



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 10:12 PM UTC

PDB ID : 6WQI / pdb_00006wqi
Title : Xanthomonas citri Methionyl-tRNA synthetase (apo)
Authors : Mercaldi, G.F.; Benedetti, C.E.
Deposited on : 2020-04-28
Resolution : 2.00 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

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

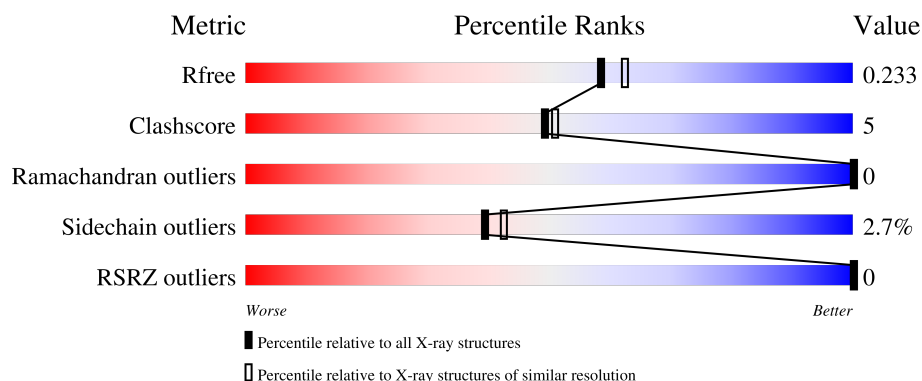
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 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	711	
1	B	711	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 8988 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 Methionine-tRNA ligase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	547	Total	C	N	O	S	0	0	0
			4244	2713	730	782	19			
1	B	547	Total	C	N	O	S	0	0	0
			4244	2713	730	782	19			

There are 34 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-16	MET	-	initiating methionine	UNP Q8PMP0
A	-15	GLY	-	expression tag	UNP Q8PMP0
A	-14	HIS	-	expression tag	UNP Q8PMP0
A	-13	HIS	-	expression tag	UNP Q8PMP0
A	-12	HIS	-	expression tag	UNP Q8PMP0
A	-11	HIS	-	expression tag	UNP Q8PMP0
A	-10	HIS	-	expression tag	UNP Q8PMP0
A	-9	HIS	-	expression tag	UNP Q8PMP0
A	-8	GLY	-	expression tag	UNP Q8PMP0
A	-7	GLU	-	expression tag	UNP Q8PMP0
A	-6	ASN	-	expression tag	UNP Q8PMP0
A	-5	LEU	-	expression tag	UNP Q8PMP0
A	-4	TYR	-	expression tag	UNP Q8PMP0
A	-3	PHE	-	expression tag	UNP Q8PMP0
A	-2	GLN	-	expression tag	UNP Q8PMP0
A	-1	GLY	-	expression tag	UNP Q8PMP0
A	0	SER	-	expression tag	UNP Q8PMP0
B	-16	MET	-	initiating methionine	UNP Q8PMP0
B	-15	GLY	-	expression tag	UNP Q8PMP0
B	-14	HIS	-	expression tag	UNP Q8PMP0
B	-13	HIS	-	expression tag	UNP Q8PMP0
B	-12	HIS	-	expression tag	UNP Q8PMP0
B	-11	HIS	-	expression tag	UNP Q8PMP0
B	-10	HIS	-	expression tag	UNP Q8PMP0
B	-9	HIS	-	expression tag	UNP Q8PMP0

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
B	-8	GLY	-	expression tag	UNP Q8PMP0
B	-7	GLU	-	expression tag	UNP Q8PMP0
B	-6	ASN	-	expression tag	UNP Q8PMP0
B	-5	LEU	-	expression tag	UNP Q8PMP0
B	-4	TYR	-	expression tag	UNP Q8PMP0
B	-3	PHE	-	expression tag	UNP Q8PMP0
B	-2	GLN	-	expression tag	UNP Q8PMP0
B	-1	GLY	-	expression tag	UNP Q8PMP0
B	0	SER	-	expression tag	UNP Q8PMP0

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Zn	0	0
			1	1		
2	B	1	Total	Zn	0	0
			1	1		

