



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 06:26 PM GMT

PDB ID : 4LN1
Title : CRYSTAL STRUCTURE OF L-lactate dehydrogenase from Bacillus cereus ATCC 14579 complexed with calcium, NYSGRC Target 029452
Authors : Malashkevich, V.N.; Bonanno, J.B.; Bhosle, R.; Toro, R.; Hillerich, B.; Gizzi, A.; Garforth, S.; Kar, A.; Chan, M.K.; Lafluer, J.; Patel, H.; Matikainen, B.; Chamala, S.; Lim, S.; Celikgil, A.; Villegas, G.; Evans, B.; Love, J.; Fiser, A.; Khafizov, K.; Seidel, R.; Almo, S.C.; New York Structural Genomics Research Consortium (NYSGRC)
Deposited on : 2013-07-11
Resolution : 1.90 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
<http://wwpdb.org/validation/2016/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 : 1.7 (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

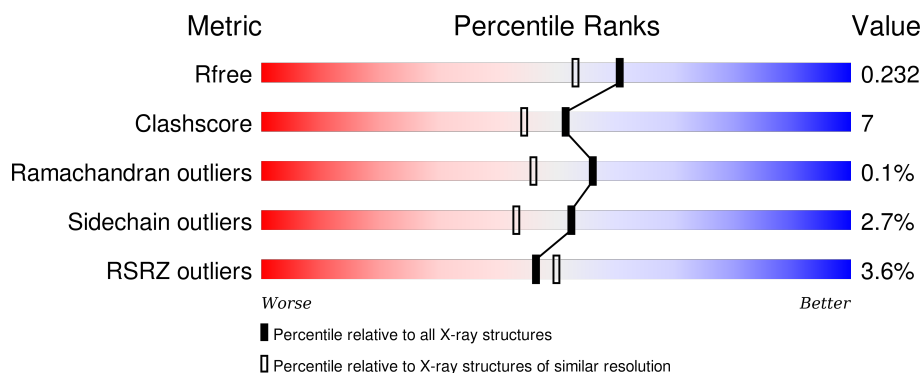
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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}	91344	4755 (1.90-1.90)
Clashscore	102246	5398 (1.90-1.90)
Ramachandran outliers	100387	5338 (1.90-1.90)
Sidechain outliers	100360	5339 (1.90-1.90)
RSRZ outliers	91569	4766 (1.90-1.90)

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 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 $\leq 5\%$. 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	336	<div> <div>2%</div> <div>76% 18% 6%</div> </div>
1	B	336	<div> <div>4%</div> <div>80% 15% •</div> </div>
1	C	336	<div> <div>3%</div> <div>74% 19% • 6%</div> </div>
1	D	336	<div> <div>5%</div> <div>77% 16% • 6%</div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard

residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	CA	A	500	-	-	-	X

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 10343 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 1.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	317	Total	C	N	O	S	Se	0	1	0
			2494	1590	415	479	3	7			
1	B	321	Total	C	N	O	S	Se	0	4	0
			2535	1617	421	486	3	8			
1	C	316	Total	C	N	O	S	Se	0	0	0
			2474	1579	409	476	3	7			
1	D	315	Total	C	N	O	S	Se	0	2	0
			2483	1584	413	476	3	7			

There are 88 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-21	MSE	-	EXPRESSION TAG	UNP Q81EP4
A	-20	HIS	-	EXPRESSION TAG	UNP Q81EP4
A	-19	HIS	-	EXPRESSION TAG	UNP Q81EP4
A	-18	HIS	-	EXPRESSION TAG	UNP Q81EP4
A	-17	HIS	-	EXPRESSION TAG	UNP Q81EP4
A	-16	HIS	-	EXPRESSION TAG	UNP Q81EP4
A	-15	HIS	-	EXPRESSION TAG	UNP Q81EP4
A	-14	SER	-	EXPRESSION TAG	UNP Q81EP4
A	-13	SER	-	EXPRESSION TAG	UNP Q81EP4
A	-12	GLY	-	EXPRESSION TAG	UNP Q81EP4
A	-11	VAL	-	EXPRESSION TAG	UNP Q81EP4
A	-10	ASP	-	EXPRESSION TAG	UNP Q81EP4
A	-9	LEU	-	EXPRESSION TAG	UNP Q81EP4
A	-8	GLY	-	EXPRESSION TAG	UNP Q81EP4
A	-7	THR	-	EXPRESSION TAG	UNP Q81EP4
A	-6	GLU	-	EXPRESSION TAG	UNP Q81EP4
A	-5	ASN	-	EXPRESSION TAG	UNP Q81EP4
A	-4	LEU	-	EXPRESSION TAG	UNP Q81EP4
A	-3	TYR	-	EXPRESSION TAG	UNP Q81EP4
A	-2	PHE	-	EXPRESSION TAG	UNP Q81EP4
A	-1	GLN	-	EXPRESSION TAG	UNP Q81EP4

