



Full wwPDB X-ray Structure Validation Report ⓘ

Dec 5, 2023 – 04:20 am GMT

PDB ID : 2J5K
Title : 2.0 Å resolution structure of the wild type malate dehydrogenase from *Haloarcula marismortui* (radiation damage series)
Authors : Fioravanti, E.; Vellieux, F.M.D.; Amara, P.; Madern, D.; Weik, M.
Deposited on : 2006-09-18
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

<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

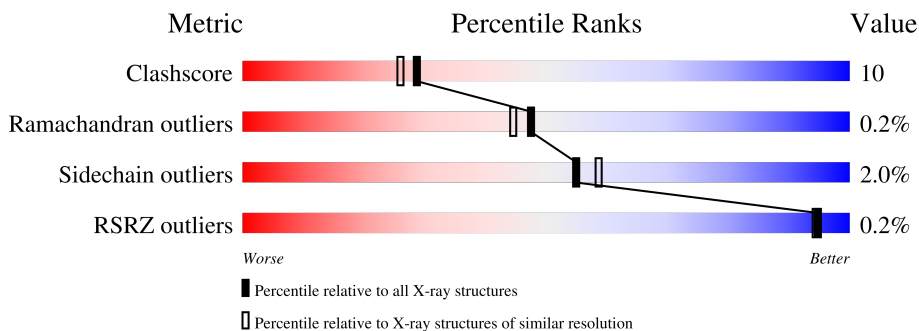
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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))
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (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 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	304	
1	B	304	
1	C	304	
1	D	304	

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	CL	A	1333	-	-	X	-
2	CL	B	1332	-	-	X	-
2	CL	C	1332	-	-	X	-
2	CL	D	1331	-	-	X	-

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 10083 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 MALATE DEHYDROGENASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	303	Total 2279	C 1407	N 391	O 477	S 4	0	0	0
1	B	298	Total 2240	C 1383	N 383	O 470	S 4	0	0	0
1	C	303	Total 2285	C 1409	N 393	O 479	S 4	0	0	0
1	D	298	Total 2248	C 1387	N 383	O 474	S 4	0	0	0

- Molecule 2 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	3	Total 3	Cl 3	0	0
2	B	2	Total 2	Cl 2	0	0
2	C	2	Total 2	Cl 2	0	0
2	D	1	Total 1	Cl 1	0	0

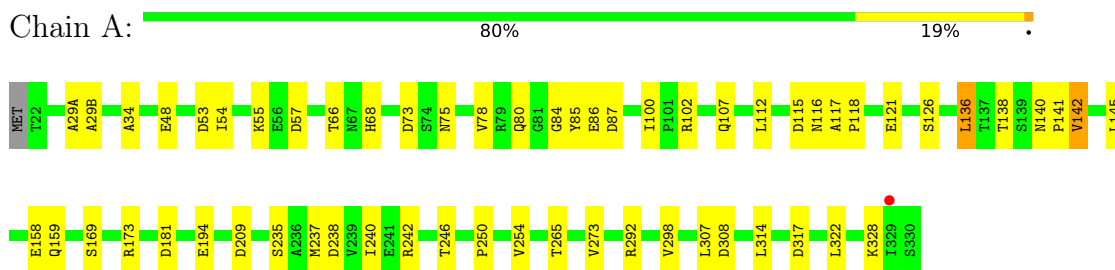
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	267	Total 267	O 267	0	0
3	B	230	Total 230	O 230	0	0
3	C	250	Total 250	O 250	0	0
3	D	276	Total 276	O 276	0	0

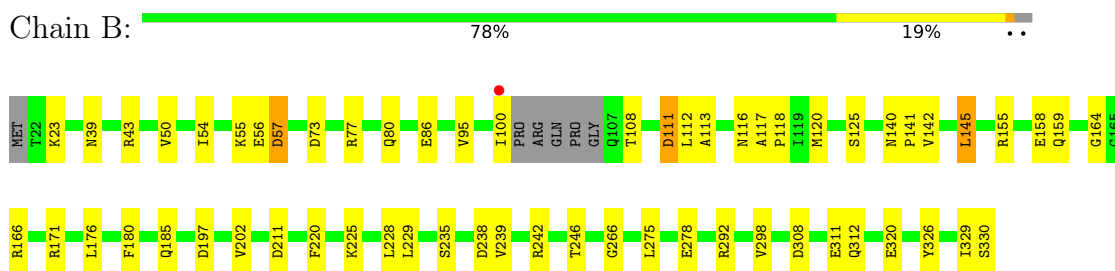
3 Residue-property plots [i](#)

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.

