



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 02:48 AM GMT

PDB ID : 2ISM
Title : Crystal structure of the putative oxidoreductase (glucose dehydrogenase) (TTHA0570) from thermus thermophilus HB8
Authors : Jeyakanthan, J.; Kanauija, S.P.; Vasuki Ranjani, C.; Sekar, K.; Ebihara, A.; Shinkai, A.; Nakagawa, N.; Shimizu, N.; Yamamoto, M.; Kuramitsu, S.; Shiro, Y.; Yokoyama, S.; RIKEN Structural Genomics/Proteomics Initiative (RSGI)
Deposited on : 2006-10-18
Resolution : 1.90 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

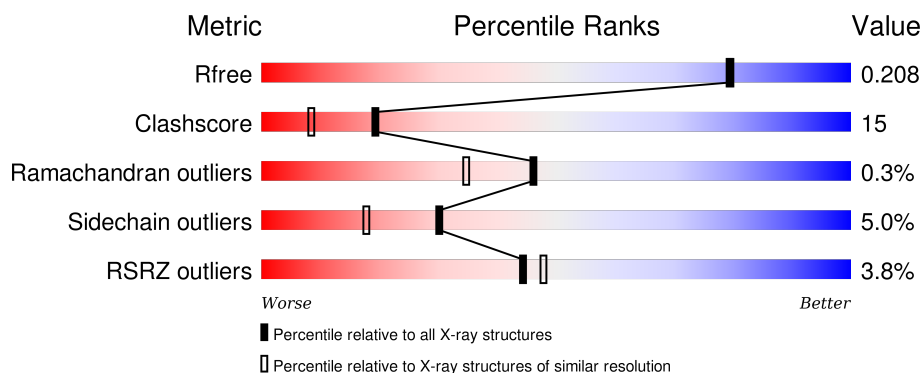
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	91344	4755 (1.90-1.90)
Clashscore	102246	5398 (1.90-1.90)
Ramachandran outliers	100387	5338 (1.90-1.90)
Sidechain outliers	100360	5339 (1.90-1.90)
RSRZ outliers	91569	4766 (1.90-1.90)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	352	<div> <div>4%</div> <div>69%</div> <div>19%</div> <div>• 7%</div> </div>
1	B	352	<div> <div>3%</div> <div>73%</div> <div>18%</div> <div>• • 5%</div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	CA	A	1002	-	-	-	X

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 5815 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 Putative oxidoreductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	327	Total	C	N	O	S	0	0	0
			2558	1621	488	447	2			
1	B	333	Total	C	N	O	S	0	0	0
			2602	1647	495	458	2			

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

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

- Molecule 3 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

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

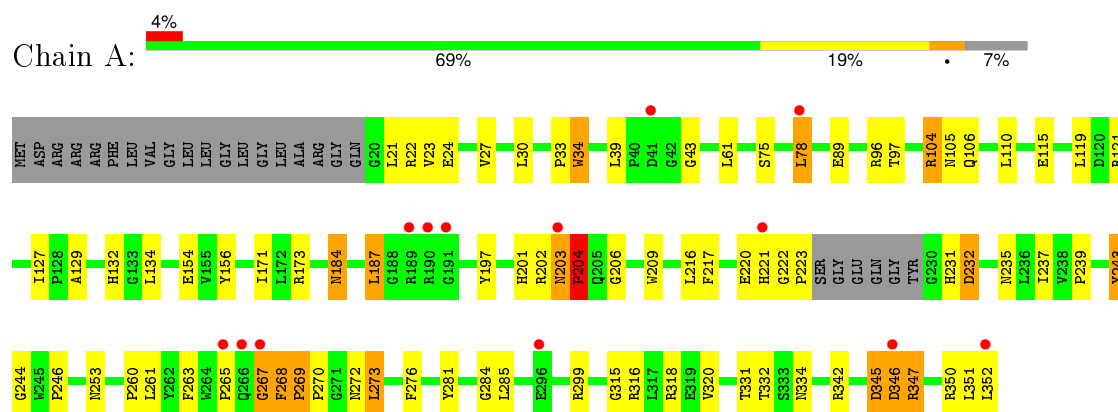
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	350	Total	O	0	0
			350	350		
4	B	301	Total	O	0	0
			301	301		

