Full wwPDB X-ray Structure Validation Report (i)

Dec 7, 2023 – 07:22 am GMT

PDB ID	:	2JF7
Title	:	Structure of Strictosidine Glucosidase
Authors	:	Barleben, L.; Panjikar, S.; Ruppert, M.; Koepke, J.; Stockigt, J.
Deposited on	:	2007-01-26
Resolution	:	2.48 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:

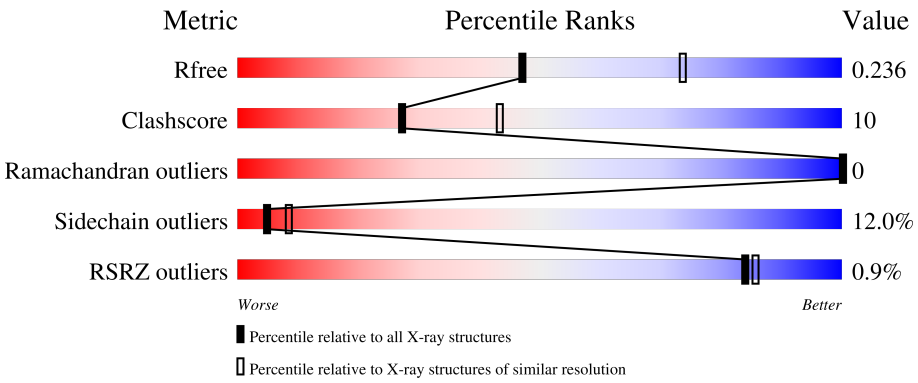
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

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

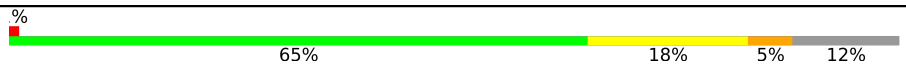
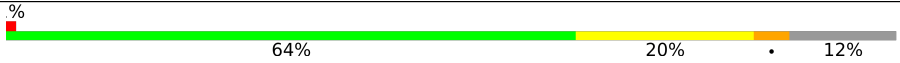
The reported resolution of this entry is 2.48 Å.

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	5857 (2.50-2.46)
Clashscore	141614	6594 (2.50-2.46)
Ramachandran outliers	138981	6469 (2.50-2.46)
Sidechain outliers	138945	6471 (2.50-2.46)
RSRZ outliers	127900	5738 (2.50-2.46)

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	532	
1	B	532	

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 8074 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 STRICTOSIDINE-O-BETA-D-GLUCOSIDASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	468	3796	2443	633	705	15	0	0	1
1	B	467	3788	2438	631	704	15	0	0	1

- Molecule 2 is water.

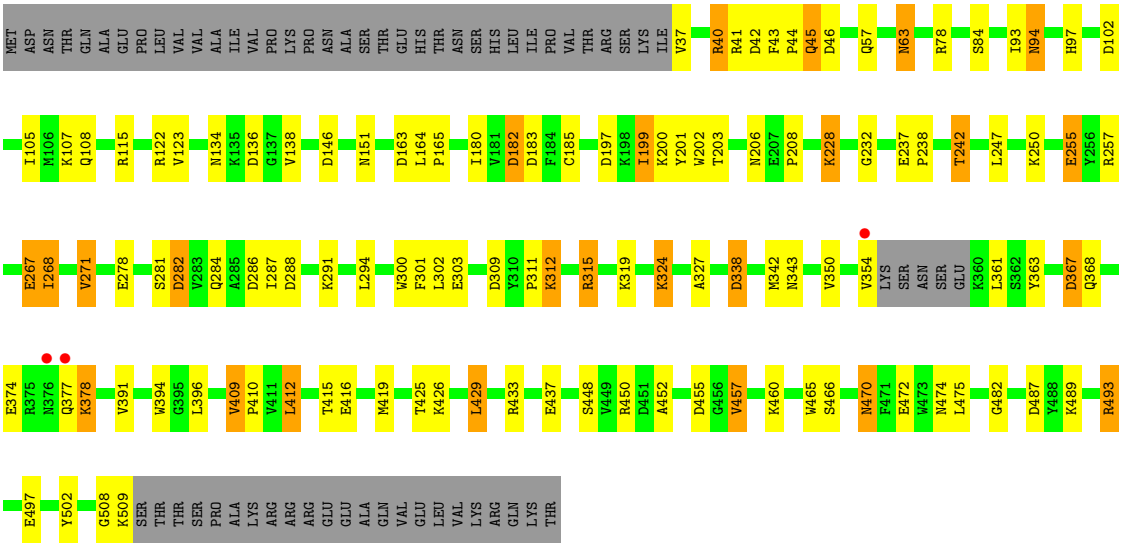
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
2	A	245	245	245	0	0
2	B	245	245	245	0	0

