



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 11:45 AM GMT

PDB ID : 3PQE  
Title : Crystal structure of L-lactate dehydrogenase from Bacillus subtilis with H171C mutation  
Authors : Zhang, Y.; Garavito, R.M.  
Deposited on : 2010-11-26  
Resolution : 2.20 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](mailto:validation@mail.wwpdb.org)  
A user guide is available at  
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>  
with specific help available everywhere you see the ⓘ symbol.

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The following versions of software and data (see [references ⓘ](#)) 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

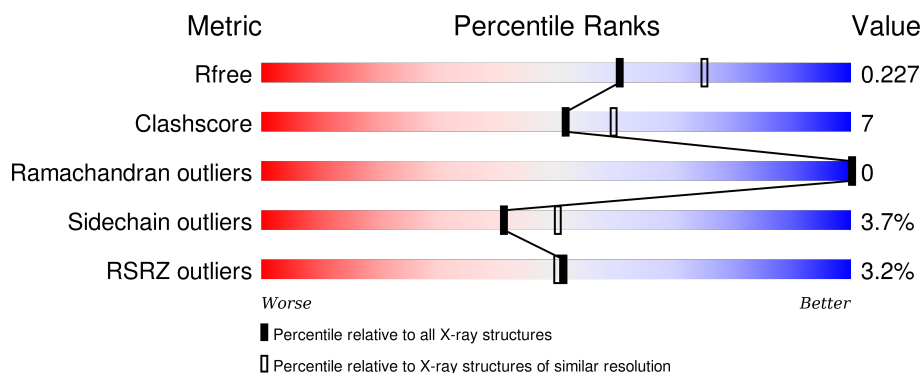
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.20 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	3774 (2.20-2.20)
Clashscore	102246	4477 (2.20-2.20)
Ramachandran outliers	100387	4404 (2.20-2.20)
Sidechain outliers	100360	4405 (2.20-2.20)
RSRZ outliers	91569	3781 (2.20-2.20)

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  $\geq 3$ , 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  $\leq 5\%$ . 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	326	<div> <div>0%</div> <div> <div></div> <div>84%</div> <div>12%</div> <div>•</div> </div> </div>
1	B	326	<div> <div>2%</div> <div> <div></div> <div>83%</div> <div>12%</div> <div>• •</div> </div> </div>
1	C	326	<div> <div>5%</div> <div> <div></div> <div>76%</div> <div>16%</div> <div>• 6%</div> </div> </div>
1	D	326	<div> <div>5%</div> <div> <div></div> <div>75%</div> <div>18%</div> <div>6%</div> </div> </div>

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 9845 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 L-lactate dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	314	Total	C	N	O	S	0	0	0
			2397	1527	401	460	9			
1	B	314	Total	C	N	O	S	0	0	0
			2397	1527	401	460	9			
1	C	305	Total	C	N	O	S	0	0	0
			2331	1488	387	447	9			
1	D	306	Total	C	N	O	S	0	0	0
			2342	1494	391	448	9			

There are 28 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	171	CYS	HIS	ENGINEERED MUTATION	UNP P13714
A	321	HIS	-	EXPRESSION TAG	UNP P13714
A	322	HIS	-	EXPRESSION TAG	UNP P13714
A	323	HIS	-	EXPRESSION TAG	UNP P13714
A	324	HIS	-	EXPRESSION TAG	UNP P13714
A	325	HIS	-	EXPRESSION TAG	UNP P13714
A	326	HIS	-	EXPRESSION TAG	UNP P13714
B	171	CYS	HIS	ENGINEERED MUTATION	UNP P13714
B	321	HIS	-	EXPRESSION TAG	UNP P13714
B	322	HIS	-	EXPRESSION TAG	UNP P13714
B	323	HIS	-	EXPRESSION TAG	UNP P13714
B	324	HIS	-	EXPRESSION TAG	UNP P13714
B	325	HIS	-	EXPRESSION TAG	UNP P13714
B	326	HIS	-	EXPRESSION TAG	UNP P13714
C	171	CYS	HIS	ENGINEERED MUTATION	UNP P13714
C	321	HIS	-	EXPRESSION TAG	UNP P13714
C	322	HIS	-	EXPRESSION TAG	UNP P13714
C	323	HIS	-	EXPRESSION TAG	UNP P13714
C	324	HIS	-	EXPRESSION TAG	UNP P13714
C	325	HIS	-	EXPRESSION TAG	UNP P13714
C	326	HIS	-	EXPRESSION TAG	UNP P13714