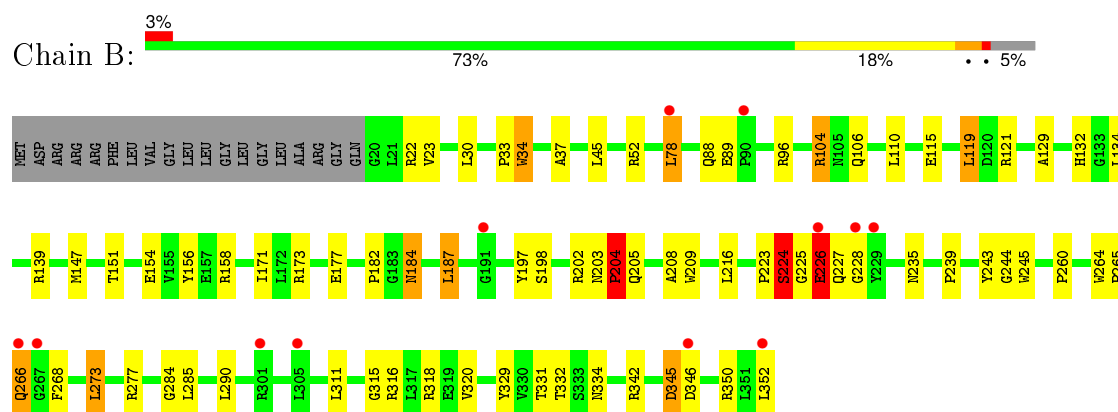
3 Residue-property plots

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: Putative oxidoreductase



• Molecule 1: Putative oxidoreductase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	36.90 Å 132.96 Å 60.78 Å 90.00° 97.18° 90.00°	Depositor
Resolution (Å)	36.61 – 1.90 36.61 – 1.90	Depositor EDS
% Data completeness (in resolution range)	97.7 (36.61-1.90) 97.9 (36.61-1.90)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.30 (at 1.89 Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.181 , 0.208 0.181 , 0.208	Depositor DCC
R_{free} test set	2236 reflections (5.01%)	DCC
Wilson B-factor (Å ²)	17.9	Xtriage
Anisotropy	0.131	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 56.9	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	1 of 45594 reflections (0.002%)	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	5815	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.75% 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

Bond lengths and bond angles in the following residue types are not validated in this section: CA, 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	0.43	1/2625 (0.0%)	0.85	15/3559 (0.4%)
1	B	0.35	1/2671 (0.0%)	0.79	8/3622 (0.2%)
All	All	0.39	2/5296 (0.0%)	0.82	23/7181 (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
1	B	0	1
All	All	0	2

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	204	PRO	N-CD	15.42	1.69	1.47
1	B	204	PRO	N-CD	5.79	1.55	1.47

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	267	GLY	N-CA-C	16.74	154.96	113.10
1	B	78	LEU	N-CA-C	9.82	137.51	111.00
1	A	268	PHE	N-CA-CB	-9.80	92.95	110.60
1	A	78	LEU	N-CA-C	9.10	135.56	111.00
1	A	203	ASN	C-N-CD	-8.44	102.03	120.60
1	A	204	PRO	N-CA-CB	7.65	112.48	103.30
1	B	226	GLU	N-CA-C	-7.61	90.46	111.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	204	PRO	CA-N-CD	-7.53	100.95	111.50
1	A	89	GLU	N-CA-C	-6.01	94.77	111.00
1	B	224	SER	N-CA-CB	-6.00	101.50	110.50
1	A	318	ARG	N-CA-C	5.88	126.86	111.00
1	B	203	ASN	N-CA-C	-5.85	95.21	111.00
1	A	244	GLY	N-CA-C	5.81	127.62	113.10
1	B	226	GLU	N-CA-CB	-5.73	100.28	110.60
1	B	244	GLY	N-CA-C	5.54	126.94	113.10
1	B	318	ARG	N-CA-C	5.49	125.83	111.00
1	A	232	ASP	CB-CG-OD2	5.26	123.03	118.30
1	A	346	ASP	CB-CG-OD2	5.22	123.00	118.30
1	A	345	ASP	CB-CG-OD2	5.21	122.99	118.30
1	B	89	GLU	N-CA-C	-5.20	96.97	111.00
1	A	201	HIS	CB-CA-C	5.08	120.56	110.40
1	A	78	LEU	CB-CA-C	-5.03	100.64	110.20
1	A	171	ILE	N-CA-C	-5.01	97.46	111.00

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	243	TYR	Sidechain
1	B	243	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	2558	0	2538	68	0
1	B	2602	0	2573	87	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	350	0	0	3	0
4	B	301	0	0	7	0
All	All	5815	0	5111	155	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 15.