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.

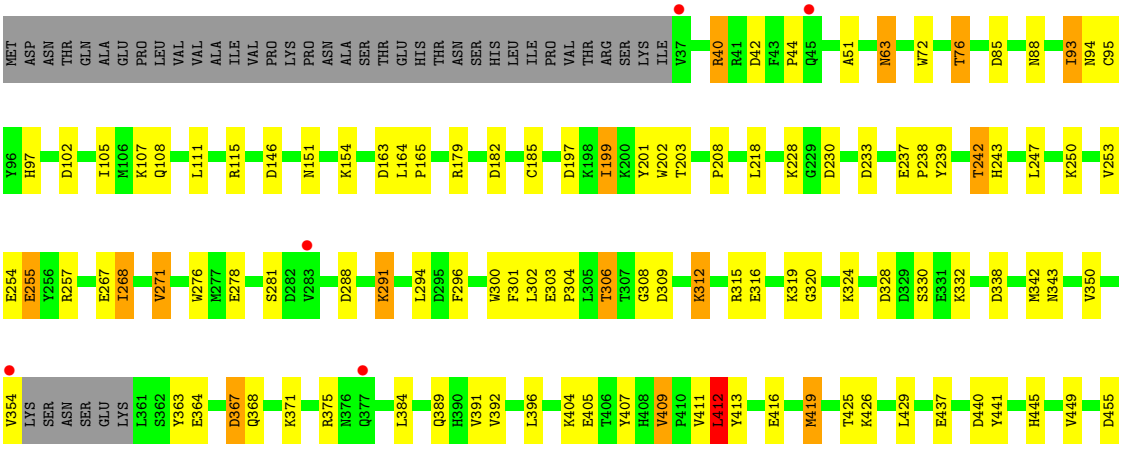
• Molecule 1: STRICTOSIDINE-O-BETA-D-GLUCOSIDASE

Chain A:



• Molecule 1: STRICTOSIDINE-O-BETA-D-GLUCOSIDASE

Chain B:



G456
V457
K460
F463
S466
F467
F468
D469
N470
N474
L475
G482
V486
K489
R493
E497
Y502
K509
SER
THR
THR
SER
PRO
ALA
LYS
ARG
ARG
ARG
GLU
GLU
ALA
GLN
VAL
GLU
LEU
VAL
LYS
ARG
GLN
LYS
THR

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 4 21 2	Depositor
Cell constants a, b, c, α, β, γ	157.63Å 157.63Å 103.59Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.48 19.98 – 2.48	Depositor EDS
% Data completeness (in resolution range)	100.0 (20.00-2.48) 100.0 (19.98-2.48)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	3.22 (at 2.47Å)	Xtriage
Refinement program	REFMAC 5.1.24	Depositor
R, R_{free}	0.185 , 0.239 0.184 , 0.236	Depositor DCC
R_{free} test set	1120 reflections (2.40%)	wwPDB-VP
Wilson B-factor (Å^2)	30.3	Xtriage
Anisotropy	0.094	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.33 , 28.3	EDS
L-test for twinning[2]	$< \|L\| > = 0.43$, $< L^2 > = 0.26$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	8074	wwPDB-VP
Average B, all atoms (Å^2)	22.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.61% 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	0.61	0/3908	0.86	18/5292 (0.3%)
1	B	0.62	0/3900	0.85	19/5283 (0.4%)
All	All	0.61	0/7808	0.85	37/10575 (0.3%)

There are no bond length outliers.

