



Full wwPDB X-ray Structure Validation Report ⓘ

Nov 8, 2022 – 08:10 PM JST

PDB ID : 5YXT
Title : Crystal structure of reducing end xylose-releasing exo-oligoxylanase
Authors : Jiang, Z.Q.; You, X.; Huang, P.; Ma, J.W.
Deposited on : 2017-12-07
Resolution : 1.88 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	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.31.2
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)	:	2.31.2

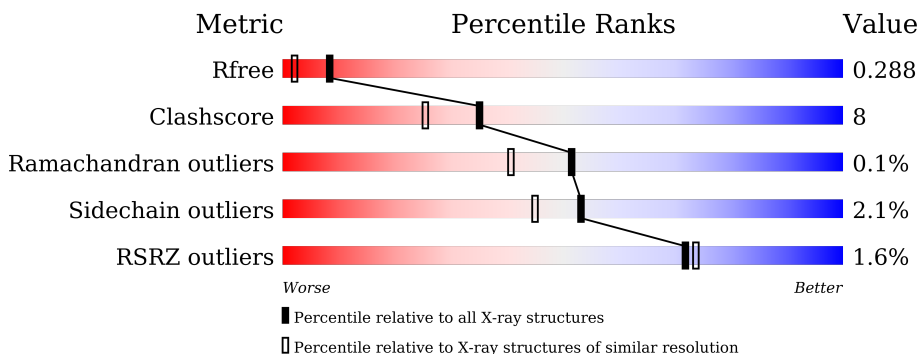
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.88 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	9470 (1.90-1.86)
Clashscore	141614	10282 (1.90-1.86)
Ramachandran outliers	138981	10152 (1.90-1.86)
Sidechain outliers	138945	10152 (1.90-1.86)
RSRZ outliers	127900	9303 (1.90-1.86)

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	 82% 16% ..
1	B	382	 83% 15% .
1	C	382	 82% 16% ..
1	D	382	 86% 14% .

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 13190 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 Reducing end xylose-releasing exo-oligoxyranase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	379	Total	C	N	O	S	0	0	0
			3038	1953	510	556	19			
1	B	379	Total	C	N	O	S	0	0	0
			3036	1952	505	560	19			
1	C	379	Total	C	N	O	S	0	0	0
			3044	1956	510	560	18			
1	D	379	Total	C	N	O	S	0	0	0
			3037	1951	510	558	18			

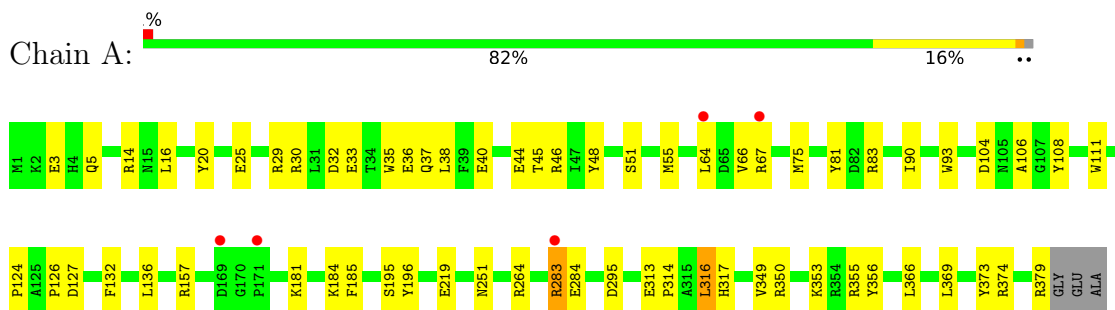
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	261	Total	O	0	0
			261	261		
2	B	264	Total	O	0	0
			264	264		
2	C	252	Total	O	0	0
			252	252		
2	D	258	Total	O	0	0
			258	258		