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Chain	Residue	Modelled	Actual	Comment	Reference
A	0	SER	-	EXPRESSION TAG	UNP Q81EP4
B	-21	MSE	-	EXPRESSION TAG	UNP Q81EP4
B	-20	HIS	-	EXPRESSION TAG	UNP Q81EP4
B	-19	HIS	-	EXPRESSION TAG	UNP Q81EP4
B	-18	HIS	-	EXPRESSION TAG	UNP Q81EP4
B	-17	HIS	-	EXPRESSION TAG	UNP Q81EP4
B	-16	HIS	-	EXPRESSION TAG	UNP Q81EP4
B	-15	HIS	-	EXPRESSION TAG	UNP Q81EP4
B	-14	SER	-	EXPRESSION TAG	UNP Q81EP4
B	-13	SER	-	EXPRESSION TAG	UNP Q81EP4
B	-12	GLY	-	EXPRESSION TAG	UNP Q81EP4
B	-11	VAL	-	EXPRESSION TAG	UNP Q81EP4
B	-10	ASP	-	EXPRESSION TAG	UNP Q81EP4
B	-9	LEU	-	EXPRESSION TAG	UNP Q81EP4
B	-8	GLY	-	EXPRESSION TAG	UNP Q81EP4
B	-7	THR	-	EXPRESSION TAG	UNP Q81EP4
B	-6	GLU	-	EXPRESSION TAG	UNP Q81EP4
B	-5	ASN	-	EXPRESSION TAG	UNP Q81EP4
B	-4	LEU	-	EXPRESSION TAG	UNP Q81EP4
B	-3	TYR	-	EXPRESSION TAG	UNP Q81EP4
B	-2	PHE	-	EXPRESSION TAG	UNP Q81EP4
B	-1	GLN	-	EXPRESSION TAG	UNP Q81EP4
B	0	SER	-	EXPRESSION TAG	UNP Q81EP4
C	-21	MSE	-	EXPRESSION TAG	UNP Q81EP4
C	-20	HIS	-	EXPRESSION TAG	UNP Q81EP4
C	-19	HIS	-	EXPRESSION TAG	UNP Q81EP4
C	-18	HIS	-	EXPRESSION TAG	UNP Q81EP4
C	-17	HIS	-	EXPRESSION TAG	UNP Q81EP4
C	-16	HIS	-	EXPRESSION TAG	UNP Q81EP4
C	-15	HIS	-	EXPRESSION TAG	UNP Q81EP4
C	-14	SER	-	EXPRESSION TAG	UNP Q81EP4
C	-13	SER	-	EXPRESSION TAG	UNP Q81EP4
C	-12	GLY	-	EXPRESSION TAG	UNP Q81EP4
C	-11	VAL	-	EXPRESSION TAG	UNP Q81EP4
C	-10	ASP	-	EXPRESSION TAG	UNP Q81EP4
C	-9	LEU	-	EXPRESSION TAG	UNP Q81EP4
C	-8	GLY	-	EXPRESSION TAG	UNP Q81EP4
C	-7	THR	-	EXPRESSION TAG	UNP Q81EP4
C	-6	GLU	-	EXPRESSION TAG	UNP Q81EP4
C	-5	ASN	-	EXPRESSION TAG	UNP Q81EP4
C	-4	LEU	-	EXPRESSION TAG	UNP Q81EP4
C	-3	TYR	-	EXPRESSION TAG	UNP Q81EP4