All (37) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	455	ASP	CB-CG-OD2	8.23	125.70	118.30
1	B	493	ARG	NE-CZ-NH2	-7.76	116.42	120.30
1	A	136	ASP	CB-CG-OD2	7.64	125.18	118.30
1	A	508	GLY	O-C-N	-7.61	110.52	122.70
1	B	197	ASP	CB-CG-OD2	7.43	124.99	118.30
1	B	163	ASP	CB-CG-OD2	7.03	124.62	118.30
1	B	182	ASP	CB-CG-OD2	6.99	124.59	118.30
1	B	42	ASP	CB-CG-OD2	6.93	124.54	118.30
1	B	102	ASP	CB-CG-OD2	6.89	124.50	118.30
1	B	328	ASP	CB-CG-OD2	6.70	124.33	118.30
1	A	338	ASP	CB-CG-OD2	6.67	124.30	118.30
1	A	182	ASP	CB-CG-OD2	6.59	124.23	118.30
1	A	163	ASP	CB-CG-OD2	6.55	124.19	118.30
1	B	309	ASP	CB-CG-OD2	6.54	124.19	118.30
1	A	487	ASP	CB-CG-OD2	6.51	124.16	118.30
1	B	233	ASP	CB-CG-OD2	6.47	124.13	118.30
1	A	309	ASP	CB-CG-OD2	6.36	124.03	118.30
1	A	46	ASP	CB-CG-OD2	6.16	123.85	118.30
1	A	197	ASP	CB-CG-OD2	6.10	123.79	118.30
1	A	455	ASP	CB-CG-OD2	5.90	123.61	118.30
1	A	42	ASP	CB-CG-OD2	5.88	123.59	118.30
1	B	469	ASP	CB-CG-OD2	5.82	123.54	118.30
1	A	286	ASP	CB-CG-OD2	5.77	123.49	118.30
1	B	288	ASP	CB-CG-OD2	5.66	123.39	118.30
1	B	367	ASP	CB-CG-OD1	5.48	123.24	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	440	ASP	CB-CG-OD2	5.42	123.18	118.30
1	A	102	ASP	CB-CG-OD2	5.39	123.15	118.30
1	A	288	ASP	CB-CG-OD2	5.27	123.04	118.30
1	B	146	ASP	CB-CG-OD2	5.27	123.04	118.30
1	B	230	ASP	CB-CG-OD2	5.25	123.02	118.30
1	B	493	ARG	NE-CZ-NH1	5.24	122.92	120.30
1	B	412	LEU	CA-CB-CG	5.22	127.30	115.30
1	A	282	ASP	CB-CG-OD2	5.20	122.98	118.30
1	A	183	ASP	CB-CG-OD2	5.18	122.97	118.30
1	A	146	ASP	CB-CG-OD2	5.18	122.96	118.30
1	B	85	ASP	CB-CG-OD2	5.04	122.84	118.30
1	A	367	ASP	CB-CG-OD2	5.02	122.82	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	3796	0	3600	72	0
1	B	3788	0	3591	77	0
2	A	245	0	0	24	2
2	B	245	0	0	14	1
All	All	8074	0	7191	149	2

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 10.

All (149) 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:123:VAL:HA	2:A:2072:HOH:O	1.50	1.12
1:B:306:THR:HG22	2:B:2175:HOH:O	1.58	1.02
1:A:115:ARG:NH1	2:A:2068:HOH:O	2.02	0.92
1:B:368:GLN:HG3	2:B:2181:HOH:O	1.71	0.88

