Full wwPDB X-ray Structure Validation Report (i)

Jun 12, 2024 – 02:54 AM EDT

PDB ID	:	1YIF
Title	:	CRYSTAL STRUCTURE OF beta-1,4-xylosidase FROM BACILLUS SUBTILIS, NEW YORK STRUCTURAL GENOMICS CONSORTIUM
Authors	:	Patskovsky, Y.; Almo, S.C.; Burley, S.K.; New York SGX Research Center for Structural Genomics (NYSGXRC)
Deposited on	:	2005-01-11
Resolution	:	1.80 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 (i) 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 (1)) were used in the production of this report:

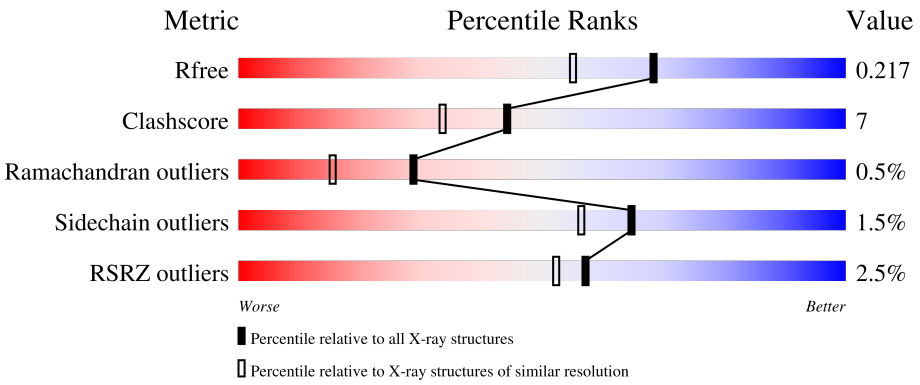
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.20.1
EDS	:	2.36.2
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36.2

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 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}	130704	5950 (1.80-1.80)
Clashscore	141614	6793 (1.80-1.80)
Ramachandran outliers	138981	6697 (1.80-1.80)
Sidechain outliers	138945	6696 (1.80-1.80)
RSRZ outliers	127900	5850 (1.80-1.80)

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 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 <=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	533	4% 87% 13% •
1	B	533	2% 86% 13% •
1	C	533	2% 84% 16% •
1	D	533	2% 84% 15% •

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 20084 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 BETA-1,4-XYLOSIDASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	532	4339	2785	728	815	11	0	0	0
1	B	532	4339	2785	728	815	11	0	0	0
1	C	532	4339	2785	728	815	11	0	0	0
1	D	532	4339	2785	728	815	11	0	0	0

- Molecule 2 is water.

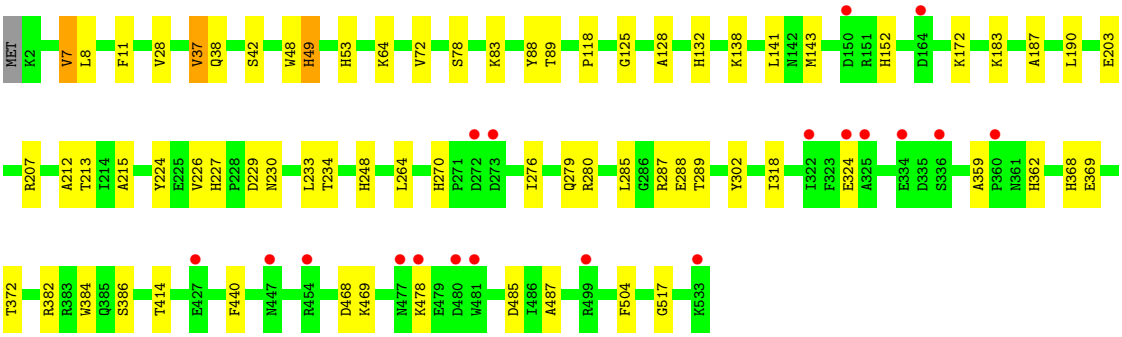
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
2	A	649	649	649	0	0
2	B	671	671	671	0	0
2	C	699	699	699	0	0
2	D	709	709	709	0	0

3 Residue-property plots (i)

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.

