



# Full wwPDB X-ray Structure Validation Report ⓘ

May 22, 2020 – 12:24 am BST

PDB ID : 2CGK  
Title : Crystal Structure of L-rhamnulose kinase from Escherichia coli in an open uncomplexed conformation.  
Authors : Grueninger, D.; Schulz, G.E.  
Deposited on : 2006-03-09  
Resolution : 2.46 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

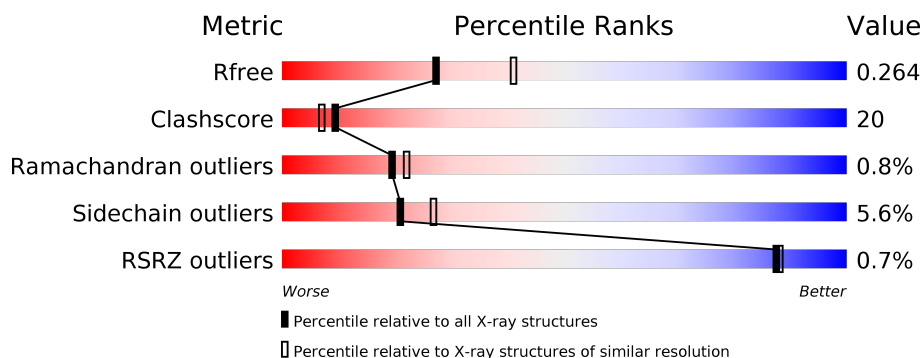
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*



The reported resolution of this entry is 2.46 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))
$R_{free}$	130704	1544 (2.48-2.44)
Clashscore	141614	1613 (2.48-2.44)
Ramachandran outliers	138981	1598 (2.48-2.44)
Sidechain outliers	138945	1598 (2.48-2.44)
RSRZ outliers	127900	1523 (2.48-2.44)

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  $\geq 3$ , 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	489	 63% 32% . .
1	B	489	 % 62% 32% . .

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 7498 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 L-RHAMNULOSE KINASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	479	Total	C	N	O	S	0	0	1
			3702	2319	657	706	20			
1	B	479	Total	C	N	O	S	0	0	1
			3702	2319	657	706	20			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	69	ALA	GLU	engineered mutation	UNP Q8X899
A	70	ALA	GLU	engineered mutation	UNP Q8X899
A	73	ALA	ARG	engineered mutation	UNP Q8X899
B	70	ALA	GLU	engineered mutation	UNP Q8X899
B	69	ALA	GLU	engineered mutation	UNP Q8X899
B	73	ALA	ARG	engineered mutation	UNP Q8X899
A	320	ALA	SER	conflict	UNP Q8X899
A	343	ASP	GLU	conflict	UNP Q8X899
A	344	GLU	THR	conflict	UNP Q8X899
A	477	LEU	ARG	conflict	UNP Q8X899
A	356	MET	THR	conflict	UNP Q8X899
B	320	ALA	SER	conflict	UNP Q8X899
B	343	ASP	GLU	conflict	UNP Q8X899
B	344	GLU	THR	conflict	UNP Q8X899
B	356	MET	THR	conflict	UNP Q8X899
B	477	LEU	ARG	conflict	UNP Q8X899

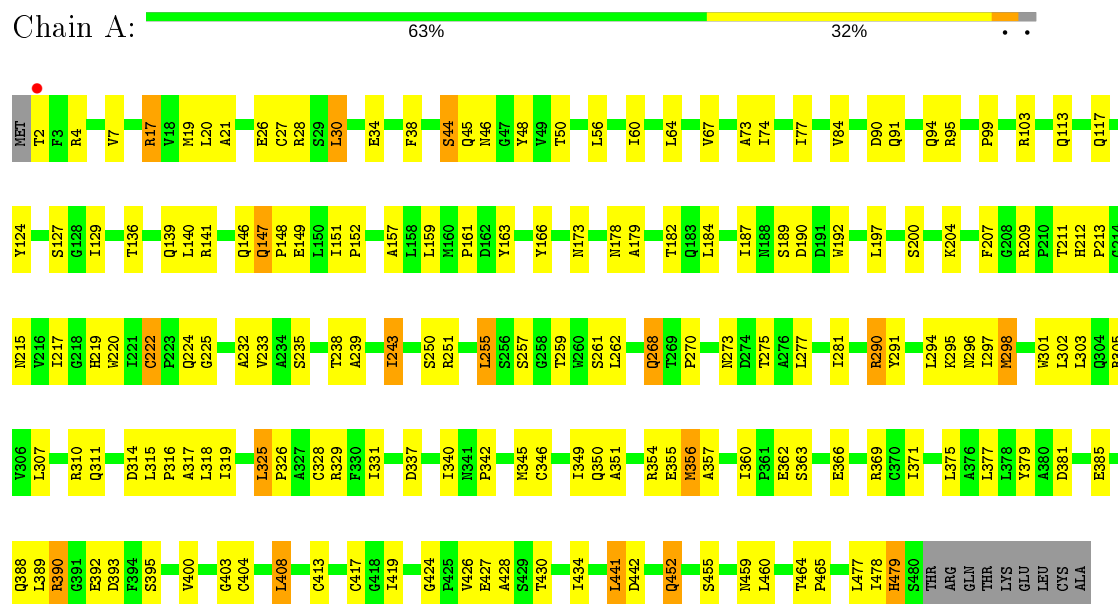
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	46	Total	O	0	0
			46	46		
2	B	48	Total	O	0	0
			48	48		

