



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 02:14 PM UTC

PDB ID : 2I5Q / pdb_00002i5q
Title : Crystal structure of Apo L-rhamnonate dehydratase from Escherichia Coli
Authors : Fedorov, A.A.; Fedorov, E.V.; Sauder, J.M.; Burley, S.K.; Gerlt, J.A.; Almo, S.C.; New York SGX Research Center for Structural Genomics (NYSGXRC)
Deposited on : 2006-08-25
Resolution : 2.10 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

<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>.

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

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

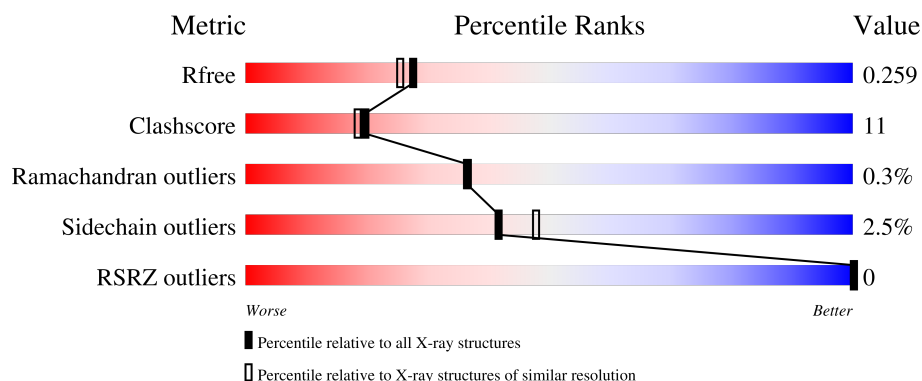
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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}	180053	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

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 ≥ 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	415	 69% 21% • 8%
1	B	415	 71% 20% • 7%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 6039 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-rhamnonate dehydratase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	381	Total	C	N	O	S	0	0	0
			2949	1875	503	545	26			
1	B	384	Total	C	N	O	S	0	0	0
			2980	1895	509	550	26			

There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	MET	-	initiating methionine	UNP Q8XE07
A	0	SER	-	cloning artifact	UNP Q8XE07
A	1	LEU	-	cloning artifact	UNP Q8XE07
A	406	GLU	-	cloning artifact	UNP Q8XE07
A	407	GLY	-	cloning artifact	UNP Q8XE07
A	408	HIS	-	expression tag	UNP Q8XE07
A	409	HIS	-	expression tag	UNP Q8XE07
A	410	HIS	-	expression tag	UNP Q8XE07
A	411	HIS	-	expression tag	UNP Q8XE07
A	412	HIS	-	expression tag	UNP Q8XE07
A	413	HIS	-	expression tag	UNP Q8XE07
B	-1	MET	-	initiating methionine	UNP Q8XE07
B	0	SER	-	cloning artifact	UNP Q8XE07
B	1	LEU	-	cloning artifact	UNP Q8XE07
B	406	GLU	-	cloning artifact	UNP Q8XE07
B	407	GLY	-	cloning artifact	UNP Q8XE07
B	408	HIS	-	expression tag	UNP Q8XE07
B	409	HIS	-	expression tag	UNP Q8XE07
B	410	HIS	-	expression tag	UNP Q8XE07
B	411	HIS	-	expression tag	UNP Q8XE07
B	412	HIS	-	expression tag	UNP Q8XE07
B	413	HIS	-	expression tag	UNP Q8XE07

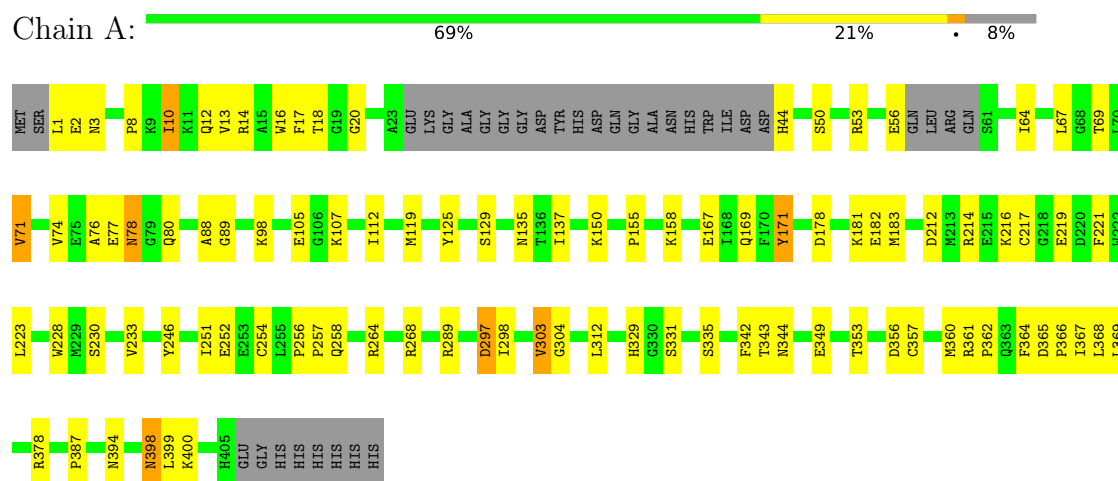
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	61	Total 61	O 61	0	0
2	B	49	Total 49	O 49	0	0