All (155) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:204:PRO:N	1:A:204:PRO:CD	1.69	1.38
1:B:266:GLN:HB3	4:B:1053:HOH:O	1.01	1.19
1:A:203:ASN:O	1:A:220:GLU:HB2	1.57	1.03
1:B:223:PRO:HB3	1:B:226:GLU:O	1.57	1.03
1:A:267:GLY:O	1:A:269:PRO:HD3	1.59	1.02
1:B:332:THR:OG1	1:B:346:ASP:HA	1.61	1.01
1:A:332:THR:OG1	1:A:346:ASP:HA	1.63	0.97
1:A:34:TRP:CE3	1:A:78:LEU:O	2.19	0.95
1:A:203:ASN:C	1:A:204:PRO:CD	2.36	0.94
1:A:342:ARG:O	1:A:345:ASP:HB2	1.68	0.93
1:B:132:HIS:HD2	1:B:134:LEU:H	1.16	0.91
1:A:132:HIS:HD2	1:A:134:LEU:H	1.13	0.90
1:B:202:ARG:HH21	1:B:226:GLU:HG2	1.34	0.89
1:B:277:ARG:HH12	1:B:352:LEU:HD12	1.42	0.85
1:A:202:ARG:HD3	1:A:222:GLY:HA2	1.59	0.83
1:B:96:ARG:HE	1:B:106:GLN:HE21	1.27	0.82
1:A:34:TRP:HE3	1:A:78:LEU:O	1.67	0.78
1:A:22:ARG:H	1:A:352:LEU:HG	1.50	0.76
1:A:132:HIS:CD2	1:A:134:LEU:H	2.03	0.75
1:B:173:ARG:HE	1:B:184:ASN:HD21	1.32	0.75
1:B:34:TRP:CE3	1:B:78:LEU:O	2.40	0.75
1:B:265:PRO:HD2	1:B:266:GLN:HE22	1.54	0.72
1:A:285:LEU:O	1:A:316:ARG:HG2	1.89	0.72
1:A:332:THR:HG1	1:A:346:ASP:HA	1.55	0.72
1:B:158:ARG:HH22	1:B:227:GLN:HB2	1.55	0.71
1:B:223:PRO:HB2	1:B:228:GLY:O	1.90	0.71
1:B:225:GLY:HA2	4:B:1299:HOH:O	1.90	0.71
1:B:226:GLU:HG3	1:B:226:GLU:O	1.90	0.71
1:B:285:LEU:O	1:B:316:ARG:HG2	1.90	0.70
1:A:96:ARG:HE	1:A:106:GLN:HE21	1.39	0.69
1:B:147:MET:HE3	1:B:182:PRO:HD2	1.75	0.69
1:A:115:GLU:HG3	4:A:1032:HOH:O	1.93	0.68
1:B:147:MET:CE	1:B:182:PRO:HD2	2.23	0.68
1:B:115:GLU:HG2	4:B:1203:HOH:O	1.93	0.67
1:B:277:ARG:NH1	1:B:352:LEU:HD12	2.09	0.67
1:B:139:ARG:HB2	1:B:205:GLN:O	1.95	0.67
1:B:22:ARG:H	1:B:352:LEU:CD2	2.08	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:227:GLN:O	1:B:227:GLN:HG2	1.94	0.66
1:B:264:TRP:HB3	1:B:266:GLN:HE21	1.62	0.65
1:A:232:ASP:CB	1:A:268:PHE:O	2.47	0.63
1:A:231:HIS:HB3	1:A:263:PHE:CZ	2.34	0.62
1:B:173:ARG:HE	1:B:184:ASN:ND2	1.97	0.62
1:B:266:GLN:NE2	1:B:266:GLN:O	2.30	0.61
1:B:342:ARG:O	1:B:345:ASP:HB2	2.01	0.61
1:A:221:HIS:CD2	1:A:232:ASP:OD1	2.54	0.60
1:A:203:ASN:CA	1:A:204:PRO:CD	2.80	0.60
1:B:96:ARG:HE	1:B:106:GLN:NE2	1.96	0.59
1:A:96:ARG:HE	1:A:106:GLN:NE2	2.00	0.59
1:A:173:ARG:HE	1:A:184:ASN:HD21	1.50	0.59
1:B:224:SER:O	1:B:225:GLY:C	2.39	0.59
1:B:332:THR:HG1	1:B:346:ASP:HA	1.65	0.58
1:B:316:ARG:H	1:B:334:ASN:HD21	1.50	0.58
1:B:264:TRP:HB3	1:B:266:GLN:O	2.04	0.58