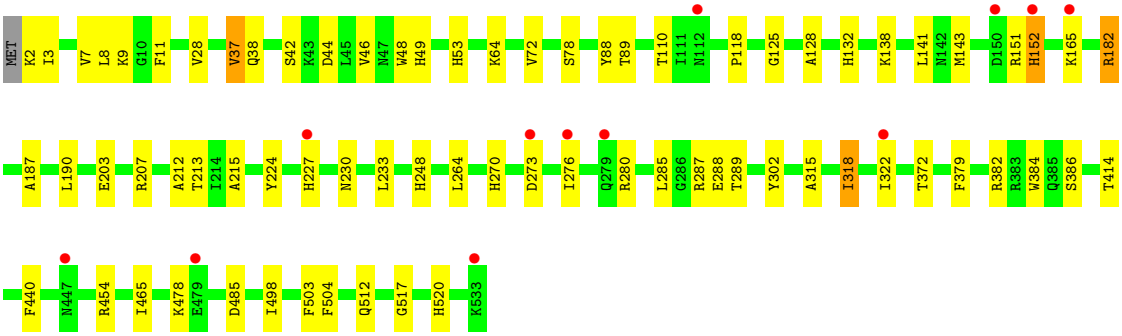
- Molecule 1: BETA-1,4-XYLOSIDASE

Chain A:



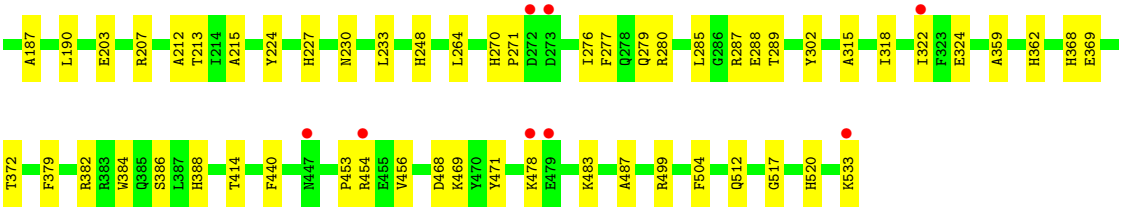
- Molecule 1: BETA-1,4-XYLOSIDASE

Chain B:



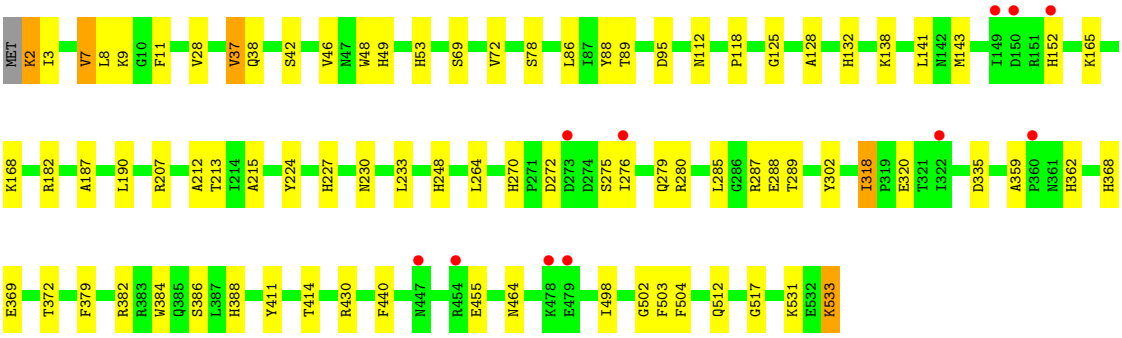
- Molecule 1: BETA-1,4-XYLOSIDASE

Chain C:



- Molecule 1: BETA-1,4-XYLOSIDASE

Chain D:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	104.23Å 104.61Å 114.60Å 90.00° 108.88° 90.00°	Depositor
Resolution (Å)	20.00 – 1.80 19.97 – 1.70	Depositor EDS
% Data completeness (in resolution range)	93.2 (20.00-1.80) 85.9 (19.97-1.70)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	0.06	Depositor
$< I/\sigma(I) >$ [1]	1.72 (at 1.70Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.205 , 0.218 0.205 , 0.217	Depositor DCC
R_{free} test set	6434 reflections (2.94%)	wwPDB-VP
Wilson B-factor (Å^2)	15.4	Xtriage
Anisotropy	0.178	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.36 , 57.3	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	20084	wwPDB-VP
Average B, all atoms (Å^2)	18.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.00% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality (i)