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-2	PHE	-	EXPRESSION TAG	UNP Q81EP4
C	-1	GLN	-	EXPRESSION TAG	UNP Q81EP4
C	0	SER	-	EXPRESSION TAG	UNP Q81EP4
D	-21	MSE	-	EXPRESSION TAG	UNP Q81EP4
D	-20	HIS	-	EXPRESSION TAG	UNP Q81EP4
D	-19	HIS	-	EXPRESSION TAG	UNP Q81EP4
D	-18	HIS	-	EXPRESSION TAG	UNP Q81EP4
D	-17	HIS	-	EXPRESSION TAG	UNP Q81EP4
D	-16	HIS	-	EXPRESSION TAG	UNP Q81EP4
D	-15	HIS	-	EXPRESSION TAG	UNP Q81EP4
D	-14	SER	-	EXPRESSION TAG	UNP Q81EP4
D	-13	SER	-	EXPRESSION TAG	UNP Q81EP4
D	-12	GLY	-	EXPRESSION TAG	UNP Q81EP4
D	-11	VAL	-	EXPRESSION TAG	UNP Q81EP4
D	-10	ASP	-	EXPRESSION TAG	UNP Q81EP4
D	-9	LEU	-	EXPRESSION TAG	UNP Q81EP4
D	-8	GLY	-	EXPRESSION TAG	UNP Q81EP4
D	-7	THR	-	EXPRESSION TAG	UNP Q81EP4
D	-6	GLU	-	EXPRESSION TAG	UNP Q81EP4
D	-5	ASN	-	EXPRESSION TAG	UNP Q81EP4
D	-4	LEU	-	EXPRESSION TAG	UNP Q81EP4
D	-3	TYR	-	EXPRESSION TAG	UNP Q81EP4
D	-2	PHE	-	EXPRESSION TAG	UNP Q81EP4
D	-1	GLN	-	EXPRESSION TAG	UNP Q81EP4
D	0	SER	-	EXPRESSION TAG	UNP Q81EP4

- Molecule 2 is CALCIUM ION (three-letter code: CA) (formula: Ca).

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

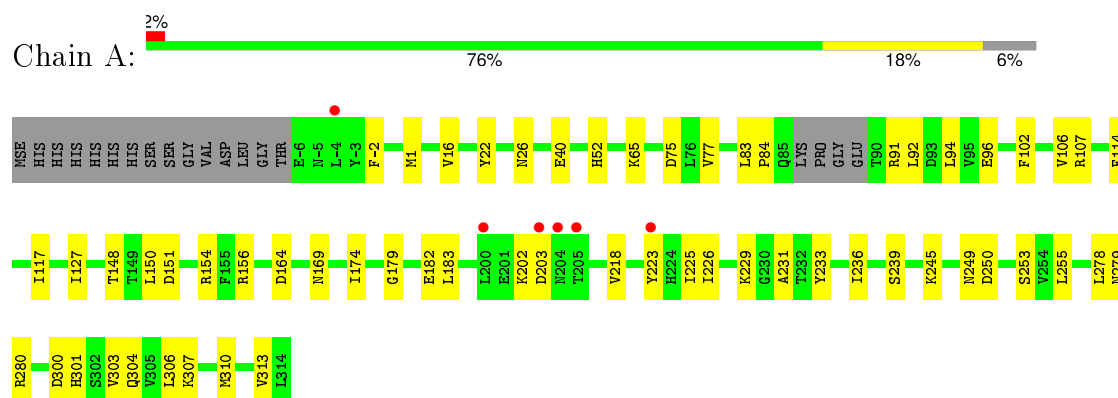
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	93	Total 93	O 93	0	0
3	B	85	Total 85	O 85	0	0
3	C	89	Total 89	O 89	0	0
3	D	84	Total 84	O 84	0	0

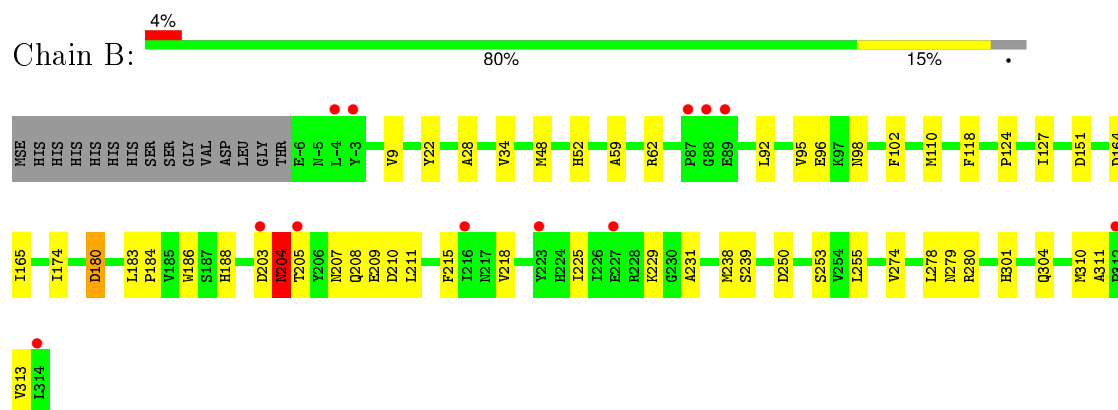
3 Residue-property plots [i](#)

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.

