



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 25, 2026 – 03:54 AM EDT

PDB ID : 1YMY / pdb_00001ymy
Title : Crystal Structure of the N-Acetylglucosamine-6-phosphate deacetylase from Escherichia coli K12
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Deposited on : 2005-01-21
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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

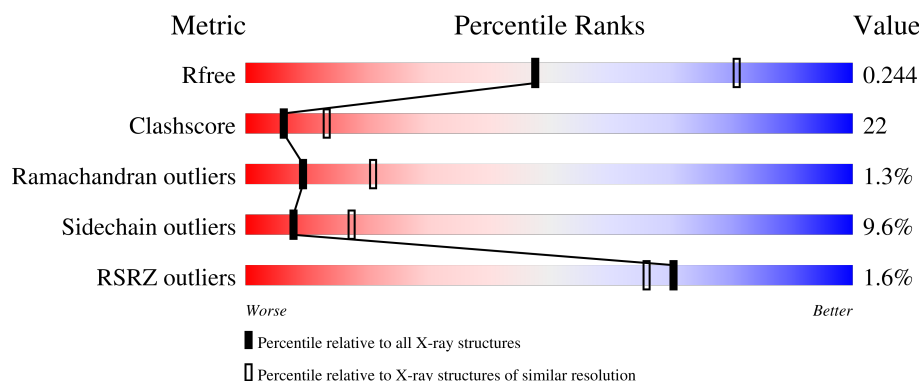
1 Overall quality at a glance

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}	180053	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. 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	382	 59% 34% 6% .
1	B	382	 51% 29% 6% . 13%

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 5306 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.

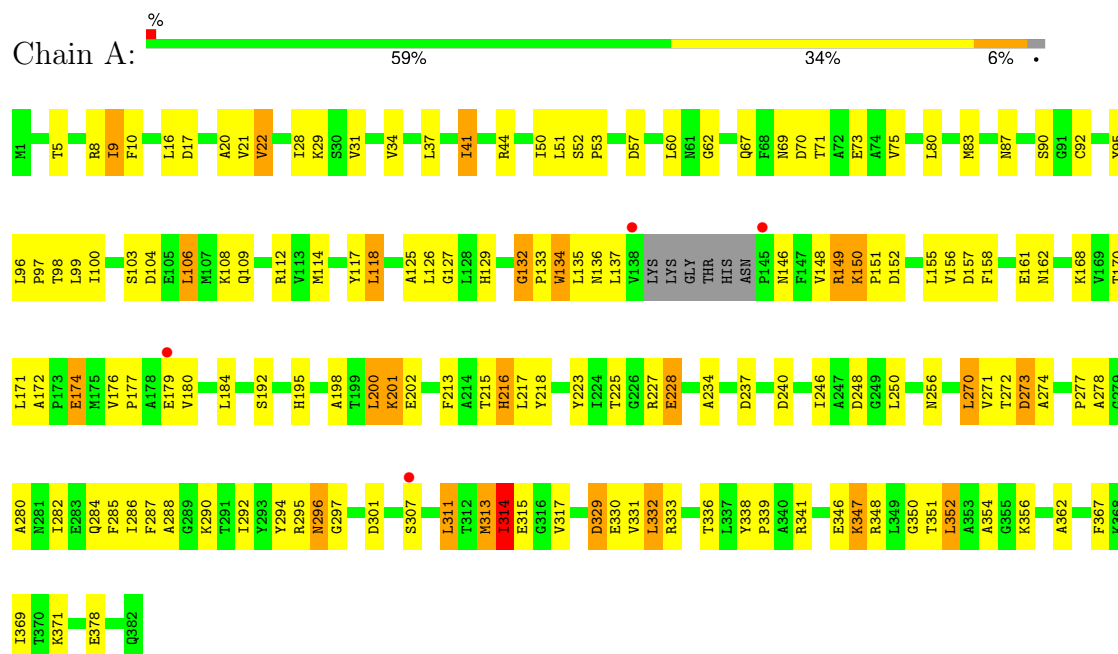
- Molecule 1 is a protein called N-acetylglucosamine-6-phosphate deacetylase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	376	Total	C	N	O	S	0	0	0
			2827	1789	483	539	16			
1	B	331	Total	C	N	O	S	0	0	0
			2479	1570	424	470	15			