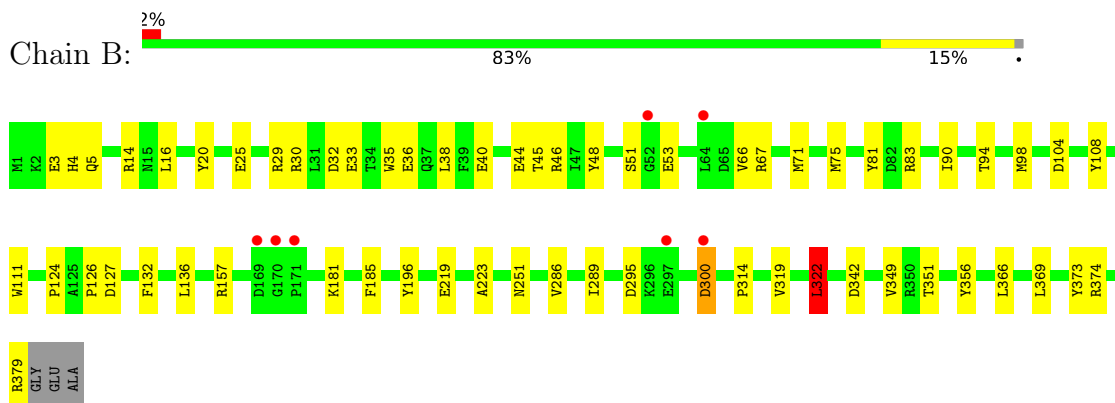
3 Residue-property plots [i](#)

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.

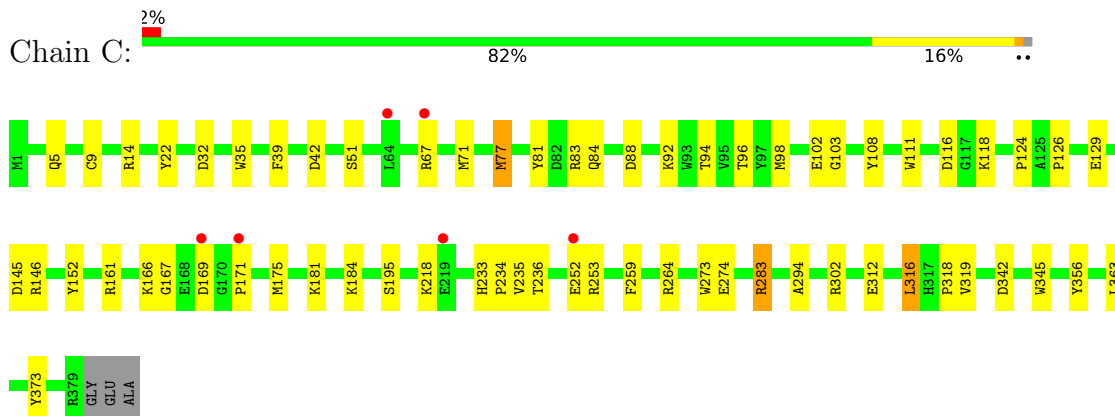
- Molecule 1: Reducing end xylose-releasing exo-oligoxylanase



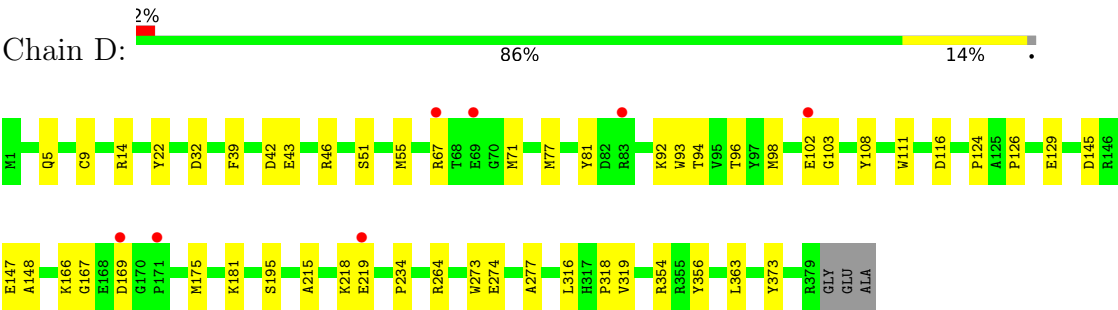
- Molecule 1: Reducing end xylose-releasing exo-oligoxylanase