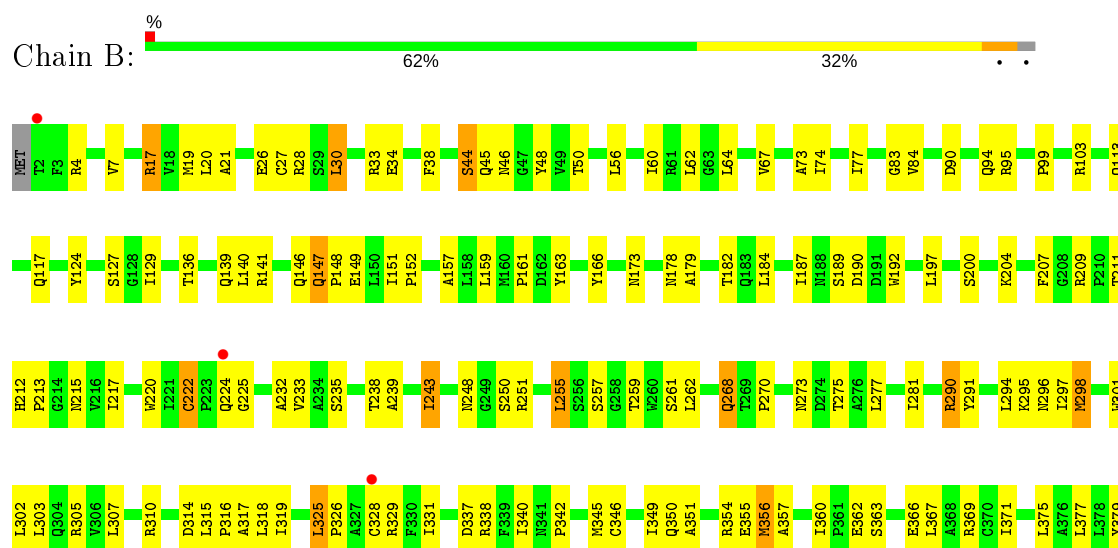
### 3 Residue-property plots

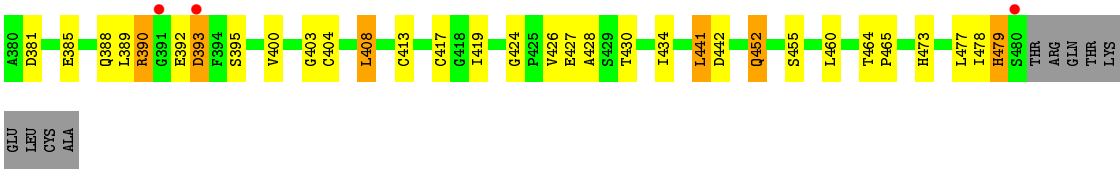
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are 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: L-RHAMNULOSE KINASE



#### • Molecule 1: L-RHAMNULOSE KINASE





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	52.26 Å   164.93 Å   51.57 Å 90.00°   93.45°   90.00°	Depositor
Resolution (Å)	44.00 – 2.46 43.67 – 2.46	Depositor EDS
% Data completeness (in resolution range)	94.7 (44.00-2.46) 95.1 (43.67-2.46)	Depositor EDS
$R_{merge}$	0.10	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.95 (at 2.45 Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.238 , 0.273 0.232 , 0.264	Depositor DCC
$R_{free}$ test set	1792 reflections (6.00%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	30.9	Xtriage
Anisotropy	0.425	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 13.8	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.44$ , $\langle L^2 \rangle = 0.26$	Xtriage
Estimated twinning fraction	0.059 for l,-k,h	Xtriage
$F_o, F_c$ correlation	0.90	EDS
Total number of atoms	7498	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 5.29% 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	0/3777	0.65	0/5147
1	B	0.40	0/3777	0.65	0/5147
All	All	0.40	0/7554	0.65	0/10294

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

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	3702	0	3613	145	3
1	B	3702	0	3613	147	4
2	A	46	0	0	0	0
2	B	48	0	0	1	0
All	All	7498	0	7226	288	5

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 20.

