



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 01:46 AM UTC

PDB ID : 3QLK / pdb_00003qlk
Title : Crystal Structure of RipA from Yersinia pestis
Authors : Torres, R.; Goulding, C.W.
Deposited on : 2011-02-02
Resolution : 3.00 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

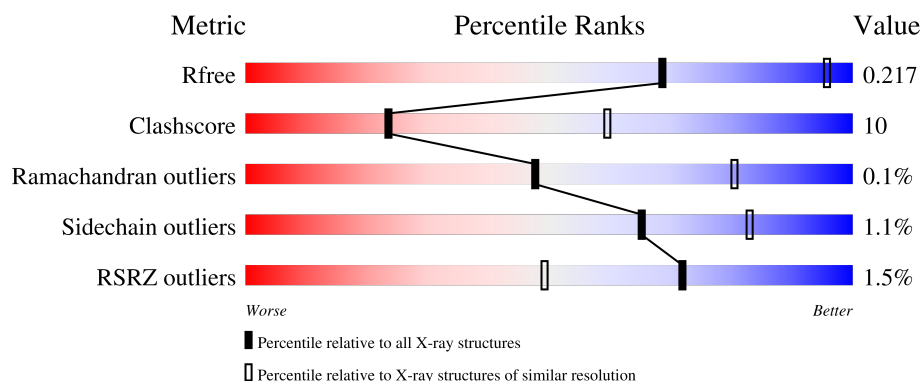
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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	2672 (3.00-3.00)
Clashscore	190562	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)
RSRZ outliers	180081	2671 (3.00-3.00)

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	455	2% 76% 18% • 5%
1	B	455	% 71% 22% • 5%

2 Entry composition [i](#)

There is only 1 type of molecule in this entry. The entry contains 6660 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 Coenzyme A transferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	431	Total	C	N	O	S	0	0	0
			3332	2092	598	626	16			
1	B	430	Total	C	N	O	S	0	0	0
			3328	2090	597	625	16			

There are 30 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-14	HIS	-	expression tag	UNP Q9ZC36
A	-13	HIS	-	expression tag	UNP Q9ZC36
A	-12	HIS	-	expression tag	UNP Q9ZC36
A	-11	HIS	-	expression tag	UNP Q9ZC36
A	-10	HIS	-	expression tag	UNP Q9ZC36
A	-9	HIS	-	expression tag	UNP Q9ZC36
A	-8	SER	-	expression tag	UNP Q9ZC36
A	-7	SER	-	expression tag	UNP Q9ZC36
A	-6	GLY	-	expression tag	UNP Q9ZC36
A	-5	LEU	-	expression tag	UNP Q9ZC36
A	-4	VAL	-	expression tag	UNP Q9ZC36
A	-3	PRO	-	expression tag	UNP Q9ZC36
A	-2	ARG	-	expression tag	UNP Q9ZC36
A	-1	GLY	-	expression tag	UNP Q9ZC36
A	0	SER	-	expression tag	UNP Q9ZC36
B	-14	HIS	-	expression tag	UNP Q9ZC36
B	-13	HIS	-	expression tag	UNP Q9ZC36
B	-12	HIS	-	expression tag	UNP Q9ZC36
B	-11	HIS	-	expression tag	UNP Q9ZC36
B	-10	HIS	-	expression tag	UNP Q9ZC36
B	-9	HIS	-	expression tag	UNP Q9ZC36
B	-8	SER	-	expression tag	UNP Q9ZC36
B	-7	SER	-	expression tag	UNP Q9ZC36
B	-6	GLY	-	expression tag	UNP Q9ZC36
B	-5	LEU	-	expression tag	UNP Q9ZC36

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-4	VAL	-	expression tag	UNP Q9ZC36
B	-3	PRO	-	expression tag	UNP Q9ZC36
B	-2	ARG	-	expression tag	UNP Q9ZC36
B	-1	GLY	-	expression tag	UNP Q9ZC36
B	0	SER	-	expression tag	UNP Q9ZC36

