



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 01:41 AM GMT

PDB ID : 2E0C
Title : crystal structure of isocitrate dehydrogenase from Sulfolobus tokodaii strain7
at 2.0 Å resolution
Authors : Kouyama, T.
Deposited on : 2006-10-06
Resolution : 2.00 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<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

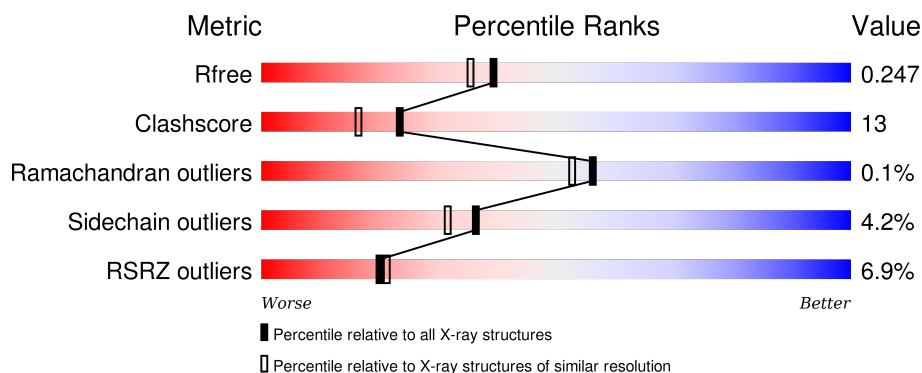
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	91344	6249 (2.00-2.00)
Clashscore	102246	7340 (2.00-2.00)
Ramachandran outliers	100387	7248 (2.00-2.00)
Sidechain outliers	100360	7247 (2.00-2.00)
RSRZ outliers	91569	6262 (2.00-2.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments 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	409	<div> <div>7%</div> <div> <div></div> <div>71%</div> <div>26%</div> <div>..</div> </div> </div>
1	B	409	<div> <div>6%</div> <div> <div></div> <div>69%</div> <div>27%</div> <div>..</div> </div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 6881 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 409aa long hypothetical NADP-dependent isocitrate dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	401	Total	C	N	O	S	0	0	0
			3209	2054	543	597	15			
1	B	401	Total	C	N	O	S	0	0	0
			3209	2054	543	597	15			

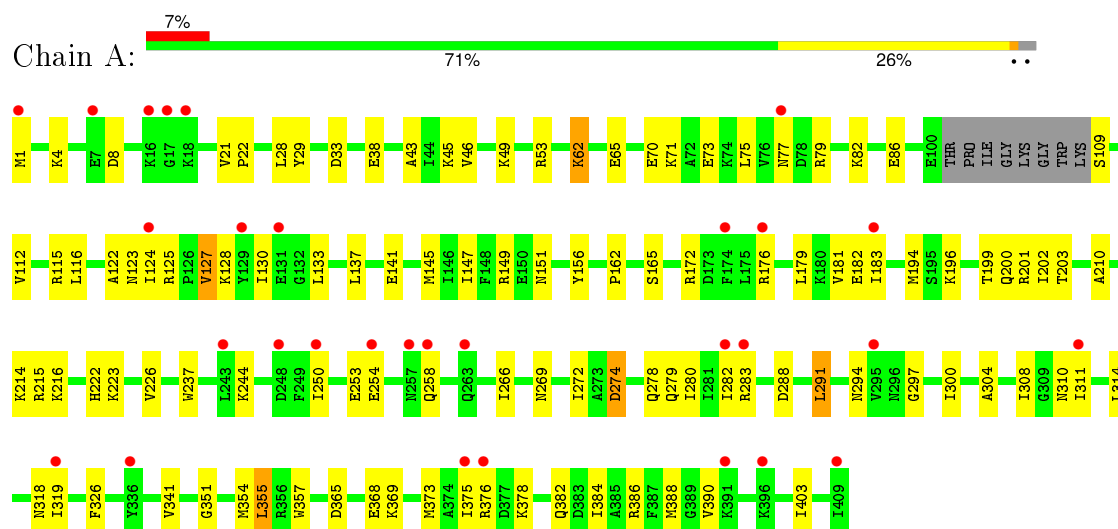
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	249	Total	O	0	0
			249	249		
2	B	214	Total	O	0	0
			214	214		

