



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 09:21 PM UTC

PDB ID : 4Q2Y / pdb_00004q2y
Title : Crystal structure of Arginyl-tRNA synthetase
Authors : Kim, H.S.; Jo, C.H.; Cha, S.Y.; Han, A.R.; Hwang, K.Y.
Deposited on : 2014-04-10
Resolution : 2.80 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

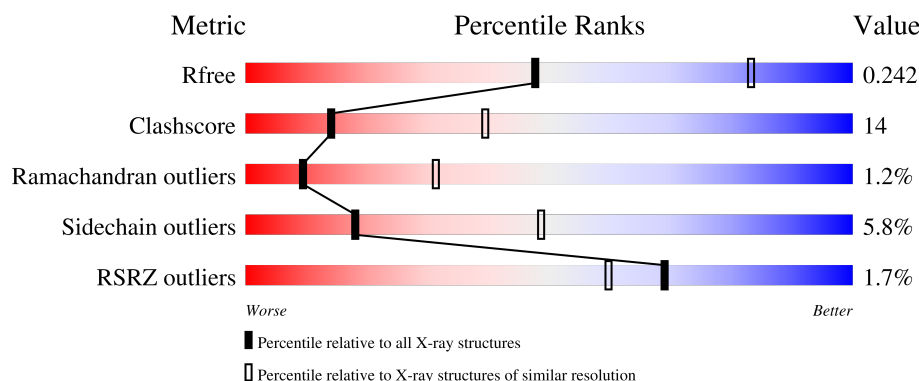
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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	607	
1	B	607	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 9333 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 Arginine-tRNA ligase, cytoplasmic.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	565	Total	C	N	O	S	0	0	0
			4533	2901	768	836	28			
1	B	577	Total	C	N	O	S	0	0	0
			4629	2960	786	856	27			

There are 40 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-18	HIS	-	expression tag	UNP P54136
A	-17	HIS	-	expression tag	UNP P54136
A	-16	HIS	-	expression tag	UNP P54136
A	-15	HIS	-	expression tag	UNP P54136
A	-14	HIS	-	expression tag	UNP P54136
A	-13	HIS	-	expression tag	UNP P54136
A	-12	SER	-	expression tag	UNP P54136
A	-11	SER	-	expression tag	UNP P54136
A	-10	GLY	-	expression tag	UNP P54136
A	-9	LEU	-	expression tag	UNP P54136
A	-8	VAL	-	expression tag	UNP P54136
A	-7	PRO	-	expression tag	UNP P54136
A	-6	ARG	-	expression tag	UNP P54136
A	-5	GLY	-	expression tag	UNP P54136
A	-4	SER	-	expression tag	UNP P54136
A	-3	HIS	-	expression tag	UNP P54136
A	-2	MET	-	expression tag	UNP P54136
A	-1	ALA	-	expression tag	UNP P54136
A	0	SER	-	expression tag	UNP P54136
A	438	ARG	HIS	engineered mutation	UNP P54136
B	-18	HIS	-	expression tag	UNP P54136
B	-17	HIS	-	expression tag	UNP P54136
B	-16	HIS	-	expression tag	UNP P54136
B	-15	HIS	-	expression tag	UNP P54136
B	-14	HIS	-	expression tag	UNP P54136

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
B	-13	HIS	-	expression tag	UNP P54136
B	-12	SER	-	expression tag	UNP P54136
B	-11	SER	-	expression tag	UNP P54136
B	-10	GLY	-	expression tag	UNP P54136
B	-9	LEU	-	expression tag	UNP P54136
B	-8	VAL	-	expression tag	UNP P54136
B	-7	PRO	-	expression tag	UNP P54136
B	-6	ARG	-	expression tag	UNP P54136
B	-5	GLY	-	expression tag	UNP P54136
B	-4	SER	-	expression tag	UNP P54136
B	-3	HIS	-	expression tag	UNP P54136
B	-2	MET	-	expression tag	UNP P54136
B	-1	ALA	-	expression tag	UNP P54136
B	0	SER	-	expression tag	UNP P54136
B	438	ARG	HIS	engineered mutation	UNP P54136