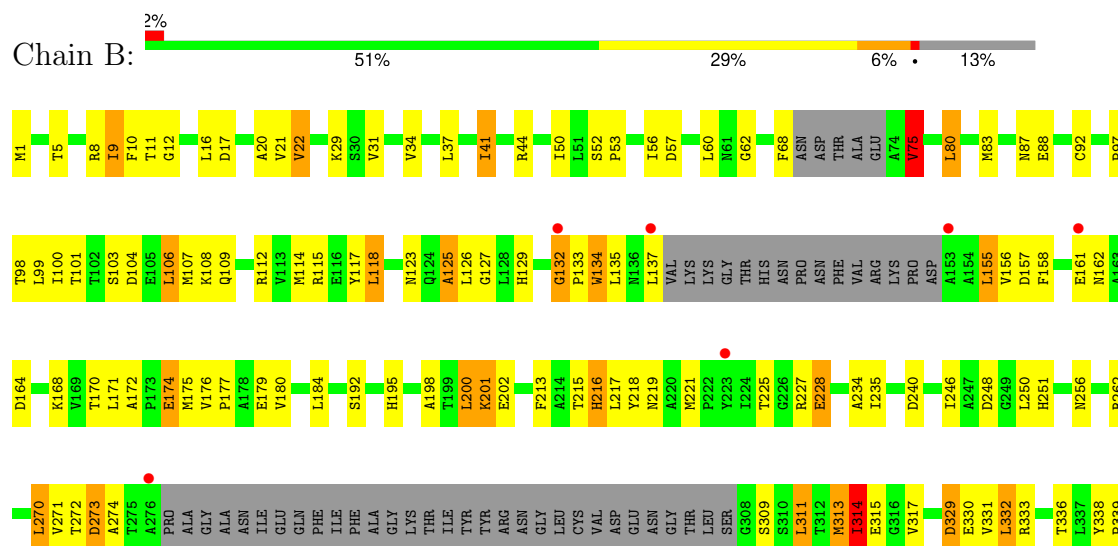
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes 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: N-acetylglucosamine-6-phosphate deacetylase



- Molecule 1: N-acetylglucosamine-6-phosphate deacetylase





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	82.23Å 114.02Å 80.22Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	25.00 – 2.60 25.00 – 2.60	Depositor EDS
% Data completeness (in resolution range)	(Not available) (25.00-2.60) 96.6 (25.00-2.60)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.10 (at 2.61Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.212 , 0.243 0.211 , 0.244	Depositor DCC
R_{free} test set	1114 reflections (4.83%)	wwPDB-VP
Wilson B-factor (Å ²)	47.2	Xtriage
Anisotropy	0.223	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 35.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.27$	Xtriage
Estimated twinning fraction	0.033 for l,-k,h	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5306	wwPDB-VP
Average B, all atoms (Å ²)	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.60% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

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	0.54	1/2873 (0.0%)	1.00	9/3900 (0.2%)
1	B	0.49	1/2515 (0.0%)	0.99	9/3410 (0.3%)
All	All	0.52	2/5388 (0.0%)	0.99	18/7310 (0.2%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	313	MET	SD-CE	-7.05	1.61	1.79
1	B	313	MET	SD-CE	-5.52	1.65	1.79

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	41	ILE	N-CA-C	6.08	117.34	108.53
1	A	132	GLY	N-CA-C	-6.03	100.04	112.34
1	A	350	GLY	N-CA-C	6.01	121.11	113.24
1	A	41	ILE	N-CA-C	6.01	117.24	108.53
1	A	126	LEU	N-CA-C	5.90	118.95	111.69
1	A	367	PHE	N-CA-C	5.90	119.83	111.92
1	B	367	PHE	N-CA-C	5.89	119.82	111.92
1	B	350	GLY	N-CA-C	5.89	120.96	113.24
1	B	132	GLY	N-CA-C	-5.82	100.48	112.34
1	B	125	ALA	N-CA-C	-5.49	101.52	109.59
1	A	280	ALA	N-CA-C	5.47	115.51	108.34
1	B	351	THR	N-CA-C	5.35	116.41	108.86
1	A	313	MET	N-CA-C	5.32	117.08	111.28
1	A	351	THR	N-CA-C	5.23	116.24	108.86
1	A	314	ILE	CB-CA-C	-5.10	105.25	112.14
1	B	313	MET	N-CA-C	5.09	116.83	111.28
1	B	314	ILE	CB-CA-C	-5.07	105.22	112.22
1	B	126	LEU	N-CA-C	5.05	118.59	112.23