1:B:224:SER:C	1:B:226:GLU:N	2.55	0.57
1:B:268:PHE:CZ	1:B:284:GLY:HA3	2.39	0.57
1:A:110:LEU:CD2	1:A:119:LEU:HD23	2.35	0.57
1:B:110:LEU:HD22	1:B:119:LEU:HA	1.86	0.56
1:B:224:SER:O	1:B:226:GLU:HB3	2.05	0.56
1:A:197:TYR:OH	1:A:239:PRO:HG3	2.05	0.56
1:A:173:ARG:HE	1:A:184:ASN:ND2	2.05	0.55
1:B:350:ARG:HD2	4:B:1116:HOH:O	2.06	0.55
1:B:158:ARG:NH2	1:B:227:GLN:HE21	2.03	0.55
1:B:34:TRP:HE3	1:B:78:LEU:O	1.87	0.55
1:B:316:ARG:H	1:B:334:ASN:ND2	2.04	0.55
1:B:30:LEU:O	1:B:346:ASP:HB2	2.07	0.54
1:B:147:MET:HE2	4:B:1082:HOH:O	2.05	0.54
1:A:232:ASP:HB3	1:A:270:PRO:HD3	1.88	0.54
1:A:203:ASN:O	1:A:220:GLU:CB	2.44	0.53
1:A:235:ASN:HD22	1:A:260:PRO:HA	1.73	0.53
1:B:110:LEU:CD2	1:B:119:LEU:HA	2.37	0.53
1:A:263:PHE:O	1:A:265:PRO:HD3	2.09	0.53
1:A:273:LEU:H	1:A:273:LEU:HD12	1.73	0.53
1:B:37:ALA:HB3	1:B:45:LEU:HB2	1.90	0.53
1:A:316:ARG:H	1:A:334:ASN:HD21	1.57	0.52
1:B:158:ARG:HH21	1:B:227:GLN:NE2	2.08	0.52
1:A:217:PHE:HE1	1:A:261:LEU:HD22	1.74	0.52
1:B:202:ARG:NE	1:B:226:GLU:OE2	2.43	0.52
1:B:202:ARG:NH2	1:B:226:GLU:HG2	2.15	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:268:PHE:O	1:A:270:PRO:HD3	2.10	0.52
1:B:277:ARG:HH22	1:B:352:LEU:HD12	1.75	0.51
1:B:158:ARG:HH21	1:B:227:GLN:HE21	1.58	0.51
1:B:52:ARG:HD3	4:B:1156:HOH:O	2.10	0.51
1:B:22:ARG:H	1:B:352:LEU:HD21	1.75	0.51
1:B:197:TYR:OH	1:B:239:PRO:HG3	2.11	0.51
1:B:264:TRP:HB3	1:B:266:GLN:NE2	2.24	0.51
1:B:22:ARG:H	1:B:352:LEU:HD23	1.75	0.50
1:B:224:SER:H	1:B:226:GLU:CG	2.24	0.50
1:A:184:ASN:HB2	1:A:187:LEU:HD13	1.93	0.50
1:A:30:LEU:HB2	1:A:346:ASP:O	2.11	0.50
1:B:235:ASN:HD22	1:B:260:PRO:HA	1.77	0.50
1:A:235:ASN:ND2	1:A:260:PRO:HA	2.26	0.49
1:B:315:GLY:HA3	1:B:334:ASN:HD21	1.77	0.49
1:A:104:ARG:HA	1:A:129:ALA:HB3	1.94	0.49
1:B:173:ARG:HH21	1:B:184:ASN:HD22	1.61	0.49
1:A:21:LEU:HA	1:A:352:LEU:HD12	1.93	0.49
1:A:350:ARG:HD2	4:A:1199:HOH:O	2.13	0.48
1:A:132:HIS:HE1	4:A:1048:HOH:O	1.96	0.48
1:B:132:HIS:CD2	1:B:134:LEU:H	2.09	0.48
1:B:139:ARG:CB	1:B:205:GLN:O	2.62	0.47
1:A:209:TRP:CE2	1:A:216:LEU:HD13	2.49	0.47
1:A:237:ILE:HA	1:A:243:TYR:OH	2.15	0.47
1:A:315:GLY:HA3	1:A:334:ASN:HD21	1.80	0.46
1:A:23:VAL:HG22	1:A:24:GLU:N	2.30	0.46
1:B:151:THR:HB	1:B:204:PRO:HG2	1.98	0.46
1:B:266:GLN:H	1:B:266:GLN:CD	2.18	0.46
1:A:217:PHE:CE1	1:A:261:LEU:HD22	2.49	0.46
