



# Full wwPDB X-ray Structure Validation Report ⓘ

May 26, 2025 – 05:42 AM EDT

PDB ID : 2E0C / pdb\_00002e0c  
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](mailto: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>.

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

MolProbity	:	4-5-2 with Phenix2.0rc1
Xtriage (Phenix)	:	2.0rc1
EDS	:	3.0
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.006 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.43.1

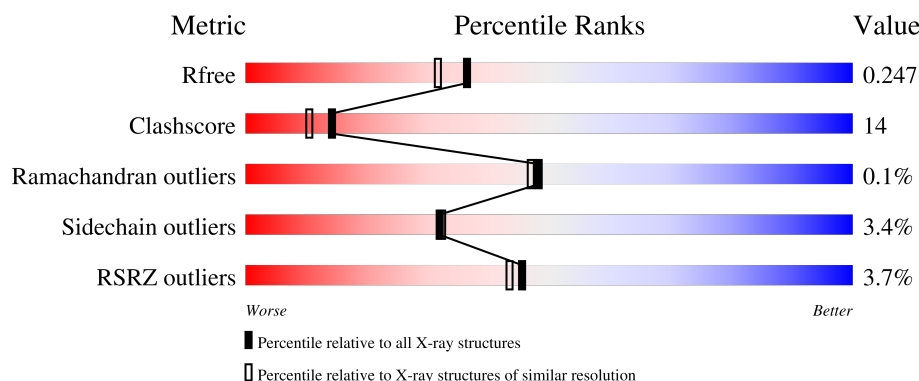
# 1 Overall quality at a glance

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))
$R_{free}$	164625	9409 (2.00-2.00)
Clashscore	180529	10737 (2.00-2.00)
Ramachandran outliers	177936	10628 (2.00-2.00)
Sidechain outliers	177891	10627 (2.00-2.00)
RSRZ outliers	164620	9409 (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 of the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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	409	
1	B	409	

## 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			

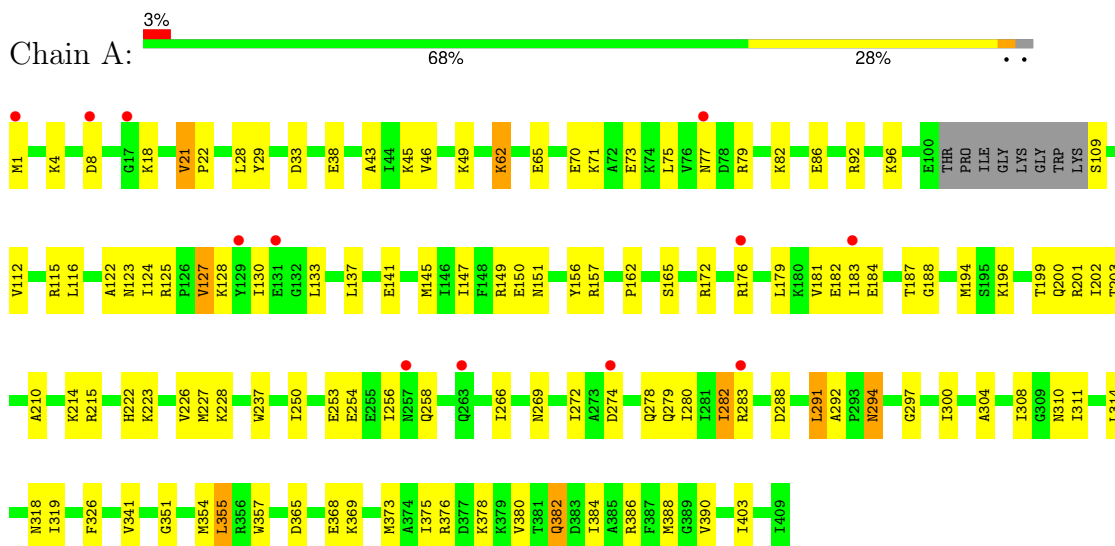
- Molecule 2 is water.