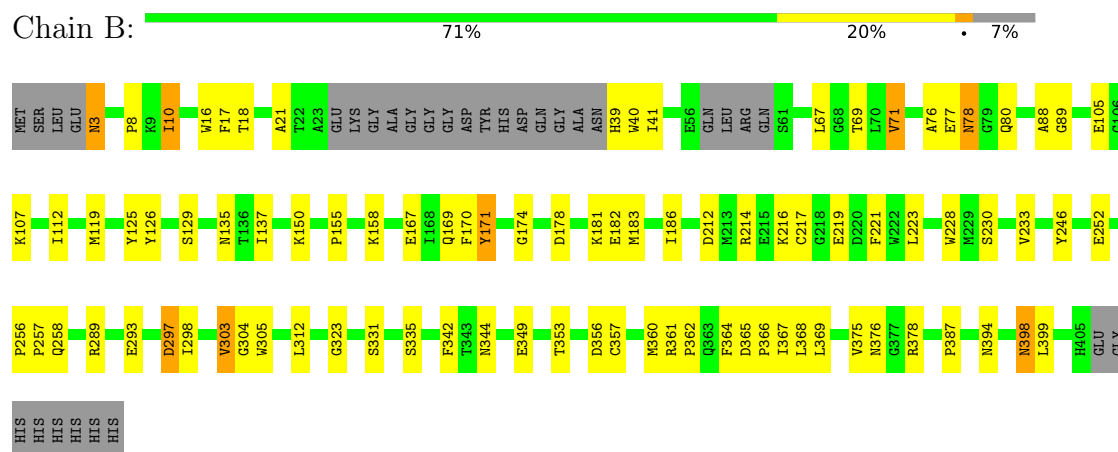
3 Residue-property plots

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: L-rhamnonate dehydratase



• Molecule 1: L-rhamnonate dehydratase



4 Data and refinement statistics

Property	Value	Source
Space group	I 4	Depositor
Cell constants a, b, c, α , β , γ	124.83Å 124.83Å 103.76Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	24.89 – 2.10 24.89 – 2.10	Depositor EDS
% Data completeness (in resolution range)	95.8 (24.89-2.10) 95.9 (24.89-2.10)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.86 (at 2.10Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.229 , 0.259 0.230 , 0.259	Depositor DCC
R_{free} test set	2309 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å ²)	24.0	Xtriage
Anisotropy	0.433	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 20.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.41$, $\langle L^2 \rangle = 0.23$	Xtriage
Estimated twinning fraction	0.237 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	6039	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.71% 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, 0.333 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	0.43	0/3020	0.93	10/4092 (0.2%)
1	B	0.42	0/3054	0.93	8/4140 (0.2%)
All	All	0.43	0/6074	0.93	18/8232 (0.2%)

There are no bond length outliers.

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	171	TYR	N-CA-C	-8.14	97.53	110.14
1	A	171	TYR	N-CA-C	-7.91	97.37	109.95
1	B	252	GLU	N-CA-C	6.58	120.56	109.76
1	A	256	PRO	N-CA-C	-6.58	104.16	110.47
1	A	252	GLU	N-CA-C	6.46	120.35	109.76
1	B	256	PRO	N-CA-C	-6.40	104.33	110.47
1	B	297	ASP	N-CA-C	6.32	118.24	111.36
1	A	297	ASP	N-CA-C	6.31	118.24	111.36
1	B	230	SER	N-CA-C	6.18	120.44	113.02
1	B	219	GLU	N-CA-C	6.14	119.24	111.69
1	A	219	GLU	N-CA-C	6.11	119.20	111.69
1	A	233	VAL	N-CA-C	6.06	116.24	110.42
1	A	343	THR	N-CA-C	-6.01	105.20	112.54
1	B	233	VAL	N-CA-C	5.91	116.09	110.42
1	A	230	SER	N-CA-C	5.68	120.20	113.16
1	B	112	ILE	N-CA-C	5.68	115.87	110.42
1	A	112	ILE	N-CA-C	5.24	115.45	110.42
1	A	329	HIS	N-CA-C	-5.04	103.88	110.53

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	2949	0	2901	74	0
1	B	2980	0	2917	67	0
2	A	61	0	0	1	0
2	B	49	0	0	2	0
All	All	6039	0	5818	135	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 11.

