



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 11:15 AM GMT

PDB ID : 3OKZ  
Title : Crystal Structure of protein gbs0355 from Streptococcus agalactiae, Northeast Structural Genomics Consortium Target SaR127  
Authors : Kuzin, A.; Lew, S.; Vorobiev, S.M.; Seetharaman, J.; Janjua, J.; Xiao, R.; Ciccosanti, C.; Wang, D.; Everett, J.K.; Nair, R.; Acton, T.B.; Rost, B.; Montelione, G.T.; Hunt, J.F.; Tong, L.; Northeast Structural Genomics Consortium (NESG)  
Deposited on : 2010-08-25  
Resolution : 2.70 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

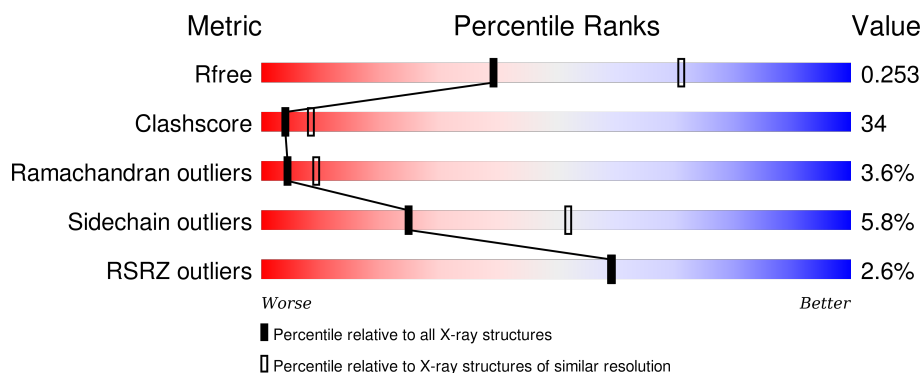
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.70 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	2103 (2.70-2.70)
Clashscore	102246	2422 (2.70-2.70)
Ramachandran outliers	100387	2382 (2.70-2.70)
Sidechain outliers	100360	2382 (2.70-2.70)
RSRZ outliers	91569	2107 (2.70-2.70)

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	306	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red, orange, yellow, green, grey);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> <span>51%</span> <span>42%</span> <span>• •</span> </div> </div>
1	B	306	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red, orange, yellow, green, grey);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> <span>40%</span> <span>51%</span> <span>6% •</span> </div> </div>

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4695 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 Putative uncharacterized protein gbs0355.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	297	Total	C	N	O	Se	0	0	0
			2327	1462	404	452	9			
1	B	295	Total	C	N	O	Se	0	0	0
			2315	1456	402	448	9			

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	41	MSE	-	expression tag	UNP Q8E703
A	339	LEU	-	expression tag	UNP Q8E703
A	340	GLU	-	expression tag	UNP Q8E703
A	341	HIS	-	expression tag	UNP Q8E703
A	342	HIS	-	expression tag	UNP Q8E703
A	343	HIS	-	expression tag	UNP Q8E703
A	344	HIS	-	expression tag	UNP Q8E703
A	345	HIS	-	expression tag	UNP Q8E703
A	346	HIS	-	expression tag	UNP Q8E703
B	41	MSE	-	expression tag	UNP Q8E703
B	339	LEU	-	expression tag	UNP Q8E703
B	340	GLU	-	expression tag	UNP Q8E703
B	341	HIS	-	expression tag	UNP Q8E703
B	342	HIS	-	expression tag	UNP Q8E703
B	343	HIS	-	expression tag	UNP Q8E703
B	344	HIS	-	expression tag	UNP Q8E703
B	345	HIS	-	expression tag	UNP Q8E703
B	346	HIS	-	expression tag	UNP Q8E703

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	37	Total	O	0	0
			37	37		

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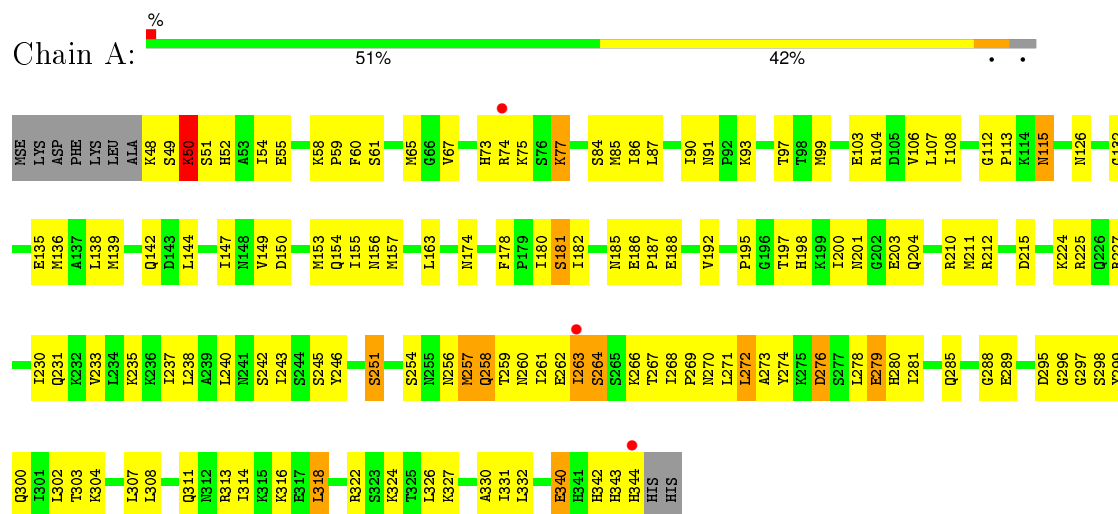
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	16	Total	O	0	0
			16	16		