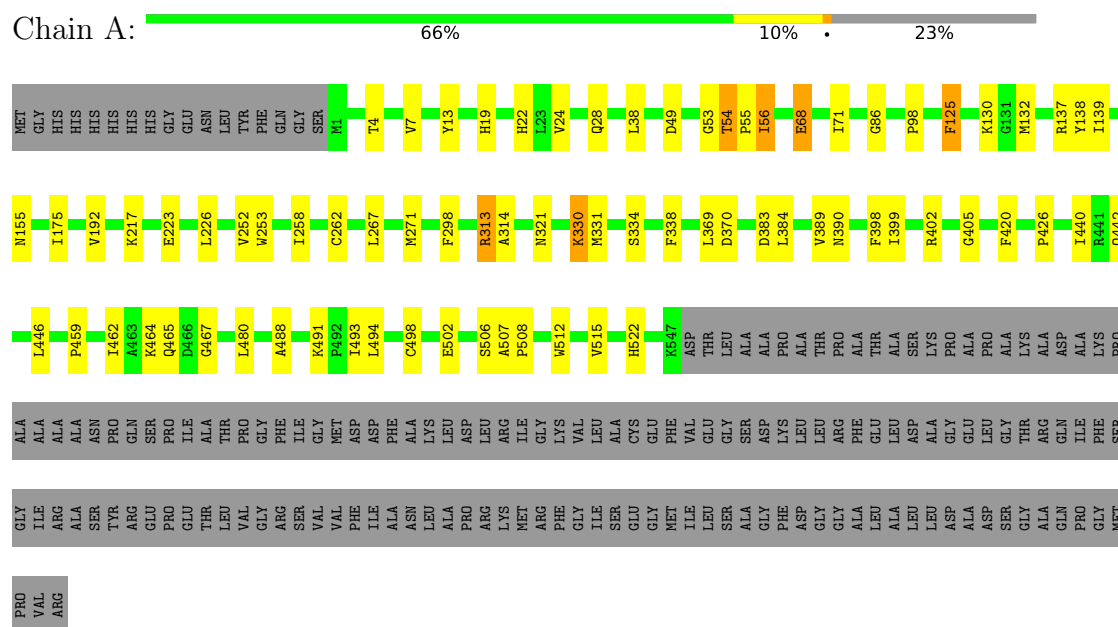
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	276	Total	O	0	0
			276	276		
3	B	222	Total	O	0	0
			222	222		

