



Full wwPDB X-ray Structure Validation Report ⓘ

Dec 8, 2023 – 01:09 am GMT

PDB ID : 2W37
Title : CRYSTAL STRUCTURE OF THE HEXAMERIC CATABOLIC ORNITHINE TRANSCARBAMYLASE FROM *Lactobacillus hilgardii*
Authors : de las Rivas, B.; Fox, G.C.; Angulo, I.; Rodriguez, H.; Munoz, R.; Mancheno, J.M.
Deposited on : 2008-11-06
Resolution : 2.10 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

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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 : 2.36
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

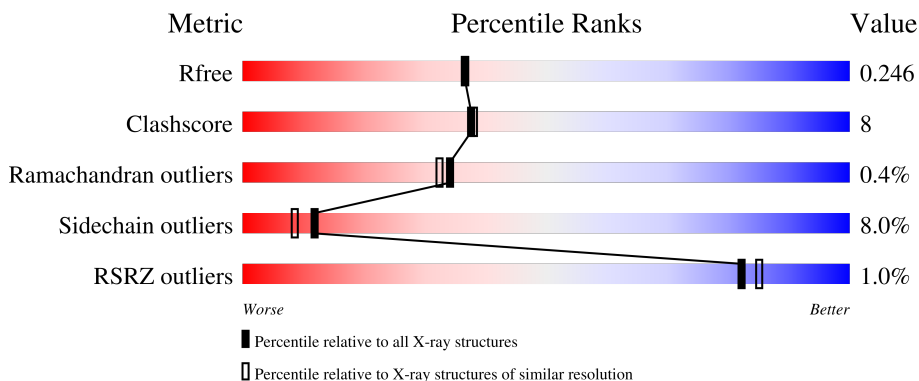
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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	5197 (2.10-2.10)
Clashscore	141614	5710 (2.10-2.10)
Ramachandran outliers	138981	5647 (2.10-2.10)
Sidechain outliers	138945	5648 (2.10-2.10)
RSRZ outliers	127900	5083 (2.10-2.10)

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	359	76% 17% • 5%
1	B	359	74% 18% • 5%
1	C	359	80% 13% • 5%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 8365 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 ORNITHINE CARBAMOYLTRANSFERASE, CATABOLIC.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	341	2677	1697	450	517	13	0	0	0
1	B	341	2677	1697	450	517	13	0	0	0
1	C	341	2677	1697	450	517	13	0	0	0

- Molecule 2 is NICKEL (II) ION (three-letter code: NI) (formula: Ni).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Ni	0	0
			1	1		

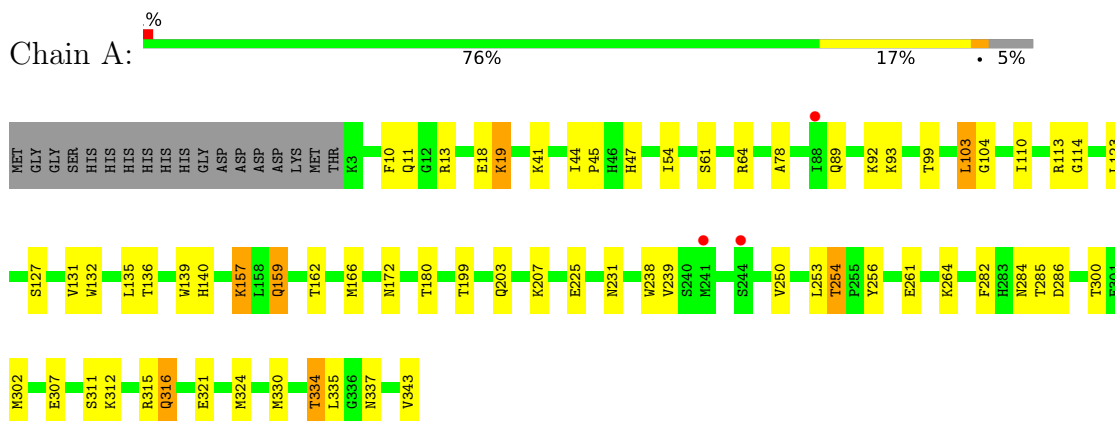
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	125	Total	O	0	0
			125	125		
3	B	110	Total	O	0	0
			110	110		
3	C	98	Total	O	0	0
			98	98		

