



# Full wwPDB X-ray Structure Validation Report ⓘ

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PDB ID : 3V3R  
Title : Crystal Structure of GES-11  
Authors : Delbruck, H.; Hoffmann, K.M.V.; Bebrone, C.  
Deposited on : 2011-12-14  
Resolution : 1.90 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

<https://www.wwpdb.org/validation/2017/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
Xtriage (Phenix)	:	1.13
EDS	:	trunk30967
Percentile statistics	:	20171227.v01 (using entries in the PDB archive December 27th 2017)
Refmac	:	5.8.0158
CCP4	:	7.0 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk30967

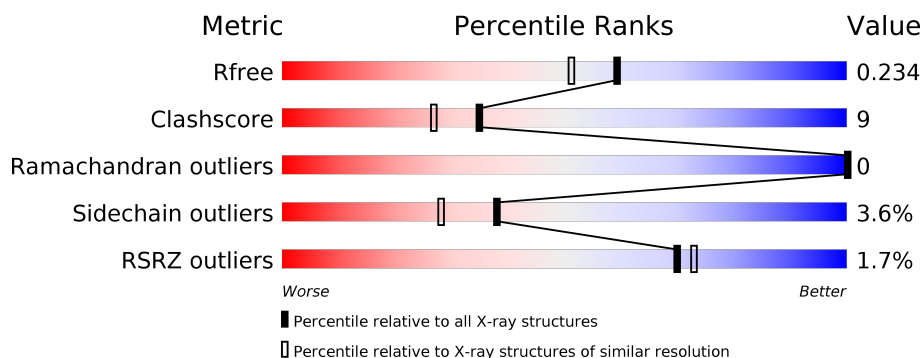
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.90 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))
$R_{free}$	111664	5502 (1.90-1.90)
Clashscore	122126	6115 (1.90-1.90)
Ramachandran outliers	120053	6048 (1.90-1.90)
Sidechain outliers	120020	6048 (1.90-1.90)
RSRZ outliers	108989	5379 (1.90-1.90)

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	287	
1	B	287	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	IOD	A	292[B]	-	-	X	-
2	IOD	A	296	-	-	X	-

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4453 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 Extended spectrum class A beta-lactamase GES-11.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	266	Total	C	N	O	S	0	13	0
			2099	1318	368	402	11			
1	B	265	Total	C	N	O	S	0	8	0
			2064	1296	364	393	11			

- Molecule 2 is IODIDE ION (three-letter code: IOD) (formula: I).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	12	Total	I	0	1
			13	13		
2	A	11	Total	I	0	1
			12	12		

- Molecule 3 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Na	0	0
			1	1		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	145	Total	O	0	0
			145	145		
4	B	119	Total	O	0	0
			119	119		



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 2 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	70.68Å 85.24Å 85.12Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 1.90 23.63 – 1.90	Depositor EDS
% Data completeness (in resolution range)	94.5 (20.00-1.90) 98.0 (23.63-1.90)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.11	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.84 (at 1.90Å)	Xtriage
Refinement program	REFMAC 5.6.0119	Depositor
R, $R_{free}$	0.192 , 0.232 0.197 , 0.234	Depositor DCC
$R_{free}$ test set	2065 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	17.8	Xtriage
Anisotropy	0.485	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.39 , 51.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.000 for -h,l,k	Xtriage
Reported twinning fraction	0.926 for H, K, L 0.074 for -H, L, K	Depositor
Outliers	1 of 40673 reflections (0.002%)	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	4453	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	21.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.59% 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, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

## 5 Model quality

### 5.1 Standard geometry

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

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	0.57	1/2165 (0.0%)	0.75	3/2925 (0.1%)
1	B	0.55	1/2121 (0.0%)	0.74	2/2867 (0.1%)
All	All	0.56	2/4286 (0.0%)	0.75	5/5792 (0.1%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	99	TRP	CD2-CE2	5.14	1.47	1.41
1	A	205	TRP	CD2-CE2	5.06	1.47	1.41

