



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 03:22 am GMT

PDB ID : 1L2F
Title : Crystal structure of NusA from *Thermotoga maritima*: a structure-based role of the N-terminal domain
Authors : Shin, D.H.; Nguyen, H.H.; Jancarik, J.; Yokota, H.; Kim, R.; Kim, S.H.; Berkeley Structural Genomics Center (BSGC)
Deposited on : 2002-02-20
Resolution : 2.50 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
Xtriage (Phenix)	:	1.9-1692
EDS	:	trunk28620
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
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)	:	recalc28949

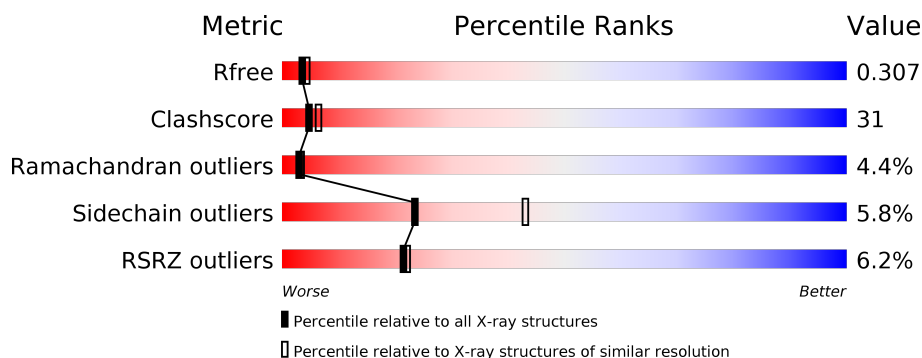
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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}	100719	3846 (2.50-2.50)
Clashscore	112137	4554 (2.50-2.50)
Ramachandran outliers	110173	4463 (2.50-2.50)
Sidechain outliers	110143	4465 (2.50-2.50)
RSRZ outliers	101464	3876 (2.50-2.50)

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	369	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 2886 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 N utilization substance protein A.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	344	Total	C	N	O	S	Se	0	0	0
			2667	1702	456	504	1	4			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MSE	MET	MODIFIED RESIDUE	UNP Q9X298
A	146	MSE	MET	MODIFIED RESIDUE	UNP Q9X298
A	208	MSE	MET	MODIFIED RESIDUE	UNP Q9X298
A	342	MSE	MET	MODIFIED RESIDUE	UNP Q9X298

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	219	Total	O	0	0
			219	219		

4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	116.19Å 116.19Å 64.63Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.05 – 2.50 50.80 – 2.48	Depositor EDS
% Data completeness (in resolution range)	99.3 (29.05-2.50) 98.4 (50.80-2.48)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.61 (at 2.48Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.223 , 0.298 0.235 , 0.307	Depositor DCC
R_{free} test set	1578 reflections (10.04%)	DCC
Wilson B-factor (Å ²)	50.6	Xtriage
Anisotropy	0.430	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 71.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	2886	wwPDB-VP
Average B, all atoms (Å ²)	60.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.32% 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](#)

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	0.56	2/2695 (0.1%)	0.76	3/3631 (0.1%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	289	PRO	N-CD	8.19	1.59	1.47
1	A	208	MSE	SE-CE	-5.48	1.63	1.95

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	289	PRO	CA-N-CD	-7.07	101.60	111.50
1	A	289	PRO	N-CA-C	5.71	126.93	112.10
1	A	273	LEU	CA-CB-CG	5.65	128.29	115.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	2667	0	2844	169	0
2	A	219	0	0	15	2
All	All	2886	0	2844	169	2

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 31.