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:425:THR:O	1:B:426:LYS:HB2	1.72	0.88
1:A:315:ARG:HG3	2:A:2168:HOH:O	1.75	0.84
1:B:93:ILE:CG1	1:B:486:VAL:HG11	2.08	0.83
1:A:271:VAL:HG11	2:A:2203:HOH:O	1.81	0.81
1:A:415:THR:HG21	2:A:2203:HOH:O	1.80	0.80
1:A:425:THR:O	1:A:426:LYS:HB2	1.81	0.79
1:A:122:ARG:O	2:A:2072:HOH:O	2.00	0.79
1:A:509:LYS:N	2:A:2245:HOH:O	2.15	0.78
1:A:105:ILE:HA	1:A:108:GLN:HE21	1.45	0.77
1:A:37:VAL:HA	2:A:2001:HOH:O	1.85	0.77
1:A:94:ASN:HD21	1:A:97:HIS:HD2	1.31	0.75
1:B:276:TRP:HE1	1:B:278:GLU:HG2	1.51	0.74
1:B:93:ILE:HG12	1:B:486:VAL:HG11	1.69	0.73
1:B:107:LYS:HE2	2:B:2065:HOH:O	1.90	0.72
1:A:367:ASP:OD1	1:A:367:ASP:O	2.07	0.71
1:A:415:THR:CB	2:A:2203:HOH:O	2.39	0.70
1:A:203:THR:HG21	1:A:271:VAL:HG13	1.76	0.68
2:A:2238:HOH:O	1:B:489:LYS:HD3	1.94	0.68
1:B:105:ILE:HA	1:B:108:GLN:HE21	1.58	0.67
1:B:76:THR:HG21	1:B:88:ASN:HB2	1.75	0.67
1:A:40:ARG:NH2	1:A:44:PRO:O	2.26	0.66
1:A:134:ASN:HB3	2:A:2072:HOH:O	1.93	0.66
1:B:276:TRP:NE1	1:B:278:GLU:HG2	2.10	0.66
1:B:409:VAL:HG21	1:B:412:LEU:HD13	1.77	0.65
1:B:306:THR:CG2	2:B:2175:HOH:O	2.28	0.65
1:B:107:LYS:HD2	1:B:151:ASN:HD22	1.62	0.65
1:A:267:GLU:HG2	1:A:338:ASP:HB3	1.77	0.64
1:B:497:GLU:HG3	2:B:2240:HOH:O	1.97	0.63
1:B:243:HIS:N	2:B:2144:HOH:O	2.32	0.63
1:A:242:THR:HG21	1:A:300:TRP:CZ2	2.35	0.61
1:B:201:TYR:HE1	2:B:2132:HOH:O	1.83	0.61
1:A:415:THR:CG2	2:A:2203:HOH:O	2.41	0.61
1:B:201:TYR:CE1	2:B:2132:HOH:O	2.50	0.60
1:A:429:LEU:HD22	1:A:433:ARG:HG3	1.83	0.59
1:A:105:ILE:HA	1:A:108:GLN:NE2	2.17	0.59
1:B:239:TYR:O	2:B:2144:HOH:O	2.17	0.59
1:A:93:ILE:O	1:A:493:ARG:NH2	2.33	0.58
1:A:242:THR:CG2	1:A:300:TRP:CZ2	2.87	0.58
1:B:76:THR:HG23	1:B:88:ASN:HB3	1.84	0.58
1:B:93:ILE:HG13	1:B:486:VAL:HG11	1.83	0.58
1:B:199:ILE:HG12	1:B:202:TRP:CE2	2.39	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:312:LYS:O	1:B:316:GLU:HG3	2.05	0.57
1:B:253:VAL:HG13	1:B:268:ILE:HD13	1.86	0.56
1:B:308:GLY:O	1:B:330:SER:OG	2.22	0.56
1:A:409:VAL:HG21	1:A:412:LEU:HD13	1.87	0.56
1:B:94:ASN:HD21	1:B:97:HIS:HD2	1.54	0.56
1:B:367:ASP:O	1:B:367:ASP:OD1	2.23	0.56
1:A:208:PRO:HB2	1:A:301:PHE:CE1	2.41	0.55
1:A:409:VAL:CG2	1:A:412:LEU:HD13	2.35	0.55
