



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 05:15 AM UTC

PDB ID : 3VTZ / pdb_00003vtz
Title : Structure of Thermoplasma volcanium aldohexose dehydrogenase
Authors : Yasutake, Y.; Nishioka, T.; Tamura, T.
Deposited on : 2012-06-12
Resolution : 2.30 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

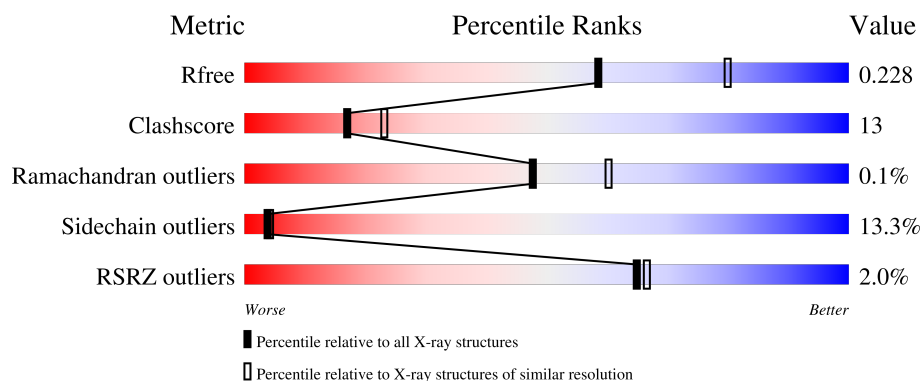
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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}	180053	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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	269	
1	B	269	
1	C	269	
1	D	269	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 7720 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 Glucose 1-dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	253	Total	C	N	O	S	0	0	0
			1912	1206	327	370	9			
1	B	253	Total	C	N	O	S	0	0	0
			1913	1208	328	368	9			
1	C	252	Total	C	N	O	S	0	0	0
			1904	1202	326	367	9			
1	D	252	Total	C	N	O	S	0	0	0
			1904	1202	326	367	9			

There are 32 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-7	MET	-	expression tag	UNP Q97CM7
A	-6	GLY	-	expression tag	UNP Q97CM7
A	-5	HIS	-	expression tag	UNP Q97CM7
A	-4	HIS	-	expression tag	UNP Q97CM7
A	-3	HIS	-	expression tag	UNP Q97CM7
A	-2	HIS	-	expression tag	UNP Q97CM7
A	-1	HIS	-	expression tag	UNP Q97CM7
A	0	HIS	-	expression tag	UNP Q97CM7
B	-7	MET	-	expression tag	UNP Q97CM7
B	-6	GLY	-	expression tag	UNP Q97CM7
B	-5	HIS	-	expression tag	UNP Q97CM7
B	-4	HIS	-	expression tag	UNP Q97CM7
B	-3	HIS	-	expression tag	UNP Q97CM7
B	-2	HIS	-	expression tag	UNP Q97CM7
B	-1	HIS	-	expression tag	UNP Q97CM7
B	0	HIS	-	expression tag	UNP Q97CM7
C	-7	MET	-	expression tag	UNP Q97CM7
C	-6	GLY	-	expression tag	UNP Q97CM7
C	-5	HIS	-	expression tag	UNP Q97CM7
C	-4	HIS	-	expression tag	UNP Q97CM7
C	-3	HIS	-	expression tag	UNP Q97CM7

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-2	HIS	-	expression tag	UNP Q97CM7
C	-1	HIS	-	expression tag	UNP Q97CM7
C	0	HIS	-	expression tag	UNP Q97CM7
D	-7	MET	-	expression tag	UNP Q97CM7
D	-6	GLY	-	expression tag	UNP Q97CM7
D	-5	HIS	-	expression tag	UNP Q97CM7
D	-4	HIS	-	expression tag	UNP Q97CM7
D	-3	HIS	-	expression tag	UNP Q97CM7
D	-2	HIS	-	expression tag	UNP Q97CM7
D	-1	HIS	-	expression tag	UNP Q97CM7
D	0	HIS	-	expression tag	UNP Q97CM7

