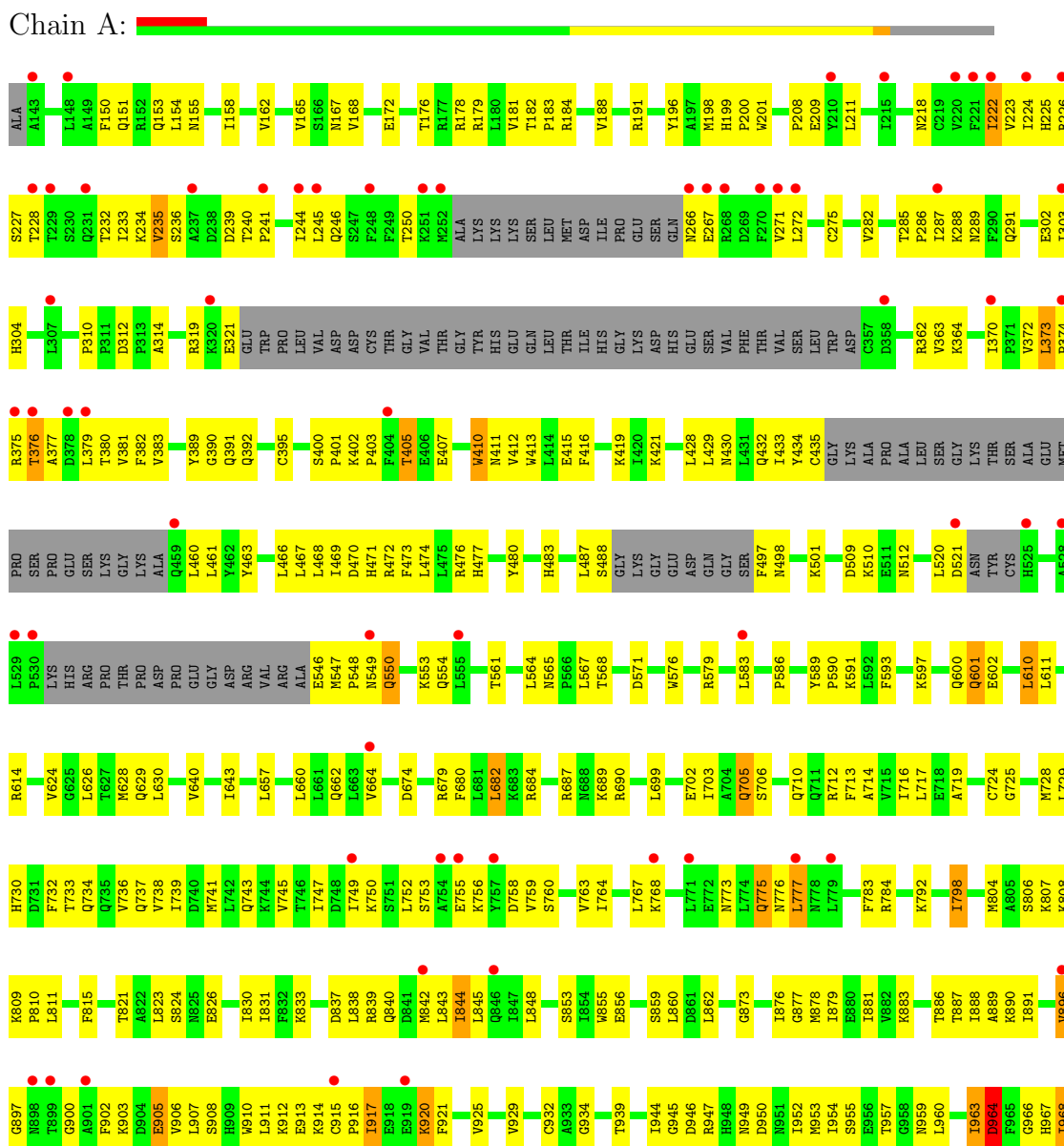


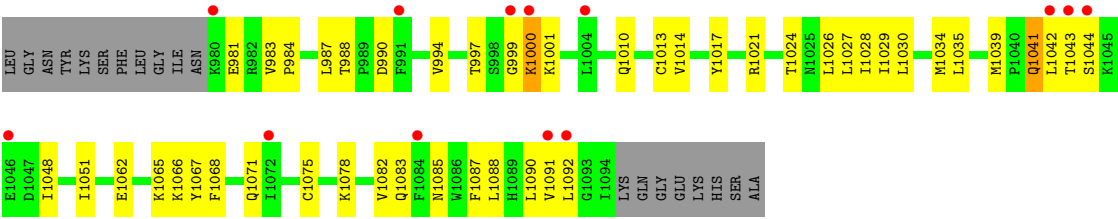
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
2	A	1	23	15	8	0	0