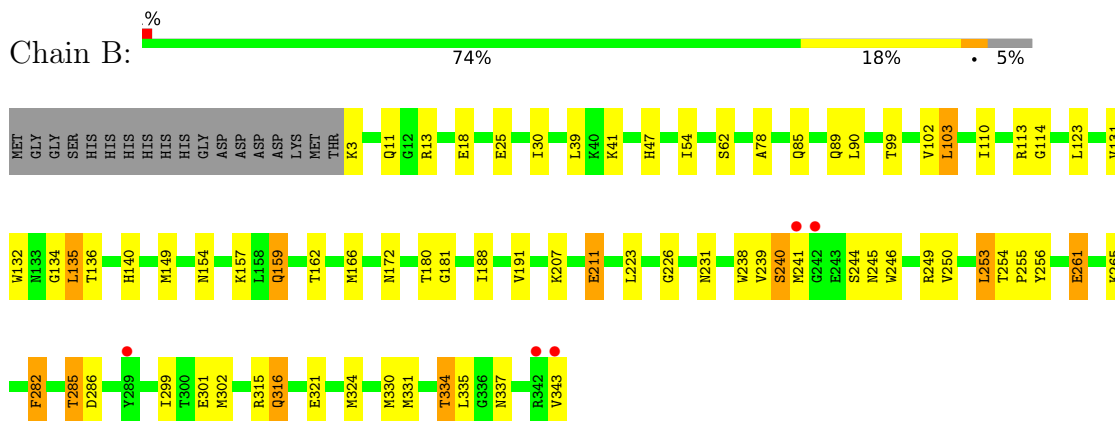
3 Residue-property plots [i](#)

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.

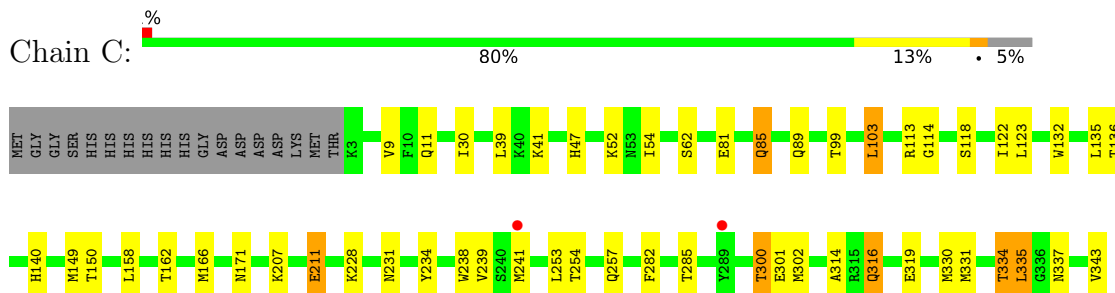
- Molecule 1: ORNITHINE CARBAMOYLTRANSFERASE, CATABOLIC



- Molecule 1: ORNITHINE CARBAMOYLTRANSFERASE, CATABOLIC



- Molecule 1: ORNITHINE CARBAMOYLTRANSFERASE, CATABOLIC



4 Data and refinement statistics i

Property	Value	Source
Space group	P 3 2 1	Depositor
Cell constants a, b, c, α , β , γ	156.81Å 156.81Å 80.62Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	69.34 – 2.10 69.33 – 2.10	Depositor EDS
% Data completeness (in resolution range)	95.7 (69.34-2.10) 82.1 (69.33-2.10)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.80 (at 2.10Å)	Xtrriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.196 , 0.244 0.198 , 0.246	Depositor DCC
R_{free} test set	2783 reflections (5.09%)	wwPDB-VP
Wilson B-factor (Å ²)	32.1	Xtrriage
Anisotropy	0.050	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 47.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	0.063 for -h,-k,l	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	8365	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.87% 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](#)

