



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 09:54 PM UTC

PDB ID : 4YF9 / pdb_00004yf9
Title : Structure of N-acylhomoserine lactone acylase MacQ
Authors : Yasutake, Y.; Kusada, H.; Kimura, N.
Deposited on : 2015-02-25
Resolution : 2.60 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

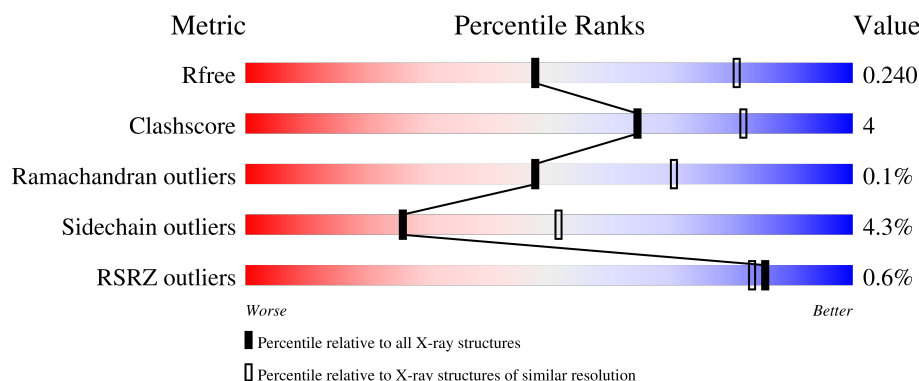
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (2.60-2.60)

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	178	87% 6% • 6%
1	D	178	87% 7% • 6%
1	G	178	% 88% 6% • 6%
1	J	178	% 81% 12% • 6%
2	B	27	4% 52% • 44%

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Mol	Chain	Length	Quality of chain
2	E	27	11% 52% 48%
2	H	27	22% 44% 11% 44%
2	K	27	48% 11% 41%
3	C	581	% 86% 12% ..
3	F	581	87% 11% ..
3	I	581	86% 11% ..
3	L	581	86% 13% ..

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 23165 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 Protein related to penicillin acylase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	167	Total	C	N	O	S	0	0	0
			1289	802	243	238	6			
1	D	168	Total	C	N	O	S	0	0	0
			1295	805	244	240	6			
1	G	167	Total	C	N	O	S	0	0	0
			1289	802	243	238	6			
1	J	167	Total	C	N	O	S	0	0	0
			1289	802	243	238	6			

- Molecule 2 is a protein called Protein related to penicillin acylase.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	B	15	Total	C	N	O	0	0	0
			110	68	22	20			
2	E	14	Total	C	N	O	0	0	0
			106	66	21	19			
2	H	15	Total	C	N	O	0	0	0
			110	68	22	20			
2	K	16	Total	C	N	O	0	0	0
			119	73	23	23			

- Molecule 3 is a protein called Protein related to penicillin acylase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	575	Total	C	N	O	S	0	0	0
			4353	2739	763	832	19			
3	F	575	Total	C	N	O	S	0	0	0
			4353	2739	763	832	19			
3	I	574	Total	C	N	O	S	0	0	0
			4344	2734	762	829	19			
3	L	575	Total	C	N	O	S	0	0	0
			4353	2739	763	832	19			

There are 32 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	574	LEU	-	expression tag	UNP A0A0A1VBK6
C	575	GLU	-	expression tag	UNP A0A0A1VBK6
C	576	HIS	-	expression tag	UNP A0A0A1VBK6
C	577	HIS	-	expression tag	UNP A0A0A1VBK6
C	578	HIS	-	expression tag	UNP A0A0A1VBK6
C	579	HIS	-	expression tag	UNP A0A0A1VBK6
C	580	HIS	-	expression tag	UNP A0A0A1VBK6
C	581	HIS	-	expression tag	UNP A0A0A1VBK6
F	574	LEU	-	expression tag	UNP A0A0A1VBK6
F	575	GLU	-	expression tag	UNP A0A0A1VBK6
F	576	HIS	-	expression tag	UNP A0A0A1VBK6
F	577	HIS	-	expression tag	UNP A0A0A1VBK6
F	578	HIS	-	expression tag	UNP A0A0A1VBK6
F	579	HIS	-	expression tag	UNP A0A0A1VBK6
F	580	HIS	-	expression tag	UNP A0A0A1VBK6
F	581	HIS	-	expression tag	UNP A0A0A1VBK6
I	574	LEU	-	expression tag	UNP A0A0A1VBK6
I	575	GLU	-	expression tag	UNP A0A0A1VBK6
I	576	HIS	-	expression tag	UNP A0A0A1VBK6
I	577	HIS	-	expression tag	UNP A0A0A1VBK6
I	578	HIS	-	expression tag	UNP A0A0A1VBK6
I	579	HIS	-	expression tag	UNP A0A0A1VBK6
I	580	HIS	-	expression tag	UNP A0A0A1VBK6
I	581	HIS	-	expression tag	UNP A0A0A1VBK6
L	574	LEU	-	expression tag	UNP A0A0A1VBK6
L	575	GLU	-	expression tag	UNP A0A0A1VBK6
L	576	HIS	-	expression tag	UNP A0A0A1VBK6
L	577	HIS	-	expression tag	UNP A0A0A1VBK6
L	578	HIS	-	expression tag	UNP A0A0A1VBK6
L	579	HIS	-	expression tag	UNP A0A0A1VBK6
L	580	HIS	-	expression tag	UNP A0A0A1VBK6
L	581	HIS	-	expression tag	UNP A0A0A1VBK6

- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	4	Total O 4 4	0	0
4	C	31	Total O 31 31	0	0
4	D	15	Total O 15 15	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	F	29	Total 29	O 29	0	0
4	G	18	Total 18	O 18	0	0
4	H	1	Total 1	O 1	0	0
4	I	33	Total 33	O 33	0	0
4	J	8	Total 8	O 8	0	0
4	L	16	Total 16	O 16	0	0

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: Protein related to penicillin acylase

Chain A: 



- Molecule 1: Protein related to penicillin acylase

Chain D: 



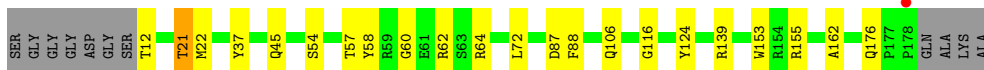
- Molecule 1: Protein related to penicillin acylase

Chain G: 



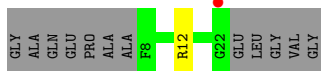
- Molecule 1: Protein related to penicillin acylase

Chain J: 

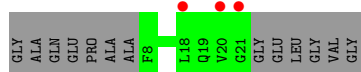


- Molecule 2: Protein related to penicillin acylase

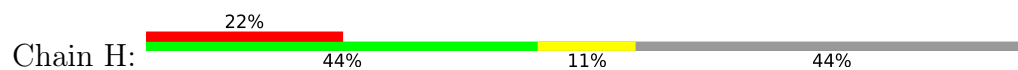
Chain B: 



- Molecule 2: Protein related to penicillin acylase



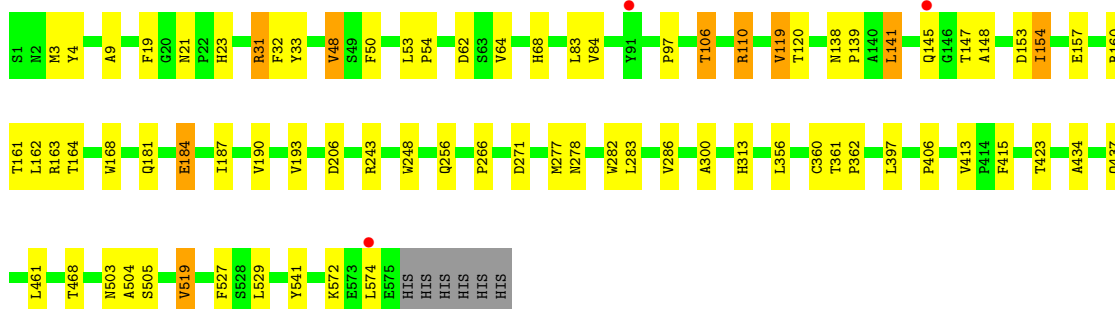
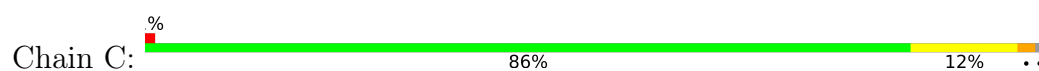
- Molecule 2: Protein related to penicillin acylase



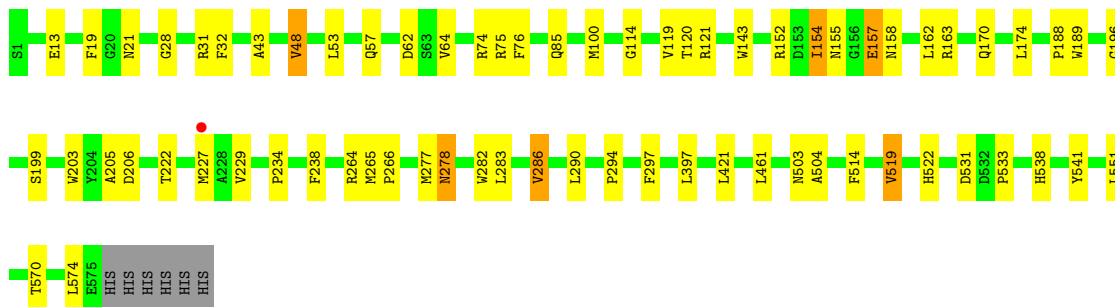
- Molecule 2: Protein related to penicillin acylase