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 3-KINASE CATALYTIC SUBUNIT





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	143.68Å 67.73Å 106.99Å 90.00° 95.94° 90.00°	Depositor
Resolution (Å)	39.30 – 2.70 39.30 – 2.60	Depositor EDS
% Data completeness (in resolution range)	99.0 (39.30-2.70) 99.0 (39.30-2.60)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.42 (at 2.61Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.272 , 0.302 0.269 , 0.288	Depositor DCC
R_{free} test set	1381 reflections (4.92%)	DCC
Wilson B-factor (Å ²)	68.2	Xtriage
Anisotropy	0.372	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 50.4	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 31426 reflections	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	6862	wwPDB-VP
Average B, all atoms (Å ²)	76.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.91% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

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

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	0.31	0/6981	0.53	0/9442

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	6839	0	6917	310	0
2	A	23	0	5	0	0
All	All	6862	0	6922	310	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 22.

All (310) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:917:ILE:H	1:A:917:ILE:HD12	1.20	1.06
1:A:745:VAL:HG12	1:A:749:ILE:HD11	1.33	1.05

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:807:LYS:HD2	1:A:807:LYS:H	1.17	1.05
1:A:887:THR:HB	1:A:890:LYS:HG3	1.53	0.88
1:A:843:LEU:HD23	1:A:1034:MET:HG3	1.54	0.87
1:A:1035:LEU:HB3	1:A:1043:THR:HG21	1.56	0.85
1:A:840:GLN:HE22	1:A:967:HIS:HA	1.43	0.83
1:A:241:PRO:HA	1:A:244:ILE:HD12	1.60	0.83
1:A:807:LYS:HD2	1:A:807:LYS:N	1.94	0.81
1:A:917:ILE:HD13	1:A:920:LYS:HD2	1.62	0.81
1:A:1041:GLN:NE2	1:A:1041:GLN:H	1.78	0.80
1:A:611:LEU:O	1:A:614:ARG:HG3	1.82	0.80
1:A:561:THR:HG21	1:A:565:ASN:ND2	1.96	0.79
1:A:760:SER:O	1:A:763:VAL:HG12	1.81	0.79
1:A:549:ASN:OD1	1:A:553:LYS:HE3	1.83	0.79
1:A:1013:CYS:HB3	1:A:1068:PHE:HE2	1.48	0.77
1:A:497:PHE:HB3	1:A:1042:LEU:HB3	1.67	0.76
1:A:944:ILE:HG22	1:A:968:ILE:HD12	1.68	0.76
1:A:576:TRP:O	1:A:579:ARG:HG3	1.87	0.75
1:A:912:LYS:HG2	1:A:921:PHE:CE1	2.21	0.75
1:A:749:ILE:HD13	1:A:811:LEU:HD21	1.68	0.75
1:A:597:LYS:HD3	1:A:600:GLN:NE2	2.02	0.74
1:A:733:THR:HG22	1:A:737:GLN:HE21	1.53	0.73
1:A:1013:CYS:HB3	1:A:1068:PHE:CE2	2.23	0.73
1:A:182:THR:HB	1:A:183:PRO:HD3	1.69	0.72
1:A:855:TRP:HB3	1:A:860:LEU:HB2	1.70	0.72
1:A:208:PRO:HD2	1:A:211:LEU:HD12	1.70	0.72
1:A:640:VAL:O	1:A:643:ILE:HG13	1.91	0.71
1:A:807:LYS:H	1:A:807:LYS:CD	1.99	0.70
1:A:747:ILE:HD11	1:A:876:ILE:HD13	1.73	0.70
1:A:921:PHE:O	1:A:925:VAL:HG23	1.91	0.70
1:A:184:ARG:O	1:A:188:VAL:HG23	1.94	0.68
1:A:777:LEU:O	1:A:777:LEU:HD23	1.94	0.68
1:A:739:ILE:O	1:A:743:GLN:HG2	1.93	0.67
1:A:222:ILE:HD13	1:A:222:ILE:H	1.59	0.67
1:A:629:GLN:HG2	1:A:1029:ILE:HD13	1.76	0.67
1:A:840:GLN:NE2	1:A:967:HIS:HA	2.09	0.67
1:A:917:ILE:HD13	1:A:920:LYS:CD	2.24	0.67
1:A:844:ILE:O	1:A:848:LEU:HD13	1.95	0.66
1:A:176:THR:HG23	1:A:674:ASP:HB2	1.78	0.66
1:A:753:SER:HB3	1:A:809:LYS:NZ	2.11	0.66
1:A:914:LYS:C	1:A:916:PRO:HD3	2.15	0.66
1:A:887:THR:HB	1:A:890:LYS:CG	2.25	0.66
1:A:908:SER:O	1:A:912:LYS:HG3	1.95	0.66