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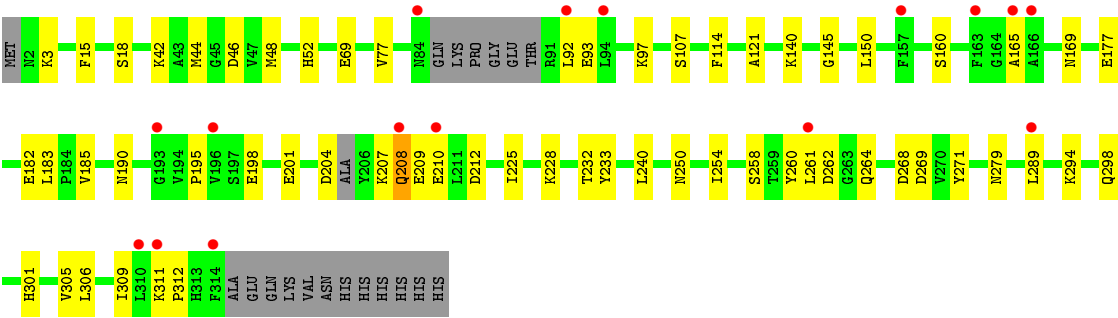
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Chain	Residue	Modelled	Actual	Comment	Reference
D	171	CYS	HIS	ENGINEERED MUTATION	UNP P13714
D	321	HIS	-	EXPRESSION TAG	UNP P13714
D	322	HIS	-	EXPRESSION TAG	UNP P13714
D	323	HIS	-	EXPRESSION TAG	UNP P13714
D	324	HIS	-	EXPRESSION TAG	UNP P13714
D	325	HIS	-	EXPRESSION TAG	UNP P13714
D	326	HIS	-	EXPRESSION TAG	UNP P13714

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	157	Total	O	0	0
			157	157		
2	B	154	Total	O	0	0
			154	154		
2	C	29	Total	O	0	0
			29	29		
2	D	38	Total	O	0	0
			38	38		





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 31	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	133.41 Å 133.41 Å 99.33 Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	26.93 – 2.20 26.93 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.5 (26.93-2.20) 99.5 (26.93-2.20)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.56 (at 2.20 Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.6 _289)	Depositor
R, $R_{free}$	0.201 , 0.235 0.194 , 0.227	Depositor DCC
$R_{free}$ test set	2995 reflections (3.00%)	DCC
Wilson B-factor (Å <sup>2</sup> )	44.3	Xtriage
Anisotropy	0.010	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	(Not available) , (Not available)	EDS
Estimated twinning fraction	0.487 for -h,-k,l 0.027 for h,-h-k,-l 0.027 for -k,-h,-l	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 99875 reflections	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	9845	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.77% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  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	0.49	0/2443	0.54	0/3311
1	B	0.49	0/2443	0.54	0/3311
1	C	0.34	0/2374	0.46	0/3215
1	D	0.33	0/2385	0.47	0/3229
All	All	0.42	0/9645	0.50	0/13066

There are no bond length outliers.

There are no bond angle outliers.

