



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 31, 2022 – 03:25 PM EDT

PDB ID : 8H50
Title : Crystal structure of carboxyspermidine dehydrogenase from *Helicobacter pylori* in space group C2221
Authors : Ko, K.Y.; Park, S.C.; Cho, S.Y.; Yoon, S.I.
Deposited on : 2022-10-11
Resolution : 2.90 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

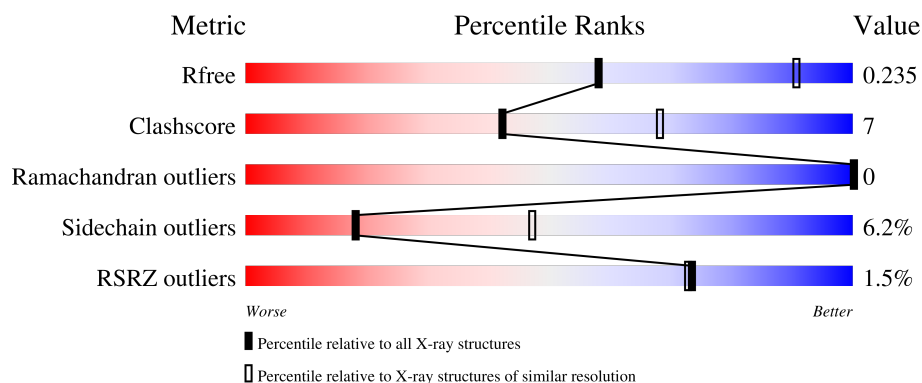
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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	1957 (2.90-2.90)
Clashscore	141614	2172 (2.90-2.90)
Ramachandran outliers	138981	2115 (2.90-2.90)
Sidechain outliers	138945	2117 (2.90-2.90)
RSRZ outliers	127900	1906 (2.90-2.90)

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	405	<div> <div>0%</div> <div> <div></div> <div>81%</div> <div>15%</div> <div>..</div> </div> </div>
1	B	405	<div> <div>0%</div> <div> <div></div> <div>76%</div> <div>18%</div> <div>..</div> </div> </div>
1	C	405	<div> <div>2%</div> <div> <div></div> <div>76%</div> <div>19%</div> <div>..</div> </div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 8970 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 Carboxyspermidine dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	391	Total	C	N	O	S	0	0	0
			2994	1912	500	559	23			
1	B	392	Total	C	N	O	S	0	0	0
			2967	1886	497	563	21			
1	C	391	Total	C	N	O	S	0	0	0
			3005	1920	495	569	21			

There are 21 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-5	GLY	-	expression tag	UNP E8QMS0
A	-4	SER	-	expression tag	UNP E8QMS0
A	-3	HIS	-	expression tag	UNP E8QMS0
A	-2	MET	-	expression tag	UNP E8QMS0
A	-1	ALA	-	expression tag	UNP E8QMS0
A	0	SER	-	expression tag	UNP E8QMS0
A	156	ARG	LYS	conflict	UNP E8QMS0
B	-5	GLY	-	expression tag	UNP E8QMS0
B	-4	SER	-	expression tag	UNP E8QMS0
B	-3	HIS	-	expression tag	UNP E8QMS0
B	-2	MET	-	expression tag	UNP E8QMS0
B	-1	ALA	-	expression tag	UNP E8QMS0
B	0	SER	-	expression tag	UNP E8QMS0
B	156	ARG	LYS	conflict	UNP E8QMS0
C	-5	GLY	-	expression tag	UNP E8QMS0
C	-4	SER	-	expression tag	UNP E8QMS0
C	-3	HIS	-	expression tag	UNP E8QMS0
C	-2	MET	-	expression tag	UNP E8QMS0
C	-1	ALA	-	expression tag	UNP E8QMS0
C	0	SER	-	expression tag	UNP E8QMS0
C	156	ARG	LYS	conflict	UNP E8QMS0