All (169) 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:288:ALA:HB3	1:A:289:PRO:HD3	1.39	1.02
1:A:86:LEU:HD23	1:A:86:LEU:H	1.21	1.02
1:A:320:LYS:HD2	2:A:2028:HOH:O	1.58	0.99
1:A:263:LYS:HE3	1:A:263:LYS:HA	1.42	0.98
1:A:83:ILE:HD12	1:A:83:ILE:H	1.32	0.95
1:A:288:ALA:CB	1:A:289:PRO:HD3	1.98	0.92
1:A:69:GLU:HB3	1:A:74:GLN:HE22	1.33	0.91
1:A:219:ILE:HG22	1:A:244:VAL:HG21	1.59	0.85
1:A:336:ILE:HD11	1:A:338:ILE:HD11	1.59	0.84
1:A:64:VAL:HA	1:A:75:ILE:HD12	1.58	0.84
1:A:4:GLY:O	1:A:7:GLU:HB3	1.82	0.80
1:A:288:ALA:CB	1:A:289:PRO:CD	2.60	0.79
1:A:288:ALA:HB3	1:A:289:PRO:CD	2.13	0.79
1:A:146:MSE:HG2	1:A:147:GLY:H	1.47	0.77
1:A:19:SER:O	1:A:22:GLU:HG2	1.85	0.77
1:A:107:ALA:O	1:A:111:LYS:HG2	1.85	0.76
1:A:298:LEU:O	1:A:299:ASP:HB2	1.84	0.76
1:A:308:LEU:HB3	1:A:341:ILE:HG23	1.66	0.76
1:A:46:GLU:HB2	1:A:61:LEU:HD11	1.69	0.74
1:A:70:ASP:H	1:A:74:GLN:NE2	1.86	0.74
1:A:159:THR:HG23	2:A:2107:HOH:O	1.87	0.73
1:A:86:LEU:CD2	1:A:86:LEU:H	2.00	0.73
1:A:68:VAL:HA	2:A:2022:HOH:O	1.89	0.71
1:A:228:GLU:HB3	1:A:231:VAL:HG22	1.72	0.69
1:A:48:VAL:HG13	1:A:57:LYS:HB2	1.75	0.68
1:A:232:ARG:HD2	1:A:323:GLN:HE21	1.59	0.68
1:A:24:ILE:HB	1:A:25:PRO:HD3	1.76	0.68
1:A:296:GLU:HB3	1:A:306:ARG:HB3	1.76	0.68
1:A:189:THR:HA	2:A:2183:HOH:O	1.93	0.68
1:A:75:ILE:HD13	1:A:80:ALA:HB2	1.76	0.67
1:A:211:GLU:HG3	2:A:2111:HOH:O	1.93	0.67
1:A:75:ILE:O	1:A:75:ILE:HD12	1.95	0.66
1:A:208:MSE:HE1	1:A:235:VAL:HG21	1.78	0.66
1:A:65:VAL:HG23	1:A:67:GLU:H	1.61	0.65
1:A:190:LYS:O	1:A:190:LYS:HD3	1.97	0.65
1:A:84:ASP:HB3	1:A:85:PRO:HD2	1.80	0.64
1:A:82:LYS:HB2	1:A:83:ILE:HD12	1.79	0.63
1:A:147:GLY:HA3	2:A:2013:HOH:O	1.97	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:146:MSE:HG2	1:A:147:GLY:N	2.14	0.62
1:A:126:GLN:HB2	1:A:192:PRO:HG3	1.81	0.62
1:A:214:GLU:H	1:A:214:GLU:CD	2.02	0.62
1:A:214:GLU:HG3	1:A:249:ALA:HB1	1.81	0.62
1:A:111:LYS:O	1:A:115:ILE:HD13	2.00	0.62
1:A:78:GLU:O	1:A:82:LYS:HG2	1.99	0.62
1:A:231:VAL:HG23	1:A:232:ARG:N	2.15	0.62
1:A:70:ASP:H	1:A:74:GLN:HE21	1.46	0.61
1:A:314:LEU:HD13	1:A:340:PRO:HG3	1.83	0.61
1:A:227:ARG:HD3	1:A:233:THR:OG1	2.00	0.61
1:A:57:LYS:HE3	2:A:2134:HOH:O	2.00	0.60
1:A:27:LEU:HD23	1:A:49:ILE:CD1	2.31	0.60
1:A:84:ASP:HB3	1:A:85:PRO:CD	2.32	0.60
1:A:137:VAL:HG23	1:A:206:GLY:HA3	1.82	0.60
1:A:165:GLU:HG3	1:A:197:SER:CB	2.31	0.59
1:A:32:VAL:O	1:A:36:ARG:HG3	2.02	0.59
1:A:219:ILE:CG2	1:A:244:VAL:HG21	2.31	0.59
1:A:11:GLN:O	1:A:15:GLU:HB2	2.02	0.59
1:A:79:GLU:O	1:A:83:ILE:HD13	2.03	0.59
