



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 02:01 PM GMT

PDB ID : 3VTZ
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
<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

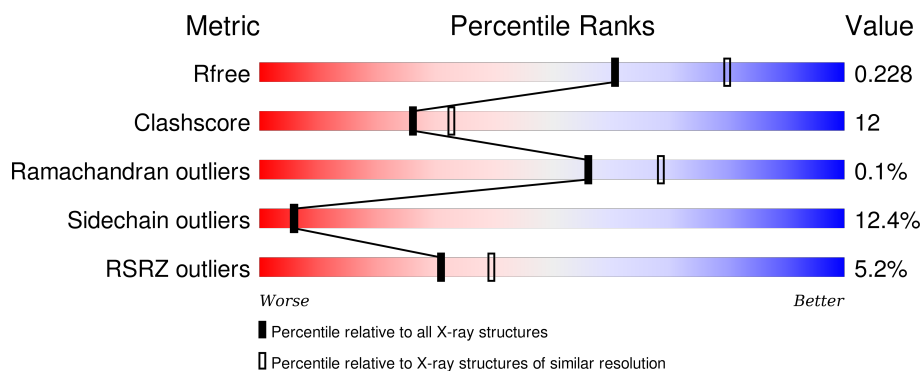
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

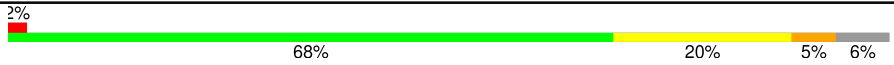



The reported resolution of this entry is 2.30 Å.

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}	91344	3852 (2.30-2.30)
Clashscore	102246	4452 (2.30-2.30)
Ramachandran outliers	100387	4410 (2.30-2.30)
Sidechain outliers	100360	4409 (2.30-2.30)
RSRZ outliers	91569	3857 (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 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	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

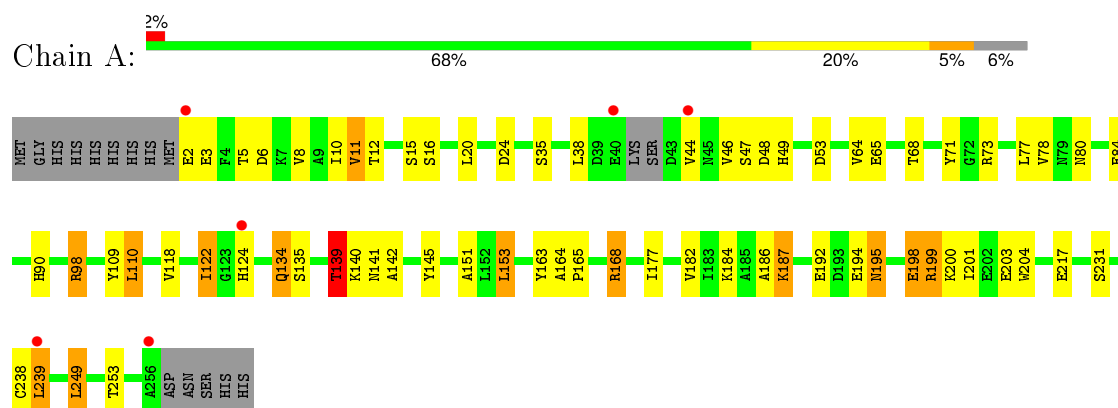
- Molecule 2 is water.