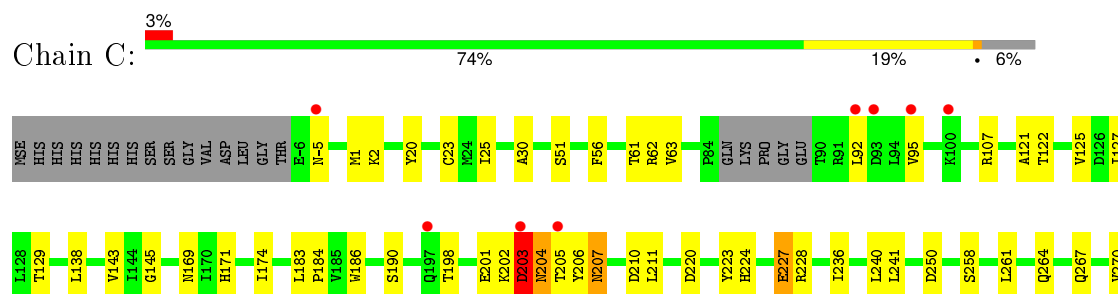
• Molecule 1: L-lactate dehydrogenase 1

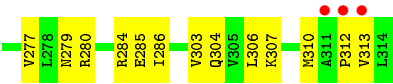


• Molecule 1: L-lactate dehydrogenase 1

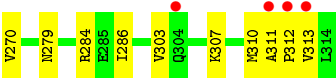
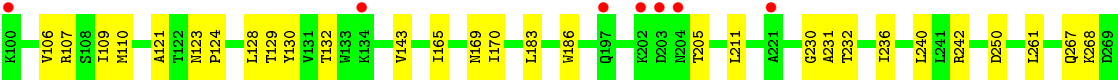
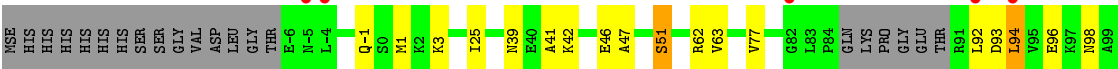
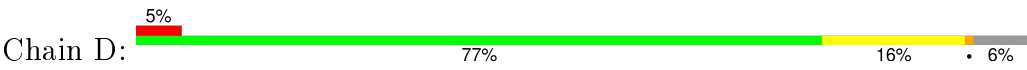


• Molecule 1: L-lactate dehydrogenase 1





● Molecule 1: L-lactate dehydrogenase 1



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	83.72Å 75.45Å 100.14Å 90.00° 89.99° 90.00°	Depositor
Resolution (Å)	42.98 – 1.90 42.97 – 1.80	Depositor EDS
% Data completeness (in resolution range)	99.6 (42.98-1.90) 99.5 (42.97-1.80)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.10 (at 1.79Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, R_{free}	0.209 , 0.241 0.210 , 0.232	Depositor DCC
R_{free} test set	4940 reflections (5.31%)	DCC
Wilson B-factor (Å ²)	29.6	Xtriage
Anisotropy	0.460	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 35.4	EDS
Estimated twinning fraction	0.514 for H, K, L 0.486 for H, -K, -L 0.468 for h,-k,-l	Xtriage
Reported twinning fraction	0.514 for H, K, L 0.486 for H, -K, -L	Depositor
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	3 of 115401 reflections (0.003%)	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	10343	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 89.69 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 3.1000e-08. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

² 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 [i](#)

5.1 Standard geometry [i](#)

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

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	0.54	1/2533 (0.0%)	0.64	0/3420
1	B	0.62	0/2588	0.66	2/3496 (0.1%)
1	C	0.70	2/2513 (0.1%)	0.67	1/3394 (0.0%)
1	D	0.55	0/2525	0.62	1/3410 (0.0%)
All	All	0.61	3/10159 (0.0%)	0.65	4/13720 (0.0%)

