



Full wwPDB X-ray Structure Validation Report ⓘ

May 21, 2020 – 04:33 am BST

PDB ID : 2DC0
Title : Crystal structure of amidase
Authors : Ohshima, T.; Sakuraba, H.; Ebihara, A.; Kanagawa, M.; Nakagawa, N.; Kuroishi, C.; Satoh, S.; Kuramitsu, S.; Yokoyama, S.; RIKEN Structural Genomics/Proteomics Initiative (RSGI)
Deposited on : 2005-12-17
Resolution : 2.00 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
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) : 2.11

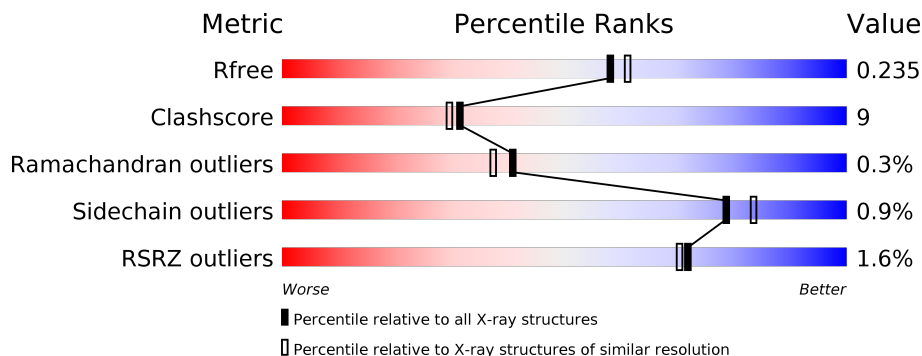
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (2.00-2.00)

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 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	434	 % 83% 16% •
1	B	434	 2% 80% 19% •

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 6879 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 probable amidase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	S	Se			
1	A	434	3282	2085	587	606	2	2	0	0	0
1	B	434	3282	2085	587	606	2	2	0	0	0

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	117	MSE	MET	MODIFIED RESIDUE	GB 55773179
A	406	MSE	MET	MODIFIED RESIDUE	GB 55773179
B	1117	MSE	MET	MODIFIED RESIDUE	GB 55773179
B	1406	MSE	MET	MODIFIED RESIDUE	GB 55773179

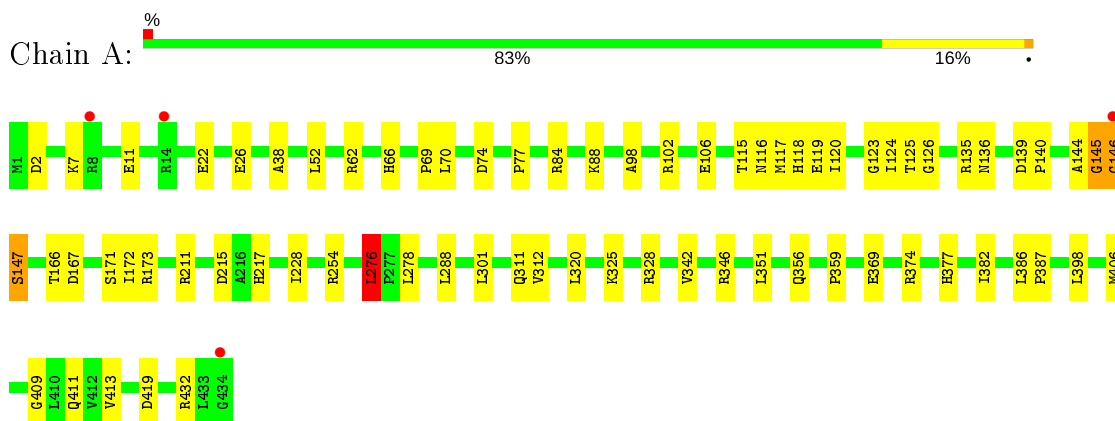
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	169	Total	O	0	0
			169	169		
2	B	146	Total	O	0	0
			146	146		