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

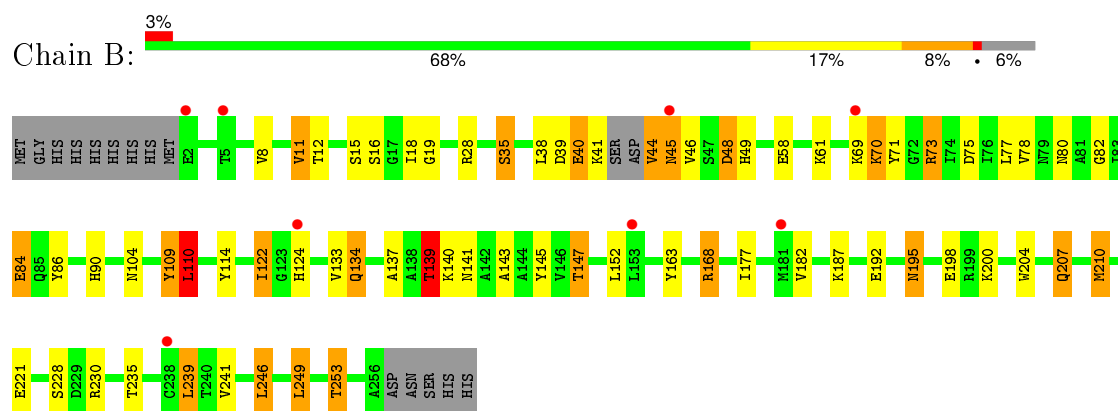
3 Residue-property plots [i](#)

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.

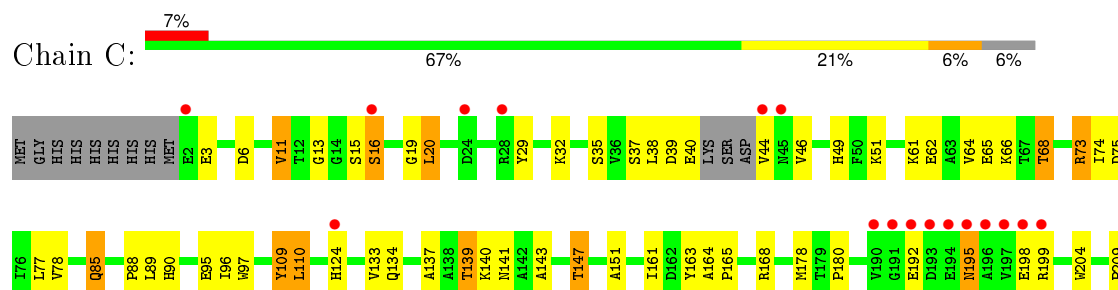
• Molecule 1: Glucose 1-dehydrogenase

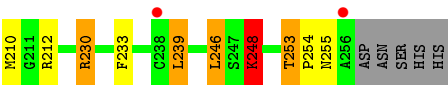


• Molecule 1: Glucose 1-dehydrogenase

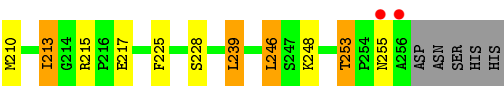
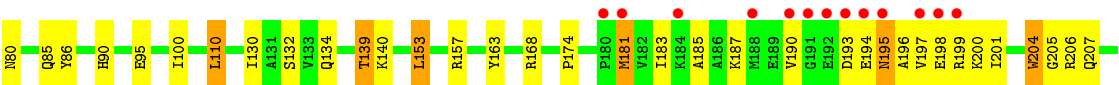


• Molecule 1: Glucose 1-dehydrogenase





● 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) 99.9 (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 (5.34%)	DCC
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 , 41.7	EDS
Estimated twinning fraction	0.017 for -h,-k,l 0.024 for h,-h-k,-l 0.023 for -k,-h,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 46331 reflections	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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.91	2/1941 (0.1%)	1.06	11/2634 (0.4%)
1	B	0.94	1/1942 (0.1%)	1.01	7/2634 (0.3%)
1	C	0.88	1/1933 (0.1%)	0.97	5/2623 (0.2%)
1	D	0.89	1/1933 (0.1%)	1.00	8/2623 (0.3%)
All	All	0.90	5/7749 (0.1%)	1.01	31/10514 (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	A	0	1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	204	TRP	CD2-CE2	6.33	1.49	1.41
1	B	204	TRP	CD2-CE2	6.00	1.48	1.41
1	A	217	GLU	CD-OE1	5.27	1.31	1.25
1	C	204	TRP	CD2-CE2	5.21	1.47	1.41
1	A	204	TRP	CD2-CE2	5.05	1.47	1.41