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	C	0	1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	206	TYR	CE1-CZ	-5.83	1.30	1.38
1	A	233	TYR	CE1-CZ	-5.45	1.31	1.38
1	C	227	GLU	CD-OE2	-5.32	1.19	1.25

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	204	ASN	N-CA-C	5.95	127.07	111.00
1	B	311	ALA	C-N-CD	5.48	139.90	128.40
1	C	227	GLU	OE1-CD-OE2	-5.31	116.92	123.30
1	D	94	LEU	CA-CB-CG	5.01	126.83	115.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	203	ASP	Peptide

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	2494	0	2485	42	0
1	B	2535	0	2537	37	0
1	C	2474	0	2465	45	0
1	D	2483	0	2476	32	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	93	0	0	5	0
3	B	85	0	0	0	0
3	C	89	0	0	4	0
3	D	84	0	0	1	0
All	All	10343	0	9963	135	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 (135) 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:306:LEU:O	1:A:310:MSE:HG3	1.68	0.92
1:C:284:ARG:NH2	1:C:285:GLU:OE2	2.06	0.87
1:C:220:ASP:O	1:C:223:TYR:HB2	1.85	0.76
1:C:202:LYS:O	1:C:203:ASP:HB2	1.86	0.76
1:B:301:HIS:HA	1:B:304:GLN:HG2	1.68	0.74
1:A:226:ILE:HG12	1:A:231:ALA:HA	1.68	0.74
1:B:164:ASP:O	1:C:284:ARG:NH1	2.22	0.73
1:D:121:ALA:HB2	1:D:240:LEU:HD21	1.70	0.72
1:B:207:ASN:OD1	1:B:207:ASN:N	2.27	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:180:ASP:OD1	1:B:180:ASP:N	2.30	0.65
1:C:92:LEU:HD21	1:C:312:PRO:HG2	1.78	0.64
1:C:207:ASN:HD21	1:C:210:ASP:CG	2.02	0.63
1:C:204:ASN:N	1:C:204:ASN:OD1	2.31	0.63
1:B:280:ARG:HB3	1:D:1:MSE:HG3	1.79	0.63
1:A:202:LYS:HG2	1:A:203:ASP:H	1.62	0.63
1:B:151:ASP:HB3	1:B:174:ILE:HD12	1.81	0.63
1:C:121:ALA:HB2	1:C:240:LEU:HD21	1.80	0.62
1:B:110[B]:MSE:HE1	1:B:118:PHE:CZ	2.35	0.62
1:C:186:TRP:CH2	1:C:211:LEU:HB3	2.36	0.61
1:D:236:ILE:O	1:D:236:ILE:HG13	2.00	0.61
1:B:250:ASP:OD1	1:B:279:ASN:HB2	2.02	0.59
1:D:96:GLU:HG2	1:D:313:VAL:HG22	1.85	0.58
1:A:75:ASP:OD1	1:A:280[A]:ARG:NH1	2.38	0.57
1:D:130:TYR:CD1	1:D:310:MSE:HE1	2.39	0.57
1:A:150:LEU:HB2	1:A:239:SER:OG	2.05	0.56
1:C:138:LEU:HD22	3:C:615:HOH:O	2.07	0.55
1:C:25:ILE:HD11	1:C:63:VAL:HG21	1.89	0.55
1:A:127:ILE:HG12	1:A:310:MSE:HG2	1.89	0.55
1:C:280:ARG:NE	3:C:658:HOH:O	2.33	0.55
1:C:145:GLY:HA3	1:C:258:SER:HB2	1.88	0.55
1:C:2:LYS:HE2	1:C:2:LYS:HA	1.88	0.55
1:C:171:HIS:HB2	1:C:190:SER:HB2	1.88	0.55
1:C:202:LYS:O	1:C:203:ASP:CB	2.54	0.55
1:B:174:ILE:CD1	1:B:218:VAL:HG11	2.37	0.55