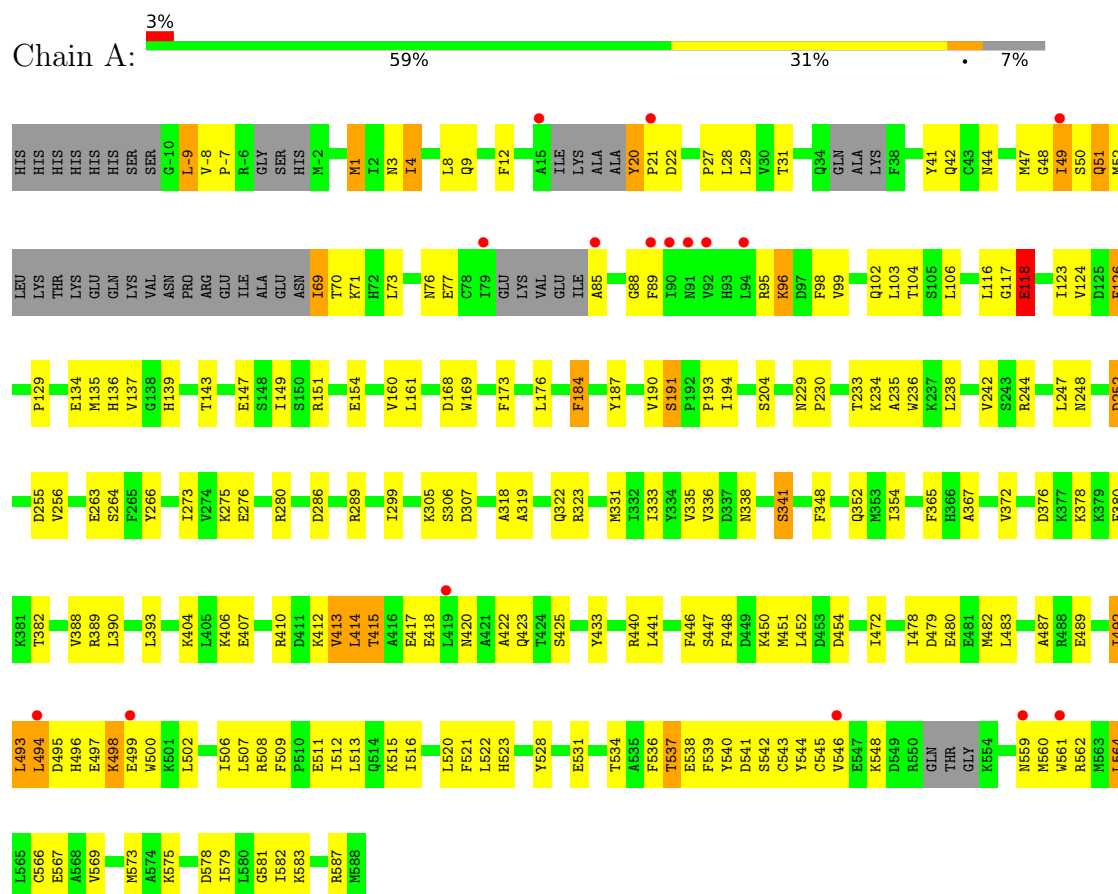
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	84	Total O 84 84	0	0
2	B	87	Total O 87 87	0	0

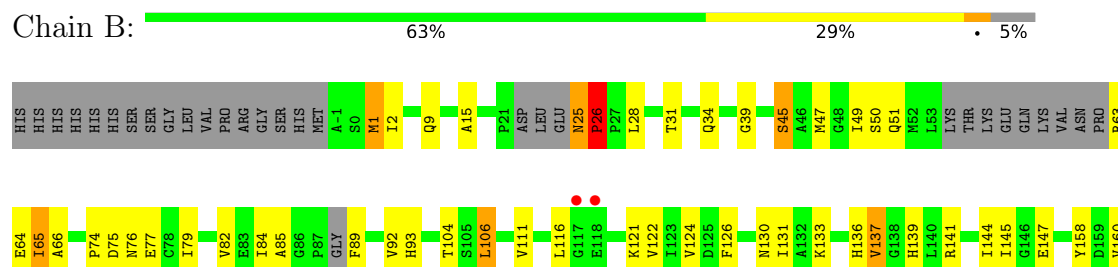
3 Residue-property plots

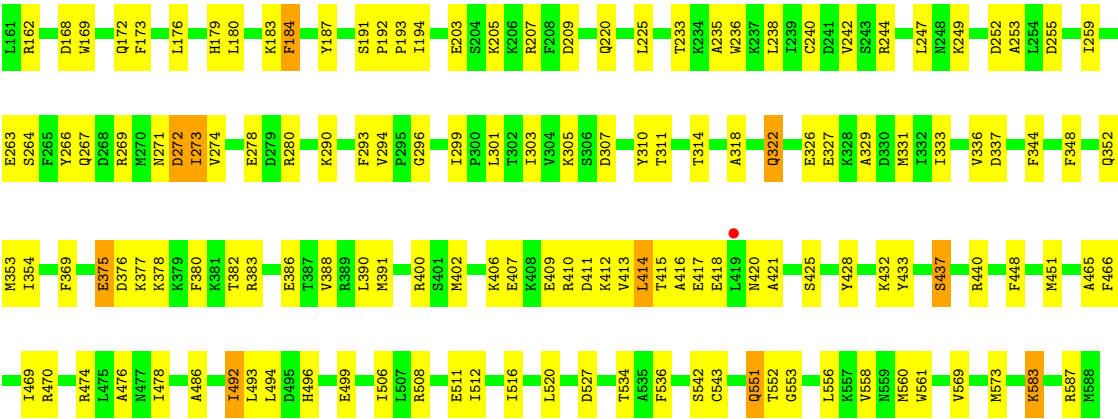
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density.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