5.1 Standard geometry (i)

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	0.32	0/4475	0.60	0/6090
1	B	0.31	0/4475	0.59	0/6090
1	C	0.32	0/4475	0.59	0/6090
1	D	0.32	0/4475	0.60	1/6090 (0.0%)
All	All	0.32	0/17900	0.59	1/24360 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	D	272	ASP	CB-CG-OD2	6.75	124.37	118.30

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	4339	0	4106	53	0
1	B	4339	0	4106	59	0
1	C	4339	0	4106	60	0
1	D	4339	0	4106	57	0
2	A	649	0	0	23	0
2	B	671	0	0	14	0
2	C	699	0	0	24	0
2	D	709	0	0	21	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	20084	0	16424	224	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 7.

All (224) 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:B:152:HIS:CD2	1:B:152:HIS:H	1.77	0.96
1:B:152:HIS:H	1:B:152:HIS:HD2	1.18	0.90
1:B:64:LYS:HD2	2:B:1117:HOH:O	1.74	0.87
1:A:468:ASP:HB3	2:A:998:HOH:O	1.77	0.83
1:A:64:LYS:HD2	2:A:946:HOH:O	1.79	0.83
1:D:464:ASN:HB3	2:D:1193:HOH:O	1.81	0.79
1:B:322:ILE:HG13	2:B:1021:HOH:O	1.84	0.77
1:D:335:ASP:HB3	2:D:1221:HOH:O	1.85	0.75
1:C:471:TYR:HB3	2:C:1212:HOH:O	1.88	0.72
1:A:478:LYS:HA	2:A:949:HOH:O	1.89	0.72
1:B:165:LYS:HG2	2:B:938:HOH:O	1.90	0.70
1:C:285:LEU:HD22	1:C:384:TRP:O	1.91	0.70
1:B:110:THR:HB	2:B:783:HOH:O	1.90	0.70
1:D:517:GLY:HA3	2:D:675:HOH:O	1.92	0.69
1:A:226:VAL:HB	2:A:1159:HOH:O	1.93	0.69
1:B:285:LEU:HD22	1:B:384:TRP:O	1.93	0.68
1:C:270:HIS:HD2	2:C:672:HOH:O	1.75	0.68
1:B:9:LYS:HE2	2:B:1080:HOH:O	1.93	0.68
1:C:2:LYS:HE2	1:C:315:ALA:H	1.59	0.67
1:A:141:LEU:HD11	1:A:190:LEU:HB2	1.76	0.67
1:B:141:LEU:HD11	1:B:190:LEU:HB2	1.76	0.67
1:A:270:HIS:HD2	2:A:762:HOH:O	1.75	0.67
1:B:318:ILE:HG21	2:B:1186:HOH:O	1.95	0.67
1:C:141:LEU:HD11	1:C:190:LEU:HB2	1.77	0.67
1:D:112:ASN:HB2	2:D:943:HOH:O	1.95	0.66
1:B:2:LYS:HE2	1:B:315:ALA:H	1.61	0.66
1:D:287:ARG:NH2	1:D:504:PHE:HB3	2.10	0.66
1:A:285:LEU:HD22	1:A:384:TRP:O	1.95	0.65
1:C:287:ARG:NH2	1:C:504:PHE:HB3	2.12	0.65
1:D:141:LEU:HD11	1:D:190:LEU:HB2	1.79	0.64
1:B:287:ARG:NH2	1:B:504:PHE:HB3	2.13	0.64
1:B:152:HIS:CD2	1:B:152:HIS:N	2.56	0.63
