Full wwPDB X-ray Structure Validation Report (i)

Feb 1, 2016 – 08:08 AM GMT

PDB ID	:	3DHU
Title	:	Crystal structure of an alpha-amylase from Lactobacillus plantarum
Authors	:	Bonanno, J.B.; Dickey, M.; Bain, K.T.; Iizuka, M.; Ozyurt, S.; Smith, D.; Wasserman, S.; Sauder, J.M.; Burley, S.K.; Almo, S.C.; New York SGX Research Center for Structural Genomics (NYSGXRC)
Deposited on	:	2008-06-18
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
http://wwpdb.org/validation/2016/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

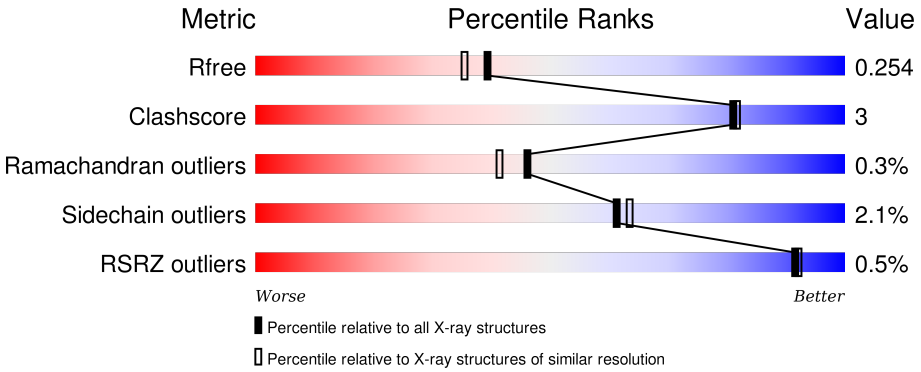
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk26865

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 2.00 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}	91344	6249 (2.00-2.00)
Clashscore	102246	7340 (2.00-2.00)
Ramachandran outliers	100387	7248 (2.00-2.00)
Sidechain outliers	100360	7247 (2.00-2.00)
RSRZ outliers	91569	6262 (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. 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	449	83% 10% 6%
1	B	449	86% 7% 7%
1	C	449	% 86% 7% • 6%
1	D	449	% 85% 7% • 7%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 14732 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 Alpha-amylase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	422	Total	C	N	O	S	0	0	0
			3361	2155	563	632	11			
1	B	418	Total	C	N	O	S	0	0	0
			3338	2142	558	627	11			
1	C	422	Total	C	N	O	S	0	0	0
			3364	2156	563	634	11			
1	D	418	Total	C	N	O	S	0	0	0
			3338	2142	558	627	11			

There are 44 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	MET	-	INSERTION	UNP Q88ZW5
A	1	SER	-	INSERTION	UNP Q88ZW5
A	2	LEU	-	INSERTION	UNP Q88ZW5
A	441	GLU	-	INSERTION	UNP Q88ZW5
A	442	GLY	-	INSERTION	UNP Q88ZW5
A	443	HIS	-	INSERTION	UNP Q88ZW5
A	444	HIS	-	INSERTION	UNP Q88ZW5
A	445	HIS	-	INSERTION	UNP Q88ZW5
A	446	HIS	-	INSERTION	UNP Q88ZW5
A	447	HIS	-	INSERTION	UNP Q88ZW5
A	448	HIS	-	INSERTION	UNP Q88ZW5
B	0	MET	-	INSERTION	UNP Q88ZW5
B	1	SER	-	INSERTION	UNP Q88ZW5
B	2	LEU	-	INSERTION	UNP Q88ZW5
B	441	GLU	-	INSERTION	UNP Q88ZW5
B	442	GLY	-	INSERTION	UNP Q88ZW5
B	443	HIS	-	INSERTION	UNP Q88ZW5
B	444	HIS	-	INSERTION	UNP Q88ZW5
B	445	HIS	-	INSERTION	UNP Q88ZW5
B	446	HIS	-	INSERTION	UNP Q88ZW5
B	447	HIS	-	INSERTION	UNP Q88ZW5