4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	197.51Å 197.51Å 61.10Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	49.72 – 3.00 49.72 – 3.00	Depositor EDS
% Data completeness (in resolution range)	96.7 (49.72-3.00) 82.7 (49.72-3.00)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.24 (at 3.01Å)	Xtriage
Refinement program	PHENIX 1.7 _650	Depositor
R, R_{free}	0.184 , 0.217 0.187 , 0.217	Depositor DCC
R_{free} test set	1998 reflections (7.24%)	wwPDB-VP
Wilson B-factor (Å ²)	48.3	Xtriage
Anisotropy	0.700	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 23.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.034 for -h,-k,l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6660	wwPDB-VP
Average B, all atoms (Å ²)	52.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.14% 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	0.76	1/3394 (0.0%)	1.10	9/4604 (0.2%)
1	B	0.75	0/3390	1.09	16/4599 (0.3%)
All	All	0.75	1/6784 (0.0%)	1.10	25/9203 (0.3%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	69	THR	CA-C	5.55	1.55	1.52

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	329	THR	N-CA-C	10.66	126.28	113.28
1	B	329	THR	N-CA-C	9.38	124.73	113.28
1	B	439	LEU	N-CA-C	-7.69	102.54	112.23
1	B	115	GLN	N-CA-C	-7.06	103.84	113.30
1	A	346	GLY	N-CA-C	-6.79	93.62	113.30
1	B	153	ARG	N-CA-C	6.11	118.01	111.36
1	A	86	VAL	N-CA-C	5.92	116.32	107.51
1	B	385	GLY	CA-C-N	5.85	125.61	119.76
1	B	385	GLY	C-N-CA	5.85	125.61	119.76
1	B	61	GLU	N-CA-C	5.77	118.23	111.02
1	B	224	GLN	N-CA-C	-5.64	99.83	109.24
1	B	89	VAL	N-CA-C	5.55	115.75	110.42
1	A	146	VAL	CA-C-N	-5.52	117.45	122.80
1	A	146	VAL	C-N-CA	-5.52	117.45	122.80
1	A	147	GLY	N-CA-C	5.48	120.47	110.71
1	B	149	ASP	CB-CA-C	-5.38	110.35	116.54
1	B	115	GLN	N-CA-CB	5.36	118.71	110.61
1	A	150	TYR	N-CA-C	-5.36	106.74	113.28
1	B	73	TYR	N-CA-C	5.26	117.70	111.33
1	B	336	HIS	CA-CB-CG	-5.25	108.56	113.80
1	A	63	ALA	N-CA-CB	-5.11	103.16	110.87

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	149	ASP	CB-CA-C	-5.07	110.71	116.54
1	B	12	THR	N-CA-CB	5.07	118.88	110.97
1	B	249	GLU	N-CA-C	-5.02	104.77	111.24
1	B	86	VAL	N-CA-C	5.00	114.97	107.51

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	3332	0	3339	63	4
1	B	3328	0	3336	77	4
All	All	6660	0	6675	136	4

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 10.

