



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 12:39 AM GMT

PDB ID : 2BAS  
Title : Crystal Structure of the Bacillus subtilis YkuI Protein, with an EAL Domain.  
Authors : Minasov, G.; Brunzelle, J.S.; Shuvalova, L.; Miller, D.J.; Collart, F.R.;  
Joachimciak, A.; Anderson, W.F.; Midwest Center for Structural Genomics  
(MCSG)  
Deposited on : 2005-10-14  
Resolution : 2.61 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](mailto: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.

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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

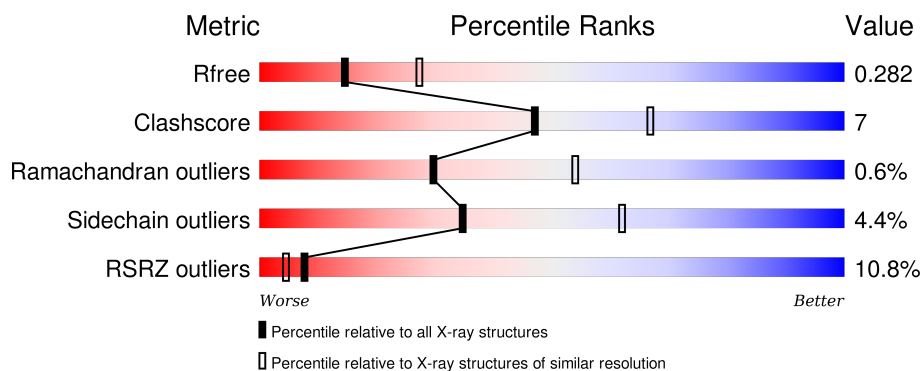
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.61 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))
$R_{free}$	91344	2700 (2.64-2.60)
Clashscore	102246	3065 (2.64-2.60)
Ramachandran outliers	100387	3015 (2.64-2.60)
Sidechain outliers	100360	3015 (2.64-2.60)
RSRZ outliers	91569	2706 (2.64-2.60)

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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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	431	
1	B	431	

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6915 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 Yku1 protein.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	402	Total	C	N	O	S	Se	0	0	0
			3351	2158	539	643	2	9			
1	B	397	Total	C	N	O	S	Se	0	0	0
			3305	2129	533	631	2	10			

There are 48 discrepancies between the modelled and reference sequences:

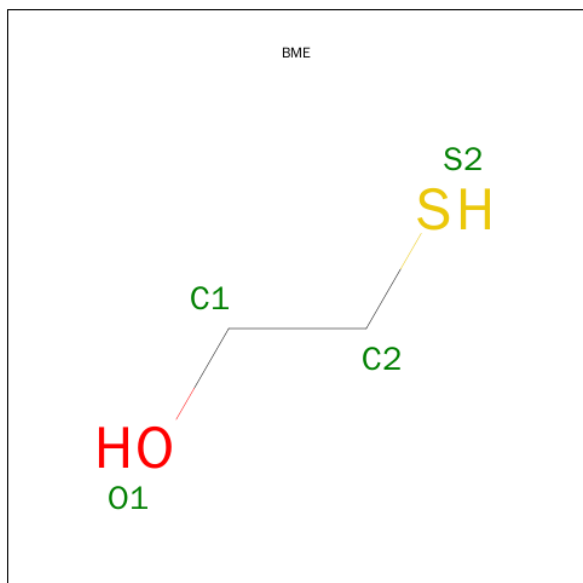
Chain	Residue	Modelled	Actual	Comment	Reference
A	-23	MSE	-	INITIATING METHIONINE	UNP O35014
A	-22	HIS	-	EXPRESSION TAG	UNP O35014
A	-21	HIS	-	EXPRESSION TAG	UNP O35014
A	-20	HIS	-	EXPRESSION TAG	UNP O35014
A	-19	HIS	-	EXPRESSION TAG	UNP O35014
A	-18	HIS	-	EXPRESSION TAG	UNP O35014
A	-17	HIS	-	EXPRESSION TAG	UNP O35014
A	-16	SER	-	CLONING ARTIFACT	UNP O35014
A	-15	SER	-	CLONING ARTIFACT	UNP O35014
A	-14	GLY	-	CLONING ARTIFACT	UNP O35014
A	-13	VAL	-	CLONING ARTIFACT	UNP O35014
A	-12	ASP	-	CLONING ARTIFACT	UNP O35014
A	-11	LEU	-	CLONING ARTIFACT	UNP O35014
A	-10	GLY	-	CLONING ARTIFACT	UNP O35014
A	-9	THR	-	CLONING ARTIFACT	UNP O35014
A	-8	GLU	-	CLONING ARTIFACT	UNP O35014
A	-7	ASN	-	CLONING ARTIFACT	UNP O35014
A	-6	LEU	-	CLONING ARTIFACT	UNP O35014
A	-5	TYR	-	CLONING ARTIFACT	UNP O35014
A	-4	PHE	-	CLONING ARTIFACT	UNP O35014
A	-3	GLN	-	CLONING ARTIFACT	UNP O35014
A	-2	SER	-	CLONING ARTIFACT	UNP O35014
A	-1	ASN	-	CLONING ARTIFACT	UNP O35014
A	0	ALA	-	CLONING ARTIFACT	UNP O35014
B	-23	MSE	-	INITIATING METHIONINE	UNP O35014