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Chain	Residue	Modelled	Actual	Comment	Reference
B	448	HIS	-	INSERTION	UNP Q88ZW5
C	0	MET	-	INSERTION	UNP Q88ZW5
C	1	SER	-	INSERTION	UNP Q88ZW5
C	2	LEU	-	INSERTION	UNP Q88ZW5
C	441	GLU	-	INSERTION	UNP Q88ZW5
C	442	GLY	-	INSERTION	UNP Q88ZW5
C	443	HIS	-	INSERTION	UNP Q88ZW5
C	444	HIS	-	INSERTION	UNP Q88ZW5
C	445	HIS	-	INSERTION	UNP Q88ZW5
C	446	HIS	-	INSERTION	UNP Q88ZW5
C	447	HIS	-	INSERTION	UNP Q88ZW5
C	448	HIS	-	INSERTION	UNP Q88ZW5
D	0	MET	-	INSERTION	UNP Q88ZW5
D	1	SER	-	INSERTION	UNP Q88ZW5
D	2	LEU	-	INSERTION	UNP Q88ZW5
D	441	GLU	-	INSERTION	UNP Q88ZW5
D	442	GLY	-	INSERTION	UNP Q88ZW5
D	443	HIS	-	INSERTION	UNP Q88ZW5
D	444	HIS	-	INSERTION	UNP Q88ZW5
D	445	HIS	-	INSERTION	UNP Q88ZW5
D	446	HIS	-	INSERTION	UNP Q88ZW5
D	447	HIS	-	INSERTION	UNP Q88ZW5
D	448	HIS	-	INSERTION	UNP Q88ZW5

- Molecule 2 is water.

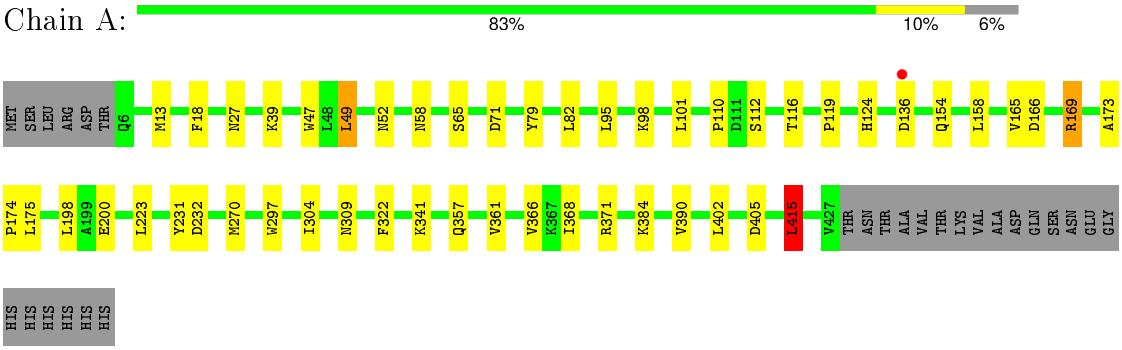
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	340	Total O 340 340	0	0
2	B	314	Total O 314 314	0	0
2	C	357	Total O 357 357	0	0
2	D	320	Total O 320 320	0	0

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 errors 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.

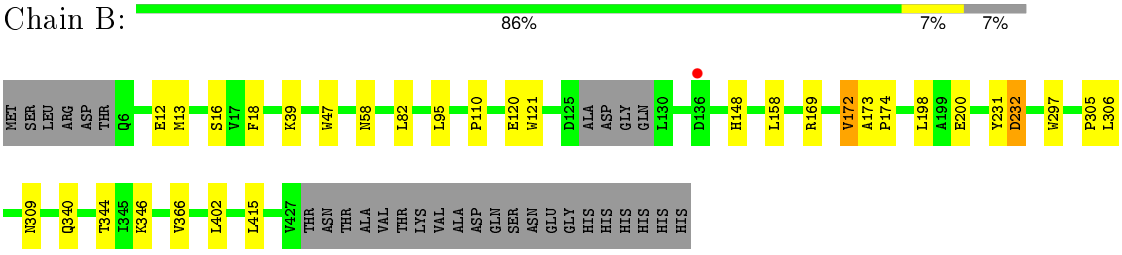
- Molecule 1: Alpha-amylase

Chain A:



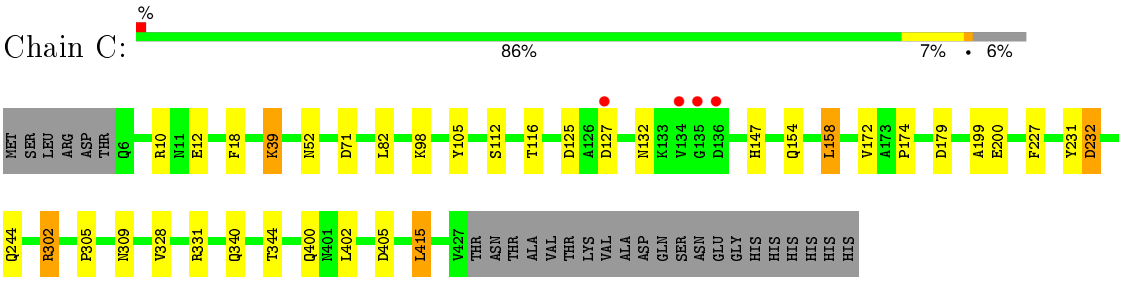
- Molecule 1: Alpha-amylase

Chain B:



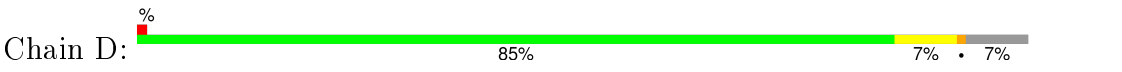
- Molecule 1: Alpha-amylase

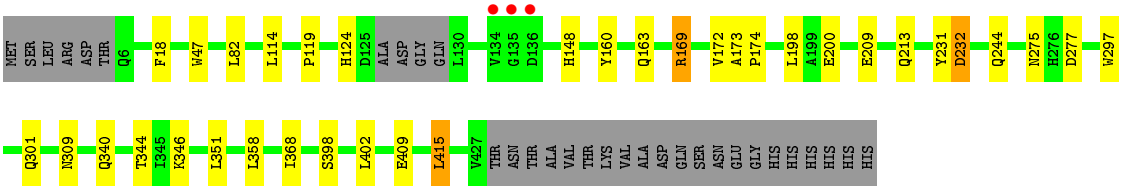
Chain C:



- Molecule 1: Alpha-amylase

Chain D:




MET SER LEU ARG ASP THR Q6 F18 W47 L82 L114 P119 H124 D125 ALA ASP GLY GLN L130 V134 G135 D136 H148 Y160 Q163 R169 V172 A173 P174 L198 A199 E200 E209 Q213 Y231 D232 Q244 N275 H276 D277 W297
Q301 N309 Q340 T344 I345 K346 L351 L358 I368 S398 L402 E409 L415 V427 THR ASN THR ALA VAL THR LYS VAL ALA ASP GLN SER ASN GLU GLY HIS HIS HIS HIS HIS HIS

4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	202.28Å 63.77Å 155.68Å 90.00° 102.93° 90.00°	Depositor
Resolution (Å)	20.00 – 2.00 24.41 – 2.00	Depositor EDS
% Data completeness (in resolution range)	98.6 (20.00-2.00) 98.6 (24.41-2.00)	Depositor EDS
R_{merge}	0.18	Depositor
R_{sym}	0.18	Depositor
$< I/\sigma(I) >$ [1]	3.10 (at 1.99Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.200 , 0.251 0.205 , 0.254	Depositor DCC
R_{free} test set	6497 reflections (5.30%)	DCC
Wilson B-factor (Å^2)	12.4	Xtriage
Anisotropy	0.489	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.39 , 44.5	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Outliers	1 of 129240 reflections (0.001%)	Xtriage
F_o,F_c correlation	0.92	EDS
Total number of atoms	14732	wwPDB-VP
Average B, all atoms (Å^2)	17.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.71% 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, 0.375 respectively for untwinned datasets, and 0.333, 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	0.65	0/3446	0.73	4/4695 (0.1%)
1	B	0.66	0/3422	0.71	2/4661 (0.0%)
1	C	0.68	0/3449	0.72	2/4699 (0.0%)
1	D	0.64	0/3422	0.71	1/4661 (0.0%)
All	All	0.66	0/13739	0.72	9/18716 (0.0%)

There are no bond length outliers.