All (136) 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:331:ALA:HB1	1:A:387:VAL:HG21	1.37	1.00
1:B:91:ARG:HD3	1:B:336:HIS:NE2	1.80	0.97
1:B:47:ALA:HB3	1:B:75:LEU:HD23	1.47	0.93
1:B:59:TYR:HE1	1:B:62:THR:HG23	1.34	0.92
1:B:91:ARG:HD3	1:B:336:HIS:CD2	2.07	0.90
1:B:76:ASN:ND2	1:B:105:VAL:HB	1.89	0.87
1:A:275:ARG:HD3	1:A:298:PHE:HB2	1.56	0.85
1:B:59:TYR:CE1	1:B:62:THR:HG23	2.10	0.85
1:B:324:LEU:O	1:B:378:ARG:HD3	1.84	0.78
1:B:327:ASP:HB3	1:B:383:ILE:HD12	1.64	0.77
1:A:236:GLY:O	1:A:239:LYS:HG2	1.84	0.75
1:B:59:TYR:CE1	1:B:62:THR:CG2	2.71	0.73
1:B:227:VAL:HG21	1:B:282:MET:HE3	1.71	0.71
1:A:90:GLU:O	1:A:94:ILE:HG12	1.91	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:112:ASN:HB3	1:B:114:HIS:CE1	2.27	0.70
1:A:73:TYR:CE1	1:A:96:ARG:HB3	2.26	0.70
1:A:230:LEU:HB3	1:A:231:PRO:HD3	1.72	0.70
1:B:47:ALA:CB	1:B:75:LEU:HD23	2.21	0.69
1:B:74:GLU:HG2	1:B:75:LEU:HD12	1.75	0.69
1:A:47:ALA:HB3	1:A:75:LEU:HD22	1.75	0.69
1:A:225:MET:HE2	1:A:320:ILE:HD12	1.75	0.69
1:A:73:TYR:OH	1:A:100:ASP:OD2	2.08	0.69
1:B:73:TYR:OH	1:B:103:ARG:NH1	2.26	0.69
1:B:70:ILE:HB	1:B:79:ILE:HG21	1.73	0.68
1:B:76:ASN:HD22	1:B:105:VAL:HB	1.57	0.68
1:B:406:LEU:HA	1:B:409:LEU:HD12	1.77	0.66
1:A:24:GLY:HA2	1:A:54:ASP:O	1.96	0.66
1:B:111:SER:HB2	1:B:119:LEU:CD1	2.27	0.65
1:A:380:ILE:HB	1:A:381:PRO:HD2	1.78	0.64
1:B:91:ARG:NH1	1:B:336:HIS:CE1	2.65	0.64
1:B:88:ALA:HA	1:B:91:ARG:CZ	2.28	0.64
1:A:110:PRO:HD2	1:B:414:ARG:HH11	1.63	0.64
1:B:91:ARG:NH1	1:B:336:HIS:NE2	2.45	0.63
1:A:321:ASN:HD22	1:A:348:GLN:HG3	1.64	0.63
1:A:412:THR:OG1	1:B:104:LYS:HE2	1.99	0.62
1:B:226:GLY:O	1:B:231:PRO:HG2	1.99	0.62
1:A:225:MET:HG2	1:A:231:PRO:HB3	1.81	0.62
1:B:134:SER:OG	1:B:143:SER:HB3	1.99	0.62
1:A:332:CYS:O	1:A:387:VAL:HG22	2.00	0.61
1:B:101:GLY:O	1:B:102:GLY:C	2.44	0.60
1:B:230:LEU:C	1:B:230:LEU:HD23	2.26	0.60
1:B:272:ASP:OD1	1:B:275:ARG:NH2	2.34	0.60
1:A:59:TYR:HE1	1:A:62:THR:HB	1.67	0.59
1:B:91:ARG:CD	1:B:336:HIS:CD2	2.85	0.59
1:B:209:ILE:HG21	1:B:230:LEU:HD11	1.83	0.58
1:B:230:LEU:HB3	1:B:231:PRO:HD3	1.84	0.58
1:A:96:ARG:O	1:A:99:GLU:HG2	2.03	0.58
1:B:227:VAL:HG21	1:B:282:MET:CE	2.33	0.57
1:A:275:ARG:NH1	1:A:295:PRO:O	2.38	0.57
1:B:236:GLY:O	1:B:239:LYS:HG2	2.04	0.57
1:A:331:ALA:CB	1:A:387:VAL:HG21	2.26	0.55
1:B:111:SER:HB2	1:B:119:LEU:HD13	1.89	0.54
1:A:73:TYR:CZ	1:A:103:ARG:NH2	2.71	0.54
1:A:133:VAL:CG1	1:A:163:VAL:HG22	2.38	0.54
1:B:416:LEU:O	1:B:419:ILE:HG13	2.09	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:74:GLU:H	1:A:74:GLU:CD	2.17	0.53
1:A:275:ARG:HD3	1:A:298:PHE:CB	2.35	0.53
1:B:323:THR:HG21	1:B:332:CYS:HB3	1.91	0.53
1:B:228:GLY:O	1:B:231:PRO:HD2	2.08	0.53
1:A:94:ILE:HD11	1:A:106:VAL:HB	1.91	0.52
1:B:225:MET:CE	1:B:231:PRO:HB3	2.39	0.52
1:B:269:LYS:HE2	1:B:271:LEU:O	2.10	0.52
1:A:134:SER:OG	1:A:143:SER:HB3	2.10	0.52
1:A:111:SER:HB2	1:A:119:LEU:CD1	2.40	0.51
1:B:34:ALA:HB1	1:B:131:HIS:CE1	2.45	0.51
1:A:336:HIS:ND1	1:A:336:HIS:N	2.59	0.51
1:A:59:TYR:CE1	1:A:62:THR:HB	2.45	0.51