- Molecule 3: Protein related to penicillin acylase

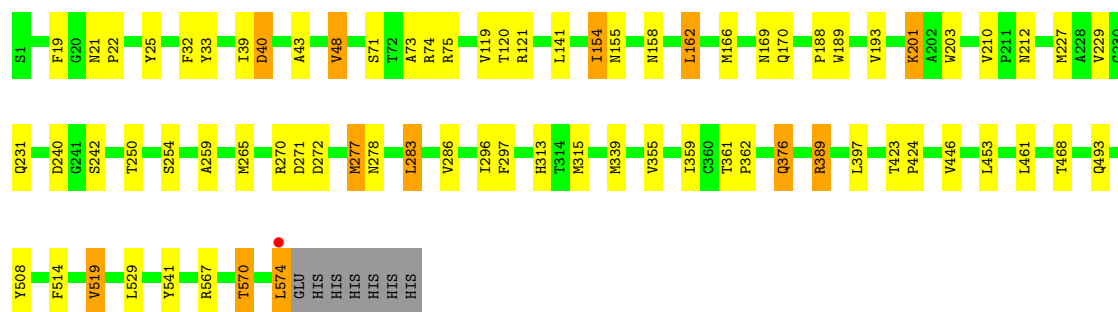


- Molecule 3: Protein related to penicillin acylase



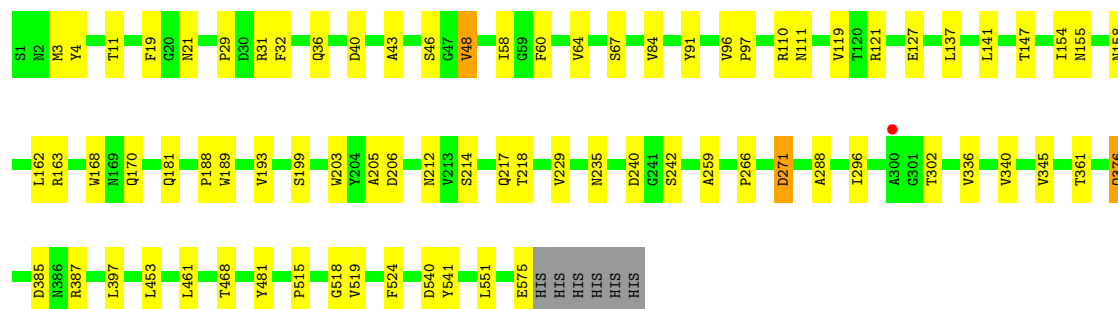
- Molecule 3: Protein related to penicillin acylase

Chain I:  86% 11% ..



• Molecule 3: Protein related to penicillin acylase

Chain L:  86% 13% ..



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	85.62Å 90.11Å 123.17Å 103.49° 104.96° 105.95°	Depositor
Resolution (Å)	38.32 – 2.60 38.32 – 2.60	Depositor EDS
% Data completeness (in resolution range)	98.7 (38.32-2.60) 98.8 (38.32-2.60)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.94 (at 2.61Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.196 , 0.241 0.196 , 0.240	Depositor DCC
R_{free} test set	4943 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å ²)	40.3	Xtriage
Anisotropy	0.026	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 28.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.004 for -h,-k,h+k+l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	23165	wwPDB-VP
Average B, all atoms (Å ²)	42.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.24% 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 [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	0.64	0/1320	0.87	0/1788
1	D	0.65	0/1326	0.85	0/1796
1	G	0.65	0/1320	0.84	0/1788
1	J	0.66	0/1320	0.84	0/1788
2	B	0.54	0/112	0.68	0/150
2	E	0.55	0/108	0.68	0/145
2	H	0.49	0/112	0.66	0/150
2	K	0.57	0/121	0.85	0/162
3	C	0.63	0/4467	0.81	0/6094
3	F	0.63	0/4467	0.81	0/6094
3	I	0.65	0/4458	0.82	0/6082
3	L	0.63	0/4467	0.83	2/6094 (0.0%)
All	All	0.64	0/23598	0.82	2/32131 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	L	288	ALA	CA-C-N	5.30	125.28	120.03
3	L	288	ALA	C-N-CA	5.30	125.28	120.03