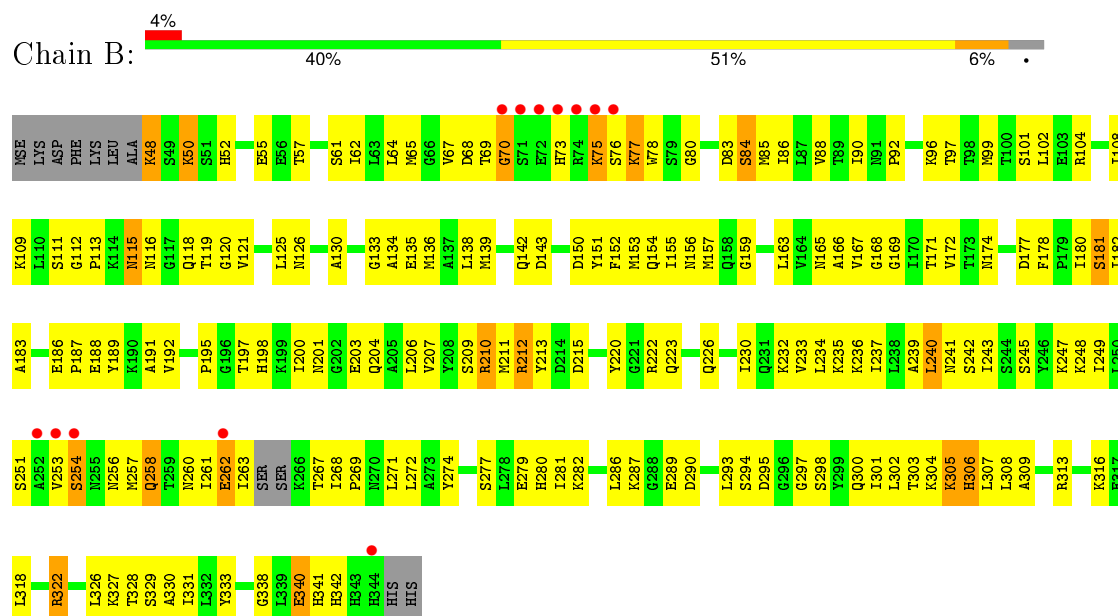
### 3 Residue-property plots

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: Putative uncharacterized protein gbs0355



- Molecule 1: Putative uncharacterized protein gbs0355



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	99.05Å 41.39Å 161.51Å 90.00° 103.44° 90.00°	Depositor
Resolution (Å)	20.00 – 2.70 29.66 – 2.71	Depositor EDS
% Data completeness (in resolution range)	79.7 (20.00-2.70) 91.6 (29.66-2.71)	Depositor EDS
$R_{merge}$	0.10	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.58 (at 2.72Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, $R_{free}$	0.204 , 0.246 0.218 , 0.253	Depositor DCC
$R_{free}$ test set	799 reflections (4.93%)	DCC
Wilson B-factor (Å <sup>2</sup> )	37.7	Xtriage
Anisotropy	0.230	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.30 , 56.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 17016 reflections	Xtriage
$F_o, F_c$ correlation	0.91	EDS
Total number of atoms	4695	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.94% 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](#)

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	0.37	0/2353	0.64	0/3153
1	B	0.32	0/2340	0.58	0/3134
All	All	0.35	0/4693	0.61	0/6287

There are no bond length outliers.

There are no bond angle outliers.

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	2327	0	2375	140	0
1	B	2315	0	2364	180	0
2	A	37	0	0	3	0
2	B	16	0	0	1	0
All	All	4695	0	4739	319	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 34.