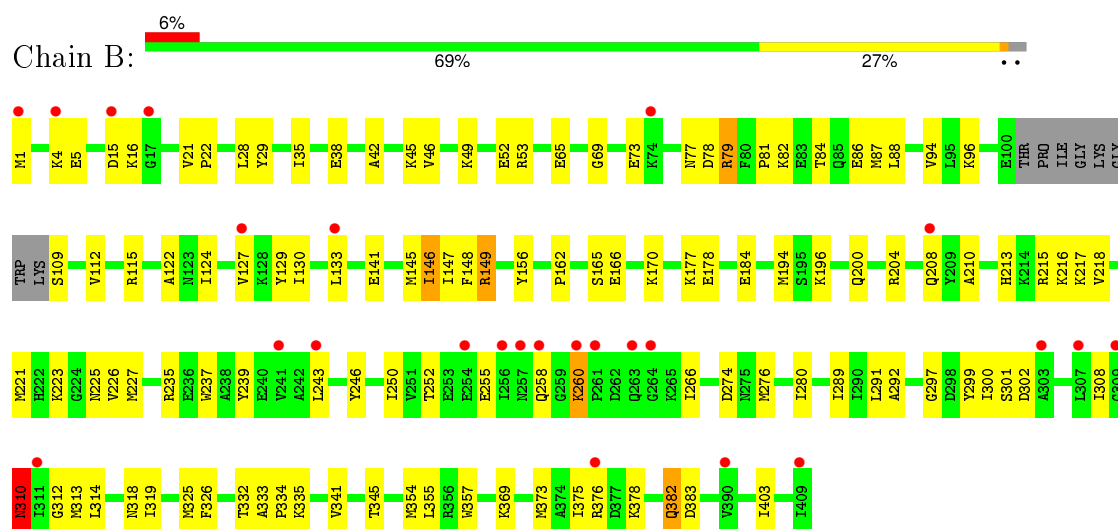
3 Residue-property plots [i](#)

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: 409aa long hypothetical NADP-dependent isocitrate dehydrogenase



- Molecule 1: 409aa long hypothetical NADP-dependent isocitrate dehydrogenase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	74.88Å 87.72Å 75.72Å 90.00° 91.37° 90.00°	Depositor
Resolution (Å)	15.00 – 2.00 45.11 – 2.00	Depositor EDS
% Data completeness (in resolution range)	99.2 (15.00-2.00) 99.2 (45.11-2.00)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.26 (at 2.00Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.221 , 0.247 0.221 , 0.247	Depositor DCC
R_{free} test set	3320 reflections (5.07%)	DCC
Wilson B-factor (Å ²)	30.4	Xtriage
Anisotropy	0.116	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 52.7	EDS
Estimated twinning fraction	0.004 for l,k,-h 0.066 for h,-k,-l 0.018 for l,-k,h	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 65657 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6881	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.85% 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 [i](#)

5.1 Standard geometry [i](#)

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	0.34	0/3264	0.61	0/4388
1	B	0.31	0/3264	0.58	0/4388
All	All	0.32	0/6528	0.60	0/8776

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3209	0	3285	79	0
1	B	3209	0	3285	97	0
2	A	249	0	0	6	0
2	B	214	0	0	8	0
All	All	6881	0	6570	173	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 (173) 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:49:LYS:HG3	1:B:403:ILE:HG23	1.45	0.97