- Molecule 1: Reducing end xylose-releasing exo-oligoxylanase



- Molecule 1: Reducing end xylose-releasing exo-oligoxylanase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	54.57Å 78.99Å 89.15Å 100.34° 95.86° 106.35°	Depositor
Resolution (Å)	22.35 – 1.88 23.95 – 1.88	Depositor EDS
% Data completeness (in resolution range)	93.0 (22.35-1.88) 93.0 (23.95-1.88)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.19 (at 1.88Å)	Xtriage
Refinement program	PHENIX (dev_2474: ???)	Depositor
R, R_{free}	0.244 , 0.289 0.244 , 0.288	Depositor DCC
R_{free} test set	1990 reflections (1.89%)	wwPDB-VP
Wilson B-factor (Å ²)	16.8	Xtriage
Anisotropy	0.002	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 53.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.36$, $\langle L^2 \rangle = 0.20$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	13190	wwPDB-VP
Average B, all atoms (Å ²)	23.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 99.26 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.8047e-12. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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	0.37	0/3144	0.54	0/4285
1	B	0.37	0/3142	0.57	3/4282 (0.1%)
1	C	0.42	0/3150	0.59	3/4292 (0.1%)
1	D	0.38	0/3143	0.54	0/4283
All	All	0.39	0/12579	0.56	6/17142 (0.0%)

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	322	LEU	CA-CB-CG	8.74	135.40	115.30
1	C	283	ARG	NE-CZ-NH1	-7.22	116.69	120.30
1	B	322	LEU	CB-CG-CD2	5.88	120.99	111.00
1	B	300	ASP	CB-CG-OD1	5.61	123.35	118.30
1	C	77	MET	CG-SD-CE	5.28	108.65	100.20
1	C	161	ARG	NE-CZ-NH1	5.20	122.90	120.30

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	3038	0	2764	52	0
1	B	3036	0	2757	49	0
1	C	3044	0	2771	53	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	3037	0	2758	41	0
2	A	261	0	0	18	1
2	B	264	0	0	19	3
2	C	252	0	0	19	2
2	D	258	0	0	12	2
All	All	13190	0	11050	190	4

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 8.

All (190) 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:C:9:CYS:SG	2:C:645:HOH:O	2.31	0.88
1:C:94:THR:HA	1:C:98:MET:HE2	1.58	0.84
1:D:9:CYS:SG	2:D:652:HOH:O	2.36	0.84
1:D:94:THR:HA	1:D:98:MET:HE2	1.60	0.83
1:B:75:MET:HE2	1:B:90:ILE:HG22	1.65	0.79
1:C:108:TYR:HA	2:C:401:HOH:O	1.85	0.76
1:A:283:ARG:NH2	1:C:345:TRP:HE1	1.83	0.76
1:A:313:GLU:OE2	2:A:402:HOH:O	2.02	0.75
1:D:108:TYR:HA	2:D:401:HOH:O	1.87	0.74
1:C:302:ARG:CZ	1:D:147:GLU:HG3	2.17	0.73
1:D:129:GLU:OE1	2:D:401:HOH:O	2.06	0.73
1:B:181:LYS:HD2	1:B:219:GLU:HB3	1.69	0.72
1:A:374:ARG:NH2	2:A:407:HOH:O	2.21	0.72
1:B:33:GLU:HB2	2:B:407:HOH:O	1.92	0.70
1:D:102:GLU:HG2	1:D:103:GLY:H	1.56	0.69
1:A:33:GLU:HB2	2:A:406:HOH:O	1.93	0.69
1:C:129:GLU:OE1	2:C:401:HOH:O	2.10	0.69
1:D:22:TYR:O	2:D:402:HOH:O	2.10	0.69
1:C:84:GLN:NE2	2:C:408:HOH:O	2.26	0.69
1:D:318:PRO:HG2	2:D:443:HOH:O	1.92	0.68
1:C:22:TYR:O	2:C:402:HOH:O	2.11	0.68
1:A:32:ASP:OD1	1:A:81:TYR:OH	2.11	0.68
1:B:32:ASP:OD1	1:B:81:TYR:OH	2.11	0.68
1:B:5:GLN:O	2:B:403:HOH:O	2.10	0.67
1:C:32:ASP:OD1	1:C:81:TYR:OH	2.10	0.67
1:A:75:MET:HE2	1:A:90:ILE:HG22	1.75	0.67
1:A:5:GLN:O	2:A:405:HOH:O	2.11	0.66
1:D:5:GLN:O	2:D:403:HOH:O	2.13	0.66