1:A:164:ASP:O	1:D:284:ARG:NH1	2.40	0.55
1:A:280[B]:ARG:HB3	1:C:1:MSE:HG3	1.90	0.54
1:B:255:LEU:HD12	1:B:278:LEU:HD22	1.90	0.53
1:D:230:GLY:O	1:D:231:ALA:HB2	2.09	0.53
1:D:46:GLU:HA	1:D:46:GLU:OE1	2.08	0.53
1:D:-1:GLN:O	1:D:3:LYS:HE3	2.09	0.53
1:B:165:ILE:CD1	1:C:277:VAL:HG21	2.38	0.53
1:A:255:LEU:HD12	1:A:278:LEU:HD22	1.91	0.52
1:B:313:VAL:HG12	1:B:313:VAL:O	2.10	0.52
1:B:110[B]:MSE:HE1	1:B:118:PHE:CE2	2.44	0.52
1:A:174:ILE:CD1	1:A:218:VAL:HG11	2.40	0.52
1:A:148:THR:HA	1:A:151:ASP:OD1	2.10	0.51
1:B:92:LEU:O	1:B:96:GLU:HB2	2.11	0.51
1:D:123[B]:ASN:OD1	1:D:124:PRO:HA	2.10	0.50
1:A:16:VAL:HG13	1:A:236:ILE:HG23	1.94	0.50
1:D:106:VAL:O	1:D:110:MSE:HG2	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:52:HIS:CE1	1:B:225:ILE:HD11	2.46	0.50
1:A:250:ASP:OD1	1:A:279:ASN:HB2	2.11	0.50
1:C:250:ASP:OD1	1:C:279:ASN:HB2	2.12	0.50
1:C:261:LEU:HD23	1:C:286:ILE:HD13	1.94	0.49
1:C:224:HIS:O	1:C:227:GLU:HB2	2.13	0.49
1:B:95:VAL:HG22	1:B:124:PRO:HG3	1.94	0.49
1:D:39:ASN:HB3	1:D:42:LYS:HB3	1.95	0.49
1:A:300:ASP:O	1:A:304:GLN:HG2	2.13	0.49
1:C:207:ASN:HD22	1:C:210:ASP:H	1.61	0.49
1:A:117:ILE:HD12	3:A:613:HOH:O	2.13	0.49
1:D:270:VAL:HG13	1:D:307:LYS:HE3	1.95	0.48
1:A:303:VAL:O	1:A:307:LYS:HG3	2.13	0.48
1:A:102:PHE:O	1:A:106:VAL:HG23	2.13	0.48
1:B:207:ASN:ND2	1:B:209:GLU:HB2	2.29	0.48
1:D:47:ALA:O	1:D:51:SER:HB2	2.13	0.48
1:D:123[A]:ASN:ND2	3:D:680:HOH:O	2.47	0.48
1:C:207:ASN:ND2	1:C:210:ASP:CG	2.67	0.48
1:B:250:ASP:OD2	1:D:62:ARG:NH2	2.47	0.48
1:A:253:SER:HA	1:D:169:ASN:ND2	2.29	0.47
1:D:261:LEU:HD23	1:D:286:ILE:HD13	1.96	0.47
1:A:301:HIS:HA	1:A:304:GLN:HE21	1.79	0.47
1:A:40:GLU:HG2	1:A:65:LYS:HE3	1.95	0.47
1:A:229:LYS:HE2	1:A:231:ALA:O	2.15	0.47
1:A:96:GLU:HG2	1:A:313:VAL:HG11	1.95	0.47
1:B:9:VAL:HG13	1:B:34:VAL:HB	1.96	0.47
1:C:30:ALA:O	1:C:61:THR:HG23	2.15	0.47
1:C:261:LEU:HD22	1:C:264:GLN:HB2	1.96	0.47
1:B:184:PRO:HG2	1:B:186:TRP:NE1	2.31	0.46
1:A:250:ASP:OD2	1:C:62:ARG:NH1	2.48	0.46
1:A:223:TYR:HD1	3:A:634:HOH:O	1.98	0.46
1:D:250:ASP:OD1	1:D:279:ASN:HB2	2.15	0.46
1:A:307:LYS:NZ	3:A:682:HOH:O	2.45	0.46
1:A:92:LEU:O	1:A:96:GLU:HG3	2.16	0.45
1:B:203:ASP:O	1:B:204:ASN:HB2	2.15	0.45
1:C:56:PHE:CD1	1:D:242:ARG:HG2	2.51	0.45
1:D:165:ILE:HG21	1:D:170:ILE:HD11	1.97	0.45
1:B:127:ILE:HG12	1:B:310:MSE:HG2	1.97	0.45
1:D:98:ASN:HB3	1:D:128:LEU:HD11	1.97	0.45