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-22	HIS	-	EXPRESSION TAG	UNP O35014
B	-21	HIS	-	EXPRESSION TAG	UNP O35014
B	-20	HIS	-	EXPRESSION TAG	UNP O35014
B	-19	HIS	-	EXPRESSION TAG	UNP O35014
B	-18	HIS	-	EXPRESSION TAG	UNP O35014
B	-17	HIS	-	EXPRESSION TAG	UNP O35014
B	-16	SER	-	CLONING ARTIFACT	UNP O35014
B	-15	SER	-	CLONING ARTIFACT	UNP O35014
B	-14	GLY	-	CLONING ARTIFACT	UNP O35014
B	-13	VAL	-	CLONING ARTIFACT	UNP O35014
B	-12	ASP	-	CLONING ARTIFACT	UNP O35014
B	-11	LEU	-	CLONING ARTIFACT	UNP O35014
B	-10	GLY	-	CLONING ARTIFACT	UNP O35014
B	-9	THR	-	CLONING ARTIFACT	UNP O35014
B	-8	GLU	-	CLONING ARTIFACT	UNP O35014
B	-7	ASN	-	CLONING ARTIFACT	UNP O35014
B	-6	LEU	-	CLONING ARTIFACT	UNP O35014
B	-5	TYR	-	CLONING ARTIFACT	UNP O35014
B	-4	PHE	-	CLONING ARTIFACT	UNP O35014
B	-3	GLN	-	CLONING ARTIFACT	UNP O35014
B	-2	SER	-	CLONING ARTIFACT	UNP O35014
B	-1	ASN	-	CLONING ARTIFACT	UNP O35014
B	0	ALA	-	CLONING ARTIFACT	UNP O35014

- Molecule 2 is BETA-MERCAPTOETHANOL (three-letter code: BME) (formula:  $C_2H_6OS$ ).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	O	S	0	0
			4	2	1	1		