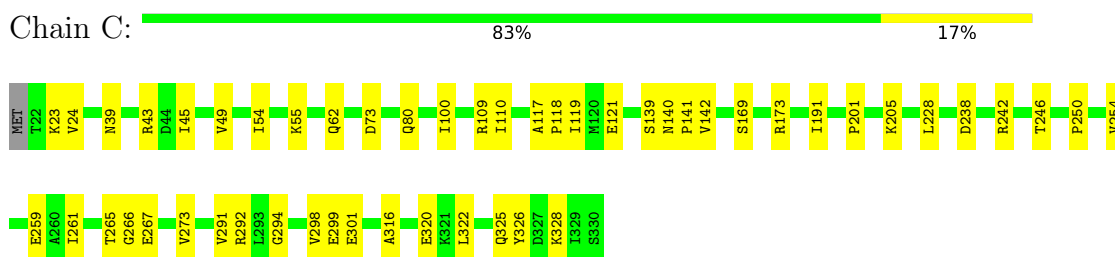
- Molecule 1: MALATE DEHYDROGENASE



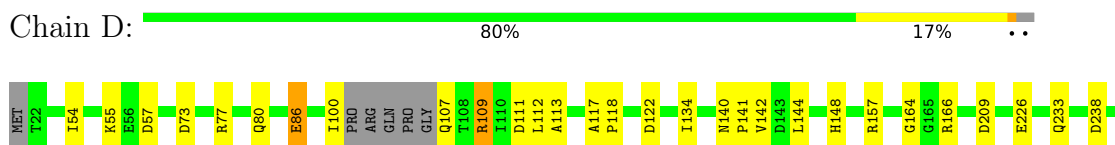
- Molecule 1: MALATE DEHYDROGENASE



- Molecule 1: MALATE DEHYDROGENASE



- Molecule 1: MALATE DEHYDROGENASE



R242
K243
G244
A245
T246
G249
P250
H256
D264
V268
K274
L275
E276
A286
V291
R292
G297
V298
E299
E301
D308
E311
K321
Q325
Y326
S330

4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	127.15Å 114.24Å 124.14Å 90.00° 93.48° 90.00°	Depositor
Resolution (Å)	20.00 – 2.00 20.00 – 2.00	Depositor EDS
% Data completeness (in resolution range)	93.8 (20.00-2.00) 93.9 (20.00-2.00)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.66 (at 2.01Å)	Xtrriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.225 , 0.270 0.230 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	17.8	Xtrriage
Anisotropy	0.671	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 42.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	10083	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.92% 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: CL

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	0.59	0/2316	0.72	0/3146
1	B	0.56	0/2274	0.73	0/3087
1	C	0.59	0/2322	0.77	0/3154
1	D	0.58	0/2282	0.74	0/3097
All	All	0.58	0/9194	0.74	0/12484

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

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	326	TYR	Sidechain

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	2279	0	2145	46	0
1	B	2240	0	2106	47	0
1	C	2285	0	2153	34	0
1	D	2248	0	2114	49	0
2	A	3	0	0	2	0
2	B	2	0	0	2	0
2	C	2	0	0	2	0
2	D	1	0	0	2	0
3	A	267	0	0	8	0
3	B	230	0	0	11	0
3	C	250	0	0	1	0
3	D	276	0	0	4	0
All	All	10083	0	8518	170	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 10.