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	168	ARG	NE-CZ-NH2	-14.41	113.10	120.30
1	A	168	ARG	NE-CZ-NH1	11.49	126.05	120.30
1	D	168	ARG	NE-CZ-NH2	-11.06	114.77	120.30
1	C	168	ARG	NE-CZ-NH2	-10.88	114.86	120.30
1	C	212	ARG	NE-CZ-NH2	6.78	123.69	120.30
1	A	168	ARG	CD-NE-CZ	6.58	132.82	123.60
1	B	168	ARG	NE-CZ-NH1	-6.08	117.26	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	139	THR	N-CA-CB	-6.05	98.81	110.30
1	A	194	GLU	OE1-CD-OE2	-5.93	116.18	123.30
1	A	98	ARG	NE-CZ-NH2	5.89	123.24	120.30
1	D	73	ARG	NE-CZ-NH1	5.88	123.24	120.30
1	D	168	ARG	CG-CD-NE	-5.83	99.57	111.80
1	B	168	ARG	NE-CZ-NH2	5.80	123.20	120.30
1	B	110	LEU	CB-CG-CD1	5.71	120.70	111.00
1	D	168	ARG	NE-CZ-NH1	5.66	123.13	120.30
1	B	249	LEU	CB-CG-CD1	5.58	120.49	111.00
1	A	53	ASP	CB-CG-OD1	5.53	123.27	118.30
1	C	248	LYS	CD-CE-NZ	-5.50	99.05	111.70
1	B	139	THR	N-CA-CB	-5.33	100.17	110.30
1	D	110	LEU	CA-CB-CG	5.28	127.44	115.30
1	C	168	ARG	CG-CD-NE	-5.25	100.78	111.80
1	D	77	LEU	CB-CG-CD2	5.16	119.78	111.00
1	C	6	ASP	CB-CG-OD1	5.13	122.92	118.30
1	B	210	MET	CG-SD-CE	5.11	108.38	100.20
1	D	246	LEU	CB-CG-CD2	5.07	119.62	111.00
1	A	24	ASP	CB-CG-OD1	5.05	122.85	118.30
1	D	248	LYS	CD-CE-NZ	-5.04	100.12	111.70
1	A	110	LEU	CB-CG-CD1	5.03	119.55	111.00
1	A	199	ARG	NE-CZ-NH2	-5.03	117.78	120.30
1	B	249	LEU	CA-CB-CG	5.02	126.86	115.30
1	A	10	ILE	CG1-CB-CG2	-5.02	100.36	111.40

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	47	0
1	B	1913	0	1953	65	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	1904	0	1940	63	0
1	D	1904	0	1940	56	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	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 12.