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	1289	0	1228	11	0
1	D	1295	0	1233	7	0
1	G	1289	0	1228	7	0
1	J	1289	0	1228	14	0
2	B	110	0	108	1	0
2	E	106	0	105	0	0
2	H	110	0	108	3	0
2	K	119	0	114	2	0
3	C	4353	0	4161	48	0
3	F	4353	0	4161	34	0
3	I	4344	0	4155	44	0
3	L	4353	0	4161	37	0
4	A	4	0	0	0	0
4	C	31	0	0	0	0
4	D	15	0	0	0	0
4	F	29	0	0	0	0
4	G	18	0	0	0	0
4	H	1	0	0	0	0
4	I	33	0	0	0	0
4	J	8	0	0	0	0
4	L	16	0	0	0	0
All	All	23165	0	21990	177	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 4.

All (177) 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)
3:I:376:GLN:HG2	3:I:453:LEU:HD12	1.60	0.83
3:C:32:PHE:HB3	3:C:48:VAL:CG2	2.21	0.70
1:G:106:GLN:HG3	1:G:153:TRP:CH2	2.27	0.69
3:F:74:ARG:HH21	3:F:234:PRO:HG2	1.57	0.69
3:C:145:GLN:H	3:C:145:GLN:CD	2.01	0.68
3:C:32:PHE:HB3	3:C:48:VAL:HG21	1.76	0.67
3:L:32:PHE:HB3	3:L:48:VAL:CG2	2.27	0.64
3:C:19:PHE:CZ	3:C:21:ASN:HB2	2.32	0.64
3:F:19:PHE:CZ	3:F:21:ASN:HB2	2.33	0.62
3:C:164:THR:OG1	3:C:184:GLU:HG2	1.99	0.62
2:B:12:ARG:HD3	3:C:139:PRO:HG2	1.81	0.61
2:H:12:ARG:NH1	1:J:176:GLN:OE1	2.33	0.61
3:C:62:ASP:HA	3:C:519:VAL:HG21	1.82	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:L:271:ASP:OD1	3:L:271:ASP:N	2.33	0.61
3:L:32:PHE:HB3	3:L:48:VAL:HG21	1.83	0.61
3:F:28:GLY:O	3:F:31:ARG:HG2	2.01	0.60
1:J:106:GLN:HG3	1:J:153:TRP:CH2	2.37	0.60
3:I:19:PHE:CZ	3:I:21:ASN:HB2	2.37	0.60
3:I:514:PHE:CE2	3:I:519:VAL:HG13	2.36	0.60
1:J:88:PHE:CZ	3:L:121:ARG:HG3	2.36	0.59
3:I:32:PHE:HB3	3:I:48:VAL:CG2	2.33	0.59
3:F:62:ASP:HA	3:F:519:VAL:HG22	1.83	0.58
1:D:12:THR:HG22	3:F:574:LEU:HD23	1.85	0.58
3:L:376:GLN:HG2	3:L:453:LEU:HD12	1.86	0.58
3:I:32:PHE:HB3	3:I:48:VAL:HG21	1.86	0.57
3:I:33:TYR:CE2	3:I:529:LEU:HD22	2.39	0.57
3:F:277:MET:O	3:F:278:ASN:HB2	2.05	0.57
3:L:168:TRP:HZ2	3:L:181:GLN:HB2	1.69	0.57
3:L:58:ILE:HG23	3:L:67:SER:HA	1.85	0.57
3:F:32:PHE:HB3	3:F:48:VAL:HG22	1.87	0.57
1:J:21:THR:HB	3:L:551:LEU:HD11	1.87	0.55
1:D:106:GLN:HG3	1:D:153:TRP:CH2	2.41	0.55
1:G:12:THR:HA	3:I:574:LEU:HB3	1.88	0.55
3:I:265:MET:HG2	3:I:297:PHE:HZ	1.73	0.54
3:C:83:LEU:HD21	3:C:148:ALA:HB3	1.91	0.53
2:H:18:LEU:HD13	3:I:74:ARG:HH22	1.73	0.53
3:I:270:ARG:HB3	3:I:272:ASP:OD1	2.09	0.53
3:L:19:PHE:CZ	3:L:21:ASN:HB2	2.44	0.52
3:C:282:TRP:HB2	3:C:300:ALA:HA	1.90	0.52
3:L:515:PRO:HD2	3:L:518:GLY:O	2.10	0.52
3:C:168:TRP:CH2	3:C:187:ILE:HD11	2.45	0.52
3:F:53:LEU:HD12	3:F:57:GLN:CD	2.33	0.52
3:F:53:LEU:HD12	3:F:57:GLN:OE1	2.10	0.52
1:A:160:ASN:HD21	3:C:161:THR:HG21	1.75	0.51
1:A:84:ILE:HD11	3:C:119:VAL:CG1	2.41	0.51
3:F:32:PHE:HB3	3:F:48:VAL:CG2	2.40	0.51
3:C:32:PHE:HB3	3:C:48:VAL:HG22	1.93	0.51
3:C:277:MET:O	3:C:278:ASN:HB2	2.11	0.51
3:I:19:PHE:CE2	3:I:21:ASN:HB2	2.45	0.51
3:C:84:VAL:HG21	3:C:97:PRO:HB3	1.92	0.51
3:L:46:SER:HB3	3:L:60:PHE:CZ	2.46	0.51
3:C:9:ALA:HB3	3:C:271:ASP:O	2.12	0.50
3:I:201:LYS:HD2	3:I:271:ASP:HB3	1.93	0.50
1:A:164:GLY:HA3	3:C:153:ASP:OD2	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:62:ASP:HA	3:C:519:VAL:CG2	2.42	0.50