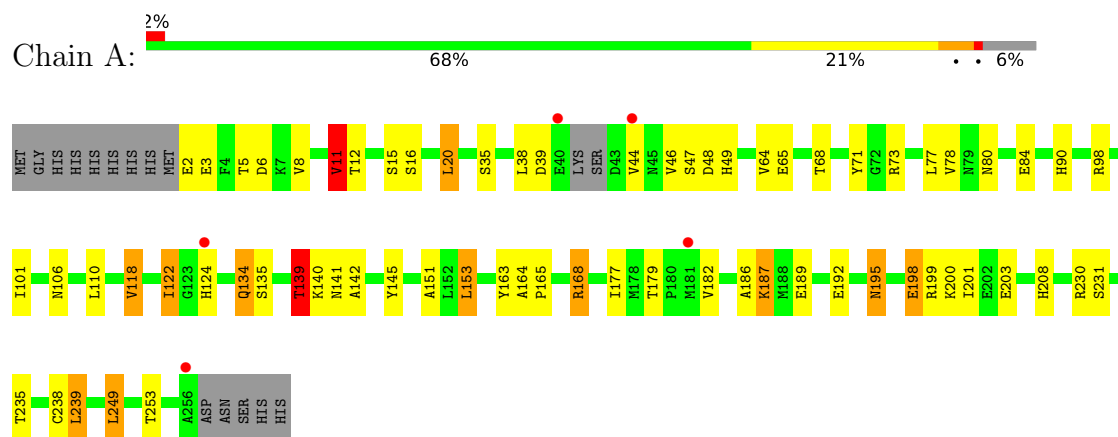
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	19	Total	O	0	0
			19	19		
2	B	25	Total	O	0	0
			25	25		
2	C	21	Total	O	0	0
			21	21		
2	D	22	Total	O	0	0
			22	22		

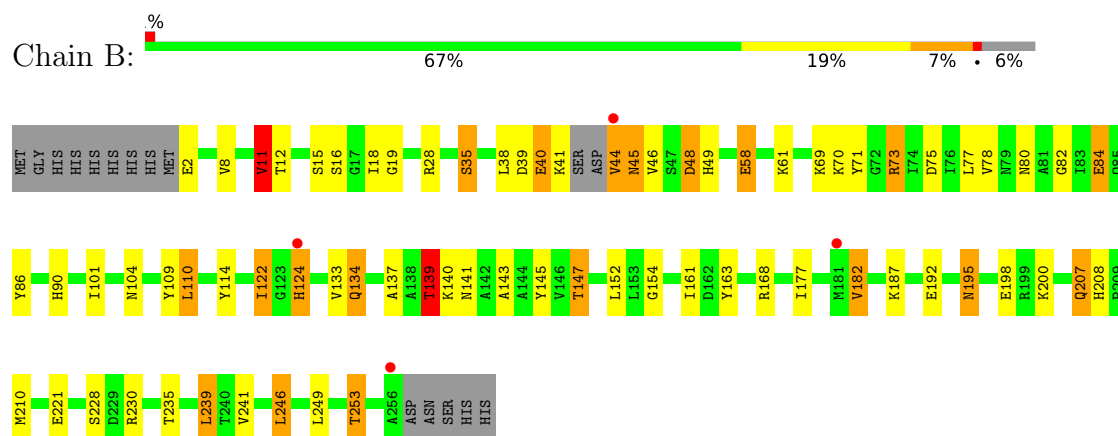
3 Residue-property plots [i](#)

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.

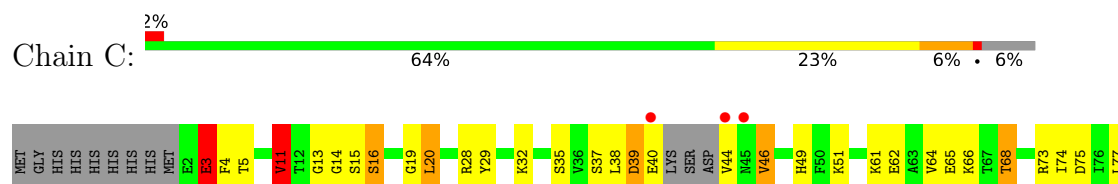
• Molecule 1: Glucose 1-dehydrogenase

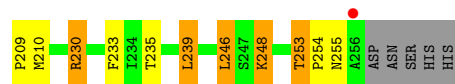
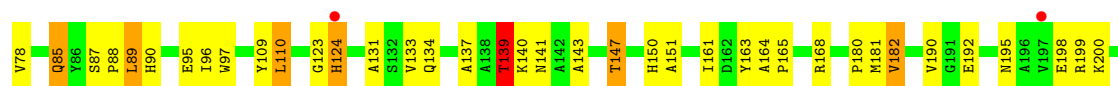