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	45	ASP	CB-CG-OD1	6.47	124.12	118.30
1	A	87	ARG	NE-CZ-NH1	5.19	122.90	120.30
1	B	87	ARG	NE-CZ-NH2	-5.15	117.72	120.30
1	B	183	ARG	NE-CZ-NH2	5.13	122.87	120.30
1	A	87	ARG	NE-CZ-NH2	-5.08	117.76	120.30

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	2099	0	2133	34	0
1	B	2064	0	2082	38	0
2	A	12	0	0	7	0
2	B	13	0	0	4	0
3	B	1	0	0	1	0
4	A	145	0	0	2	0
4	B	119	0	0	4	0
All	All	4453	0	4215	75	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 9.

All (75) 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:261[A]:LYS:HE3	2:A:296:IOD:I	2.22	1.09
1:A:33[B]:ARG:HG3	1:A:33[B]:ARG:HH11	1.19	1.06
1:A:33[B]:ARG:CG	1:A:33[B]:ARG:HH11	1.69	1.04
2:B:294[B]:IOD:I	4:B:389:HOH:O	2.49	1.01
2:B:299:IOD:I	4:B:356:HOH:O	2.48	0.98
2:B:290:IOD:I	3:B:300:NA:NA	2.07	0.98
1:A:145[B]:GLN:HG2	2:A:288:IOD:I	2.34	0.97
1:B:95[A]:MET:HE3	4:B:352:HOH:O	1.65	0.94
1:B:95[A]:MET:HE1	1:B:131:ASN:HB2	1.57	0.86
1:B:57:ALA:HB3	2:B:291:IOD:I	2.51	0.81
1:A:33[B]:ARG:NH1	1:A:33[B]:ARG:HG3	1.89	0.81
1:B:95[A]:MET:CE	1:B:131:ASN:HB2	2.11	0.81
1:B:70:LEU:HD13	1:B:143[A]:MET:HE1	1.64	0.78
1:B:70:LEU:HB2	1:B:143[A]:MET:HE3	1.65	0.77
1:A:78:ILE:HD13	1:A:83:GLU:HB3	1.64	0.77
1:B:137:ILE:HD12	1:B:142:ALA:HB3	1.69	0.75
1:A:70:LEU:HD13	1:A:143[A]:MET:HE3	1.70	0.74
1:B:266[B]:GLU:OE1	1:B:266[B]:GLU:HA	1.87	0.73
1:B:50:ILE:HD11	1:B:179:ILE:HD11	1.69	0.72
1:B:50:ILE:CD1	1:B:179:ILE:HD11	2.20	0.71
1:B:70:LEU:HD13	1:B:143[A]:MET:CE	2.20	0.71
1:A:64:SER:OG	2:A:292[B]:IOD:I	2.79	0.69
1:B:70:LEU:CD1	1:B:143[A]:MET:HE1	2.23	0.69
1:A:142:ALA:O	1:A:145[B]:GLN:HG3	1.92	0.69
1:A:245:LYS:NZ	4:A:358:HOH:O	2.29	0.65
1:A:33[B]:ARG:HG2	1:A:33[B]:ARG:HH11	1.63	0.62
1:A:53:GLY:HA3	1:A:56:MET:CE	2.31	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:73:LEU:HD23	1:B:73:LEU:C	2.20	0.61
1:A:70:LEU:HA	1:A:143[A]:MET:HE1	1.83	0.60
1:A:53:GLY:HA3	1:A:56:MET:HE3	1.82	0.60
1:A:36:ALA:HB1	1:A:261[A]:LYS:HD3	1.84	0.58
1:A:20:GLU:OE2	1:A:21[B]:LYS:HG3	2.04	0.58
1:A:73:LEU:HD13	1:A:143[A]:MET:HE2	1.87	0.56