3:F:514:PHE:CE2	3:F:519:VAL:HG13	2.46	0.50
1:G:162:ALA:HB3	3:I:189:TRP:CE2	2.47	0.50
3:L:203:TRP:CZ2	3:L:205:ALA:HB2	2.47	0.50
3:F:100:MET:HE1	3:F:143:TRP:CH2	2.46	0.49
3:I:277:MET:O	3:I:278:ASN:HB2	2.12	0.49
3:L:206:ASP:HB3	3:L:266:PRO:HG2	1.95	0.49
3:L:155:ASN:HA	3:L:158:ASN:HB3	1.94	0.49
3:C:413:VAL:HG22	3:I:231:GLN:NE2	2.27	0.49
3:I:193:VAL:HA	3:I:203:TRP:O	2.13	0.49
3:F:206:ASP:HB3	3:F:266:PRO:HG2	1.95	0.48
1:G:16:GLU:HG3	3:I:570:THR:HG22	1.95	0.48
3:F:533:PRO:HA	3:F:538:HIS:CD2	2.48	0.48
3:C:4:TYR:HB2	3:C:19:PHE:HB3	1.94	0.48
3:L:32:PHE:HB3	3:L:48:VAL:HG22	1.95	0.48
3:I:75:ARG:HD2	3:I:188:PRO:O	2.14	0.48
3:I:240:ASP:OD1	3:I:242:SER:HB2	2.14	0.48
1:A:160:ASN:HD21	3:C:161:THR:CG2	2.27	0.48
3:C:160:ARG:HA	3:C:163:ARG:HD3	1.95	0.48
3:C:154:ILE:HG23	3:C:248:TRP:HZ3	1.78	0.47
3:I:155:ASN:HA	3:I:158:ASN:HB3	1.96	0.47
3:I:210:VAL:O	3:I:259:ALA:HA	2.14	0.47
3:C:110:ARG:HE	3:C:110:ARG:HB2	1.56	0.47
3:C:434:ALA:HA	3:C:437:GLN:OE1	2.13	0.47
3:I:43:ALA:HB2	3:I:170:GLN:HG2	1.95	0.47
3:C:3:MET:HE3	3:C:193:VAL:HG23	1.95	0.47
1:D:32:TRP:CD1	1:D:109:LYS:HD3	2.50	0.47
1:G:88:PHE:CZ	3:I:121:ARG:HG3	2.50	0.47
3:L:193:VAL:HA	3:L:203:TRP:O	2.15	0.47
1:D:101:ARG:NH1	3:F:157:GLU:OE1	2.48	0.47
1:J:22:MET:HA	3:L:540:ASP:OD2	2.14	0.47
3:L:188:PRO:HB2	3:L:189:TRP:CD1	2.50	0.47
3:C:356:LEU:O	3:C:360:CYS:HB2	2.15	0.46
1:G:156:ILE:HG21	3:I:162:LEU:HG	1.98	0.46
3:L:36:GLN:HG3	3:L:46:SER:HB2	1.97	0.46
3:I:22:PRO:HD2	3:I:508:TYR:O	2.16	0.46
3:L:345:VAL:HG22	3:L:481:TYR:CZ	2.51	0.45
3:C:23:HIS:HA	3:C:505:SER:O	2.17	0.45
3:F:188:PRO:HB2	3:F:189:TRP:CD1	2.51	0.45
3:C:157:GLU:HB2	3:C:256:GLN:NE2	2.32	0.45
3:I:389:ARG:HH11	3:I:389:ARG:HG3	1.81	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:K:22:GLY:O	2:K:23:GLU:HB3	2.16	0.45
2:H:21:GLY:HA2	3:I:25:TYR:OH	2.16	0.45
3:I:166:MET:O	3:I:169:ASN:HB2	2.17	0.45
3:I:315:MET:HB3	3:I:339:MET:HE2	1.99	0.45
3:C:154:ILE:HG23	3:C:248:TRP:CZ3	2.52	0.45
3:C:415:PHE:HA	3:C:423:THR:HB	1.98	0.45
1:J:37:TYR:HA	1:J:116:GLY:O	2.17	0.45
3:L:336:VAL:O	3:L:340:VAL:HG13	2.17	0.45
3:I:283:LEU:O	3:I:313:HIS:CE1	2.70	0.44
1:J:58:TYR:CE2	1:J:155:ARG:HG2	2.52	0.44
1:A:107:PRO:HD3	3:C:163:ARG:HG2	1.99	0.44
1:A:72:LEU:HD23	1:A:81:PRO:O	2.18	0.44
1:D:56:LEU:HD11	1:D:143:TRP:CB	2.48	0.44
1:D:88:PHE:CE1	3:F:121:ARG:HG2	2.53	0.44
3:I:250:THR:HG23	3:I:254:SER:HB2	1.99	0.44
3:C:283:LEU:O	3:C:313:HIS:CE1	2.71	0.44
3:F:43:ALA:HB2	3:F:170:GLN:HG2	1.99	0.43
3:F:286:VAL:HG21	3:F:421:LEU:HD11	2.00	0.43
3:C:503:ASN:O	3:C:504:ALA:C	2.59	0.43
3:F:203:TRP:CZ2	3:F:205:ALA:HB2	2.53	0.43
3:C:406:PRO:HB3	3:F:114:GLY:HA3	1.99	0.43
3:C:206:ASP:HB3	3:C:266:PRO:HG2	2.01	0.43
3:I:567:ARG:HH21	3:I:567:ARG:HG3	1.84	0.43
1:J:54:SER:HB3	3:L:29:PRO:HB3	2.00	0.43
1:J:57:THR:OG1	1:J:62:ARG:NH1	2.49	0.43
1:D:47:ASN:HA	3:F:531:ASP:OD1	2.18	0.43
3:I:39:ILE:O	3:I:40:ASP:C	2.60	0.43
3:F:100:MET:HE1	3:F:143:TRP:CZ2	2.54	0.43