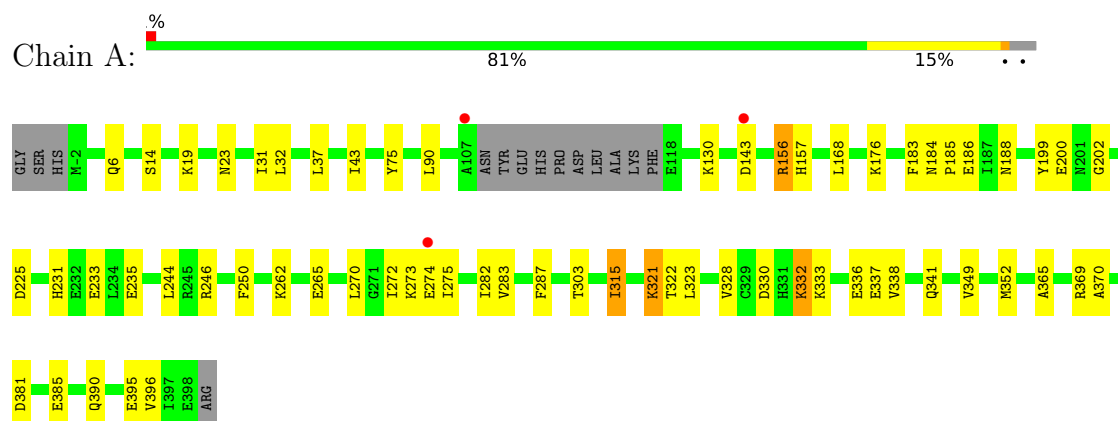
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total 1	O 1	0	0
2	B	1	Total 1	O 1	0	0
2	C	2	Total 2	O 2	0	0

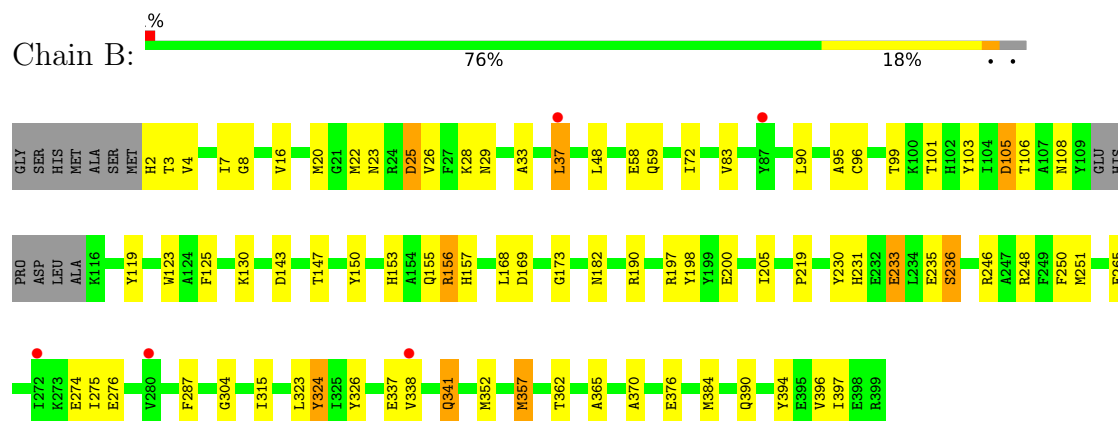
3 Residue-property plots

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.

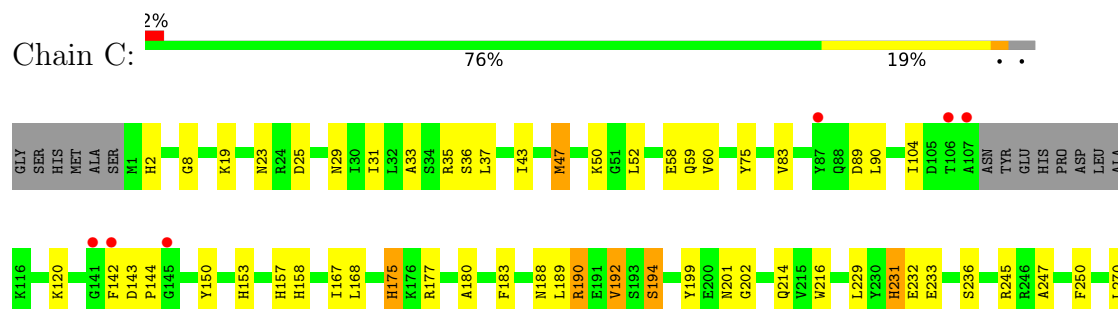
• Molecule 1: Carboxyspermidine dehydrogenase