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	2827	0	2861	127	0
1	B	2479	0	2527	112	0
All	All	5306	0	5388	237	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 22.

All (237) 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:201:LYS:HE3	1:A:201:LYS:H	1.08	1.07
1:B:201:LYS:HE3	1:B:201:LYS:H	1.08	1.06
1:B:272:THR:HG23	1:B:313:MET:HE2	1.40	0.99
1:A:272:THR:HG22	1:A:311:LEU:HD13	1.43	0.98
1:A:201:LYS:HE3	1:A:201:LYS:N	1.85	0.92
1:A:272:THR:HG23	1:A:313:MET:HE2	1.51	0.92
1:A:62:GLY:HA2	1:A:83:MET:HE3	1.52	0.90
1:B:201:LYS:HE3	1:B:201:LYS:N	1.86	0.89
1:B:272:THR:HG22	1:B:311:LEU:HD13	1.58	0.84
1:B:62:GLY:HA2	1:B:83:MET:HE3	1.60	0.82
1:B:114:MET:HE2	1:B:118:LEU:HD13	1.61	0.80
1:A:114:MET:HE2	1:A:118:LEU:HD13	1.63	0.80
1:A:149:ARG:HG2	1:A:149:ARG:HH11	1.46	0.79
1:A:152:ASP:O	1:A:156:VAL:HG23	1.85	0.76
1:A:157:ASP:O	1:A:161:GLU:HG3	1.86	0.76
1:A:5:THR:HG22	1:A:20:ALA:HB2	1.68	0.75
1:B:5:THR:HG22	1:B:20:ALA:HB2	1.68	0.75
1:A:69:ASN:ND2	1:A:100:ILE:H	1.83	0.75
1:B:157:ASP:O	1:B:161:GLU:HG3	1.87	0.74
1:B:272:THR:C	1:B:274:ALA:H	1.95	0.74
1:A:69:ASN:HD21	1:A:100:ILE:HG12	1.54	0.73
1:A:287:PHE:HB3	1:A:292:ILE:HD12	1.71	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:272:THR:HG23	1:B:313:MET:CE	2.20	0.72
1:A:129:HIS:CE1	1:A:168:LYS:HE3	2.25	0.72
1:A:272:THR:C	1:A:274:ALA:H	1.97	0.71
1:B:272:THR:O	1:B:274:ALA:N	2.24	0.70
1:A:69:ASN:ND2	1:A:100:ILE:HG12	2.07	0.69
1:B:60:LEU:CD1	1:B:274:ALA:HB3	2.22	0.68
1:B:129:HIS:HE1	1:B:170:THR:OG1	1.76	0.67
1:A:69:ASN:HD21	1:A:100:ILE:H	1.41	0.66
1:A:272:THR:O	1:A:274:ALA:N	2.26	0.66
1:A:103:SER:HB2	1:A:106:LEU:HB2	1.77	0.66
1:A:129:HIS:HE1	1:A:170:THR:OG1	1.80	0.65
1:A:277:PRO:HD3	1:A:282:ILE:HD11	1.77	0.64
1:B:103:SER:HB2	1:B:106:LEU:HB2	1.80	0.64
1:B:68:PHE:CE1	1:B:80:LEU:HD13	2.33	0.64
1:B:248:ASP:HB3	1:B:311:LEU:HG	1.78	0.64
1:B:5:THR:HG22	1:B:20:ALA:CB	2.28	0.63
1:A:62:GLY:CA	1:A:83:MET:HE3	2.29	0.63
1:B:60:LEU:HD12	1:B:274:ALA:HB3	1.80	0.63
1:B:129:HIS:CE1	1:B:168:LYS:HE3	2.35	0.62
1:B:272:THR:C	1:B:274:ALA:N	2.58	0.61
1:B:251:HIS:CE1	1:B:309:SER:HB3	2.35	0.61
1:A:5:THR:HG22	1:A:20:ALA:CB	2.28	0.61