1:A:287:ARG:NH2	1:A:504:PHE:HB3	2.13	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:276:ILE:O	1:A:280:ARG:HD2	2.00	0.62
1:C:533:LYS:HG2	2:C:1211:HOH:O	1.98	0.61
1:C:454:ARG:HG3	2:C:992:HOH:O	1.99	0.61
1:D:270:HIS:HD2	2:D:615:HOH:O	1.84	0.61
1:A:83:LYS:HE2	2:A:939:HOH:O	2.01	0.60
1:C:276:ILE:O	1:C:280:ARG:HD2	2.01	0.60
1:C:324:GLU:HG2	2:C:1001:HOH:O	2.02	0.60
1:B:44:ASP:HB2	2:B:963:HOH:O	2.02	0.60
1:C:382:ARG:HD2	2:C:577:HOH:O	2.01	0.60
1:C:112:ASN:HB3	2:C:883:HOH:O	2.02	0.59
1:C:64:LYS:HD2	2:C:818:HOH:O	2.02	0.59
1:C:322:ILE:HG13	2:C:945:HOH:O	2.03	0.59
1:C:388:HIS:HD2	2:C:994:HOH:O	1.85	0.59
1:D:37:VAL:HG11	1:D:86:LEU:HD21	1.86	0.58
1:C:483:LYS:HG3	2:C:1212:HOH:O	2.03	0.58
1:D:285:LEU:HD22	1:D:384:TRP:O	2.03	0.57
1:D:165:LYS:HG3	2:D:891:HOH:O	2.04	0.57
1:D:276:ILE:O	1:D:280:ARG:HD2	2.05	0.57
1:B:270:HIS:HD2	2:B:557:HOH:O	1.88	0.57
1:B:227:HIS:CE1	1:B:302:TYR:CZ	2.93	0.56
1:D:165:LYS:HG2	2:D:823:HOH:O	2.05	0.56
1:D:455:GLU:H	1:D:455:GLU:CD	2.08	0.56
1:D:78:SER:OG	1:D:132:HIS:HE1	1.89	0.56
1:A:517:GLY:HA3	2:A:1135:HOH:O	2.04	0.56
1:D:531:LYS:HE2	2:D:922:HOH:O	2.06	0.56
1:B:182:ARG:HD3	1:C:277:PHE:CD2	2.42	0.55
1:B:273:ASP:O	1:C:152:HIS:CE1	2.60	0.54
1:C:517:GLY:HA3	2:C:1014:HOH:O	2.08	0.54
1:B:276:ILE:O	1:B:280:ARG:HD2	2.08	0.54
1:B:187:ALA:HB2	1:B:248:HIS:CE1	2.43	0.53
1:A:226:VAL:HG22	2:A:660:HOH:O	2.08	0.53
1:C:78:SER:OG	1:C:132:HIS:HE1	1.91	0.53
1:A:152:HIS:HB3	2:A:721:HOH:O	2.09	0.53
1:C:499:ARG:HG3	2:C:942:HOH:O	2.09	0.53
1:D:9:LYS:HE2	2:D:899:HOH:O	2.09	0.52
1:A:230:ASN:CB	2:A:1159:HOH:O	2.57	0.52
1:C:227:HIS:HE1	1:C:302:TYR:CE1	2.27	0.52
1:C:280:ARG:HD3	2:C:762:HOH:O	2.08	0.52
1:D:187:ALA:HB2	1:D:248:HIS:CE1	2.45	0.52
1:B:78:SER:OG	1:B:132:HIS:HE1	1.93	0.52
1:B:212:ALA:HB3	1:B:233:LEU:HB3	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:227:HIS:CE1	1:B:302:TYR:CE1	2.98	0.51
1:A:212:ALA:HB3	1:A:233:LEU:HB3	1.92	0.51
1:B:382:ARG:HD2	2:B:612:HOH:O	2.10	0.51
1:A:213:THR:CG2	1:A:230:ASN:HD21	2.22	0.51
1:C:8:LEU:HB2	1:C:289:THR:HB	1.91	0.51