1:A:280[A]:ARG:HB3	1:C:1:MSE:HG3	1.98	0.45
1:C:129:THR:HG23	1:C:143:VAL:HG12	1.99	0.45
1:A:26:ASN:ND2	1:B:238:MSE:HE1	2.30	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:267:GLN:HG3	1:D:303:VAL:HG21	1.99	0.44
1:C:236:ILE:HG13	1:C:236:ILE:O	2.18	0.44
1:B:165:ILE:HD12	1:C:277:VAL:CG2	2.47	0.44
1:A:179:GLY:O	1:A:182:GLU:HG2	2.18	0.44
1:A:245:LYS:NZ	3:A:647:HOH:O	2.51	0.43
1:B:253:SER:HA	1:C:169:ASN:ND2	2.33	0.43
1:B:208:GLN:HA	1:B:211:LEU:HD12	2.00	0.43
1:C:122:THR:O	1:C:125:VAL:HA	2.19	0.43
1:A:156:ARG:NE	3:A:675:HOH:O	2.51	0.43
1:D:311:ALA:HB3	1:D:312:PRO:HD3	2.00	0.43
1:D:205:THR:HG22	1:D:205:THR:O	2.18	0.43
1:B:184:PRO:HG3	1:B:215:PHE:CG	2.53	0.43
1:B:62:ARG:NH1	1:D:250:ASP:OD2	2.50	0.43
1:B:229:LYS:HE2	1:B:231:ALA:O	2.19	0.43
1:C:241:LEU:HD13	1:C:241:LEU:C	2.39	0.43
1:D:186:TRP:CH2	1:D:211:LEU:HB3	2.54	0.42
1:D:25:ILE:HD11	1:D:63:VAL:HG21	2.01	0.42
1:D:77:VAL:HG11	1:D:109:ILE:HD13	2.02	0.42
1:C:95:VAL:HG12	1:C:313:VAL:HG11	2.01	0.42
1:A:-2:PHE:HA	1:A:1:MSE:HE3	2.01	0.42
1:A:225:ILE:CG1	1:B:48:MSE:HE3	2.49	0.42
1:A:225:ILE:HG12	1:B:48:MSE:HE3	2.02	0.42
1:B:165:ILE:HD12	1:C:277:VAL:HG21	2.01	0.42
1:B:28:ALA:CB	1:B:59:ALA:HB3	2.50	0.42
1:C:198:THR:O	1:C:201:GLU:HB2	2.20	0.42
1:C:228:ARG:HB3	1:D:41:ALA:O	2.20	0.42
1:C:270:VAL:HG13	1:C:307:LYS:HE3	2.01	0.42
1:A:250:ASP:OD2	1:C:62:ARG:NH2	2.53	0.42
1:A:77:VAL:HG23	1:A:114:PHE:CE1	2.55	0.41
1:A:16:VAL:HG13	1:A:236:ILE:CG2	2.50	0.41
1:D:106:VAL:HG21	1:D:132:THR:HG23	2.02	0.41
1:A:83:LEU:HA	1:A:84:PRO:HD2	1.78	0.41
1:A:245:LYS:O	1:A:249:ASN:HB2	2.21	0.41
1:C:174:ILE:HG13	1:C:184:PRO:HA	2.02	0.41
1:C:127:ILE:HG12	1:C:306:LEU:HD11	2.02	0.41
1:B:98:ASN:O	1:B:102:PHE:HB2	2.20	0.41
1:C:267:GLN:HG3	1:C:303:VAL:HG21	2.03	0.40
1:A:91:ARG:O	1:A:94:LEU:HB3	2.21	0.40
1:A:154:ARG:HG2	1:B:52:HIS:CG	2.56	0.40
1:B:188:HIS:HB3	3:C:660:HOH:O	2.20	0.40
1:B:203:ASP:O	1:B:204:ASN:CB	2.69	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:20:TYR:O	1:C:23:CYS:HB2	2.21	0.40
1:C:171:HIS:HB3	3:C:682:HOH:O	2.22	0.40
1:D:129:THR:HG23	1:D:143:VAL:HG12	2.04	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	314/336 (94%)	303 (96%)	11 (4%)	0	100	100
1	B	323/336 (96%)	313 (97%)	10 (3%)	0	100	100
1	C	312/336 (93%)	300 (96%)	11 (4%)	1 (0%)	46	35
1	D	313/336 (93%)	299 (96%)	14 (4%)	0	100	100
All	All	1262/1344 (94%)	1215 (96%)	46 (4%)	1 (0%)	56	46