1:A:34:VAL:O	1:A:37:LEU:HB2	2.00	0.60
1:B:347:LYS:NZ	1:B:347:LYS:HB3	2.16	0.60
1:A:294:TYR:CZ	1:A:297:GLY:HA2	2.36	0.60
1:A:272:THR:C	1:A:274:ALA:N	2.59	0.60
1:A:150:LYS:HB2	1:A:150:LYS:NZ	2.17	0.60
1:A:132:GLY:HA2	1:A:135:LEU:HG	1.84	0.60
1:B:314:ILE:HD12	1:B:315:GLU:H	1.66	0.60
1:A:286:ILE:HD12	1:A:286:ILE:N	2.16	0.59
1:A:60:LEU:HD12	1:A:274:ALA:HB3	1.83	0.59
1:B:34:VAL:O	1:B:37:LEU:HB2	2.01	0.59
1:A:192:SER:HA	1:A:213:PHE:O	2.02	0.59
1:B:115:ARG:NH1	1:B:164:ASP:OD2	2.36	0.58
1:A:172:ALA:HB1	1:A:174:GLU:OE2	2.04	0.58
1:A:347:LYS:NZ	1:A:347:LYS:HB3	2.19	0.58
1:B:172:ALA:HB1	1:B:174:GLU:OE2	2.03	0.58
1:A:314:ILE:HD12	1:A:315:GLU:H	1.69	0.58
1:A:272:THR:HG23	1:A:313:MET:CE	2.32	0.58
1:B:109:GLN:HG3	1:B:112:ARG:NH2	2.19	0.58
1:B:132:GLY:HA2	1:B:135:LEU:HG	1.86	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:117:TYR:CE1	1:A:125:ALA:HB3	2.39	0.57
1:B:329:ASP:OD2	1:B:329:ASP:N	2.37	0.57
1:A:329:ASP:OD2	1:A:329:ASP:N	2.37	0.57
1:B:192:SER:HA	1:B:213:PHE:O	2.04	0.56
1:A:57:ASP:CG	1:A:313:MET:HE3	2.32	0.55
1:A:71:THR:HB	1:A:73:GLU:CD	2.31	0.55
1:A:176:VAL:HG12	1:A:177:PRO:O	2.06	0.55
1:B:176:VAL:HG12	1:B:177:PRO:O	2.07	0.54
1:B:117:TYR:CE1	1:B:125:ALA:HB3	2.43	0.54
1:A:87:ASN:HB3	1:A:92:CYS:HB3	1.89	0.54
1:A:150:LYS:N	1:A:151:PRO:HD2	2.23	0.54
1:A:137:LEU:HD22	1:A:137:LEU:N	2.23	0.53
1:A:90:SER:HA	1:A:369:ILE:HG13	1.90	0.53
1:B:137:LEU:HD21	1:B:195:HIS:HB3	1.90	0.53
1:A:272:THR:HG22	1:A:311:LEU:CD1	2.29	0.53
1:A:52:SER:HB2	1:A:53:PRO:HD2	1.90	0.53
1:A:60:LEU:CD1	1:A:274:ALA:HB3	2.38	0.53
1:B:52:SER:HB2	1:B:53:PRO:HD2	1.89	0.53
1:A:71:THR:HB	1:A:73:GLU:OE2	2.09	0.52
1:A:149:ARG:HG2	1:A:149:ARG:NH1	2.19	0.52
1:A:347:LYS:HB3	1:A:347:LYS:HZ3	1.75	0.52
1:A:109:GLN:HG3	1:A:112:ARG:NH2	2.24	0.52
1:B:87:ASN:HB3	1:B:92:CYS:HB3	1.91	0.52
1:A:133:PRO:O	1:A:135:LEU:N	2.41	0.51
1:A:218:TYR:CE1	1:A:256:ASN:HB3	2.45	0.51
1:B:129:HIS:CE1	1:B:170:THR:OG1	2.59	0.51
1:A:114:MET:CE	1:A:118:LEU:HD13	2.37	0.51
1:A:338:TYR:HB2	1:A:339:PRO:HD3	1.93	0.51
1:A:129:HIS:CE1	1:A:170:THR:OG1	2.62	0.51
1:A:156:VAL:HG21	1:A:180:VAL:HG22	1.92	0.51
1:A:198:ALA:HA	1:A:202:GLU:OE2	2.10	0.51