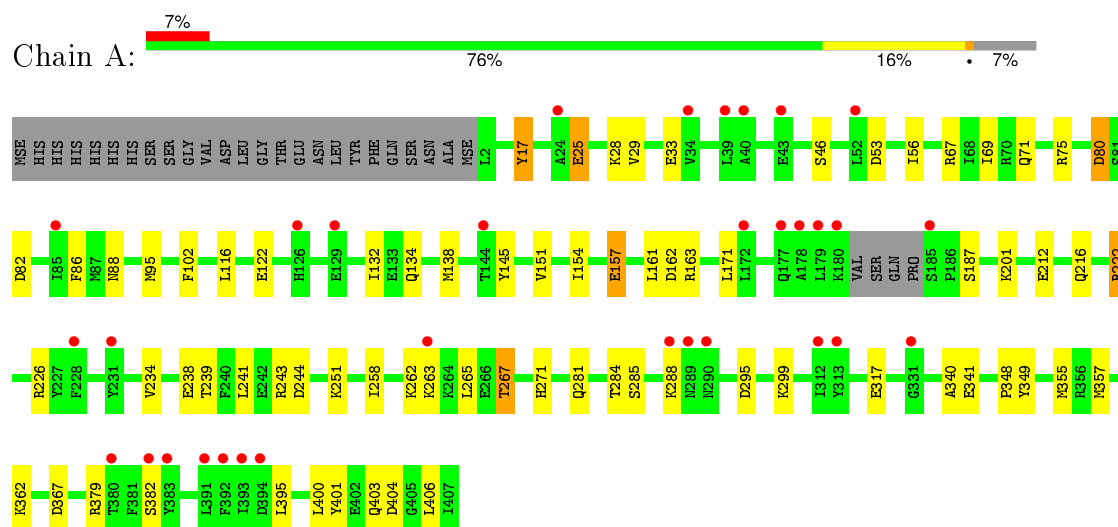
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	148	Total	O	0	0
			148	148		
3	B	107	Total	O	0	0
			107	107		

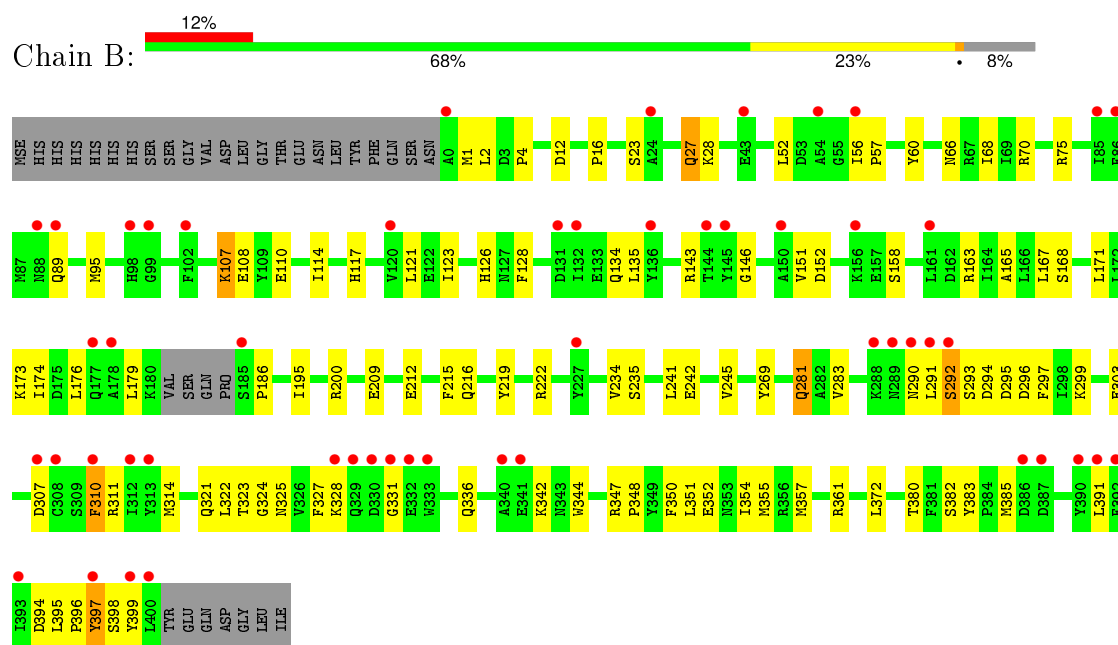
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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: YkuI protein