1:A:64:VAL:HG13	1:A:91:SER:H	1.66	0.59
1:A:207:LEU:HD21	1:A:264:GLU:HG2	1.86	0.58
1:A:100:LYS:O	1:A:101:ASN:HB2	2.01	0.58
1:A:68:VAL:O	1:A:68:VAL:HG12	2.04	0.58
1:A:79:GLU:OE1	1:A:79:GLU:HA	2.03	0.58
1:A:83:ILE:H	1:A:83:ILE:CD1	2.10	0.57
1:A:212:ILE:HD13	1:A:250:CYS:HA	1.86	0.57
1:A:275:TRP:HZ3	1:A:277:ASP:OD2	1.87	0.57
1:A:92:ILE:O	1:A:93:VAL:HG23	2.04	0.56
1:A:308:LEU:HA	1:A:339:LYS:O	2.05	0.56
1:A:99:VAL:O	1:A:99:VAL:HG13	2.06	0.56
1:A:130:TYR:O	1:A:185:VAL:HG21	2.07	0.55
1:A:231:VAL:CG2	1:A:232:ARG:N	2.70	0.54
1:A:298:LEU:HD12	2:A:2166:HOH:O	2.06	0.54
1:A:112:GLN:O	1:A:116:GLN:HG3	2.07	0.54
1:A:244:VAL:HG23	1:A:244:VAL:O	2.08	0.54
1:A:59:TYR:HB3	1:A:94:LYS:HB3	1.89	0.54
1:A:2:ASN:O	1:A:3:ILE:C	2.46	0.54
1:A:312:THR:HG23	2:A:2079:HOH:O	2.07	0.54
1:A:149:TRP:NE1	1:A:160:ARG:HD3	2.23	0.54
1:A:336:ILE:HD11	1:A:338:ILE:CD1	2.36	0.54
1:A:166:TRP:CD1	1:A:170:GLU:HG2	2.43	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:74:GLN:O	1:A:75:ILE:HG23	2.07	0.53
1:A:48:VAL:CG1	1:A:59:TYR:HE2	2.20	0.53
1:A:228:GLU:HB3	1:A:231:VAL:CG2	2.38	0.53
1:A:275:TRP:CZ3	1:A:277:ASP:OD2	2.61	0.53
1:A:86:LEU:HD23	1:A:86:LEU:N	2.05	0.53
1:A:343:ASN:O	1:A:344:LEU:HD23	2.09	0.52
1:A:214:GLU:CG	1:A:249:ALA:HB1	2.38	0.52
1:A:181:TYR:O	1:A:196:VAL:HA	2.10	0.52
1:A:3:ILE:HB	2:A:2039:HOH:O	2.09	0.52
1:A:64:VAL:HA	1:A:75:ILE:CD1	2.35	0.52
1:A:288:ALA:HB1	1:A:289:PRO:HD3	1.91	0.51
1:A:27:LEU:HD23	1:A:49:ILE:HD13	1.91	0.51
1:A:165:GLU:HG3	1:A:197:SER:OG	2.11	0.51
1:A:2:ASN:HB2	2:A:2024:HOH:O	2.11	0.50
1:A:137:VAL:CG2	1:A:206:GLY:HA3	2.40	0.50
1:A:297:ILE:HG13	1:A:297:ILE:O	2.11	0.50
1:A:64:VAL:HG12	1:A:91:SER:O	2.12	0.50
1:A:144:ARG:NH1	2:A:2011:HOH:O	2.45	0.50
1:A:241:ASP:O	1:A:244:VAL:HG22	2.12	0.50
1:A:38:ASN:HB3	1:A:105:ILE:HD13	1.93	0.50
1:A:188:THR:O	1:A:190:LYS:N	2.40	0.49
1:A:308:LEU:HB3	1:A:341:ILE:CG2	2.40	0.49
1:A:204:VAL:HG12	1:A:208:MSE:HE2	1.94	0.48
1:A:232:ARG:HD2	1:A:323:GLN:NE2	2.26	0.48
1:A:213:PRO:HD2	1:A:214:GLU:OE2	2.13	0.48
1:A:38:ASN:CG	1:A:105:ILE:HG21	2.34	0.48
1:A:344:LEU:O	1:A:344:LEU:HG	2.13	0.48
1:A:134:LYS:HE2	1:A:184:ASP:OD1	2.14	0.48
1:A:299:ASP:C	1:A:301:GLU:H	2.16	0.47
1:A:94:LYS:N	1:A:94:LYS:HD3	2.29	0.47
1:A:245:ASP:OD1	1:A:247:ILE:N	2.48	0.47
1:A:341:ILE:HD12	1:A:341:ILE:O	2.15	0.47
1:A:156:LYS:HD3	1:A:156:LYS:H	1.80	0.46
1:A:253:GLU:HG2	1:A:256:SER:OG	2.15	0.46
1:A:163:LYS:HE2	1:A:166:TRP:CZ2	2.50	0.46
1:A:232:ARG:HH11	1:A:323:GLN:HE21	1.64	0.46