• Molecule 1: Arginine-tRNA ligase, cytoplasmic





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	77.62Å 105.23Å 172.26Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	44.90 – 2.80 44.90 – 2.80	Depositor EDS
% Data completeness (in resolution range)	95.5 (44.90-2.80) 95.5 (44.90-2.80)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.00 (at 2.81Å)	Xtriage
Refinement program	PHENIX 1.8.2_1309	Depositor
R, R_{free}	0.208 , 0.268 (Not available) , 0.242	Depositor DCC
R_{free} test set	1712 reflections (4.82%)	wwPDB-VP
Wilson B-factor (Å ²)	44.8	Xtriage
Anisotropy	0.417	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 60.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.27$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	9333	wwPDB-VP
Average B, all atoms (Å ²)	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 30.96 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.1991e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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	0.62	0/4616	1.00	9/6214 (0.1%)
1	B	0.66	1/4714 (0.0%)	1.03	21/6349 (0.3%)
All	All	0.64	1/9330 (0.0%)	1.02	30/12563 (0.2%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	26	PRO	CA-C	5.06	1.56	1.52

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	25	ASN	CA-C-N	14.04	134.84	120.38
1	B	25	ASN	C-N-CA	14.04	134.84	120.38
1	A	118	GLU	N-CA-C	7.68	118.40	108.34
1	B	50	SER	N-CA-C	-7.28	103.55	113.30
1	A	252	ASP	N-CA-C	-7.08	103.49	111.14
1	A	126	PHE	N-CA-C	6.77	118.09	108.74
1	A	567	GLU	N-CA-C	-6.58	102.03	110.19
1	B	273	ILE	N-CA-C	-6.27	104.16	111.00
1	A	184	PHE	CA-C-N	6.23	126.70	119.47
1	A	184	PHE	C-N-CA	6.23	126.70	119.47
1	A	244	ARG	N-CA-C	-5.92	104.83	111.28
1	B	160	VAL	N-CA-C	5.92	116.63	107.99
1	B	26	PRO	CA-C-N	5.78	140.87	127.00
1	B	26	PRO	C-N-CA	5.78	140.87	127.00
1	B	85	ALA	N-CA-C	5.67	117.60	107.80
1	B	26	PRO	N-CA-C	5.65	117.60	110.70
1	B	369	PHE	N-CA-C	5.58	117.89	109.23
1	B	583	LYS	CA-C-N	5.54	125.31	119.76
1	B	583	LYS	C-N-CA	5.54	125.31	119.76
1	B	26	PRO	C-N-CD	-5.43	108.64	120.60
1	B	476	ALA	N-CA-C	-5.43	104.25	112.99

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	191	SER	CA-C-N	-5.36	114.86	120.38
1	B	191	SER	C-N-CA	-5.36	114.86	120.38
1	A	117	GLY	N-CA-C	-5.31	105.90	113.86
1	B	184	PHE	CA-C-N	5.16	125.45	119.47
1	B	184	PHE	C-N-CA	5.16	125.45	119.47
1	B	259	ILE	CB-CA-C	-5.10	105.79	111.35
1	A	335	VAL	N-CA-C	5.09	111.66	106.21
1	B	45	SER	N-CA-C	-5.01	107.13	113.20
1	B	2	ILE	N-CA-C	5.01	115.19	107.77

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	4533	0	4561	144	0
1	B	4629	0	4666	120	0
2	A	84	0	0	11	0
2	B	87	0	0	14	0
All	All	9333	0	9227	262	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 (262) 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:A:578:ASP:OD1	2:A:604:HOH:O	1.91	0.89
1:A:502:LEU:HD12	1:A:536:PHE:HA	1.61	0.83
1:B:440:ARG:NH1	1:B:520:LEU:O	2.11	0.83
1:A:137:VAL:HG12	1:A:372:VAL:HG21	1.58	0.83
1:B:75:ASP:OD2	2:B:623:HOH:O	1.95	0.82
1:A:406:LYS:HD2	1:A:410:ARG:HD3	1.61	0.80
1:B:496:HIS:ND1	1:B:543:CYS:SG	2.54	0.79