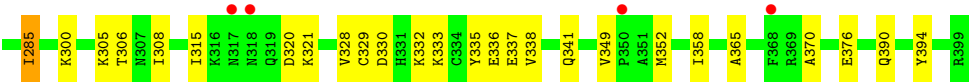


• Molecule 1: Carboxyspermidine dehydrogenase



• Molecule 1: Carboxyspermidine dehydrogenase





4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	152.16Å 160.65Å 162.02Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.90 – 2.90 29.90 – 2.88	Depositor EDS
% Data completeness (in resolution range)	99.8 (29.90-2.90) 98.8 (29.90-2.88)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.79 (at 2.90Å)	Xtriage
Refinement program	PHENIX 1.15.2_3472	Depositor
R, R_{free}	0.195 , 0.234 0.194 , 0.235	Depositor DCC
R_{free} test set	2292 reflections (5.15%)	wwPDB-VP
Wilson B-factor (Å ²)	62.7	Xtriage
Anisotropy	0.897	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 50.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	8970	wwPDB-VP
Average B, all atoms (Å ²)	77.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.30% 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	0.36	0/3058	0.57	0/4157
1	B	0.36	0/3031	0.59	2/4128 (0.0%)
1	C	0.36	0/3067	0.57	0/4165
All	All	0.36	0/9156	0.58	2/12450 (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	105	ASP	CB-CG-OD1	5.35	123.11	118.30
1	B	248	ARG	NE-CZ-NH1	-5.06	117.77	120.30

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	2994	0	2871	34	0
1	B	2967	0	2765	47	0
1	C	3005	0	2889	43	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	2	0	0	0	0
All	All	8970	0	8525	121	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 7.