1:A:229:ASP:HB2	2:A:1001:HOH:O	2.10	0.51
1:D:382:ARG:HD2	2:D:634:HOH:O	2.09	0.51
1:A:172:LYS:HD3	2:A:1170:HOH:O	2.10	0.51
1:C:212:ALA:HB3	1:C:233:LEU:HB3	1.93	0.51
1:C:213:THR:CG2	1:C:230:ASN:HD21	2.24	0.51
1:D:212:ALA:HB3	1:D:233:LEU:HB3	1.93	0.51
1:A:324:GLU:HB2	2:A:1040:HOH:O	2.10	0.51
1:D:213:THR:CG2	1:D:230:ASN:HD21	2.23	0.51
1:A:8:LEU:HB2	1:A:289:THR:HB	1.92	0.50
1:B:227:HIS:HE1	1:B:302:TYR:CE1	2.29	0.50
1:B:125:GLY:HA3	1:B:143:MET:O	2.10	0.50
1:A:280:ARG:HD3	2:A:763:HOH:O	2.12	0.50
1:C:119:ILE:HG13	2:C:937:HOH:O	2.12	0.50
1:A:187:ALA:HB2	1:A:248:HIS:CE1	2.47	0.50
1:A:78:SER:OG	1:A:132:HIS:HE1	1.94	0.49
1:B:8:LEU:HB2	1:B:289:THR:HB	1.94	0.49
1:C:44:ASP:HB2	2:C:946:HOH:O	2.11	0.49
1:A:207:ARG:HB2	2:A:591:HOH:O	2.12	0.49
1:A:132:HIS:HD2	1:A:138:LYS:NZ	2.11	0.49
1:B:213:THR:CG2	1:B:230:ASN:HD21	2.25	0.49
1:D:168:LYS:HE3	2:D:1222:HOH:O	2.11	0.49
1:D:8:LEU:HB2	1:D:289:THR:HB	1.95	0.49
1:D:95:ASP:HB2	2:D:996:HOH:O	2.13	0.48
1:B:38:GLN:HG2	1:B:53:HIS:NE2	2.29	0.48
1:D:279:GLN:HB3	2:D:1237:HOH:O	2.13	0.48
1:A:215:ALA:HB1	1:A:224:TYR:HB3	1.96	0.48
1:C:2:LYS:HE2	1:C:315:ALA:N	2.27	0.48
1:D:320:GLU:HG3	2:D:884:HOH:O	2.14	0.48
1:B:88:TYR:OH	1:B:118:PRO:HB3	2.14	0.48
1:A:152:HIS:HA	2:A:880:HOH:O	2.13	0.47
1:A:42:SER:HB2	1:A:48:TRP:CD2	2.50	0.47
1:D:533:LYS:HB3	1:D:533:LYS:NZ	2.29	0.47
1:A:37:VAL:O	1:A:53:HIS:HA	2.14	0.47
1:A:382:ARG:HD2	2:A:699:HOH:O	2.13	0.47
1:D:132:HIS:HD2	1:D:138:LYS:NZ	2.12	0.47
1:C:42:SER:HB2	1:C:48:TRP:CD2	2.50	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:279:GLN:HB3	2:C:619:HOH:O	2.13	0.47
1:B:182:ARG:HD3	1:C:277:PHE:CE2	2.49	0.47
1:C:128:ALA:HA	1:C:141:LEU:O	2.15	0.47
1:D:227:HIS:HE1	1:D:302:TYR:CE1	2.33	0.47
1:C:187:ALA:HB2	1:C:248:HIS:CE1	2.49	0.47
1:D:125:GLY:HA3	1:D:143:MET:O	2.14	0.47
1:C:125:GLY:HA3	1:C:143:MET:O	2.15	0.47
1:B:517:GLY:HA3	2:B:556:HOH:O	2.14	0.46
1:B:3:ILE:HG12	1:B:46:VAL:HG22	1.96	0.46
1:B:38:GLN:HG2	1:B:53:HIS:HE2	1.80	0.46
1:B:454:ARG:O	1:B:454:ARG:HG3	2.15	0.46
1:B:478:LYS:O	1:B:478:LYS:HG2	2.14	0.46