All (170) 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:238:ASP:HA	1:C:242:ARG:HD2	1.55	0.88
1:A:181:ASP:HB2	3:A:2146:HOH:O	1.74	0.88
1:D:291:VAL:HG13	1:D:298:VAL:HG13	1.57	0.86
1:D:86:GLU:CD	1:D:86:GLU:H	1.85	0.80
1:B:100:ILE:HG12	1:B:116:ASN:OD1	1.84	0.77
1:B:108:THR:OG1	1:B:111:ASP:HB2	1.86	0.76
1:A:250:PRO:O	1:A:254:VAL:HG23	1.89	0.72
1:A:100:ILE:HG12	1:A:116:ASN:OD1	1.90	0.71
1:A:140:ASN:HB2	3:A:2116:HOH:O	1.93	0.69
1:C:62:GLN:NE2	1:D:243:LYS:HE3	2.07	0.69
1:D:100:ILE:O	1:D:100:ILE:HG13	1.92	0.68
1:A:75:ASN:OD1	1:B:185:GLN:HG3	1.95	0.67
2:C:1332:CL:CL	3:D:2058:HOH:O	2.49	0.67
1:B:246:THR:HG21	3:B:2108:HOH:O	1.93	0.67
1:B:308:ASP:OD1	1:B:311:GLU:HG3	1.94	0.67
1:C:191:ILE:HD13	1:C:201:PRO:HA	1.75	0.66
1:D:117:ALA:HB3	1:D:118:PRO:HD3	1.78	0.65
1:A:209:ASP:OD2	1:D:292:ARG:NH2	2.30	0.65
1:C:39:ASN:O	1:C:43:ARG:HG2	1.96	0.65
1:A:102:ARG:HB2	1:A:112:LEU:HD22	1.79	0.64
1:C:140:ASN:HA	1:C:142:VAL:N	2.12	0.64
1:B:95:VAL:HG13	3:B:2088:HOH:O	1.97	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:141:PRO:HB2	3:B:2091:HOH:O	1.99	0.63
1:C:266:GLY:O	1:C:292:ARG:HG3	1.98	0.63
1:B:100:ILE:HG13	1:B:100:ILE:O	1.99	0.62
1:D:122:ASP:HB3	3:D:2103:HOH:O	1.99	0.61
1:D:113:ALA:HB2	1:D:141:PRO:HG3	1.81	0.61
1:C:110:ILE:HD11	1:C:325:GLN:OE1	2.01	0.61
1:D:246:THR:HG22	1:D:250:PRO:HD3	1.82	0.60
1:B:54:ILE:HD11	1:B:55:LYS:HE3	1.84	0.60
1:B:86:GLU:H	1:B:86:GLU:CD	2.05	0.59
1:A:328:LYS:O	1:A:328:LYS:HD3	2.03	0.59
1:D:107:GLN:HE21	1:D:111:ASP:CB	2.16	0.58
1:D:238:ASP:O	1:D:242:ARG:HB2	2.04	0.58
1:D:246:THR:HG22	1:D:250:PRO:CD	2.34	0.57
1:B:326:TYR:O	1:B:330:SER:HB2	2.05	0.57
1:B:320:GLU:HG3	3:B:2222:HOH:O	2.04	0.57
1:A:169:SER:O	1:A:173:ARG:HG3	2.04	0.57
1:A:314:LEU:O	1:A:317:ASP:HB2	2.05	0.57
1:A:29(B):ALA:HB3	1:A:53:ASP:OD2	2.05	0.56
1:A:102:ARG:HD2	1:A:107:GLN:O	2.05	0.56
1:C:141:PRO:HB2	3:C:2114:HOH:O	2.05	0.56
1:C:316:ALA:O	1:C:320:GLU:HG3	2.06	0.56
1:D:134:ILE:HD13	1:D:297:GLY:CA	2.35	0.56
1:B:56:GLU:HA	3:B:2027:HOH:O	2.06	0.55
1:D:291:VAL:HG11	1:D:298:VAL:HG22	1.88	0.55
1:D:291:VAL:CG1	1:D:298:VAL:HG13	2.33	0.55
1:C:62:GLN:CD	1:D:243:LYS:HE3	2.27	0.55
1:D:166:ARG:HH11	1:D:166:ARG:HG3	1.71	0.54
3:B:2015:HOH:O	1:D:264:ASP:HB3	2.07	0.54
1:A:140:ASN:OD1	1:A:141:PRO:HA	2.07	0.54
1:D:274:LYS:HA	1:D:286:ALA:HB2	1.89	0.54
1:D:109:ARG:CG	1:D:109:ARG:HH21	2.21	0.54
1:D:268:VAL:HA	1:D:291:VAL:O	2.07	0.54
1:A:209:ASP:CG	1:D:292:ARG:HH22	2.10	0.54
1:A:117:ALA:O	1:A:121:GLU:HG3	2.09	0.53
1:D:166:ARG:HG3	1:D:166:ARG:NH1	2.24	0.53
1:A:86:GLU:H	1:A:86:GLU:CD	2.13	0.52
1:B:23:LYS:HE3	1:B:50:VAL:HG21	1.90	0.52
