



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 06:06 PM GMT

PDB ID : 4KJS
Title : Structure of native YfkE
Authors : Wu, M.; Tong, S.; Zheng, L.
Deposited on : 2013-05-03
Resolution : 3.05 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

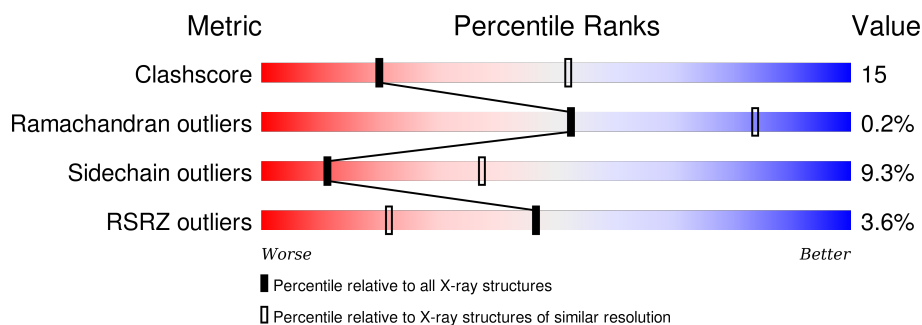
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.05 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))
Clashscore	102246	1303 (3.10-3.02)
Ramachandran outliers	100387	1254 (3.10-3.02)
Sidechain outliers	100360	1254 (3.10-3.02)
RSRZ outliers	91569	1197 (3.10-3.02)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	351	<div> <div>3%</div> <div>64%</div> <div>24%</div> <div>• • 9%</div> </div>
1	B	351	<div> <div>3%</div> <div>69%</div> <div>17%</div> <div>5% • 9%</div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4800 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 cation exchanger YfkE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	320	Total	C	N	O	S	0	0	0
			2391	1596	369	412	14			
1	B	320	Total	C	N	O	S	0	0	0
			2391	1596	369	412	14			

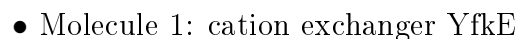
There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	77	MET	LEU	CONFLICT	UNP O34840
A	116	ALA	LYS	CONFLICT	UNP O34840
B	77	MET	LEU	CONFLICT	UNP O34840
B	116	ALA	LYS	CONFLICT	UNP O34840

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	8	Total	O	0	0
			8	8		
2	B	10	Total	O	0	0
			10	10		

- Molecule 1: cation exchanger YfkE



4 Data and refinement statistics

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, α , β , γ	168.50 Å 168.50 Å 94.17 Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	50.01 – 3.05 47.59 – 3.05	Depositor EDS
% Data completeness (in resolution range)	99.8 (50.01-3.05) 99.9 (47.59-3.05)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.06 (at 3.07 Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.216 , 0.262 (Not available) , (Not available)	Depositor DCC
R_{free} test set	NotAvailable	DCC
Wilson B-factor (Å ²)	83.9	Xtriage
Anisotropy	0.175	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 66.3	EDS
Estimated twinning fraction	0.486 for H, K, L 0.514 for K, H, -L 0.036 for h,-h-k,-l	Xtriage
Reported twinning fraction	0.486 for H, K, L 0.514 for K, H, -L	Depositor
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 18969 reflections	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	4800	wwPDB-VP
Average B, all atoms (Å ²)	83.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.04% 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, 0.375 respectively for untwinned datasets, and 0.333, 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	0.46	2/2438 (0.1%)	0.62	6/3317 (0.2%)
1	B	0.50	1/2438 (0.0%)	0.60	2/3317 (0.1%)
All	All	0.48	3/4876 (0.1%)	0.61	8/6634 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2
1	B	0	3
All	All	0	5

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	237	TRP	CD2-CE2	5.15	1.47	1.41
1	B	237	TRP	CD2-CE2	5.14	1.47	1.41
1	A	24	TRP	CD2-CE2	5.05	1.47	1.41