1:D:42:SER:HB2	1:D:48:TRP:CD2	2.51	0.46
1:D:388:HIS:HD2	2:D:1153:HOH:O	1.96	0.46
1:D:7:VAL:HG12	1:D:8:LEU:HG	1.98	0.46
1:B:273:ASP:O	1:C:152:HIS:HE1	1.98	0.46
1:C:28:VAL:CG2	1:C:38:GLN:HB2	2.46	0.46
1:C:37:VAL:O	1:C:53:HIS:HA	2.16	0.46
1:D:88:TYR:OH	1:D:118:PRO:HB3	2.16	0.46
1:B:2:LYS:HE2	1:B:315:ALA:N	2.30	0.46
1:D:207:ARG:HB2	2:D:586:HOH:O	2.16	0.45
1:A:203:GLU:OE1	1:A:213:THR:HG21	2.16	0.45
1:A:226:VAL:HG23	2:A:1021:HOH:O	2.16	0.45
1:A:227:HIS:CE1	1:A:302:TYR:CE1	3.03	0.45
1:C:207:ARG:HB2	2:C:665:HOH:O	2.17	0.45
1:A:227:HIS:HE1	1:A:302:TYR:CE1	2.34	0.45
1:D:128:ALA:HA	1:D:141:LEU:O	2.17	0.45
1:A:38:GLN:HG2	1:A:53:HIS:HE2	1.82	0.45
1:B:215:ALA:HB1	1:B:224:TYR:HB3	1.99	0.45
1:C:533:LYS:HE3	2:C:722:HOH:O	2.17	0.45
1:A:28:VAL:CG2	1:A:38:GLN:HB2	2.47	0.45
1:A:88:TYR:OH	1:A:118:PRO:HB3	2.17	0.45
1:C:264:LEU:HA	1:C:288:GLU:O	2.17	0.45
1:A:125:GLY:HA3	1:A:143:MET:O	2.17	0.44
1:C:7:VAL:HG12	1:C:8:LEU:HG	2.00	0.44
1:C:203:GLU:OE1	1:C:213:THR:HG21	2.18	0.44
1:C:215:ALA:HB1	1:C:224:TYR:HB3	1.99	0.44
1:A:38:GLN:HG2	1:A:53:HIS:NE2	2.32	0.44
1:A:324:GLU:CB	2:A:1040:HOH:O	2.66	0.44
1:D:38:GLN:HG2	1:D:53:HIS:NE2	2.32	0.44
1:D:318:ILE:HD11	2:D:837:HOH:O	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:379:PHE:CD1	1:D:512:GLN:HB2	2.52	0.44
1:C:469:LYS:HD3	1:C:487:ALA:HB1	2.00	0.44
1:D:215:ALA:HB1	1:D:224:TYR:HB3	2.00	0.44
1:A:7:VAL:HG12	1:A:8:LEU:HG	1.99	0.44
1:B:132:HIS:HD2	1:B:138:LYS:NZ	2.16	0.44
1:D:28:VAL:O	1:D:37:VAL:HG23	2.17	0.44
1:A:279:GLN:HB3	2:A:592:HOH:O	2.17	0.44
1:A:469:LYS:HD3	1:A:487:ALA:HB1	2.00	0.44
1:B:285:LEU:CD2	1:B:384:TRP:O	2.65	0.44
1:D:270:HIS:HE1	2:D:995:HOH:O	2.01	0.44
1:C:76:CYS:HB3	1:C:87:ILE:HB	2.00	0.43
1:A:230:ASN:CG	2:A:1159:HOH:O	2.57	0.43
1:C:132:HIS:HD2	1:C:138:LYS:NZ	2.16	0.43
1:D:498:ILE:HD11	1:D:503:PHE:HA	1.99	0.43
1:B:151:ARG:HG3	2:B:861:HOH:O	2.17	0.43
1:B:128:ALA:HA	1:B:141:LEU:O	2.18	0.43
1:B:227:HIS:HE1	1:B:302:TYR:CZ	2.34	0.43
1:B:379:PHE:CD1	1:B:512:GLN:HB2	2.54	0.43
1:A:128:ALA:HA	1:A:141:LEU:O	2.17	0.42
1:B:42:SER:HB2	1:B:48:TRP:CD2	2.54	0.42