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, 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: 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, $\alpha$ , $\beta$ , $\gamma$	74.88Å 87.72Å 75.72Å 90.00° 91.37° 90.00°	Depositor
Resolution (Å)	15.00 – 2.00 15.00 – 2.00	Depositor EDS
% Data completeness (in resolution range)	99.2 (15.00-2.00) 99.0 (15.00-2.00)	Depositor EDS
$R_{merge}$	0.06	Depositor
$R_{sym}$	0.06	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.26 (at 2.00Å)	Xtriage
Refinement program	CNS	Depositor
R, $R_{free}$	0.221 , 0.247 0.221 , 0.247	Depositor DCC
$R_{free}$ test set	3320 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	30.4	Xtriage
Anisotropy	0.116	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.39 , 49.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.004 for l,k,-h 0.066 for h,-k,-l 0.018 for l,-k,h	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	6881	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	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.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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	0.40	0/3264	0.95	12/4388 (0.3%)
1	B	0.37	0/3264	0.91	14/4388 (0.3%)
All	All	0.39	0/6528	0.93	26/8776 (0.3%)

There are no bond length outliers.

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	21	VAL	N-CA-C	8.75	117.57	107.77
1	B	187	THR	N-CA-C	7.48	121.12	109.07
1	A	150	GLU	N-CA-C	-7.31	99.00	109.96
1	B	21	VAL	N-CA-C	7.31	115.95	107.77
1	A	188	GLY	N-CA-C	-7.10	102.43	112.37
1	A	228	LYS	N-CA-C	6.71	118.59	111.28
1	A	292	ALA	N-CA-C	6.59	115.90	108.25
1	B	188	GLY	N-CA-C	-6.49	103.28	112.37
1	A	187	THR	N-CA-C	6.40	119.33	108.90
1	A	294	ASN	N-CA-C	6.38	118.81	111.02
1	B	150	GLU	N-CA-C	-6.38	100.40	109.96
1	B	292	ALA	N-CA-C	6.14	115.37	108.25
1	B	228	LYS	N-CA-C	6.10	118.43	111.11
1	B	350	ALA	N-CA-C	-6.01	104.80	111.36
1	B	294	ASN	N-CA-C	5.91	118.23	111.02
1	A	92	ARG	N-CA-C	5.78	123.11	110.80
1	A	157	ARG	N-CA-C	-5.62	106.26	113.01
1	A	382	GLN	N-CA-C	5.44	118.71	111.75
1	B	92	ARG	N-CA-C	5.43	122.37	110.80
1	A	380	VAL	N-CA-C	5.42	116.51	108.65
1	B	96	LYS	N-CA-C	5.30	117.54	108.90
1	B	189	ILE	N-CA-C	5.27	115.71	108.12
1	B	216	LYS	N-CA-C	5.27	119.98	113.50
1	B	380	VAL	N-CA-C	5.25	116.46	108.80
1	A	96	LYS	N-CA-C	5.01	117.07	108.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	382	GLN	N-CA-C	5.01	117.39	111.33

There are no chirality outliers.

There are no planarity outliers.

## 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	3209	0	3285	84	0
1	B	3209	0	3285	101	0
2	A	249	0	0	6	0
2	B	214	0	0	8	0
All	All	6881	0	6570	182	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 14.

