



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 08:21 AM UTC

PDB ID : 2J5Q / pdb_00002j5q
Title : 2.15 Å resolution structure of the wild type malate dehydrogenase from *Haloarcula marismortui* after first radiation burn (radiation damage series)
Authors : Fioravanti, E.; Vellieux, F.M.D.; Amara, P.; Madern, D.; Weik, M.
Deposited on : 2006-09-19
Resolution : 2.15 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	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

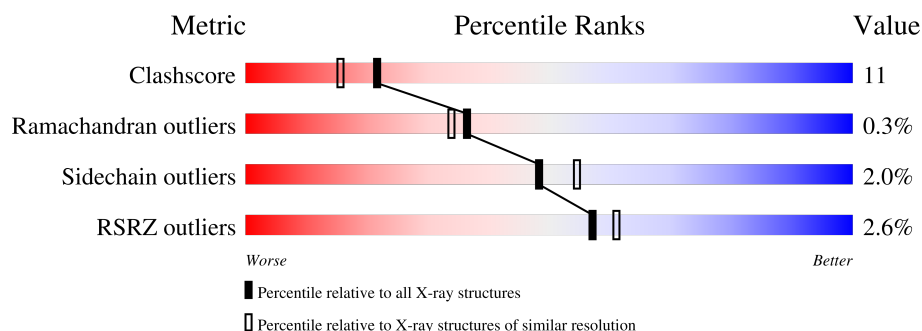
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.15 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	190562	2159 (2.16-2.16)
Ramachandran outliers	187476	2134 (2.16-2.16)
Sidechain outliers	187428	2133 (2.16-2.16)
RSRZ outliers	180081	2059 (2.16-2.16)

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	3% 75% 23% .
1	B	304	5% 73% 22% ..
1	C	304	79% 19% .
1	D	304	3% 75% 20% ..

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

There are 3 unique types of molecules in this entry. The entry contains 10082 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
1	A	303	Total	C	N	O	S	0	0	0
			2279	1407	391	477	4			
1	B	298	Total	C	N	O	S	0	0	0
			2239	1382	383	470	4			
1	C	303	Total	C	N	O	S	0	0	0
			2285	1409	393	479	4			
1	D	298	Total	C	N	O	S	0	0	0
			2248	1387	383	474	4			

- Molecule 2 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

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

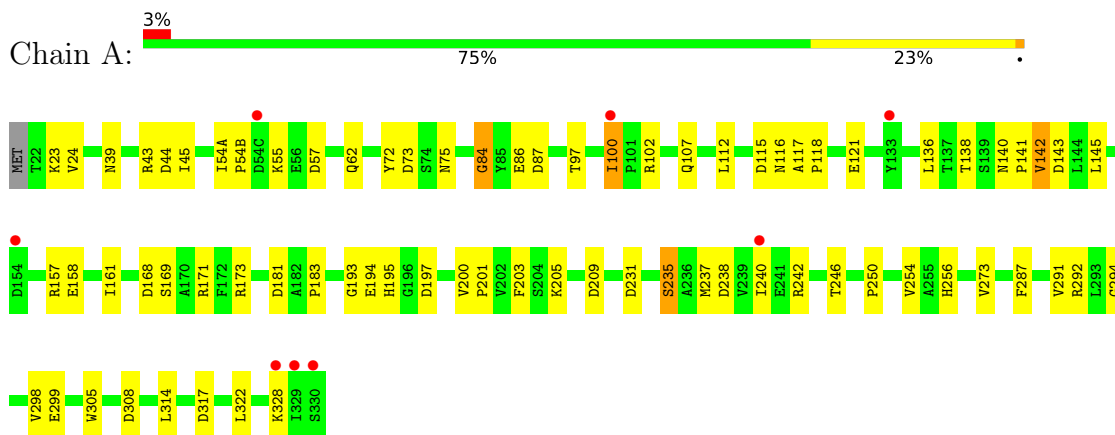
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	264	Total	O	0	0
			264	264		
3	B	229	Total	O	0	0
			229	229		
3	C	251	Total	O	0	0
			251	251		
3	D	279	Total	O	0	0
			279	279		