#### • Molecule 1: YkuI protein



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	46.34Å 125.27Å 167.96Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	24.04 – 2.61 24.01 – 2.61	Depositor EDS
% Data completeness (in resolution range)	100.0 (24.04-2.61) 98.1 (24.01-2.61)	Depositor EDS
$R_{merge}$	0.04	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.33 (at 2.60Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, $R_{free}$	0.203 , 0.287 0.202 , 0.282	Depositor DCC
$R_{free}$ test set	1505 reflections (5.26%)	DCC
Wilson B-factor (Å <sup>2</sup> )	64.9	Xtriage
Anisotropy	0.129	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.28 , 65.9	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.29$	Xtriage
Outliers	0 of 30096 reflections	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	6915	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	65.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.71% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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 [i](#)

### 5.1 Standard geometry [i](#)

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

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	0.51	2/3418 (0.1%)	0.63	0/4601
1	B	0.68	7/3370 (0.2%)	0.72	3/4535 (0.1%)
All	All	0.60	9/6788 (0.1%)	0.67	3/9136 (0.0%)

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	295	ASP	CG-OD1	13.20	1.55	1.25
1	B	296	ASP	CG-OD1	10.49	1.49	1.25
1	B	107	LYS	CE-NZ	8.46	1.70	1.49
1	B	296	ASP	CG-OD2	7.78	1.43	1.25
1	B	292	SER	C-O	6.96	1.36	1.23
1	B	292	SER	C-N	6.23	1.48	1.34
1	A	244	ASP	CG-OD1	5.64	1.38	1.25
1	B	303	GLU	CD-OE2	5.47	1.31	1.25
1	A	244	ASP	CG-OD2	5.39	1.37	1.25

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	296	ASP	CB-CG-OD2	-17.09	102.92	118.30
1	B	296	ASP	CB-CG-OD1	6.87	124.48	118.30
1	B	295	ASP	CB-CG-OD1	-5.86	113.02	118.30

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	3351	0	3244	42	0
1	B	3305	0	3210	66	0
2	A	4	0	5	0	0
3	A	148	0	0	1	0
3	B	107	0	0	3	0
All	All	6915	0	6459	98	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 7.