All (288) 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:33:ARG:HD2	2:B:2006:HOH:O	1.51	1.09

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:217:ILE:HD11	1:A:232:ALA:HB2	1.48	0.96
1:B:217:ILE:HD11	1:B:232:ALA:HB2	1.48	0.96
1:A:281:ILE:HD13	1:A:389:LEU:HD22	1.54	0.89
1:B:281:ILE:HD13	1:B:389:LEU:HD22	1.54	0.89
1:B:187:ILE:HD11	1:B:291:TYR:CE1	2.13	0.83
1:A:187:ILE:HD11	1:A:291:TYR:CE1	2.13	0.82
1:A:19:MET:SD	1:A:34:GLU:HG2	2.19	0.82
1:B:19:MET:SD	1:B:34:GLU:HG2	2.19	0.82
1:B:4:ARG:HH21	1:B:441:LEU:HG	1.43	0.81
1:A:4:ARG:HH21	1:A:441:LEU:HG	1.43	0.81
1:A:302:LEU:HB3	1:A:371:ILE:HD12	1.64	0.79
1:A:91:GLN:HG2	1:B:48:TYR:CZ	2.17	0.79
1:B:302:LEU:HB3	1:B:371:ILE:HD12	1.64	0.78
1:B:90:ASP:OD2	1:B:94:GLN:HB2	1.89	0.73
1:A:90:ASP:OD2	1:A:94:GLN:HB2	1.89	0.72
1:B:354:ARG:HG2	1:B:360:ILE:HD11	1.70	0.72
1:B:297:ILE:HG21	1:B:379:TYR:HE1	1.53	0.72
1:B:239:ALA:O	1:B:243:ILE:HD13	1.89	0.72
1:A:239:ALA:O	1:A:243:ILE:HD13	1.89	0.72
1:A:297:ILE:HG21	1:A:379:TYR:HE1	1.53	0.72
1:A:354:ARG:HG2	1:A:360:ILE:HD11	1.70	0.72
1:B:328:CYS:HA	1:B:331:ILE:HD11	1.72	0.72
1:A:328:CYS:HA	1:A:331:ILE:HD11	1.72	0.71
1:B:268:GLN:HG3	1:B:290:ARG:CZ	2.21	0.70
1:A:268:GLN:HG3	1:A:290:ARG:CZ	2.21	0.70
1:A:297:ILE:HG21	1:A:379:TYR:CE1	2.28	0.68
1:B:297:ILE:HG21	1:B:379:TYR:CE1	2.28	0.68
1:A:281:ILE:CD1	1:A:389:LEU:HD22	2.23	0.68
1:B:281:ILE:CD1	1:B:389:LEU:HD22	2.24	0.68
1:A:408:LEU:O	1:A:408:LEU:HD12	1.95	0.67
1:B:408:LEU:O	1:B:408:LEU:HD12	1.95	0.67
1:A:220:TRP:CZ3	1:A:222:CYS:SG	2.89	0.66
1:A:28:ARG:HH21	1:A:452:GLN:NE2	1.94	0.66
1:B:28:ARG:HH21	1:B:452:GLN:NE2	1.94	0.66
1:B:220:TRP:CZ3	1:B:222:CYS:SG	2.89	0.65
1:A:315:LEU:HB3	1:A:316:PRO:HD3	1.77	0.65
1:B:319:ILE:HG23	1:B:408:LEU:HD23	1.77	0.65
1:A:351:ALA:O	1:A:355:GLU:HG3	1.97	0.65
1:B:351:ALA:O	1:B:355:GLU:HG3	1.97	0.65
1:B:315:LEU:HB3	1:B:316:PRO:HD3	1.77	0.64
1:B:27:CYS:O	1:B:28:ARG:HB3	1.98	0.64