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:406:LYS:HD2	1:B:410:ARG:HD3	1.66	0.77
1:B:474:ARG:NH1	2:B:685:HOH:O	2.12	0.77
1:B:130:ASN:HD21	1:B:172:GLN:HE22	1.35	0.75
1:A:20:TYR:OH	1:A:69:ILE:O	2.04	0.75
1:B:179:HIS:ND1	2:B:683:HOH:O	2.21	0.73
1:B:207:ARG:NH1	2:B:683:HOH:O	2.19	0.72
1:A:495:ASP:N	1:A:499:GLU:OE2	2.23	0.71
1:B:205:LYS:NZ	1:B:209:ASP:OD2	2.25	0.70
1:A:508:ARG:NE	2:A:621:HOH:O	2.25	0.70
1:A:499:GLU:HB3	2:A:601:HOH:O	1.92	0.70
1:A:508:ARG:NH2	1:A:531:GLU:OE2	2.23	0.70
1:B:527:ASP:OD2	2:B:620:HOH:O	2.09	0.70
1:B:194:ILE:HD12	1:B:242:VAL:HG21	1.72	0.69
1:A:27:PRO:HG3	1:A:48:GLY:HA3	1.75	0.69
1:A:496:HIS:ND1	1:A:543:CYS:SG	2.66	0.68
1:A:275:LYS:NZ	2:A:660:HOH:O	2.26	0.68
1:A:413:VAL:HG12	1:A:418:GLU:HB2	1.76	0.68
1:A:338:ASN:O	1:A:341:SER:OG	2.11	0.67
1:A:331:MET:HE2	1:A:333:ILE:HD11	1.76	0.67
1:B:496:HIS:HB3	1:B:499:GLU:HG2	1.75	0.67
1:A:29:LEU:HD23	1:A:44:ASN:HB2	1.75	0.67
1:B:122:VAL:HG22	1:B:331:MET:HB3	1.75	0.67
1:B:272:ASP:OD1	1:B:272:ASP:N	2.29	0.66
1:A:478:ILE:HD13	1:A:560:MET:HE1	1.76	0.66
1:A:266:TYR:HE1	1:A:322:GLN:HG2	1.60	0.65
1:B:486:ALA:HB2	1:B:560:MET:HE3	1.79	0.65
1:B:496:HIS:HD1	1:B:543:CYS:HG	1.42	0.65
1:A:96:LYS:HE3	1:A:96:LYS:O	1.95	0.65
1:B:9:GLN:HG2	1:B:28:LEU:HB3	1.80	0.64
1:B:433:TYR:O	1:B:437:SER:OG	2.08	0.64
1:A:448:PHE:HD1	1:A:451:MET:HE2	1.61	0.64
1:A:450:LYS:NZ	2:A:663:HOH:O	2.25	0.64
1:A:506:ILE:HD11	1:A:569:VAL:HG23	1.79	0.64
1:A:149:ILE:HD13	1:A:520:LEU:HD13	1.79	0.64
1:A:42:GLN:NE2	2:A:627:HOH:O	2.30	0.64
1:A:480:GLU:HA	1:A:483:LEU:HD12	1.80	0.64
1:B:414:LEU:HD22	1:B:417:GLU:H	1.64	0.63
1:B:1:MET:HE3	1:B:106:LEU:HD22	1.80	0.63
1:B:252:ASP:O	2:B:650:HOH:O	2.16	0.63
1:B:587:ARG:O	2:B:686:HOH:O	2.15	0.62
1:B:380:PHE:CE2	1:B:388:VAL:HG21	2.34	0.62

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:508:ARG:O	1:B:511:GLU:HG2	1.99	0.62
1:B:409:GLU:O	1:B:413:VAL:HG22	1.99	0.62
1:A:478:ILE:HG21	1:A:560:MET:HE1	1.81	0.61
1:A:194:ILE:HD12	1:A:242:VAL:HG21	1.83	0.60
1:A:404:LYS:HE2	1:A:452:LEU:O	2.02	0.60
1:B:173:PHE:HD1	1:B:176:LEU:HD12	1.67	0.60
1:A:51:GLN:HG2	1:A:52:MET:N	2.17	0.59
1:B:124:VAL:HG12	1:B:126:PHE:HD1	1.66	0.59
1:A:96:LYS:HD2	1:A:500:TRP:HZ3	1.68	0.59
1:B:269:ARG:NH1	1:B:326:GLU:OE2	2.35	0.59
1:B:47:MET:SD	1:B:47:MET:N	2.76	0.58
1:B:512:ILE:O	1:B:516:ILE:HG13	2.03	0.58
1:A:77:GLU:O	1:A:95:ARG:NH2	2.38	0.57
1:A:96:LYS:HD2	1:A:500:TRP:CZ3	2.39	0.57
1:A:20:TYR:HB2	1:A:21:PRO:HD3	1.87	0.57
1:A:124:VAL:HG12	1:A:126:PHE:HD1	1.68	0.57
1:A:234:LYS:HE2	1:A:238:LEU:HD21	1.87	0.57
1:A:498:LYS:HG2	1:A:538:GLU:O	2.05	0.57
1:B:269:ARG:O	1:B:273:ILE:HG12	2.05	0.57
1:B:382:THR:HG22	1:B:383:ARG:H	1.70	0.56
1:B:411:ASP:OD2	2:B:622:HOH:O	2.18	0.56
1:B:249:LYS:HB3	1:B:391:MET:HE2	1.86	0.56
1:A:404:LYS:NZ	1:A:454:ASP:OD1	2.39	0.56
1:A:420:ASN:O	1:A:423:GLN:N	2.39	0.56
1:B:428:TYR:O	1:B:432:LYS:HG3	2.07	0.55
1:B:63:ARG:NH2	2:B:672:HOH:O	2.40	0.54
1:B:121:LYS:NZ	1:B:327:GLU:OE1	2.36	0.54
1:A:69:ILE:O	1:A:71:LYS:N	2.39	0.54
1:B:506:ILE:HD11	1:B:569:VAL:HG23	1.90	0.54
1:A:506:ILE:HG22	1:A:507:LEU:HD12	1.90	0.54
1:B:273:ILE:HG13	1:B:318:ALA:HA	1.90	0.54
1:A:173:PHE:HD1	1:A:176:LEU:HD12	1.74	0.53
1:A:184:PHE:CE1	1:A:193:PRO:HD2	2.42	0.53
1:B:15:ALA:HB2	1:B:74:PRO:HD3	1.90	0.53
1:A:414:LEU:HG	1:A:417:GLU:HB2	1.91	0.53
1:A:169:TRP:CE3	1:A:264:SER:HB3	2.43	0.53
1:A:447:SER:HB3	1:A:450:LYS:HB2	1.90	0.53
1:A:508:ARG:O	1:A:511:GLU:HG2	2.09	0.53
1:B:448:PHE:CD1	1:B:451:MET:HE2	2.43	0.53
1:A:187:TYR:CE2	1:A:235:ALA:HB2	2.44	0.53
1:B:407:GLU:OE1	1:B:407:GLU:N	2.42	0.53