All (319) 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:304:LYS:HG2	1:A:332:LEU:HD13	1.33	1.09

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:99:MSE:HE1	1:A:230:ILE:HG21	1.45	0.98
1:B:268:ILE:HA	1:B:271:LEU:HD12	1.43	0.97
1:B:240:LEU:HD13	1:B:245:SER:HB3	1.48	0.94
1:A:257:MSE:HE1	1:A:259:THR:HG22	1.52	0.91
1:B:109:LYS:HE3	1:B:120:GLY:HA2	1.53	0.90
1:A:156:ASN:HD21	1:A:256:ASN:HD22	1.18	0.90
1:B:156:ASN:HD21	1:B:256:ASN:HD22	0.95	0.88
1:B:156:ASN:HD21	1:B:256:ASN:ND2	1.71	0.87
1:B:113:PRO:HD2	1:B:136:MSE:HE3	1.57	0.86
1:B:52:HIS:HA	1:B:55:GLU:OE1	1.75	0.86
1:B:210:ARG:HB3	1:B:210:ARG:HH11	1.43	0.84
1:B:181:SER:HB2	1:B:192:VAL:HG12	1.61	0.82
1:B:75:LYS:HZ2	1:B:75:LYS:HB3	1.44	0.81
1:B:156:ASN:ND2	1:B:256:ASN:HD22	1.77	0.80
1:B:316:LYS:HA	1:B:322:ARG:HH21	1.47	0.80
1:B:135:GLU:O	1:B:139:MSE:HG3	1.81	0.80
1:A:157:MSE:HG3	2:A:427:HOH:O	1.81	0.79
1:A:156:ASN:ND2	1:A:256:ASN:HD22	1.81	0.78
1:A:251:SER:HA	1:A:254:SER:HB2	1.65	0.78
1:A:316:LYS:HG3	1:A:322:ARG:HH21	1.48	0.78
1:A:261:ILE:HG22	1:A:263:ILE:H	1.48	0.78
1:A:240:LEU:HD13	1:A:245:SER:HB3	1.68	0.76
1:B:99:MSE:HE1	1:B:230:ILE:HG21	1.67	0.75
1:A:267:THR:HG22	1:A:271:LEU:HD12	1.68	0.75
1:A:156:ASN:HD21	1:A:256:ASN:ND2	1.85	0.74
1:A:243:ILE:HG12	1:A:268:ILE:HD12	1.70	0.74
1:B:302:LEU:HD13	1:B:307:LEU:HD13	1.69	0.74
1:B:313:ARG:HH11	1:B:313:ARG:HG3	1.53	0.74
1:B:316:LYS:HG3	1:B:322:ARG:HH21	1.52	0.73
1:A:174:ASN:HD22	1:A:195:PRO:HA	1.52	0.73
1:A:65:MSE:HE3	1:A:85:MSE:SE	2.39	0.73
1:B:104:ARG:HB2	1:B:126:ASN:HB2	1.69	0.73
1:A:304:LYS:CG	1:A:332:LEU:HD13	2.14	0.72
1:B:108:ILE:HG13	1:B:330:ALA:HA	1.69	0.72
1:A:155:ILE:HG12	1:A:257:MSE:HB2	1.70	0.72
1:B:167:VAL:HB	1:B:236:LYS:HG2	1.73	0.71
1:B:113:PRO:HG2	1:B:116:ASN:HB2	1.72	0.71
1:B:61:SER:HB3	1:B:318:LEU:HD11	1.73	0.71
1:A:238:LEU:HD11	1:A:278:LEU:HB2	1.73	0.70
1:A:326:LEU:HD12	1:A:327:LYS:H	1.57	0.70
1:A:201:ASN:H	1:A:204:GLN:HE21	1.39	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:316:LYS:HG3	1:B:322:ARG:NH2	2.06	0.69
1:B:268:ILE:HG22	1:B:269:PRO:HD3	1.74	0.69
1:B:112:GLY:HA3	1:B:136:MSE:HE2	1.73	0.69
1:A:227:ARG:NH2	1:A:285:GLN:HB2	2.08	0.69
1:B:236:LYS:O	1:B:239:ALA:HB3	1.93	0.68
1:B:174:ASN:HB3	1:B:195:PRO:HA	1.74	0.68
1:A:331:ILE:N	1:A:331:ILE:HD12	2.10	0.67
1:A:298:SER:HB3	1:B:298:SER:HB3	1.75	0.67
1:B:322:ARG:H	1:B:322:ARG:HD2	1.60	0.66
1:B:322:ARG:HD2	1:B:322:ARG:N	2.10	0.66
1:A:60:PHE:CZ	1:A:90:ILE:HD13	2.30	0.66
1:A:233:VAL:O	1:A:237:ILE:HG13	1.96	0.66
1:B:340:GLU:O	1:B:341:HIS:HB2	1.95	0.65
1:B:73:HIS:O	1:B:77:LYS:HD2	1.95	0.65
1:B:50:LYS:HD2	1:B:50:LYS:H	1.62	0.65
1:B:57:THR:O	1:B:57:THR:HG22	1.97	0.65
1:B:268:ILE:O	1:B:272:LEU:HD13	1.98	0.64
1:A:142:GLN:HG2	1:A:149:VAL:HG23	1.79	0.64
1:B:115:ASN:C	1:B:115:ASN:HD22	2.01	0.64