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:149:ARG:HG3	1:A:297:GLY:HA3	1.47	0.95
1:A:162:PRO:HG2	1:A:165:SER:HB3	1.50	0.91
1:A:49:LYS:HG3	1:A:403:ILE:HG23	1.54	0.88
1:B:149:ARG:HG3	1:B:297:GLY:HA3	1.58	0.86
1:B:124:ILE:HG23	1:B:146:ILE:HD11	1.59	0.84
1:B:162:PRO:HG2	1:B:165:SER:HB3	1.61	0.81
1:A:311:ILE:HD12	1:A:314:LEU:HD12	1.62	0.80
1:B:28:LEU:HB2	1:B:94:VAL:HG22	1.63	0.79
1:B:5:GLU:HG2	1:B:87:MET:HE1	1.67	0.77
1:A:223:LYS:HG3	1:A:226:VAL:HG22	1.68	0.76
1:B:124:ILE:CG2	1:B:146:ILE:HD11	2.16	0.75
1:A:250:ILE:HG22	1:A:266:ILE:HB	1.67	0.75
1:A:210:ALA:HA	1:A:215:ARG:HB2	1.71	0.73
1:A:172:ARG:HG2	1:A:183:ILE:HD12	1.73	0.70
1:A:282:ILE:HG13	1:A:283:ARG:N	2.09	0.68
1:B:109:SER:HB3	1:B:112:VAL:HG12	1.76	0.66
1:B:82:LYS:O	1:B:86:GLU:HG3	1.95	0.66
1:B:223:LYS:HG3	1:B:226:VAL:HG22	1.78	0.66
1:B:341:VAL:CG1	1:B:382:GLN:HG2	2.25	0.66
1:B:223:LYS:HE3	1:B:274:ASP:OD2	1.97	0.65
1:B:29:TYR:CZ	1:B:65:GLU:HB2	2.32	0.65
1:A:82:LYS:O	1:A:86:GLU:HG3	1.98	0.64
1:A:279:GLN:HE22	1:A:282:ILE:HD11	1.63	0.64
1:B:382:GLN:HG3	1:B:383:ASP:N	2.13	0.64
1:A:311:ILE:HG22	2:A:650:HOH:O	1.96	0.64
1:A:201:ARG:HD3	2:A:548:HOH:O	1.98	0.63
1:B:88:LEU:CD2	1:B:94:VAL:HG21	2.28	0.63
1:B:28:LEU:CB	1:B:94:VAL:HG22	2.28	0.62
1:A:176:ARG:NH1	1:A:182:GLU:HG2	2.14	0.62
1:B:96:LYS:HD3	1:B:325:MET:SD	2.41	0.61
1:B:130:ILE:HB	1:B:133:LEU:HD12	1.84	0.60
1:B:146:ILE:HD13	1:B:147:ILE:N	2.17	0.59
1:A:130:ILE:HB	1:A:133:LEU:HD12	1.84	0.59
1:A:388:MET:HG3	1:A:390:VAL:HG22	1.86	0.57
1:B:69:GLY:HA2	2:B:623:HOH:O	2.05	0.57
1:B:210:ALA:HA	1:B:215:ARG:HB2	1.86	0.56
1:A:127:VAL:HG13	1:A:145:MET:HG3	1.86	0.56
1:B:221:MET:HB3	1:B:292:ALA:HB2	1.87	0.56
1:B:146:ILE:CG2	1:B:289:ILE:HG12	2.36	0.56
1:B:38:GLU:HG3	2:B:475:HOH:O	2.05	0.56
1:B:115:ARG:HD3	2:B:513:HOH:O	2.06	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:172:ARG:HA	1:A:183:ILE:HD11	1.88	0.55
1:B:146:ILE:HD12	1:B:148:PHE:CE2	2.42	0.55
1:B:79:ARG:HD3	1:B:79:ARG:H	1.72	0.55
1:A:45:LYS:HG2	1:A:403:ILE:HD11	1.88	0.54
1:B:213:HIS:HE1	2:B:489:HOH:O	1.91	0.54
1:A:382:GLN:O	1:A:386:ARG:HD3	2.07	0.54
1:B:166:GLU:O	1:B:170:LYS:HG3	2.08	0.54
1:B:335:LYS:HD3	1:B:335:LYS:H	1.73	0.54
1:A:151:ASN:HD22	1:A:318:ASN:HD21	1.57	0.53
1:A:29:TYR:CZ	1:A:65:GLU:HB2	2.43	0.53
1:B:280:ILE:HG21	1:B:300:ILE:HD11	1.91	0.53
1:B:88:LEU:HD21	1:B:94:VAL:HG21	1.90	0.53
1:B:4:LYS:HD2	2:B:545:HOH:O	2.08	0.53
1:B:88:LEU:HD23	1:B:94:VAL:HG21	1.91	0.53
1:A:43:ALA:O	1:A:46:VAL:HG22	2.09	0.53
1:B:335:LYS:N	1:B:335:LYS:HD3	2.23	0.52
1:A:123:ASN:ND2	2:A:456:HOH:O	2.43	0.52