• Molecule 1: Glucose 1-dehydrogenase

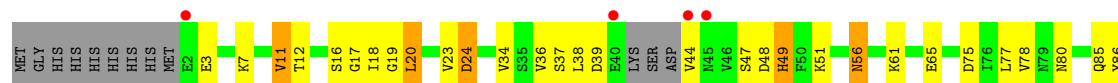


• Molecule 1: Glucose 1-dehydrogenase





● Molecule 1: Glucose 1-dehydrogenase



4 Data and refinement statistics

Property	Value	Source
Space group	P 31	Depositor
Cell constants a, b, c, α , β , γ	110.02Å 110.02Å 77.17Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	44.79 – 2.30 44.79 – 2.30	Depositor EDS
% Data completeness (in resolution range)	99.8 (44.79-2.30) 99.8 (44.79-2.30)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.34 (at 2.29Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.167 , 0.232 0.164 , 0.228	Depositor DCC
R_{free} test set	2347 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	32.2	Xtriage
Anisotropy	0.025	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.41 , 41.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.017 for -h,-k,l 0.024 for h,-h-k,-l 0.023 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	7720	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.77% 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

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	1.22	2/1941 (0.1%)	1.29	13/2634 (0.5%)
1	B	1.27	5/1942 (0.3%)	1.27	6/2634 (0.2%)
1	C	1.18	5/1933 (0.3%)	1.27	12/2623 (0.5%)
1	D	1.20	6/1933 (0.3%)	1.27	7/2623 (0.3%)
All	All	1.22	18/7749 (0.2%)	1.28	38/10514 (0.4%)

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	A	0	1

All (18) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	208	HIS	CG-CD2	7.41	1.44	1.35
1	D	131	ALA	CA-C	7.18	1.57	1.53
1	C	124	HIS	CG-CD2	6.67	1.43	1.35
1	B	208	HIS	CG-CD2	6.54	1.43	1.35
1	D	150	HIS	CG-CD2	6.53	1.43	1.35
1	D	150	HIS	ND1-CE1	6.47	1.39	1.32
1	D	116	ILE	C-O	-6.31	1.19	1.24
1	B	124	HIS	CG-CD2	6.19	1.42	1.35
1	C	150	HIS	CG-CD2	6.00	1.42	1.35
1	B	208	HIS	CE1-NE2	5.91	1.38	1.32
1	B	124	HIS	CE1-NE2	5.74	1.38	1.32
1	A	168	ARG	CD-NE	-5.73	1.38	1.46
1	C	124	HIS	CE1-NE2	5.34	1.37	1.32
1	D	124	HIS	CG-CD2	5.11	1.41	1.35
1	C	124	HIS	ND1-CE1	5.10	1.37	1.32
1	D	49	HIS	CG-CD2	5.09	1.41	1.35

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	150	HIS	ND1-CE1	5.05	1.37	1.32
1	B	161	ILE	C-O	-5.01	1.17	1.24

All (38) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	208	HIS	N-CA-C	-7.86	100.17	110.07
1	A	139	THR	N-CA-CB	-7.66	98.81	110.45
1	D	24	ASP	N-CA-C	6.84	118.81	111.36
1	A	168	ARG	NE-CZ-NH2	-6.78	113.10	119.20
1	B	139	THR	N-CA-CB	-6.76	100.17	110.45
1	B	154	GLY	N-CA-C	-6.58	104.35	112.77
1	D	131	ALA	N-CA-C	-6.49	102.52	108.75
1	C	14	GLY	N-CA-C	6.49	121.80	114.67
1	A	208	HIS	N-CA-C	-6.45	101.95	110.07
1	A	118	VAL	N-CA-C	-6.33	104.33	110.72
1	A	189	GLU	CA-C-N	-6.27	114.39	121.85
1	A	189	GLU	C-N-CA	-6.27	114.39	121.85
1	B	11	VAL	CB-CA-C	6.05	118.58	110.42
1	A	168	ARG	CD-NE-CZ	6.01	132.82	124.40
1	C	131	ALA	N-CA-C	-6.00	103.86	108.78
1	D	139	THR	N-CA-CB	-5.79	101.65	110.45
1	A	11	VAL	CB-CA-C	5.75	118.19	110.42
1	D	168	ARG	CG-CD-NE	-5.65	99.57	112.00
1	C	89	LEU	N-CA-C	5.60	117.83	111.11
1	B	139	THR	N-CA-C	-5.55	103.06	110.55
1	A	179	THR	CA-C-N	-5.54	113.96	119.56
1	A	179	THR	C-N-CA	-5.54	113.96	119.56
1	C	46	VAL	N-CA-C	-5.54	103.95	110.21
1	A	106	ASN	N-CA-C	5.51	117.29	111.28
1	D	18	ILE	N-CA-C	5.47	116.25	110.72
1	C	139	THR	N-CA-CB	-5.34	102.18	110.46
1	C	87	SER	CA-C-N	-5.32	115.10	120.21
1	C	87	SER	C-N-CA	-5.32	115.10	120.21
1	D	130	ILE	CB-CA-C	-5.29	106.19	111.80
1	C	123	GLY	N-CA-C	5.27	120.76	114.48
1	C	39	ASP	N-CA-C	5.22	121.93	110.80
1	A	101	ILE	N-CA-C	-5.22	105.62	110.53
1	A	39	ASP	N-CA-C	5.19	117.51	110.35
1	C	11	VAL	CB-CA-C	5.14	118.07	110.77
1	B	101	ILE	N-CA-C	-5.12	105.55	110.72
1	C	3	GLU	N-CA-C	5.11	117.59	111.71