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	415	LEU	CA-CB-CG	8.58	135.03	115.30
1	A	169	ARG	NE-CZ-NH1	8.33	124.47	120.30
1	D	415	LEU	CA-CB-CG	7.83	133.31	115.30
1	C	415	LEU	CA-CB-CG	7.09	131.60	115.30
1	B	415	LEU	CA-CB-CG	6.65	130.59	115.30
1	A	169	ARG	NE-CZ-NH2	-6.45	117.07	120.30
1	B	13	MET	CG-SD-CE	-6.27	90.17	100.20
1	C	302	ARG	NE-CZ-NH1	6.11	123.36	120.30
1	A	13	MET	CG-SD-CE	-5.37	91.61	100.20

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	3361	0	3243	31	0
1	B	3338	0	3224	17	0
1	C	3364	0	3245	21	0
1	D	3338	0	3224	19	0
2	A	340	0	0	4	0
2	B	314	0	0	1	0
2	C	357	0	0	6	0
2	D	320	0	0	2	0
All	All	14732	0	12936	88	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 3.

All (88) 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:D:244:GLN:NE2	2:D:552:HOH:O	2.04	0.91
1:D:47:TRP:CZ2	1:D:169:ARG:HD3	2.07	0.90
1:B:169:ARG:HG3	1:B:198:LEU:HD23	1.57	0.86
1:A:47:TRP:CZ2	1:A:169:ARG:HD3	2.13	0.83
1:C:112:SER:O	1:C:116:THR:HG23	1.78	0.81
1:C:244:GLN:NE2	2:C:567:HOH:O	2.05	0.79
1:D:169:ARG:HG3	1:D:198:LEU:HD23	1.66	0.78
1:A:169:ARG:HG3	1:A:198:LEU:HD23	1.69	0.73
1:B:47:TRP:CZ2	1:B:169:ARG:HD3	2.23	0.72
1:C:125:ASP:OD1	1:C:127:ASP:OD1	2.08	0.71
1:C:328:VAL:O	1:C:331:ARG:NH1	2.23	0.71
1:D:148:HIS:HB2	2:D:736:HOH:O	1.95	0.65
1:D:200:GLU:HA	1:D:231:TYR:CD1	2.33	0.64
1:A:169:ARG:HG3	1:A:198:LEU:CD2	2.30	0.61
1:D:398:SER:HB2	1:D:409:GLU:OE2	2.00	0.61
1:D:160:TYR:O	1:D:163:GLN:HG3	2.01	0.61
1:C:340:GLN:O	1:C:344:THR:HG23	2.01	0.60
1:D:119:PRO:O	1:D:124:HIS:HE1	1.86	0.58
1:A:101:LEU:HG	1:A:165:VAL:HG11	1.86	0.58
1:A:58:ASN:HD22	1:A:110:PRO:HG2	1.70	0.56
1:C:39:LYS:C	1:C:39:LYS:HD3	2.26	0.56
1:B:169:ARG:HG3	1:B:198:LEU:CD2	2.33	0.56
1:D:47:TRP:CE2	1:D:169:ARG:HD3	2.41	0.56
1:C:200:GLU:HA	1:C:231:TYR:CD1	2.41	0.55
1:A:341:LYS:NZ	1:A:405:ASP:OD2	2.36	0.54
1:B:148:HIS:O	2:B:609:HOH:O	2.18	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:39:LYS:HE3	1:A:95:LEU:O	2.08	0.54
1:B:39:LYS:HE3	1:B:95:LEU:O	2.09	0.53
1:C:344:THR:HG21	2:C:621:HOH:O	2.09	0.53
1:D:209:GLU:O	1:D:213:GLN:HG2	2.10	0.52
1:C:179:ASP:HB3	2:C:753:HOH:O	2.09	0.52
1:A:119:PRO:O	1:A:124:HIS:HE1	1.93	0.51
1:D:169:ARG:HG3	1:D:198:LEU:CD2	2.40	0.50
1:A:270:MET:HA	1:A:304:ILE:O	2.12	0.49
1:A:112:SER:O	1:A:116:THR:HG23	2.12	0.49
1:A:98:LYS:HB3	1:A:166:ASP:HB2	1.95	0.48
1:A:384:LYS:NZ	2:A:782:HOH:O	2.41	0.48
1:A:174:PRO:HG3	1:A:223:LEU:HD11	1.94	0.48