1:A:110:PRO:HD2	1:B:414:ARG:NH1	2.25	0.50
1:B:327:ASP:OD1	1:B:330:GLY:HA2	2.11	0.50
1:A:321:ASN:ND2	1:A:348:GLN:HG3	2.27	0.50
1:A:383:ILE:HD13	1:A:387:VAL:HG23	1.94	0.50
1:A:206:TYR:HE2	1:A:229:ALA:HB3	1.75	0.49
1:A:84:MET:HG2	1:A:109:VAL:HB	1.93	0.49
1:A:67:GLY:HA2	1:A:71:LEU:HB2	1.94	0.49
1:A:230:LEU:HB3	1:A:231:PRO:CD	2.42	0.49
1:B:59:TYR:HE1	1:B:62:THR:CG2	2.09	0.49
1:A:90:GLU:HA	1:A:93:LEU:HD12	1.94	0.49
1:B:91:ARG:CD	1:B:336:HIS:NE2	2.66	0.48
1:B:203:ILE:CD1	1:B:205:GLU:HB2	2.44	0.48
1:A:194:LEU:HD12	1:A:288:TYR:CG	2.49	0.47
1:B:24:GLY:HA2	1:B:54:ASP:O	2.14	0.47
1:B:328:LEU:HD13	1:B:434:ALA:HB2	1.96	0.47
1:B:197:MET:SD	1:B:227:VAL:CG1	3.03	0.46
1:B:327:ASP:OD1	1:B:327:ASP:C	2.59	0.46
1:B:415:ALA:O	1:B:419:ILE:HG23	2.15	0.46
1:A:401:PHE:CE2	1:A:423:HIS:HA	2.51	0.46
1:A:116:ALA:N	1:A:117:PRO:CD	2.79	0.46
1:B:76:ASN:HD21	1:B:105:VAL:HB	1.73	0.46
1:A:382:ARG:HB2	1:A:437:MET:HE1	1.98	0.46
1:A:250:VAL:HG21	1:A:282:MET:HE2	1.98	0.45
1:B:397:ILE:HD11	1:B:418:ILE:HG12	1.99	0.45
1:B:27:LEU:HD12	1:B:128:THR:HB	1.99	0.45
1:B:255:LEU:HA	1:B:255:LEU:HD23	1.71	0.45
1:A:133:VAL:HG11	1:A:163:VAL:HG22	1.98	0.45
1:B:327:ASP:HA	1:B:380:ILE:O	2.16	0.45
1:A:73:TYR:CE2	1:A:103:ARG:NH2	2.85	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:423:HIS:CD2	1:A:424:PRO:HD2	2.52	0.45
1:B:372:ALA:O	1:B:375:THR:OG1	2.33	0.45
1:A:269:LYS:HE3	1:A:271:LEU:O	2.16	0.45
1:A:71:LEU:HD22	1:A:81:PRO:HB3	1.99	0.45
1:A:94:ILE:CD1	1:A:106:VAL:HB	2.47	0.45
1:B:413:GLU:HB3	1:B:417:ARG:NH1	2.32	0.44
1:B:423:HIS:CG	1:B:424:PRO:HD2	2.53	0.44
1:A:2:ASP:OD2	1:A:5:ALA:HB2	2.18	0.44
1:A:34:ALA:HB1	1:A:131:HIS:CE1	2.53	0.43
1:A:249:GLU:O	1:A:279:THR:HG22	2.18	0.43
1:A:104:LYS:HE2	1:B:412:THR:OG1	2.18	0.43
1:B:197:MET:SD	1:B:227:VAL:HG11	2.58	0.43
1:B:248:THR:O	1:B:278:PHE:HA	2.18	0.43
1:A:327:ASP:HA	1:A:380:ILE:O	2.19	0.43
1:A:56:ARG:HD3	1:A:82:TYR:CZ	2.54	0.42
1:A:133:VAL:HG13	1:A:163:VAL:HG22	2.00	0.42
1:B:27:LEU:O	1:B:57:VAL:HA	2.19	0.42
1:B:73:TYR:HH	1:B:103:ARG:NH1	2.17	0.42
1:B:3:ILE:HD12	1:B:136:MET:HE3	2.02	0.42
1:B:88:ALA:HA	1:B:91:ARG:NH1	2.35	0.42
1:A:206:TYR:CE2	1:A:229:ALA:HB3	2.53	0.42
1:B:180:HIS:CE1	1:B:182:SER:HG	2.28	0.41
1:B:271:LEU:HD11	1:B:313:GLN:HG3	2.01	0.41
1:A:133:VAL:HG22	1:A:163:VAL:HG13	2.00	0.41
1:A:104:LYS:HB3	1:A:104:LYS:HE3	1.84	0.41
1:B:75:LEU:HD12	1:B:75:LEU:N	2.36	0.41
1:B:161:PHE:HE2	1:B:179:ILE:HG21	1.86	0.41
1:B:327:ASP:CB	1:B:383:ILE:HD12	2.42	0.41
1:B:35:GLU:OE1	1:B:63:ALA:HB3	2.21	0.41
1:A:248:THR:O	1:A:278:PHE:HA	2.21	0.41
1:B:323:THR:HG22	1:B:334:SER:OG	2.21	0.41
1:A:63:ALA:HB1	1:A:195:ILE:HD13	2.03	0.41
1:A:133:VAL:HG11	1:A:163:VAL:CG2	2.50	0.41
1:B:216:LEU:O	1:B:218:PRO:HD3	2.21	0.41
1:B:249:GLU:O	1:B:279:THR:HG22	2.21	0.40
1:A:36:PRO:HA	1:A:37:PRO:HD2	2.02	0.40
1:A:382:ARG:HB2	1:A:437:MET:CE	2.51	0.40
1:B:156:ARG:CZ	1:B:179:ILE:HD13	2.51	0.40
1:B:84:MET:HG2	1:B:109:VAL:HB	2.04	0.40
1:B:101:GLY:O	1:B:103:ARG:N	2.54	0.40