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	168	ARG	CG-CD-NE	-5.10	100.78	112.00
1	D	195	ASN	N-CA-C	5.04	117.51	111.71

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	44	VAL	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	1912	0	1944	48	0
1	B	1913	0	1953	62	0
1	C	1904	0	1940	62	0
1	D	1904	0	1940	61	0
2	A	19	0	0	1	0
2	B	25	0	0	0	0
2	C	21	0	0	0	0
2	D	22	0	0	0	0
All	All	7720	0	7777	195	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 13.

All (195) 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:235:THR:HG21	1:C:210:MET:HE2	1.37	1.05
1:C:230:ARG:HH11	1:C:230:ARG:HG2	1.15	1.04
1:A:164:ALA:HB1	1:D:210:MET:HE2	1.45	0.98
1:C:230:ARG:HH11	1:C:230:ARG:CG	1.78	0.96
1:B:168:ARG:HG2	1:C:210:MET:HE1	1.46	0.94
1:C:230:ARG:HG2	1:C:230:ARG:NH1	1.87	0.86

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:143:ALA:O	1:B:147:THR:HG23	1.76	0.85
1:B:235:THR:CG2	1:C:210:MET:HE2	2.06	0.85
1:C:134:GLN:NE2	1:C:139:THR:HB	1.93	0.83
1:A:163:TYR:OH	1:C:90:HIS:HD2	1.62	0.82
1:B:163:TYR:OH	1:D:90:HIS:HD2	1.64	0.80
1:C:35:SER:HB3	1:C:49:HIS:HD2	1.46	0.80
1:C:143:ALA:O	1:C:147:THR:HG23	1.82	0.80
1:B:90:HIS:HD2	1:D:163:TYR:OH	1.66	0.79
1:B:58:GLU:H	1:B:58:GLU:CD	1.93	0.77
1:B:134:GLN:HE22	1:B:139:THR:H	1.34	0.76
1:A:187:LYS:HE2	1:A:192:GLU:OE2	1.84	0.75
1:A:134:GLN:HE22	1:A:139:THR:H	1.35	0.74
1:D:134:GLN:NE2	1:D:139:THR:HB	2.02	0.73
1:A:235:THR:HG21	1:D:210:MET:HE3	1.71	0.72
1:B:15:SER:HB3	1:B:41:LYS:HG2	1.71	0.72
1:A:15:SER:HA	1:A:20:LEU:HD13	1.72	0.71
1:B:239:LEU:HD11	1:C:239:LEU:HD11	1.72	0.71
1:A:164:ALA:CB	1:D:210:MET:HE2	2.21	0.70
1:C:134:GLN:HE22	1:C:139:THR:HB	1.58	0.69
1:C:140:LYS:O	1:D:253:THR:HG21	1.92	0.69
1:C:35:SER:HB2	1:C:46:VAL:HG11	1.75	0.68
1:A:139:THR:HG23	1:A:140:LYS:O	1.94	0.68
1:B:139:THR:HG23	1:B:140:LYS:O	1.94	0.68
1:C:62:GLU:O	1:C:66:LYS:HG3	1.95	0.67
1:B:15:SER:CB	1:B:41:LYS:HG2	2.25	0.66
1:B:133:VAL:HG12	1:B:246:LEU:HD22	1.77	0.66
1:C:16:SER:O	1:C:20:LEU:HB2	1.95	0.66
1:A:35:SER:HB2	1:A:49:HIS:HD2	1.61	0.66
1:C:253:THR:HG21	1:D:140:LYS:O	1.97	0.65
1:D:16:SER:HB2	1:D:19:GLY:H	1.62	0.65
1:D:197:VAL:O	1:D:201:ILE:HG12	1.97	0.64
1:B:35:SER:OG	1:B:49:HIS:HD2	1.82	0.63
1:A:84:GLU:HG3	1:A:145:TYR:CD1	2.33	0.63
1:C:133:VAL:HG12	1:C:246:LEU:HD22	1.81	0.62