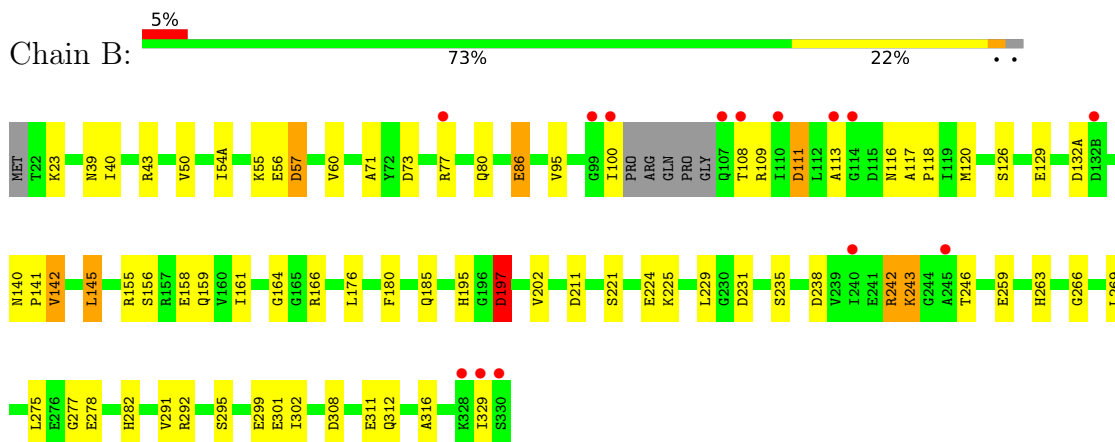
3 Residue-property plots

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.

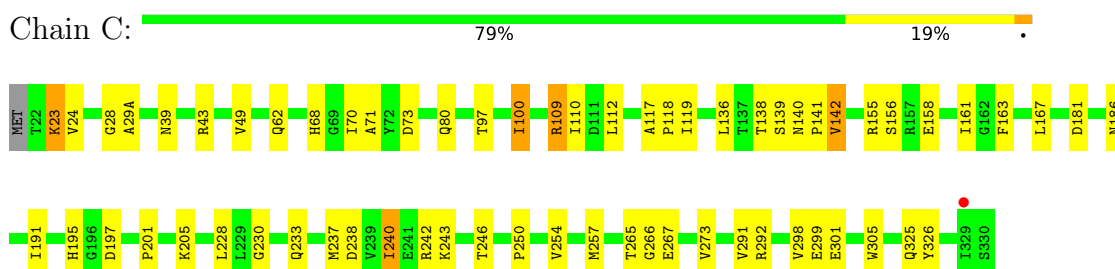
• Molecule 1: MALATE DEHYDROGENASE



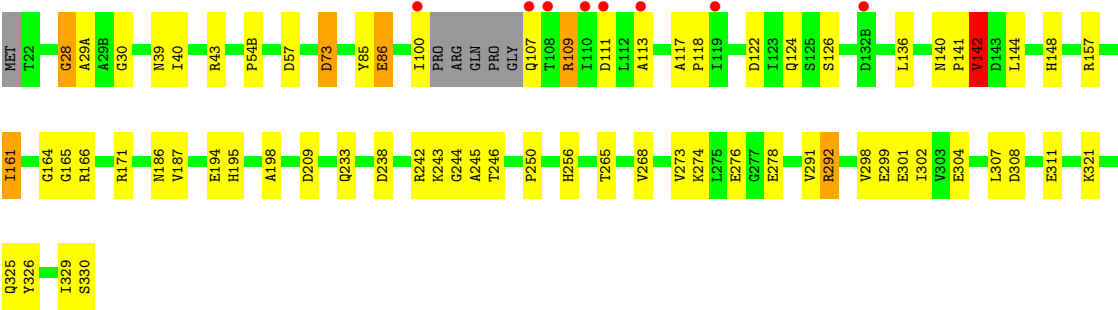
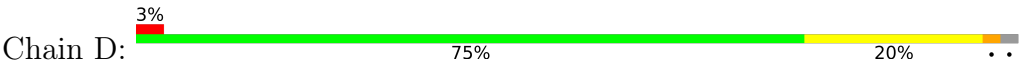
• Molecule 1: MALATE DEHYDROGENASE