1:B:218:TYR:CE1	1:B:256:ASN:HB3	2.45	0.51
1:A:287:PHE:HB3	1:A:292:ILE:CD1	2.39	0.51
1:A:129:HIS:ND1	1:A:168:LYS:HG3	2.26	0.51
1:B:133:PRO:O	1:B:135:LEU:N	2.41	0.50
1:A:75:VAL:HG11	1:A:106:LEU:CD2	2.41	0.50
1:A:295:ARG:O	1:A:296:ASN:HB2	2.11	0.50
1:B:114:MET:CE	1:B:118:LEU:HD13	2.36	0.50
1:B:75:VAL:HG11	1:B:106:LEU:CD2	2.41	0.50
1:B:158:PHE:CE1	1:B:162:ASN:ND2	2.79	0.50
1:A:248:ASP:HB3	1:A:311:LEU:HG	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:336:THR:O	1:B:339:PRO:HD2	2.12	0.50
1:B:271:VAL:HA	1:B:313:MET:HE1	1.94	0.49
1:A:287:PHE:O	1:A:288:ALA:HB3	2.13	0.49
1:A:294:TYR:CE1	1:A:297:GLY:HA2	2.47	0.49
1:A:41:ILE:HD12	1:A:41:ILE:O	2.12	0.49
1:B:1:MET:HE2	1:B:41:ILE:CG2	2.42	0.49
1:B:200:LEU:HD12	1:B:234:ALA:HB2	1.94	0.49
1:A:146:ASN:HA	1:A:177:PRO:HA	1.95	0.49
1:B:75:VAL:HG11	1:B:106:LEU:HD21	1.95	0.49
1:A:277:PRO:HD3	1:A:285:PHE:CE1	2.48	0.49
1:A:237:ASP:HA	1:B:262:ARG:HH22	1.77	0.49
1:B:198:ALA:HA	1:B:202:GLU:OE2	2.12	0.49
1:A:171:LEU:HD12	1:A:171:LEU:C	2.38	0.49
1:A:271:VAL:HA	1:A:313:MET:HE1	1.94	0.49
1:A:158:PHE:CE1	1:A:162:ASN:ND2	2.81	0.48
1:A:104:ASP:O	1:A:108:LYS:HG3	2.14	0.48
1:B:346:GLU:O	1:B:356:LYS:HE3	2.14	0.48
1:B:104:ASP:OD1	1:B:155:LEU:HD23	2.14	0.48
1:B:201:LYS:N	1:B:201:LYS:CE	2.71	0.48
1:A:149:ARG:HB3	1:A:151:PRO:HD2	1.95	0.48
1:B:123:ASN:CG	1:B:357:VAL:HG21	2.40	0.47
1:B:156:VAL:HG21	1:B:180:VAL:HG22	1.94	0.47
1:B:338:TYR:HB2	1:B:339:PRO:HD3	1.96	0.47
1:B:336:THR:C	1:B:339:PRO:HD2	2.40	0.47
1:B:171:LEU:C	1:B:171:LEU:HD12	2.38	0.47
1:A:225:THR:OG1	1:A:228:GLU:HG2	2.15	0.47
1:B:270:LEU:HD22	1:B:271:VAL:N	2.30	0.47
1:A:317:VAL:HG13	1:A:331:VAL:HG11	1.97	0.47
1:A:8:ARG:NH2	1:A:17:ASP:OD1	2.48	0.47
1:A:10:PHE:CE1	1:A:332:LEU:HB3	2.49	0.47
1:A:362:ALA:HB3	1:A:371:LYS:HG3	1.96	0.47
1:A:28:ILE:O	1:A:354:ALA:HA	2.15	0.46
1:A:135:LEU:HD22	1:A:195:HIS:HB2	1.96	0.46
1:A:98:THR:OG1	1:A:129:HIS:HD2	1.97	0.46
1:A:9:ILE:HD12	1:A:21:VAL:HG23	1.97	0.46
1:B:104:ASP:O	1:B:108:LYS:HG3	2.16	0.46
1:A:317:VAL:HG13	1:A:331:VAL:CG1	2.46	0.46
1:B:227:ARG:HG2	1:B:227:ARG:HH21	1.80	0.46
1:A:149:ARG:NH1	1:A:149:ARG:CG	2.79	0.46