1:A:279:GLN:NE2	1:A:282:ILE:HD11	2.24	0.52
1:A:147:ILE:HD13	1:A:300:ILE:HG23	1.92	0.52
1:B:334:PRO:HD2	1:B:335:LYS:HZ3	1.74	0.52
1:B:196:LYS:O	1:B:200:GLN:HG3	2.10	0.51
1:B:308:ILE:CD1	1:B:314:LEU:HD21	2.41	0.51
1:A:122:ALA:O	1:A:318:ASN:HA	2.10	0.51
1:A:33:ASP:OD2	1:A:70:GLU:HG2	2.10	0.51
1:B:146:ILE:HD13	1:B:146:ILE:C	2.31	0.51
1:A:196:LYS:O	1:A:200:GLN:HG3	2.11	0.51
1:A:116:LEU:HD11	2:A:637:HOH:O	2.11	0.51
1:B:45:LYS:HG2	1:B:403:ILE:HD11	1.93	0.51
1:B:146:ILE:HD12	1:B:148:PHE:CD2	2.45	0.50
1:B:204:ARG:O	1:B:208:GLN:HG3	2.11	0.50
1:B:318:ASN:C	1:B:319:ILE:HD12	2.31	0.50
1:A:124:ILE:CG1	1:A:319:ILE:HD13	2.42	0.50
1:B:5:GLU:HG2	1:B:87:MET:CE	2.41	0.49
1:A:176:ARG:HH11	1:A:182:GLU:HG2	1.76	0.49
1:B:258:GLN:HE21	1:B:260:LYS:NZ	2.11	0.49
1:B:299:TYR:HD2	2:B:414:HOH:O	1.94	0.49
1:A:124:ILE:HG13	1:A:319:ILE:HD13	1.94	0.49
1:A:365:ASP:O	1:A:369:LYS:HG3	2.14	0.48
1:A:226:VAL:HG21	1:B:156:TYR:CE1	2.47	0.48
1:A:215:ARG:HD2	1:A:288:ASP:OD2	2.13	0.48
1:B:79:ARG:N	1:B:79:ARG:HD3	2.29	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:156:TYR:CE1	1:B:226:VAL:HG21	2.48	0.48
1:B:127:VAL:HG13	1:B:145:MET:HG3	1.94	0.47
1:A:21:VAL:HA	1:A:22:PRO:HD3	1.81	0.47
1:B:313:MET:HE3	1:B:345:THR:HG21	1.96	0.47
1:A:253:GLU:HB2	1:A:269:ASN:HB3	1.96	0.47
1:B:124:ILE:HG13	1:B:319:ILE:HD13	1.96	0.47
1:A:272:ILE:N	1:A:272:ILE:HD12	2.30	0.47
1:B:127:VAL:HG13	1:B:145:MET:CG	2.45	0.47
1:A:373:MET:HG2	1:A:376:ARG:NH2	2.29	0.47
1:B:109:SER:HB3	1:B:112:VAL:CG1	2.43	0.47
1:A:109:SER:OG	1:A:112:VAL:HG12	2.15	0.47
1:B:124:ILE:CG1	1:B:319:ILE:HD13	2.45	0.46
1:A:319:ILE:HD12	1:A:319:ILE:N	2.29	0.46
1:B:319:ILE:HD12	1:B:319:ILE:N	2.30	0.46
1:B:81:PRO:HG2	1:B:84:THR:OG1	2.16	0.46
1:A:141:GLU:H	1:A:141:GLU:CD	2.19	0.46
1:B:373:MET:O	1:B:376:ARG:HG2	2.15	0.46
1:A:199:THR:O	1:A:202:ILE:HG22	2.16	0.46
1:A:378:LYS:HG3	1:A:388:MET:SD	2.56	0.46
1:B:333:ALA:HA	1:B:335:LYS:HZ3	1.81	0.46
1:B:177:LYS:HE3	1:B:178:GLU:OE2	2.16	0.46
1:B:109:SER:CB	1:B:112:VAL:HG12	2.45	0.46
1:A:351:GLY:O	1:A:355:LEU:HD22	2.16	0.45
1:B:250:ILE:HG22	1:B:266:ILE:HB	1.99	0.45
1:B:310:ASN:HD22	1:B:312:GLY:H	1.65	0.45
1:A:38:GLU:HG3	2:A:600:HOH:O	2.17	0.45
1:A:45:LYS:HG2	1:A:403:ILE:CD1	2.47	0.44
1:B:308:ILE:HD11	1:B:314:LEU:HD21	1.99	0.44
1:B:276:MET:O	1:B:280:ILE:HG12	2.17	0.44
1:A:73:GLU:OE1	1:A:79:ARG:HD3	2.18	0.44
1:B:15:ASP:OD2	1:B:16:LYS:N	2.47	0.44
1:A:73:GLU:HG3	1:A:79:ARG:HH21	1.82	0.44
1:B:333:ALA:HA	1:B:335:LYS:NZ	2.32	0.44
1:A:115:ARG:HD3	1:A:318:ASN:ND2	2.33	0.44
1:B:326:PHE:CE2	1:B:354:MET:HA	2.53	0.44
1:A:149:ARG:HD3	1:A:294:ASN:HA	2.00	0.44