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:30:ARG:O	2:A:406:HOH:O	2.13	0.66
1:C:5:GLN:O	2:C:403:HOH:O	2.14	0.66
1:B:181:LYS:HE2	1:B:223:ALA:HB2	1.76	0.66
1:B:319:VAL:HA	1:B:322:LEU:HD13	1.78	0.66
1:C:318:PRO:HG2	2:C:446:HOH:O	1.95	0.66
1:B:374:ARG:NH2	2:B:406:HOH:O	2.18	0.65
1:D:32:ASP:OD1	1:D:81:TYR:OH	2.13	0.64
1:D:116:ASP:OD2	2:D:404:HOH:O	2.15	0.63
1:A:181:LYS:HD3	1:A:219:GLU:HB3	1.80	0.63
1:C:67:ARG:HD3	1:C:111:TRP:HE3	1.63	0.63
1:C:35:TRP:CZ2	1:C:83:ARG:HD3	2.34	0.62
1:B:71:MET:HG3	1:B:98:MET:CE	2.29	0.62
1:C:234:PRO:HB3	2:C:626:HOH:O	2.00	0.62
1:A:16:LEU:HD22	1:A:369:LEU:HD22	1.80	0.61
1:A:44:GLU:HG2	2:A:415:HOH:O	2.00	0.61
1:C:218:LYS:HG2	2:C:455:HOH:O	2.00	0.61
1:D:42:ASP:C	1:D:46:ARG:HD3	2.20	0.61
1:B:16:LEU:HD22	1:B:369:LEU:HD22	1.80	0.61
1:C:102:GLU:HG2	1:C:103:GLY:N	2.16	0.61
1:C:302:ARG:NH2	1:D:147:GLU:HG3	2.17	0.60
1:B:94:THR:HA	1:B:98:MET:HE2	1.82	0.59
1:A:316:LEU:HD12	1:A:317:HIS:CE1	2.37	0.59
1:D:215:ALA:O	1:D:219:GLU:HG2	2.02	0.59
1:B:66:VAL:N	2:B:415:HOH:O	2.35	0.59
1:A:66:VAL:N	2:A:413:HOH:O	2.35	0.59
1:C:312:GLU:OE2	2:C:404:HOH:O	2.16	0.59
1:A:283:ARG:NH2	1:C:345:TRP:NE1	2.52	0.58
1:C:92:LYS:O	1:C:96:THR:HG23	2.04	0.57
1:D:92:LYS:O	1:D:96:THR:HG23	2.05	0.57
1:C:166:LYS:NZ	2:C:419:HOH:O	2.37	0.56
1:C:233:HIS:CD2	1:C:235:VAL:H	2.22	0.56
1:D:102:GLU:HG2	1:D:103:GLY:N	2.20	0.56
1:A:25:GLU:OE2	1:A:29:ARG:NH2	2.39	0.56
1:C:102:GLU:HG2	1:C:103:GLY:H	1.70	0.56
1:B:53:GLU:OE1	2:B:405:HOH:O	2.18	0.56
1:C:146:ARG:NH2	1:C:152:TYR:OH	2.39	0.55
1:D:166:LYS:NZ	2:D:418:HOH:O	2.38	0.55
1:C:342:ASP:OD2	2:C:405:HOH:O	2.18	0.55
1:A:283:ARG:HD2	2:C:423:HOH:O	2.07	0.55
1:D:55:MET:HE3	1:D:93:TRP:CD1	2.41	0.55
1:C:252:GLU:HG2	1:C:253:ARG:N	2.22	0.55