*Continued on next page...*



*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:319:ILE:HG23	1:A:408:LEU:HD23	1.77	0.64
1:A:27:CYS:O	1:A:28:ARG:HB3	1.98	0.64
1:A:305:ARG:NH2	1:A:342:PRO:O	2.30	0.64
1:B:305:ARG:NH2	1:B:342:PRO:O	2.30	0.64
1:A:262:LEU:HD23	1:A:296:ASN:HA	1.79	0.63
1:A:315:LEU:O	1:A:318:LEU:HB3	1.99	0.63
1:B:315:LEU:O	1:B:318:LEU:HB3	1.99	0.63
1:B:262:LEU:HD23	1:B:296:ASN:HA	1.79	0.63
1:A:91:GLN:HG2	1:B:48:TYR:CE2	2.34	0.62
1:A:325:LEU:HD21	1:A:366:GLU:HG2	1.80	0.62
1:A:296:ASN:O	1:A:297:ILE:HD12	2.00	0.62
1:B:296:ASN:O	1:B:297:ILE:HD12	2.00	0.62
1:A:238:THR:HG21	1:A:262:LEU:CD1	2.30	0.61
1:B:187:ILE:HD11	1:B:291:TYR:HE1	1.64	0.61
1:B:238:THR:HG21	1:B:262:LEU:CD1	2.30	0.61
1:B:325:LEU:HD21	1:B:366:GLU:HG2	1.80	0.61
1:B:345:MET:O	1:B:349:ILE:HG12	2.00	0.61
1:B:316:PRO:O	1:B:317:ALA:HB3	1.99	0.61
1:A:345:MET:O	1:A:349:ILE:HG12	2.00	0.61
1:A:316:PRO:O	1:A:317:ALA:HB3	1.99	0.61
1:A:7:VAL:HB	1:A:77:ILE:HD13	1.83	0.61
1:B:250:SER:O	1:B:390:ARG:NH2	2.27	0.61
1:B:7:VAL:HB	1:B:77:ILE:HD13	1.83	0.61
1:A:326:PRO:O	1:A:369:ARG:HD3	2.01	0.61
1:A:250:SER:O	1:A:390:ARG:NH2	2.27	0.60
1:A:259:THR:HG23	1:A:301:TRP:HB2	1.83	0.60
1:B:326:PRO:O	1:B:369:ARG:HD3	2.01	0.60
1:A:147:GLN:HA	1:A:147:GLN:HE21	1.66	0.60
1:B:259:THR:HG23	1:B:301:TRP:HB2	1.83	0.60
1:B:147:GLN:HE21	1:B:147:GLN:HA	1.66	0.60
1:B:136:THR:HB	1:B:182:THR:HB	1.83	0.59
1:A:136:THR:HB	1:A:182:THR:HB	1.83	0.59
1:A:30:LEU:O	1:A:30:LEU:HD23	2.03	0.59
1:B:140:LEU:HD12	1:B:184:LEU:HD11	1.84	0.59
1:B:30:LEU:HD23	1:B:30:LEU:O	2.03	0.58
1:A:140:LEU:HD12	1:A:184:LEU:HD11	1.84	0.58
1:B:220:TRP:HZ3	1:B:222:CYS:SG	2.27	0.58
1:A:220:TRP:HZ3	1:A:222:CYS:SG	2.27	0.58
1:B:273:ASN:OD1	1:B:275:THR:HB	2.04	0.57
1:B:30:LEU:C	1:B:30:LEU:HD23	2.24	0.57
1:A:273:ASN:OD1	1:A:275:THR:HB	2.04	0.57

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:152:PRO:HB3	1:B:44:SER:CB	2.34	0.57
1:A:173:ASN:CB	1:A:217:ILE:HD12	2.34	0.57
1:B:173:ASN:CB	1:B:217:ILE:HD12	2.34	0.57
1:A:30:LEU:C	1:A:30:LEU:HD23	2.24	0.57
1:A:152:PRO:HB3	1:B:44:SER:HB3	1.86	0.56
1:A:316:PRO:C	1:A:318:LEU:H	2.07	0.56
1:A:67:VAL:CG1	1:A:74:ILE:HD11	2.35	0.56
1:B:67:VAL:CG1	1:B:74:ILE:HD11	2.35	0.56
1:B:316:PRO:C	1:B:318:LEU:H	2.07	0.56
1:A:187:ILE:HD11	1:A:291:TYR:HE1	1.64	0.56
1:B:346:CYS:O	1:B:350:GLN:HG3	2.05	0.55
1:A:346:CYS:O	1:A:350:GLN:HG3	2.06	0.55
1:B:4:ARG:NH2	1:B:441:LEU:HG	2.19	0.55
1:A:4:ARG:NH2	1:A:441:LEU:HG	2.19	0.55
1:A:290:ARG:HE	1:A:290:ARG:HA	1.71	0.55
1:B:290:ARG:HA	1:B:290:ARG:HE	1.71	0.54
1:B:255:LEU:HD12	1:B:255:LEU:C	2.28	0.54
1:A:255:LEU:HD12	1:A:255:LEU:C	2.28	0.54
1:A:268:GLN:HG3	1:A:290:ARG:NH1	2.22	0.54
1:B:455:SER:HA	1:B:460:LEU:HD11	1.90	0.54
1:A:455:SER:HA	1:A:460:LEU:HD11	1.90	0.54
1:B:268:GLN:HG3	1:B:290:ARG:NH1	2.22	0.54
1:A:173:ASN:HB2	1:A:217:ILE:HD12	1.88	0.53
1:A:67:VAL:HG11	1:A:74:ILE:HD11	1.91	0.53
1:B:173:ASN:HB2	1:B:217:ILE:HD12	1.88	0.53
1:B:67:VAL:HG11	1:B:74:ILE:HD11	1.91	0.53
1:A:316:PRO:O	1:A:318:LEU:N	2.39	0.53
1:B:477:LEU:C	1:B:477:LEU:HD12	2.30	0.53
1:A:477:LEU:C	1:A:477:LEU:HD12	2.30	0.53
1:B:124:TYR:CG	1:B:277:LEU:HD13	2.44	0.53
1:A:124:TYR:CG	1:A:277:LEU:HD13	2.44	0.52
1:B:45:GLN:O	1:B:46:ASN:HB2	2.09	0.52
1:A:17:ARG:CB	1:A:17:ARG:HH11	2.22	0.52
1:B:161:PRO:CB	1:B:235:SER:HB2	2.39	0.52
1:B:17:ARG:HH11	1:B:17:ARG:CB	2.22	0.52
1:A:45:GLN:O	1:A:46:ASN:HB2	2.09	0.52
1:A:161:PRO:CB	1:A:235:SER:HB2	2.39	0.52
1:B:316:PRO:O	1:B:318:LEU:N	2.39	0.52
1:A:127:SER:O	1:A:129:ILE:HD13	2.09	0.51
1:B:127:SER:O	1:B:129:ILE:HD13	2.09	0.51
1:A:379:TYR:HD2	1:A:413:CYS:SG	2.33	0.51