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:244:ILE:HD11	1:A:287:ILE:CG2	2.26	0.65
1:A:550:GLN:O	1:A:550:GLN:NE2	2.24	0.65
1:A:312:ASP:OD2	1:A:314:ALA:HB3	1.96	0.65
1:A:1000:LYS:HE3	1:A:1000:LYS:HA	1.79	0.65
1:A:466:LEU:HD11	1:A:476:ARG:HD3	1.77	0.64
1:A:917:ILE:H	1:A:917:ILE:CD1	1.98	0.64
1:A:470:ASP:OD2	1:A:474:LEU:HB2	1.98	0.64
1:A:546:GLU:HG3	1:A:547:MET:N	2.12	0.64
1:A:402:LYS:HB3	1:A:403:PRO:HD2	1.79	0.64
1:A:903:LYS:HB2	1:A:906:VAL:HG23	1.80	0.64
1:A:501:LYS:NZ	1:A:501:LYS:HB3	2.13	0.64
1:A:201:TRP:NE1	1:A:291:GLN:HG3	2.12	0.63
1:A:747:ILE:HD11	1:A:876:ILE:CD1	2.28	0.63
1:A:983:VAL:HG22	1:A:984:PRO:HD2	1.81	0.63
1:A:917:ILE:N	1:A:917:ILE:HD12	2.04	0.62
1:A:597:LYS:HD3	1:A:600:GLN:HE22	1.64	0.62
1:A:939:THR:CG2	1:A:945:GLY:HA2	2.30	0.62
1:A:487:LEU:HD23	1:A:488:SER:N	2.14	0.62
1:A:583:LEU:HD22	1:A:610:LEU:CD2	2.30	0.62
1:A:891:ILE:HG22	1:A:906:VAL:HG12	1.82	0.61
1:A:705:GLN:HE21	1:A:873:GLY:C	2.03	0.61
1:A:564:LEU:HD11	1:A:1048:ILE:HG22	1.83	0.61
1:A:983:VAL:CG2	1:A:984:PRO:HD2	2.31	0.61
1:A:472:ARG:O	1:A:473:PHE:HB2	2.00	0.61
1:A:561:THR:HG21	1:A:565:ASN:HD22	1.66	0.60
1:A:432:GLN:HB3	1:A:460:LEU:HD11	1.83	0.60
1:A:239:ASP:O	1:A:287:ILE:HG23	2.00	0.60
1:A:246:GLN:O	1:A:250:THR:HG23	2.02	0.60
1:A:564:LEU:HD21	1:A:1048:ILE:HG21	1.83	0.60
1:A:275:CYS:HB3	1:A:823:LEU:HD13	1.83	0.60
1:A:1035:LEU:HA	1:A:1039:MET:HG2	1.82	0.60
1:A:712:ARG:O	1:A:716:ILE:HD13	2.02	0.60
1:A:376:THR:O	1:A:376:THR:HG23	2.00	0.59
1:A:886:THR:HG22	1:A:887:THR:N	2.17	0.59
1:A:303:ILE:HD12	1:A:303:ILE:N	2.17	0.59
1:A:764:ILE:HG22	1:A:768:LYS:HE2	1.83	0.59
1:A:209:GLU:HB3	1:A:859:SER:HB3	1.85	0.59
1:A:165:VAL:O	1:A:165:VAL:HG12	2.03	0.59
1:A:947:ARG:NH1	1:A:964:ASP:O	2.27	0.59
1:A:302:GLU:HB2	1:A:304:HIS:CE1	2.38	0.58
1:A:567:LEU:HD21	1:A:591:LYS:HD2	1.85	0.58
1:A:887:THR:HG22	1:A:889:ALA:H	1.68	0.58