1:B:122:ALA:O	1:B:318:ASN:HA	2.17	0.44
1:B:129:TYR:CD1	1:B:141:GLU:HA	2.53	0.44
1:A:4:LYS:N	1:A:4:LYS:HD3	2.33	0.43
1:A:179:LEU:HB3	1:A:181:VAL:HG23	2.00	0.43
1:B:78:ASP:O	1:B:81:PRO:HD3	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:217:LYS:HG2	1:B:218:VAL:N	2.33	0.43
1:B:4:LYS:N	1:B:4:LYS:HD3	2.33	0.43
1:A:223:LYS:HD2	1:A:223:LYS:HA	1.76	0.43
1:A:71:LYS:O	1:A:75:LEU:HD13	2.19	0.43
1:A:254:GLU:O	1:A:258:GLN:HB2	2.18	0.43
1:B:73:GLU:O	1:B:77:ASN:HA	2.18	0.43
1:A:308:ILE:HD13	1:A:314:LEU:HD21	2.00	0.43
1:B:369:LYS:HE3	1:B:373:MET:SD	2.59	0.43
1:B:21:VAL:HA	1:B:22:PRO:HD3	1.81	0.43
1:B:313:MET:CE	1:B:345:THR:HG21	2.48	0.43
1:B:127:VAL:CG1	1:B:145:MET:HG3	2.49	0.42
1:A:310:ASN:ND2	1:A:384:ILE:HD11	2.34	0.42
1:A:279:GLN:HE21	1:A:283:ARG:HE	1.66	0.42
1:B:196:LYS:HA	1:B:237:TRP:CE3	2.55	0.42
1:B:149:ARG:HG3	1:B:297:GLY:CA	2.39	0.42
1:B:223:LYS:HE2	1:B:225:ASN:OD1	2.19	0.42
1:B:133:LEU:HG	1:B:375:ILE:HG21	2.02	0.42
1:B:146:ILE:O	1:B:146:ILE:HG23	2.19	0.42
1:A:8:ASP:HB3	1:A:62:LYS:HE3	2.02	0.42
1:A:308:ILE:CD1	1:A:314:LEU:HD21	2.49	0.42
1:A:341:VAL:HG13	1:A:382:GLN:HG2	2.02	0.42
1:A:214:LYS:NZ	1:A:214:LYS:HB2	2.34	0.42
1:A:196:LYS:HA	1:A:237:TRP:CE3	2.55	0.42
1:A:222:HIS:O	1:A:272:ILE:HA	2.20	0.41
1:A:326:PHE:CE2	1:A:354:MET:HA	2.55	0.41
1:A:278:GLN:HG3	1:B:302:ASP:HB3	2.02	0.41
1:B:246:TYR:O	1:B:250:ILE:HG13	2.20	0.41
1:B:35:ILE:HG12	1:B:332:THR:O	2.20	0.41
1:B:235:ARG:NH1	1:B:235:ARG:HB3	2.35	0.41
1:A:128:LYS:HD3	1:A:368:GLU:OE2	2.19	0.41
1:A:73:GLU:O	1:A:77:ASN:HA	2.19	0.41
1:B:226:VAL:HG23	1:B:227:MET:HG3	2.03	0.41
1:B:130:ILE:HG21	1:B:375:ILE:HD12	2.03	0.41
1:B:313:MET:HE1	1:B:375:ILE:HD11	2.03	0.41
1:B:252:THR:OG1	1:B:255:GLU:HG3	2.20	0.41
1:A:210:ALA:CA	1:A:215:ARG:HB2	2.47	0.41
1:A:223:LYS:HD2	1:A:272:ILE:HG23	2.03	0.41
1:B:42:ALA:O	1:B:46:VAL:HG23	2.21	0.41
1:A:130:ILE:HG21	1:A:375:ILE:HD12	2.03	0.41
1:B:239:TYR:O	1:B:243:LEU:HD13	2.21	0.41
1:A:1:MET:HE3	2:A:481:HOH:O	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:166:GLU:HG3	2:B:508:HOH:O	2.20	0.40
1:B:310:ASN:HA	2:B:606:HOH:O	2.21	0.40
1:A:280:ILE:HD11	1:A:304:ALA:HB2	2.03	0.40
1:A:203:THR:HA	1:A:291:LEU:HD11	2.03	0.40
1:B:147:ILE:CD1	1:B:301:SER:HA	2.52	0.40
1:A:223:LYS:HE3	1:A:274:ASP:OD2	2.22	0.40
1:A:125:ARG:HB2	1:A:147:ILE:HB	2.03	0.40
1:B:225:ASN:ND2	1:B:226:VAL:HG13	2.37	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	397/409 (97%)	382 (96%)	15 (4%)	0	100	100
1	B	397/409 (97%)	371 (94%)	25 (6%)	1 (0%)	46	41
All	All	794/818 (97%)	753 (95%)	40 (5%)	1 (0%)	56	53