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:30:ARG:HD3	2:B:407:HOH:O	2.06	0.54
1:D:147:GLU:OE1	1:D:148:ALA:N	2.32	0.54
1:B:286:VAL:HA	1:B:289:ILE:HG22	1.88	0.54
1:C:77:MET:CE	1:C:363:LEU:HD23	2.38	0.54
1:D:218:LYS:HG2	2:D:456:HOH:O	2.07	0.54
1:D:77:MET:CE	1:D:363:LEU:HD23	2.38	0.54
1:B:83:ARG:NH1	2:B:418:HOH:O	2.41	0.54
1:B:35:TRP:CZ2	1:B:83:ARG:HD3	2.42	0.54
1:A:83:ARG:NH1	2:A:419:HOH:O	2.41	0.53
1:A:283:ARG:HH22	1:C:345:TRP:HE1	1.53	0.53
1:C:67:ARG:HD2	1:C:111:TRP:O	2.08	0.53
1:A:136:LEU:HG	2:A:631:HOH:O	2.08	0.53
1:A:30:ARG:HD3	2:A:406:HOH:O	2.08	0.53
1:C:145:ASP:O	2:C:407:HOH:O	2.19	0.53
1:A:35:TRP:CZ2	1:A:83:ARG:HD3	2.43	0.52
1:D:43:GLU:N	1:D:46:ARG:NH1	2.58	0.52
1:C:67:ARG:HD3	1:C:111:TRP:CE3	2.43	0.52
1:A:283:ARG:HG2	1:A:284:GLU:N	2.25	0.52
1:B:71:MET:HG3	1:B:98:MET:HE1	1.89	0.52
1:B:157:ARG:NH1	2:B:402:HOH:O	2.07	0.52
1:A:30:ARG:HA	2:A:406:HOH:O	2.08	0.51
1:C:81:TYR:HD2	1:C:83:ARG:HD2	1.75	0.51
1:C:88:ASP:OD2	1:C:146:ARG:NH2	2.35	0.51
1:C:71:MET:HG3	1:C:98:MET:CE	2.41	0.51
1:A:66:VAL:HG23	2:A:413:HOH:O	2.11	0.50
1:B:4:HIS:HE1	2:B:627:HOH:O	1.95	0.50
1:D:316:LEU:HD11	1:D:354:ARG:HE	1.76	0.50
1:B:44:GLU:HG3	2:B:416:HOH:O	2.11	0.49
1:D:234:PRO:HB3	2:D:634:HOH:O	2.13	0.49
1:B:71:MET:HG3	1:B:98:MET:HE3	1.94	0.49
1:A:83:ARG:NH2	2:A:425:HOH:O	2.46	0.49
1:D:145:ASP:O	2:D:406:HOH:O	2.20	0.49
1:B:32:ASP:O	1:B:36:GLU:HG3	2.12	0.49
1:D:43:GLU:CA	1:D:46:ARG:CZ	2.91	0.49
1:A:157:ARG:NH1	2:A:404:HOH:O	2.10	0.49
1:A:295:ASP:OD1	1:A:295:ASP:N	2.41	0.49
1:B:295:ASP:OD1	1:B:295:ASP:N	2.41	0.49
1:D:71:MET:HG3	1:D:98:MET:CE	2.43	0.48
1:C:259:PHE:CE2	1:C:316:LEU:HD23	2.49	0.48
1:B:30:ARG:HG3	1:B:366:LEU:HD21	1.96	0.48
1:B:66:VAL:HG23	2:B:415:HOH:O	2.13	0.48