3:I:73:ALA:HA	3:I:296:ILE:O	2.19	0.43
3:L:4:TYR:HB2	3:L:19:PHE:HB3	2.00	0.43
3:L:385:ASP:OD2	3:L:387:ARG:NH1	2.50	0.43
3:F:264:ARG:NH1	3:F:294:PRO:HG3	2.33	0.43
1:A:76:SER:C	1:A:78:LEU:H	2.27	0.43
3:I:265:MET:HG2	3:I:297:PHE:CZ	2.53	0.42
3:L:3:MET:HE3	3:L:193:VAL:HG23	2.01	0.42
3:C:33:TYR:CE2	3:C:529:LEU:HD22	2.54	0.42
3:C:138:ASN:HB3	3:C:141:LEU:HD22	2.01	0.42
3:F:155:ASN:HA	3:F:158:ASN:HB3	2.01	0.42
3:F:265:MET:HG2	3:F:297:PHE:HZ	1.84	0.42
1:J:162:ALA:HB3	3:L:189:TRP:CE2	2.54	0.42
3:C:361:THR:HA	3:C:362:PRO:HA	1.85	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:F:76:PHE:HA	3:F:152:ARG:O	2.20	0.42
2:K:16:PRO:HD3	3:L:137:LEU:HD13	2.02	0.42
3:C:53:LEU:HA	3:C:54:PRO:HD3	1.93	0.42
3:C:68:HIS:HB3	3:C:190:VAL:HB	2.01	0.42
1:A:54:SER:O	1:A:57:THR:HB	2.20	0.42
1:A:84:ILE:HD11	3:C:119:VAL:HG13	2.00	0.42
3:C:31:ARG:HG3	3:C:527:PHE:CE1	2.54	0.42
3:C:168:TRP:HZ2	3:C:181:GLN:HB2	1.83	0.42
3:I:361:THR:HA	3:I:362:PRO:HA	1.89	0.42
3:I:423:THR:HA	3:I:424:PRO:C	2.45	0.42
1:A:84:ILE:CD1	3:C:119:VAL:HG11	2.50	0.41
3:F:522:HIS:HA	3:F:551:LEU:O	2.21	0.41
1:G:110:LEU:HG	3:I:166:MET:SD	2.61	0.41
3:F:174:LEU:HB2	3:F:196:GLY:HA3	2.03	0.41
3:C:106:THR:HG23	3:C:120:THR:HB	2.02	0.41
3:I:355:VAL:O	3:I:359:ILE:HG12	2.21	0.41
1:J:45:GLN:HA	1:J:124:TYR:CZ	2.55	0.41
1:J:60:GLY:HA2	1:J:87:ASP:HA	2.02	0.41
3:F:75:ARG:HB3	3:F:154:ILE:HD11	2.02	0.41
3:F:282:TRP:HD1	3:F:290:LEU:O	2.03	0.41
3:L:212:ASN:HB2	3:L:259:ALA:C	2.46	0.41
3:L:214:SER:OG	3:L:217:GLN:HB2	2.20	0.41
3:C:160:ARG:HB3	3:C:184:GLU:HG3	2.02	0.41
3:F:503:ASN:O	3:F:504:ALA:C	2.64	0.41
3:I:212:ASN:HB2	3:I:259:ALA:C	2.45	0.41
1:J:54:SER:CB	3:L:29:PRO:HB3	2.51	0.41
3:L:524:PHE:HA	3:L:541:TYR:CE1	2.56	0.41
3:F:32:PHE:CD1	3:F:48:VAL:HG13	2.56	0.41
3:F:222:THR:HA	3:F:238:PHE:O	2.21	0.41
3:I:154:ILE:H	3:I:154:ILE:HG13	1.73	0.41
1:J:12:THR:N	3:L:575:GLU:OE2	2.54	0.41
3:L:240:ASP:OD1	3:L:242:SER:HB2	2.20	0.41
3:I:32:PHE:CD1	3:I:48:VAL:HG13	2.56	0.41
3:L:43:ALA:HB2	3:L:170:GLN:HG2	2.02	0.40
3:C:50:PHE:HB2	3:C:53:LEU:HD12	2.03	0.40
3:L:91:TYR:OH	3:L:127:GLU:HG2	2.21	0.40
1:A:19:ARG:HD3	1:A:23:GLY:O	2.21	0.40
3:I:32:PHE:HB3	3:I:48:VAL:HG22	2.02	0.40
3:I:446:VAL:HG11	3:I:453:LEU:HD23	2.02	0.40
3:L:96:VAL:HA	3:L:97:PRO:HD3	1.97	0.40
3:L:235:ASN:HB3	3:L:296:ILE:HA	2.02	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	165/178 (93%)	158 (96%)	7 (4%)	0	100	100
1	D	166/178 (93%)	159 (96%)	7 (4%)	0	100	100
1	G	165/178 (93%)	160 (97%)	5 (3%)	0	100	100
1	J	165/178 (93%)	158 (96%)	6 (4%)	1 (1%)	21	42
2	B	13/27 (48%)	12 (92%)	1 (8%)	0	100	100
2	E	12/27 (44%)	11 (92%)	1 (8%)	0	100	100
2	H	13/27 (48%)	11 (85%)	2 (15%)	0	100	100
2	K	14/27 (52%)	14 (100%)	0	0	100	100
3	C	573/581 (99%)	547 (96%)	26 (4%)	0	100	100
3	F	573/581 (99%)	545 (95%)	27 (5%)	1 (0%)	43	66
3	I	572/581 (98%)	545 (95%)	26 (4%)	1 (0%)	43	66
3	L	573/581 (99%)	549 (96%)	23 (4%)	1 (0%)	43	66
All	All	3004/3144 (96%)	2869 (96%)	131 (4%)	4 (0%)	48	70