1:C:112:ASN:CB	2:C:883:HOH:O	2.63	0.42
1:C:9:LYS:HE2	2:C:1054:HOH:O	2.18	0.42
1:A:359:ALA:O	1:A:362:HIS:HB2	2.19	0.42
1:D:69:SER:HB2	1:D:502:GLY:HA3	2.01	0.42
1:C:271:PRO:HB3	2:C:980:HOH:O	2.19	0.42
1:A:49:HIS:HA	2:A:620:HOH:O	2.19	0.42
1:D:3:ILE:HG12	1:D:46:VAL:HG22	2.02	0.42
1:D:72:VAL:HA	1:D:89:THR:O	2.19	0.42
1:D:411:TYR:HE2	2:D:901:HOH:O	2.03	0.42
1:B:454:ARG:HA	2:B:551:HOH:O	2.19	0.42
1:D:38:GLN:HG2	1:D:53:HIS:CD2	2.54	0.42
1:D:227:HIS:CE1	1:D:302:TYR:CE1	3.08	0.42
1:B:38:GLN:HG2	1:B:53:HIS:CD2	2.54	0.42
1:B:498:ILE:HD11	1:B:503:PHE:HA	2.02	0.42
1:C:359:ALA:O	1:C:362:HIS:HB2	2.20	0.42
1:D:372:THR:HA	1:D:414:THR:HB	2.02	0.42
1:C:72:VAL:HA	1:C:89:THR:O	2.20	0.41
1:B:203:GLU:OE1	1:B:213:THR:HG21	2.21	0.41
1:B:372:THR:HA	1:B:414:THR:HB	2.01	0.41
1:B:465:ILE:N	1:B:465:ILE:HD12	2.35	0.41
1:C:478:LYS:HG2	1:C:478:LYS:O	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:520:HIS:HE1	2:C:579:HOH:O	2.03	0.41
1:D:38:GLN:HG2	1:D:53:HIS:HE2	1.86	0.41
1:A:264:LEU:HA	1:A:288:GLU:O	2.21	0.41
1:C:468:ASP:HB3	2:C:980:HOH:O	2.20	0.41
1:D:359:ALA:O	1:D:362:HIS:HB2	2.21	0.41
1:A:372:THR:HA	1:A:414:THR:HB	2.01	0.41
1:C:372:THR:HA	1:C:414:THR:HB	2.02	0.41
1:B:37:VAL:O	1:B:53:HIS:HA	2.21	0.41
1:B:143:MET:C	1:B:143:MET:SD	2.99	0.41
1:B:520:HIS:HE1	2:B:604:HOH:O	2.02	0.41
1:C:88:TYR:OH	1:C:118:PRO:HB3	2.21	0.41
1:D:2:LYS:HA	2:D:1087:HOH:O	2.21	0.41
1:A:368:HIS:HB3	1:A:369:GLU:H	1.65	0.41
1:D:368:HIS:HB3	1:D:369:GLU:H	1.62	0.41
1:D:430:ARG:HB2	2:D:640:HOH:O	2.21	0.41
1:B:264:LEU:HA	1:B:288:GLU:O	2.20	0.41
1:C:368:HIS:HB3	1:C:369:GLU:H	1.66	0.41
1:B:28:VAL:CG2	1:B:38:GLN:HB2	2.51	0.41
1:A:183:LYS:HB2	1:D:275:SER:HB2	2.02	0.40
1:A:183:LYS:HB3	2:A:633:HOH:O	2.21	0.40
1:D:28:VAL:CG2	1:D:38:GLN:HB2	2.52	0.40
1:B:72:VAL:HA	1:B:89:THR:O	2.22	0.40
1:C:379:PHE:CD1	1:C:512:GLN:HB2	2.56	0.40
1:A:72:VAL:HA	1:A:89:THR:O	2.21	0.40
1:B:207:ARG:HB2	2:B:617:HOH:O	2.21	0.40
1:C:453:PRO:HB2	1:C:456:VAL:HG23	2.03	0.40
1:D:264:LEU:HA	1:D:288:GLU:O	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles (i)