1:B:75:VAL:HG22	1:B:75:VAL:O	2.15	0.46
1:B:347:LYS:HB3	1:B:347:LYS:HZ3	1.81	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:67:GLN:HE22	1:A:287:PHE:HB2	1.82	0.45
1:A:223:TYR:CD2	1:B:225:THR:HG22	2.50	0.45
1:B:50:ILE:N	1:B:50:ILE:HD12	2.31	0.45
1:B:107:MET:HE1	1:B:155:LEU:HD21	1.97	0.45
1:A:22:VAL:HG22	1:A:29:LYS:HB3	1.97	0.45
1:A:217:LEU:O	1:A:218:TYR:HB2	2.16	0.45
1:A:292:ILE:HG12	1:A:307:SER:OG	2.16	0.45
1:B:135:LEU:O	1:B:175:MET:HG3	2.17	0.45
1:A:9:ILE:HD12	1:A:21:VAL:CG2	2.47	0.45
1:A:75:VAL:HG11	1:A:106:LEU:HD21	1.98	0.45
1:B:8:ARG:NH2	1:B:17:ASP:OD1	2.50	0.45
1:B:98:THR:OG1	1:B:129:HIS:HD2	1.99	0.45
1:A:70:ASP:O	1:A:71:THR:C	2.58	0.45
1:A:246:ILE:CD1	1:A:273:ASP:HB2	2.47	0.45
1:A:62:GLY:HA2	1:A:83:MET:CE	2.34	0.45
1:A:75:VAL:O	1:A:75:VAL:HG22	2.16	0.45
1:B:11:THR:HA	1:B:352:LEU:HB2	1.98	0.45
1:B:362:ALA:HB3	1:B:371:LYS:HG3	1.97	0.45
1:B:41:ILE:HD12	1:B:41:ILE:O	2.16	0.45
1:B:217:LEU:O	1:B:218:TYR:HB2	2.16	0.45
1:A:114:MET:HE2	1:A:118:LEU:CD1	2.40	0.45
1:A:129:HIS:ND1	1:A:168:LYS:HE3	2.32	0.45
1:B:57:ASP:CG	1:B:313:MET:HE3	2.42	0.45
1:A:50:ILE:N	1:A:50:ILE:HD12	2.31	0.44
1:A:227:ARG:HG2	1:A:227:ARG:HH21	1.82	0.44
1:B:9:ILE:HD12	1:B:21:VAL:HG23	1.99	0.44
1:B:225:THR:OG1	1:B:228:GLU:HG2	2.17	0.44
1:B:22:VAL:HG22	1:B:29:LYS:HB3	1.98	0.44
1:B:129:HIS:ND1	1:B:168:LYS:HE3	2.32	0.44
1:B:133:PRO:HG2	1:B:134:TRP:H	1.81	0.44
1:A:215:THR:HG22	1:A:216:HIS:CD2	2.53	0.44
1:A:346:GLU:O	1:A:356:LYS:HE3	2.17	0.44
1:B:114:MET:HE2	1:B:118:LEU:CD1	2.40	0.44
1:B:129:HIS:HA	1:B:168:LYS:O	2.18	0.44
1:A:90:SER:O	1:A:314:ILE:HG13	2.16	0.44
1:A:129:HIS:HA	1:A:168:LYS:O	2.18	0.44
1:A:290:LYS:HE2	1:A:301:ASP:OD1	2.17	0.44
1:A:148:VAL:HG23	1:A:152:ASP:HB2	2.00	0.44
1:A:171:LEU:HD12	1:A:171:LEU:O	2.18	0.44
1:B:215:THR:HG22	1:B:216:HIS:CD2	2.53	0.44
1:B:330:GLU:HA	1:B:333:ARG:HH11	1.83	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:201:LYS:N	1:A:201:LYS:CE	2.70	0.43
1:B:10:PHE:CE1	1:B:12:GLY:HA2	2.53	0.43
1:B:347:LYS:NZ	1:B:347:LYS:CB	2.82	0.43
1:B:10:PHE:CE1	1:B:332:LEU:HB3	2.54	0.43
1:B:129:HIS:ND1	1:B:168:LYS:HG3	2.34	0.43
1:A:133:PRO:HG2	1:A:134:TRP:H	1.83	0.43