• Molecule 1: MALATE DEHYDROGENASE



● Molecule 1: MALATE DEHYDROGENASE



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	126.98Å 114.32Å 124.12Å 90.00° 93.47° 90.00°	Depositor
Resolution (Å)	20.00 – 2.15 20.00 – 2.15	Depositor EDS
% Data completeness (in resolution range)	95.5 (20.00-2.15) 95.5 (20.00-2.15)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.90 (at 2.15Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.220 , 0.267 0.225 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	18.3	Xtriage
Anisotropy	0.688	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 58.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.27$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	10082	wwPDB-VP
Average B, all atoms (Å ²)	22.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.45% 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

5.1 Standard geometry

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	0.71	0/2316	1.05	5/3146 (0.2%)
1	B	0.67	0/2273	1.05	11/3085 (0.4%)
1	C	0.71	0/2322	1.09	15/3154 (0.5%)
1	D	0.69	0/2282	1.08	12/3097 (0.4%)
All	All	0.69	0/9193	1.07	43/12482 (0.3%)

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.

All (43) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	142	VAL	N-CA-C	9.68	120.53	110.36
1	C	142	VAL	N-CA-C	9.37	120.20	110.36
1	D	142	VAL	N-CA-C	8.07	118.85	110.62
1	B	243	LYS	N-CA-C	-7.35	103.35	111.36
1	D	161	ILE	N-CA-C	7.21	118.67	108.36
1	B	197	ASP	N-CA-C	7.05	119.81	111.71
1	B	142	VAL	N-CA-C	6.83	117.59	110.62
1	C	100	ILE	N-CA-C	-6.68	102.88	108.63
1	A	161	ILE	N-CA-C	6.53	117.31	108.17
1	A	84	GLY	N-CA-C	-6.30	103.94	112.57
1	B	242	ARG	N-CA-C	6.29	117.94	111.14
1	B	71	ALA	N-CA-C	6.24	118.89	111.71

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	243	LYS	N-CA-C	-6.00	104.67	112.23
1	D	28	GLY	N-CA-C	-5.99	106.38	115.00
1	C	161	ILE	N-CA-C	5.98	117.19	108.58
1	C	240	ILE	N-CA-C	5.88	116.66	110.72
1	D	186	ASN	N-CA-C	-5.86	105.79	113.17
1	D	73	ASP	N-CA-C	5.81	121.81	114.31
1	C	230	GLY	N-CA-C	-5.79	105.48	112.49
1	D	292	ARG	N-CA-C	-5.78	99.87	109.46
1	D	187	VAL	N-CA-C	5.67	116.28	107.99
1	B	316	ALA	N-CA-C	-5.61	105.25	111.36
1	C	71	ALA	N-CA-C	5.53	118.07	111.71
1	C	70	ILE	N-CA-C	5.50	117.40	112.17
1	B	80	GLN	N-CA-C	-5.50	102.74	110.50
1	C	112	LEU	N-CA-C	-5.49	105.20	111.07
1	A	197	ASP	N-CA-C	5.43	117.01	111.14
1	B	295	SER	N-CA-C	5.40	118.66	111.75
1	C	29(A)	ALA	N-CA-C	5.38	118.95	112.38
1	C	195	HIS	N-CA-C	-5.35	103.24	110.68
1	D	195	HIS	N-CA-C	-5.26	99.40	108.56
1	B	195	HIS	N-CA-C	-5.26	99.37	108.52
1	B	161	ILE	N-CA-C	5.25	115.86	108.36
1	D	304	GLU	N-CA-C	5.20	117.57	108.52
1	C	28	GLY	N-CA-C	-5.17	107.54	116.01
1	C	197	ASP	N-CA-C	5.16	117.64	111.71
1	A	157	ARG	N-CA-C	5.13	118.31	111.75
1	D	329	ILE	N-CA-C	-5.13	107.71	112.43
1	D	40	ILE	N-CA-C	-5.11	105.62	110.42
1	C	23	LYS	N-CA-C	5.07	117.66	109.40
1	C	186	ASN	N-CA-C	-5.04	106.72	112.92
1	D	265	THR	N-CA-C	5.03	117.66	111.82
1	B	231	ASP	N-CA-C	-5.00	105.71	111.07

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 ⓘ

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	52	0
1	B	2239	0	2105	50	0
1	C	2285	0	2153	40	0
1	D	2248	0	2114	55	0
2	A	3	0	0	3	0
2	B	2	0	0	3	0
2	C	2	0	0	2	0
2	D	1	0	0	2	0
3	A	264	0	0	5	0
3	B	229	0	0	8	0
3	C	251	0	0	4	0
3	D	279	0	0	8	0
All	All	10082	0	8517	190	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 11.