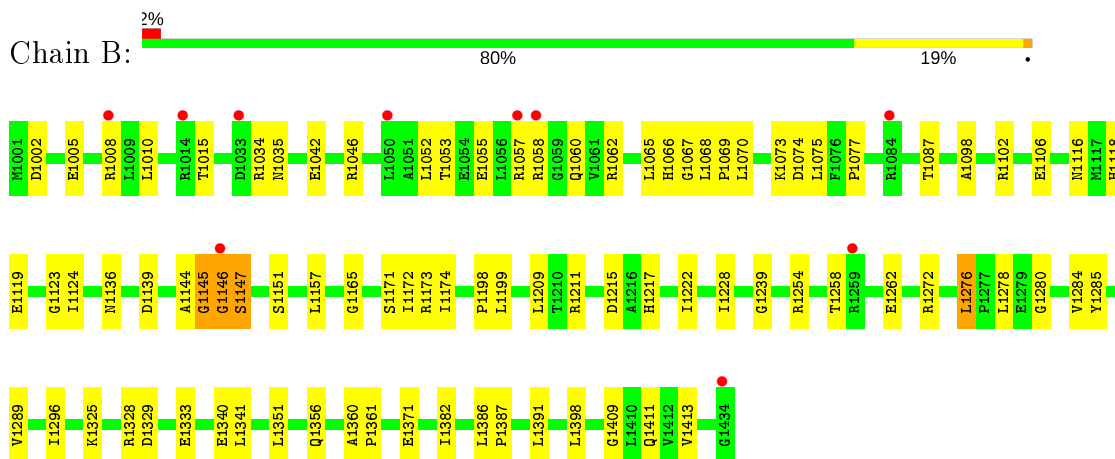
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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: probable amidase



- Molecule 1: probable amidase



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	81.34Å 81.34Å 313.26Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.13 – 2.00 46.36 – 2.00	Depositor EDS
% Data completeness (in resolution range)	(Not available) (46.13-2.00) 96.5 (46.36-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	18.05 (at 2.00Å)	Xtrriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.220 , 0.240 0.218 , 0.235	Depositor DCC
R_{free} test set	4883 reflections (6.99%)	wwPDB-VP
Wilson B-factor (Å ²)	21.1	Xtrriage
Anisotropy	0.470	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 41.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6879	wwPDB-VP
Average B, all atoms (Å ²)	23.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.85% 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 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	0.48	5/3342 (0.1%)	0.65	7/4533 (0.2%)
1	B	0.42	3/3342 (0.1%)	0.64	6/4533 (0.1%)
All	All	0.45	8/6684 (0.1%)	0.64	13/9066 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	B	0	1
All	All	0	2

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	144	ALA	C-N	-14.78	1.06	1.33
1	B	1144	ALA	C-N	-11.77	1.11	1.33
1	A	146	GLY	C-N	9.01	1.54	1.34
1	B	1146	GLY	C-N	9.00	1.54	1.34
1	A	115	THR	C-N	-6.21	1.19	1.34
1	A	117	MSE	C-N	-5.74	1.20	1.34
1	A	145	GLY	C-N	5.43	1.42	1.33
1	B	1145	GLY	C-N	5.42	1.42	1.33

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	1147	SER	O-C-N	-7.60	110.54	122.70
1	A	147	SER	O-C-N	-7.29	111.03	122.70
1	A	146	GLY	N-CA-C	5.84	127.71	113.10
1	B	1146	GLY	N-CA-C	5.84	127.69	113.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	276	LEU	CB-CA-C	5.71	121.06	110.20
1	B	1276	LEU	CB-CA-C	5.71	121.04	110.20
1	B	1147	SER	CA-C-N	5.35	128.98	117.20
1	A	147	SER	CA-C-N	5.28	128.82	117.20
1	A	115	THR	O-C-N	-5.28	114.25	122.70
1	A	145	GLY	O-C-N	-5.17	114.41	123.20
1	B	1145	GLY	O-C-N	-5.15	114.44	123.20
1	B	1147	SER	N-CA-C	-5.10	97.24	111.00
1	A	147	SER	N-CA-C	-5.09	97.27	111.00

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	145	GLY	Mainchain
1	B	1145	GLY	Mainchain

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	3282	0	3383	57	0
1	B	3282	0	3380	63	0
2	A	169	0	0	1	0
2	B	146	0	0	3	0
All	All	6879	0	6763	120	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 9.