1:D:183:ILE:HG22	1:D:187:LYS:HE2	1.81	0.62
1:D:134:GLN:NE2	1:D:139:THR:H	1.98	0.61
1:C:230:ARG:CG	1:C:230:ARG:NH1	2.48	0.61
1:A:141:ASN:HD22	1:B:253:THR:CG2	2.14	0.60
1:B:90:HIS:CD2	1:D:163:TYR:OH	2.54	0.60
1:A:165:PRO:HA	1:D:210:MET:CE	2.32	0.60
1:B:210:MET:HE2	1:C:164:ALA:HB1	1.84	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:182:VAL:HG12	1:A:201:ILE:HD12	1.85	0.59
1:D:39:ASP:O	1:D:49:HIS:HE1	1.84	0.59
1:D:86:TYR:OH	1:D:181:MET:HE1	2.02	0.59
1:D:61:LYS:O	1:D:65:GLU:HG3	2.02	0.59
1:D:85:GLN:HB3	1:D:100:ILE:HD11	1.84	0.59
1:D:181:MET:O	1:D:181:MET:HE2	2.04	0.58
1:B:163:TYR:OH	1:D:90:HIS:CD2	2.53	0.58
1:B:210:MET:HE3	1:C:235:THR:HG21	1.85	0.58
1:B:39:ASP:O	1:B:49:HIS:HE1	1.86	0.57
1:D:196:ALA:HA	1:D:199:ARG:CZ	2.34	0.57
1:A:153:LEU:HG	1:A:238:CYS:SG	2.44	0.57
1:A:35:SER:OG	1:A:46:VAL:HG11	2.04	0.57
1:A:90:HIS:HD2	1:C:163:TYR:OH	1.88	0.57
1:C:16:SER:HB3	1:C:19:GLY:H	1.70	0.56
1:C:35:SER:HB3	1:C:49:HIS:CD2	2.34	0.56
1:D:16:SER:O	1:D:20:LEU:HB2	2.05	0.56
1:D:134:GLN:HE22	1:D:139:THR:HB	1.70	0.56
1:A:239:LEU:HD11	1:D:239:LEU:HD11	1.89	0.55
1:B:61:LYS:HG3	1:B:114:TYR:CE1	2.42	0.55
1:A:253:THR:HG21	1:B:139:THR:CG2	2.37	0.55
1:D:205:GLY:N	1:D:213:ILE:HD12	2.21	0.55
1:D:85:GLN:CB	1:D:100:ILE:HD11	2.37	0.54
1:B:84:GLU:HG2	1:B:145:TYR:CD1	2.42	0.54
1:A:35:SER:CB	1:A:49:HIS:HD2	2.21	0.54
1:B:45:ASN:H	1:B:45:ASN:ND2	2.05	0.54
1:C:209:PRO:HG3	1:C:248:LYS:HD3	1.91	0.53
1:D:85:GLN:HB3	1:D:100:ILE:CD1	2.39	0.53
1:A:11:VAL:HB	1:A:78:VAL:HB	1.91	0.53
1:D:205:GLY:CA	1:D:213:ILE:HD12	2.39	0.53
1:A:35:SER:HB2	1:A:49:HIS:CD2	2.43	0.52
1:B:11:VAL:HB	1:B:78:VAL:HB	1.91	0.52
1:D:196:ALA:HA	1:D:199:ARG:NH2	2.23	0.52
1:B:235:THR:CG2	1:C:210:MET:CE	2.84	0.52
1:C:97:TRP:HH2	1:C:147:THR:HG21	1.75	0.52
1:A:16:SER:O	1:A:20:LEU:HB2	2.10	0.52
1:B:73:ARG:HD2	1:B:75:ASP:OD2	2.09	0.52
1:A:195:ASN:HB2	1:A:198:GLU:HG2	1.92	0.52
1:D:215:ARG:HB3	1:D:217:GLU:OE2	2.09	0.52
1:C:195:ASN:O	1:C:199:ARG:HD3	2.09	0.52
1:A:249:LEU:HD13	1:B:137:ALA:HB3	1.92	0.51
1:B:39:ASP:HB3	1:B:41:LYS:H	1.75	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:47:SER:HB2	2:A:319:HOH:O	2.10	0.51
1:B:210:MET:HE2	1:C:165:PRO:HA	1.91	0.51
1:B:109:TYR:CE2	1:B:110:LEU:HD13	2.46	0.51
1:A:135:SER:HB3	1:A:153:LEU:HD22	1.91	0.51
1:C:44:VAL:HG12	1:C:44:VAL:O	2.11	0.51