1:A:39:LEU:HD12	1:A:43:GLY:HA3	1.97	0.45
1:A:22:ARG:N	1:A:352:LEU:HG	2.27	0.45
1:B:223:PRO:CB	1:B:226:GLU:O	2.46	0.45
1:A:316:ARG:H	1:A:334:ASN:ND2	2.15	0.45
1:A:223:PRO:HB2	1:A:231:HIS:CD2	2.52	0.45
1:B:104:ARG:HA	1:B:129:ALA:HB3	1.98	0.45
1:A:27:VAL:O	1:A:347:ARG:HB3	2.16	0.45
1:B:223:PRO:HG3	1:B:245:TRP:CZ3	2.52	0.45
1:B:273:LEU:HD12	1:B:273:LEU:H	1.82	0.45
1:A:204:PRO:HA	1:A:220:GLU:CB	2.46	0.44
1:B:320:VAL:HA	1:B:329:TYR:O	2.18	0.44
1:A:33:PRO:HG2	1:A:331:THR:HB	1.99	0.44
1:A:222:GLY:O	1:A:223:PRO:C	2.54	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:221:HIS:CD2	1:A:222:GLY:O	2.71	0.44
1:B:121:ARG:NH2	1:B:177:GLU:O	2.51	0.43
1:A:351:LEU:HD12	1:A:351:LEU:N	2.33	0.43
1:B:277:ARG:NH2	1:B:352:LEU:HD12	2.33	0.43
1:B:316:ARG:N	1:B:334:ASN:HD21	2.16	0.43
1:B:202:ARG:HH21	1:B:226:GLU:CG	2.19	0.43
1:A:273:LEU:N	1:A:273:LEU:HD12	2.32	0.43
1:A:221:HIS:HD2	1:A:232:ASP:OD1	1.98	0.43
1:A:75:SER:HB3	1:A:97:THR:OG1	2.19	0.42
1:B:208:ALA:CB	1:B:273:LEU:HD13	2.48	0.42
1:B:208:ALA:HB2	1:B:273:LEU:HD13	2.00	0.42
1:A:173:ARG:HH21	1:A:184:ASN:HD22	1.66	0.42
1:B:23:VAL:HG21	1:B:311:LEU:HD11	2.01	0.42
1:B:265:PRO:HD2	1:B:266:GLN:NE2	2.30	0.42
1:B:158:ARG:NH2	1:B:227:GLN:NE2	2.67	0.42
1:B:30:LEU:HB2	1:B:346:ASP:O	2.19	0.42
1:B:96:ARG:HH21	1:B:106:GLN:HE22	1.68	0.42
1:B:33:PRO:HG2	1:B:331:THR:HB	2.01	0.41
1:B:197:TYR:CZ	1:B:239:PRO:HG3	2.55	0.41
1:B:235:ASN:ND2	1:B:260:PRO:HA	2.35	0.41
1:A:75:SER:HB3	1:A:97:THR:HG1	1.85	0.41
1:A:276:PHE:HB3	1:A:281:TYR:CD2	2.55	0.41
1:A:206:GLY:HA2	1:A:272:ASN:OD1	2.20	0.41
1:B:132:HIS:HE1	4:B:1079:HOH:O	2.03	0.41
1:A:268:PHE:CZ	1:A:284:GLY:HA3	2.56	0.41
1:B:171:ILE:HB	1:B:198:SER:OG	2.21	0.41
1:A:299:ARG:HG2	1:A:299:ARG:HH11	1.86	0.41
1:B:173:ARG:HH21	1:B:184:ASN:ND2	2.19	0.41
1:B:290:LEU:HB3	1:B:311:LEU:HB3	2.02	0.41
1:A:320:VAL:O	1:A:320:VAL:HG13	2.21	0.41
1:B:187:LEU:HA	1:B:187:LEU:HD12	1.86	0.41
1:B:34:TRP:HB3	1:B:78:LEU:O	2.21	0.40
1:A:105:ASN:HB2	1:A:127:ILE:HB	2.04	0.40
1:A:187:LEU:HA	1:A:187:LEU:HD12	1.89	0.40
1:A:232:ASP:CG	1:A:268:PHE:O	2.60	0.40
1:B:209:TRP:CE2	1:B:216:LEU:HD13	2.57	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	323/352 (92%)	307 (95%)	15 (5%)	1 (0%)	46	35
1	B	331/352 (94%)	314 (95%)	16 (5%)	1 (0%)	46	35
All	All	654/704 (93%)	621 (95%)	31 (5%)	2 (0%)	46	35