All (120) 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:38:ALA:HB2	1:A:116:ASN:HD22	1.28	0.95
1:B:1058:ARG:HG3	1:B:1060:GLN:HG3	1.51	0.90

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:38:ALA:HB2	1:A:116:ASN:ND2	1.87	0.89
1:B:1062:ARG:H	1:B:1066:HIS:HD2	1.26	0.82
1:A:62:ARG:H	1:A:66:HIS:HD2	1.30	0.78
1:B:1055:GLU:HA	1:B:1058:ARG:HD3	1.68	0.73
1:A:38:ALA:CB	1:A:116:ASN:ND2	2.52	0.72
1:B:1058:ARG:NE	1:B:1060:GLN:HE21	1.89	0.70
1:A:124:ILE:C	1:A:124:ILE:HD12	2.11	0.70
1:B:1198:PRO:HB2	1:B:1296:ILE:HD13	1.74	0.69
1:A:124:ILE:HG21	1:A:382:ILE:HD13	1.77	0.67
1:B:1002:ASP:HB2	1:B:1215:ASP:OD2	1.95	0.66
1:B:1356:GLN:NE2	1:B:1409:GLY:H	1.95	0.65
1:B:1124:ILE:HG21	1:B:1382:ILE:HD13	1.79	0.64
1:A:7:LYS:O	1:A:11:GLU:HG3	1.98	0.64
1:A:356:GLN:NE2	1:A:409:GLY:H	1.96	0.63
1:A:38:ALA:HA	1:A:116:ASN:HD21	1.63	0.63
1:B:1258:THR:O	1:B:1262:GLU:HG3	1.98	0.63
1:B:1217:HIS:HE1	1:B:1228:ILE:O	1.81	0.62
1:A:369:GLU:HG2	1:A:374:ARG:HH12	1.65	0.61
1:A:118:HIS:HD1	1:A:126:GLY:H	1.47	0.61
1:A:124:ILE:CD1	1:A:124:ILE:C	2.69	0.61
1:B:1329:ASP:O	1:B:1333:GLU:HG3	2.01	0.61
1:B:1062:ARG:H	1:B:1066:HIS:CD2	2.14	0.61
1:B:1356:GLN:HE21	1:B:1409:GLY:H	1.48	0.61
1:A:382:ILE:HG23	1:A:386:LEU:HD23	1.83	0.60
1:B:1254:ARG:HD3	1:B:1371:GLU:OE1	2.01	0.60
1:A:369:GLU:OE2	1:A:374:ARG:HG2	2.03	0.59
1:A:62:ARG:H	1:A:66:HIS:CD2	2.16	0.58
1:B:1055:GLU:HA	1:B:1058:ARG:CD	2.30	0.58
1:B:1382:ILE:HG23	1:B:1386:LEU:HD23	1.85	0.58
1:A:124:ILE:HD11	1:A:377:HIS:CE1	2.39	0.58
1:A:124:ILE:CD1	1:A:377:HIS:HE1	2.16	0.57
1:B:1035:ASN:O	1:B:1116:ASN:OD1	2.23	0.56
1:A:38:ALA:CA	1:A:116:ASN:ND2	2.68	0.56
1:A:217:HIS:HE1	1:A:228:ILE:O	1.87	0.56
1:A:2:ASP:HB2	1:A:215:ASP:OD2	2.08	0.54
1:B:1102:ARG:O	1:B:1106:GLU:HG3	2.08	0.54
1:B:1199:LEU:O	1:B:1296:ILE:HD12	2.08	0.53
1:B:1052:LEU:HD13	1:B:1067:GLY:HA2	1.91	0.53
1:A:356:GLN:HE21	1:A:409:GLY:H	1.55	0.53
1:B:1034:ARG:NH1	1:B:1136:ASN:O	2.42	0.52
1:A:22:GLU:O	1:A:26:GLU:HG3	2.09	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:301:LEU:HD23	1:A:320:LEU:HD11	1.91	0.52
1:A:38:ALA:HA	1:A:116:ASN:ND2	2.25	0.52
1:A:118:HIS:HB3	1:A:147:SER:HB2	1.92	0.52
1:A:77:PRO:HD2	1:A:98:ALA:HB2	1.92	0.52
1:B:1053:THR:HG22	1:B:1057:ARG:NH1	2.25	0.51
1:B:1010:LEU:HD23	1:B:1015:THR:HG23	1.92	0.51
1:B:1055:GLU:O	1:B:1058:ARG:HG2	2.10	0.50
1:A:369:GLU:HG2	1:A:374:ARG:NH1	2.25	0.50
1:B:1042:GLU:HB3	1:B:1046:ARG:NH1	2.27	0.50
1:A:166:THR:O	1:A:172:ILE:HG13	2.12	0.50
1:B:1296:ILE:HD11	2:B:102:HOH:O	2.12	0.50
1:A:276:LEU:HG	1:A:278:LEU:HG	1.94	0.49
1:B:1087:THR:HB	1:B:1119:GLU:HG3	1.95	0.49
1:B:1073:LYS:HE3	1:B:1147:SER:O	2.13	0.48
1:A:124:ILE:HD12	1:A:125:THR:N	2.28	0.48
1:A:38:ALA:CB	1:A:116:ASN:HD22	2.11	0.48
1:B:1341:LEU:HD22	1:B:1391:LEU:HD12	1.96	0.48
1:A:167:ASP:HB2	1:A:172:ILE:HB	1.96	0.47