All (190) 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	1.47	0.92
1:D:86:GLU:CD	1:D:86:GLU:H	1.84	0.83
1:B:100:ILE:HG12	1:B:116:ASN:OD1	1.85	0.77
1:D:100:ILE:O	1:D:100:ILE:HG13	1.84	0.76
1:D:291:VAL:HG13	1:D:298:VAL:HG13	1.67	0.76
1:A:181:ASP:HB2	3:A:2142:HOH:O	1.88	0.74
1:D:109:ARG:CG	1:D:109:ARG:HH21	2.01	0.74
1:C:62:GLN:NE2	1:D:243:LYS:HE3	2.03	0.72
2:A:1333:CL:CL	3:B:2037:HOH:O	2.45	0.72
1:A:75:ASN:OD1	1:B:185:GLN:HG3	1.90	0.71
1:C:140:ASN:HA	1:C:142:VAL:N	2.07	0.69
1:A:100:ILE:HG12	1:A:116:ASN:OD1	1.91	0.69
1:B:308:ASP:OD1	1:B:311:GLU:HG3	1.92	0.68
2:C:1332:CL:CL	3:D:2057:HOH:O	2.48	0.67
1:D:109:ARG:HH21	1:D:109:ARG:HG3	1.60	0.67

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:246:THR:HG21	3:B:2109:HOH:O	1.94	0.66
1:B:95:VAL:HG13	3:B:2091:HOH:O	1.93	0.66
1:A:102:ARG:HB2	1:A:112:LEU:HD22	1.78	0.66
1:C:39:ASN:O	1:C:43:ARG:HG2	1.96	0.66
1:A:86:GLU:H	1:A:86:GLU:CD	2.05	0.65
1:D:308:ASP:OD1	1:D:311:GLU:HG3	1.96	0.65
1:D:117:ALA:HB3	1:D:118:PRO:HD3	1.79	0.65
1:A:328:LYS:O	1:A:328:LYS:HD3	1.96	0.64
1:B:86:GLU:H	1:B:86:GLU:CD	2.06	0.64
1:C:62:GLN:CD	1:D:243:LYS:HE3	2.22	0.64
1:D:109:ARG:HH21	1:D:109:ARG:CB	2.11	0.64
1:A:140:ASN:HB2	3:A:2114:HOH:O	1.99	0.63
1:D:321:LYS:O	1:D:325:GLN:HG3	1.98	0.63
1:B:166:ARG:NH2	2:B:1332:CL:CL	2.69	0.63
1:B:108:THR:OG1	1:B:111:ASP:HB2	1.99	0.62
1:C:191:ILE:HD13	1:C:201:PRO:HA	1.82	0.61
1:B:113:ALA:HB2	1:B:141:PRO:HG2	1.81	0.60
1:D:238:ASP:O	1:D:242:ARG:HB2	2.01	0.60
1:A:250:PRO:O	1:A:254:VAL:HG23	2.01	0.60
1:B:77:ARG:HA	1:B:77:ARG:HE	1.67	0.59
1:A:194:GLU:HB2	1:A:322:LEU:HD11	1.83	0.59
1:B:56:GLU:O	1:B:60:VAL:HG23	2.03	0.58
1:C:265:THR:OG1	1:C:267:GLU:HG3	2.04	0.58
1:B:141:PRO:HB2	3:B:2095:HOH:O	2.04	0.58
1:C:110:ILE:HD11	1:C:325:GLN:OE1	2.05	0.57
1:B:100:ILE:HG13	1:B:100:ILE:O	2.04	0.57
1:A:209:ASP:OD2	1:D:292:ARG:NH2	2.39	0.56
1:D:144:LEU:HG	3:D:2127:HOH:O	2.06	0.56
1:D:157:ARG:NE	1:D:276:GLU:OE1	2.40	0.55
1:A:237:MET:CE	1:A:240:ILE:HD12	2.37	0.55
1:C:233:GLN:HE21	1:C:233:GLN:HA	1.72	0.54
1:C:140:ASN:HA	1:C:141:PRO:C	2.32	0.54
1:A:314:LEU:O	1:A:317:ASP:HB2	2.08	0.54
1:D:122:ASP:HB3	3:D:2101:HOH:O	2.08	0.54
1:A:102:ARG:HD2	1:A:107:GLN:O	2.08	0.54
1:B:109:ARG:HD3	3:B:2222:HOH:O	2.09	0.53
1:A:308:ASP:OD1	1:A:308:ASP:C	2.51	0.53
1:A:205:LYS:HE3	1:A:305:TRP:HB3	1.89	0.53
1:C:100:ILE:HD11	1:C:119:ILE:HD11	1.91	0.53
1:C:273:VAL:HG21	1:C:298:VAL:HG21	1.89	0.53
1:D:268:VAL:HA	1:D:291:VAL:O	2.08	0.52