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	116	ALA	N-CA-CB	8.23	121.62	110.10
1	A	328	ASP	CB-CA-C	8.03	126.45	110.40
1	A	329	SER	N-CA-CB	-8.02	98.47	110.50
1	A	115	TYR	N-CA-C	7.08	130.10	111.00
1	A	326	ASP	N-CA-C	6.80	129.36	111.00
1	A	115	TYR	CB-CA-C	-6.76	96.88	110.40
1	B	120	PHE	N-CA-CB	-5.81	100.15	110.60
1	B	24	TRP	N-CA-C	-5.00	97.50	111.00

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	297	PHE	Peptide
1	A	299	THR	Peptide
1	B	120	PHE	Peptide
1	B	24	TRP	Peptide
1	B	297	PHE	Peptide

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	2391	0	2509	73	0
1	B	2391	0	2509	78	0
2	A	8	0	0	0	0
2	B	10	0	0	0	0
All	All	4800	0	5018	145	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 15.

All (145) 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:299:THR:HG21	1:B:82:GLU:O	1.05	1.21
1:B:20:SER:O	1:B:23:HIS:CD2	1.96	1.18
1:B:122:ILE:HG13	1:B:123:HIS:H	1.08	1.14
1:B:120:PHE:HA	1:B:328:ASP:OD2	1.45	1.12
1:A:299:THR:CG2	1:B:82:GLU:O	2.01	1.09
1:B:122:ILE:HG13	1:B:123:HIS:N	1.60	1.08
1:B:301:MET:O	1:B:301:MET:HG3	1.52	1.03
1:B:20:SER:O	1:B:23:HIS:NE2	1.91	1.01
1:A:299:THR:HG21	1:B:82:GLU:C	1.83	0.97
1:A:295:ILE:HD12	1:A:300:SER:HB3	1.47	0.96
1:B:120:PHE:CA	1:B:328:ASP:OD2	2.16	0.93
1:A:295:ILE:CD1	1:A:300:SER:HB3	1.99	0.92