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:302:GLU:HB2	1:A:304:HIS:HE1	1.68	0.58
1:A:200:PRO:HG3	1:A:282:VAL:HG23	1.84	0.58
1:A:939:THR:HG23	1:A:944:ILE:HG13	1.83	0.58
1:A:732:PHE:O	1:A:736:VAL:HG23	2.03	0.58
1:A:550:GLN:HE21	1:A:550:GLN:C	2.06	0.58
1:A:624:VAL:O	1:A:628:MET:HG2	2.03	0.58
1:A:946:ASP:O	1:A:947:ARG:HG2	2.04	0.57
1:A:831:ILE:HD11	1:A:881:ILE:HD11	1.86	0.57
1:A:810:PRO:HG3	1:A:833:LYS:HE3	1.85	0.57
1:A:235:VAL:HG12	1:A:239:ASP:OD2	2.03	0.57
1:A:843:LEU:CD2	1:A:1034:MET:HG3	2.29	0.57
1:A:981:GLU:OE1	1:A:1078:LYS:HE3	2.05	0.57
1:A:1048:ILE:O	1:A:1051:ILE:HG22	2.04	0.57
1:A:752:LEU:O	1:A:753:SER:HB3	2.05	0.57
1:A:233:ILE:N	1:A:233:ILE:HD12	2.20	0.57
1:A:477:HIS:O	1:A:480:TYR:HE1	1.88	0.56
1:A:725:GLY:O	1:A:729:LEU:HB2	2.05	0.56
1:A:753:SER:HB3	1:A:809:LYS:HZ2	1.70	0.56
1:A:755:GLU:HG2	1:A:755:GLU:O	2.05	0.56
1:A:469:ILE:N	1:A:469:ILE:HD12	2.19	0.56
1:A:568:THR:H	1:A:571:ASP:HB2	1.69	0.56
1:A:773:ASN:HA	1:A:776:ASN:ND2	2.21	0.56
1:A:703:ILE:HD11	1:A:714:ALA:N	2.21	0.56
1:A:429:LEU:HB2	1:A:468:LEU:HD21	1.88	0.56
1:A:839:ARG:HA	1:A:842:MET:HE2	1.89	0.55
1:A:910:TRP:O	1:A:914:LYS:HG2	2.06	0.55
1:A:749:ILE:CD1	1:A:811:LEU:HD21	2.35	0.55
1:A:498:ASN:ND2	1:A:1042:LEU:HD23	2.21	0.55
1:A:724:CYS:HB2	1:A:728:MET:HE3	1.88	0.55
1:A:741:MET:O	1:A:745:VAL:HG23	2.06	0.55
1:A:745:VAL:O	1:A:749:ILE:HG13	2.06	0.55
1:A:963:ILE:O	1:A:964:ASP:C	2.44	0.55
1:A:903:LYS:HB3	1:A:905:GLU:OE1	2.06	0.55
1:A:432:GLN:HB3	1:A:460:LEU:CD1	2.36	0.55
1:A:244:ILE:HD11	1:A:287:ILE:HG21	1.88	0.54
1:A:915:CYS:HB3	1:A:920:LYS:HG2	1.88	0.54
1:A:939:THR:HG22	1:A:945:GLY:HA2	1.88	0.54
1:A:806:SER:HB3	1:A:810:PRO:HD3	1.90	0.54
1:A:364:LYS:HE2	1:A:411:ASN:OD1	2.06	0.54
1:A:831:ILE:HB	1:A:879:ILE:HB	1.90	0.54
1:A:997:THR:HG23	1:A:1001:LYS:O	2.07	0.54
1:A:689:LYS:HG2	1:A:728:MET:CE	2.38	0.54

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:662:GLN:HE21	1:A:1030:LEU:HD22	1.74	0.53
1:A:1035:LEU:HD12	1:A:1048:ILE:HD13	1.91	0.53
1:A:509:ASP:OD2	1:A:512:ASN:HB2	2.08	0.53
1:A:826:GLU:HB3	1:A:883:LYS:HE2	1.90	0.53
1:A:225:HIS:CE1	1:A:823:LEU:HD21	2.43	0.53
1:A:837:ASP:OD1	1:A:839:ARG:HB2	2.09	0.53
1:A:840:GLN:O	1:A:844:ILE:HD13	2.09	0.53
1:A:743:GLN:O	1:A:747:ILE:HD13	2.08	0.52
1:A:657:LEU:HD11	1:A:690:ARG:HD3	1.91	0.52
1:A:381:VAL:HG12	1:A:435:CYS:HA	1.91	0.52
1:A:1014:VAL:HG11	1:A:1065:LYS:HG3	1.91	0.52
1:A:245:LEU:HD21	1:A:272:LEU:HG	1.90	0.52
1:A:240:THR:O	1:A:244:ILE:HG13	2.10	0.52
1:A:705:GLN:NE2	1:A:873:GLY:O	2.43	0.52
1:A:266:ASN:CG	1:A:267:GLU:H	2.12	0.52
1:A:749:ILE:HD13	1:A:767:LEU:HD23	1.91	0.52
1:A:887:THR:HG22	1:A:889:ALA:N	2.25	0.52
1:A:821:THR:HG22	1:A:821:THR:O	2.10	0.51
1:A:232:THR:C	1:A:233:ILE:HD12	2.31	0.51
1:A:287:ILE:HD12	1:A:288:LYS:N	2.25	0.51
1:A:1041:GLN:CD	1:A:1041:GLN:H	2.14	0.51
1:A:1010:GLN:O	1:A:1014:VAL:HG23	2.11	0.51
1:A:1088:LEU:HD23	1:A:1092:LEU:HD12	1.92	0.51
1:A:383:VAL:HG22	1:A:433:ILE:CD1	2.41	0.51
1:A:375:ARG:O	1:A:377:ALA:N	2.44	0.51
1:A:1062:GLU:O	1:A:1066:LYS:HD3	2.10	0.50
1:A:912:LYS:HE2	1:A:921:PHE:CZ	2.47	0.50
1:A:1024:THR:O	1:A:1028:ILE:HG12	2.11	0.50
1:A:228:THR:O	1:A:228:THR:HG22	2.10	0.50
1:A:470:ASP:HB3	1:A:476:ARG:NH2	2.27	0.50
1:A:932:CYS:HA	1:A:960:LEU:HD23	1.92	0.50
1:A:554:GLN:HA	1:A:554:GLN:NE2	2.26	0.50
1:A:1041:GLN:N	1:A:1041:GLN:NE2	2.54	0.50
1:A:602:GLU:H	1:A:602:GLU:CD	2.15	0.50
1:A:184:ARG:HH22	1:A:321:GLU:CD	2.15	0.50
1:A:784:ARG:HG2	1:A:784:ARG:HH11	1.77	0.49
1:A:660:LEU:O	1:A:664:VAL:HG23	2.11	0.49
1:A:162:VAL:HG12	1:A:714:ALA:HB1	1.94	0.49
1:A:896:VAL:HG12	1:A:897:GLY:N	2.27	0.49
1:A:497:PHE:HB3	1:A:1042:LEU:HD22	1.93	0.49
1:A:939:THR:HG21	1:A:945:GLY:HA2	1.94	0.49
1:A:379:LEU:HB3	1:A:435:CYS:SG	2.53	0.49