All (98) 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:107:LYS:CE	1:B:107:LYS:NZ	1.70	1.53
1:B:385:MSE:HE2	1:B:391:LEU:HB2	1.48	0.93
1:A:263:LYS:O	1:A:267:THR:HG22	1.71	0.91
1:B:212:GLU:H	1:B:216:GLN:NE2	1.69	0.89
1:B:321:GLN:HE21	1:B:344:TRP:HE1	1.26	0.80
1:B:242:GLU:O	1:B:245:VAL:HG23	1.81	0.80
1:B:152:ASP:OD1	1:B:173:LYS:HD3	1.84	0.77
1:B:212:GLU:H	1:B:216:GLN:HE22	1.35	0.72
1:B:328:LYS:HE2	1:B:331:GLY:HA2	1.75	0.69
1:B:66:ASN:O	1:B:70:ARG:HG3	1.94	0.68
1:A:379:ARG:NH1	1:A:406:LEU:O	2.27	0.67
1:A:340:ALA:O	1:A:341:GLU:HB2	1.97	0.65
1:A:161:LEU:CD1	1:B:195:ILE:HG13	2.28	0.64
1:A:234:VAL:HG11	1:A:241:LEU:HD11	1.78	0.64
1:B:350:PHE:CE1	1:B:354:ILE:HD11	2.33	0.63
1:B:70:ARG:NH2	1:B:108:GLU:OE2	2.34	0.60
1:B:357:MSE:SE	1:B:382:SER:HB3	2.50	0.60
1:B:143:ARG:NH2	1:B:168:SER:O	2.34	0.60
1:B:327:PHE:HB2	1:B:336:GLN:NE2	2.17	0.60
1:A:95:MSE:HE2	1:A:134:GLN:HG3	1.82	0.60
1:A:271:HIS:HE1	1:A:403:GLN:HE21	1.51	0.59
1:B:126:HIS:HB2	1:B:158:SER:HB2	1.84	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:311:ARG:NH1	1:B:394:ASP:OD2	2.32	0.56
1:B:117:HIS:HA	1:B:146:GLY:O	2.07	0.55
1:A:262:LYS:HD3	1:B:351:LEU:HD22	1.90	0.54
1:B:350:PHE:HE1	1:B:354:ILE:HD11	1.72	0.54
1:A:348:PRO:O	1:A:349:TYR:HB2	2.08	0.53
1:A:212:GLU:H	1:A:216:GLN:NE2	2.06	0.53
1:B:57:PRO:HG2	1:B:60:TYR:CD1	2.43	0.53
1:A:116:LEU:HD13	1:A:145:TYR:HB3	1.89	0.53
1:A:25:GLU:HB3	1:B:342:LYS:HB3	1.90	0.53
1:A:82:ASP:CG	1:A:243:ARG:HH12	2.12	0.53
1:B:57:PRO:HG2	1:B:60:TYR:HD1	1.74	0.52
1:B:215:PHE:O	1:B:219:TYR:HD1	1.92	0.52
1:A:357:MSE:SE	1:A:382:SER:HB3	2.59	0.52
1:B:173:LYS:HE3	1:B:209:GLU:OE1	2.10	0.52
1:B:110:GLU:HA	1:B:114:ILE:O	2.10	0.52
1:B:89:GLN:HG3	1:B:121:LEU:HD11	1.92	0.52
1:B:4:PRO:HB3	1:B:56:ILE:HD12	1.92	0.51
1:B:344:TRP:CE3	1:B:347:ARG:HD2	2.46	0.51
1:A:95:MSE:SE	1:A:138:MSE:HG2	2.61	0.51
1:B:128:PHE:CZ	1:B:135:LEU:HB2	2.46	0.50
1:B:242:GLU:O	1:B:245:VAL:CG2	2.57	0.50
1:A:281:GLN:O	1:A:285:SER:HB3	2.11	0.50
1:A:271:HIS:HE1	1:A:403:GLN:NE2	2.09	0.50
1:B:397:TYR:HB3	1:B:399:TYR:HB2	1.94	0.50
1:A:258:ILE:O	1:A:262:LYS:HG2	2.13	0.49
1:B:107:LYS:NZ	3:B:437:HOH:O	2.46	0.49
1:A:53:ASP:HB3	1:A:56:ILE:HD12	1.93	0.49
1:B:174:ILE:HD11	1:B:195:ILE:HG21	1.95	0.48
1:A:284:THR:O	1:A:288:LYS:HB2	2.13	0.48
1:B:1:MSE:HG2	1:B:2:LEU:H	1.77	0.48
1:B:294:ASP:HB3	1:B:314:MSE:CE	2.43	0.48
1:B:241:LEU:HB3	1:B:245:VAL:HG21	1.96	0.48
1:A:151:VAL:HG12	1:A:154:ILE:HD11	1.95	0.48
1:A:17:TYR:CE1	1:A:238:GLU:HB3	2.48	0.47
1:A:33:GLU:HA	1:A:86:PHE:HB2	1.96	0.47
1:A:88:ASN:HA	1:A:122:GLU:HB2	1.96	0.47
1:A:67:ARG:HG3	1:A:71:GLN:HE21	1.79	0.47
1:B:310:PHE:HB2	1:B:396:PRO:HD3	1.96	0.46
1:B:151:VAL:HG13	3:B:433:HOH:O	2.15	0.46
1:B:383:TYR:HB3	1:B:391:LEU:HB3	1.97	0.46
1:A:82:ASP:OD2	1:A:243:ARG:NH1	2.49	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:234:VAL:HG11	1:B:241:LEU:HD11	1.98	0.46
1:A:161:LEU:HD13	1:B:195:ILE:HG13	1.96	0.46
1:A:161:LEU:HD12	1:B:195:ILE:HG13	1.98	0.46
1:B:291:LEU:HD13	1:B:297:PHE:HA	1.99	0.45
1:B:16:PRO:HB3	1:B:68:ILE:HG23	1.98	0.45
1:A:157:GLU:H	1:A:157:GLU:HG2	1.27	0.45
1:A:151:VAL:CG1	1:A:154:ILE:HD11	2.46	0.45
1:A:132:ILE:HG12	1:A:163:ARG:NH1	2.33	0.44
1:B:281:GLN:HB3	1:B:281:GLN:HE21	1.61	0.44
1:B:23:SER:O	1:B:27:GLN:HA	2.18	0.44
1:A:222:ARG:HB2	3:A:565:HOH:O	2.17	0.43
1:A:75:ARG:HH12	1:A:239:THR:HG23	1.84	0.43
1:B:347:ARG:HA	1:B:348:PRO:HD3	1.93	0.43
1:A:80:ASP:N	1:A:80:ASP:OD2	2.52	0.43
1:B:292:SER:OG	1:B:293:SER:N	2.52	0.43
1:A:28:LYS:HB2	1:A:243:ARG:O	2.19	0.43
1:A:226:ARG:HH12	1:B:372:LEU:HB3	1.84	0.43
1:B:325:ASN:HB2	1:B:336:GLN:HB2	2.00	0.42
1:B:163:ARG:O	1:B:167:LEU:HG	2.19	0.42
1:A:295:ASP:O	1:A:299:LYS:HG3	2.20	0.42
1:B:95:MSE:HE2	1:B:134:GLN:HG2	2.02	0.42
1:B:234:VAL:HG12	1:B:235:SER:O	2.20	0.41
1:A:367:ASP:OD1	1:B:200:ARG:NH2	2.52	0.41
1:A:69:ILE:HD13	1:A:102:PHE:HE1	1.86	0.41
1:A:355:MSE:HE1	1:B:269:TYR:HB2	2.02	0.41
1:B:283:VAL:HG22	1:B:297:PHE:HZ	1.86	0.41
1:B:176:LEU:O	1:B:179:LEU:HB3	2.20	0.41
1:B:310:PHE:HB3	1:B:394:ASP:O	2.21	0.41
1:B:395:LEU:HD23	1:B:396:PRO:HD2	2.03	0.41
1:A:187:SER:HB3	1:B:165:ALA:HB1	2.03	0.41
1:B:297:PHE:CE2	1:B:385:MSE:HE1	2.55	0.41
1:B:28:LYS:HA	3:B:439:HOH:O	2.21	0.41
1:B:107:LYS:CD	1:B:107:LYS:NZ	2.71	0.40
1:A:265:LEU:CD2	1:B:352:GLU:HG3	2.51	0.40
1:B:323:THR:HG22	1:B:324:GLY:O	2.21	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	398/431 (92%)	379 (95%)	19 (5%)	0	100	100
1	B	393/431 (91%)	365 (93%)	23 (6%)	5 (1%)	15	29
All	All	791/862 (92%)	744 (94%)	42 (5%)	5 (1%)	30	54