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:124:VAL:HG12	1:A:126:PHE:CD1	2.43	0.53
1:A:425:SER:OG	1:A:583:LYS:O	2.20	0.53
1:B:400:ARG:HD3	2:B:630:HOH:O	2.09	0.52
1:A:47:MET:SD	1:A:47:MET:N	2.82	0.52
1:A:417:GLU:OE2	1:A:587:ARG:HD3	2.09	0.52
1:A:487:ALA:HA	1:A:564:LEU:HD21	1.91	0.52
1:A:561:TRP:CE2	1:A:562:ARG:HG3	2.45	0.52
1:A:234:LYS:O	1:A:238:LEU:HG	2.09	0.52
1:B:34:GLN:H	1:B:34:GLN:CD	2.14	0.52
1:A:561:TRP:NE1	1:A:562:ARG:HG3	2.25	0.52
1:A:27:PRO:HD2	1:A:49:ILE:HG23	1.91	0.52
1:B:448:PHE:HD1	1:B:451:MET:HE2	1.75	0.52
1:A:496:HIS:HB3	1:A:499:GLU:HG2	1.92	0.52
1:A:540:TYR:HE1	1:A:545:CYS:HB3	1.75	0.52
1:A:543:CYS:O	1:A:562:ARG:NH1	2.42	0.52
1:A:447:SER:CB	1:A:450:LYS:HB2	2.41	0.51
1:A:143:THR:HG22	1:A:256:VAL:HG21	1.92	0.51
1:A:441:LEU:HD12	1:B:293:PHE:HE1	1.75	0.51
1:A:380:PHE:CE2	1:A:388:VAL:HG21	2.45	0.51
1:A:559:ASN:HB3	1:A:562:ARG:HB2	1.92	0.51
1:B:465:ALA:O	1:B:469:ILE:HG13	2.11	0.51
1:B:348:PHE:O	1:B:352:GLN:HG3	2.11	0.51
1:A:413:VAL:HG13	1:A:417:GLU:HB3	1.92	0.50
1:B:233:THR:O	1:B:236:TRP:HB3	2.11	0.50
1:A:129:PRO:HA	1:A:139:HIS:CE1	2.46	0.50
1:A:393:LEU:HD11	1:A:446:PHE:CE1	2.47	0.50
1:A:440:ARG:NH1	1:A:520:LEU:O	2.44	0.50
1:A:20:TYR:OH	1:A:69:ILE:HG13	2.12	0.50
1:A:76:ASN:OD1	1:A:77:GLU:N	2.43	0.50
1:A:516:ILE:HD13	1:A:522:LEU:HA	1.93	0.50
1:A:508:ARG:HD2	2:A:621:HOH:O	2.13	0.49
1:A:48:GLY:O	1:A:50:SER:N	2.45	0.49
1:A:184:PHE:O	1:A:187:TYR:HB3	2.13	0.49
1:A:116:LEU:HB2	1:A:118:GLU:O	2.12	0.49
1:A:348:PHE:O	1:A:352:GLN:HG3	2.13	0.49
1:A:489:GLU:OE1	2:A:649:HOH:O	2.19	0.49
1:B:413:VAL:HG23	1:B:418:GLU:HB2	1.95	0.49
1:A:380:PHE:CE1	1:A:446:PHE:HB3	2.49	0.48
1:B:173:PHE:CD1	1:B:176:LEU:HD12	2.47	0.48
1:B:280:ARG:HH11	1:B:354:ILE:HG12	1.78	0.48
1:B:187:TYR:CE2	1:B:235:ALA:HB2	2.48	0.48