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:372:VAL:HG12	1:A:373:LEU:N	2.28	0.49
1:A:165:VAL:HG13	1:A:168:VAL:HG21	1.95	0.49
1:A:808:LYS:HD2	1:A:833:LYS:HE2	1.95	0.49
1:A:705:GLN:OE1	1:A:839:ARG:NE	2.46	0.48
1:A:303:ILE:HD12	1:A:303:ILE:H	1.77	0.48
1:A:806:SER:HB3	1:A:810:PRO:CD	2.43	0.48
1:A:747:ILE:HD12	1:A:747:ILE:N	2.29	0.48
1:A:223:VAL:HG12	1:A:225:HIS:CD2	2.48	0.48
1:A:165:VAL:HG13	1:A:168:VAL:CG2	2.43	0.48
1:A:853:SER:O	1:A:856:GLU:HB3	2.13	0.48
1:A:1067:TYR:O	1:A:1071:GLN:HG2	2.12	0.48
1:A:784:ARG:HG2	1:A:784:ARG:NH1	2.29	0.48
1:A:241:PRO:HA	1:A:244:ILE:CD1	2.40	0.48
1:A:200:PRO:HG3	1:A:282:VAL:CG2	2.44	0.48
1:A:1017:TYR:O	1:A:1021:ARG:HG3	2.14	0.48
1:A:589:TYR:HB2	1:A:590:PRO:HD3	1.95	0.48
1:A:167:ASN:O	1:A:168:VAL:HG13	2.14	0.47
1:A:370:ILE:O	1:A:370:ILE:HG23	2.14	0.47
1:A:886:THR:HG22	1:A:887:THR:H	1.79	0.47
1:A:188:VAL:O	1:A:191:ARG:HG2	2.15	0.47
1:A:419:LYS:HD3	1:A:421:LYS:HE2	1.95	0.47
1:A:549:ASN:O	1:A:553:LYS:HG3	2.15	0.47
1:A:428:LEU:HD23	1:A:467:LEU:HD23	1.95	0.47
1:A:756:LYS:HD3	1:A:756:LYS:HA	1.75	0.47
1:A:363:VAL:HG12	1:A:416:PHE:HE1	1.79	0.47
1:A:767:LEU:HD13	1:A:767:LEU:C	2.35	0.47
1:A:222:ILE:HD13	1:A:222:ILE:N	2.27	0.47
1:A:236:SER:HG	1:A:239:ASP:CG	2.18	0.47
1:A:501:LYS:HZ3	1:A:501:LYS:HB3	1.80	0.47
1:A:586:PRO:HA	1:A:589:TYR:CE1	2.50	0.47
1:A:702:GLU:O	1:A:706:SER:HB3	2.15	0.47
1:A:955:SER:C	1:A:957:THR:H	2.17	0.47
1:A:405:THR:HG23	1:A:407:GLU:O	2.15	0.47
1:A:233:ILE:HG22	1:A:234:LYS:N	2.30	0.47
1:A:568:THR:N	1:A:571:ASP:HB2	2.29	0.47
1:A:179:ARG:NH2	1:A:679:ARG:HH22	2.13	0.47
1:A:1087:PHE:O	1:A:1091:VAL:HB	2.15	0.47
1:A:662:GLN:HG3	1:A:1030:LEU:HD22	1.98	0.46
1:A:887:THR:HG21	1:A:950:ASP:OD1	2.16	0.46
1:A:750:LYS:HE3	1:A:808:LYS:HA	1.97	0.46
1:A:876:ILE:HG22	1:A:877:GLY:N	2.30	0.46
1:A:629:GLN:HG2	1:A:1029:ILE:CD1	2.45	0.46