1:A:270:LEU:HD22	1:A:271:VAL:N	2.33	0.43
1:B:37:LEU:HD22	1:B:41:ILE:CD1	2.49	0.43
1:B:97:PRO:HD2	1:B:127:GLY:O	2.19	0.43
1:B:332:LEU:HD12	1:B:332:LEU:HA	1.87	0.43
1:B:221:MET:HE1	1:B:235:ILE:CD1	2.49	0.43
1:B:118:LEU:HD12	1:B:118:LEU:HA	1.85	0.43
1:B:8:ARG:HG2	1:B:8:ARG:HH21	1.84	0.42
1:B:134:TRP:HB3	1:B:176:VAL:CG2	2.48	0.42
1:A:136:ASN:C	1:A:136:ASN:HD22	2.26	0.42
1:B:371:LYS:HE3	1:B:378:GLU:OE1	2.19	0.42
1:A:330:GLU:HA	1:A:333:ARG:HH11	1.84	0.42
1:A:200:LEU:HD12	1:A:234:ALA:HB2	2.01	0.42
1:A:106:LEU:HA	1:A:106:LEU:HD23	1.86	0.42
1:A:336:THR:C	1:A:339:PRO:HD2	2.45	0.42
1:B:106:LEU:HD23	1:B:106:LEU:HA	1.86	0.42
1:B:171:LEU:HD12	1:B:171:LEU:O	2.19	0.42
1:B:133:PRO:HG2	1:B:134:TRP:N	2.35	0.42
1:B:317:VAL:HG13	1:B:331:VAL:CG1	2.50	0.42
1:B:347:LYS:O	1:B:347:LYS:HG3	2.19	0.42
1:B:10:PHE:O	1:B:352:LEU:HD22	2.19	0.41
1:A:346:GLU:C	1:A:348:ARG:H	2.28	0.41
1:B:9:ILE:HD12	1:B:21:VAL:CG2	2.49	0.41
1:B:88:GLU:OE2	1:B:379:VAL:HG21	2.20	0.41
1:B:215:THR:O	1:B:216:HIS:C	2.62	0.41
1:A:9:ILE:HD11	1:A:16:LEU:HD12	2.01	0.41
1:A:51:LEU:HD13	1:A:352:LEU:HD21	2.02	0.41
1:B:9:ILE:HD11	1:B:16:LEU:HD12	2.02	0.41
1:A:215:THR:O	1:A:216:HIS:C	2.63	0.41
1:B:346:GLU:C	1:B:348:ARG:H	2.29	0.41
1:A:347:LYS:NZ	1:A:347:LYS:CB	2.84	0.41
1:B:135:LEU:HD22	1:B:195:HIS:HB2	2.02	0.41
1:B:317:VAL:HG13	1:B:331:VAL:HG11	2.03	0.41
1:B:56:ILE:HD12	1:B:340:ALA:HB2	2.03	0.41
1:B:219:ASN:OD1	1:B:251:HIS:HB3	2.21	0.41
1:A:97:PRO:HD2	1:A:127:GLY:O	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:347:LYS:O	1:A:347:LYS:HG3	2.20	0.41
1:B:366:ASP:OD2	1:B:368:LYS:HE2	2.20	0.40
1:A:295:ARG:O	1:A:296:ASN:CB	2.69	0.40
1:A:8:ARG:HH21	1:A:8:ARG:HG2	1.87	0.40
1:A:37:LEU:HD22	1:A:41:ILE:CD1	2.51	0.40
1:B:1:MET:HE2	1:B:41:ILE:HG21	2.02	0.40
1:A:95:TYR:O	1:A:96:LEU:HD12	2.21	0.40
1:A:371:LYS:HE3	1:A:378:GLU:OE1	2.21	0.40
1:B:100:ILE:O	1:B:101:THR:HB	2.21	0.40
1:B:246:ILE:CD1	1:B:273:ASP:HB2	2.51	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	372/382 (97%)	341 (92%)	26 (7%)	5 (1%)	9	21
1	B	323/382 (85%)	299 (93%)	20 (6%)	4 (1%)	10	23
All	All	695/764 (91%)	640 (92%)	46 (7%)	9 (1%)	9	21