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:375:GLU:C	1:B:377:LYS:H	2.22	0.48
1:B:474:ARG:NH2	2:B:626:HOH:O	2.45	0.48
1:A:20:TYR:CD1	1:A:20:TYR:N	2.81	0.48
1:A:492:ILE:HA	1:A:561:TRP:HB2	1.95	0.47
1:B:469:ILE:HD13	1:B:569:VAL:HG11	1.95	0.47
1:B:169:TRP:HB2	1:B:263:GLU:HG3	1.96	0.47
1:B:238:LEU:O	1:B:242:VAL:HG23	2.14	0.47
1:A:1:MET:HE3	1:A:106:LEU:HD12	1.95	0.47
1:A:168:ASP:H	1:A:263:GLU:HG2	1.80	0.47
1:A:498:LYS:NZ	1:A:538:GLU:OE1	2.47	0.47
1:A:407:GLU:N	1:A:407:GLU:OE1	2.47	0.47
1:A:508:ARG:CD	2:A:621:HOH:O	2.61	0.47
1:B:322:GLN:OE1	1:B:327:GLU:HG2	2.15	0.47
1:A:20:TYR:N	1:A:20:TYR:HD1	2.12	0.47
1:A:420:ASN:O	1:A:422:ALA:N	2.48	0.47
1:A:545:CYS:SG	1:A:546:VAL:N	2.87	0.47
1:A:9:GLN:HG2	1:A:28:LEU:HB3	1.97	0.47
1:A:376:ASP:OD1	1:A:378:LYS:HG2	2.14	0.47
1:A:540:TYR:C	1:A:542:SER:HA	2.40	0.47
1:B:77:GLU:H	1:B:77:GLU:CD	2.22	0.47
1:B:240:CYS:HB3	1:B:244:ARG:HH21	1.80	0.47
1:A:440:ARG:HH12	1:A:520:LEU:HB2	1.79	0.46
1:A:286:ASP:OD2	1:A:289:ARG:NH2	2.36	0.46
1:A:509:PHE:CE2	1:A:513:LEU:HD11	2.51	0.46
1:B:64:GLU:O	1:B:66:ALA:N	2.48	0.46
1:A:493:LEU:HG	1:A:561:TRP:CE2	2.50	0.46
1:A:447:SER:OG	1:A:450:LYS:HB2	2.16	0.46
1:B:266:TYR:HE1	1:B:322:GLN:HG2	1.79	0.46
1:A:77:GLU:O	1:A:98:PHE:HB2	2.16	0.46
1:A:512:ILE:HG13	1:A:528:TYR:CD1	2.51	0.46
1:B:294:VAL:HG13	1:B:353:MET:SD	2.56	0.46
1:B:168:ASP:OD2	1:B:264:SER:OG	2.26	0.46
1:B:169:TRP:NE1	1:B:267:GLN:OE1	2.45	0.46
1:B:296:GLY:O	2:B:656:HOH:O	2.21	0.46
1:B:144:ILE:HD13	1:B:432:LYS:HA	1.97	0.45
1:B:331:MET:HE2	1:B:333:ILE:HD11	1.98	0.45
1:B:130:ASN:ND2	1:B:172:GLN:HE22	2.10	0.45
1:B:414:LEU:H	1:B:414:LEU:HD13	1.81	0.45
1:A:135:MET:HE3	1:A:139:HIS:HB3	1.98	0.45
1:B:253:ALA:HB1	1:B:428:TYR:CZ	2.52	0.45
1:B:425:SER:OG	1:B:583:LYS:O	2.19	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:8:LEU:O	1:A:12:PHE:HD1	1.99	0.45
1:A:123:ILE:HG12	1:A:161:LEU:HB3	1.97	0.45
1:A:307:ASP:OD1	1:A:307:ASP:N	2.48	0.45
1:A:276:GLU:OE2	1:A:280:ARG:NH2	2.50	0.45
1:A:433:TYR:HB2	1:A:582:ILE:HD11	1.98	0.45
1:B:274:VAL:HG22	1:B:303:ILE:HG22	1.99	0.45
1:B:556:LEU:HD23	1:B:556:LEU:HA	1.75	0.45
1:B:131:ILE:C	1:B:133:LYS:H	2.25	0.44
1:B:220:GLN:HG2	2:B:602:HOH:O	2.18	0.44
1:A:136:HIS:HD1	1:A:136:HIS:H	1.64	0.44
1:A:273:ILE:HD13	1:A:273:ILE:HA	1.77	0.44
1:A:521:PHE:HB3	1:A:523:HIS:CE1	2.52	0.44
1:A:48:GLY:C	1:A:50:SER:H	2.26	0.44
1:A:136:HIS:HE1	1:A:139:HIS:CD2	2.35	0.44
1:A:548:LYS:HB2	1:A:548:LYS:HE3	1.76	0.44
1:A:99:VAL:O	1:A:103:LEU:HD12	2.18	0.44
1:A:496:HIS:CD2	1:A:497:GLU:H	2.35	0.44
1:A:515:LYS:HA	1:A:515:LYS:HD2	1.74	0.44
1:B:82:VAL:HG12	1:B:92:VAL:HG22	1.99	0.44
1:B:116:LEU:HD11	1:B:158:TYR:CD1	2.53	0.44