Bond lengths and bond angles in the following residue types are not validated in this section:
NI

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	0.57	0/2730	0.67	1/3682 (0.0%)
1	B	0.58	0/2730	0.67	1/3682 (0.0%)
1	C	0.57	0/2730	0.65	0/3682
All	All	0.57	0/8190	0.66	2/11046 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	324	MET	CG-SD-CE	-5.57	91.29	100.20
1	A	324	MET	CG-SD-CE	-5.14	91.97	100.20

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	2677	0	2638	43	0
1	B	2677	0	2638	48	0
1	C	2677	0	2638	32	0
2	A	1	0	0	0	0
3	A	125	0	0	7	0
3	B	110	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	C	98	0	0	2	0
All	All	8365	0	7914	122	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 8.

All (122) 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:132:TRP:HE1	1:B:334:THR:HG21	1.11	1.13
1:A:132:TRP:HE1	1:A:334:THR:HG21	1.05	1.13
1:C:132:TRP:HE1	1:C:334:THR:HG21	1.09	1.08
1:A:250:VAL:O	1:A:254:THR:HG22	1.61	1.00
1:A:159:GLN:HE21	1:A:159:GLN:H	1.18	0.88
1:A:132:TRP:NE1	1:A:334:THR:HG21	1.90	0.86
1:B:132:TRP:NE1	1:B:334:THR:HG21	1.92	0.85
1:C:132:TRP:NE1	1:C:334:THR:HG21	1.91	0.83
1:C:314:ALA:HB1	1:C:316:GLN:HE22	1.43	0.83
1:C:54:ILE:HD13	1:C:331:MET:CE	2.13	0.79
1:A:162:THR:H	1:A:231:ASN:HD22	1.32	0.77
1:A:159:GLN:HE21	1:A:159:GLN:N	1.82	0.77
1:B:250:VAL:O	1:B:254:THR:HG23	1.86	0.76
1:C:54:ILE:HD13	1:C:331:MET:HE3	1.66	0.76
1:A:172:ASN:HB2	3:A:2084:HOH:O	1.87	0.75
1:B:246:TRP:CZ3	1:B:250:VAL:HG21	2.27	0.69
1:A:132:TRP:HE1	1:A:334:THR:CG2	1.96	0.69
1:C:314:ALA:CB	1:C:316:GLN:HE22	2.05	0.69
1:C:330:MET:O	1:C:334:THR:HB	1.92	0.68
1:A:18:GLU:HG3	1:A:180:THR:OG1	1.95	0.67
1:B:162:THR:H	1:B:231:ASN:HD22	1.44	0.66
1:B:330:MET:O	1:B:334:THR:HB	1.95	0.65
1:A:159:GLN:H	1:A:159:GLN:NE2	1.91	0.65
1:A:203:GLN:O	1:A:207:LYS:HG3	1.96	0.65
1:A:54:ILE:HD12	1:A:78:ALA:HB1	1.78	0.65
1:C:118:SER:O	1:C:122:ILE:HG13	1.97	0.65
1:B:47:HIS:HE1	3:B:2044:HOH:O	1.81	0.62
1:B:261:GLU:OE2	1:B:265:LYS:HE2	1.99	0.62
1:C:47:HIS:HE1	3:C:2031:HOH:O	1.84	0.61
1:A:330:MET:O	1:A:334:THR:HB	2.01	0.61
1:A:250:VAL:O	1:A:254:THR:CG2	2.44	0.60