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	134	TRP
1	A	278	ALA
1	B	134	TRP
1	A	216	HIS
1	B	216	HIS
1	A	273	ASP
1	A	296	ASN
1	B	273	ASP

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Mol	Chain	Res	Type
1	B	75	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	300/305 (98%)	272 (91%)	28 (9%)	8	18
1	B	263/305 (86%)	237 (90%)	26 (10%)	7	16
All	All	563/610 (92%)	509 (90%)	54 (10%)	8	17

All (54) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	ILE
1	A	22	VAL
1	A	31	VAL
1	A	44	ARG
1	A	80	LEU
1	A	99	LEU
1	A	106	LEU
1	A	118	LEU
1	A	149	ARG
1	A	150	LYS
1	A	155	LEU
1	A	174	GLU
1	A	179	GLU
1	A	184	LEU
1	A	200	LEU
1	A	201	LYS
1	A	228	GLU
1	A	240	ASP
1	A	250	LEU
1	A	270	LEU
1	A	284	GLN
1	A	311	LEU

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Mol	Chain	Res	Type
1	A	314	ILE
1	A	329	ASP
1	A	332	LEU
1	A	341	ARG
1	A	347	LYS
1	A	352	LEU
1	B	9	ILE
1	B	22	VAL
1	B	31	VAL
1	B	44	ARG
1	B	75	VAL
1	B	80	LEU
1	B	99	LEU
1	B	106	LEU
1	B	118	LEU
1	B	155	LEU
1	B	174	GLU
1	B	179	GLU
1	B	184	LEU
1	B	200	LEU
1	B	201	LYS
1	B	228	GLU
1	B	240	ASP
1	B	250	LEU
1	B	270	LEU
1	B	311	LEU
1	B	314	ILE
1	B	329	ASP
1	B	332	LEU
1	B	341	ARG
1	B	347	LYS
1	B	352	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	6	GLN
1	A	67	GLN
1	A	69	ASN
1	A	84	GLN
1	A	109	GLN
1	A	129	HIS

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Mol	Chain	Res	Type
1	A	136	ASN
1	A	284	GLN
1	A	377	ASN
1	B	6	GLN
1	B	67	GLN
1	B	84	GLN
1	B	109	GLN
1	B	129	HIS
1	B	136	ASN
1	B	195	HIS
1	B	377	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

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, 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	376/382 (98%)	-0.18	4 (1%) 78 74	24, 41, 77, 107	0
1	B	331/382 (86%)	-0.01	7 (2%) 63 58	29, 52, 81, 112	0
All	All	707/764 (92%)	-0.10	11 (1%) 70 66	24, 47, 78, 112	0

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	276	ALA	3.9
1	B	132	GLY	3.6
1	A	307	SER	2.6
1	A	138	VAL	2.5
1	B	350	GLY	2.5
1	B	137	LEU	2.3
1	B	161	GLU	2.2
1	A	145	PRO	2.2
1	B	223	TYR	2.2
1	A	179	GLU	2.1
1	B	153	ALA	2.1

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 oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.