1:B:210:MET:HE2	1:C:164:ALA:CB	2.41	0.50
1:A:140:LYS:O	1:B:253:THR:HG21	2.11	0.50
1:B:58:GLU:CD	1:B:58:GLU:N	2.68	0.50
1:B:8:VAL:HG21	1:B:71:TYR:HB3	1.93	0.50
1:C:181:MET:HE2	1:C:182:VAL:HG23	1.93	0.50
1:A:151:ALA:HB2	1:C:151:ALA:HB2	1.93	0.50
1:B:230:ARG:NH1	1:C:29:TYR:OH	2.45	0.50
1:A:141:ASN:HD22	1:B:253:THR:HG22	1.77	0.49
1:A:253:THR:HG21	1:B:139:THR:HG23	1.93	0.49
1:A:164:ALA:HB1	1:D:210:MET:CE	2.32	0.49
1:D:80:ASN:HD22	1:D:130:ILE:HG13	1.76	0.49
1:C:13:GLY:O	1:C:16:SER:HB2	2.13	0.49
1:C:35:SER:CB	1:C:49:HIS:HD2	2.21	0.49
1:C:109:TYR:CE2	1:C:110:LEU:HD13	2.47	0.49
1:D:195:ASN:HA	1:D:198:GLU:OE1	2.12	0.49
1:A:5:THR:O	1:A:6:ASP:HB2	2.12	0.49
1:C:89:LEU:C	1:C:89:LEU:HD23	2.38	0.49
1:A:164:ALA:CB	1:D:210:MET:CE	2.91	0.49
1:A:68:THR:HG21	1:A:118:VAL:HG21	1.93	0.49
1:B:35:SER:OG	1:B:49:HIS:CD2	2.64	0.49
1:D:11:VAL:HB	1:D:78:VAL:HB	1.94	0.48
1:C:73:ARG:HD3	1:C:75:ASP:OD1	2.13	0.48
1:D:196:ALA:CB	1:D:199:ARG:NH2	2.76	0.48
1:C:40:GLU:OE1	1:C:51:LYS:HD2	2.13	0.48
1:B:82:GLY:HA2	1:B:104:ASN:OD1	2.14	0.48
1:C:37:SER:O	1:C:51:LYS:HA	2.13	0.48
1:D:95:GLU:O	1:D:95:GLU:OE2	2.31	0.48
1:C:3:GLU:O	1:C:3:GLU:HG3	2.12	0.47
1:D:37:SER:O	1:D:51:LYS:HA	2.14	0.47
1:D:12:THR:O	1:D:80:ASN:HB3	2.15	0.47
1:D:206:ARG:HB2	1:D:206:ARG:NH1	2.30	0.47
1:B:143:ALA:O	1:B:147:THR:CG2	2.57	0.47
1:B:195:ASN:HD22	1:B:195:ASN:H	1.62	0.47
1:D:56:ASN:C	1:D:56:ASN:HD22	2.22	0.47
1:D:132:SER:HA	1:D:174:PRO:HD2	1.96	0.47
1:C:195:ASN:HB2	1:C:198:GLU:OE2	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:196:ALA:HB2	1:D:199:ARG:HH22	1.80	0.47
1:B:134:GLN:HE22	1:B:139:THR:HB	1.80	0.46
1:C:39:ASP:O	1:C:40:GLU:HB2	2.15	0.46
1:C:64:VAL:O	1:C:68:THR:HG23	2.15	0.46
1:D:193:ASP:O	1:D:194:GLU:C	2.58	0.46
1:A:186:ALA:HB2	1:A:201:ILE:CD1	2.47	0.45
1:D:34:VAL:HG22	1:D:48:ASP:HB2	1.98	0.45
1:B:187:LYS:NZ	1:B:192:GLU:OE2	2.40	0.45
1:C:11:VAL:HB	1:C:78:VAL:HB	1.98	0.45
1:C:64:VAL:O	1:C:68:THR:CG2	2.64	0.45
1:D:195:ASN:O	1:D:198:GLU:HG2	2.16	0.45
1:A:195:ASN:ND2	1:A:199:ARG:HD3	2.32	0.45
1:C:88:PRO:HD3	1:C:141:ASN:OD1	2.17	0.45
1:D:195:ASN:HA	1:D:198:GLU:CD	2.41	0.45
1:D:204:TRP:HB2	1:D:213:ILE:HD11	1.98	0.45
1:B:40:GLU:O	1:B:41:LYS:C	2.60	0.45
1:A:163:TYR:OH	1:C:90:HIS:CD2	2.53	0.45
1:B:70:LYS:HG2	1:B:71:TYR:CE2	2.52	0.45
1:C:195:ASN:OD1	1:C:195:ASN:C	2.59	0.45