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:586:PRO:HA	1:A:589:TYR:CD1	2.50	0.46
1:A:949:ASN:HB2	1:A:1083:GLN:NE2	2.31	0.46
1:A:783:PHE:N	1:A:792:LYS:HE2	2.30	0.46
1:A:687:ARG:O	1:A:687:ARG:HG2	2.16	0.45
1:A:463:TYR:CE1	1:A:501:LYS:HA	2.52	0.45
1:A:151:GLN:HG3	1:A:155:ASN:HD21	1.80	0.45
1:A:643:ILE:C	1:A:643:ILE:HD12	2.36	0.45
1:A:362:ARG:HB3	1:A:415:GLU:HG3	1.99	0.45
1:A:196:TYR:OH	1:A:728:MET:HE2	2.16	0.45
1:A:954:ILE:HD12	1:A:959:ASN:C	2.37	0.45
1:A:389:TYR:O	1:A:392:GLN:HG2	2.17	0.45
1:A:593:PHE:CZ	1:A:611:LEU:HD21	2.52	0.45
1:A:903:LYS:HB2	1:A:906:VAL:CG2	2.45	0.45
1:A:201:TRP:CE2	1:A:291:GLN:HG3	2.52	0.45
1:A:946:ASP:C	1:A:947:ARG:HG2	2.37	0.45
1:A:734:GLN:O	1:A:738:VAL:HG23	2.16	0.45
1:A:939:THR:HG22	1:A:945:GLY:CA	2.47	0.45
1:A:844:ILE:CD1	1:A:844:ILE:N	2.79	0.45
1:A:763:VAL:HG13	1:A:764:ILE:N	2.32	0.44
1:A:271:VAL:HG22	1:A:272:LEU:N	2.32	0.44
1:A:1014:VAL:CG1	1:A:1065:LYS:HG3	2.47	0.44
1:A:286:PRO:HD2	1:A:289:ASN:HD22	1.80	0.44
1:A:1000:LYS:HE3	1:A:1000:LYS:CA	2.48	0.44
1:A:703:ILE:HD11	1:A:713:PHE:CB	2.46	0.44
1:A:838:LEU:HD12	1:A:877:GLY:HA3	1.99	0.44
1:A:730:HIS:O	1:A:734:GLN:HG2	2.17	0.44
1:A:181:VAL:HG13	1:A:184:ARG:NH2	2.33	0.44
1:A:380:THR:O	1:A:435:CYS:HA	2.18	0.44
1:A:150:PHE:CE1	1:A:319:ARG:HD3	2.52	0.44
1:A:287:ILE:C	1:A:287:ILE:HD12	2.38	0.44
1:A:699:LEU:O	1:A:703:ILE:HG12	2.18	0.44
1:A:815:PHE:CE1	1:A:830:ILE:HD12	2.52	0.44
1:A:172:GLU:HG3	1:A:471:HIS:CG	2.53	0.43
1:A:498:ASN:C	1:A:498:ASN:OD1	2.57	0.43
1:A:824:SER:OG	1:A:826:GLU:HG3	2.17	0.43
1:A:804:MET:HE2	1:A:831:ILE:HG12	2.00	0.43
1:A:583:LEU:HD22	1:A:610:LEU:HD21	2.00	0.43
1:A:804:MET:C	1:A:806:SER:H	2.22	0.43
1:A:888:ILE:HA	1:A:891:ILE:HD12	2.00	0.43
1:A:410:TRP:HB3	1:A:412:VAL:HG23	1.99	0.43
1:A:862:LEU:HD12	1:A:934:GLY:HA2	2.00	0.43
1:A:1044:SER:O	1:A:1048:ILE:HG12	2.19	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:382:PHE:CE1	1:A:434:TYR:HB2	2.54	0.43
1:A:225:HIS:HE1	1:A:823:LEU:HD21	1.82	0.43
1:A:364:LYS:HD2	1:A:413:TRP:CE2	2.54	0.43
1:A:886:THR:CG2	1:A:887:THR:N	2.82	0.43
1:A:953:MET:HG3	1:A:963:ILE:HD13	2.01	0.42
1:A:419:LYS:CD	1:A:421:LYS:HE2	2.49	0.42
1:A:990:ASP:O	1:A:994:VAL:HG23	2.19	0.42