1:A:325:LYS:HD3	1:A:328:ARG:HH22	1.78	0.47
1:B:1340:GLU:HG3	2:B:303:HOH:O	2.14	0.47
1:A:301:LEU:HD23	1:A:320:LEU:CD1	2.44	0.47
1:B:1052:LEU:O	1:B:1066:HIS:HE1	1.98	0.47
1:A:118:HIS:CD2	1:A:123:GLY:HA3	2.50	0.47
1:B:1386:LEU:HD12	1:B:1387:PRO:N	2.30	0.46
1:B:1077:PRO:HD2	1:B:1098:ALA:HB2	1.98	0.46
1:B:1198:PRO:HB2	1:B:1296:ILE:CD1	2.44	0.46
1:A:325:LYS:HD3	1:A:328:ARG:NH2	2.31	0.46
1:A:359:PRO:HB2	1:A:406:MSE:HB2	1.96	0.46
1:B:1211:ARG:HD2	2:B:81:HOH:O	2.14	0.46
1:B:1280:GLY:O	1:B:1284:VAL:HG23	2.16	0.46
1:A:211:ARG:HD2	2:A:475:HOH:O	2.16	0.46
1:B:1053:THR:HG22	1:B:1057:ARG:HH12	1.81	0.46
1:B:1174:ILE:HD11	1:B:1382:ILE:HD12	1.97	0.46
1:B:1285:TYR:CE1	1:B:1289:VAL:HG21	2.51	0.46
1:B:1136:ASN:HB3	1:B:1139:ASP:O	2.16	0.46
1:A:52:LEU:O	1:A:66:HIS:HE1	1.99	0.45
1:A:84:ARG:HH21	1:A:88:LYS:NZ	2.15	0.45
1:B:1069:PRO:C	1:B:1070:LEU:HD12	2.37	0.45
1:B:1075:LEU:C	1:B:1075:LEU:HD12	2.37	0.45
1:B:1065:LEU:HA	1:B:1068:LEU:HD12	1.98	0.45
1:B:1146:GLY:O	1:B:1171:SER:HA	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:69:PRO:C	1:A:70:LEU:HD12	2.37	0.45
1:A:386:LEU:HD12	1:A:387:PRO:N	2.33	0.44
1:B:1173:ARG:HH11	1:B:1411:GLN:NE2	2.16	0.44
1:B:1165:GLY:C	1:B:1172:ILE:HD11	2.37	0.44
1:A:288:LEU:HD23	1:A:288:LEU:C	2.38	0.43
1:A:167:ASP:HB2	1:A:172:ILE:HD12	2.01	0.43
1:B:1118:HIS:HB3	1:B:1147:SER:HB2	2.00	0.43
1:A:119:GLU:HB3	1:A:312:VAL:HG21	2.01	0.42
1:B:1008:ARG:HH11	1:B:1008:ARG:HG3	1.84	0.42
1:A:102:ARG:O	1:A:106:GLU:HG3	2.19	0.42
1:A:351:LEU:HB2	1:A:413:VAL:HB	2.01	0.42
1:A:136:ASN:HB3	1:A:139:ASP:O	2.20	0.42
1:A:146:GLY:O	1:A:171:SER:HA	2.20	0.42
1:B:1239:GLY:HA2	1:B:1272:ARG:O	2.19	0.42
1:B:1058:ARG:NE	1:B:1060:GLN:NE2	2.64	0.42
1:B:1351:LEU:HB2	1:B:1413:VAL:HB	2.02	0.42
1:A:124:ILE:CD1	1:A:377:HIS:CE1	2.98	0.41
1:B:1118:HIS:CD2	1:B:1123:GLY:HA3	2.55	0.41
1:B:1151:SER:HB3	1:B:1209:LEU:HD12	2.01	0.41
1:A:124:ILE:HG21	1:A:382:ILE:CD1	2.49	0.41
1:A:173:ARG:HH11	1:A:411:GLN:NE2	2.18	0.41
1:B:1124:ILE:HG21	1:B:1382:ILE:CD1	2.49	0.41
1:A:432:ARG:HH11	1:A:432:ARG:HG3	1.85	0.41
1:B:1034:ARG:HG3	1:B:1157:LEU:CD1	2.51	0.41
1:B:1360:ALA:HA	1:B:1361:PRO:HD3	1.87	0.41
1:B:1065:LEU:HG	1:B:1222:ILE:HD13	2.02	0.41
1:B:1034:ARG:HG3	1:B:1157:LEU:HD11	2.03	0.40
1:A:342:VAL:O	1:A:346:ARG:HG3	2.21	0.40
1:B:1147:SER:H	1:B:1171:SER:CB	2.35	0.40
1:A:276:LEU:HD21	1:A:278:LEU:HD11	2.03	0.40
1:B:1002:ASP:OD1	1:B:1005:GLU:HB2	2.22	0.40
1:B:1325:LYS:HD3	1:B:1328:ARG:HH22	1.86	0.40
1:B:1386:LEU:HG	1:B:1387:PRO:HD3	2.03	0.40
1:A:374:ARG:HH11	1:A:374:ARG:HG2	1.86	0.40
1:A:135:ARG:HB3	1:A:140:PRO:HA	2.03	0.40
1:B:1276:LEU:HG	1:B:1278:LEU:HG	2.02	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	432/434 (100%)	416 (96%)	14 (3%)	2 (0%)	29	23
1	B	432/434 (100%)	422 (98%)	9 (2%)	1 (0%)	47	44
All	All	864/868 (100%)	838 (97%)	23 (3%)	3 (0%)	41	37