All (135) 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:3:ASN:HD22	1:B:3:ASN:N	1.71	0.87
1:B:303:VAL:HG13	1:B:312:LEU:HD22	1.61	0.82
1:A:303:VAL:HG13	1:A:312:LEU:HD22	1.61	0.81
1:B:214:ARG:HD2	1:B:246:TYR:O	1.82	0.80
1:A:18:THR:HG23	1:A:399:LEU:HD23	1.63	0.80
1:A:214:ARG:HD2	1:A:246:TYR:O	1.83	0.79
1:A:169:GLN:HG2	1:A:378:ARG:HH11	1.49	0.76
1:A:214:ARG:HG2	1:A:223:LEU:HD12	1.68	0.75
1:B:169:GLN:HG2	1:B:378:ARG:HH11	1.50	0.74
1:B:364:PHE:O	1:B:367:ILE:HG22	1.88	0.73
1:A:364:PHE:O	1:A:367:ILE:HG22	1.89	0.73
1:B:16:TRP:HB2	1:B:71:VAL:HG13	1.71	0.72
1:B:214:ARG:HG3	1:B:246:TYR:HB3	1.72	0.72
1:B:214:ARG:HG2	1:B:223:LEU:HD12	1.70	0.71
1:A:214:ARG:HG3	1:A:246:TYR:HB3	1.73	0.69
1:B:367:ILE:HG23	1:B:368:LEU:HD13	1.73	0.69
1:B:39:HIS:ND1	1:B:41:ILE:HG22	2.08	0.69
1:A:129:SER:HB3	1:B:258:GLN:HE21	1.57	0.69
1:A:367:ILE:HG23	1:A:368:LEU:HD13	1.75	0.68
1:A:18:THR:HG22	1:A:398:ASN:O	1.94	0.68
1:A:8:PRO:HB2	1:A:77:GLU:HG2	1.76	0.67
1:B:16:TRP:HB2	1:B:71:VAL:CG1	2.25	0.67
1:A:18:THR:HG22	1:A:398:ASN:C	2.21	0.65

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:258:GLN:HE21	1:B:129:SER:HB3	1.63	0.64
1:B:398:ASN:HD22	1:B:398:ASN:N	1.95	0.62
1:A:18:THR:CG2	1:A:399:LEU:HD23	2.29	0.62
1:A:98:LYS:HB3	1:B:21:ALA:HB1	1.81	0.62
1:A:12:GLN:OE1	1:A:14:ARG:HD2	2.01	0.61
1:A:16:TRP:HB2	1:A:71:VAL:CG1	2.30	0.61
1:A:169:GLN:HG2	1:A:378:ARG:NH1	2.16	0.60
1:A:398:ASN:HD22	1:A:398:ASN:N	1.96	0.60
1:B:360:MET:HA	1:B:360:MET:HE2	1.83	0.60
1:B:3:ASN:N	1:B:3:ASN:ND2	2.45	0.60
1:A:360:MET:HA	1:A:360:MET:HE2	1.84	0.59
1:B:323:GLY:HA2	2:B:448:HOH:O	2.02	0.58
1:B:169:GLN:HG2	1:B:378:ARG:NH1	2.18	0.58
1:A:10:ILE:HD13	1:A:74:VAL:HG13	1.86	0.58
1:B:78:ASN:ND2	1:B:80:GLN:H	2.02	0.57
1:B:331:SER:HB2	1:B:335:SER:OG	2.04	0.57
1:A:16:TRP:HB2	1:A:71:VAL:HG13	1.87	0.57
1:B:10:ILE:HG23	1:B:107:LYS:O	2.05	0.56
1:B:18:THR:OG1	1:B:399:LEU:HD23	2.06	0.56
1:A:50:SER:HA	1:A:56:GLU:HA	1.88	0.56
1:B:40:TRP:CH2	1:B:174:GLY:HA2	2.42	0.55
1:A:78:ASN:ND2	1:A:80:GLN:H	2.04	0.55
1:A:155:PRO:HD2	1:A:158:LYS:HD2	1.90	0.54
1:B:303:VAL:HG22	2:B:430:HOH:O	2.07	0.53
1:B:364:PHE:HB2	1:B:368:LEU:HD22	1.90	0.53
1:A:12:GLN:HG2	1:A:13:VAL:N	2.24	0.53
1:A:78:ASN:HD22	1:A:78:ASN:H	1.57	0.53
1:A:212:ASP:OD1	1:A:216:LYS:HE3	2.09	0.53
1:A:331:SER:HB2	1:A:335:SER:OG	2.09	0.52
1:A:2:GLU:HG3	1:A:3:ASN:N	2.25	0.51
1:B:289:ARG:C	1:B:289:ARG:HD3	2.36	0.51
1:A:364:PHE:HB2	1:A:368:LEU:HD22	1.92	0.51
1:B:10:ILE:O	1:B:10:ILE:HG13	2.10	0.51
1:B:155:PRO:HD2	1:B:158:LYS:HD2	1.92	0.51
1:B:78:ASN:HD22	1:B:78:ASN:H	1.59	0.50
1:A:289:ARG:C	1:A:289:ARG:HD3	2.37	0.50
1:B:178:ASP:O	1:B:182:GLU:HG3	2.12	0.50
1:A:178:ASP:O	1:A:182:GLU:HG3	2.12	0.49
1:A:129:SER:HB3	1:B:258:GLN:NE2	2.25	0.49
1:A:17:PHE:HA	1:A:69:THR:O	2.11	0.49
1:A:181:LYS:HE3	1:A:217:CYS:O	2.13	0.49