1:A:312:LYS:HD3	2:A:2166:HOH:O	2.06	0.55
1:B:208:PRO:HB2	1:B:301:PHE:CE1	2.41	0.55
1:A:232:GLY:HA3	2:A:2140:HOH:O	2.07	0.54
1:A:412:LEU:HD23	1:A:457:VAL:HG23	1.89	0.54
1:B:72:TRP:O	1:B:76:THR:HB	2.08	0.54
1:A:107:LYS:HD2	1:A:151:ASN:HD22	1.73	0.54
1:B:76:THR:CG2	1:B:88:ASN:HB2	2.39	0.53
1:A:303:GLU:HB3	1:A:311:PRO:HD3	1.91	0.53
1:B:242:THR:HG21	1:B:300:TRP:CZ2	2.43	0.53
1:A:63:ASN:HD22	1:A:63:ASN:H	1.57	0.53
1:B:242:THR:CG2	1:B:300:TRP:CZ2	2.92	0.52
1:A:203:THR:CG2	1:A:271:VAL:HG13	2.39	0.52
1:B:154:LYS:HE2	2:B:2096:HOH:O	2.10	0.52
1:A:327:ALA:HB3	2:A:2174:HOH:O	2.09	0.52
1:B:76:THR:CG2	1:B:88:ASN:CB	2.88	0.52
1:A:281:SER:HA	2:A:2157:HOH:O	2.09	0.51
1:A:312:LYS:NZ	2:A:2168:HOH:O	2.39	0.51
1:A:206:ASN:HA	1:A:271:VAL:HG22	1.93	0.51
1:A:94:ASN:HD21	1:A:97:HIS:CD2	2.19	0.51
1:A:303:GLU:OE2	1:A:303:GLU:HA	2.10	0.51
1:A:271:VAL:CG1	2:A:2203:HOH:O	2.51	0.50
1:A:94:ASN:ND2	1:A:97:HIS:HD2	2.05	0.50
1:B:303:GLU:HB3	1:B:304:PRO:HD3	1.94	0.50
1:A:242:THR:HG21	1:A:300:TRP:CH2	2.46	0.50
1:B:202:TRP:O	1:B:268:ILE:HA	2.12	0.49
1:A:278:GLU:HG3	1:A:394:TRP:HH2	1.77	0.49
1:B:115:ARG:HG3	1:B:463:PHE:CD2	2.48	0.49
1:B:242:THR:HG21	1:B:300:TRP:CH2	2.47	0.49
1:A:78:ARG:CG	1:A:78:ARG:HH21	2.25	0.48
1:A:378:LYS:NZ	2:A:2191:HOH:O	2.47	0.48
1:A:312:LYS:HA	1:A:312:LYS:HD2	1.61	0.48
1:B:253:VAL:HG13	1:B:268:ILE:CD1	2.43	0.47
1:A:470:ASN:ND2	1:A:472:GLU:OE1	2.46	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:468:PHE:HB2	2:B:2224:HOH:O	2.14	0.47
1:B:470:ASN:HD22	1:B:470:ASN:C	2.18	0.47
1:A:199:ILE:HD12	1:A:201:TYR:H	1.79	0.47
1:B:107:LYS:CE	2:B:2065:HOH:O	2.56	0.47
1:B:199:ILE:HG12	1:B:202:TRP:CZ2	2.50	0.47
1:B:237:GLU:N	1:B:238:PRO:CD	2.78	0.47
1:A:164:LEU:HD12	1:A:165:PRO:HD2	1.97	0.47
1:B:218:LEU:HD21	1:B:364:GLU:HG3	1.97	0.46
1:B:115:ARG:NH2	1:B:203:THR:OG1	2.48	0.46
1:B:267:GLU:HG2	1:B:338:ASP:HB3	1.98	0.46
1:B:412:LEU:HD23	1:B:457:VAL:HG23	1.98	0.46
1:B:409:VAL:CG2	1:B:412:LEU:HD13	2.43	0.46
1:A:367:ASP:OD1	1:A:367:ASP:C	2.55	0.46
1:A:433:ARG:NH1	1:A:497:GLU:OE2	2.48	0.46
1:B:445:HIS:O	1:B:449:VAL:HG23	2.16	0.45
1:B:154:LYS:CE	2:B:2096:HOH:O	2.63	0.45
