



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 5, 2026 – 02:39 PM UTC

PDB ID : 4FMA / pdb_00004fma
Title : EspG structure
Authors : Shao, F.; Zhu, Y.; Hu, L.
Deposited on : 2012-06-16
Resolution : 2.15 Å(reported)

This is a wwPDB X-ray Structure Validation Summary Report for a publicly released PDB entry.

We welcome your comments at 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.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

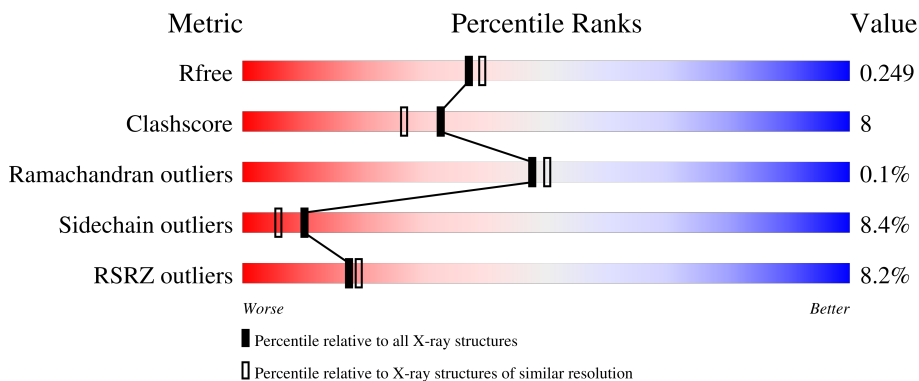
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.15 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}	180053	2057 (2.16-2.16)
Clashscore	190562	2159 (2.16-2.16)
Ramachandran outliers	187476	2134 (2.16-2.16)
Sidechain outliers	187428	2133 (2.16-2.16)
RSRZ outliers	180081	2059 (2.16-2.16)

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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. 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	 3% 85% 12% .
1	B	351	 7% 87% 11% .
1	C	351	 7% 82% 14% ..
1	D	351	 8% 80% 16% ..
1	E	351	 5% 80% 16% ..

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Mol	Chain	Length	Quality of chain
1	F	351	
1	G	351	
1	H	351	
1	I	351	
1	J	351	
1	K	351	
1	L	351	

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 34644 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 EspG protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	351	2719	1679	477	547	16	0	0	0
1	B	351	2719	1679	477	547	16	0	0	0
1	C	346	2671	1653	470	532	16	0	0	0
1	D	341	2636	1632	459	529	16	0	0	0
1	E	349	2696	1666	475	539	16	0	0	0
1	F	344	2658	1645	467	530	16	0	0	0
1	G	343	2627	1630	464	517	16	0	0	0
1	H	347	2678	1657	472	533	16	0	0	0
1	I	348	2658	1644	466	532	16	0	0	0
1	J	341	2624	1626	459	523	16	0	0	0
1	K	344	2639	1635	464	524	16	0	0	0
1	L	348	2673	1655	470	532	16	0	0	0

- Molecule 2 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

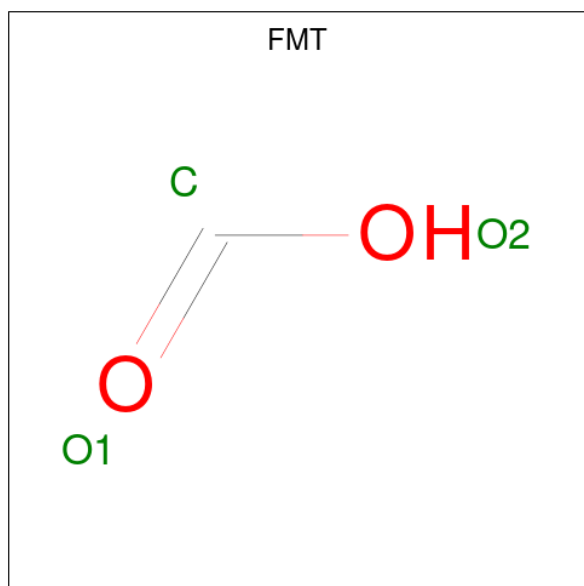
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total 2	Mg 2	0	0
2	B	1	Total 1	Mg 1	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	C	2	Total 2	Mg 2	0	0
2	D	3	Total 3	Mg 3	0	0
2	E	1	Total 1	Mg 1	0	0
2	F	2	Total 2	Mg 2	0	0
2	H	1	Total 1	Mg 1	0	0
2	I	2	Total 2	Mg 2	0	0
2	J	2	Total 2	Mg 2	0	0
2	L	3	Total 3	Mg 3	0	0

- Molecule 3 is FORMIC ACID (CCD ID: FMT) (formula: CH₂O₂).