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	I	40	ASP
3	L	40	ASP
1	J	139	ARG
3	F	278	ASN

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	125/130 (96%)	124 (99%)	1 (1%)	73	88
1	D	126/130 (97%)	122 (97%)	4 (3%)	34	62
1	G	125/130 (96%)	121 (97%)	4 (3%)	34	62
1	J	125/130 (96%)	122 (98%)	3 (2%)	43	70
2	B	11/17 (65%)	11 (100%)	0	100	100
2	E	11/17 (65%)	11 (100%)	0	100	100
2	H	11/17 (65%)	11 (100%)	0	100	100
2	K	12/17 (71%)	12 (100%)	0	100	100
3	C	446/452 (99%)	426 (96%)	20 (4%)	24	50
3	F	446/452 (99%)	426 (96%)	20 (4%)	24	50
3	I	445/452 (98%)	422 (95%)	23 (5%)	21	44
3	L	446/452 (99%)	422 (95%)	24 (5%)	20	42
All	All	2329/2396 (97%)	2230 (96%)	99 (4%)	26	51

All (99) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	72	LEU
3	C	31	ARG
3	C	48	VAL
3	C	64	VAL
3	C	106	THR
3	C	110	ARG
3	C	119	VAL
3	C	141	LEU
3	C	147	THR
3	C	154	ILE
3	C	162	LEU
3	C	184	GLU
3	C	243	ARG
3	C	286	VAL