1:C:169:SER:O	1:C:173:ARG:HG3	2.10	0.52
1:A:238:ASP:O	1:A:242:ARG:HB2	2.09	0.52
1:B:266:GLY:O	1:B:292:ARG:HG3	2.10	0.52
1:C:261:ILE:HG23	1:C:294:GLY:O	2.10	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:140:ASN:HA	1:B:142:VAL:N	2.25	0.51
1:C:291:VAL:HB	1:C:298:VAL:HG13	1.92	0.51
1:C:265:THR:OG1	1:C:267:GLU:HG3	2.11	0.51
1:B:77:ARG:HA	1:B:77:ARG:HE	1.76	0.51
1:B:166:ARG:NH2	2:B:1332:CL:CL	2.77	0.51
1:D:308:ASP:OD1	1:D:311:GLU:HG3	2.11	0.51
1:C:246:THR:O	1:C:246:THR:HG23	2.11	0.51
1:B:57:ASP:HB2	3:B:2021:HOH:O	2.12	0.49
1:C:45:ILE:HG21	1:C:259:GLU:HG3	1.94	0.49
1:A:237:MET:CE	1:A:240:ILE:HD12	2.42	0.49
1:C:100:ILE:HD11	1:C:119:ILE:HD11	1.94	0.49
1:D:144:LEU:HG	3:D:2129:HOH:O	2.13	0.49
1:B:39:ASN:O	1:B:43:ARG:HG2	2.13	0.49
1:C:250:PRO:O	1:C:254:VAL:HG23	2.12	0.49
1:D:140:ASN:HA	1:D:142:VAL:N	2.28	0.48
1:D:242:ARG:NH2	1:D:242:ARG:HG3	2.28	0.48
1:C:140:ASN:HA	1:C:141:PRO:C	2.34	0.48
1:D:54:ILE:HD11	1:D:55:LYS:HE3	1.96	0.48
1:A:54:ILE:HD11	1:A:55:LYS:HE3	1.95	0.48
1:D:107:GLN:HE21	1:D:111:ASP:HB2	1.77	0.48
1:C:73:ASP:OD2	2:D:1331:CL:CL	2.69	0.48
1:A:117:ALA:N	1:A:118:PRO:HD2	2.29	0.48
2:A:1333:CL:CL	1:B:73:ASP:OD2	2.70	0.47
1:A:235:SER:O	1:A:238:ASP:HB2	2.15	0.47
1:C:273:VAL:HG21	1:C:298:VAL:HG21	1.95	0.47
1:A:142:VAL:HG22	3:A:2119:HOH:O	2.14	0.47
1:C:117:ALA:HB3	1:C:118:PRO:HD3	1.97	0.47
1:C:100:ILE:HG13	1:C:119:ILE:CD1	2.44	0.47
1:A:194:GLU:HB2	1:A:322:LEU:HD11	1.97	0.47
2:A:1333:CL:CL	3:B:2037:HOH:O	2.58	0.46
1:B:176:LEU:O	1:B:180:PHE:HD1	1.99	0.46
1:A:68:HIS:CG	1:B:171:ARG:HG2	2.50	0.46
1:A:246:THR:HG21	3:A:2139:HOH:O	2.14	0.46
1:C:24:VAL:O	1:C:49:VAL:HA	2.16	0.46
1:B:238:ASP:O	1:B:242:ARG:HB2	2.16	0.45
1:A:141:PRO:HB2	3:A:2125:HOH:O	2.15	0.45
2:C:1332:CL:CL	1:D:73:ASP:OD2	2.71	0.45
1:A:328:LYS:HD3	1:A:328:LYS:C	2.37	0.45
1:B:220:PHE:O	1:B:225:LYS:HE3	2.17	0.45
1:C:299:GLU:O	1:C:301:GLU:HB3	2.15	0.45
1:D:157:ARG:NE	1:D:276:GLU:OE1	2.49	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:113:ALA:HB2	1:B:141:PRO:HG2	1.98	0.45
1:B:211:ASP:OD2	1:C:205:LYS:NZ	2.40	0.45
1:B:235:SER:O	1:B:238:ASP:HB2	2.17	0.45
1:D:242:ARG:HG3	1:D:242:ARG:HH21	1.82	0.45
1:B:117:ALA:N	1:B:118:PRO:HD2	2.32	0.44
1:B:202:VAL:O	1:B:202:VAL:HG12	2.17	0.44
1:D:109:ARG:HH21	1:D:109:ARG:HB3	1.82	0.44
1:C:54:ILE:HG12	1:C:55:LYS:HD2	1.99	0.44
1:D:100:ILE:HD11	1:D:112:LEU:HD12	1.99	0.44
1:B:275:LEU:HD22	1:B:278:GLU:HB2	2.00	0.44
1:D:326:TYR:O	1:D:330:SER:HB2	2.17	0.44
1:B:312:GLN:HG2	3:B:2070:HOH:O	2.17	0.44
1:A:73:ASP:OD2	2:B:1332:CL:CL	2.73	0.44
1:B:23:LYS:HE3	1:B:50:VAL:CG2	2.48	0.44