1:C:239:VAL:HG22	1:C:302:MET:HE1	1.83	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:99:THR:HG22	1:A:103:LEU:HD22	1.84	0.59
1:B:249:ARG:O	1:B:253:LEU:HB2	2.03	0.58
1:C:52:LYS:HE2	1:C:335:LEU:HD22	1.85	0.58
1:B:54:ILE:HD13	1:B:331:MET:CE	2.33	0.58
1:A:157:LYS:C	3:A:2082:HOH:O	2.41	0.58
1:C:207:LYS:O	1:C:211:GLU:HG2	2.04	0.57
1:C:238:TRP:HE3	1:C:302:MET:HE3	1.69	0.57
1:C:300:THR:HG23	1:C:301:GLU:HG3	1.87	0.57
1:B:99:THR:HG22	1:B:103:LEU:HD22	1.87	0.57
1:B:114:GLY:O	1:B:136:THR:HA	2.06	0.56
1:A:316:GLN:H	1:A:316:GLN:NE2	2.04	0.56
1:C:162:THR:H	1:C:231:ASN:HD22	1.53	0.56
1:A:11:GLN:HE22	1:A:337:ASN:H	1.55	0.55
1:B:223:LEU:HD11	1:B:255:PRO:O	2.05	0.55
1:A:157:LYS:N	3:A:2082:HOH:O	2.39	0.55
1:B:316:GLN:NE2	1:B:316:GLN:H	2.05	0.55
1:C:99:THR:HG22	1:C:103:LEU:HD22	1.89	0.54
1:B:282:PHE:C	1:B:282:PHE:HD2	2.11	0.54
1:A:239:VAL:HG22	1:A:302:MET:HE1	1.90	0.53
1:B:166:MET:HG2	1:B:256:TYR:O	2.08	0.53
1:B:282:PHE:C	1:B:282:PHE:CD2	2.82	0.53
1:C:314:ALA:HB1	1:C:316:GLN:NE2	2.20	0.52
1:B:238:TRP:HE3	1:B:302:MET:CE	2.22	0.52
1:A:11:GLN:NE2	1:A:337:ASN:H	2.07	0.52
1:B:207:LYS:O	1:B:211:GLU:CG	2.58	0.52
1:B:30:ILE:HG23	1:B:149:MET:HA	1.92	0.52
1:C:316:GLN:NE2	1:C:316:GLN:H	2.08	0.51
1:A:238:TRP:CH2	1:A:254:THR:HA	2.45	0.50
3:B:2043:HOH:O	1:C:81:GLU:HG2	2.12	0.50
1:B:11:GLN:NE2	1:B:337:ASN:H	2.10	0.50
1:B:207:LYS:O	1:B:211:GLU:HG2	2.12	0.49
1:B:154:ASN:OD1	1:B:315:ARG:HD2	2.13	0.49
1:B:240:SER:O	1:B:241:MET:HG2	2.13	0.49
1:B:18:GLU:HG3	1:B:180:THR:OG1	2.13	0.48
1:A:13:ARG:HD3	3:A:2004:HOH:O	2.14	0.48
1:B:159:GLN:HE21	1:B:159:GLN:H	1.61	0.48
1:C:257:GLN:HG3	1:C:302:MET:O	2.13	0.48
1:B:181:GLY:HA3	1:B:188:ILE:HD11	1.96	0.48
1:A:199:THR:O	1:A:203:GLN:HG3	2.14	0.47
1:B:11:GLN:HE22	1:B:337:ASN:H	1.62	0.47
1:A:312:LYS:O	1:A:315:ARG:NH2	2.47	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:41:LYS:NZ	1:A:321:GLU:OE2	2.45	0.47
1:B:54:ILE:HD13	1:B:331:MET:HE3	1.96	0.47
1:C:30:ILE:HG23	1:C:149:MET:HA	1.96	0.47
1:A:47:HIS:HE1	3:A:2053:HOH:O	1.96	0.47
1:C:238:TRP:HB2	1:C:302:MET:HE2	1.97	0.46