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:230:ARG:HH11	1:C:230:ARG:HG2	1.15	1.04
1:A:165:PRO:HA	1:D:210:MET:HE1	1.43	0.98
1:C:230:ARG:CG	1:C:230:ARG:HH11	1.78	0.96
1:C:230:ARG:HG2	1:C:230:ARG:NH1	1.87	0.86
1:B:143:ALA:O	1:B:147:THR:HG23	1.76	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:C:143:ALA:O	1:C:147:THR:HG23	1.82	0.80
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:B:90:HIS:HD2	1:D:163:TYR:OH	1.66	0.79
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:B:235:THR:CG2	1:C:210:MET:HE2	2.18	0.74
1:D:134:GLN:NE2	1:D:139:THR:HB	2.02	0.73
1:B:58:GLU:H	1:B:58:GLU:CD	1.93	0.73
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: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:B:139:THR:HG23	1:B:140:LYS:O	1.94	0.68
1:A:139:THR:HG23	1:A: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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:B:235:THR:HG21	1:C:210:MET:CE	2.29	0.62
1:D:183:ILE:HG22	1:D:187:LYS:HE2	1.81	0.62
1:C:133:VAL:HG12	1:C:246:LEU:HD22	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:164:ALA:HB1	1:D:210:MET:CE	2.32	0.59
1:A:182:VAL:HG12	1:A:201:ILE:HD12	1.85	0.59
1:B:235:THR:HG22	1:C:210:MET:HE2	1.85	0.59
1:D:39:ASP:O	1:D:49:HIS:HE1	1.84	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:B:235:THR:HG21	1:C:210:MET:HE2	1.85	0.58
1:B:163:TYR:OH	1:D:90:HIS:CD2	2.53	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:D:181:MET:HE1	1:D:185:ALA:HB2	1.86	0.57
1:A:90:HIS:HD2	1:C:163:TYR:OH	1.88	0.57
1:C:35:SER:HB3	1:C:49:HIS:CD2	2.34	0.56
1:C:16:SER:HB3	1:C:19:GLY:H	1.70	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:168:ARG:HG2	1:C:210:MET:CE	2.37	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:45:ASN:H	1:B:45:ASN:ND2	2.05	0.54
1:B:168:ARG:HG2	1:C:210:MET:HE3	1.90	0.53
1:A:164:ALA:HB1	1:D:210:MET:HE1	1.89	0.53
1:D:85:GLN:HB3	1:D:100:ILE:CD1	2.39	0.53
1:C:209:PRO:HG3	1:C:248:LYS:HD3	1.91	0.53
1:D:205:GLY:CA	1:D:213:ILE:HD12	2.39	0.53
1:A:11:VAL:HB	1:A:78:VAL:HB	1.91	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: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:D:215:ARG:HB3	1:D:217:GLU:OE2	2.09	0.52
1:A:195:ASN:HB2	1:A:198:GLU:HG2	1.92	0.52
1:C:195:ASN:O	1:C:199:ARG:HD3	2.09	0.52
1:B:39:ASP:HB3	1:B:41:LYS:H	1.75	0.51
1:A:249:LEU:HD13	1:B:137:ALA:HB3	1.92	0.51
1:A:47:SER:HB2	2:A:319:HOH:O	2.10	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:235:THR:CG2	1:C:210:MET:CE	2.84	0.50
1:B:210:MET:HE2	1:C:164:ALA:HB1	1.94	0.50
1:A:140:LYS:O	1:B:253:THR:HG21	2.11	0.50
1:B:8:VAL:HG21	1:B:71:TYR:HB3	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:253:THR:HG21	1:B:139:THR:HG23	1.93	0.49
1:A:141:ASN:HD22	1:B:253:THR:HG22	1.77	0.49
1:D:80:ASN:HD22	1:D:130:ILE:HG13	1.76	0.49
1:C:35:SER:CB	1:C:49:HIS:HD2	2.21	0.49
1:C:13:GLY:O	1:C:16:SER:HB2	2.13	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:A:164:ALA:CB	1:D:210:MET:CE	2.91	0.49
1:B:35:SER:OG	1:B:49:HIS:CD2	2.64	0.49
1:A:68:THR:HG21	1:A:118:VAL:HG21	1.93	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:C:3:GLU:O	1:C:3:GLU:HG3	2.12	0.48
1:C:37:SER:O	1:C:51:LYS:HA	2.13	0.48
1:B:82:GLY:HA2	1:B:104:ASN:OD1	2.14	0.48
1:D:95:GLU:OE2	1:D:95:GLU:O	2.31	0.48
1:B:235:THR:HG21	1:C:210:MET:HE1	1.96	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:D:132:SER:HA	1:D:174:PRO:HD2	1.96	0.47
1:B:195:ASN:HD22	1:B:195:ASN:H	1.62	0.47
1:D:196:ALA:HB2	1:D:199:ARG:HH22	1.80	0.47
1:C:195:ASN:HB2	1:C:198:GLU:OE2	2.15	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:A:164:ALA:HB1	1:D:210:MET:HE2	1.98	0.45
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:D:195:ASN:O	1:D:198:GLU:HG2	2.16	0.45
1:C:64:VAL:O	1:C:68:THR:CG2	2.64	0.45
1:C:11:VAL:HB	1:C:78:VAL:HB	1.98	0.45
1:D:204:TRP:HB2	1:D:213:ILE:HD11	1.98	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: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:61:LYS:O	1:C:65:GLU:HG3	2.17	0.45
1:B:187:LYS:NZ	1:B:192:GLU:OE2	2.40	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: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:89:LEU:C	1:C:89:LEU:HD23	2.38	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:B:210:MET:HE2	1:C:164:ALA:CB	2.48	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:118:VAL:O	1:A:122:ILE:HG13	2.18	0.44
1:A:164:ALA:CB	1:D:210:MET:HE1	2.48	0.43
1:A:139:THR:HG22	1:A:142:ALA:HB3	1.99	0.43
1:B:210:MET:HE1	1:C:165:PRO:HA	2.00	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:D:206:ARG:HH11	1:D:206:ARG:CB	2.32	0.43
1:B:195:ASN:HD22	1:B:195:ASN:N	2.14	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
1:D:56:ASN:HD22	1:D:56:ASN:C	2.22	0.42
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:A:165:PRO:HA	1:D:210:MET:CE	2.32	0.42
1:D:12:THR:HA	1:D:36:VAL:HB	2.01	0.41
1:B:44:VAL:HB	1:B:45:ASN:H	1.47	0.41
1:B:16:SER:HB2	1:B:19:GLY:H	1.85	0.41
1:B:207:GLN:HG2	1:B:207:GLN:H	1.56	0.41
1:D:193:ASP:O	1:D:194:GLU:C	2.58	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:B:210:MET:CE	1:C:165:PRO:HA	2.51	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:C:195:ASN:HA	1:C:198:GLU:HG2	2.03	0.41
1:C:85:GLN:HG2	1:C:96:ILE:HG21	2.02	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: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:B:58:GLU:CD	1:B:58:GLU:N	2.68	0.40
1:B:40:GLU:O	1:B:41:LYS:C	2.60	0.40
1:D:195:ASN:HA	1:D:198:GLU:CD	2.41	0.40
1:B:48:ASP:OD2	1:B:70:LYS:HE3	2.21	0.40
1:B:28:ARG:NH2	1:B:221:GLU:OE2	2.55	0.40
1:C:195:ASN:C	1:C:195:ASN:OD1	2.59	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