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:117:ARG:O	1:A:118:GLN:HG2	1.69	0.91
1:A:118:GLN:HG3	1:A:329:SER:HB3	1.51	0.91
1:A:118:GLN:HA	1:A:118:GLN:HE21	1.36	0.88
1:A:117:ARG:C	1:A:118:GLN:HG2	1.94	0.88
1:B:323:ILE:O	1:B:326:ASP:OD1	1.95	0.85
1:B:122:ILE:CG1	1:B:123:HIS:H	1.89	0.85
1:B:25:PRO:O	1:B:29:LEU:HD22	1.76	0.84
1:A:108:PHE:CE2	1:A:328:ASP:O	2.32	0.82
1:B:298:PRO:HD2	1:B:299:THR:H	1.45	0.81
1:B:120:PHE:CD1	1:B:120:PHE:N	2.48	0.80
1:A:117:ARG:O	1:A:118:GLN:CG	2.30	0.79
1:B:20:SER:O	1:B:23:HIS:HD2	1.67	0.78
1:B:326:ASP:CG	1:B:327:GLY:N	2.37	0.78
1:A:114:LYS:O	1:A:115:TYR:HB2	1.83	0.78
1:B:20:SER:HG	1:B:219:TYR:HH	0.80	0.78
1:B:301:MET:O	1:B:301:MET:CG	2.30	0.77
1:B:295:ILE:HG22	1:B:300:SER:HB2	1.66	0.75
1:A:301:MET:HG3	1:A:301:MET:O	1.86	0.75
1:A:116:ALA:O	1:A:117:ARG:HG2	1.89	0.72
1:B:96:VAL:HB	1:B:279:THR:HG23	1.71	0.72
1:A:173:ALA:O	1:A:177:LYS:N	2.24	0.71
1:A:299:THR:HG23	1:B:83:GLY:HA3	1.73	0.70
1:B:26:SER:HB3	1:B:227:THR:HG23	1.73	0.70
1:A:299:THR:CG2	1:B:82:GLU:C	2.54	0.69
1:A:173:ALA:O	1:A:177:LYS:C	2.31	0.69
1:A:229:HIS:O	1:A:233:GLU:HG3	1.93	0.68
1:B:317:VAL:O	1:B:321:ILE:HG12	1.94	0.67
1:A:172:ALA:O	1:A:176:PHE:HD1	1.77	0.67
1:A:164:ILE:HD13	1:A:168:LEU:HD11	1.77	0.66
1:B:326:ASP:OD1	1:B:327:GLY:N	2.30	0.65
1:A:117:ARG:O	1:A:118:GLN:NE2	2.30	0.64
1:A:164:ILE:HD13	1:A:168:LEU:CD1	2.27	0.63
1:B:160:ILE:HD11	1:B:295:ILE:HG21	1.81	0.63
1:B:24:TRP:CE3	1:B:25:PRO:HD3	2.34	0.62
1:B:298:PRO:CD	1:B:299:THR:H	2.13	0.62
1:A:164:ILE:CD1	1:A:168:LEU:HD11	2.29	0.62
1:B:121:ASN:O	1:B:122:ILE:C	2.37	0.62
1:A:67:PHE:HA	1:A:70:ALA:HB2	1.82	0.61
1:B:25:PRO:HG2	1:B:28:VAL:HG22	1.82	0.61
1:B:222:GLU:O	1:B:226:HIS:HD2	1.85	0.60
1:A:326:ASP:OD1	1:A:327:GLY:N	2.34	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:164:ILE:O	1:A:168:LEU:HG	2.03	0.59
1:A:297:PHE:CG	1:A:297:PHE:O	2.56	0.59
1:A:240:LEU:O	1:A:244:VAL:HG23	2.03	0.59
1:A:117:ARG:O	1:A:118:GLN:CD	2.40	0.59
1:B:108:PHE:CZ	1:B:328:ASP:O	2.56	0.59
1:B:30:PHE:CD2	1:B:231:VAL:HG21	2.38	0.58
1:A:118:GLN:HA	1:A:118:GLN:NE2	2.14	0.58
1:B:26:SER:HB3	1:B:227:THR:CG2	2.33	0.58
1:A:229:HIS:O	1:A:233:GLU:CG	2.51	0.58
1:A:339:ALA:O	1:A:343:ILE:HG12	2.02	0.58
1:B:299:THR:HG23	1:B:300:SER:N	2.19	0.57
1:B:118:GLN:HG3	1:B:329:SER:HB3	1.87	0.57
1:B:327:GLY:O	1:B:328:ASP:HB2	2.04	0.56
1:B:67:PHE:HA	1:B:70:ALA:HB2	1.86	0.56
1:B:297:PHE:CG	1:B:297:PHE:O	2.59	0.56
1:A:141:VAL:HA	1:A:310:LEU:HD11	1.87	0.55
1:A:99:ASN:HA	1:A:103:VAL:HG22	1.88	0.55
1:A:229:HIS:CE1	1:A:233:GLU:OE2	2.60	0.55
1:A:216:VAL:O	1:A:220:ILE:HG12	2.07	0.54
1:B:120:PHE:N	1:B:328:ASP:OD2	2.41	0.54
1:A:13:VAL:HB	1:A:14:PRO:HD3	1.90	0.54
1:A:167:ILE:HG23	1:A:288:PRO:HB2	1.90	0.53
1:B:268:MET:SD	1:B:329:SER:HB2	2.49	0.53
1:A:118:GLN:HE21	1:A:118:GLN:CA	2.15	0.53
1:B:278:SER:O	1:B:281:GLN:HG2	2.09	0.52
1:A:295:ILE:HD11	1:A:300:SER:HB3	1.86	0.52
1:B:114:LYS:HE3	1:B:115:TYR:CD1	2.45	0.52
1:A:144:GLU:HG2	1:A:310:LEU:HD22	1.91	0.52