1:C:238:TRP:CE3	1:C:302:MET:HE3	2.50	0.46
1:A:166:MET:HG2	1:A:256:TYR:O	2.16	0.46
1:A:284:ASN:OD1	1:A:286:ASP:HB2	2.16	0.46
1:C:54:ILE:CD1	1:C:331:MET:CE	2.89	0.46
1:C:9:VAL:HG21	1:C:343:VAL:HB	1.97	0.45
1:B:13:ARG:HH22	1:B:25:GLU:CD	2.19	0.45
1:B:299:ILE:HG22	1:B:301:GLU:O	2.16	0.45
1:B:254:THR:N	1:B:255:PRO:HD2	2.31	0.45
1:B:261:GLU:O	1:B:265:LYS:HG3	2.17	0.45
1:A:316:GLN:H	1:A:316:GLN:HE21	1.63	0.45
1:B:25:GLU:HA	1:B:343:VAL:HG21	1.99	0.45
1:B:238:TRP:CD1	1:B:253:LEU:HG	2.52	0.45
1:B:191:VAL:HG22	1:B:226:GLY:HA3	1.98	0.45
1:C:114:GLY:O	1:C:136:THR:HA	2.17	0.45
1:A:64:ARG:NE	1:B:102:VAL:HG11	2.32	0.44
1:B:41:LYS:NZ	1:B:321:GLU:OE2	2.50	0.44
1:B:285:THR:CG2	1:B:285:THR:O	2.65	0.44
1:C:11:GLN:NE2	1:C:337:ASN:H	2.16	0.44
1:B:159:GLN:H	1:B:159:GLN:NE2	2.17	0.43
1:B:207:LYS:O	1:B:211:GLU:HG3	2.18	0.43
1:B:238:TRP:CE3	1:B:302:MET:CE	3.00	0.43
1:B:238:TRP:HE3	1:B:302:MET:HE3	1.83	0.43
1:A:114:GLY:O	1:A:136:THR:HA	2.18	0.43
1:B:54:ILE:HD12	1:B:78:ALA:HB1	2.01	0.43
1:C:11:GLN:HE22	1:C:337:ASN:H	1.66	0.43
1:B:245:ASN:O	1:B:249:ARG:HG3	2.18	0.43
1:A:307:GLU:O	1:A:311:SER:HB3	2.18	0.42
1:A:64:ARG:HD2	3:A:2105:HOH:O	2.19	0.42
1:B:110:ILE:O	1:B:131:VAL:HA	2.19	0.42
1:A:104:GLY:HA3	1:A:127:SER:O	2.20	0.42
1:A:110:ILE:O	1:A:131:VAL:HA	2.20	0.42
1:C:158:LEU:HA	3:C:2054:HOH:O	2.19	0.42
1:A:238:TRP:HE3	1:A:302:MET:HE3	1.85	0.41
1:B:134:GLY:O	1:B:135:LEU:CB	2.68	0.41
1:A:44:ILE:HA	1:A:45:PRO:HD3	1.94	0.41
1:A:238:TRP:CE3	1:A:302:MET:HE3	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:85:GLN:H	1:C:85:GLN:HE21	1.67	0.41
1:C:150:THR:HG21	1:C:319:GLU:HB2	2.03	0.41
1:C:238:TRP:CH2	1:C:254:THR:HA	2.56	0.41
1:A:54:ILE:HD12	1:A:78:ALA:CB	2.50	0.41
1:A:19:LYS:HG2	1:A:139:TRP:CH2	2.56	0.41
1:A:172:ASN:CB	3:A:2084:HOH:O	2.58	0.41
1:B:238:TRP:CE3	1:B:302:MET:HE2	2.56	0.40
1:A:10:PHE:O	1:A:13:ARG:HB2	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	339/359 (94%)	326 (96%)	12 (4%)	1 (0%)	41	41
1	B	339/359 (94%)	323 (95%)	14 (4%)	2 (1%)	25	21
1	C	339/359 (94%)	327 (96%)	11 (3%)	1 (0%)	41	41
All	All	1017/1077 (94%)	976 (96%)	37 (4%)	4 (0%)	34	32