1:B:384:LEU:HD12	1:B:389:GLN:HG2	1.98	0.45
1:A:363:TYR:CD2	1:A:363:TYR:C	2.89	0.45
1:A:228:LYS:HA	2:A:2137:HOH:O	2.17	0.45
1:B:63:ASN:HD22	1:B:63:ASN:H	1.65	0.45
1:B:250:LYS:O	1:B:254:GLU:HG3	2.17	0.45
1:A:378:LYS:HE3	1:A:378:LYS:HB3	1.56	0.45
1:B:343:ASN:CG	1:B:416:GLU:HB2	2.37	0.45
1:B:419:MET:SD	1:B:437:GLU:HG2	2.57	0.45
1:B:276:TRP:CD1	1:B:278:GLU:HG2	2.52	0.45
1:B:411:VAL:HG13	1:B:460:LYS:HG3	1.98	0.45
1:B:51:ALA:HB3	1:B:111:LEU:HD21	1.98	0.45
1:B:291:LYS:HZ2	1:B:291:LYS:HG2	1.44	0.44
1:B:199:ILE:O	1:B:199:ILE:HG13	2.14	0.44
1:B:76:THR:HG23	1:B:88:ASN:CB	2.46	0.44
1:A:45:GLN:HE21	1:A:45:GLN:HA	1.82	0.44
1:A:324:LYS:H	1:A:324:LYS:HG2	1.59	0.44
1:A:415:THR:HB	2:A:2203:HOH:O	2.09	0.44
1:A:182:ASP:OD1	2:A:2107:HOH:O	2.21	0.43
1:B:296:PHE:CG	1:B:367:ASP:HB3	2.53	0.43
1:B:203:THR:HG21	1:B:271:VAL:HG13	1.99	0.43
1:A:40:ARG:HH21	1:A:43:PHE:HB2	1.84	0.43
1:B:164:LEU:HD12	1:B:165:PRO:HD2	2.01	0.43
1:A:465:TRP:HA	1:A:466:SER:HA	1.77	0.43
1:B:425:THR:O	1:B:426:LYS:CB	2.47	0.43
1:A:237:GLU:N	1:A:238:PRO:CD	2.82	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:489:LYS:NZ	2:A:2237:HOH:O	2.50	0.43
1:A:452:ALA:O	1:A:457:VAL:HG13	2.19	0.42
1:B:95:CYS:HB3	2:B:2062:HOH:O	2.19	0.42
1:A:242:THR:HG23	1:A:300:TRP:CZ2	2.54	0.42
1:A:466:SER:O	1:A:482:GLY:HA2	2.20	0.42
1:A:199:ILE:HD11	1:A:202:TRP:CD1	2.55	0.42
1:A:57:GLN:O	1:A:470:ASN:HB2	2.20	0.42
1:B:306:THR:HG21	1:B:407:TYR:HB3	2.02	0.42
1:A:343:ASN:CG	1:A:416:GLU:HB2	2.40	0.42
1:A:409:VAL:HA	1:A:410:PRO:HD3	1.91	0.42
1:B:40:ARG:NH2	1:B:44:PRO:O	2.53	0.42
1:B:367:ASP:OD1	1:B:367:ASP:C	2.58	0.42
1:A:450:ARG:HD2	2:A:2003:HOH:O	2.20	0.41
1:B:413:TYR:CE2	1:B:460:LYS:HB2	2.54	0.41
1:B:242:THR:HG23	1:B:300:TRP:CZ2	2.55	0.41
1:B:185:CYS:SG	1:B:255:GLU:HG2	2.61	0.41
1:A:185:CYS:SG	1:A:255:GLU:HG2	2.60	0.41
1:A:202:TRP:O	1:A:268:ILE:HA	2.21	0.41
1:B:466:SER:O	1:B:482:GLY:HA2	2.20	0.41
1:B:319:LYS:HB3	1:B:320:GLY:H	1.65	0.40
1:B:441:TYR:CZ	1:B:445:HIS:CE1	3.10	0.40
1:B:363:TYR:CD2	1:B:363:TYR:C	2.94	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:2074:HOH:O	2:A:2216:HOH:O[3_655]	1.72	0.48
2:A:2047:HOH:O	2:B:2110:HOH:O[3_655]	1.97	0.23