1:A:245:ILE:HD11	1:A:345:ALA:HA	1.91	0.51
1:B:118:GLN:HE22	1:B:266:ASN:HA	1.76	0.50
1:A:35:VAL:HA	1:A:38:ILE:HD12	1.93	0.50
1:B:108:PHE:CE2	1:B:328:ASP:O	2.66	0.49
1:B:99:ASN:ND2	1:B:279:THR:OG1	2.45	0.49
1:B:24:TRP:CD2	1:B:25:PRO:HD3	2.47	0.49
1:A:82:GLU:HB3	1:A:84:LEU:HD13	1.94	0.49
1:A:330:ASN:HB3	1:A:333:GLU:HB2	1.94	0.49
1:B:298:PRO:CD	1:B:299:THR:N	2.76	0.49
1:B:297:PHE:N	1:B:297:PHE:CD1	2.81	0.49
1:A:82:GLU:OE1	1:A:225:VAL:HG12	2.12	0.49
1:A:128:ASN:HB3	1:A:276:VAL:HG11	1.95	0.48
1:B:13:VAL:HG13	1:B:14:PRO:HD3	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:285:PHE:O	1:A:289:VAL:HG23	2.15	0.47
1:A:24:TRP:HA	1:A:25:PRO:HA	1.78	0.47
1:A:298:PRO:HG2	1:A:299:THR:H	1.80	0.47
1:A:298:PRO:C	1:A:299:THR:HG22	2.36	0.46
1:B:325:ASN:N	1:B:325:ASN:OD1	2.46	0.46
1:B:122:ILE:CG1	1:B:123:HIS:N	2.41	0.46
1:B:76:SER:HB3	1:B:88:VAL:HG13	1.98	0.46
1:A:95:SER:HB3	1:A:282:ILE:HD13	1.96	0.46
1:A:173:ALA:O	1:A:177:LYS:CA	2.63	0.46
1:A:317:VAL:O	1:A:321:ILE:HG13	2.16	0.45
1:B:29:LEU:HA	1:B:29:LEU:HD12	1.66	0.45
1:B:129:SER:HB3	1:B:321:ILE:HG23	1.99	0.45
1:A:69:ASN:HD21	1:A:281:GLN:HG3	1.82	0.45
1:B:119:GLU:C	1:B:328:ASP:OD2	2.55	0.44
1:A:270:ILE:O	1:A:274:ILE:HG12	2.17	0.44
1:B:323:ILE:HD12	1:B:323:ILE:HA	1.72	0.44
1:B:128:ASN:HB3	1:B:276:VAL:HG11	1.99	0.44
1:B:28:VAL:HG23	1:B:29:LEU:N	2.32	0.44
1:B:20:SER:C	1:B:23:HIS:HD2	2.20	0.43
1:B:118:GLN:HB2	1:B:329:SER:HB3	1.99	0.43
1:A:34:CYS:O	1:A:38:ILE:HG13	2.18	0.43
1:B:252:ASN:O	1:B:256:HIS:HD2	2.01	0.43
1:A:87:ILE:HD11	1:A:239:GLU:HB3	2.00	0.43
1:A:76:SER:HB3	1:A:88:VAL:HG13	2.00	0.43
1:B:292:ILE:O	1:B:295:ILE:HG12	2.18	0.43
1:B:156:LEU:HD13	1:B:300:SER:CB	2.48	0.43
1:A:301:MET:CG	1:A:301:MET:O	2.61	0.42
1:B:307:LEU:HB3	1:B:308:PRO:CD	2.49	0.42
1:B:114:LYS:HE3	1:B:115:TYR:HD1	1.83	0.42
1:A:85:THR:OG1	1:A:301:MET:HB2	2.19	0.42
1:A:287:ALA:HB3	1:A:288:PRO:HD3	2.02	0.42
1:A:96:VAL:HB	1:A:279:THR:HG23	2.02	0.42
1:B:281:GLN:O	1:B:285:PHE:HB3	2.19	0.42
1:A:297:PHE:HA	1:A:298:PRO:HA	1.80	0.42
1:A:301:MET:HA	1:A:302:PRO:HD3	1.80	0.42
1:A:282:ILE:HA	1:A:286:VAL:HB	2.01	0.42
1:B:108:PHE:N	1:B:108:PHE:CD1	2.87	0.41
1:A:80:LEU:HB2	1:A:297:PHE:CE1	2.55	0.41
1:A:156:LEU:HD22	1:A:300:SER:HB2	2.03	0.41
1:B:34:CYS:O	1:B:38:ILE:HG13	2.21	0.41
1:B:108:PHE:HE2	1:B:323:ILE:HG21	1.86	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:100:LEU:HD22	1:B:320:MET:HG2	2.03	0.41
1:A:25:PRO:HD2	1:A:28:VAL:HB	2.01	0.41
1:B:282:ILE:HD13	1:B:286:VAL:HG21	2.03	0.41
1:B:120:PHE:HA	1:B:328:ASP:CG	2.31	0.41
1:A:262:MET:HG3	1:A:267:LYS:HB2	2.03	0.41
1:B:108:PHE:HE2	1:B:323:ILE:CG2	2.35	0.40
1:A:161:GLY:O	1:A:164:ILE:HG22	2.21	0.40
1:A:299:THR:CG2	1:B:83:GLY:HA3	2.46	0.40
1:B:98:GLY:O	1:B:102:LEU:HB3	2.20	0.40
1:B:118:GLN:CG	1:B:329:SER:HB3	2.49	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	314/351 (90%)	295 (94%)	19 (6%)	0	100	100
1	B	314/351 (90%)	298 (95%)	15 (5%)	1 (0%)	46	78
All	All	628/702 (90%)	593 (94%)	34 (5%)	1 (0%)	52	84