1:A:967:HIS:O	1:A:968:ILE:O	2.38	0.42
1:A:150:PHE:O	1:A:153:GLN:HG2	2.20	0.42
1:A:497:PHE:CB	1:A:1042:LEU:HD22	2.49	0.42
1:A:905:GLU:N	1:A:905:GLU:OE1	2.37	0.42
1:A:1083:GLN:HA	1:A:1083:GLN:NE2	2.34	0.42
1:A:285:THR:HG22	1:A:289:ASN:HB2	2.02	0.42
1:A:223:VAL:HG12	1:A:225:HIS:NE2	2.34	0.42
1:A:400:SER:HA	1:A:401:PRO:HD3	1.94	0.42
1:A:745:VAL:HG12	1:A:749:ILE:CD1	2.24	0.42
1:A:225:HIS:N	1:A:225:HIS:CD2	2.88	0.42
1:A:178:ARG:O	1:A:181:VAL:HG23	2.19	0.42
1:A:753:SER:HB3	1:A:809:LYS:HZ1	1.81	0.42
1:A:568:THR:H	1:A:571:ASP:CB	2.31	0.42
1:A:435:CYS:HB2	1:A:461:LEU:HD12	2.02	0.42
1:A:154:LEU:O	1:A:158:ILE:HG13	2.20	0.42
1:A:967:HIS:O	1:A:968:ILE:C	2.58	0.41
1:A:798:ILE:HG13	1:A:798:ILE:H	1.69	0.41
1:A:497:PHE:CB	1:A:1042:LEU:HB3	2.45	0.41
1:A:952:ILE:CG2	1:A:960:LEU:HD11	2.50	0.41
1:A:162:VAL:CG1	1:A:714:ALA:HB1	2.50	0.41
1:A:999:GLY:C	1:A:1001:LYS:H	2.23	0.41
1:A:266:ASN:CG	1:A:267:GLU:N	2.72	0.41
1:A:896:VAL:HG12	1:A:897:GLY:H	1.85	0.41
1:A:831:ILE:O	1:A:878:MET:HA	2.21	0.41
1:A:601:GLN:HG3	1:A:602:GLU:N	2.35	0.41
1:A:430:ASN:OD1	1:A:432:GLN:HG3	2.20	0.41
1:A:165:VAL:O	1:A:165:VAL:CG1	2.69	0.41
1:A:373:LEU:HB2	1:A:374:PRO:CD	2.50	0.41
1:A:271:VAL:HB	1:A:310:PRO:HG3	2.03	0.41
1:A:682:LEU:HD11	1:A:719:ALA:HB1	2.03	0.41
1:A:630:LEU:CD1	1:A:643:ILE:HD11	2.51	0.41
1:A:1088:LEU:O	1:A:1092:LEU:HB2	2.21	0.41
1:A:390:GLY:O	1:A:391:GLN:HB3	2.21	0.41
1:A:900:GLY:O	1:A:902:PHE:N	2.47	0.41
1:A:224:ILE:HD12	1:A:272:LEU:CD2	2.51	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:775:GLN:O	1:A:776:ASN:C	2.57	0.40
1:A:373:LEU:HB2	1:A:374:PRO:HD2	2.02	0.40
1:A:987:LEU:HB3	1:A:1075:CYS:HB3	2.03	0.40
1:A:831:ILE:CD1	1:A:881:ILE:HD11	2.50	0.40
1:A:703:ILE:HD13	1:A:710:GLN:HA	2.02	0.40
1:A:988:THR:HG21	1:A:1083:GLN:HG3	2.03	0.40
1:A:520:LEU:O	1:A:521:ASP:C	2.59	0.40
1:A:198:MET:O	1:A:199:HIS:C	2.60	0.40
1:A:467:LEU:O	1:A:476:ARG:HD2	2.21	0.40
1:A:1082:VAL:O	1:A:1085:ASN:HB2	2.21	0.40
1:A:680:PHE:O	1:A:684:ARG:HG2	2.22	0.40
1:A:925:VAL:O	1:A:929:VAL:HG23	2.21	0.40
1:A:483:HIS:CD2	1:A:510:LYS:HG2	2.57	0.40
1:A:983:VAL:HG22	1:A:984:PRO:CD	2.50	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone ⓘ