All (121) 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:144:PRO:HB3	1:C:233:GLU:HB3	1.80	0.62
1:A:176:LYS:NZ	1:A:303:THR:OG1	2.32	0.62
1:B:72:ILE:HG21	1:B:99:THR:HG21	1.83	0.61
1:A:330:ASP:HB3	1:A:333:LYS:HB2	1.82	0.60
1:A:200:GLU:HG3	1:A:246:ARG:HG3	1.83	0.59
1:A:19:LYS:HG3	1:A:349:VAL:HG22	1.85	0.58
1:A:156:ARG:HG3	1:A:157:HIS:ND1	2.18	0.58
1:C:332:LYS:O	1:C:336:GLU:HG3	2.03	0.58
1:B:357:MET:HG3	1:B:362:THR:HB	1.85	0.58
1:A:168:LEU:HD23	1:A:250:PHE:HB2	1.85	0.57
1:A:365:ALA:HB1	1:A:370:ALA:HB2	1.85	0.57
1:B:16:VAL:O	1:B:20:MET:HB2	2.06	0.56
1:A:156:ARG:NH2	1:A:381:ASP:OD1	2.39	0.56
1:B:37:LEU:HD23	1:B:59:GLN:HB2	1.86	0.56
1:B:168:LEU:HD23	1:B:250:PHE:HB2	1.88	0.56
1:B:156:ARG:HD3	1:B:157:HIS:CE1	2.41	0.55
1:B:236:SER:OG	1:B:376:GLU:OE1	2.21	0.55
1:A:287:PHE:CE2	1:B:275:ILE:HD11	2.42	0.55
1:C:142:PHE:CE1	1:C:308:ILE:HB	2.42	0.54
1:C:19:LYS:HG3	1:C:349:VAL:HG22	1.89	0.54
1:C:190:ARG:O	1:C:194:SER:OG	2.26	0.54
1:C:330:ASP:HB3	1:C:333:LYS:HB2	1.89	0.53
1:A:143:ASP:OD2	1:A:143:ASP:N	2.40	0.53
1:B:200:GLU:HG3	1:B:246:ARG:HG3	1.91	0.53
1:A:23:ASN:OD1	1:A:390:GLN:NE2	2.37	0.53
1:C:270:LEU:HA	1:C:285:ILE:HD12	1.91	0.52
1:C:175:HIS:CD2	1:C:180:ALA:HB3	2.44	0.52
1:C:236:SER:OG	1:C:376:GLU:OE1	2.27	0.52
1:C:143:ASP:O	1:C:231:HIS:HE1	1.92	0.51
1:B:23:ASN:OD1	1:B:390:GLN:NE2	2.38	0.51
1:C:365:ALA:HB1	1:C:370:ALA:HB2	1.92	0.51
1:B:99:THR:HG23	1:B:101:THR:OG1	2.09	0.51
1:B:103:TYR:OH	1:B:105:ASP:HB3	2.11	0.50
1:B:143:ASP:OD2	1:B:143:ASP:N	2.44	0.50
1:C:315:ILE:HD13	1:C:320:ASP:HA	1.94	0.50
1:C:104:ILE:HD11	1:C:358:ILE:HD12	1.94	0.49
1:A:130:LYS:HE3	1:A:370:ALA:O	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:323:LEU:HD11	1:A:396:VAL:CG1	2.42	0.49
1:B:143:ASP:HB2	1:B:233:GLU:OE2	2.13	0.49
1:A:184:ASN:OD1	1:A:186:GLU:HG3	2.12	0.49
1:A:274:GLU:HG2	1:A:283:VAL:HG22	1.95	0.49
1:C:300:LYS:HB3	1:C:335:TYR:CZ	2.48	0.49
1:A:185:PRO:HA	1:A:188:ASN:HD22	1.78	0.48
1:B:337:GLU:HG2	1:B:338:VAL:HG13	1.94	0.48
1:B:72:ILE:HD13	1:B:99:THR:HG21	1.96	0.48
1:C:43:ILE:HG22	1:C:47:MET:HE2	1.96	0.48
1:A:381:ASP:O	1:A:385:GLU:HG3	2.13	0.48
1:C:168:LEU:HD13	1:C:216:TRP:CE2	2.49	0.47
1:B:197:ARG:HB3	1:B:230:TYR:CD2	2.49	0.47
1:C:143:ASP:OD2	1:C:143:ASP:N	2.46	0.47
1:A:14:SER:HA	1:A:43:ILE:HD13	1.97	0.47
1:C:188:ASN:O	1:C:192:VAL:HG13	2.15	0.46
1:B:4:VAL:HG21	1:B:20:MET:SD	2.56	0.46
1:C:8:GLY:HA3	1:C:83:VAL:HG23	1.98	0.46
1:C:337:GLU:HG2	1:C:338:VAL:HG13	1.96	0.46
1:A:287:PHE:HE2	1:B:275:ILE:HD11	1.79	0.46
1:B:25:ASP:OD2	1:B:26:VAL:HG23	2.15	0.46
1:A:332:LYS:O	1:A:336:GLU:HG3	2.16	0.45
1:C:232:GLU:H	1:C:232:GLU:CD	2.20	0.45
1:C:25:ASP:OD1	1:C:25:ASP:N	2.46	0.45
1:A:31:ILE:HG12	1:A:75:TYR:CE1	2.52	0.45
1:A:199:TYR:OH	1:A:202:GLY:HA2	2.15	0.45
1:A:321:LYS:HG3	1:A:322:THR:N	2.31	0.45
1:C:305:LYS:HA	1:C:329:CYS:O	2.17	0.45
1:A:337:GLU:HG2	1:A:338:VAL:HG13	1.98	0.44