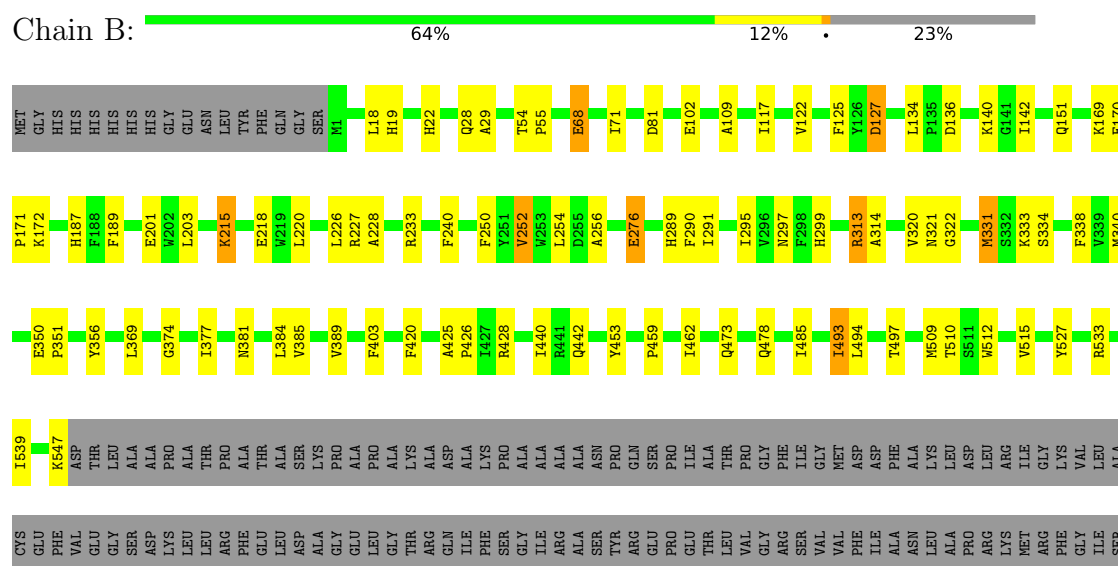
3 Residue-property plots

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: Methionine-tRNA ligase



• Molecule 1: Methionine-tRNA ligase



GLU	GLY	MET	ILE	LEU	SER	ALA	GLY	PHE	ASP	GLY	GLY	ALA	LEU	ALA	LEU	LEU	ASP	ASP	ALA	ASP	SER	GLY	ALA	GLN	PRO	GLY	MET	PRO	VAL	ARG
-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	74.02Å 95.25Å 84.08Å 90.00° 90.52° 90.00°	Depositor
Resolution (Å)	47.83 – 2.00 47.83 – 2.00	Depositor EDS
% Data completeness (in resolution range)	94.3 (47.83-2.00) 94.3 (47.83-2.00)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.44 (at 2.00Å)	Xtriage
Refinement program	REFMAC 5.8.0258	Depositor
R, R_{free}	0.185 , 0.230 0.190 , 0.233	Depositor DCC
R_{free} test set	3766 reflections (4.78%)	wwPDB-VP
Wilson B-factor (Å ²)	26.8	Xtriage
Anisotropy	0.885	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 29.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	0.340 for h,-k,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	8988	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.48% 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](#)

Bond lengths and bond angles in the following residue types are not validated in this section: ZN

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	1.11	5/4356 (0.1%)	1.43	10/5921 (0.2%)
1	B	1.14	5/4356 (0.1%)	1.49	11/5921 (0.2%)
All	All	1.12	10/8712 (0.1%)	1.46	21/11842 (0.2%)

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	7	VAL	C-O	6.49	1.31	1.24
1	A	488	ALA	C-O	6.39	1.31	1.24
1	B	297	ASN	C-O	6.24	1.31	1.24
1	A	53	GLY	C-O	6.14	1.28	1.23
1	A	98	PRO	C-O	-6.02	1.16	1.24
1	B	289	HIS	C-O	6.01	1.30	1.24
1	B	127	ASP	C-O	5.63	1.30	1.23
1	A	498	CYS	C-O	-5.30	1.17	1.24
1	B	322	GLY	C-O	5.12	1.28	1.23
1	B	240	PHE	C-O	5.00	1.30	1.24

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	374	GLY	CA-C-N	6.40	128.86	120.28
1	B	374	GLY	C-N-CA	6.40	128.86	120.28
1	A	389	VAL	CA-C-N	6.14	128.79	120.38
1	A	389	VAL	C-N-CA	6.14	128.79	120.38
1	A	4	THR	CA-CB-OG1	-6.07	100.49	109.60
1	B	127	ASP	O-C-N	5.89	130.17	122.68
1	B	170	GLU	CA-C-N	5.88	125.81	119.76
1	B	170	GLU	C-N-CA	5.88	125.81	119.76
1	A	465	GLN	CA-C-N	5.59	127.71	120.44

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	465	GLN	C-N-CA	5.59	127.71	120.44
1	B	331	MET	CA-C-N	5.53	130.80	123.00
1	B	331	MET	C-N-CA	5.53	130.80	123.00
1	B	510	THR	CA-CB-OG1	-5.35	101.57	109.60
1	A	383	ASP	CA-CB-CG	5.34	117.94	112.60
1	A	258	ILE	CA-C-N	5.33	125.90	119.98
1	A	258	ILE	C-N-CA	5.33	125.90	119.98
1	B	493	ILE	CA-C-O	-5.33	115.01	120.71
1	B	109	ALA	CA-C-O	-5.21	115.35	120.82
1	B	136	ASP	CA-CB-CG	5.18	117.78	112.60
1	A	480	LEU	CA-C-N	5.10	127.54	120.29
1	A	480	LEU	C-N-CA	5.10	127.54	120.29

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	4244	0	4130	43	0
1	B	4244	0	4130	48	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	276	0	0	5	0
3	B	222	0	0	1	0
All	All	8988	0	8260	91	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 5.

All (91) 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:331:MET:HE2	1:B:338:PHE:HA	1.68	0.76
1:A:19:HIS:H	1:A:22:HIS:HD2	1.38	0.70