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	135	LEU
1	B	135	LEU
1	C	135	LEU
1	B	157	LYS

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	287/302 (95%)	264 (92%)	23 (8%)	12	8
1	B	287/302 (95%)	263 (92%)	24 (8%)	11	7
1	C	287/302 (95%)	265 (92%)	22 (8%)	13	9
All	All	861/906 (95%)	792 (92%)	69 (8%)	12	8

All (69) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	19	LYS
1	A	61	SER
1	A	89	GLN
1	A	92	LYS
1	A	93	LYS
1	A	103	LEU
1	A	113	ARG
1	A	123	LEU
1	A	140	HIS
1	A	157	LYS
1	A	159	GLN
1	A	225	GLU
1	A	253	LEU
1	A	254	THR
1	A	261	GLU
1	A	264	LYS
1	A	282	PHE
1	A	285	THR
1	A	300	THR
1	A	316	GLN
1	A	334	THR
1	A	335	LEU
1	A	343	VAL
1	B	3	LYS
1	B	39	LEU
1	B	62	SER

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Mol	Chain	Res	Type
1	B	85	GLN
1	B	89	GLN
1	B	90	LEU
1	B	103	LEU
1	B	113	ARG
1	B	123	LEU
1	B	140	HIS
1	B	159	GLN
1	B	172	ASN
1	B	211	GLU
1	B	239	VAL
1	B	240	SER
1	B	244	SER
1	B	253	LEU
1	B	261	GLU
1	B	282	PHE
1	B	285	THR
1	B	286	ASP
1	B	316	GLN
1	B	334	THR
1	B	335	LEU
1	C	39	LEU
1	C	41	LYS
1	C	62	SER
1	C	85	GLN
1	C	89	GLN
1	C	103	LEU
1	C	113	ARG
1	C	123	LEU
1	C	140	HIS
1	C	166	MET
1	C	171	ASN
1	C	211	GLU
1	C	228	LYS
1	C	234	TYR
1	C	241	MET
1	C	253	LEU
1	C	282	PHE
1	C	285	THR
1	C	300	THR
1	C	316	GLN
1	C	334	THR

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Mol	Chain	Res	Type
1	C	335	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	11	GLN
1	A	47	HIS
1	A	159	GLN
1	A	203	GLN
1	A	231	ASN
1	A	316	GLN
1	B	11	GLN
1	B	47	HIS
1	B	159	GLN
1	B	203	GLN
1	B	231	ASN
1	B	272	GLN
1	B	316	GLN
1	C	11	GLN
1	C	47	HIS
1	C	85	GLN
1	C	89	GLN
1	C	231	ASN
1	C	316	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

Of 1 ligands modelled in this entry, 1 is monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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 [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	341/359 (94%)	-0.13	3 (0%) 84 86	19, 30, 47, 64	0
1	B	341/359 (94%)	-0.10	5 (1%) 73 77	19, 31, 48, 64	0
1	C	341/359 (94%)	-0.19	2 (0%) 89 91	19, 31, 47, 62	0
All	All	1023/1077 (94%)	-0.14	10 (0%) 82 85	19, 31, 47, 64	0

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	241	MET	4.5
1	A	241	MET	4.2
1	B	242	GLY	3.8
1	B	241	MET	3.4
1	A	88	ILE	2.9
1	A	244	SER	2.8
1	B	343	VAL	2.8
1	B	342	ARG	2.2
1	B	289	TYR	2.1
1	C	289	TYR	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](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	NI	A	1344	1/1	1.00	0.13	23,23,23,23	0

6.5 Other polymers [i](#)

There are no such residues in this entry.