1:B:414:LEU:HB2	1:B:415:THR:H	1.61	0.44
1:A:85:ALA:O	1:A:89:PHE:HB2	2.18	0.44
1:A:229:ASN:HB2	2:A:632:HOH:O	2.18	0.44
1:B:106:LEU:HD13	1:B:111:VAL:HG22	2.00	0.44
1:B:303:ILE:O	1:B:314:THR:OG1	2.35	0.43
1:B:271:ASN:OD1	1:B:310:TYR:OH	2.10	0.43
1:B:247:LEU:HD23	1:B:247:LEU:HA	1.67	0.43
1:B:486:ALA:CB	1:B:560:MET:HE3	2.47	0.43
1:A:187:TYR:HA	1:A:191:SER:HB2	2.00	0.43
1:B:147:GLU:OE2	1:B:162:ARG:NH1	2.48	0.43
1:A:147:GLU:HB2	1:A:256:VAL:HG22	2.00	0.43
1:A:273:ILE:HG13	1:A:318:ALA:HA	1.99	0.43
1:A:537:THR:O	1:A:541:ASP:HB2	2.19	0.43
1:A:319:ALA:O	1:A:323:ARG:HG3	2.18	0.43
1:A:154:GLU:HG3	1:A:160:VAL:HG21	2.00	0.43
1:A:233:THR:O	1:A:236:TRP:HB3	2.19	0.43
1:B:376:ASP:C	1:B:378:LYS:H	2.27	0.43
1:B:136:HIS:O	1:B:139:HIS:HB2	2.18	0.42
1:A:472:ILE:HD12	1:A:566:CYS:SG	2.60	0.42
1:B:466:PHE:HB2	1:B:573:MET:HE1	2.01	0.42
1:A:134:GLU:OE2	1:A:389:ARG:HG3	2.19	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:414:LEU:HB2	1:A:415:THR:H	1.51	0.42
1:B:307:ASP:N	1:B:307:ASP:OD1	2.53	0.42
1:B:551:GLN:O	1:B:553:GLY:N	2.52	0.42
1:B:244:ARG:HE	1:B:244:ARG:HB2	1.64	0.42
1:A:280:ARG:HH21	1:A:354:ILE:HG12	1.85	0.42
1:A:479:ASP:OD1	1:A:482:MET:HG3	2.20	0.42
1:B:415:THR:OG1	1:B:416:ALA:N	2.53	0.42
1:A:4:ILE:HD11	1:A:41:TYR:CE2	2.55	0.41
1:A:49:ILE:O	1:A:51:GLN:HA	2.20	0.41
1:A:102:GLN:HB2	1:A:507:LEU:HD21	2.02	0.41
1:A:493:LEU:HG	1:A:561:TRP:NE1	2.35	0.41
1:B:84:ILE:HA	1:B:89:PHE:O	2.20	0.41
1:B:411:ASP:OD1	1:B:412:LYS:N	2.52	0.41
1:A:248:ASN:O	1:A:252:ASP:N	2.51	0.41
1:A:578:ASP:HA	2:A:604:HOH:O	2.21	0.41
1:B:45:SER:O	1:B:49:ILE:HG13	2.20	0.41
1:B:278:GLU:OE2	1:B:290:LYS:NZ	2.27	0.41
1:A:98:PHE:O	1:A:102:GLN:HG2	2.21	0.41
1:A:365:PHE:HE1	1:A:367:ALA:HA	1.84	0.41
1:B:180:LEU:HD23	1:B:235:ALA:HB1	2.03	0.41
1:B:382:THR:HG22	1:B:383:ARG:N	2.33	0.41
1:B:420:ASN:O	1:B:421:ALA:C	2.63	0.41
1:A:151:ARG:NH1	1:A:581:GLY:HA2	2.34	0.41
1:B:137:VAL:HA	1:B:390:LEU:HD21	2.01	0.41
1:B:470:ARG:NH1	2:B:654:HOH:O	2.53	0.41
1:A:390:LEU:HD23	1:A:390:LEU:HA	1.90	0.41
1:B:39:GLY:HA2	1:B:93:HIS:CD2	2.56	0.41
1:B:336:VAL:HG11	1:B:344:PHE:CE2	2.56	0.41
1:A:136:HIS:CE1	1:A:139:HIS:CG	3.09	0.41
1:A:573:MET:HE2	1:A:573:MET:HB3	1.78	0.41
1:B:192:PRO:HA	1:B:193:PRO:HD3	1.90	0.41
1:B:25:ASN:HA	1:B:26:PRO:HD2	1.62	0.41
1:B:124:VAL:HG12	1:B:126:PHE:CD1	2.52	0.41
1:B:402:MET:SD	1:B:418:GLU:OE2	2.80	0.41
1:B:413:VAL:CG2	1:B:418:GLU:HB2	2.50	0.41
1:A:494:LEU:HA	1:A:494:LEU:HD22	1.85	0.40
1:B:121:LYS:HB3	1:B:329:ALA:HA	2.03	0.40
1:B:469:ILE:HG12	1:B:536:PHE:CE2	2.57	0.40
1:B:492:ILE:HA	1:B:561:TRP:HB2	2.02	0.40
1:A:-8:VAL:HG13	1:A:-7:PRO:HD2	2.02	0.40
1:B:76:ASN:OD1	1:B:79:ILE:HB	2.21	0.40