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:156:SER:OG	1:C:158:GLU:OE1	2.28	0.52
1:D:113:ALA:HB2	1:D:141:PRO:HG3	1.91	0.52
1:D:166:ARG:HH11	1:D:166:ARG:HG3	1.74	0.52
1:B:57:ASP:HB2	3:B:2019:HOH:O	2.10	0.52
1:B:155:ARG:HB3	1:B:159:GLN:HE21	1.74	0.52
1:A:140:ASN:HB3	3:A:2118:HOH:O	2.10	0.52
1:A:169:SER:O	1:A:173:ARG:HG3	2.10	0.52
1:B:235:SER:O	1:B:238:ASP:HB2	2.11	0.51
1:B:23:LYS:HE3	1:B:50:VAL:HG21	1.92	0.51
1:B:197:ASP:HB3	3:B:2224:HOH:O	2.10	0.51
1:D:246:THR:HG22	1:D:250:PRO:CD	2.40	0.51
1:C:117:ALA:HB3	1:C:118:PRO:HD3	1.92	0.51
1:D:109:ARG:HG3	1:D:109:ARG:NH2	2.23	0.51
1:D:166:ARG:HG3	1:D:166:ARG:NH1	2.26	0.51
1:D:256:HIS:HB2	2:D:1331:CL:CL	2.48	0.51
1:A:183:PRO:CD	1:D:292:ARG:HH12	2.24	0.51
1:A:237:MET:HE1	1:A:240:ILE:HD12	1.93	0.51
1:C:291:VAL:HB	1:C:298:VAL:HG13	1.93	0.51
1:C:233:GLN:HA	1:C:233:GLN:NE2	2.27	0.50
1:A:84:GLY:N	1:A:87:ASP:OD1	2.35	0.50
1:B:39:ASN:O	1:B:43:ARG:HG2	2.12	0.50
1:C:73:ASP:OD2	2:D:1331:CL:CL	2.66	0.50
1:A:273:VAL:HG11	1:A:298:VAL:HG21	1.93	0.50
1:C:110:ILE:HG13	3:C:2086:HOH:O	2.11	0.50
1:A:328:LYS:HD3	1:A:328:LYS:C	2.37	0.50
1:B:140:ASN:HA	1:B:142:VAL:N	2.27	0.49
1:C:246:THR:O	1:C:246:THR:HG23	2.12	0.49
1:A:44:ASP:OD1	1:A:75:ASN:N	2.39	0.49
1:A:54(A):ILE:HG12	1:A:55:LYS:HD2	1.95	0.49
1:C:266:GLY:O	1:C:292:ARG:HG3	2.12	0.49
1:A:256:HIS:ND1	2:A:1333:CL:CL	2.82	0.49
1:C:181:ASP:O	1:C:181:ASP:CG	2.56	0.49
1:C:68:HIS:CE1	1:D:171:ARG:HG2	2.47	0.49
1:D:291:VAL:HG11	1:D:298:VAL:HG22	1.95	0.49
1:B:155:ARG:HB3	1:B:159:GLN:NE2	2.28	0.49
1:A:117:ALA:N	1:A:118:PRO:HD2	2.27	0.48
1:D:140:ASN:HA	1:D:142:VAL:N	2.28	0.48
1:C:24:VAL:O	1:C:49:VAL:HA	2.14	0.48
1:B:299:GLU:O	1:B:301:GLU:HB3	2.11	0.48
1:A:291:VAL:HB	1:A:298:VAL:HG13	1.94	0.48
1:A:117:ALA:O	1:A:121:GLU:HG3	2.14	0.48