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	204	PRO
1	B	204	PRO

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	257/275 (94%)	244 (95%)	13 (5%)	29	17
1	B	261/275 (95%)	248 (95%)	13 (5%)	30	18
All	All	518/550 (94%)	492 (95%)	26 (5%)	30	18

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	34	TRP
1	A	61	LEU
1	A	104	ARG
1	A	121	ARG

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Mol	Chain	Res	Type
1	A	154	GLU
1	A	156	TYR
1	A	184	ASN
1	A	187	LEU
1	A	246	PRO
1	A	253	ASN
1	A	269	PRO
1	A	273	LEU
1	A	347	ARG
1	B	34	TRP
1	B	88	GLN
1	B	104	ARG
1	B	119	LEU
1	B	154	GLU
1	B	156	TYR
1	B	184	ASN
1	B	187	LEU
1	B	224	SER
1	B	226	GLU
1	B	266	GLN
1	B	273	LEU
1	B	345	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	105	ASN
1	A	106	GLN
1	A	132	HIS
1	A	184	ASN
1	A	221	HIS
1	A	231	HIS
1	A	235	ASN
1	A	334	ASN
1	B	88	GLN
1	B	105	ASN
1	B	106	GLN
1	B	132	HIS
1	B	184	ASN
1	B	221	HIS
1	B	227	GLN
1	B	235	ASN

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Mol	Chain	Res	Type
1	B	266	GLN
1	B	334	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 4 ligands modelled in this entry, 4 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	327/352 (92%)	-0.05	13 (3%) 42 46	9, 20, 46, 72	0
1	B	333/352 (94%)	-0.03	12 (3%) 46 50	9, 19, 45, 79	0
All	All	660/704 (93%)	-0.04	25 (3%) 44 48	9, 19, 46, 79	0

All (25) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	226	GLU	4.9
1	B	266	GLN	4.6
1	A	267	GLY	4.4
1	A	352	LEU	4.4
1	B	228	GLY	3.8
1	B	305	LEU	3.7
1	A	78	LEU	3.2
1	A	346	ASP	3.2
1	B	229	TYR	3.1
1	B	301	ARG	3.0
1	A	221	HIS	2.8
1	A	265	PRO	2.8
1	B	191	GLY	2.7
1	A	203	ASN	2.6
1	B	78	LEU	2.6
1	A	191	GLY	2.6
1	B	267	GLY	2.5
1	A	296	GLU	2.5
1	B	352	LEU	2.5
1	B	346	ASP	2.5
1	B	90	PRO	2.2
1	A	41	ASP	2.1
1	A	266	GLN	2.1
1	A	190	ARG	2.0

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Mol	Chain	Res	Type	RSRZ
1	A	189	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 carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
2	CA	A	1002	1/1	0.68	0.28	10.48	82,82,82,82	0
2	CA	B	1001	1/1	0.99	0.05	-1.32	21,21,21,21	0
3	CL	A	1003	1/1	0.98	0.06	-1.58	24,24,24,24	0
3	CL	B	1004	1/1	0.98	0.04	-3.08	23,23,23,23	0

6.5 Other polymers [i](#)

There are no such residues in this entry.