All (4) symmetry-related close contacts are listed below. The label for Atom-2 includes the sym-

metry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:99:GLU:OE2	1:B:78:ARG:NH1[3_445]	1.54	0.66
1:A:99:GLU:OE2	1:B:78:ARG:NH2[3_445]	1.76	0.44
1:A:99:GLU:OE2	1:B:78:ARG:CZ[3_445]	1.77	0.43
1:A:100:ASP:OD1	1:B:78:ARG:NE[3_445]	1.78	0.42

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	427/455 (94%)	414 (97%)	13 (3%)	0	100	100
1	B	426/455 (94%)	411 (96%)	14 (3%)	1 (0%)	43	76
All	All	853/910 (94%)	825 (97%)	27 (3%)	1 (0%)	48	80

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	102	GLY

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	359/379 (95%)	356 (99%)	3 (1%)	73	86
1	B	359/379 (95%)	354 (99%)	5 (1%)	59	80

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	718/758 (95%)	710 (99%)	8 (1%)	65 83

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	163	VAL
1	A	199	VAL
1	A	336	HIS
1	B	103	ARG
1	B	201	SER
1	B	275	ARG
1	B	279	THR
1	B	440	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	173	GLN
1	A	316	ASN
1	B	107	ASN
1	B	131	HIS
1	B	316	ASN
1	B	325	GLN
1	B	333	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	431/455 (94%)	-0.18	7 (1%)	70 47	30, 47, 69, 89	0
1	B	430/455 (94%)	0.00	6 (1%)	73 51	34, 52, 84, 105	0
All	All	861/910 (94%)	-0.09	13 (1%)	72 49	30, 50, 80, 105	0

All (13) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	336	HIS	3.9
1	B	386	PRO	3.6
1	A	62	THR	3.0
1	B	114	HIS	2.8
1	A	438	HIS	2.8
1	B	226	GLY	2.6
1	B	242	ASN	2.5
1	B	1	MET	2.4
1	A	386	PRO	2.4
1	B	201	SER	2.3
1	A	348	GLN	2.3
1	A	100	ASP	2.2
1	A	85	PHE	2.1

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no 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.