All (182) 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
1:A:149:ARG:HG3	1:A:297:GLY:HA3	1.47	0.95
1:B:5:GLU:HG2	1:B:87:MET:HE1	1.51	0.92
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:A:223:LYS:HG3	1:A:226:VAL:HG22	1.68	0.76
1:A:250:ILE:HG22	1:A:266:ILE:HB	1.67	0.75
1:B:124:ILE:CG2	1:B:146:ILE:HD11	2.16	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:29:TYR:CZ	1:B:65:GLU:HB2	2.32	0.65
1:B:223:LYS:HE3	1:B:274:ASP:OD2	1.97	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:A:130:ILE:HB	1:A:133:LEU:HD12	1.84	0.59
1:B:146:ILE:HD13	1:B:147:ILE:N	2.17	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:A:127:VAL:HG13	1:A:145:MET:HG3	1.86	0.56
1:B:210:ALA:HA	1:B:215:ARG:HB2	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:146:ILE:HD13	1:B:146:ILE:C	2.31	0.55
1:B:318:ASN:C	1:B:319:ILE:HD12	2.31	0.55
1:B:115:ARG:HD3	2:B:513:HOH:O	2.06	0.55
1:A:172:ARG:HA	1:A:183:ILE:HD11	1.88	0.55
1:B:79:ARG:HD3	1:B:79:ARG:H	1.72	0.55
1:B:146:ILE:HD12	1:B:148:PHE:CE2	2.42	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:29:TYR:CZ	1:A:65:GLU:HB2	2.43	0.53
1:A:151:ASN:HD22	1:A:318:ASN:HD21	1.57	0.53
1:B:221:MET:HE2	1:B:296:ASN:HB3	1.90	0.53
1:B:280:ILE:HG21	1:B:300:ILE:HD11	1.91	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:4:LYS:HD2	2:B:545:HOH:O	2.08	0.53
1:B:88:LEU:HD21	1:B:94:VAL:HG21	1.90	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:HD3	1:B:335:LYS:N	2.23	0.52
1:A:123:ASN:ND2	2:A:456:HOH:O	2.43	0.52
1:A:147:ILE:HD13	1:A:300:ILE:HG23	1.92	0.52
1:A:279:GLN:NE2	1:A:282:ILE:HD11	2.24	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:B:334:PRO:HD2	1:B:335:LYS:HZ3	1.75	0.51
1:A:33:ASP:OD2	1:A:70:GLU:HG2	2.10	0.51
1:A:122:ALA:O	1:A:318:ASN:HA	2.10	0.51
1:A:141:GLU:H	1:A:141:GLU:CD	2.19	0.51
1:A:116:LEU:HD11	2:A:637:HOH:O	2.11	0.51
1:A:196:LYS:O	1:A:200:GLN:HG3	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:A:223:LYS:HB3	1:A:227:MET:HE3	1.93	0.50
1:B:204:ARG:O	1:B:208:GLN:HG3	2.11	0.50
1:A:124:ILE:CG1	1:A:319:ILE:HD13	2.42	0.50
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:B:333:ALA:HA	1:B:335:LYS:HZ3	1.78	0.48
1:A:365:ASP:O	1:A:369:LYS:HG3	2.14	0.48
1:A:215:ARG:HD2	1:A:288:ASP:OD2	2.13	0.48
1:A:226:VAL:HG21	1:B:156:TYR:CE1	2.47	0.48
1:B:79:ARG:HD3	1:B:79:ARG:N	2.29	0.48
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:A:253:GLU:HB2	1:A:269:ASN:HB3	1.96	0.47
1:A:272:ILE:HD12	1:A:272:ILE:N	2.30	0.47
1:B:124:ILE:HG13	1:B:319:ILE:HD13	1.96	0.47
1:A:373:MET:HG2	1:A:376:ARG:NH2	2.29	0.47
1:B:127:VAL:HG13	1:B:145:MET:CG	2.45	0.47
1:A:109:SER:OG	1:A:112:VAL:HG12	2.15	0.47
1:B:109:SER:HB3	1:B:112:VAL:CG1	2.43	0.47
1:A:319:ILE:HD12	1:A:319:ILE:N	2.29	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:124:ILE:CG1	1:B:319:ILE:HD13	2.45	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: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: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:B:313:MET:HE1	1:B:375:ILE:HD11	1.98	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:A:184:GLU:H	1:A:184:GLU:CD	2.25	0.44
1:B:256:ILE:C	1:B:258:GLN:H	2.25	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: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:B:333:ALA:HA	1:B:335:LYS:NZ	2.32	0.44
1:A:149:ARG:HD3	1:A:294:ASN:HA	2.00	0.44
1:A:318:ASN:C	1:A:319:ILE:HD12	2.43	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:256:ILE:C	1:A:258:GLN:H	2.26	0.43
1:A:179:LEU:HB3	1:A:181:VAL:HG23	2.00	0.43
1:A:250:ILE:HD12	1:A:250:ILE:C	2.43	0.43
1:B:78:ASP:O	1:B:81:PRO:HD3	2.18	0.43
1:B:313:MET:HE2	1:B:345:THR:HG21	1.99	0.43
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:71:LYS:O	1:A:75:LEU:HD13	2.19	0.43
1:A:223:LYS:HD2	1:A:223:LYS:HA	1.76	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:21:VAL:HA	1:B:22:PRO:HD3	1.81	0.43
1:B:369:LYS:HE3	1:B:373:MET:SD	2.59	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:133:LEU:HG	1:B:375:ILE:HG21	2.02	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:A:8:ASP:HB3	1:A:62:LYS:HE3	2.02	0.42
1:B:146:ILE:O	1:B:146:ILE:HG23	2.19	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:HB2	1:A:214:LYS:NZ	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:278:GLN:HG3	1:B:302:ASP:HB3	2.02	0.41
1:A:326:PHE:CE2	1:A:354:MET:HA	2.55	0.41
1:B:35:ILE:HG12	1:B:332:THR:O	2.20	0.41
1:B:235:ARG:HB3	1:B:235:ARG:NH1	2.35	0.41
1:B:246:TYR:O	1:B:250:ILE:HG13	2.20	0.41
1:A:128:LYS:HD3	1:A:368:GLU:OE2	2.19	0.41
1:B:64:LEU:C	1:B:64:LEU:HD13	2.45	0.41
1:A:73:GLU:O	1:A:77:ASN:HA	2.19	0.41
1:B:130:ILE:HG21	1:B:375:ILE:HD12	2.03	0.41
1:B:226:VAL:HG23	1:B:227:MET:HG3	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:B:313:MET:HE1	1:B:375:ILE:CD1	2.51	0.41
1:A:1:MET:HE3	2:A:481:HOH:O	2.21	0.40
1:B:178:GLU:C	1:B:179:LEU:HD22	2.47	0.40
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:203:THR:HA	1:A:291:LEU:HD11	2.03	0.40
1:A:280:ILE:HD11	1:A:304:ALA:HB2	2.03	0.40
1:A:125:ARG:HB2	1:A:147:ILE:HB	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:223:LYS:HE3	1:A:274:ASP:OD2	2.22	0.40
1:B:147:ILE:CD1	1:B:301:SER:HA	2.52	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%)	37	35
All	All	794/818 (97%)	753 (95%)	40 (5%)	1 (0%)	48	47