1:B:50:ILE:HD13	1:B:50:ILE:N	2.20	0.56
1:B:95[A]:MET:HE1	1:B:128:GLY:O	2.06	0.56
1:B:95[A]:MET:CE	4:B:352:HOH:O	2.38	0.53
1:B:95[A]:MET:CE	1:B:131:ASN:CB	2.86	0.53
1:A:24:PHE:CD1	1:A:281:LEU:HD22	2.45	0.52
1:A:70:LEU:HD13	1:A:143[A]:MET:CE	2.39	0.52
1:A:64:SER:CB	2:A:292[B]:IOD:I	3.29	0.51
1:B:95[A]:MET:HE1	1:B:131:ASN:CB	2.36	0.50
1:A:59:ARG:NH2	4:A:389:HOH:O	2.30	0.50
1:B:73:LEU:HD23	1:B:73:LEU:O	2.11	0.49
1:A:143[A]:MET:HE2	1:A:143[A]:MET:HA	1.94	0.49
1:B:68:PHE:CZ	1:B:206:LEU:HD11	2.48	0.49
1:B:50:ILE:HD12	1:B:179:ILE:HD11	1.95	0.47
1:B:44:VAL:CG2	1:B:252:ALA:HB3	2.44	0.47
1:A:73:LEU:HD13	1:A:143[A]:MET:CE	2.44	0.47
1:A:70:LEU:CD1	1:A:143[A]:MET:CE	2.94	0.46
1:A:73:LEU:HD12	1:A:146:TYR:HB2	1.98	0.46
1:B:68:PHE:HB3	1:B:69:PRO:HD3	1.99	0.45
1:B:154:VAL:HG11	1:B:177:THR:HG23	1.98	0.45
1:B:167:ASN:O	1:B:235:ASN:ND2	2.49	0.45
1:B:262:LEU:HD13	1:B:266[B]:GLU:HB3	1.99	0.44
1:A:70:LEU:HD12	1:A:143[A]:MET:HE1	2.00	0.43
1:A:125:SER:OG	2:A:292[B]:IOD:I	2.95	0.43
1:A:64:SER:HB2	2:A:292[B]:IOD:I	2.89	0.43
1:B:137:ILE:HD12	1:B:142:ALA:CB	2.43	0.43
1:B:259:ALA:HB1	1:B:262:LEU:HG	1.99	0.43
1:B:95[A]:MET:HE3	1:B:131:ASN:HB2	1.97	0.42
1:B:215:THR:O	1:B:218:ALA:HB3	2.19	0.42
1:A:142:ALA:HA	1:A:145[B]:GLN:HG3	2.00	0.42
1:A:78:ILE:HD13	1:A:83:GLU:CB	2.42	0.42
1:A:228:GLU:OE1	1:A:240:ASP:OD1	2.38	0.42
1:B:221:PRO:HD2	1:B:224:TRP:CD2	2.55	0.42
1:A:261[A]:LYS:CE	2:A:296:IOD:I	3.15	0.42
1:B:64:SER:HB3	1:B:67:LYS:HD2	2.02	0.42
1:A:142:ALA:HA	1:A:145[B]:GLN:CG	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:270:LEU:O	1:B:274:VAL:HG23	2.20	0.41
1:A:68:PHE:C	1:A:68:PHE:CD1	2.94	0.41
1:B:44:VAL:HG23	1:B:252:ALA:HB3	2.01	0.41
1:B:95[A]:MET:HE3	1:B:131:ASN:CB	2.51	0.41
1:B:166:ASP:O	1:B:234:ALA:HB3	2.21	0.41
1:B:143[B]:MET:HG2	1:B:157:LEU:HD22	2.03	0.40
1:B:77[B]:ARG:NH1	1:B:77[B]:ARG:HB3	2.37	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	277/287 (96%)	267 (96%)	10 (4%)	0	100	100
1	B	271/287 (94%)	263 (97%)	8 (3%)	0	100	100
All	All	548/574 (96%)	530 (97%)	18 (3%)	0	100	100

There are no Ramachandran outliers to report.