1:A:232:ARG:HB3	1:A:327:LEU:CD1	2.46	0.46
1:A:203:PHE:HB3	1:A:227:ARG:NH2	2.30	0.46
1:A:265:LEU:HB3	1:A:268:GLU:HB3	1.97	0.45
1:A:318:ILE:HG22	1:A:322:GLY:HA2	1.98	0.45
1:A:343:ASN:C	1:A:344:LEU:HD23	2.37	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:232:ARG:HB3	1:A:327:LEU:HD11	1.98	0.45
1:A:64:VAL:HG11	1:A:91:SER:HB2	1.97	0.45
1:A:214:GLU:HG3	1:A:249:ALA:CB	2.46	0.45
1:A:203:PHE:HB3	1:A:227:ARG:HH21	1.81	0.45
1:A:277:ASP:O	1:A:279:PRO:HD3	2.16	0.45
1:A:48:VAL:HG11	1:A:59:TYR:HE2	1.80	0.45
1:A:286:ALA:C	1:A:288:ALA:H	2.19	0.45
1:A:64:VAL:CG1	1:A:91:SER:H	2.30	0.45
1:A:245:ASP:OD1	1:A:245:ASP:C	2.56	0.44
1:A:70:ASP:OD1	1:A:72:ALA:HB3	2.17	0.44
1:A:239:SER:C	1:A:241:ASP:H	2.21	0.44
1:A:282:LEU:HD22	1:A:332:THR:HG22	1.99	0.44
1:A:228:GLU:O	1:A:231:VAL:HG22	2.18	0.44
1:A:211:GLU:HG2	1:A:260:ALA:HB3	1.99	0.44
1:A:221:GLU:OE2	1:A:223:LYS:HE3	2.17	0.43
1:A:62:LEU:HB3	1:A:75:ILE:HG13	1.99	0.43
1:A:232:ARG:CD	1:A:323:GLN:HE21	2.30	0.43
1:A:39:PHE:CE2	1:A:102:PHE:HB2	2.53	0.43
1:A:125:LYS:O	1:A:129:LYS:HD3	2.19	0.42
1:A:180:VAL:HB	1:A:196:VAL:HB	2.01	0.42
1:A:251:ILE:HG23	1:A:258:ILE:HD11	2.01	0.42
1:A:288:ALA:HB1	1:A:289:PRO:CD	2.46	0.42
1:A:60:GLN:HE21	1:A:97:LEU:HD21	1.85	0.42
1:A:34:ALA:CB	1:A:109:THR:HG21	2.50	0.42
1:A:5:LEU:C	1:A:5:LEU:HD13	2.40	0.42
1:A:205:ILE:HA	1:A:208:MSE:HE3	2.02	0.42
1:A:187:LYS:HD2	2:A:2069:HOH:O	2.20	0.42
1:A:149:TRP:CD1	1:A:160:ARG:HD3	2.55	0.41
1:A:300:LYS:CB	1:A:300:LYS:NZ	2.83	0.41
1:A:34:ALA:HB2	1:A:109:THR:HG21	2.02	0.41
1:A:156:LYS:HD3	1:A:156:LYS:N	2.35	0.41
1:A:339:LYS:HA	1:A:340:PRO:HD3	1.81	0.41
1:A:153:ARG:HH11	1:A:153:ARG:HG3	1.85	0.41
1:A:86:LEU:O	1:A:86:LEU:HG	2.20	0.41
1:A:292:VAL:HG13	1:A:307:VAL:CG1	2.50	0.41
1:A:159:THR:HG21	1:A:194:ILE:O	2.20	0.41
1:A:200:VAL:HG22	2:A:2197:HOH:O	2.20	0.41
1:A:122:GLU:O	1:A:126:GLN:HG2	2.21	0.41
1:A:188:THR:C	1:A:190:LYS:H	2.23	0.41
1:A:223:LYS:HA	1:A:223:LYS:HE2	2.02	0.41
1:A:64:VAL:HA	1:A:75:ILE:O	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:302:ASN:O	1:A:302:ASN:OD1	2.39	0.41
1:A:83:ILE:HG13	2:A:2080:HOH:O	2.21	0.41
1:A:135:GLY:HA2	1:A:181:TYR:CZ	2.56	0.41
1:A:189:THR:O	1:A:190:LYS:HB2	2.20	0.41
1:A:97:LEU:O	1:A:99:VAL:HG12	2.21	0.40
1:A:163:LYS:HA	1:A:166:TRP:CE3	2.56	0.40
1:A:236:ALA:HA	1:A:273:LEU:HD22	2.03	0.40
1:A:292:VAL:HG13	1:A:307:VAL:HG13	2.03	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:2103:HOH:O	2:A:2103:HOH:O[7_556]	1.79	0.41
2:A:2094:HOH:O	2:A:2094:HOH:O[7_556]	2.11	0.09