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%)	327 (97%)	10 (3%)	36	37
1	B	337/343 (98%)	324 (96%)	13 (4%)	27	27

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
All	All	674/686 (98%)	651 (97%)	23 (3%)	32	32

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	18	LYS
1	A	28	LEU
1	A	62	LYS
1	A	127	VAL
1	A	137	LEU
1	A	194	MET
1	A	282	ILE
1	A	291	LEU
1	A	355	LEU
1	A	357	TRP
1	B	28	LEU
1	B	52	GLU
1	B	127	VAL
1	B	146	ILE
1	B	149	ARG
1	B	184	GLU
1	B	194	MET
1	B	235	ARG
1	B	291	LEU
1	B	355	LEU
1	B	357	TRP
1	B	378	LYS
1	B	382	GLN

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

Mol	Chain	Res	Type
1	A	85	GLN
1	A	123	ASN
1	A	151	ASN
1	A	258	GLN
1	A	263	GLN
1	A	279	GLN
1	A	310	ASN
1	A	361	ASN
1	B	23	ASN

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Mol	Chain	Res	Type
1	B	85	GLN
1	B	123	ASN
1	B	139	HIS
1	B	151	ASN
1	B	213	HIS
1	B	258	GLN
1	B	263	GLN
1	B	310	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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	401/409 (98%)	0.04	12 (2%) 52 51	16, 29, 46, 66	0
1	B	401/409 (98%)	0.20	18 (4%) 39 37	19, 32, 54, 83	0
All	All	802/818 (98%)	0.12	30 (3%) 45 43	16, 31, 51, 83	0

All (30) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	309	GLY	4.5
1	B	1	MET	4.1
1	A	17	GLY	3.6
1	B	256	ILE	3.5
1	A	257	ASN	3.4
1	B	15	ASP	3.3
1	B	264	GLY	3.3
1	A	129	TYR	3.3
1	B	257	ASN	3.2
1	B	259	GLY	3.1
1	A	176	ARG	3.0
1	A	283	ARG	2.9
1	A	1	MET	2.6
1	A	77	ASN	2.6
1	B	8	ASP	2.5
1	B	274	ASP	2.4
1	B	286	GLU	2.3
1	A	131	GLU	2.3
1	B	133	LEU	2.3
1	B	258	GLN	2.3
1	B	243	LEU	2.2
1	A	263	GLN	2.2
1	A	274	ASP	2.2
1	A	183	ILE	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	280	ILE	2.2
1	A	8	ASP	2.2
1	B	263	GLN	2.1
1	B	311	ILE	2.1
1	B	310	ASN	2.1
1	B	17	GLY	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 monosaccharides 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.