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:95:ARG:HH21	1:B:99:PRO:HD3	1.75	0.51
1:A:298:MET:SD	1:A:302:LEU:HD13	2.51	0.51
1:B:298:MET:SD	1:B:302:LEU:HD13	2.51	0.51
1:A:159:LEU:HB3	1:A:178:ASN:ND2	2.26	0.51
1:A:21:ALA:HB2	1:A:434:ILE:HD13	1.93	0.51
1:B:21:ALA:HB2	1:B:434:ILE:HD13	1.93	0.51
1:B:251:ARG:HD3	1:B:392:GLU:OE1	2.11	0.51
1:A:251:ARG:HD3	1:A:392:GLU:OE1	2.11	0.51
1:A:95:ARG:HH21	1:A:99:PRO:HD3	1.75	0.51
1:B:379:TYR:HD2	1:B:413:CYS:SG	2.33	0.51
1:A:295:LYS:O	1:A:297:ILE:HD13	2.11	0.50
1:B:159:LEU:HB3	1:B:178:ASN:ND2	2.26	0.50
1:B:187:ILE:HG12	1:B:270:PRO:HG3	1.94	0.50
1:B:295:LYS:O	1:B:297:ILE:HD13	2.11	0.50
1:B:314:ASP:OD1	1:B:316:PRO:O	2.28	0.50
1:A:187:ILE:HG12	1:A:270:PRO:HG3	1.94	0.50
1:A:314:ASP:OD1	1:A:316:PRO:O	2.28	0.50
1:B:360:ILE:N	1:B:360:ILE:HD12	2.25	0.50
1:A:360:ILE:HD12	1:A:360:ILE:N	2.25	0.50
1:B:141:ARG:HA	1:B:200:SER:O	2.11	0.50
1:A:141:ARG:HA	1:A:200:SER:O	2.11	0.50
1:B:151:ILE:HB	1:B:152:PRO:HD3	1.94	0.49
1:A:173:ASN:CG	1:A:217:ILE:HD12	2.32	0.49
1:A:259:THR:CG2	1:A:301:TRP:HB2	2.42	0.49
1:B:44:SER:HA	1:B:48:TYR:O	2.12	0.49
1:B:173:ASN:CG	1:B:217:ILE:HD12	2.32	0.49
1:A:151:ILE:HB	1:A:152:PRO:HD3	1.94	0.49
1:A:255:LEU:HA	1:A:262:LEU:O	2.12	0.49
1:A:44:SER:HA	1:A:48:TYR:O	2.12	0.49
1:B:259:THR:CG2	1:B:301:TRP:HB2	2.42	0.49
1:B:316:PRO:C	1:B:318:LEU:N	2.66	0.49
1:A:316:PRO:C	1:A:318:LEU:N	2.66	0.49
1:B:255:LEU:HA	1:B:262:LEU:O	2.12	0.49
1:A:310:ARG:NH2	1:A:362:GLU:O	2.46	0.49
1:B:310:ARG:NH2	1:B:362:GLU:O	2.46	0.49
1:B:262:LEU:CD2	1:B:296:ASN:HA	2.44	0.48
1:A:262:LEU:CD2	1:A:296:ASN:HA	2.44	0.48
1:A:238:THR:HG21	1:A:262:LEU:HD13	1.94	0.48
1:A:441:LEU:O	1:A:442:ASP:HB3	2.13	0.48
1:B:441:LEU:O	1:B:442:ASP:HB3	2.13	0.48
1:B:328:CYS:CA	1:B:331:ILE:HD11	2.41	0.48