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:139:SER:O	1:C:142:VAL:HA	2.14	0.48
1:B:238:ASP:O	1:B:242:ARG:HB2	2.14	0.48
1:A:140:ASN:OD1	1:A:141:PRO:HA	2.15	0.47
1:D:54(B):PRO:HD2	3:D:2034:HOH:O	2.13	0.47
1:A:73:ASP:OD2	2:B:1332:CL:CL	2.69	0.47
1:B:259:GLU:HG3	1:B:263:HIS:CD2	2.50	0.47
1:A:238:ASP:O	1:A:242:ARG:HB2	2.14	0.47
1:D:246:THR:HG22	1:D:250:PRO:HD3	1.96	0.47
1:D:43:ARG:NH1	3:D:2014:HOH:O	2.48	0.47
1:B:269:LEU:O	1:B:291:VAL:HG22	2.15	0.47
1:A:143:ASP:OD1	1:A:287:PHE:HA	2.15	0.46
1:A:39:ASN:O	1:A:43:ARG:HG2	2.15	0.46
1:D:86:GLU:CD	1:D:86:GLU:N	2.62	0.46
1:B:120:MET:HG3	1:B:145:LEU:HG	1.97	0.46
1:D:136:LEU:HA	1:D:161:ILE:O	2.14	0.46
1:B:176:LEU:O	1:B:180:PHE:HD1	1.99	0.46
1:C:163:PHE:HD1	1:C:257:MET:SD	2.38	0.46
1:D:291:VAL:CG1	1:D:298:VAL:HG13	2.40	0.46
1:A:171:ARG:HH22	1:A:246:THR:HG21	1.79	0.46
1:D:242:ARG:NH2	1:D:242:ARG:HG3	2.30	0.46
1:A:294:GLY:HA3	1:A:299:GLU:OE2	2.16	0.46
1:A:23:LYS:HG2	1:A:24:VAL:N	2.31	0.46
1:B:202:VAL:O	1:B:202:VAL:HG12	2.15	0.46
1:A:72:TYR:HB2	2:B:1332:CL:CL	2.52	0.46
1:B:275:LEU:HD22	1:B:278:GLU:HB2	1.98	0.46
2:C:1332:CL:CL	1:D:73:ASP:OD2	2.71	0.45
1:D:107:GLN:HE21	1:D:111:ASP:CB	2.30	0.45
1:A:100:ILE:HD11	1:A:115:ASP:HB3	1.98	0.45
1:C:155:ARG:HG3	3:C:2100:HOH:O	2.16	0.45
1:D:124:GLN:NE2	3:D:2107:HOH:O	2.50	0.45
1:B:141:PRO:O	1:B:145:LEU:HB2	2.16	0.45
1:A:193:GLY:HA3	1:A:200:VAL:HG23	1.99	0.45
1:D:326:TYR:O	1:D:330:SER:HB2	2.17	0.45
1:C:205:LYS:HE3	1:C:305:TRP:HB3	1.99	0.45
1:D:117:ALA:HB1	1:D:148:HIS:CD2	2.52	0.45
1:B:259:GLU:HG3	1:B:263:HIS:HD2	1.81	0.45
1:C:238:ASP:O	1:C:242:ARG:HB2	2.17	0.45
1:C:237:MET:HE2	1:C:240:ILE:HD12	1.99	0.44
1:D:28:GLY:O	1:D:30:GLY:N	2.42	0.44
1:A:107:GLN:NE2	1:A:115:ASP:OD2	2.48	0.44
1:D:244:GLY:O	1:D:245:ALA:HB2	2.18	0.44