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	464/532 (87%)	444 (96%)	20 (4%)	0	100	100
1	B	463/532 (87%)	443 (96%)	20 (4%)	0	100	100
All	All	927/1064 (87%)	887 (96%)	40 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	396/456 (87%)	344 (87%)	52 (13%)	4	6
1	B	395/456 (87%)	352 (89%)	43 (11%)	6	11
All	All	791/912 (87%)	696 (88%)	95 (12%)	5	8

All (95) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	40	ARG
1	A	41	ARG
1	A	45	GLN
1	A	63	ASN
1	A	84	SER
1	A	94	ASN
1	A	138	VAL
1	A	180	ILE
1	A	199	ILE
1	A	200	LYS
1	A	228	LYS
1	A	242	THR
1	A	247	LEU
1	A	250	LYS
1	A	255	GLU
1	A	257	ARG
1	A	267	GLU
1	A	268	ILE

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Mol	Chain	Res	Type
1	A	271	VAL
1	A	282	ASP
1	A	284	GLN
1	A	287	ILE
1	A	291	LYS
1	A	294	LEU
1	A	302	LEU
1	A	312	LYS
1	A	315	ARG
1	A	319	LYS
1	A	324	LYS
1	A	342	MET
1	A	350	VAL
1	A	354	VAL
1	A	361	LEU
1	A	368	GLN
1	A	374	GLU
1	A	377	GLN
1	A	378	LYS
1	A	391	VAL
1	A	396	LEU
1	A	409	VAL
1	A	412	LEU
1	A	419	MET
1	A	429	LEU
1	A	437	GLU
1	A	448	SER
1	A	457	VAL
1	A	460	LYS
1	A	470	ASN
1	A	474	ASN
1	A	475	LEU
1	A	493	ARG
1	A	502	TYR
1	B	40	ARG
1	B	63	ASN
1	B	76	THR
1	B	93	ILE
1	B	179	ARG
1	B	199	ILE
1	B	228	LYS
1	B	242	THR

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Mol	Chain	Res	Type
1	B	247	LEU
1	B	255	GLU
1	B	257	ARG
1	B	268	ILE
1	B	271	VAL
1	B	281	SER
1	B	291	LYS
1	B	294	LEU
1	B	302	LEU
1	B	306	THR
1	B	312	LYS
1	B	315	ARG
1	B	324	LYS
1	B	332	LYS
1	B	342	MET
1	B	350	VAL
1	B	354	VAL
1	B	371	LYS
1	B	375	ARG
1	B	391	VAL
1	B	392	VAL
1	B	396	LEU
1	B	404	LYS
1	B	405	GLU
1	B	409	VAL
1	B	412	LEU
1	B	419	MET
1	B	429	LEU
1	B	457	VAL
1	B	470	ASN
1	B	474	ASN
1	B	475	LEU
1	B	489	LYS
1	B	493	ARG
1	B	502	TYR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (27) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	45	GLN
1	A	57	GLN
1	A	63	ASN

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Mol	Chain	Res	Type
1	A	77	GLN
1	A	94	ASN
1	A	97	HIS
1	A	108	GLN
1	A	151	ASN
1	A	178	HIS
1	A	214	ASN
1	A	264	GLN
1	A	470	ASN
1	A	474	ASN
1	B	45	GLN
1	B	57	GLN
1	B	63	ASN
1	B	77	GLN
1	B	94	ASN
1	B	97	HIS
1	B	108	GLN
1	B	151	ASN
1	B	214	ASN
1	B	264	GLN
1	B	284	GLN
1	B	389	GLN
1	B	470	ASN
1	B	474	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	468/532 (87%)	-0.63	3 (0%)	89	90	17, 21, 26, 30	0
1	B	467/532 (87%)	-0.64	5 (1%)	80	82	17, 21, 26, 30	0
All	All	935/1064 (87%)	-0.63	8 (0%)	84	86	17, 21, 26, 30	0

All (8) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	377	GLN	3.3
1	A	377	GLN	2.9
1	B	37	VAL	2.7
1	A	354	VAL	2.6
1	A	376	ASN	2.4
1	B	283	VAL	2.3
1	B	45	GLN	2.3
1	B	354	VAL	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 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.