1:C:61:LYS:O	1:C:65:GLU:HG3	2.17	0.45
1:D:204:TRP:CB	1:D:213:ILE:HD11	2.47	0.45
1:D:205:GLY:HA2	1:D:213:ILE:HD12	1.99	0.45
1:A:177:ILE:O	1:A:182:VAL:HG21	2.17	0.44
1:D:16:SER:O	1:D:17:GLY:C	2.60	0.44
1:A:165:PRO:HA	1:D:210:MET:HE2	1.99	0.44
1:C:134:GLN:HA	1:C:137:ALA:O	2.17	0.44
1:A:64:VAL:O	1:A:68:THR:HG23	2.17	0.44
1:C:143:ALA:O	1:C:147:THR:CG2	2.62	0.44
1:C:39:ASP:O	1:C:40:GLU:CB	2.66	0.44
1:C:254:PRO:HG2	1:D:86:TYR:CD2	2.53	0.44
1:A:118:VAL:O	1:A:122:ILE:HG13	2.18	0.44
1:A:139:THR:HG22	1:A:142:ALA:HB3	1.99	0.43
1:D:80:ASN:CG	1:D:80:ASN:O	2.60	0.43
1:A:168:ARG:NH2	1:A:231:SER:OG	2.51	0.43
1:D:7:LYS:HA	1:D:75:ASP:OD2	2.19	0.43
1:B:134:GLN:NE2	1:B:139:THR:H	2.10	0.43
1:B:73:ARG:HB3	1:B:122:ILE:HD11	2.00	0.43
1:B:195:ASN:HD22	1:B:195:ASN:N	2.14	0.43
1:D:206:ARG:CB	1:D:206:ARG:HH11	2.32	0.43
1:B:35:SER:HB3	1:B:46:VAL:HG11	2.01	0.43
1:A:8:VAL:HG21	1:A:71:TYR:HB3	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:68:THR:HG22	1:C:74:ILE:HD11	2.02	0.42
1:B:134:GLN:NE2	1:B:139:THR:HB	2.35	0.42
1:C:39:ASP:HB3	1:C:40:GLU:H	1.53	0.42
1:D:12:THR:HA	1:D:36:VAL:HB	2.01	0.41
1:B:16:SER:HB2	1:B:19:GLY:H	1.85	0.41
1:B:44:VAL:HB	1:B:45:ASN:H	1.47	0.41
1:B:207:GLN:H	1:B:207:GLN:HG2	1.56	0.41
1:B:86:TYR:HA	1:B:141:ASN:O	2.20	0.41
1:A:253:THR:HG21	1:B:139:THR:HG21	2.03	0.41
1:D:217:GLU:H	1:D:217:GLU:CD	2.27	0.41
1:A:134:GLN:NE2	1:A:139:THR:H	2.10	0.41
1:D:153:LEU:O	1:D:157:ARG:HG3	2.20	0.41
1:B:12:THR:O	1:B:80:ASN:HB3	2.21	0.41
1:B:241:VAL:HG12	1:C:233:PHE:HB3	2.03	0.41
1:C:85:GLN:HG2	1:C:96:ILE:HG21	2.02	0.41
1:C:195:ASN:HA	1:C:198:GLU:HG2	2.03	0.41
1:D:89:LEU:C	1:D:89:LEU:HD23	2.45	0.41
1:D:225:PHE:O	1:D:228:SER:HB3	2.21	0.41
1:A:12:THR:O	1:A:80:ASN:HB3	2.20	0.41
1:C:4:PHE:O	1:C:5:THR:C	2.64	0.40
1:A:80:ASN:O	1:A:80:ASN:CG	2.64	0.40
1:B:28:ARG:NH2	1:B:221:GLU:OE2	2.55	0.40
1:B:48:ASP:OD2	1:B:70:LYS:HE3	2.21	0.40
1:B:210:MET:CE	1:C:165:PRO:HA	2.51	0.40
1:B:177:ILE:O	1:B:182:VAL:HG21	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	249/269 (93%)	236 (95%)	13 (5%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	249/269 (93%)	237 (95%)	12 (5%)	0	100	100
1	C	248/269 (92%)	233 (94%)	14 (6%)	1 (0%)	30	38
1	D	248/269 (92%)	238 (96%)	10 (4%)	0	100	100
All	All	994/1076 (92%)	944 (95%)	49 (5%)	1 (0%)	48	60