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:257:PRO:HG2	1:A:258:GLN:NE2	2.28	0.49
1:B:181:LYS:HE3	1:B:217:CYS:O	2.13	0.49
1:B:17:PHE:HA	1:B:69:THR:O	2.13	0.49
1:A:78:ASN:HD21	1:A:80:GLN:HB2	1.78	0.48
1:B:258:GLN:CD	1:B:258:GLN:H	2.22	0.48
1:B:78:ASN:HD22	1:B:78:ASN:N	2.12	0.48
1:A:78:ASN:HD22	1:A:78:ASN:N	2.12	0.48
1:B:8:PRO:HB2	1:B:77:GLU:HG2	1.96	0.48
1:B:212:ASP:OD1	1:B:216:LYS:HE3	2.12	0.47
1:B:398:ASN:N	1:B:398:ASN:ND2	2.62	0.47
1:A:258:GLN:H	1:A:258:GLN:CD	2.22	0.47
1:B:171:TYR:CZ	1:B:349:GLU:HB2	2.49	0.47
1:A:44:HIS:O	1:A:44:HIS:ND1	2.46	0.47
1:A:171:TYR:CZ	1:A:349:GLU:HB2	2.50	0.47
1:B:257:PRO:HG2	1:B:258:GLN:NE2	2.29	0.47
1:A:12:GLN:CD	1:A:14:ARG:HD2	2.40	0.47
1:B:78:ASN:HD21	1:B:80:GLN:HB2	1.78	0.47
1:B:362:PRO:HB2	1:B:365:ASP:HB2	1.96	0.47
1:B:356:ASP:OD1	1:B:356:ASP:O	2.33	0.47
1:A:214:ARG:NH2	1:A:221:PHE:O	2.39	0.47
1:A:258:GLN:NE2	1:B:129:SER:HB3	2.29	0.47
1:A:362:PRO:HB2	1:A:365:ASP:HB2	1.97	0.46
1:A:297:ASP:C	1:A:298:ILE:HG13	2.41	0.46
1:A:214:ARG:HG2	1:A:223:LEU:CD1	2.44	0.46
1:A:398:ASN:N	1:A:398:ASN:ND2	2.63	0.46
1:A:10:ILE:HA	1:A:76:ALA:HA	1.98	0.45
1:B:135:ASN:HA	1:B:304:GLY:O	2.16	0.45
1:B:367:ILE:HG23	1:B:368:LEU:CD1	2.44	0.45
1:B:214:ARG:NH2	1:B:221:PHE:O	2.38	0.45
1:B:10:ILE:CG1	1:B:105:GLU:HA	2.47	0.45
1:A:10:ILE:HG23	1:A:107:LYS:O	2.17	0.45
1:A:76:ALA:HB3	1:A:78:ASN:ND2	2.31	0.45
1:B:76:ALA:HB3	1:B:78:ASN:ND2	2.32	0.45
1:A:356:ASP:O	1:A:356:ASP:OD1	2.35	0.44
1:A:12:GLN:CG	1:A:13:VAL:N	2.80	0.44
1:B:67:LEU:O	1:B:89:GLY:HA3	2.18	0.44
1:A:20:GLY:HA3	2:A:438:HOH:O	2.17	0.44
1:A:125:TYR:O	1:B:305:TRP:HZ3	2.01	0.44
1:A:135:ASN:HA	1:A:304:GLY:O	2.18	0.44
1:A:356:ASP:O	1:A:357:CYS:HB2	2.17	0.44
1:A:53:ARG:HG3	1:A:53:ARG:HH11	1.82	0.44