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:17:ARG:HH11	1:B:17:ARG:HB2	1.79	0.48
1:A:303:LEU:O	1:A:307:LEU:HG	2.13	0.48
1:A:404:CYS:SG	1:A:426:VAL:HG22	2.54	0.48
1:B:238:THR:HG21	1:B:262:LEU:HD13	1.94	0.48
1:B:303:LEU:O	1:B:307:LEU:HG	2.13	0.48
1:B:404:CYS:SG	1:B:426:VAL:HG22	2.54	0.48
1:A:17:ARG:HB2	1:A:17:ARG:HH11	1.79	0.48
1:A:328:CYS:CA	1:A:331:ILE:HD11	2.41	0.47
1:A:157:ALA:HB3	1:A:207:PHE:CE1	2.50	0.47
1:B:157:ALA:HB3	1:B:207:PHE:CE1	2.50	0.47
1:B:159:LEU:HD11	1:B:179:ALA:HA	1.96	0.47
1:A:159:LEU:HD11	1:A:179:ALA:HA	1.96	0.47
1:A:238:THR:HG21	1:A:262:LEU:HD12	1.96	0.47
1:A:26:GLU:OE1	1:A:26:GLU:HA	2.14	0.47
1:A:295:LYS:NZ	1:A:385:GLU:OE2	2.48	0.47
1:B:26:GLU:OE1	1:B:26:GLU:HA	2.14	0.47
1:B:60:ILE:O	1:B:64:LEU:HG	2.15	0.47
1:A:60:ILE:O	1:A:64:LEU:HG	2.15	0.46
1:B:238:THR:HG21	1:B:262:LEU:HD12	1.96	0.46
1:B:354:ARG:CG	1:B:360:ILE:HD11	2.41	0.46
1:B:295:LYS:NZ	1:B:385:GLU:OE2	2.48	0.46
1:B:103:ARG:O	1:B:103:ARG:HG3	2.16	0.46
1:A:103:ARG:HG3	1:A:103:ARG:O	2.16	0.46
1:A:4:ARG:HH21	1:A:441:LEU:CG	2.22	0.46
1:B:297:ILE:CG2	1:B:379:TYR:HE1	2.25	0.46
1:B:4:ARG:HH21	1:B:441:LEU:CG	2.22	0.46
1:A:103:ARG:HG2	1:A:103:ARG:NH1	2.31	0.45
1:A:354:ARG:CG	1:A:360:ILE:HD11	2.41	0.45
1:B:103:ARG:NH1	1:B:103:ARG:HG2	2.31	0.45
1:B:395:SER:HA	1:B:419:ILE:HD11	1.98	0.45
1:A:297:ILE:CG2	1:A:379:TYR:HE1	2.25	0.45
1:A:395:SER:HA	1:A:419:ILE:HD11	1.98	0.45
1:B:173:ASN:ND2	1:B:215:ASN:HD21	2.15	0.45
1:A:173:ASN:ND2	1:A:215:ASN:HD21	2.15	0.45
1:B:192:TRP:HB3	1:B:197:LEU:HD11	1.98	0.45
1:B:408:LEU:C	1:B:408:LEU:HD12	2.37	0.45
1:A:257:SER:OG	1:A:403:GLY:HA3	2.17	0.45
1:A:408:LEU:C	1:A:408:LEU:HD12	2.37	0.44
1:A:192:TRP:HB3	1:A:197:LEU:HD11	1.98	0.44
1:A:28:ARG:HH21	1:A:452:GLN:HE21	1.63	0.44
1:A:45:GLN:HG3	1:A:50:THR:OG1	2.18	0.44