### 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	224/227 (99%)	215 (96%)	9 (4%)	34	24

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	216/227 (95%)	210 (97%)	6 (3%)	47	39
All	All	440/454 (97%)	425 (97%)	15 (3%)	38	30

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	47	GLN
1	A	68	PHE
1	A	76	GLU
1	A	119	GLN
1	A	228	GLU
1	A	238	ARG
1	A	245	LYS
1	A	248	GLU
1	A	262	LEU
1	B	68	PHE
1	B	125	SER
1	B	145	GLN
1	B	183	ARG
1	B	228	GLU
1	B	249	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	127	ASN
1	B	280	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

Of 26 ligands modelled in this entry, 26 are monoatomic - leaving 0 for Mogul analysis.

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

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, 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	266/287 (92%)	0.27	1 (0%) 92 93	13, 19, 29, 36	6 (2%)
1	B	265/287 (92%)	0.37	8 (3%) 50 53	14, 21, 30, 36	1 (0%)
All	All	531/574 (92%)	0.32	9 (1%) 70 73	13, 20, 29, 36	7 (1%)

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	197[A]	THR	2.9
1	B	261	LYS	2.8
1	B	19	SER	2.6
1	B	47[A]	GLN	2.5
1	B	249	ARG	2.4
1	B	248	GLU	2.3
1	B	145	GLN	2.3
1	B	70	LEU	2.1
1	B	99	TRP	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](#)

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. 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	B-factors(Å <sup>2</sup> )	Q<0.9
2	IOD	A	298	1/1	0.94	0.05	53,53,53,53	1
3	NA	B	300	1/1	0.96	0.18	16,16,16,16	0
2	IOD	A	296	1/1	0.98	0.04	38,38,38,38	1
2	IOD	B	295	1/1	0.99	0.03	32,32,32,32	1
2	IOD	A	297	1/1	0.99	0.04	41,41,41,41	1
2	IOD	B	298	1/1	0.99	0.03	35,35,35,35	1
2	IOD	A	295	1/1	0.99	0.06	30,30,30,30	1
2	IOD	A	289	1/1	0.99	0.02	37,37,37,37	0
2	IOD	B	293	1/1	0.99	0.05	22,22,22,22	1
2	IOD	B	299	1/1	0.99	0.13	52,52,52,52	1
2	IOD	B	297	1/1	0.99	0.03	27,27,27,27	1
2	IOD	B	296	1/1	0.99	0.03	33,33,33,33	1
2	IOD	A	294	1/1	0.99	0.06	37,37,37,37	1
2	IOD	A	293	1/1	0.99	0.02	29,29,29,29	1
2	IOD	A	291	1/1	0.99	0.02	31,31,31,31	1
2	IOD	B	294[A]	1/1	1.00	0.05	23,23,23,23	1
2	IOD	B	294[B]	1/1	1.00	0.05	34,34,34,34	1
2	IOD	A	290	1/1	1.00	0.03	24,24,24,24	1
2	IOD	B	291	1/1	1.00	0.02	21,21,21,21	1
2	IOD	B	289	1/1	1.00	0.02	22,22,22,22	0
2	IOD	B	292	1/1	1.00	0.02	23,23,23,23	1
2	IOD	A	292[A]	1/1	1.00	0.05	26,26,26,26	1
2	IOD	B	288	1/1	1.00	0.08	42,42,42,42	1
2	IOD	B	290	1/1	1.00	0.03	17,17,17,17	1
2	IOD	A	292[B]	1/1	1.00	0.05	27,27,27,27	1
2	IOD	A	288	1/1	1.00	0.05	15,15,15,15	1

## 6.5 Other polymers

There are no such residues in this entry.