1:B:316:LYS:HA	1:B:322:ARG:NH2	2.11	0.64
1:A:326:LEU:HD12	1:A:327:LYS:N	2.13	0.64
1:B:64:LEU:HD23	1:B:65:MSE:N	2.13	0.64
1:A:188:GLU:H	1:A:188:GLU:CD	2.01	0.63
1:B:268:ILE:CG2	1:B:269:PRO:HD3	2.27	0.62
1:B:235:LYS:NZ	1:B:235:LYS:HB3	2.13	0.62
1:A:86:ILE:HD12	1:A:86:ILE:N	2.14	0.62
1:A:201:ASN:OD1	1:A:204:GLN:HG3	1.97	0.62
1:A:322:ARG:N	1:A:322:ARG:HD2	2.14	0.62
1:B:171:THR:HG23	1:B:198:HIS:O	1.99	0.62
1:B:340:GLU:HB2	1:B:342:HIS:CD2	2.35	0.62
1:A:263:ILE:O	1:A:263:ILE:HG22	1.99	0.61
1:B:97:THR:HB	1:B:281:ILE:HD12	1.80	0.61
1:A:257:MSE:HE1	1:A:259:THR:CG2	2.27	0.61
1:B:308:LEU:HD23	1:B:326:LEU:HB2	1.82	0.61
1:A:307:LEU:O	1:A:311:GLN:HG3	1.99	0.61
1:B:167:VAL:HB	1:B:236:LYS:CG	2.31	0.61
1:B:280:HIS:C	1:B:281:ILE:HD13	2.21	0.61
1:B:67:VAL:O	1:B:156:ASN:HA	2.00	0.60
1:A:211:MSE:HB3	1:A:225:ARG:NH1	2.16	0.60
1:A:113:PRO:HD2	1:A:136:MSE:HE3	1.82	0.60
1:A:316:LYS:HG3	1:A:322:ARG:NH2	2.16	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:50:LYS:N	1:B:50:LYS:HD2	2.16	0.60
1:B:174:ASN:ND2	1:B:180:ILE:HG13	2.17	0.60
1:A:174:ASN:ND2	1:A:195:PRO:HA	2.16	0.59
1:B:90:ILE:N	1:B:90:ILE:HD12	2.17	0.59
1:A:262:GLU:O	1:A:264:SER:N	2.36	0.59
1:A:174:ASN:HD22	1:A:195:PRO:CA	2.16	0.59
1:A:201:ASN:ND2	1:A:203:GLU:H	2.01	0.59
1:B:80:GLY:O	1:B:130:ALA:HA	2.03	0.59
1:B:293:LEU:C	1:B:295:ASP:H	2.07	0.58
1:A:316:LYS:CG	1:A:322:ARG:HH21	2.16	0.58
1:A:242:SER:HB3	1:A:245:SER:HB2	1.84	0.58
1:B:174:ASN:HD21	1:B:180:ILE:HG13	1.69	0.58
1:B:48:LYS:HD3	1:B:48:LYS:N	2.19	0.58
1:B:83:ASP:O	1:B:84:SER:HB2	2.03	0.58
1:A:201:ASN:HD21	1:A:203:GLU:HB2	1.67	0.58
1:B:62:ILE:HA	1:B:151:TYR:O	2.04	0.58
1:B:211:MSE:HE1	1:B:215:ASP:HB2	1.86	0.57
1:B:83:ASP:HA	1:B:104:ARG:HH22	1.70	0.57
1:B:50:LYS:CD	1:B:50:LYS:H	2.18	0.57
1:B:83:ASP:HA	1:B:104:ARG:NH2	2.20	0.57
1:B:116:ASN:HB3	1:B:119:THR:OG1	2.05	0.57
1:A:67:VAL:O	1:A:156:ASN:HA	2.04	0.57
1:B:268:ILE:CA	1:B:271:LEU:HD12	2.27	0.57
1:A:87:LEU:HD12	1:A:314:ILE:HD13	1.87	0.57
1:B:268:ILE:HG13	1:B:272:LEU:CD1	2.35	0.56
1:A:276:ASP:O	1:A:279:GLU:HB2	2.03	0.56
1:B:99:MSE:HE1	1:B:230:ILE:CG2	2.34	0.56
1:A:52:HIS:HA	1:A:55:GLU:OE1	2.06	0.56
1:A:268:ILE:HG12	1:A:269:PRO:HD3	1.87	0.56
1:B:263:ILE:HA	1:B:267:THR:CB	2.36	0.56
1:B:169:GLY:O	1:B:236:LYS:HE3	2.06	0.55
1:B:209:SER:HB2	1:B:226:GLN:HG2	1.88	0.55
1:A:49:SER:C	1:A:51:SER:H	2.10	0.55
1:B:86:ILE:HD12	1:B:86:ILE:N	2.22	0.55
1:B:201:ASN:OD1	1:B:204:GLN:HG3	2.07	0.55
1:A:58:LYS:CD	1:A:59:PRO:HD2	2.36	0.55
1:B:268:ILE:HA	1:B:271:LEU:CD1	2.29	0.55
1:A:201:ASN:HD22	1:A:203:GLU:H	1.54	0.55
1:B:293:LEU:O	1:B:295:ASP:N	2.38	0.55
1:B:108:ILE:HA	1:B:331:ILE:HD13	1.89	0.55
1:B:331:ILE:N	1:B:331:ILE:HD12	2.22	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:186:GLU:HB3	1:A:188:GLU:OE1	2.07	0.55