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:136:LEU:HG	2:B:637:HOH:O	2.14	0.47
1:B:30:ARG:HA	2:B:407:HOH:O	2.13	0.47
1:A:30:ARG:NH1	2:A:427:HOH:O	2.47	0.47
1:A:32:ASP:O	1:A:36:GLU:HG3	2.14	0.47
1:C:233:HIS:HD2	1:C:236:THR:H	1.62	0.47
1:A:30:ARG:HG3	1:A:366:LEU:HD21	1.97	0.47
1:D:77:MET:HE2	1:D:363:LEU:HD23	1.96	0.47
1:D:316:LEU:HD11	1:D:354:ARG:NE	2.30	0.47
1:D:39:PHE:HE1	1:D:77:MET:HE2	1.78	0.47
1:C:273:TRP:NE1	1:C:283:ARG:NH1	2.63	0.47
1:C:77:MET:HE3	1:C:363:LEU:HD23	1.97	0.46
1:A:64:LEU:O	1:A:64:LEU:HG	2.16	0.46
1:B:44:GLU:HG2	1:B:45:THR:HG23	1.97	0.46
1:C:294:ALA:HB1	2:C:444:HOH:O	2.15	0.46
1:D:126:PRO:HG3	1:D:175:MET:HG2	1.96	0.46
1:C:169:ASP:OD2	1:C:169:ASP:N	2.47	0.46
1:B:83:ARG:NH1	2:B:432:HOH:O	2.48	0.46
1:B:342:ASP:OD2	2:B:408:HOH:O	2.21	0.46
1:A:181:LYS:CD	1:A:219:GLU:HB3	2.44	0.46
1:A:283:ARG:CD	2:C:423:HOH:O	2.63	0.46
1:A:316:LEU:HD12	1:A:317:HIS:NE2	2.31	0.46
1:B:25:GLU:OE2	1:B:29:ARG:NH2	2.49	0.46
1:B:30:ARG:NH1	2:B:433:HOH:O	2.49	0.45
1:D:169:ASP:N	1:D:169:ASP:OD1	2.48	0.45
1:B:75:MET:HE2	1:B:90:ILE:CG2	2.41	0.45
1:C:319:VAL:HG23	2:C:446:HOH:O	2.15	0.45
1:B:67:ARG:HA	1:B:111:TRP:O	2.16	0.45
1:C:39:PHE:HE1	1:C:77:MET:HE2	1.81	0.45
1:B:75:MET:HE3	1:B:132:PHE:HE1	1.82	0.45
1:B:75:MET:CE	1:B:132:PHE:HE1	2.30	0.45
1:C:126:PRO:HG3	1:C:175:MET:HG2	1.98	0.45
1:C:166:LYS:HD3	1:C:167:GLY:H	1.81	0.45
1:A:67:ARG:HA	1:A:111:TRP:O	2.16	0.45
1:C:14:ARG:NH1	1:C:274:GLU:OE1	2.50	0.45
1:B:3:GLU:HG3	1:B:14:ARG:HD3	1.99	0.44
1:A:3:GLU:HG3	1:A:14:ARG:HD3	1.99	0.44
1:D:319:VAL:HG23	2:D:443:HOH:O	2.16	0.44
1:C:67:ARG:HA	1:C:111:TRP:O	2.18	0.44
1:D:166:LYS:HD3	1:D:167:GLY:H	1.83	0.43
1:C:77:MET:HE3	1:C:363:LEU:HB3	2.01	0.43
1:D:14:ARG:NH1	1:D:274:GLU:OE1	2.52	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:92:LYS:HA	1:D:92:LYS:HD3	1.85	0.43
1:D:67:ARG:HA	1:D:111:TRP:O	2.19	0.43
1:B:126:PRO:HD2	1:B:185:PHE:O	2.19	0.43
1:A:46:ARG:NE	1:A:48:TYR:O	2.52	0.42
1:B:16:LEU:HD21	1:B:20:TYR:CE2	2.54	0.42
1:B:300:ASP:O	1:B:300:ASP:OD1	2.37	0.42
1:A:106:ALA:N	2:A:436:HOH:O	2.51	0.42
1:A:40:GLU:OE2	1:A:83:ARG:NH1	2.52	0.42
1:D:71:MET:HG3	1:D:98:MET:HE1	2.02	0.42
1:C:184:LYS:NZ	2:C:413:HOH:O	2.31	0.42
1:D:273:TRP:CE3	1:D:277:ALA:HA	2.55	0.42
1:A:16:LEU:HD21	1:A:20:TYR:CE2	2.55	0.42
1:A:75:MET:HE1	1:A:132:PHE:HE1	1.84	0.42
1:C:42:ASP:HB2	2:C:415:HOH:O	2.19	0.42
1:B:38:LEU:HD21	1:B:349:VAL:HG21	2.01	0.42
1:D:195:SER:HB2	1:D:264:ARG:HB2	2.02	0.42
1:A:55:MET:HE1	1:A:93:TRP:HA	2.03	0.41
1:A:75:MET:HE1	1:A:132:PHE:CE1	2.55	0.41
1:A:127:ASP:HB3	1:A:196:TYR:CD1	2.56	0.41
1:A:184:LYS:NZ	2:A:403:HOH:O	2.09	0.41
1:C:116:ASP:OD2	1:C:118:LYS:HG3	2.21	0.41
1:B:30:ARG:O	2:B:407:HOH:O	2.21	0.41
1:B:314:PRO:HB3	2:B:600:HOH:O	2.20	0.41
1:A:38:LEU:HD21	1:A:349:VAL:HG21	2.02	0.41
1:A:350:ARG:HD3	1:A:355:ARG:HA	2.01	0.41
1:C:273:TRP:CZ2	1:C:283:ARG:HD2	2.54	0.41
1:D:108:TYR:HB3	1:D:124:PRO:HB3	2.02	0.41
1:A:126:PRO:HD2	1:A:185:PHE:O	2.20	0.41
1:A:195:SER:HB2	1:A:264:ARG:HB2	2.02	0.41
1:B:40:GLU:OE1	1:B:83:ARG:NH1	2.53	0.41
1:C:195:SER:HB2	1:C:264:ARG:HB2	2.02	0.41
1:A:314:PRO:HB3	2:A:590:HOH:O	2.21	0.41
1:B:108:TYR:HB3	1:B:124:PRO:HB3	2.03	0.41
1:A:37:GLN:HB3	1:A:45:THR:OG1	2.21	0.40
1:B:351:THR:HG22	2:B:588:HOH:O	2.21	0.40
1:C:108:TYR:HB3	1:C:124:PRO:HB3	2.02	0.40
1:A:108:TYR:HB3	1:A:124:PRO:HB3	2.04	0.40
1:B:46:ARG:NE	1:B:48:TYR:O	2.54	0.40
1:B:127:ASP:HB3	1:B:196:TYR:CD1	2.57	0.40