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	307	ASP
1	B	310	PHE
1	B	186	PRO
1	B	397	TYR
1	B	290	ASN

### 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	363/378 (96%)	345 (95%)	18 (5%)	30	55
1	B	358/378 (95%)	344 (96%)	14 (4%)	39	67
All	All	721/756 (95%)	689 (96%)	32 (4%)	35	62

All (32) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	17	TYR

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Mol	Chain	Res	Type
1	A	25	GLU
1	A	29	VAL
1	A	46	SER
1	A	80	ASP
1	A	157	GLU
1	A	162	ASP
1	A	171	LEU
1	A	201	LYS
1	A	222	ARG
1	A	251	LYS
1	A	267	THR
1	A	317	GLU
1	A	362	LYS
1	A	395	LEU
1	A	400	LEU
1	A	401	TYR
1	A	404	ASP
1	B	12	ASP
1	B	27	GLN
1	B	52	LEU
1	B	75	ARG
1	B	123	ILE
1	B	171	LEU
1	B	222	ARG
1	B	281	GLN
1	B	299	LYS
1	B	322	LEU
1	B	355	MSE
1	B	361	ARG
1	B	380	THR
1	B	398	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	71	GLN
1	A	134	GLN
1	A	216	GLN
1	A	218	GLN
1	A	256	GLN
1	A	271	HIS
1	A	290	ASN