5.3.1 Protein backbone

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
1	B	249/269 (93%)	237 (95%)	12 (5%)	0	100	100
1	C	248/269 (92%)	233 (94%)	14 (6%)	1 (0%)	39	48
1	D	248/269 (92%)	238 (96%)	10 (4%)	0	100	100
All	All	994/1076 (92%)	944 (95%)	49 (5%)	1 (0%)	56	68

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%)	7	7
1	B	208/223 (93%)	178 (86%)	30 (14%)	4	4
1	C	207/223 (93%)	181 (87%)	26 (13%)	5	5
1	D	207/223 (93%)	184 (89%)	23 (11%)	8	8
All	All	830/892 (93%)	727 (88%)	103 (12%)	6	6

All (103) 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	38	LEU
1	A	48	ASP
1	A	65	GLU
1	A	73	ARG
1	A	77	LEU
1	A	98	ARG
1	A	109	TYR
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	184	LYS
1	A	187	LYS
1	A	195	ASN
1	A	198	GLU
1	A	200	LYS
1	A	203	GLU
1	A	239	LEU
1	A	249	LEU
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	69	LYS
1	B	70	LYS
1	B	73	ARG
1	B	77	LEU
1	B	84	GLU
1	B	109	TYR
1	B	110	LEU
1	B	122	ILE
1	B	124	HIS
1	B	134	GLN
1	B	139	THR