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	1074	ASP
1	A	74	ASP
1	A	120	ILE

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	337/335 (101%)	332 (98%)	5 (2%)	65	69
1	B	337/335 (101%)	336 (100%)	1 (0%)	92	95
All	All	674/670 (101%)	668 (99%)	6 (1%)	78	83

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	254	ARG
1	A	276	LEU
1	A	311	GLN

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Mol	Chain	Res	Type
1	A	398	LEU
1	A	419	ASP
1	B	1398	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	66	HIS
1	A	116	ASN
1	A	217	HIS
1	A	235	ASN
1	A	311	GLN
1	A	356	GLN
1	A	411	GLN
1	B	1060	GLN
1	B	1066	HIS
1	B	1116	ASN
1	B	1217	HIS
1	B	1356	GLN
1	B	1411	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 carbohydrates 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](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	B	2
1	A	2

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	B	1405:GLY	C	1406:MSE	N	1.62
1	A	115:THR	C	116:ASN	N	1.19
1	B	1144:ALA	C	1145:GLY	N	1.11
1	A	144:ALA	C	145:GLY	N	1.06

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	432/434 (99%)	-0.31	4 (0%) 84 83	13, 20, 34, 46	0
1	B	432/434 (99%)	-0.11	10 (2%) 60 59	13, 22, 37, 47	0
All	All	864/868 (99%)	-0.21	14 (1%) 72 70	13, 21, 36, 47	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	434	GLY	5.6
1	B	1434	GLY	4.3
1	B	1033	ASP	3.8
1	B	1014	ARG	3.1
1	B	1146	GLY	3.0
1	B	1057	ARG	3.0
1	B	1058	ARG	2.8
1	B	1259	ARG	2.6
1	A	8	ARG	2.6
1	A	146	GLY	2.5
1	B	1050	LEU	2.4
1	A	14	ARG	2.3
1	B	1008	ARG	2.3
1	B	1084	ARG	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 carbohydrates in this entry.

6.4 Ligands

There are no ligands in this entry.

6.5 Other polymers

There are no such residues in this entry.