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	203	ASP

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	270/277 (98%)	266 (98%)	4 (2%)	72	69
1	B	276/277 (100%)	268 (97%)	8 (3%)	50	40
1	C	268/277 (97%)	259 (97%)	9 (3%)	44	33
1	D	269/277 (97%)	261 (97%)	8 (3%)	48	38
All	All	1083/1108 (98%)	1054 (97%)	29 (3%)	52	43

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	TYR
1	A	107	ARG
1	A	169	ASN
1	A	183	LEU
1	B	22	TYR
1	B	180	ASP
1	B	183	LEU
1	B	204	ASN
1	B	205	THR
1	B	210	ASP
1	B	239	SER
1	B	274	VAL
1	C	-5	ASN
1	C	51	SER
1	C	107	ARG
1	C	183	LEU
1	C	204	ASN
1	C	205	THR
1	C	207	ASN
1	C	304	GLN
1	C	310	MSE
1	D	51	SER
1	D	92	LEU
1	D	93	ASP
1	D	94	LEU
1	D	107	ARG
1	D	183	LEU
1	D	232	THR
1	D	268	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	178	HIS
1	A	304	GLN
1	C	104	GLN
1	C	169	ASN
1	C	207	ASN
1	D	169	ASN

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](#)

Of 6 ligands modelled in this entry, 6 are 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 [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, 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	310/336 (92%)	0.07	6 (1%) 70 73	20, 37, 61, 92	0
1	B	314/336 (93%)	0.15	12 (3%) 44 48	22, 39, 68, 92	0
1	C	309/336 (91%)	0.10	11 (3%) 46 50	23, 40, 70, 92	0
1	D	308/336 (91%)	0.21	16 (5%) 31 34	23, 39, 70, 98	0
All	All	1241/1344 (92%)	0.13	45 (3%) 46 50	20, 39, 68, 98	0

All (45) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	203	ASP	9.9
1	A	205	THR	6.2
1	D	94	LEU	6.0
1	B	205	THR	5.5
1	D	204	ASN	4.7
1	B	87	PRO	4.4
1	A	204	ASN	4.3
1	C	311	ALA	4.0
1	A	223	TYR	3.9
1	C	-5	ASN	3.7
1	B	203	ASP	3.7
1	D	312	PRO	3.6
1	D	311	ALA	3.5
1	C	93	ASP	3.5
1	C	312	PRO	3.0
1	D	100	LYS	3.0
1	B	89	GLU	2.9
1	D	92	LEU	2.9
1	C	95	VAL	2.9
1	B	216	ILE	2.9
1	C	100	LYS	2.8

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Mol	Chain	Res	Type	RSRZ
1	D	313	VAL	2.6
1	D	197	GLN	2.6
1	C	313	VAL	2.6
1	A	203	ASP	2.5
1	B	314	LEU	2.5
1	D	82	GLY	2.5
1	D	-4	LEU	2.5
1	C	92	LEU	2.5
1	D	-5	ASN	2.4
1	B	88	GLY	2.4
1	C	203	ASP	2.3
1	D	202	LYS	2.3
1	C	205	THR	2.3
1	A	200	LEU	2.2
1	B	-4	LEU	2.2
1	D	304	GLN	2.2
1	B	223	TYR	2.2
1	A	-4	LEU	2.1
1	D	134	LYS	2.1
1	C	197	GLN	2.1
1	B	227	GLU	2.1
1	B	-3	TYR	2.0
1	B	312	PRO	2.0
1	D	221	ALA	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](#)

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(\AA^2)	Q<0.9
2	CA	A	500	1/1	0.99	0.11	2.51	34,34,34,34	0
2	CA	D	501	1/1	0.96	0.14	0.89	43,43,43,43	0
2	CA	B	500	1/1	0.98	0.11	0.68	34,34,34,34	0
2	CA	B	501	1/1	0.94	0.12	0.09	40,40,40,40	0
2	CA	C	501	1/1	0.98	0.08	-0.88	37,37,37,37	0
2	CA	A	501	1/1	0.98	0.08	-1.43	37,37,37,37	0

6.5 Other polymers [i](#)

There are no such residues in this entry.