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	327	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	254/280 (91%)	234 (92%)	20 (8%)	15	46
1	B	254/280 (91%)	227 (89%)	27 (11%)	8	30
All	All	508/560 (91%)	461 (91%)	47 (9%)	11	37

All (47) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	PHE
1	A	51	LEU
1	A	64	ASN
1	A	97	LEU
1	A	100	LEU
1	A	118	GLN
1	A	120	PHE
1	A	122	ILE
1	A	137	ILE
1	A	164	ILE
1	A	174	LEU
1	A	177	LYS
1	A	224	LEU
1	A	290	LEU
1	A	293	CYS
1	A	299	THR
1	A	300	SER
1	A	307	LEU
1	A	328	ASP
1	A	332	PHE
1	B	13	VAL
1	B	26	SER
1	B	29	LEU
1	B	44	MET
1	B	49	GLU
1	B	82	GLU
1	B	96	VAL
1	B	113	LEU
1	B	114	LYS
1	B	120	PHE
1	B	142	ILE
1	B	156	LEU
1	B	220	ILE

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Mol	Chain	Res	Type
1	B	222	GLU
1	B	227	THR
1	B	229	HIS
1	B	266	ASN
1	B	281	GLN
1	B	297	PHE
1	B	299	THR
1	B	300	SER
1	B	301	MET
1	B	307	LEU
1	B	314	VAL
1	B	323	ILE
1	B	325	ASN
1	B	350	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	64	ASN
1	A	69	ASN
1	A	118	GLN
1	A	229	HIS
1	A	281	GLN
1	A	325	ASN
1	B	99	ASN
1	B	118	GLN
1	B	226	HIS
1	B	266	ASN
1	B	330	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 [i](#)

There are no carbohydrates 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 ⓘ

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	320/351 (91%)	-0.15	11 (3%)	49 23	52, 81, 107, 128	0
1	B	320/351 (91%)	-0.21	12 (3%)	44 20	50, 83, 111, 126	0
All	All	640/702 (91%)	-0.18	23 (3%)	46 21	50, 82, 110, 128	0

All (23) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	175	TYR	6.2
1	B	63	LEU	4.6
1	A	175	TYR	4.0
1	A	264	PHE	3.9
1	A	208	ILE	3.7
1	B	4	ILE	3.6
1	B	5	PHE	3.4
1	B	6	PHE	3.4
1	A	261	ILE	3.4
1	B	265	LYS	3.3
1	B	176	PHE	3.2
1	A	204	LYS	2.9
1	B	124	ASP	2.9
1	B	264	PHE	2.8
1	A	265	LYS	2.8
1	B	43	TYR	2.8
1	A	124	ASP	2.7
1	B	62	LEU	2.7
1	A	219	TYR	2.6
1	B	126	ARG	2.5
1	A	223	ASN	2.1
1	A	4	ILE	2.1
1	A	22	MET	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 carbohydrates 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.