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Mol	Chain	Res	Type
3	C	397	LEU
3	C	461	LEU
3	C	468	THR
3	C	519	VAL
3	C	541	TYR
3	C	572	LYS
3	C	574	LEU
1	D	12	THR
1	D	64	ARG
1	D	72	LEU
1	D	176	GLN
3	F	13	GLU
3	F	48	VAL
3	F	64	VAL
3	F	85	GLN
3	F	119	VAL
3	F	120	THR
3	F	154	ILE
3	F	157	GLU
3	F	162	LEU
3	F	163	ARG
3	F	199	SER
3	F	227	MET
3	F	229	VAL
3	F	283	LEU
3	F	286	VAL
3	F	397	LEU
3	F	461	LEU
3	F	519	VAL
3	F	541	TYR
3	F	570	THR
1	G	12	THR
1	G	72	LEU
1	G	97	GLU
1	G	176	GLN
3	I	48	VAL
3	I	71	SER
3	I	119	VAL
3	I	120	THR
3	I	141	LEU
3	I	154	ILE
3	I	162	LEU

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Mol	Chain	Res	Type
3	I	201	LYS
3	I	227	MET
3	I	229	VAL
3	I	277	MET
3	I	283	LEU
3	I	286	VAL
3	I	376	GLN
3	I	389	ARG
3	I	397	LEU
3	I	461	LEU
3	I	468	THR
3	I	493	GLN
3	I	519	VAL
3	I	541	TYR
3	I	570	THR
3	I	574	LEU
1	J	21	THR
1	J	64	ARG
1	J	72	LEU
3	L	11	THR
3	L	31	ARG
3	L	48	VAL
3	L	64	VAL
3	L	84	VAL
3	L	110	ARG
3	L	111	ASN
3	L	119	VAL
3	L	141	LEU
3	L	147	THR
3	L	154	ILE
3	L	162	LEU
3	L	163	ARG
3	L	199	SER
3	L	218	THR
3	L	229	VAL
3	L	271	ASP
3	L	302	THR
3	L	361	THR
3	L	376	GLN
3	L	397	LEU
3	L	461	LEU
3	L	468	THR

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Mol	Chain	Res	Type
3	L	519	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	160	ASN
1	A	167	ASN
3	C	21	ASN
3	C	269	GLN
3	C	368	ASN
3	C	563	ASN
1	D	65	HIS
1	D	92	HIS
1	D	145	GLN
3	F	21	ASN
3	F	111	ASN
3	F	448	GLN
3	I	21	ASN
3	I	111	ASN
3	I	231	GLN
3	I	493	GLN
1	J	65	HIS
3	L	111	ASN
3	L	191	ASN
3	L	235	ASN
3	L	305	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 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 ⓘ

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	167/178 (93%)	-0.27	0 100 100	31, 46, 62, 82	0
1	D	168/178 (94%)	-0.50	0 100 100	26, 37, 52, 70	0
1	G	167/178 (93%)	-0.53	1 (0%) 85 83	27, 36, 49, 57	0
1	J	167/178 (93%)	-0.40	1 (0%) 85 83	29, 38, 57, 77	0
2	B	15/27 (55%)	0.86	1 (6%) 24 19	58, 65, 77, 94	0
2	E	14/27 (51%)	1.35	3 (21%) 2 2	65, 79, 82, 85	0
2	H	15/27 (55%)	1.26	6 (40%) 1 0	50, 68, 80, 81	0
2	K	16/27 (59%)	0.52	0 100 100	59, 68, 74, 85	0
3	C	575/581 (98%)	-0.38	3 (0%) 87 85	24, 39, 69, 88	0
3	F	575/581 (98%)	-0.37	1 (0%) 91 90	26, 40, 64, 80	0
3	I	574/581 (98%)	-0.50	1 (0%) 91 90	21, 36, 56, 75	0
3	L	575/581 (98%)	-0.10	1 (0%) 91 90	28, 48, 71, 96	0
All	All	3028/3144 (96%)	-0.33	18 (0%) 85 83	21, 41, 67, 96	0

All (18) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	E	21	GLY	4.6
2	E	20	VAL	3.2
2	H	20	VAL	2.8
2	H	21	GLY	2.7
2	H	17	SER	2.5
3	C	574	LEU	2.5
3	I	574	LEU	2.4
2	H	18	LEU	2.4
3	F	227	MET	2.2
3	C	91	TYR	2.2
1	G	12	THR	2.2

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Mol	Chain	Res	Type	RSRZ
2	H	19	GLN	2.2
2	E	18	LEU	2.1
1	J	178	PRO	2.1
2	B	22	GLY	2.1
3	C	145	GLN	2.0
2	H	22	GLY	2.0
3	L	300	ALA	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 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.