1:A:243:ILE:HA	1:A:246:TYR:CD2	2.42	0.55
1:B:108:ILE:HD11	1:B:329:SER:OG	2.07	0.55
1:A:201:ASN:HD21	1:A:203:GLU:CB	2.20	0.54
1:B:254:SER:O	1:B:257:MSE:HB3	2.08	0.54
1:B:90:ILE:HG21	1:B:274:TYR:HB3	1.87	0.54
1:B:191:ALA:O	1:B:207:VAL:HG21	2.08	0.54
1:B:313:ARG:NH1	1:B:313:ARG:HG3	2.19	0.54
1:A:65:MSE:HE3	1:A:85:MSE:CG	2.38	0.54
1:A:331:ILE:CD1	1:A:331:ILE:N	2.71	0.54
1:A:108:ILE:HG13	1:A:330:ALA:HA	1.90	0.54
1:B:243:ILE:O	1:B:247:LYS:HG3	2.08	0.54
1:A:262:GLU:C	1:A:264:SER:H	2.12	0.54
1:B:257:MSE:SE	1:B:258:GLN:N	2.91	0.53
1:A:243:ILE:HD13	1:A:272:LEU:HD11	1.88	0.53
1:A:224:LYS:HD2	1:A:227:ARG:NH1	2.22	0.53
1:B:75:LYS:O	1:B:75:LYS:HD2	2.08	0.53
1:B:316:LYS:CG	1:B:322:ARG:HH21	2.20	0.53
1:B:73:HIS:HB2	1:B:77:LYS:O	2.09	0.53
1:A:181:SER:HB2	1:A:192:VAL:HG12	1.89	0.53
1:A:261:ILE:HG22	1:A:263:ILE:N	2.21	0.53
1:A:112:GLY:HA3	1:A:136:MSE:CE	2.39	0.53
1:A:108:ILE:HG13	1:A:330:ALA:HB2	1.91	0.53
1:A:211:MSE:HE1	1:A:215:ASP:HB2	1.91	0.52
1:A:308:LEU:HD23	1:A:326:LEU:HB2	1.91	0.52
1:A:200:ILE:HB	1:A:204:GLN:HB2	1.89	0.52
1:A:108:ILE:HG13	1:A:330:ALA:CA	2.40	0.52
1:B:115:ASN:C	1:B:115:ASN:ND2	2.61	0.52
1:B:68:ASP:O	1:B:70:GLY:N	2.39	0.52
1:B:257:MSE:SE	1:B:258:GLN:H	2.43	0.52
1:A:261:ILE:HG22	1:A:262:GLU:N	2.25	0.51
1:A:201:ASN:ND2	1:A:203:GLU:N	2.59	0.51
1:B:109:LYS:HE3	1:B:120:GLY:CA	2.33	0.51
1:A:103:GLU:HB3	1:A:106:VAL:HG23	1.91	0.51
1:A:54:ILE:HA	1:A:274:TYR:OH	2.11	0.51
1:A:316:LYS:HD2	2:A:432:HOH:O	2.10	0.51
1:A:200:ILE:HA	1:A:204:GLN:NE2	2.26	0.51
1:B:290:ASP:OD1	1:B:300:GLN:NE2	2.44	0.51
1:A:84:SER:OG	1:A:86:ILE:HD11	2.11	0.51
1:B:182:ILE:HD11	1:B:189:TYR:HB2	1.93	0.51
1:A:135:GLU:O	1:A:139:MSE:HG3	2.11	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:316:LYS:CA	1:B:322:ARG:HH21	2.20	0.50
1:A:112:GLY:HA3	1:A:136:MSE:HE2	1.92	0.50
1:A:211:MSE:HB3	1:A:225:ARG:HH12	1.76	0.50
1:A:50:LYS:N	1:A:50:LYS:HE2	2.26	0.50
1:B:85:MSE:HE1	1:B:125:LEU:HG	1.93	0.50
1:A:316:LYS:HA	1:A:322:ARG:NH2	2.27	0.50
1:B:240:LEU:CD1	1:B:249:ILE:HD11	2.42	0.50
1:B:262:GLU:O	1:B:263:ILE:HB	2.11	0.50
1:B:261:ILE:HG22	1:B:262:GLU:N	2.27	0.50
1:A:267:THR:HG22	1:A:271:LEU:CD1	2.38	0.49
1:B:154:GLN:HG3	1:B:258:GLN:HB3	1.94	0.49
1:A:316:LYS:HA	1:A:322:ARG:HH21	1.78	0.49
1:B:182:ILE:HG23	1:B:207:VAL:HG13	1.94	0.49
1:B:109:LYS:CE	1:B:120:GLY:HA2	2.35	0.49
1:A:154:GLN:HE21	1:A:258:GLN:NE2	2.10	0.49
1:B:112:GLY:HA3	1:B:136:MSE:CE	2.41	0.49
1:B:75:LYS:NZ	1:B:75:LYS:HB3	2.22	0.49
1:B:263:ILE:HA	1:B:267:THR:HG21	1.95	0.49
1:A:61:SER:O	1:A:150:ASP:HB2	2.11	0.49
1:B:143:ASP:HB3	1:B:329:SER:HB3	1.95	0.48
1:B:92:PRO:HD3	1:B:274:TYR:CE2	2.47	0.48
1:B:212:ARG:HB2	1:B:222:ARG:HD2	1.94	0.48
1:B:239:ALA:O	1:B:241:ASN:N	2.46	0.48
1:B:108:ILE:HG13	1:B:330:ALA:CA	2.41	0.48
1:B:295:ASP:C	1:B:297:GLY:H	2.16	0.48