1:D:86:GLU:CD	1:D:86:GLU:N	2.62	0.44
1:A:145:LEU:HD23	1:A:145:LEU:HA	1.76	0.43
1:C:117:ALA:O	1:C:121:GLU:HG3	2.17	0.43
1:A:292:ARG:NH1	1:D:209:ASP:OD2	2.48	0.43
1:D:117:ALA:HB1	1:D:148:HIS:CD2	2.53	0.43
1:D:244:GLY:O	1:D:245:ALA:HB2	2.19	0.43
1:D:77:ARG:NH2	3:D:2062:HOH:O	2.52	0.43
1:B:225:LYS:O	1:B:229:LEU:HG	2.19	0.42
1:D:299:GLU:O	1:D:301:GLU:HB3	2.19	0.42
1:B:120:MET:HG3	1:B:145:LEU:HG	2.01	0.42
1:A:273:VAL:HG11	1:A:298:VAL:HG21	2.01	0.42
1:B:140:ASN:HA	1:B:141:PRO:C	2.40	0.42
1:B:141:PRO:O	1:B:145:LEU:HB2	2.20	0.42
1:A:107:GLN:NE2	1:A:115:ASP:OD2	2.49	0.42
1:B:54:ILE:HG12	1:B:55:LYS:HD2	2.02	0.42
1:B:158:GLU:HB2	1:B:298:VAL:HB	2.02	0.42
1:D:109:ARG:HH21	1:D:109:ARG:CB	2.33	0.42
1:D:109:ARG:HG3	1:D:109:ARG:NH2	2.34	0.42
1:A:100:ILE:HD11	1:A:115:ASP:HB3	2.02	0.42
1:B:43:ARG:HG3	3:B:2178:HOH:O	2.18	0.42
1:A:85:TYR:CG	1:A:126:SER:HB3	2.55	0.42
1:C:139:SER:O	1:C:142:VAL:HA	2.20	0.42
1:B:239:VAL:CG1	1:B:246:THR:HA	2.50	0.41
1:D:321:LYS:O	1:D:325:GLN:HG3	2.20	0.41
1:B:108:THR:HG1	1:B:111:ASP:HB2	1.82	0.41
1:B:155:ARG:HB3	1:B:159:GLN:HE21	1.85	0.41
1:A:66:THR:HG22	1:A:78:VAL:HG21	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:249:GLY:N	1:D:250:PRO:HD2	2.36	0.41
1:A:237:MET:HE1	1:A:240:ILE:HD12	2.01	0.41
1:B:100:ILE:HD11	1:B:112:LEU:HD12	2.01	0.41
1:D:109:ARG:HH21	1:D:109:ARG:HG3	1.85	0.41
1:A:54:ILE:HG12	1:A:55:LYS:HD2	2.03	0.41
1:A:159:GLN:CG	3:A:2113:HOH:O	2.68	0.41
1:B:228:LEU:HG	3:B:2157:HOH:O	2.21	0.41
1:A:84:GLY:N	1:A:87:ASP:OD1	2.41	0.41
1:A:265:THR:HA	3:A:2216:HOH:O	2.20	0.41
1:C:23:LYS:HG2	1:C:24:VAL:N	2.34	0.41
1:C:322:LEU:HA	1:C:322:LEU:HD23	1.86	0.41
1:C:140:ASN:HD22	1:C:141:PRO:HA	1.85	0.41
1:A:29(A):ALA:O	1:A:34:ALA:HB2	2.21	0.41
1:D:109:ARG:CG	1:D:109:ARG:NH2	2.83	0.41
1:D:256:HIS:HB2	2:D:1331:CL:CL	2.58	0.41
1:B:329:ILE:O	1:B:329:ILE:HG13	2.20	0.40
1:A:138:THR:O	1:A:142:VAL:HG13	2.22	0.40
1:A:308:ASP:C	1:A:308:ASP:OD1	2.59	0.40
1:C:325:GLN:O	1:C:328:LYS:HB2	2.21	0.40
1:D:113:ALA:HB2	1:D:141:PRO:CG	2.49	0.40
1:A:158:GLU:H	1:A:158:GLU:CD	2.24	0.40
1:C:228:LEU:HD23	1:C:228:LEU:HA	1.75	0.40
1:A:136:LEU:HD23	1:A:136:LEU:HA	1.78	0.40
1:A:307:LEU:HD12	3:A:2102:HOH:O	2.20	0.40
1:B:100:ILE:HG12	1:B:116:ASN:CG	2.39	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	301/304 (99%)	292 (97%)	9 (3%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	294/304 (97%)	283 (96%)	10 (3%)	1 (0%)	41	37
1	C	301/304 (99%)	290 (96%)	11 (4%)	0	100	100
1	D	294/304 (97%)	283 (96%)	10 (3%)	1 (0%)	41	37
All	All	1190/1216 (98%)	1148 (96%)	40 (3%)	2 (0%)	47	44