1:A:173:ALA:N	1:A:174:PRO:CD	2.77	0.48
1:B:305:PRO:HD3	1:B:346:LYS:HE2	1.94	0.48
1:B:340:GLN:O	1:B:344:THR:HG23	2.14	0.48
1:A:175:LEU:HD12	1:A:175:LEU:N	2.28	0.48
1:C:132:ASN:C	2:C:778:HOH:O	2.52	0.47
1:A:27:ASN:HA	1:A:79:TYR:HA	1.95	0.47
1:D:173:ALA:N	1:D:174:PRO:HD2	2.28	0.47
1:B:173:ALA:N	1:B:174:PRO:CD	2.77	0.47
1:B:47:TRP:CE2	1:B:169:ARG:HD3	2.50	0.47
1:C:98:LYS:HD3	2:C:517:HOH:O	2.14	0.47
1:B:39:LYS:HD3	1:B:39:LYS:O	2.15	0.46
1:A:65:SER:HB2	1:A:322:PHE:CE1	2.50	0.46
1:A:58:ASN:ND2	1:A:110:PRO:HG2	2.31	0.45
1:A:297:TRP:CZ3	1:A:366:VAL:HG11	2.52	0.45
1:C:147:HIS:HE1	2:C:541:HOH:O	1.99	0.45
1:B:58:ASN:HD22	1:B:110:PRO:HG2	1.81	0.45
1:B:172:VAL:C	1:B:174:PRO:HD2	2.37	0.45
1:A:371:ARG:NH1	2:A:724:HOH:O	2.47	0.45
1:A:47:TRP:CE2	1:A:169:ARG:HD3	2.50	0.44
1:A:47:TRP:CH2	1:A:169:ARG:HD3	2.52	0.44
1:A:49:LEU:HB2	2:A:464:HOH:O	2.18	0.44
1:A:119:PRO:O	1:A:124:HIS:CE1	2.71	0.44
1:D:114:LEU:HA	1:D:114:LEU:HD23	1.64	0.44
1:B:16:SER:HB2	1:B:306:LEU:HD11	2.00	0.44
1:A:357:GLN:HG2	2:A:457:HOH:O	2.18	0.44
1:D:358:LEU:HD23	1:D:368:ILE:HG12	1.99	0.43
1:C:172:VAL:C	1:C:174:PRO:HD2	2.38	0.43
1:B:200:GLU:HA	1:B:231:TYR:CD1	2.52	0.43
1:D:172:VAL:C	1:D:174:PRO:HD2	2.38	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:39:LYS:O	1:C:39:LYS:HD3	2.18	0.43
1:D:340:GLN:O	1:D:344:THR:HG23	2.17	0.43
1:C:12:GLU:O	1:C:305:PRO:HD2	2.18	0.43
1:B:120:GLU:HG2	1:B:121:TRP:CD1	2.54	0.42
1:C:10:ARG:CZ	1:C:302:ARG:HD3	2.49	0.42
1:D:297:TRP:O	1:D:301:GLN:HG3	2.19	0.42
1:A:390:VAL:CG2	1:A:415:LEU:HD22	2.50	0.42
1:A:297:TRP:CZ3	1:A:368:ILE:HD11	2.55	0.41
1:C:105:TYR:CE2	1:C:154:GLN:HB3	2.55	0.41
1:B:12:GLU:O	1:B:305:PRO:HD2	2.20	0.41
1:D:346:LYS:HA	1:D:351:LEU:HG	2.03	0.41
1:C:400:GLN:HE21	1:C:405:ASP:HA	1.84	0.41
1:A:200:GLU:HA	1:A:231:TYR:CD1	2.55	0.41
1:C:52:ASN:HB3	1:C:71:ASP:O	2.21	0.41
1:A:52:ASN:HB3	1:A:71:ASP:O	2.21	0.41
1:A:154:GLN:OE1	1:A:154:GLN:HA	2.21	0.41
1:A:175:LEU:CD1	1:A:175:LEU:N	2.84	0.40
1:D:275:ASN:OD1	1:D:277:ASP:HB2	2.22	0.40
1:C:105:TYR:OH	1:C:158:LEU:HD13	2.21	0.40
1:C:199:ALA:HB2	1:C:227:PHE:CG	2.56	0.40
1:B:297:TRP:CZ3	1:B:366:VAL:HG11	2.57	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	420/449 (94%)	407 (97%)	12 (3%)	1 (0%)	52	48
1	B	414/449 (92%)	402 (97%)	10 (2%)	2 (0%)	34	26
1	C	420/449 (94%)	406 (97%)	13 (3%)	1 (0%)	52	48
1	D	414/449 (92%)	401 (97%)	12 (3%)	1 (0%)	52	48