1:A:313:ARG:HH11	1:A:313:ARG:HG3	1.77	0.48
1:B:277:SER:C	1:B:279:GLU:H	2.17	0.48
1:B:111:SER:O	1:B:136:MSE:HG3	2.13	0.48
1:B:268:ILE:N	1:B:269:PRO:CD	2.77	0.47
1:B:61:SER:HA	1:B:88:VAL:O	2.13	0.47
1:A:108:ILE:HG13	1:A:330:ALA:CB	2.44	0.47
1:B:166:ALA:C	1:B:168:GLY:H	2.17	0.47
1:B:243:ILE:CD1	1:B:268:ILE:HG12	2.44	0.47
1:A:163:LEU:HG	1:A:233:VAL:HG11	1.96	0.47
1:B:243:ILE:HD11	1:B:268:ILE:HG21	1.97	0.47
1:A:270:ASN:O	1:A:273:ALA:HB3	2.15	0.47
1:B:96:LYS:HD2	1:B:282:LYS:HD2	1.96	0.47
1:B:163:LEU:HG	1:B:233:VAL:HG11	1.97	0.47
1:A:235:LYS:HB3	1:A:235:LYS:NZ	2.30	0.47
1:B:253:VAL:O	1:B:254:SER:C	2.53	0.47
1:A:279:GLU:HB3	1:A:280:HIS:CE1	2.50	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:58:LYS:HD2	1:A:59:PRO:HD2	1.97	0.47
1:B:248:LYS:O	1:B:251:SER:HB3	2.15	0.47
1:A:240:LEU:CD1	1:A:245:SER:HB3	2.43	0.47
1:B:76:SER:O	1:B:78:TRP:N	2.47	0.47
1:B:326:LEU:HD12	1:B:327:LYS:H	1.80	0.47
1:A:261:ILE:CG2	1:A:262:GLU:N	2.78	0.47
1:B:97:THR:HB	1:B:281:ILE:CD1	2.45	0.47
1:B:200:ILE:HB	1:B:204:GLN:HB2	1.96	0.47
1:B:305:LYS:O	1:B:306:HIS:C	2.52	0.46
1:A:295:ASP:C	1:A:297:GLY:H	2.19	0.46
1:A:203:GLU:HG3	2:A:412:HOH:O	2.15	0.46
1:A:147:ILE:HD11	1:A:318:LEU:HD23	1.97	0.46
1:A:340:GLU:OE1	1:A:340:GLU:N	2.38	0.46
1:A:58:LYS:HD2	1:A:59:PRO:CD	2.46	0.46
1:B:295:ASP:C	1:B:297:GLY:N	2.69	0.46
1:B:263:ILE:HA	1:B:267:THR:HB	1.97	0.46
1:A:288:GLY:CA	1:A:302:LEU:HD23	2.44	0.46
1:A:180:ILE:HG23	1:A:211:MSE:HG3	1.98	0.46
1:A:73:HIS:HB3	1:A:77:LYS:HG2	1.97	0.46
1:B:50:LYS:N	1:B:50:LYS:CD	2.78	0.46
1:B:212:ARG:HD3	1:B:213:TYR:CE1	2.50	0.46
1:A:242:SER:HB3	1:A:245:SER:CB	2.45	0.45
1:B:92:PRO:HD3	1:B:274:TYR:CD2	2.50	0.45
1:B:309:ALA:O	1:B:313:ARG:HB2	2.16	0.45
1:B:177:ASP:O	1:B:195:PRO:HG2	2.16	0.45
1:B:326:LEU:HD12	1:B:327:LYS:N	2.31	0.45
1:B:268:ILE:HG13	1:B:272:LEU:HD13	1.98	0.45
1:B:68:ASP:C	1:B:70:GLY:H	2.18	0.45
1:A:340:GLU:HB2	1:A:342:HIS:CD2	2.52	0.45
1:B:75:LYS:NZ	1:B:76:SER:OG	2.50	0.45
1:B:235:LYS:HZ2	1:B:235:LYS:HB3	1.80	0.44
1:B:134:ALA:O	1:B:135:GLU:C	2.56	0.44
1:B:178:PHE:CD1	1:B:178:PHE:N	2.85	0.44
1:A:104:ARG:HB2	1:A:126:ASN:HB2	2.00	0.44
1:B:171:THR:O	1:B:232:LYS:NZ	2.44	0.44
1:B:101:SER:HB3	1:B:223:GLN:NE2	2.32	0.44
1:A:107:LEU:HD13	1:A:299:TYR:CD2	2.52	0.44
1:B:78:TRP:HH2	1:B:155:ILE:HA	1.81	0.44
1:B:220:TYR:OH	1:B:287:LYS:HB2	2.17	0.44
1:B:230:ILE:O	1:B:234:LEU:HB2	2.18	0.44
1:B:242:SER:HB3	1:B:245:SER:OG	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:240:LEU:HD12	1:B:249:ILE:HD11	2.00	0.44
1:B:186:GLU:HB3	1:B:188:GLU:OE1	2.17	0.44
1:A:156:ASN:ND2	1:A:256:ASN:HB3	2.32	0.44
1:A:58:LYS:CE	1:A:59:PRO:HD2	2.48	0.44
1:A:178:PHE:N	1:A:178:PHE:CD1	2.85	0.44
1:B:197:THR:O	1:B:198:HIS:ND1	2.51	0.43
1:A:115:ASN:O	1:A:115:ASN:ND2	2.50	0.43
1:B:301:ILE:CG2	1:B:333:TYR:HA	2.48	0.43