5.3.1 Protein backbone (i)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	530/533 (99%)	498 (94%)	28 (5%)	4 (1%)	19	7
1	B	530/533 (99%)	500 (94%)	27 (5%)	3 (1%)	25	12
1	C	530/533 (99%)	500 (94%)	28 (5%)	2 (0%)	34	21
1	D	530/533 (99%)	498 (94%)	30 (6%)	2 (0%)	34	21
All	All	2120/2132 (99%)	1996 (94%)	113 (5%)	11 (0%)	29	15

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	485	ASP
1	B	485	ASP
1	A	11	PHE
1	B	11	PHE
1	B	37	VAL
1	C	11	PHE
1	D	11	PHE
1	D	37	VAL
1	A	37	VAL
1	C	37	VAL
1	A	234	THR

5.3.2 Protein sidechains (i)

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	466/467 (100%)	461 (99%)	5 (1%)	73	68
1	B	466/467 (100%)	459 (98%)	7 (2%)	65	56
1	C	466/467 (100%)	459 (98%)	7 (2%)	65	56
1	D	466/467 (100%)	457 (98%)	9 (2%)	57	46
All	All	1864/1868 (100%)	1836 (98%)	28 (2%)	65	56

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	VAL
1	A	49	HIS
1	A	318	ILE
1	A	386	SER
1	A	440	PHE
1	B	7	VAL
1	B	49	HIS
1	B	152	HIS
1	B	182	ARG
1	B	318	ILE
1	B	386	SER
1	B	440	PHE
1	C	7	VAL
1	C	18	CYS
1	C	49	HIS
1	C	112	ASN
1	C	318	ILE
1	C	386	SER
1	C	440	PHE
1	D	2	LYS
1	D	7	VAL
1	D	49	HIS
1	D	152	HIS
1	D	182	ARG
1	D	318	ILE
1	D	386	SER
1	D	440	PHE
1	D	533	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	132	HIS
1	A	270	HIS
1	A	361	ASN
1	A	424	HIS
1	B	132	HIS
1	B	152	HIS
1	B	270	HIS
1	B	361	ASN
1	B	388	HIS
1	B	448	ASN
1	C	132	HIS

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Mol	Chain	Res	Type
1	C	152	HIS
1	C	270	HIS
1	C	361	ASN
1	D	132	HIS
1	D	270	HIS
1	D	278	GLN
1	D	361	ASN
1	D	447	ASN
1	D	448	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 monosaccharides 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	532/533 (99%)	-0.09	19 (3%)	42	37	9, 15, 29, 45	0
1	B	532/533 (99%)	-0.08	12 (2%)	60	56	8, 15, 28, 43	0
1	C	532/533 (99%)	-0.12	12 (2%)	60	56	8, 14, 27, 42	0
1	D	532/533 (99%)	-0.15	11 (2%)	63	59	8, 14, 27, 40	0
All	All	2128/2132 (99%)	-0.11	54 (2%)	57	52	8, 14, 28, 45	0

All (54) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	322	ILE	6.3
1	C	478	LYS	6.3
1	D	150	ASP	5.4
1	C	150	ASP	5.2
1	B	273	ASP	4.9
1	A	478	LYS	4.7
1	D	152	HIS	4.6
1	D	273	ASP	4.5
1	B	276	ILE	4.2
1	C	272	ASP	4.2
1	B	152	HIS	3.8
1	B	150	ASP	3.5
1	C	165	LYS	3.5
1	A	150	ASP	3.3
1	A	427	GLU	3.2
1	A	447	ASN	3.2
1	C	447	ASN	3.2
1	A	477	ASN	3.1
1	D	447	ASN	3.1
1	A	480	ASP	2.9
1	A	325	ALA	2.9

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Mol	Chain	Res	Type	RSRZ
1	C	111	ILE	2.9
1	A	324	GLU	2.8
1	A	273	ASP	2.7
1	C	479	GLU	2.7
1	B	279	GLN	2.7
1	A	272	ASP	2.7
1	B	165	LYS	2.6
1	C	152	HIS	2.6
1	A	322	ILE	2.6
1	A	164	ASP	2.6
1	A	533	LYS	2.5
1	B	533	LYS	2.5
1	C	273	ASP	2.5
1	C	454	ARG	2.5
1	D	454	ARG	2.4
1	A	336	SER	2.4
1	D	360	PRO	2.4
1	C	322	ILE	2.4
1	A	499	ARG	2.3
1	D	276	ILE	2.3
1	A	334	GLU	2.2
1	B	479	GLU	2.2
1	B	447	ASN	2.2
1	A	481	TRP	2.2
1	A	454	ARG	2.2
1	D	478	LYS	2.1
1	D	149	ILE	2.1
1	C	533	LYS	2.1
1	D	479	GLU	2.1
1	D	322	ILE	2.1
1	A	360	PRO	2.1
1	B	227	HIS	2.0
1	B	112	ASN	2.0

6.2 Non-standard residues in protein, DNA, RNA chains (i)

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

6.3 Carbohydrates (i)

There are no monosaccharides 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.