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Mol	Chain	Res	Type
1	A	403	GLN
1	B	19	GLN
1	B	45	GLN
1	B	89	GLN
1	B	117	HIS
1	B	216	GLN
1	B	218	GLN
1	B	223	ASN
1	B	229	GLN
1	B	248	GLN
1	B	255	HIS
1	B	274	GLN
1	B	281	GLN
1	B	321	GLN
1	B	325	ASN
1	B	336	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 carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

1 ligand is modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. 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  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
2	BME	A	501	1	3,3,3	0.28	0	2,2,2	0.21	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	BME	A	501	1	-	0/1/1/1	0/0/0/0

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 ⓘ

### 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	393/431 (91%)	0.52	32 (8%) 15 10	44, 64, 75, 88	0
1	B	387/431 (89%)	0.79	52 (13%) 4 2	48, 66, 78, 88	0
All	All	780/862 (90%)	0.65	84 (10%) 8 4	44, 65, 77, 88	0

All (84) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	0	ALA	5.8
1	B	391	LEU	5.4
1	A	393	ILE	4.9
1	A	289	ASN	4.7
1	B	400	LEU	4.6
1	B	292	SER	4.5
1	B	330	ASP	4.5
1	B	290	ASN	4.5
1	B	312	ILE	4.4
1	B	313	TYR	4.3
1	A	180	LYS	4.2
1	B	393	ILE	4.1
1	A	43	GLU	4.0
1	B	333	TRP	4.0
1	A	185	SER	3.9
1	B	397	TYR	3.9
1	B	392	PHE	3.9
1	B	98	HIS	3.8
1	A	392	PHE	3.6
1	A	129	GLU	3.6
1	B	329	GLN	3.5
1	B	185	SER	3.5
1	B	289	ASN	3.5
1	A	313	TYR	3.5

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Mol	Chain	Res	Type	RSRZ
1	B	86	PHE	3.3
1	B	291	LEU	3.3
1	B	177	GLN	3.3
1	A	39	LEU	3.3
1	A	52	LEU	3.3
1	A	144	THR	3.3
1	B	132	ILE	3.2
1	B	331	GLY	3.2
1	B	332	GLU	3.2
1	B	43	GLU	3.1
1	B	144	THR	3.1
1	B	54	ALA	3.1
1	B	156	LYS	3.0
1	B	328	LYS	3.0
1	B	340	ALA	2.9
1	B	99	GLY	2.9
1	B	120	VAL	2.9
1	B	89	GLN	2.8
1	A	288	LYS	2.8
1	A	177	GLN	2.7
1	A	312	ILE	2.7
1	A	391	LEU	2.6
1	B	145	TYR	2.6
1	A	179	LEU	2.6
1	B	56	ILE	2.6
1	B	24	ALA	2.6
1	A	172	LEU	2.6
1	B	399	TYR	2.6
1	B	131	ASP	2.5
1	B	387	ASP	2.5
1	A	382	SER	2.5
1	B	341	GLU	2.4
1	A	34	VAL	2.4
1	A	231	TYR	2.4
1	B	307	ASP	2.4
1	B	88	ASN	2.4
1	B	102	PHE	2.4
1	A	331	GLY	2.4
1	B	178	ALA	2.3
1	A	290	ASN	2.3
1	A	40	ALA	2.3
1	A	178	ALA	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	85	ILE	2.3
1	B	227	TYR	2.2
1	B	310	PHE	2.2
1	A	85	ILE	2.2
1	B	288	LYS	2.2
1	B	386	ASP	2.2
1	B	390	TYR	2.2
1	A	383	TYR	2.1
1	B	136	TYR	2.1
1	A	380	THR	2.1
1	A	126	HIS	2.1
1	A	24	ALA	2.1
1	A	263	LYS	2.1
1	B	150	ALA	2.1
1	B	161	LEU	2.1
1	B	308	CYS	2.1
1	A	228	PHE	2.0
1	A	394	ASP	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](#)

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> 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	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
2	BME	A	501	4/4	0.93	0.23	1.86	88,92,92,93	0

## 6.5 Other polymers

There are no such residues in this entry.