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	342/369 (93%)	298 (87%)	29 (8%)	15 (4%)	3 3

All (15) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	3	ILE
1	A	82	LYS
1	A	190	LYS
1	A	288	ALA
1	A	299	ASP
1	A	2	ASN
1	A	85	PRO

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Mol	Chain	Res	Type
1	A	87	ALA
1	A	342	MSE
1	A	7	GLU
1	A	83	ILE
1	A	88	GLU
1	A	99	VAL
1	A	104	ARG
1	A	253	GLU

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	291/310 (94%)	274 (94%)	17 (6%)	23	43

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	37	LYS
1	A	64	VAL
1	A	86	LEU
1	A	109	THR
1	A	121	LEU
1	A	190	LYS
1	A	200	VAL
1	A	243	ASN
1	A	258	ILE
1	A	263	LYS
1	A	273	LEU
1	A	282	LEU
1	A	298	LEU
1	A	320	LYS
1	A	335	LYS
1	A	336	ILE
1	A	343	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	2	ASN
1	A	11	GLN
1	A	38	ASN
1	A	41	ASN
1	A	44	ASN
1	A	52	ASN
1	A	60	GLN
1	A	74	GLN
1	A	217	ASN
1	A	240	ASN
1	A	243	ASN
1	A	302	ASN
1	A	323	GLN

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 carbohydrates 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	340/369 (92%)	0.39	21 (6%)	21 22	28, 53, 101, 130	0

All (21) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	83	ILE	5.1
1	A	343	ASN	4.6
1	A	78	GLU	4.1
1	A	344	LEU	3.6
1	A	35	TYR	2.8
1	A	39	PHE	2.7
1	A	5	LEU	2.7
1	A	14	GLU	2.7
1	A	288	ALA	2.7
1	A	84	ASP	2.6
1	A	69	GLU	2.4
1	A	92	ILE	2.4
1	A	97	LEU	2.4
1	A	2	ASN	2.4
1	A	297	ILE	2.3
1	A	57	LYS	2.2
1	A	111	LYS	2.2
1	A	82	LYS	2.1
1	A	85	PRO	2.1
1	A	37	LYS	2.0
1	A	253	GLU	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.