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	1668/1796 (93%)	1616 (97%)	47 (3%)	5 (0%)	46 41

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	232	ASP
1	D	232	ASP
1	C	232	ASP
1	A	361	VAL
1	B	172	VAL

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	352/384 (92%)	343 (97%)	9 (3%)	54 54
1	B	351/384 (91%)	345 (98%)	6 (2%)	68 71
1	C	353/384 (92%)	345 (98%)	8 (2%)	58 60
1	D	351/384 (91%)	344 (98%)	7 (2%)	63 65
All	All	1407/1536 (92%)	1377 (98%)	30 (2%)	61 63

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	18	PHE
1	A	49	LEU
1	A	82	LEU
1	A	136	ASP
1	A	158	LEU
1	A	232	ASP
1	A	309	ASN
1	A	402	LEU
1	A	415	LEU
1	B	18	PHE

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Mol	Chain	Res	Type
1	B	82	LEU
1	B	158	LEU
1	B	232	ASP
1	B	309	ASN
1	B	402	LEU
1	C	18	PHE
1	C	39	LYS
1	C	82	LEU
1	C	158	LEU
1	C	232	ASP
1	C	309	ASN
1	C	402	LEU
1	C	415	LEU
1	D	18	PHE
1	D	82	LEU
1	D	169	ARG
1	D	232	ASP
1	D	309	ASN
1	D	402	LEU
1	D	415	LEU

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

Mol	Chain	Res	Type
1	A	8	GLN
1	A	58	ASN
1	A	124	HIS
1	A	244	GLN
1	B	8	GLN
1	B	58	ASN
1	B	244	GLN
1	C	118	HIS
1	C	124	HIS
1	C	132	ASN
1	C	147	HIS
1	C	400	GLN
1	D	118	HIS
1	D	124	HIS
1	D	132	ASN
1	D	244	GLN
1	D	278	ASN

5.3.3 RNA

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains

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

5.5 Carbohydrates

There are no carbohydrates in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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, 95^{th} 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	422/449 (93%)	-0.34	1 (0%)	95	95	7, 15, 27, 36	0
1	B	418/449 (93%)	-0.33	1 (0%)	95	95	8, 15, 27, 33	0
1	C	422/449 (93%)	-0.32	4 (0%)	85	86	7, 14, 26, 41	0
1	D	418/449 (93%)	-0.25	3 (0%)	89	89	8, 16, 27, 48	0
All	All	1680/1796 (93%)	-0.31	9 (0%)	91	92	7, 15, 27, 48	0

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	134	VAL	5.5
1	C	134	VAL	5.0
1	D	136	ASP	4.7
1	D	135	GLY	4.5
1	C	136	ASP	3.9
1	B	136	ASP	3.8
1	C	135	GLY	3.7
1	A	136	ASP	2.8
1	C	127	ASP	2.2

6.2 Non-standard residues in protein, DNA, RNA chains

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

6.3 Carbohydrates

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.