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	828/961 (86%)	752 (91%)	67 (8%)	9 (1%)	21 49

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	896	VAL
1	A	966	GLY
1	A	758	ASP
1	A	777	LEU
1	A	964	ASP
1	A	963	ILE
1	A	376	THR
1	A	798	ILE
1	A	759	VAL

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	761/857 (89%)	728 (96%)	33 (4%)	40 72

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	218	ASN
1	A	222	ILE
1	A	226	ARG
1	A	227	SER
1	A	235	VAL
1	A	373	LEU
1	A	395	CYS
1	A	405	THR
1	A	410	TRP
1	A	548	PRO
1	A	550	GLN
1	A	601	GLN
1	A	610	LEU
1	A	626	LEU
1	A	682	LEU
1	A	705	GLN
1	A	717	LEU
1	A	775	GLN
1	A	844	ILE
1	A	845	LEU
1	A	905	GLU
1	A	907	LEU
1	A	911	LEU
1	A	913	GLU
1	A	917	ILE
1	A	920	LYS
1	A	964	ASP
1	A	968	ILE
1	A	1000	LYS
1	A	1026	LEU
1	A	1027	LEU

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Mol	Chain	Res	Type
1	A	1041	GLN
1	A	1090	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	299	ASN
1	A	304	HIS
1	A	396	GLN
1	A	459	GLN
1	A	550	GLN
1	A	554	GLN
1	A	565	ASN
1	A	600	GLN
1	A	601	GLN
1	A	639	ASN
1	A	708	HIS
1	A	710	GLN
1	A	737	GLN
1	A	776	ASN
1	A	778	ASN
1	A	840	GLN
1	A	922	GLN
1	A	951	ASN
1	A	1007	GLN
1	A	1023	HIS
1	A	1041	GLN
1	A	1083	GLN
1	A	1089	HIS

5.3.3 RNA ⓘ

There are no RNA chains 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	MYC	A	2095	-	24,25,25	1.95	5 (20%)	35,38,38	1.62	6 (17%)

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	MYC	A	2095	-	-	0/4/4/4	0/1/3/3

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	2095	MYC	C9-C3	4.38	1.47	1.40
2	A	2095	MYC	C14-C11	3.93	1.51	1.46
2	A	2095	MYC	C1-C6	3.65	1.45	1.39
2	A	2095	MYC	C15-C16	3.04	1.43	1.38
2	A	2095	MYC	O12-C11	2.81	1.37	1.35

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	2095	MYC	C14-C11-C10	4.20	128.35	120.38
2	A	2095	MYC	C10-C9-C3	-4.15	116.86	121.41
2	A	2095	MYC	O12-C11-C10	-2.71	120.34	121.54
2	A	2095	MYC	C19-C18-C17	2.62	122.13	120.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	2095	MYC	C14-C19-C18	-2.58	119.90	121.03
2	A	2095	MYC	C18-C17-C16	-2.55	118.21	119.56

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

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	844/961 (87%)	0.57	76 (9%) 10 9	27, 72, 128, 184	0

All (76) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	898	ASN	7.6
1	A	375	ARG	6.8
1	A	529	LEU	6.2
1	A	245	LEU	5.7
1	A	248	PHE	5.6
1	A	919	GLU	5.3
1	A	777	LEU	5.1
1	A	252	MET	4.9
1	A	404	PHE	4.8
1	A	530	PRO	4.6
1	A	148	LEU	4.1
1	A	754	ALA	4.0
1	A	307	LEU	3.9
1	A	270	PHE	3.8
1	A	525	HIS	3.7
1	A	272	LEU	3.6
1	A	266	ASN	3.6
1	A	379	LEU	3.5
1	A	549	ASN	3.5
1	A	226	ARG	3.5
1	A	376	THR	3.4
1	A	757	TYR	3.4
1	A	268	ARG	3.4
1	A	755	GLU	3.3
1	A	251	LYS	3.3
1	A	244	ILE	3.2
1	A	224	ILE	3.2

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Mol	Chain	Res	Type	RSRZ
1	A	374	PRO	3.2
1	A	271	VAL	3.1
1	A	143	ALA	3.1
1	A	528	ALA	3.1
1	A	222	ILE	3.0
1	A	228	THR	2.9
1	A	1092	LEU	2.9
1	A	1091	VAL	2.9
1	A	267	GLU	2.9
1	A	220	VAL	2.8
1	A	896	VAL	2.8
1	A	999	GLY	2.8
1	A	229	THR	2.8
1	A	521	ASP	2.7
1	A	664	VAL	2.7
1	A	320	LYS	2.7
1	A	1043	THR	2.7
1	A	749	ILE	2.6
1	A	459	GLN	2.6
1	A	991	PHE	2.6
1	A	1084	PHE	2.6
1	A	231	GLN	2.6
1	A	1042	LEU	2.5
1	A	779	LEU	2.5
1	A	287	ILE	2.5
1	A	768	LYS	2.4
1	A	303	ILE	2.4
1	A	358	ASP	2.4
1	A	370	ILE	2.4
1	A	555	LEU	2.3
1	A	378	ASP	2.3
1	A	215	ILE	2.2
1	A	846	GLN	2.2
1	A	210	TYR	2.2
1	A	583	LEU	2.2
1	A	1000	LYS	2.2
1	A	842	MET	2.2
1	A	980	LYS	2.2
1	A	1004	LEU	2.2
1	A	901	ALA	2.1
1	A	1072	ILE	2.1
1	A	915	CYS	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	1044	SER	2.1
1	A	237	ALA	2.1
1	A	221	PHE	2.1
1	A	771	LEU	2.1
1	A	241	PRO	2.1
1	A	899	THR	2.0
1	A	1046	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 ⓘ

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	RSR	LLDF	B-factors(Å ²)	Q<0.9
2	MYC	A	2095	23/23	0.25	1.02	93,93,93,93	0

6.5 Other polymers ⓘ

There are no such residues in this entry.