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	2397	0	2384	31	0
1	B	2397	0	2384	35	0
1	C	2331	0	2315	40	0
1	D	2342	0	2328	40	0
2	A	157	0	0	6	0
2	B	154	0	0	4	0
2	C	29	0	0	1	0
2	D	38	0	0	0	0
All	All	9845	0	9411	129	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 (129) 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:156:ARG:HG2	1:B:156:ARG:HH11	1.12	1.13
1:B:85:GLN:HB3	1:B:94:LEU:HD22	1.43	1.00
1:B:156:ARG:NH1	1:B:156:ARG:HG2	1.82	0.90
1:A:48:MET:CE	1:D:228:LYS:HD2	2.08	0.84
1:A:159:LEU:N	2:A:462:HOH:O	2.07	0.81
1:C:311:LYS:HG3	1:C:312:PRO:HD3	1.63	0.80
1:C:209:GLU:HA	1:C:212:ASP:HB2	1.67	0.76
1:C:254:ILE:HD12	1:D:169:ASN:HB3	1.70	0.74
1:C:81:ALA:O	1:C:122:THR:HG21	1.89	0.72
1:A:156:ARG:C	2:A:462:HOH:O	2.28	0.71
1:A:48:MET:HE2	1:D:228:LYS:HD2	1.70	0.71
1:A:48:MET:HE1	1:D:228:LYS:HD2	1.72	0.70
1:A:158:MET:N	2:A:462:HOH:O	2.25	0.70
1:B:48:MET:CE	1:C:228:LYS:HD2	2.23	0.68
1:B:156:ARG:CG	1:B:156:ARG:HH11	1.97	0.68
1:A:52:HIS:CE1	1:D:150:LEU:HD21	2.30	0.66
1:A:155:PHE:O	2:A:462:HOH:O	2.12	0.66
1:B:162:TYR:HE1	1:B:210:GLU:HG2	1.62	0.65
1:B:48:MET:HE1	1:C:228:LYS:HD2	1.77	0.65
1:D:160:SER:HB2	1:D:165:ALA:O	1.97	0.64
1:D:294:LYS:HE3	1:D:298:GLN:HE22	1.62	0.64
1:C:160:SER:HB2	1:C:165:ALA:O	1.99	0.62
1:C:301:HIS:O	1:C:305:VAL:HG23	2.00	0.61
1:B:190:ASN:C	2:B:478:HOH:O	2.39	0.61
1:C:140:LYS:HD2	1:C:260:TYR:CD2	2.36	0.60
1:A:188:HIS:HD2	1:B:193:GLY:O	1.83	0.60
1:A:85:GLN:HB2	1:A:94:LEU:HD22	1.83	0.60
1:D:140:LYS:HD2	1:D:260:TYR:CD2	2.37	0.59
1:D:177:GLU:HB2	1:D:306:LEU:HD11	1.85	0.59
1:A:161:GLU:HG2	1:A:161:GLU:O	2.04	0.58
1:D:301:HIS:O	1:D:305:VAL:HG23	2.03	0.58
1:B:52:HIS:CE1	1:C:150:LEU:HD21	2.39	0.58
1:D:93:GLU:O	1:D:97:LYS:HG3	2.03	0.58
1:B:44:MET:O	1:B:48:MET:HE3	2.03	0.58
1:A:178:HIS:HD2	2:A:372:HOH:O	1.87	0.57
1:A:193:GLY:O	1:B:188:HIS:HD2	1.88	0.57
1:B:204:ASP:O	1:B:207:LYS:HE2	2.06	0.55
1:D:262:ASP:HA	1:D:268:ASP:HA	1.89	0.55
1:A:214:ILE:O	1:A:218:VAL:HG23	2.07	0.54
1:C:169:ASN:ND2	1:D:254:ILE:H	2.05	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:154:ARG:O	1:A:158:MET:HG3	2.08	0.54
1:D:305:VAL:O	1:D:309:ILE:HG13	2.07	0.54