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:146:GLN:C	1:B:148:PRO:HD3	2.37	0.44
1:B:337:ASP:O	1:B:340:ILE:HG12	2.16	0.44
1:A:262:LEU:HD21	1:A:296:ASN:OD1	2.18	0.44
1:B:45:GLN:HG3	1:B:50:THR:OG1	2.18	0.44
1:A:146:GLN:C	1:A:148:PRO:HD3	2.37	0.44
1:A:337:ASP:O	1:A:340:ILE:HG12	2.16	0.44
1:B:182:THR:OG1	1:B:184:LEU:HB2	2.17	0.44
1:B:257:SER:OG	1:B:403:GLY:HA3	2.17	0.44
1:B:28:ARG:HH21	1:B:452:GLN:HE21	1.63	0.44
1:A:182:THR:OG1	1:A:184:LEU:HB2	2.18	0.44
1:B:356:MET:CE	1:B:356:MET:HA	2.48	0.44
1:B:113:GLN:O	1:B:117:GLN:HG3	2.16	0.44
1:B:262:LEU:HD21	1:B:296:ASN:OD1	2.18	0.44
1:A:356:MET:CE	1:A:356:MET:HA	2.48	0.44
1:A:113:GLN:O	1:A:117:GLN:HG3	2.16	0.44
1:B:371:ILE:O	1:B:375:LEU:HG	2.18	0.43
1:A:371:ILE:O	1:A:375:LEU:HG	2.18	0.43
1:B:262:LEU:HD22	1:B:294:LEU:HD13	1.99	0.43
1:B:189:SER:O	1:B:190:ASP:HB2	2.18	0.43
1:A:233:VAL:O	1:A:235:SER:N	2.52	0.43
1:A:262:LEU:HD22	1:A:294:LEU:HD13	1.99	0.43
1:B:233:VAL:O	1:B:235:SER:N	2.52	0.43
1:B:328:CYS:HA	1:B:331:ILE:CD1	2.45	0.43
1:B:7:VAL:HG22	1:B:20:LEU:HD23	2.00	0.43
1:B:338:ARG:NH1	1:B:355:GLU:OE1	2.36	0.43
1:A:189:SER:O	1:A:190:ASP:HB2	2.18	0.43
1:A:84:VAL:HG12	1:A:139:GLN:NE2	2.34	0.43
1:B:296:ASN:C	1:B:297:ILE:HD12	2.38	0.43
1:A:296:ASN:C	1:A:297:ILE:HD12	2.38	0.42
1:A:224:GLN:HE21	1:A:224:GLN:HA	1.84	0.42
1:B:224:GLN:HA	1:B:224:GLN:HE21	1.85	0.42
1:A:328:CYS:HA	1:A:331:ILE:CD1	2.45	0.42
1:B:19:MET:HG3	1:B:430:THR:HG21	2.01	0.42
1:A:149:GLU:O	1:A:152:PRO:HD2	2.19	0.42
1:A:7:VAL:HG22	1:A:20:LEU:HD23	2.00	0.42
1:B:149:GLU:O	1:B:152:PRO:HD2	2.19	0.42
1:B:84:VAL:HG12	1:B:139:GLN:NE2	2.34	0.42
1:A:19:MET:HG3	1:A:430:THR:HG21	2.02	0.42
1:A:316:PRO:O	1:A:317:ALA:CB	2.64	0.42
1:B:464:THR:HG23	1:B:465:PRO:HD2	2.02	0.42
1:A:73:ALA:C	1:A:74:ILE:HD13	2.40	0.42

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:73:ALA:C	1:B:74:ILE:HD13	2.40	0.42
1:A:464:THR:HG23	1:A:465:PRO:HD2	2.02	0.42
1:A:238:THR:CG2	1:A:262:LEU:HD12	2.50	0.41
1:B:238:THR:CG2	1:B:262:LEU:HD12	2.50	0.41
1:A:103:ARG:HG2	1:A:103:ARG:HH11	1.86	0.41
1:A:363:SER:OG	1:A:366:GLU:HG3	2.20	0.41
1:B:377:LEU:O	1:B:381:ASP:OD2	2.39	0.41
1:A:377:LEU:O	1:A:381:ASP:OD2	2.39	0.41
1:A:7:VAL:HB	1:A:77:ILE:CD1	2.50	0.41
1:A:163:TYR:O	1:A:166:TYR:HB3	2.21	0.41
1:A:38:PHE:HE2	1:A:56:LEU:HD22	1.86	0.41
1:B:103:ARG:HH11	1:B:103:ARG:HG2	1.86	0.41
1:B:224:GLN:HA	1:B:224:GLN:NE2	2.36	0.41
1:A:400:VAL:HG12	1:A:424:GLY:O	2.21	0.41
1:B:20:LEU:C	1:B:20:LEU:HD13	2.42	0.41
1:B:38:PHE:HE2	1:B:56:LEU:HD22	1.86	0.41
1:A:20:LEU:HD13	1:A:20:LEU:C	2.41	0.41
1:A:224:GLN:NE2	1:A:224:GLN:HA	2.36	0.41
1:A:261:SER:O	1:A:262:LEU:HD23	2.21	0.41
1:A:400:VAL:HB	1:A:428:ALA:HB2	2.03	0.41
1:B:212:HIS:HB3	1:B:213:PRO:HD2	2.03	0.41
1:B:363:SER:OG	1:B:366:GLU:HG3	2.20	0.41
1:A:212:HIS:HB3	1:A:213:PRO:HD2	2.03	0.41
1:A:270:PRO:HD3	1:A:291:TYR:CE2	2.55	0.41
1:B:163:TYR:O	1:B:166:TYR:HB3	2.21	0.41
1:B:17:ARG:HH11	1:B:17:ARG:CG	2.35	0.41
1:A:17:ARG:CG	1:A:17:ARG:HH11	2.35	0.40
1:B:261:SER:O	1:B:262:LEU:HD23	2.21	0.40
1:B:400:VAL:HB	1:B:428:ALA:HB2	2.03	0.40
1:B:7:VAL:HB	1:B:77:ILE:CD1	2.50	0.40
1:A:73:ALA:O	1:A:74:ILE:HD13	2.22	0.40
1:B:73:ALA:O	1:B:74:ILE:HD13	2.22	0.40
1:A:478:ILE:HG13	1:A:479:HIS:ND1	2.36	0.40
1:B:270:PRO:HD3	1:B:291:TYR:CE2	2.55	0.40
1:B:367:LEU:HD23	1:B:367:LEU:HA	1.91	0.40
1:B:400:VAL:HG12	1:B:424:GLY:O	2.21	0.40
1:B:478:ILE:HG13	1:B:479:HIS:ND1	2.36	0.40
1:A:360:ILE:HG22	1:A:362:GLU:HG3	2.03	0.40
1:A:417:CYS:HB2	1:A:419:ILE:HG22	2.03	0.40
1:B:83:GLY:O	1:B:84:VAL:HB	2.22	0.40
1:B:417:CYS:HB2	1:B:419:ILE:HG22	2.03	0.40