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:331:MET:HE2	1:A:338:PHE:HA	1.78	0.65
1:B:19:HIS:H	1:B:22:HIS:HD2	1.42	0.65
1:B:377:ILE:HD11	1:B:539:ILE:HG13	1.81	0.63
1:A:369:LEU:HD23	1:A:369:LEU:C	2.28	0.58
1:A:19:HIS:H	1:A:22:HIS:CD2	2.22	0.57
1:A:13:TYR:CZ	1:A:56:ILE:HD12	2.39	0.57
1:A:267:LEU:HG	1:A:271:MET:HE2	1.86	0.56
1:B:203:LEU:HD12	1:B:220:LEU:HD11	1.87	0.55
1:B:291:ILE:O	1:B:320:VAL:HA	2.06	0.55
1:B:389:VAL:HG13	1:B:527:TYR:OH	2.08	0.54
1:A:426:PRO:HB2	1:A:442:GLN:HE22	1.71	0.54
1:B:389:VAL:CG1	1:B:527:TYR:OH	2.56	0.54
1:B:428:ARG:HD3	1:B:512:TRP:CD1	2.43	0.53
1:B:19:HIS:HA	1:B:340:MET:HA	1.91	0.53
1:A:369:LEU:HD23	1:A:370:ASP:N	2.24	0.53
1:A:28:GLN:HE22	1:A:321:ASN:HB3	1.74	0.53
1:B:440:ILE:HD11	1:B:493:ILE:HD13	1.91	0.52
1:A:192:VAL:HG12	1:A:226:LEU:HD13	1.93	0.51
1:B:381:ASN:O	1:B:385:VAL:HB	2.10	0.51
1:A:155:ASN:ND2	3:A:808:HOH:O	2.40	0.51
1:A:384:LEU:HD21	1:A:494:LEU:HD11	1.91	0.51
1:B:220:LEU:HD23	1:B:226:LEU:HD11	1.92	0.51
1:A:390:ASN:HA	3:A:902:HOH:O	2.11	0.50
1:A:420:PHE:CE2	1:A:446:LEU:HB3	2.47	0.49
1:B:276:GLU:H	1:B:276:GLU:CD	2.19	0.49
1:A:426:PRO:HB2	1:A:442:GLN:NE2	2.28	0.49
1:B:68:GLU:HA	1:B:71:ILE:HG12	1.94	0.49
1:A:330:LYS:HE3	3:A:1032:HOH:O	2.13	0.48
1:B:189:PHE:HB3	1:B:228:ALA:HB1	1.95	0.48
1:B:142:ILE:HB	1:B:169:LYS:HB2	1.94	0.48
1:A:54:THR:N	1:A:55:PRO:CD	2.77	0.48
1:B:122:VAL:CG2	1:B:187:HIS:CD2	2.96	0.47
1:A:506:SER:O	1:A:522:HIS:CE1	2.67	0.47
1:B:127:ASP:HB2	1:B:134:LEU:HD21	1.96	0.47
1:A:506:SER:O	1:A:522:HIS:HE1	1.98	0.47
1:B:29:ALA:HB2	1:B:290:PHE:CG	2.50	0.47
1:A:467:GLY:HA2	3:A:921:HOH:O	2.14	0.46
1:A:54:THR:CG2	1:A:132:MET:HE1	2.45	0.46
1:A:253:TRP:O	1:A:298:PHE:HZ	1.98	0.46
1:B:385:VAL:HG11	1:B:533:ARG:HG3	1.97	0.46
1:A:130:LYS:HD3	1:A:175:ILE:HG21	1.97	0.46