All (4) 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:B:635:HOH:O	2:C:463:HOH:O[1_565]	1.89	0.31
2:B:528:HOH:O	2:C:629:HOH:O[1_455]	2.12	0.08
2:A:632:HOH:O	2:D:500:HOH:O[1_665]	2.13	0.07
2:B:567:HOH:O	2:D:469:HOH:O[1_554]	2.17	0.03

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	377/382 (99%)	369 (98%)	8 (2%)	0	100	100
1	B	377/382 (99%)	369 (98%)	8 (2%)	0	100	100
1	C	377/382 (99%)	364 (97%)	12 (3%)	1 (0%)	41	30
1	D	377/382 (99%)	366 (97%)	11 (3%)	0	100	100
All	All	1508/1528 (99%)	1468 (97%)	39 (3%)	1 (0%)	51	41

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	171	PRO

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	293/309 (95%)	284 (97%)	9 (3%)	40	29
1	B	295/309 (96%)	288 (98%)	7 (2%)	49	39

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	295/309 (96%)	290 (98%)	5 (2%)	60	54
1	D	293/309 (95%)	289 (99%)	4 (1%)	67	62
All	All	1176/1236 (95%)	1151 (98%)	25 (2%)	53	45

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	51	SER
1	A	104	ASP
1	A	251	ASN
1	A	283	ARG
1	A	316	LEU
1	A	353	LYS
1	A	356	TYR
1	A	373	TYR
1	A	379	ARG
1	B	51	SER
1	B	104	ASP
1	B	251	ASN
1	B	322	LEU
1	B	356	TYR
1	B	373	TYR
1	B	379	ARG
1	C	51	SER
1	C	181	LYS
1	C	316	LEU
1	C	356	TYR
1	C	373	TYR
1	D	51	SER
1	D	181	LYS
1	D	356	TYR
1	D	373	TYR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	28	GLN
1	C	233	HIS
1	D	5	GLN

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

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	379/382 (99%)	-0.36	5 (1%) 77 79	13, 21, 34, 54	0
1	B	379/382 (99%)	-0.35	7 (1%) 68 70	13, 21, 35, 53	0
1	C	379/382 (99%)	-0.34	6 (1%) 72 74	13, 21, 34, 55	0
1	D	379/382 (99%)	-0.36	7 (1%) 68 70	13, 21, 35, 55	0
All	All	1516/1528 (99%)	-0.35	25 (1%) 72 74	13, 21, 35, 55	0

All (25) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	171	PRO	6.3
1	B	171	PRO	5.8
1	C	64	LEU	5.1
1	C	252	GLU	4.2
1	B	64	LEU	4.0
1	C	171	PRO	4.0
1	D	171	PRO	4.0
1	A	64	LEU	3.9
1	C	67	ARG	3.8
1	A	169	ASP	3.8
1	B	169	ASP	3.8
1	D	67	ARG	3.7
1	C	169	ASP	3.4
1	D	169	ASP	3.4
1	B	300	ASP	3.3
1	A	283	ARG	2.9
1	B	297	GLU	2.7
1	D	219	GLU	2.6
1	D	69	GLU	2.4
1	A	67	ARG	2.3
1	C	219	GLU	2.3

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	B	52	GLY	2.3
1	D	83	ARG	2.2
1	B	170	GLY	2.2
1	D	102	GLU	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.