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	310	ASN

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	337/343 (98%)	325 (96%)	12 (4%)	42	39
1	B	337/343 (98%)	321 (95%)	16 (5%)	32	27
All	All	674/686 (98%)	646 (96%)	28 (4%)	36	31

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	28	LEU
1	A	53	ARG
1	A	62	LYS
1	A	127	VAL
1	A	137	LEU
1	A	194	MET
1	A	216	LYS
1	A	244	LYS
1	A	274	ASP
1	A	291	LEU
1	A	355	LEU
1	A	357	TRP
1	B	1	MET
1	B	52	GLU
1	B	53	ARG
1	B	79	ARG
1	B	146	ILE
1	B	149	ARG
1	B	184	GLU
1	B	194	MET
1	B	216	LYS
1	B	260	LYS
1	B	291	LEU
1	B	310	ASN
1	B	355	LEU
1	B	357	TRP
1	B	378	LYS
1	B	382	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	85	GLN
1	A	123	ASN

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Mol	Chain	Res	Type
1	A	151	ASN
1	A	208	GLN
1	A	258	GLN
1	A	263	GLN
1	A	279	GLN
1	A	310	ASN
1	B	85	GLN
1	B	123	ASN
1	B	213	HIS
1	B	258	GLN
1	B	263	GLN
1	B	310	ASN
1	B	318	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 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	401/409 (98%)	0.49	30 (7%) 17 18	16, 29, 46, 66	0
1	B	401/409 (98%)	0.52	25 (6%) 24 25	19, 32, 54, 83	0
All	All	802/818 (98%)	0.50	55 (6%) 20 21	16, 31, 51, 83	0

All (55) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	264	GLY	5.7
1	B	257	ASN	5.4
1	B	256	ILE	5.3
1	B	1	MET	4.4
1	B	263	GLN	4.1
1	B	309	GLY	3.9
1	B	260	LYS	3.7
1	B	254	GLU	3.6
1	B	243	LEU	3.6
1	A	17	GLY	3.5
1	A	376	ARG	3.4
1	A	1	MET	3.4
1	A	129	TYR	3.3
1	A	263	GLN	3.2
1	A	257	ASN	3.1
1	A	254	GLU	3.1
1	B	258	GLN	3.1
1	A	283	ARG	3.0
1	A	243	LEU	3.0
1	A	131	GLU	2.8
1	B	311	ILE	2.8
1	A	183	ILE	2.8
1	B	241	VAL	2.8
1	B	133	LEU	2.8

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Mol	Chain	Res	Type	RSRZ
1	B	15	ASP	2.7
1	A	258	GLN	2.6
1	B	208	GLN	2.6
1	A	409	ILE	2.6
1	A	176	ARG	2.6
1	B	390	VAL	2.5
1	B	127	VAL	2.5
1	A	18	LYS	2.5
1	B	409	ILE	2.5
1	B	4	LYS	2.4
1	A	311	ILE	2.4
1	A	250	ILE	2.4
1	A	336	TYR	2.4
1	A	7	GLU	2.4
1	A	282	ILE	2.3
1	B	17	GLY	2.3
1	A	319	ILE	2.3
1	A	375	ILE	2.3
1	A	77	ASN	2.2
1	A	174	PHE	2.2
1	A	16	LYS	2.2
1	A	248	ASP	2.2
1	B	303	ALA	2.2
1	B	307	LEU	2.1
1	A	396	LYS	2.1
1	B	376	ARG	2.1
1	A	124	ILE	2.1
1	A	295	VAL	2.1
1	A	391	LYS	2.0
1	B	74	LYS	2.0
1	B	261	PRO	2.0

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

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

6.3 Carbohydrates ⓘ

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.