1:B:301:ILE:HG21	1:B:333:TYR:HA	2.00	0.43
1:A:182:ILE:HD12	1:A:186:GLU:OE2	2.18	0.43
1:B:187:PRO:HB2	2:B:410:HOH:O	2.18	0.43
1:B:237:ILE:HG12	1:B:249:ILE:HD13	2.01	0.43
1:A:262:GLU:C	1:A:264:SER:N	2.71	0.43
1:B:102:LEU:HD22	1:B:286:LEU:HD23	2.01	0.43
1:B:155:ILE:HG22	1:B:159:GLY:HA3	2.00	0.43
1:B:191:ALA:CB	1:B:203:GLU:HG3	2.48	0.43
1:B:303:THR:HG22	1:B:333:TYR:CB	2.48	0.43
1:B:73:HIS:ND1	1:B:73:HIS:O	2.49	0.43
1:B:67:VAL:O	1:B:157:MSE:HE3	2.19	0.42
1:B:307:LEU:CD2	1:B:328:THR:HG21	2.49	0.42
1:A:154:GLN:HE21	1:A:258:GLN:HE21	1.67	0.42
1:A:97:THR:HB	1:A:281:ILE:HD13	2.01	0.42
1:B:113:PRO:HD2	1:B:136:MSE:CE	2.37	0.42
1:B:206:LEU:CD1	1:B:210:ARG:HD2	2.49	0.42
1:A:58:LYS:HE3	1:A:59:PRO:HD2	2.01	0.42
1:A:268:ILE:N	1:A:269:PRO:CD	2.81	0.42
1:A:153:MSE:HE2	1:A:257:MSE:SE	2.69	0.42
1:A:65:MSE:HG2	1:A:85:MSE:HG2	2.00	0.42
1:A:138:LEU:HD11	1:A:149:VAL:HB	2.01	0.42
1:B:118:GLN:O	1:B:121:VAL:HB	2.18	0.42
1:A:201:ASN:ND2	1:A:203:GLU:HB2	2.33	0.42
1:A:266:LYS:O	1:A:270:ASN:ND2	2.53	0.42
1:B:240:LEU:HB3	1:B:245:SER:CB	2.50	0.42
1:B:90:ILE:CG2	1:B:274:TYR:HB3	2.50	0.42
1:A:103:GLU:HG2	1:A:300:GLN:HG3	2.02	0.42
1:A:324:LYS:N	1:A:324:LYS:HD2	2.34	0.42
1:B:331:ILE:N	1:B:331:ILE:CD1	2.82	0.42
1:B:166:ALA:C	1:B:168:GLY:N	2.72	0.42
1:A:91:ASN:OD1	1:A:93:LYS:HB3	2.19	0.42
1:B:109:LYS:HG2	1:B:120:GLY:O	2.19	0.42
1:A:295:ASP:O	1:A:297:GLY:N	2.53	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:163:LEU:HD13	1:B:253:VAL:HG11	2.01	0.41
1:B:253:VAL:HG23	1:B:254:SER:N	2.35	0.41
1:A:227:ARG:HB3	1:A:231:GLN:HE21	1.85	0.41
1:A:108:ILE:HD11	1:A:144:LEU:HB2	2.02	0.41
1:B:210:ARG:CB	1:B:210:ARG:HH11	2.24	0.41
1:A:343:HIS:O	1:A:344:HIS:ND1	2.54	0.41
1:B:280:HIS:O	1:B:281:ILE:HD13	2.19	0.41
1:A:268:ILE:O	1:A:272:LEU:HD13	2.21	0.41
1:B:174:ASN:HD21	1:B:180:ILE:H	1.68	0.41
1:B:183:ALA:HA	1:B:186:GLU:O	2.20	0.41
1:A:156:ASN:HD21	1:A:256:ASN:CB	2.33	0.41
1:A:85:MSE:C	1:A:86:ILE:HD12	2.40	0.41
1:A:211:MSE:SE	1:A:212:ARG:N	3.04	0.41
1:A:263:ILE:O	1:A:263:ILE:CG2	2.69	0.41
1:A:65:MSE:HE3	1:A:85:MSE:HG2	2.03	0.41
1:B:48:LYS:N	1:B:48:LYS:CD	2.83	0.41
1:B:165:ASN:O	1:B:168:GLY:N	2.44	0.41
1:B:239:ALA:C	1:B:241:ASN:N	2.71	0.41
1:A:322:ARG:H	1:A:322:ARG:HD2	1.83	0.41
1:B:313:ARG:CG	1:B:313:ARG:NH1	2.81	0.41
1:A:197:THR:C	1:A:198:HIS:ND1	2.75	0.41
1:B:153:MSE:HG2	1:B:154:GLN:N	2.36	0.40
1:B:78:TRP:CD2	1:B:154:GLN:NE2	2.90	0.40
1:A:201:ASN:N	1:A:204:GLN:HE21	2.12	0.40
1:B:263:ILE:HA	1:B:267:THR:CG2	2.51	0.40
1:B:138:LEU:O	1:B:142:GLN:HG3	2.21	0.40
1:A:74:ARG:O	1:A:75:LYS:C	2.60	0.40
1:A:316:LYS:CB	1:A:322:ARG:HH21	2.34	0.40
1:A:210:ARG:NH1	1:A:210:ARG:HG2	2.36	0.40
1:B:234:LEU:O	1:B:237:ILE:N	2.54	0.40
1:B:239:ALA:O	1:B:240:LEU:C	2.60	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles ⓘ