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:168:ASP:CG	1:A:195:HIS:HD1	2.26	0.44
1:A:292:ARG:NH1	1:D:209:ASP:OD2	2.49	0.44
1:A:102:ARG:HA	1:A:112:LEU:HD13	2.00	0.43
1:C:136:LEU:HD23	1:C:254:VAL:HG13	2.01	0.43
1:C:237:MET:CE	1:C:240:ILE:HD12	2.48	0.43
1:B:140:ASN:HA	1:B:141:PRO:C	2.44	0.43
1:C:141:PRO:HB2	3:C:2114:HOH:O	2.17	0.43
1:B:211:ASP:OD2	1:C:205:LYS:NZ	2.43	0.42
1:C:299:GLU:O	1:C:301:GLU:HB3	2.19	0.42
1:A:201:PRO:HB2	1:A:203:PHE:CE2	2.54	0.42
1:D:242:ARG:HG3	1:D:242:ARG:HH21	1.84	0.42
1:B:312:GLN:HG2	3:B:2085:HOH:O	2.19	0.42
1:D:109:ARG:HH21	1:D:109:ARG:HB3	1.84	0.42
1:D:157:ARG:HD2	1:D:274:LYS:O	2.19	0.42
1:A:97:THR:HG22	1:A:138:THR:OG1	2.19	0.42
1:A:145:LEU:HA	1:A:145:LEU:HD23	1.76	0.42
1:A:158:GLU:H	1:A:158:GLU:CD	2.26	0.42
1:B:126:SER:O	1:B:129:GLU:HG2	2.20	0.42
1:B:277:GLY:N	1:B:282:HIS:O	2.41	0.42
2:A:1333:CL:CL	1:B:73:ASP:OD2	2.75	0.42
1:B:23:LYS:HE3	1:B:50:VAL:CG2	2.50	0.42
1:B:54(A):ILE:HG12	1:B:55:LYS:HD2	2.01	0.42
1:D:39:ASN:O	1:D:43:ARG:HG2	2.20	0.42
1:B:77:ARG:HA	1:B:77:ARG:NE	2.31	0.42
1:B:109:ARG:HH11	1:B:109:ARG:HG3	1.84	0.42
1:B:117:ALA:N	1:B:118:PRO:HD2	2.34	0.42
1:D:165:GLY:HA3	3:D:2161:HOH:O	2.19	0.42
1:D:299:GLU:O	1:D:301:GLU:HB3	2.18	0.42
1:D:307:LEU:HB3	1:D:311:GLU:HB2	2.02	0.41
1:C:228:LEU:HA	1:C:228:LEU:HD23	1.82	0.41
1:B:278:GLU:HG2	1:B:302:ILE:HD12	2.01	0.41
1:B:225:LYS:O	1:B:229:LEU:HG	2.21	0.41
1:C:109:ARG:HD3	3:C:2153:HOH:O	2.20	0.41
1:D:117:ALA:N	1:D:118:PRO:CD	2.83	0.41
1:A:62:GLN:NE2	1:B:243:LYS:HE3	2.36	0.41
1:A:235:SER:O	1:A:238:ASP:HB2	2.21	0.41
1:B:329:ILE:O	1:B:329:ILE:HG13	2.20	0.41
1:B:221:SER:OG	1:B:224:GLU:HG3	2.20	0.41
1:A:142:VAL:HG22	3:A:2116:HOH:O	2.19	0.41
1:B:266:GLY:O	1:B:292:ARG:HG3	2.21	0.41
1:C:97:THR:HG22	1:C:138:THR:OG1	2.21	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:250:PRO:O	1:C:254:VAL:HG23	2.21	0.41
1:D:85:TYR:CG	1:D:126:SER:HB3	2.57	0.41
1:C:23:LYS:HG2	1:C:24:VAL:N	2.35	0.40
1:D:43:ARG:CZ	3:D:2015:HOH:O	2.68	0.40
1:D:273:VAL:HG11	1:D:298:VAL:HG21	2.02	0.40
1:A:54(A):ILE:HA	1:A:54(B):PRO:HD3	1.95	0.40
1:A:231:ASP:CG	3:A:2193:HOH:O	2.64	0.40
1:C:110:ILE:CD1	1:C:325:GLN:HG2	2.52	0.40
1:D:278:GLU:CD	1:D:302:ILE:HD12	2.46	0.40
1:B:156:SER:OG	1:B:158:GLU:CD	2.65	0.40
1:D:194:GLU:OE1	1:D:198:ALA:HB3	2.21	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%)	291 (97%)	10 (3%)	0	100	100
1	B	294/304 (97%)	282 (96%)	10 (3%)	2 (1%)	18	13
1	C	301/304 (99%)	289 (96%)	12 (4%)	0	100	100
1	D	294/304 (97%)	283 (96%)	9 (3%)	2 (1%)	18	13
All	All	1190/1216 (98%)	1145 (96%)	41 (3%)	4 (0%)	36	34