1:B:162:TYR:CE1	1:B:210:GLU:HG2	2.41	0.53
1:D:311:LYS:HB2	1:D:312:PRO:HD3	1.90	0.52
1:D:225:ILE:HG21	1:D:232:THR:HG22	1.90	0.52
1:C:177:GLU:O	1:C:182:GLU:HB3	2.10	0.52
1:B:156:ARG:HD3	2:B:390:HOH:O	2.10	0.52
1:D:77:VAL:HG23	1:D:114:PHE:CE1	2.45	0.52
1:D:225:ILE:CG2	1:D:232:THR:HG22	2.40	0.52
1:D:185:VAL:HG11	1:D:289:LEU:HD13	1.92	0.52
1:D:44:MET:O	1:D:48:MET:HG3	2.11	0.51
1:D:294:LYS:HE3	1:D:298:GLN:NE2	2.24	0.50
1:A:2:ASN:HB3	1:A:5:VAL:O	2.11	0.50
1:A:45:GLY:HA2	1:A:48:MET:HE3	1.93	0.50
1:A:140:LYS:HD2	1:A:260:TYR:CD2	2.46	0.50
1:D:260:TYR:OH	1:D:269:ASP:HA	2.12	0.50
1:C:305:VAL:O	1:C:309:ILE:HG13	2.12	0.49
1:A:204:ASP:O	1:A:207:LYS:HE2	2.11	0.49
1:D:260:TYR:CZ	1:D:269:ASP:HA	2.47	0.49
1:C:92:LEU:O	1:C:96:GLU:HG3	2.13	0.49
1:C:169:ASN:HB2	1:D:254:ILE:HD12	1.95	0.49
1:D:250:ASN:OD1	1:D:279:ASN:HB2	2.13	0.49
1:B:157:PHE:CE2	1:B:161:GLU:OE2	2.66	0.49
1:C:157:PHE:CE2	1:C:161:GLU:OE1	2.66	0.49
1:D:177:GLU:O	1:D:182:GLU:HB3	2.13	0.48
1:B:275:PRO:HB2	1:B:287:THR:HB	1.94	0.48
1:C:155:PHE:CE2	1:C:172:ALA:HB1	2.48	0.48
1:C:311:LYS:CG	1:C:312:PRO:HD3	2.39	0.48
1:B:140:LYS:HD2	1:B:260:TYR:CD2	2.48	0.48
1:B:214:ILE:O	1:B:218:VAL:HG23	2.14	0.48
1:A:250:ASN:OD1	1:A:279:ASN:HB2	2.13	0.48
1:B:150:LEU:HD21	1:C:52:HIS:CE1	2.49	0.47
1:A:158:MET:SD	2:A:457:HOH:O	2.60	0.47
1:D:195:PRO:HG2	1:D:198:GLU:HG3	1.97	0.47
1:A:150:LEU:HD21	1:D:52:HIS:CE1	2.49	0.47
1:C:77:VAL:HG23	1:C:114:PHE:CE1	2.50	0.47
1:A:40:LYS:HE3	1:A:40:LYS:HB3	1.80	0.47
1:B:177:GLU:O	1:B:177:GLU:HG2	2.12	0.47
1:A:25:ILE:HG12	1:A:61:VAL:HG11	1.95	0.47
1:B:202:LYS:HB2	1:B:202:LYS:NZ	2.30	0.47
1:A:158:MET:O	1:A:161:GLU:C	2.53	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:280:ARG:NH2	1:D:3:LYS:HD3	2.30	0.46
1:A:77:VAL:HG23	1:A:114:PHE:CE1	2.51	0.46
1:B:196:VAL:O	1:B:200:VAL:HG23	2.16	0.46
1:C:154:ARG:NH1	2:C:348:HOH:O	2.46	0.46
1:D:177:GLU:HG3	1:D:306:LEU:HD21	1.96	0.45
1:B:77:VAL:HG23	1:B:114:PHE:CE1	2.51	0.45
1:D:42:LYS:HA	1:D:42:LYS:HD2	1.82	0.45
1:D:145:GLY:HA3	1:D:258:SER:HB2	1.99	0.44
1:B:250:ASN:OD1	1:B:279:ASN:HB2	2.17	0.44
1:C:207:LYS:N	1:C:207:LYS:HD2	2.32	0.44
1:C:40:LYS:HE3	1:C:41:GLU:OE1	2.18	0.44