All (5) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:62:LEU:CD1	1:B:392:GLU:CA[1_556]	1.87	0.33
1:A:459:ASN:N	1:B:473:HIS:CE1[2_645]	1.93	0.27
1:A:2:THR:CA	1:B:248:ASN:OD1[2_545]	2.06	0.14
1:B:62:LEU:CD1	1:B:393:ASP:N[1_556]	2.07	0.13
1:A:219:HIS:NE2	1:A:311:GLN:OE1[1_455]	2.09	0.11

## 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	477/489 (98%)	450 (94%)	23 (5%)	4 (1%)	19	22
1	B	477/489 (98%)	450 (94%)	23 (5%)	4 (1%)	19	22
All	All	954/978 (98%)	900 (94%)	46 (5%)	8 (1%)	19	22

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	357	ALA
1	B	357	ALA
1	A	225	GLY
1	B	225	GLY
1	A	298	MET
1	B	298	MET
1	A	393	ASP
1	B	393	ASP

### 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	396/406 (98%)	374 (94%)	22 (6%)	21	27
1	B	396/406 (98%)	374 (94%)	22 (6%)	21	27
All	All	792/812 (98%)	748 (94%)	44 (6%)	21	27

All (44) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	17	ARG
1	A	30	LEU
1	A	44	SER
1	A	147	GLN
1	A	204	LYS
1	A	209	ARG
1	A	211	THR
1	A	222	CYS
1	A	243	ILE
1	A	255	LEU
1	A	268	GLN
1	A	290	ARG
1	A	325	LEU
1	A	329	ARG
1	A	356	MET
1	A	388	GLN
1	A	390	ARG
1	A	408	LEU
1	A	427	GLU
1	A	441	LEU
1	A	452	GLN
1	A	479	HIS
1	B	17	ARG
1	B	30	LEU
1	B	44	SER
1	B	147	GLN
1	B	204	LYS
1	B	209	ARG
1	B	211	THR
1	B	222	CYS
1	B	243	ILE

*Continued on next page...*



*Continued from previous page...*

Mol	Chain	Res	Type
1	B	255	LEU
1	B	268	GLN
1	B	290	ARG
1	B	325	LEU
1	B	329	ARG
1	B	356	MET
1	B	388	GLN
1	B	390	ARG
1	B	408	LEU
1	B	427	GLU
1	B	441	LEU
1	B	452	GLN
1	B	479	HIS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (37) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	40	ASN
1	A	43	HIS
1	A	65	ASN
1	A	91	GLN
1	A	94	GLN
1	A	113	GLN
1	A	116	GLN
1	A	130	GLN
1	A	147	GLN
1	A	173	ASN
1	A	224	GLN
1	A	268	GLN
1	A	280	ASN
1	A	311	GLN
1	A	323	GLN
1	A	388	GLN
1	A	411	GLN
1	A	452	GLN
1	B	40	ASN
1	B	43	HIS
1	B	45	GLN
1	B	65	ASN
1	B	91	GLN
1	B	94	GLN
1	B	113	GLN

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type
1	B	116	GLN
1	B	130	GLN
1	B	147	GLN
1	B	173	ASN
1	B	224	GLN
1	B	268	GLN
1	B	280	ASN
1	B	311	GLN
1	B	323	GLN
1	B	388	GLN
1	B	411	GLN
1	B	452	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 [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

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	479/489 (97%)	-0.27	1 (0%) 95 95	9, 31, 56, 86	0
1	B	479/489 (97%)	-0.20	6 (1%) 77 76	9, 31, 56, 86	0
All	All	958/978 (97%)	-0.24	7 (0%) 87 88	9, 31, 56, 86	0

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	480	SER	3.4
1	A	2	THR	2.7
1	B	224	GLN	2.3
1	B	391	GLY	2.3
1	B	2	THR	2.2
1	B	328	CYS	2.2
1	B	393	ASP	2.1

### 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

There are no such residues in this entry.