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	164	GLY
1	D	164	GLY

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	237/245 (97%)	232 (98%)	5 (2%)	53	57
1	B	233/245 (95%)	227 (97%)	6 (3%)	46	48
1	C	239/245 (98%)	237 (99%)	2 (1%)	81	86
1	D	235/245 (96%)	229 (97%)	6 (3%)	46	48
All	All	944/980 (96%)	925 (98%)	19 (2%)	55	58

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	48	GLU
1	A	57	ASP
1	A	80	GLN
1	A	136	LEU
1	A	142	VAL
1	B	57	ASP
1	B	80	GLN
1	B	111	ASP
1	B	125	SER

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Mol	Chain	Res	Type
1	B	145	LEU
1	B	197	ASP
1	C	80	GLN
1	C	109	ARG
1	D	57	ASP
1	D	80	GLN
1	D	86	GLU
1	D	109	ARG
1	D	226	GLU
1	D	233	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	80	GLN
1	A	130	HIS
1	A	233	GLN
1	A	325	GLN
1	B	80	GLN
1	B	159	GLN
1	B	185	GLN
1	B	199	GLN
1	C	80	GLN
1	C	140	ASN
1	C	159	GLN
1	C	199	GLN
1	C	233	GLN
1	D	80	GLN
1	D	107	GLN
1	D	233	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 [i](#)

Of 8 ligands modelled in this entry, 8 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 [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	303/304 (99%)	-0.45	1 (0%) 94 93	11, 19, 32, 54	0
1	B	298/304 (98%)	-0.42	1 (0%) 94 93	8, 20, 37, 50	0
1	C	303/304 (99%)	-0.48	0 100 100	9, 17, 29, 46	0
1	D	298/304 (98%)	-0.43	0 100 100	10, 20, 34, 54	0
All	All	1202/1216 (98%)	-0.45	2 (0%) 95 94	8, 19, 33, 54	0

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	100	ILE	3.2
1	A	329	ILE	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	CL	B	1332	1/1	0.93	0.36	64,64,64,64	0
2	CL	D	1331	1/1	0.97	0.14	36,36,36,36	0
2	CL	A	1333	1/1	0.99	0.15	31,31,31,31	0
2	CL	B	1331	1/1	0.99	0.07	20,20,20,20	0
2	CL	A	1331	1/1	0.99	0.09	18,18,18,18	0
2	CL	C	1332	1/1	0.99	0.09	26,26,26,26	0
2	CL	A	1332	1/1	0.99	0.07	23,23,23,23	0
2	CL	C	1331	1/1	1.00	0.09	11,11,11,11	0

6.5 Other polymers [i](#)

There are no such residues in this entry.