1:C:262:ASP:HA	1:C:268:ASP:HA	2.00	0.44
1:C:177:GLU:HG3	1:C:306:LEU:HD21	1.99	0.44
1:C:260:TYR:OH	1:C:269:ASP:HA	2.18	0.43
1:C:157:PHE:HE2	1:C:161:GLU:OE1	2.00	0.43
1:D:208:GLN:HE21	1:D:208:GLN:HA	1.83	0.43
1:B:42:LYS:NZ	2:B:435:HOH:O	2.51	0.43
1:C:133:TRP:HA	1:C:143:VAL:HG21	2.00	0.43
1:B:48:MET:HE2	1:C:228:LYS:HD2	1.99	0.43
1:C:260:TYR:CZ	1:C:269:ASP:HA	2.53	0.43
1:C:106:VAL:O	1:C:110:MET:HG2	2.18	0.43
1:A:209:GLU:HA	1:A:209:GLU:OE2	2.18	0.43
1:A:159:LEU:HA	1:A:159:LEU:HD12	1.73	0.42
1:D:260:TYR:HB2	1:D:271:TYR:CZ	2.54	0.42
1:B:204:ASP:HA	2:B:342:HOH:O	2.19	0.42
1:C:213:GLN:HA	1:C:216:ASP:OD2	2.20	0.42
1:D:15:PHE:HB3	1:D:233:TYR:CD2	2.54	0.42
1:D:261:LEU:HD22	1:D:264:GLN:HB2	2.00	0.42
1:C:44:MET:O	1:C:48:MET:HG3	2.19	0.42
1:C:260:TYR:HB2	1:C:271:TYR:CZ	2.55	0.42
1:D:145:GLY:CA	1:D:258:SER:HB2	2.49	0.42
1:A:275:PRO:HB2	1:A:287:THR:HB	2.01	0.42
1:A:157:PHE:CD2	1:A:157:PHE:C	2.93	0.42
1:C:169:ASN:HD21	1:D:254:ILE:H	1.67	0.41
1:B:61:VAL:O	1:B:61:VAL:HG13	2.19	0.41
1:B:40:LYS:HE3	1:B:40:LYS:HB3	1.90	0.41
1:C:309:ILE:HG13	1:C:309:ILE:H	1.65	0.41
1:B:15:PHE:HB3	1:B:233:TYR:CD2	2.55	0.41
1:B:207:LYS:HD3	1:B:207:LYS:N	2.36	0.41
1:B:168:GLN:HG2	1:B:168:GLN:H	1.65	0.41
1:B:202:LYS:CB	1:B:202:LYS:NZ	2.84	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:204:ASP:O	1:C:207:LYS:HE3	2.20	0.41
1:C:166:ALA:HB1	1:C:168:GLN:OE1	2.21	0.41
1:C:73:ASP:OD1	1:C:73:ASP:N	2.53	0.41
1:D:209:GLU:HA	1:D:212:ASP:HB3	2.03	0.40
1:D:121:ALA:HB2	1:D:240:LEU:HD21	2.03	0.40
1:A:61:VAL:O	1:A:61:VAL:HG13	2.21	0.40
1:C:42:LYS:HB3	1:C:42:LYS:HE2	1.90	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	312/326 (96%)	301 (96%)	11 (4%)	0	100	100
1	B	312/326 (96%)	301 (96%)	11 (4%)	0	100	100
1	C	299/326 (92%)	284 (95%)	15 (5%)	0	100	100
1	D	300/326 (92%)	291 (97%)	9 (3%)	0	100	100
All	All	1223/1304 (94%)	1177 (96%)	46 (4%)	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	255/267 (96%)	248 (97%)	7 (3%)	52	64
1	B	255/267 (96%)	247 (97%)	8 (3%)	47	59
1	C	249/267 (93%)	239 (96%)	10 (4%)	38	47
1	D	250/267 (94%)	238 (95%)	12 (5%)	31	37
All	All	1009/1068 (94%)	972 (96%)	37 (4%)	41	50