1:A:275:ILE:HD11	1:B:287:PHE:CZ	2.53	0.44
1:A:315:ILE:HA	1:A:315:ILE:HD13	1.70	0.44
1:C:35:ARG:O	1:C:59:GLN:NE2	2.50	0.44
1:B:147:THR:OG1	1:B:233:GLU:OE1	2.30	0.44
1:C:332:LYS:HD3	1:C:332:LYS:HA	1.73	0.44
1:C:33:ALA:HA	1:C:58:GLU:O	2.17	0.44
1:C:189:LEU:HD13	1:C:270:LEU:HD13	1.99	0.44
1:A:244:LEU:HD23	1:A:244:LEU:HA	1.78	0.44
1:C:50:LYS:HB2	1:C:52:LEU:HG	2.00	0.44
1:B:8:GLY:HA3	1:B:83:VAL:HG23	1.99	0.44
1:B:150:TYR:OH	1:B:394:TYR:HB3	2.18	0.43
1:A:272:ILE:HD12	1:A:273:LYS:N	2.34	0.43
1:B:169:ASP:O	1:B:251:MET:HA	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:153:HIS:CE1	1:C:157:HIS:HD2	2.37	0.43
1:B:119:TYR:HB3	1:B:123:TRP:CE2	2.54	0.43
1:C:89:ASP:OD2	1:C:90:LEU:N	2.51	0.43
1:C:167:ILE:O	1:C:250:PHE:N	2.47	0.43
1:B:275:ILE:HD12	1:B:275:ILE:O	2.18	0.43
1:B:341:GLN:H	1:B:341:GLN:HG3	1.68	0.43
1:C:23:ASN:OD1	1:C:390:GLN:NE2	2.47	0.43
1:C:47:MET:HG3	1:C:52:LEU:HB2	2.01	0.43
1:A:185:PRO:HA	1:A:188:ASN:ND2	2.34	0.43
1:B:219:PRO:HG2	1:B:326:TYR:CE1	2.53	0.43
1:C:150:TYR:OH	1:C:394:TYR:HB3	2.19	0.43
1:B:130:LYS:HG3	1:B:370:ALA:O	2.19	0.42
1:B:150:TYR:HD1	1:B:384:MET:HE1	1.83	0.42
1:B:83:VAL:HG12	1:B:106:THR:CG2	2.48	0.42
1:B:2:HIS:O	1:B:28:LYS:N	2.50	0.42
1:B:324:TYR:HB3	1:B:397:ILE:HB	2.01	0.42
1:C:37:LEU:HD23	1:C:37:LEU:HA	1.86	0.42
1:B:95:ALA:O	1:B:99:THR:HG22	2.19	0.42
1:C:158:HIS:O	1:C:321:LYS:HD3	2.19	0.42
1:A:37:LEU:HD23	1:A:37:LEU:HA	1.84	0.42
1:B:90:LEU:HD22	1:B:125:PHE:HE2	1.84	0.42
1:B:96:CYS:HA	1:B:99:THR:HG22	2.02	0.42
1:C:31:ILE:HG12	1:C:75:TYR:CD1	2.55	0.42
1:C:199:TYR:OH	1:C:202:GLY:HA2	2.19	0.42
1:B:365:ALA:HB1	1:B:370:ALA:CB	2.50	0.42
1:C:59:GLN:HG3	1:C:60:VAL:N	2.35	0.42
1:A:265:GLU:HB2	1:A:270:LEU:HD12	2.02	0.42
1:A:323:LEU:HD11	1:A:396:VAL:HG13	2.00	0.41
1:B:33:ALA:HA	1:B:58:GLU:O	2.19	0.41
1:B:265:GLU:O	1:B:265:GLU:HG2	2.20	0.41
1:C:31:ILE:HG12	1:C:75:TYR:CE1	2.55	0.41
1:B:153:HIS:HB2	1:B:384:MET:HE2	2.02	0.41
1:B:198:TYR:CE1	1:B:205:ILE:HB	2.55	0.41
1:C:199:TYR:HB3	1:C:247:ALA:HB3	2.02	0.41
1:B:3:THR:HA	1:B:29:ASN:O	2.20	0.41
1:C:306:THR:O	1:C:328:VAL:HA	2.21	0.41
1:A:6:GLN:NE2	1:A:32:LEU:HD21	2.36	0.41
1:B:22:MET:HE1	1:B:337:GLU:HG2	2.03	0.40
1:B:173:GLY:HA2	1:B:304:GLY:HA3	2.03	0.40
1:A:19:LYS:HD3	1:A:19:LYS:HA	1.89	0.40
1:B:103:TYR:CZ	1:B:105:ASP:HB3	2.56	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:323:LEU:HD11	1:B:396:VAL:CG1	2.52	0.40
1:C:190:ARG:HH11	1:C:190:ARG:HD3	1.72	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	387/405 (96%)	374 (97%)	13 (3%)	0	100	100
1	B	388/405 (96%)	378 (97%)	10 (3%)	0	100	100
1	C	387/405 (96%)	374 (97%)	13 (3%)	0	100	100
All	All	1162/1215 (96%)	1126 (97%)	36 (3%)	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	306/346 (88%)	289 (94%)	17 (6%)	21	52
1	B	295/346 (85%)	275 (93%)	20 (7%)	16	42
1	C	308/346 (89%)	289 (94%)	19 (6%)	18	47
All	All	909/1038 (88%)	853 (94%)	56 (6%)	18	47