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:356:ASP:O	1:B:357:CYS:HB2	2.18	0.44
1:A:10:ILE:HD12	1:A:12:GLN:N	2.33	0.43
1:A:183:MET:CE	1:A:360:MET:HE3	2.48	0.43
1:B:10:ILE:HA	1:B:76:ALA:HA	1.99	0.43
1:B:150:LYS:HA	1:B:387:PRO:HB3	1.99	0.43
1:A:353:THR:OG1	1:A:361:ARG:HD3	2.18	0.43
1:A:167:GLU:OE1	1:A:378:ARG:HD3	2.18	0.43
1:B:119:MET:HE3	1:B:137:ILE:HG23	2.01	0.43
1:B:167:GLU:OE1	1:B:378:ARG:HD3	2.19	0.43
1:A:67:LEU:O	1:A:89:GLY:HA3	2.18	0.43
1:B:297:ASP:C	1:B:298:ILE:HG13	2.43	0.43
1:B:369:LEU:HG	1:B:394:ASN:HB2	2.01	0.43
1:A:10:ILE:HG13	1:A:105:GLU:HA	2.01	0.42
1:A:119:MET:HE3	1:A:137:ILE:HG23	2.00	0.42
1:B:353:THR:OG1	1:B:361:ARG:HD3	2.19	0.42
1:A:150:LYS:HA	1:A:387:PRO:HB3	2.00	0.42
1:B:170:PHE:O	1:B:186:ILE:HG12	2.20	0.42
1:B:375:VAL:O	1:B:376:ASN:HB2	2.20	0.42
1:A:369:LEU:HG	1:A:394:ASN:HB2	2.02	0.42
1:A:367:ILE:HG23	1:A:368:LEU:CD1	2.46	0.42
1:B:183:MET:CE	1:B:360:MET:HE3	2.49	0.42
1:A:251:ILE:HG12	1:A:254:CYS:SG	2.60	0.42
1:B:39:HIS:CE1	1:B:41:ILE:HG22	2.54	0.42
1:B:289:ARG:NE	1:B:293:GLU:OE2	2.43	0.42
1:A:16:TRP:HA	1:A:400:LYS:O	2.20	0.41
1:B:125:TYR:CZ	1:B:126:TYR:HD2	2.39	0.41
1:A:264:ARG:O	1:A:268:ARG:HG3	2.21	0.40
1:A:342:PHE:C	1:A:344:ASN:N	2.76	0.40
1:A:64:ILE:O	1:A:64:ILE:HG13	2.20	0.40
1:B:342:PHE:C	1:B:344:ASN:N	2.77	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	375/415 (90%)	360 (96%)	14 (4%)	1 (0%)	36	36
1	B	378/415 (91%)	362 (96%)	15 (4%)	1 (0%)	36	36
All	All	753/830 (91%)	722 (96%)	29 (4%)	2 (0%)	36	36

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	88	ALA
1	B	88	ALA

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	318/344 (92%)	310 (98%)	8 (2%)	42	48
1	B	321/344 (93%)	313 (98%)	8 (2%)	42	48
All	All	639/688 (93%)	623 (98%)	16 (2%)	42	48

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	LEU
1	A	10	ILE
1	A	71	VAL
1	A	78	ASN
1	A	228	TRP
1	A	303	VAL
1	A	366	PRO
1	A	398	ASN
1	B	3	ASN
1	B	10	ILE
1	B	71	VAL

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	78	ASN
1	B	228	TRP
1	B	303	VAL
1	B	366	PRO
1	B	398	ASN

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

Mol	Chain	Res	Type
1	A	65	ASN
1	A	78	ASN
1	A	121	ASN
1	A	231	GLN
1	A	258	GLN
1	A	269	ASN
1	A	398	ASN
1	B	65	ASN
1	B	78	ASN
1	B	121	ASN
1	B	231	GLN
1	B	258	GLN
1	B	269	ASN
1	B	329	HIS
1	B	398	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 oligosaccharides in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

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 [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, 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	381/415 (91%)	-1.45	0 100 100	12, 26, 49, 69	0
1	B	384/415 (92%)	-1.40	0 100 100	16, 29, 51, 59	0
All	All	765/830 (92%)	-1.43	0 100 100	12, 27, 50, 69	0

There are no RSRZ outliers to report.

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 oligosaccharides 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.