All (37) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	86	LYS
1	A	159	LEU
1	A	160	SER
1	A	169	ASN
1	A	183	LEU
1	A	202	LYS
1	A	229	LYS
1	B	5	VAL
1	B	156	ARG
1	B	160	SER
1	B	177	GLU
1	B	183	LEU
1	B	201	GLU
1	B	207	LYS
1	B	208	GLN
1	C	42	LYS
1	C	46	ASP
1	C	107	SER
1	C	122	THR
1	C	154	ARG
1	C	183	LEU
1	C	190	ASN
1	C	207	LYS
1	C	210	GLU
1	C	308	ASN
1	D	18	SER
1	D	46	ASP
1	D	69	GLU
1	D	92	LEU
1	D	107	SER
1	D	183	LEU
1	D	190	ASN
1	D	201	GLU

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Mol	Chain	Res	Type
1	D	204	ASP
1	D	207	LYS
1	D	208	GLN
1	D	210	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	52	HIS
1	A	213	GLN
1	B	6	ASN
1	B	52	HIS
1	C	169	ASN
1	C	203	ASN
1	D	52	HIS
1	D	188	HIS
1	D	203	ASN
1	D	208	GLN
1	D	249	HIS

### 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](#)

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<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	314/326 (96%)	0.03	4 (1%) 79 78	27, 40, 65, 81	0
1	B	314/326 (96%)	0.03	5 (1%) 74 73	27, 40, 66, 81	0
1	C	305/326 (93%)	0.28	15 (4%) 33 33	39, 63, 92, 103	0
1	D	306/326 (93%)	0.27	16 (5%) 31 30	39, 64, 93, 110	0
All	All	1239/1304 (95%)	0.15	40 (3%) 51 50	27, 51, 86, 110	0

All (40) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	94	LEU	4.5
1	C	193	GLY	4.5
1	C	94	LEU	4.4
1	A	205	ALA	4.2
1	C	314	PHE	3.9
1	D	193	GLY	3.8
1	D	314	PHE	3.7
1	D	163	PHE	3.6
1	C	163	PHE	3.5
1	D	210	GLU	3.2
1	A	211	LEU	3.0
1	D	208	GLN	3.0
1	B	205	ALA	3.0
1	D	310	LEU	3.0
1	D	157	PHE	3.0
1	C	310	LEU	2.9
1	D	92	LEU	2.8
1	D	311	LYS	2.8
1	C	208	GLN	2.8
1	C	157	PHE	2.7
1	D	166	ALA	2.7

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Mol	Chain	Res	Type	RSRZ
1	B	168	GLN	2.7
1	C	306	LEU	2.6
1	D	261	LEU	2.6
1	D	289	LEU	2.5
1	C	311	LYS	2.5
1	D	84	ASN	2.5
1	C	166	ALA	2.4
1	C	210	GLU	2.4
1	C	300	LEU	2.4
1	B	214	ILE	2.4
1	C	2	ASN	2.4
1	B	211	LEU	2.3
1	C	84	ASN	2.3
1	D	165	ALA	2.2
1	A	214	ILE	2.1
1	C	261	LEU	2.1
1	A	163	PHE	2.0
1	D	196	VAL	2.0
1	B	163	PHE	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 [i](#)

There are no ligands in this entry.

## 6.5 Other polymers [i](#)

There are no such residues in this entry.