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	255	ASN

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	208/223 (93%)	184 (88%)	24 (12%)	5	6
1	B	208/223 (93%)	177 (85%)	31 (15%)	3	3
1	C	207/223 (93%)	180 (87%)	27 (13%)	4	4
1	D	207/223 (93%)	179 (86%)	28 (14%)	4	4
All	All	830/892 (93%)	720 (87%)	110 (13%)	4	4

All (110) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	GLU
1	A	3	GLU
1	A	11	VAL
1	A	20	LEU
1	A	38	LEU
1	A	48	ASP
1	A	65	GLU
1	A	73	ARG
1	A	77	LEU
1	A	98	ARG

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Mol	Chain	Res	Type
1	A	110	LEU
1	A	122	ILE
1	A	124	HIS
1	A	134	GLN
1	A	139	THR
1	A	153	LEU
1	A	187	LYS
1	A	195	ASN
1	A	198	GLU
1	A	200	LYS
1	A	203	GLU
1	A	230	ARG
1	A	239	LEU
1	A	249	LEU
1	B	2	GLU
1	B	11	VAL
1	B	18	ILE
1	B	35	SER
1	B	38	LEU
1	B	40	GLU
1	B	44	VAL
1	B	45	ASN
1	B	48	ASP
1	B	58	GLU
1	B	69	LYS
1	B	73	ARG
1	B	77	LEU
1	B	84	GLU
1	B	110	LEU
1	B	122	ILE
1	B	124	HIS
1	B	134	GLN
1	B	139	THR
1	B	147	THR
1	B	152	LEU
1	B	182	VAL
1	B	195	ASN
1	B	198	GLU
1	B	200	LYS
1	B	207	GLN
1	B	228	SER
1	B	239	LEU

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Mol	Chain	Res	Type
1	B	246	LEU
1	B	249	LEU
1	B	253	THR
1	C	3	GLU
1	C	11	VAL
1	C	15	SER
1	C	16	SER
1	C	20	LEU
1	C	28	ARG
1	C	32	LYS
1	C	38	LEU
1	C	68	THR
1	C	77	LEU
1	C	85	GLN
1	C	95	GLU
1	C	110	LEU
1	C	124	HIS
1	C	139	THR
1	C	147	THR
1	C	161	ILE
1	C	180	PRO
1	C	182	VAL
1	C	190	VAL
1	C	192	GLU
1	C	200	LYS
1	C	230	ARG
1	C	239	LEU
1	C	246	LEU
1	C	248	LYS
1	C	253	THR
1	D	3	GLU
1	D	11	VAL
1	D	20	LEU
1	D	23	VAL
1	D	24	ASP
1	D	38	LEU
1	D	44	VAL
1	D	47	SER
1	D	56	ASN
1	D	77	LEU
1	D	110	LEU
1	D	139	THR

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Mol	Chain	Res	Type
1	D	153	LEU
1	D	161	ILE
1	D	181	MET
1	D	190	VAL
1	D	195	ASN
1	D	199	ARG
1	D	200	LYS
1	D	207	GLN
1	D	209	PRO
1	D	213	ILE
1	D	217	GLU
1	D	239	LEU
1	D	246	LEU
1	D	251	ILE
1	D	253	THR
1	D	255	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (26) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	49	HIS
1	A	85	GLN
1	A	90	HIS
1	A	134	GLN
1	A	141	ASN
1	A	195	ASN
1	B	45	ASN
1	B	49	HIS
1	B	90	HIS
1	B	134	GLN
1	B	195	ASN
1	C	45	ASN
1	C	49	HIS
1	C	80	ASN
1	C	85	GLN
1	C	90	HIS
1	C	134	GLN
1	C	207	GLN
1	C	255	ASN
1	D	49	HIS
1	D	56	ASN
1	D	80	ASN

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Mol	Chain	Res	Type
1	D	90	HIS
1	D	134	GLN
1	D	207	GLN
1	D	255	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 oligosaccharides 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, 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	253/269 (94%)	-0.42	5 (1%) 65 66	21, 32, 54, 85	0
1	B	253/269 (94%)	-0.37	4 (1%) 70 72	21, 32, 52, 82	0
1	C	252/269 (93%)	-0.25	6 (2%) 59 62	22, 35, 66, 87	0
1	D	252/269 (93%)	-0.24	5 (1%) 65 66	23, 35, 70, 91	0
All	All	1010/1076 (93%)	-0.32	20 (1%) 65 66	21, 33, 61, 91	0

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	44	VAL	6.9
1	D	256	ALA	5.6
1	D	44	VAL	5.4
1	C	256	ALA	5.0
1	B	44	VAL	4.6
1	C	44	VAL	4.1
1	B	256	ALA	3.8
1	C	124	HIS	3.4
1	B	124	HIS	3.4
1	D	45	ASN	2.7
1	B	181	MET	2.7
1	D	40	GLU	2.5
1	C	197	VAL	2.4
1	C	40	GLU	2.3
1	A	124	HIS	2.3
1	A	40	GLU	2.3
1	A	256	ALA	2.2
1	D	2	GLU	2.2
1	A	181	MET	2.0
1	C	45	ASN	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.