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Mol	Chain	Res	Type
1	B	147	THR
1	B	152	LEU
1	B	195	ASN
1	B	198	GLU
1	B	200	LYS
1	B	207	GLN
1	B	228	SER
1	B	239	LEU
1	B	246	LEU
1	B	249	LEU
1	B	253	THR
1	C	11	VAL
1	C	15	SER
1	C	16	SER
1	C	20	LEU
1	C	32	LYS
1	C	38	LEU
1	C	68	THR
1	C	73	ARG
1	C	77	LEU
1	C	85	GLN
1	C	95	GLU
1	C	109	TYR
1	C	110	LEU
1	C	124	HIS
1	C	139	THR
1	C	147	THR
1	C	161	ILE
1	C	178	MET
1	C	180	PRO
1	C	192	GLU
1	C	195	ASN
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	24	ASP
1	D	38	LEU

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Mol	Chain	Res	Type
1	D	44	VAL
1	D	47	SER
1	D	56	ASN
1	D	70	LYS
1	D	77	LEU
1	D	110	LEU
1	D	139	THR
1	D	153	LEU
1	D	181	MET
1	D	190	VAL
1	D	195	ASN
1	D	200	LYS
1	D	207	GLN
1	D	213	ILE
1	D	239	LEU
1	D	246	LEU
1	D	253	THR
1	D	255	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) such sidechains are listed below:

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

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Mol	Chain	Res	Type
1	D	134	GLN
1	D	207	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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%)	-0.00	6 (2%)	62 71	21, 32, 54, 85	0
1	B	253/269 (94%)	0.08	8 (3%)	51 60	21, 32, 52, 82	0
1	C	252/269 (93%)	0.21	19 (7%)	17 24	22, 35, 66, 87	0
1	D	252/269 (93%)	0.28	20 (7%)	15 22	23, 35, 70, 91	0
All	All	1010/1076 (93%)	0.14	53 (5%)	31 39	21, 33, 61, 91	0

All (53) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	192	GLU	5.4
1	C	197	VAL	5.3
1	A	44	VAL	4.9
1	C	256	ALA	4.8
1	D	198	GLU	4.7
1	D	194	GLU	4.6
1	D	193	ASP	4.6
1	C	194	GLU	4.6
1	D	195	ASN	4.4
1	D	2	GLU	4.2
1	D	256	ALA	3.8
1	D	44	VAL	3.8
1	B	2	GLU	3.6
1	B	5	THR	3.5
1	C	45	ASN	3.4
1	D	197	VAL	3.4
1	D	184	LYS	3.3
1	C	190	VAL	3.3
1	D	199	ARG	3.2
1	C	199	ARG	3.2
1	D	190	VAL	3.1

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Mol	Chain	Res	Type	RSRZ
1	A	40	GLU	3.0
1	D	255	ASN	2.9
1	C	28	ARG	2.9
1	D	191	GLY	2.9
1	C	196	ALA	2.8
1	C	198	GLU	2.8
1	B	238	CYS	2.8
1	B	45	ASN	2.7
1	C	191	GLY	2.7
1	C	193	ASP	2.6
1	C	195	ASN	2.6
1	C	44	VAL	2.6
1	A	256	ALA	2.5
1	D	45	ASN	2.4
1	A	2	GLU	2.4
1	D	180	PRO	2.4
1	B	124	HIS	2.3
1	B	153	LEU	2.3
1	D	188	MET	2.3
1	B	181	MET	2.3
1	C	238	CYS	2.3
1	C	24	ASP	2.3
1	C	2	GLU	2.2
1	D	181	MET	2.2
1	C	124	HIS	2.2
1	B	69	LYS	2.1
1	A	239	LEU	2.1
1	D	24	ASP	2.1
1	C	192	GLU	2.0
1	A	124	HIS	2.0
1	D	28	ARG	2.0
1	C	16	SER	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.