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:313:ARG:HD3	1:A:314:ALA:O	2.15	0.46
1:A:130:LYS:HD3	1:A:175:ILE:CG2	2.46	0.46
1:B:28:GLN:NE2	3:B:812:HOH:O	2.49	0.46
1:B:117:ILE:HA	1:B:189:PHE:O	2.16	0.46
1:B:313:ARG:HD3	1:B:314:ALA:O	2.17	0.45
1:B:18:LEU:HD12	1:B:81:ASP:HB2	1.99	0.45
1:B:512:TRP:O	1:B:515:VAL:HG22	2.16	0.45
1:A:137:ARG:HD2	1:A:138:TYR:CE2	2.52	0.45
1:B:28:GLN:HE22	1:B:321:ASN:HB3	1.82	0.45
1:A:125:PHE:HD2	1:A:139:ILE:HD13	1.81	0.44
1:B:295:ILE:HD12	1:B:299:HIS:CE1	2.52	0.44
1:B:459:PRO:HA	1:B:462:ILE:HD12	1.99	0.44
1:A:223:GLU:O	1:A:223:GLU:HG3	2.16	0.44
1:A:24:VAL:O	1:A:28:GLN:HG3	2.19	0.43
1:B:453:TYR:CZ	1:B:478:GLN:HG2	2.54	0.43
1:B:54:THR:N	1:B:55:PRO:CD	2.82	0.43
1:A:459:PRO:HA	1:A:462:ILE:HD12	2.01	0.43
1:B:494:LEU:HB3	1:B:497:THR:HB	2.01	0.43
1:B:356:TYR:CE1	1:B:384:LEU:HD21	2.54	0.42
1:A:440:ILE:HD11	1:A:493:ILE:HD13	2.01	0.42
1:B:215:LYS:HE3	1:B:218:GLU:OE1	2.19	0.42
1:B:356:TYR:CD1	1:B:384:LEU:HD21	2.54	0.42
1:A:398:PHE:O	1:A:402:ARG:HB2	2.19	0.42
1:A:399:ILE:HG22	1:A:405:GLY:HA2	2.01	0.42
1:A:384:LEU:HD23	1:A:384:LEU:HA	1.82	0.42
1:A:491:LYS:HE2	1:A:502:GLU:OE2	2.20	0.42
1:B:171:PRO:O	1:B:172:LYS:HD3	2.19	0.41
1:B:509:MET:HE3	1:B:512:TRP:HA	2.02	0.41
1:A:86:GLY:HA3	3:A:950:HOH:O	2.20	0.41
1:B:140:LYS:HA	1:B:151:GLN:O	2.20	0.41
1:B:250:PHE:CD1	1:B:254:LEU:HD23	2.55	0.41
1:B:252:VAL:HG13	1:B:256:ALA:HB2	2.02	0.41
1:B:425:ALA:HB3	1:B:426:PRO:CD	2.50	0.41
1:B:331:MET:HB3	1:B:338:PHE:HA	2.03	0.41
1:B:68:GLU:H	1:B:68:GLU:CD	2.29	0.41
1:B:233:ARG:HD3	1:B:233:ARG:HA	1.90	0.41
1:B:334:SER:O	1:B:547:LYS:HD2	2.20	0.41
1:A:369:LEU:C	1:A:369:LEU:CD2	2.93	0.41
1:A:38:LEU:HD21	1:A:491:LYS:HD3	2.03	0.41
1:B:403:PHE:CZ	1:B:473:GLN:HB2	2.55	0.41
1:A:68:GLU:HA	1:A:71:ILE:HG12	2.02	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:125:PHE:HD1	1:A:125:PHE:HA	1.72	0.41
1:A:271:MET:HE3	1:A:271:MET:HB2	1.99	0.41
1:B:350:GLU:HA	1:B:351:PRO:HD2	1.99	0.41
1:A:512:TRP:O	1:A:515:VAL:HG22	2.21	0.40
1:B:426:PRO:HB2	1:B:442:GLN:HE22	1.85	0.40
1:A:507:ALA:N	1:A:508:PRO:HD3	2.36	0.40
1:B:420:PHE:HB3	1:B:485:ILE:HG21	2.03	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	545/711 (77%)	533 (98%)	12 (2%)	0	100	100
1	B	545/711 (77%)	532 (98%)	13 (2%)	0	100	100
All	All	1090/1422 (77%)	1065 (98%)	25 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	432/553 (78%)	420 (97%)	12 (3%)	38	41

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	432/553 (78%)	421 (98%)	11 (2%)	42	45
All	All	864/1106 (78%)	841 (97%)	23 (3%)	39	42

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

Mol	Chain	Res	Type
1	A	49	ASP
1	A	54	THR
1	A	56	ILE
1	A	68	GLU
1	A	125	PHE
1	A	217	LYS
1	A	252	VAL
1	A	262	CYS
1	A	313	ARG
1	A	330	LYS
1	A	334	SER
1	A	464	LYS
1	B	68	GLU
1	B	102	GLU
1	B	125	PHE
1	B	201	GLU
1	B	215	LYS
1	B	227	ARG
1	B	252	VAL
1	B	276	GLU
1	B	313	ARG
1	B	333	LYS
1	B	369	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	28	GLN
1	A	96	ASN
1	A	442	GLN
1	B	28	GLN
1	B	96	ASN
1	B	321	ASN
1	B	390	ASN
1	B	442	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 [i](#)

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	547/711 (76%)	-1.13	0 100 100	23, 33, 48, 68	0
1	B	547/711 (76%)	-1.10	0 100 100	25, 34, 52, 70	0
All	All	1094/1422 (76%)	-1.12	0 100 100	23, 33, 50, 70	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](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q < 0.9’ lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	ZN	A	701	1/1	1.00	0.01	32,32,32,32	0
2	ZN	B	701	1/1	1.00	0.02	34,34,34,34	0

6.5 Other polymers [i](#)

There are no such residues in this entry.