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:141:ARG:O	1:B:145:ILE:HG13	2.21	0.40
1:A:441:LEU:HD12	1:B:293:PHE:CE1	2.55	0.40
1:A:575:LYS:O	1:A:579:ILE:HG13	2.22	0.40
1:B:116:LEU:HD21	1:B:158:TYR:CE1	2.56	0.40
1:B:183:LYS:NZ	1:B:203:GLU:OE1	2.52	0.40
1:A:137:VAL:HG11	1:A:372:VAL:HG11	2.03	0.40
1:B:169:TRP:CH2	1:B:225:LEU:HG	2.56	0.40
1:B:180:LEU:HD12	1:B:184:PHE:HD2	1.86	0.40
1:B:169:TRP:CB	1:B:263:GLU:HG3	2.52	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	551/607 (91%)	490 (89%)	51 (9%)	10 (2%)	6	23
1	B	569/607 (94%)	520 (91%)	45 (8%)	4 (1%)	18	47
All	All	1120/1214 (92%)	1010 (90%)	96 (9%)	14 (1%)	9	31

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	49	ILE
1	A	70	THR
1	B	26	PRO
1	B	65	ILE
1	A	-9	LEU
1	B	552	THR
1	A	22	ASP
1	A	191	SER
1	A	544	TYR

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	542	SER
1	A	88	GLY
1	A	190	VAL
1	A	415	THR
1	A	230	PRO

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	494/530 (93%)	462 (94%)	32 (6%)	15	43
1	B	503/530 (95%)	477 (95%)	26 (5%)	21	53
All	All	997/1060 (94%)	939 (94%)	58 (6%)	18	49

All (58) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	-9	LEU
1	A	1	MET
1	A	3	ASN
1	A	4	ILE
1	A	20	TYR
1	A	31	THR
1	A	51	GLN
1	A	69	ILE
1	A	73	LEU
1	A	96	LYS
1	A	104	THR
1	A	118	GLU
1	A	204	SER
1	A	247	LEU
1	A	255	ASP
1	A	299	ILE
1	A	305	LYS
1	A	306	SER

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	336	VAL
1	A	341	SER
1	A	382	THR
1	A	412	LYS
1	A	413	VAL
1	A	414	LEU
1	A	492	ILE
1	A	493	LEU
1	A	494	LEU
1	A	498	LYS
1	A	534	THR
1	A	537	THR
1	A	539	PHE
1	A	564	LEU
1	B	1	MET
1	B	31	THR
1	B	51	GLN
1	B	65	ILE
1	B	104	THR
1	B	106	LEU
1	B	137	VAL
1	B	255	ASP
1	B	272	ASP
1	B	299	ILE
1	B	301	LEU
1	B	305	LYS
1	B	311	THR
1	B	322	GLN
1	B	337	ASP
1	B	375	GLU
1	B	386	GLU
1	B	414	LEU
1	B	437	SER
1	B	478	ILE
1	B	492	ILE
1	B	493	LEU
1	B	494	LEU
1	B	534	THR
1	B	551	GLN
1	B	558	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	93	HIS
1	A	165	HIS
1	A	172	GLN
1	A	559	ASN
1	B	172	GLN
1	B	284	GLN
1	B	340	GLN
1	B	366	HIS
1	B	514	GLN
1	B	559	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 [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	565/607 (93%)	0.09	16 (2%) 55 45	18, 52, 86, 113	0
1	B	577/607 (95%)	-0.31	3 (0%) 87 82	15, 39, 74, 89	0
All	All	1142/1214 (94%)	-0.11	19 (1%) 69 60	15, 44, 81, 113	0

All (19) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	117	GLY	5.1
1	A	94	LEU	4.6
1	A	15	ALA	3.9
1	A	419	LEU	3.8
1	A	494	LEU	3.4
1	A	21	PRO	3.4
1	A	92	VAL	3.2
1	B	419	LEU	3.2
1	A	85	ALA	3.0
1	A	89	PHE	3.0
1	A	561	TRP	2.8
1	A	79	ILE	2.8
1	A	499	GLU	2.5
1	A	49	ILE	2.4
1	B	118	GLU	2.3
1	A	546	VAL	2.3
1	A	90	ILE	2.3
1	A	91	ASN	2.0
1	A	559	ASN	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.