### 5.3.1 Protein backbone ⓘ

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	295/306 (96%)	261 (88%)	28 (10%)	6 (2%)	9	24
1	B	291/306 (95%)	237 (81%)	39 (13%)	15 (5%)	2	4
All	All	586/612 (96%)	498 (85%)	67 (11%)	21 (4%)	4	9

All (21) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	77	LYS
1	B	254	SER
1	B	294	SER
1	A	263	ILE
1	A	264	SER
1	B	69	THR
1	B	305	LYS
1	B	84	SER
1	B	150	ASP
1	B	240	LEU
1	B	306	HIS
1	B	338	GLY
1	A	50	LYS
1	A	132	GLY
1	B	133	GLY
1	B	304	LYS
1	A	296	GLY
1	B	262	GLU
1	B	70	GLY
1	B	172	VAL
1	A	187	PRO

### 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	260/258 (101%)	243 (94%)	17 (6%)	21	46
1	B	258/258 (100%)	245 (95%)	13 (5%)	30	60
All	All	518/516 (100%)	488 (94%)	30 (6%)	25	52

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	48	LYS
1	A	50	LYS
1	A	77	LYS
1	A	115	ASN
1	A	181	SER
1	A	185	ASN
1	A	251	SER
1	A	257	MSE
1	A	258	GLN
1	A	260	ASN
1	A	272	LEU
1	A	276	ASP
1	A	279	GLU
1	A	289	GLU
1	A	303	THR
1	A	318	LEU
1	A	340	GLU
1	B	48	LYS
1	B	50	LYS
1	B	75	LYS
1	B	115	ASN
1	B	152	PHE
1	B	181	SER
1	B	210	ARG
1	B	212	ARG
1	B	258	GLN
1	B	260	ASN
1	B	289	GLU
1	B	322	ARG
1	B	340	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (24) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	142	GLN
1	A	148	ASN
1	A	185	ASN
1	A	201	ASN
1	A	204	GLN
1	A	226	GLN
1	A	231	GLN
1	A	256	ASN
1	A	258	GLN
1	A	260	ASN
1	A	270	ASN
1	A	311	GLN
1	A	312	ASN
1	B	115	ASN
1	B	142	GLN
1	B	148	ASN
1	B	154	GLN
1	B	226	GLN
1	B	231	GLN
1	B	256	ASN
1	B	258	GLN
1	B	260	ASN
1	B	311	GLN
1	B	342	HIS

### 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 [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, 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	288/306 (94%)	-0.25	3 (1%) 84 85	6, 28, 68, 115	0
1	B	286/306 (93%)	0.29	12 (4%) 40 39	17, 49, 99, 409	0
All	All	574/612 (93%)	0.02	15 (2%) 59 59	6, 38, 88, 409	0

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	73	HIS	8.9
1	B	71	SER	7.3
1	B	72	GLU	7.2
1	B	74	ARG	6.1
1	B	75	LYS	4.9
1	B	76	SER	4.8
1	B	253	VAL	3.6
1	B	344	HIS	3.1
1	A	344	HIS	2.9
1	B	262	GLU	2.5
1	B	252	ALA	2.4
1	A	74	ARG	2.4
1	A	263	ILE	2.2
1	B	254	SER	2.1
1	B	70	GLY	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 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.