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	132(A)	ASP
1	B	164	GLY
1	D	164	GLY
1	D	29(A)	ALA

5.3.2 Protein sidechains ⓘ

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%)	47	52
1	B	233/245 (95%)	227 (97%)	6 (3%)	40	43
1	C	239/245 (98%)	236 (99%)	3 (1%)	61	68
1	D	235/245 (96%)	230 (98%)	5 (2%)	47	52
All	All	944/980 (96%)	925 (98%)	19 (2%)	48	54

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

Mol	Chain	Res	Type
1	A	45	ILE
1	A	57	ASP
1	A	100	ILE
1	A	136	LEU
1	A	235	SER
1	B	40	ILE
1	B	57	ASP
1	B	86	GLU
1	B	111	ASP
1	B	145	LEU
1	B	197	ASP
1	C	80	GLN
1	C	109	ARG
1	C	167	LEU
1	D	57	ASP
1	D	86	GLU
1	D	109	ARG
1	D	142	VAL
1	D	233	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	80	GLN
1	A	185	GLN
1	A	233	GLN
1	A	325	GLN
1	B	80	GLN
1	B	140	ASN
1	B	159	GLN
1	C	80	GLN
1	C	159	GLN
1	C	199	GLN
1	C	233	GLN
1	C	325	GLN
1	D	80	GLN
1	D	107	GLN
1	D	131	ASN
1	D	159	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 oligosaccharides 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

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 ⓘ

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	303/304 (99%)	0.51	8 (2%) 57 61	9, 20, 32, 55	0
1	B	298/304 (98%)	0.57	14 (4%) 36 41	10, 20, 39, 51	0
1	C	303/304 (99%)	0.44	1 (0%) 90 91	8, 18, 30, 46	0
1	D	298/304 (98%)	0.59	8 (2%) 56 60	10, 20, 36, 54	0
All	All	1202/1216 (98%)	0.53	31 (2%) 57 61	8, 20, 34, 55	0

All (31) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	100	ILE	5.7
1	B	100	ILE	5.4
1	A	330	SER	3.4
1	B	110	ILE	3.4
1	B	330	SER	3.3
1	D	110	ILE	3.3
1	A	329	ILE	3.2
1	D	107	GLN	3.2
1	C	329	ILE	2.9
1	A	328	LYS	2.8
1	B	329	ILE	2.8
1	D	108	THR	2.7
1	B	132(B)	ASP	2.6
1	B	108	THR	2.6
1	D	132(B)	ASP	2.6
1	A	100	ILE	2.5
1	B	245	ALA	2.3
1	B	114	GLY	2.3
1	A	54(C)	ASP	2.3
1	D	113	ALA	2.3
1	D	111	ASP	2.3

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	B	328	LYS	2.3
1	D	119	ILE	2.3
1	B	240	ILE	2.2
1	A	154	ASP	2.2
1	B	107	GLN	2.2
1	A	240	ILE	2.2
1	B	99	GLY	2.1
1	A	133	TYR	2.1
1	B	113	ALA	2.0
1	B	77	ARG	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 oligosaccharides 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($\text{\AA}^2$)	Q<0.9
2	CL	B	1332	1/1	0.85	0.45	68,68,68,68	0
2	CL	A	1333	1/1	0.94	0.28	37,37,37,37	0
2	CL	D	1331	1/1	0.94	0.28	41,41,41,41	0
2	CL	C	1332	1/1	0.97	0.26	31,31,31,31	0
2	CL	B	1331	1/1	0.97	0.11	21,21,21,21	0
2	CL	A	1332	1/1	0.98	0.12	24,24,24,24	0
2	CL	A	1331	1/1	0.98	0.07	22,22,22,22	0
2	CL	C	1331	1/1	0.99	0.11	13,13,13,13	0

6.5 Other polymers [i](#)

There are no such residues in this entry.