All (56) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	90	LEU
1	A	156	ARG
1	A	183	PHE
1	A	225	ASP
1	A	231	HIS
1	A	233	GLU
1	A	235	GLU
1	A	262	LYS
1	A	282	ILE
1	A	315	ILE
1	A	321	LYS
1	A	328	VAL
1	A	332	LYS
1	A	341	GLN
1	A	352	MET
1	A	369	ARG
1	A	395	GLU
1	B	7	ILE
1	B	25	ASP
1	B	37	LEU
1	B	48	LEU
1	B	108	ASN
1	B	155	GLN
1	B	156	ARG
1	B	182	ASN
1	B	190	ARG
1	B	231	HIS
1	B	233	GLU
1	B	235	GLU
1	B	236	SER
1	B	274	GLU
1	B	276	GLU
1	B	315	ILE
1	B	324	TYR
1	B	341	GLN
1	B	352	MET
1	B	357	MET
1	C	2	HIS
1	C	29	ASN
1	C	36	SER
1	C	47	MET
1	C	120	LYS

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Mol	Chain	Res	Type
1	C	175	HIS
1	C	177	ARG
1	C	183	PHE
1	C	190	ARG
1	C	192	VAL
1	C	194	SER
1	C	201	ASN
1	C	214	GLN
1	C	229	LEU
1	C	231	HIS
1	C	245	ARG
1	C	285	ILE
1	C	341	GLN
1	C	352	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	2	HIS
1	A	188	ASN
1	A	266	ASN
1	B	155	GLN
1	B	266	ASN
1	C	2	HIS
1	C	153	HIS
1	C	266	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 monosaccharides 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 [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	391/405 (96%)	-0.05	3 (0%) 86 86	40, 73, 100, 124	0
1	B	392/405 (96%)	-0.03	5 (1%) 77 77	59, 84, 115, 130	0
1	C	391/405 (96%)	-0.06	10 (2%) 56 52	41, 72, 95, 131	0
All	All	1174/1215 (96%)	-0.05	18 (1%) 73 73	40, 77, 107, 131	0

All (18) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	107	ALA	5.7
1	A	107	ALA	3.9
1	B	37	LEU	3.3
1	C	87	TYR	2.6
1	C	318	ASN	2.6
1	C	141	GLY	2.4
1	C	142	PHE	2.4
1	C	368	PHE	2.4
1	C	350	PRO	2.4
1	B	87	TYR	2.3
1	B	338	VAL	2.2
1	C	317	ASN	2.2
1	B	272	ILE	2.2
1	A	274	GLU	2.1
1	C	106	THR	2.1
1	A	143	ASP	2.1
1	C	145	GLY	2.0
1	B	280	VAL	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.