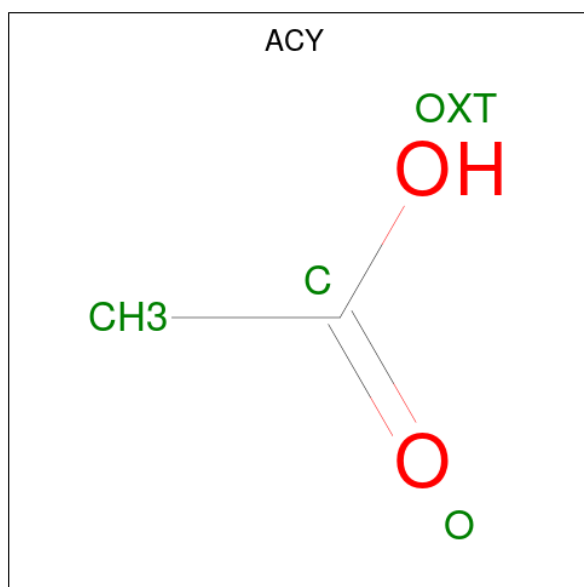
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total 3	C 1	O 2	0	0
3	A	1	Total 3	C 1	O 2	0	0
3	B	1	Total 3	C 1	O 2	0	0

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	D	1	Total	C	O	0	0
			3	1	2		
3	L	1	Total	C	O	0	0
			3	1	2		
3	L	1	Total	C	O	0	0
			3	1	2		

- Molecule 4 is ACETIC ACID (CCD ID: ACY) (formula: C₂H₄O₂).



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

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	281	Total	O	0	0
			281	281		
5	B	287	Total	O	0	0
			287	287		
5	C	251	Total	O	0	0
			251	251		
5	D	227	Total	O	0	0
			227	227		
5	E	243	Total	O	0	0
			243	243		

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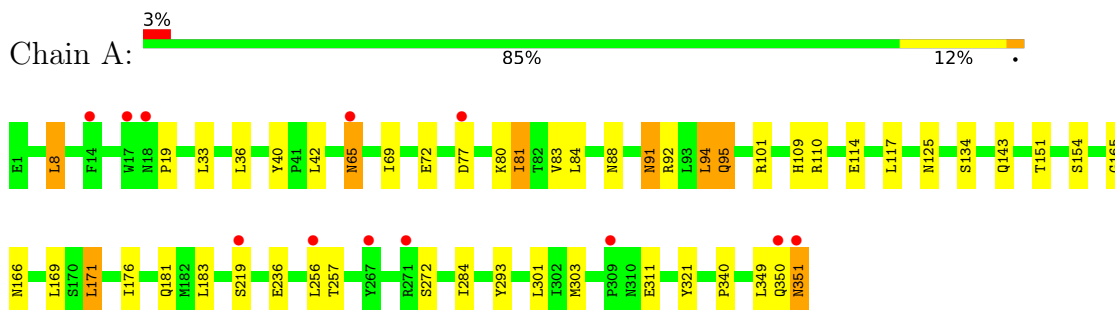
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	F	240	Total 240	O 240	0	0
5	G	164	Total 164	O 164	0	0
5	H	196	Total 196	O 196	0	0
5	I	153	Total 153	O 153	0	0
5	J	152	Total 152	O 152	0	0
5	K	184	Total 184	O 184	0	0
5	L	227	Total 227	O 227	0	0

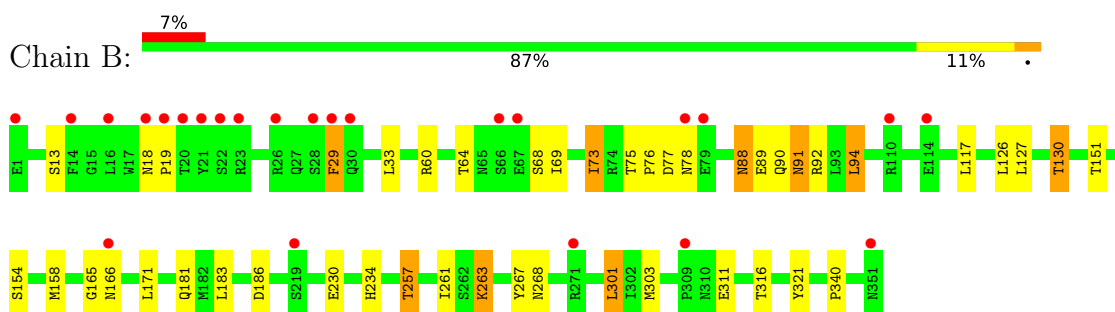
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes 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.

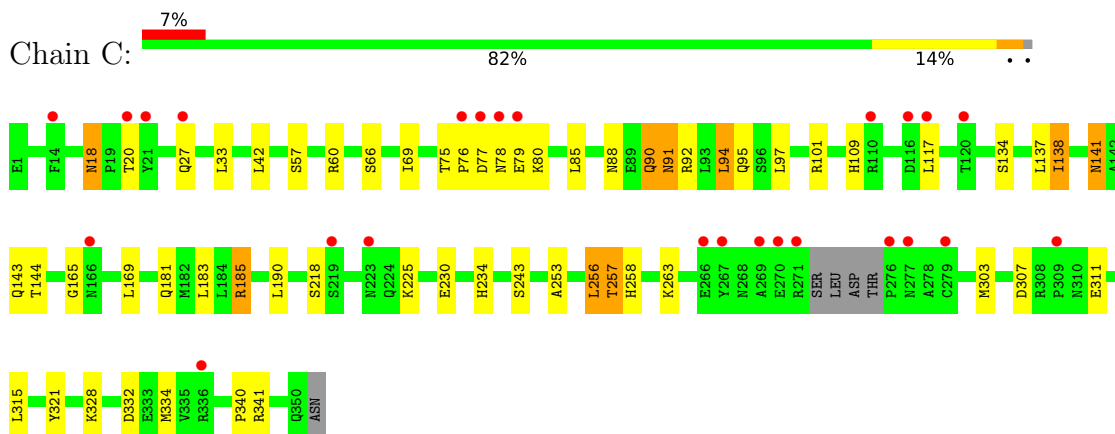
- Molecule 1: EspG protein



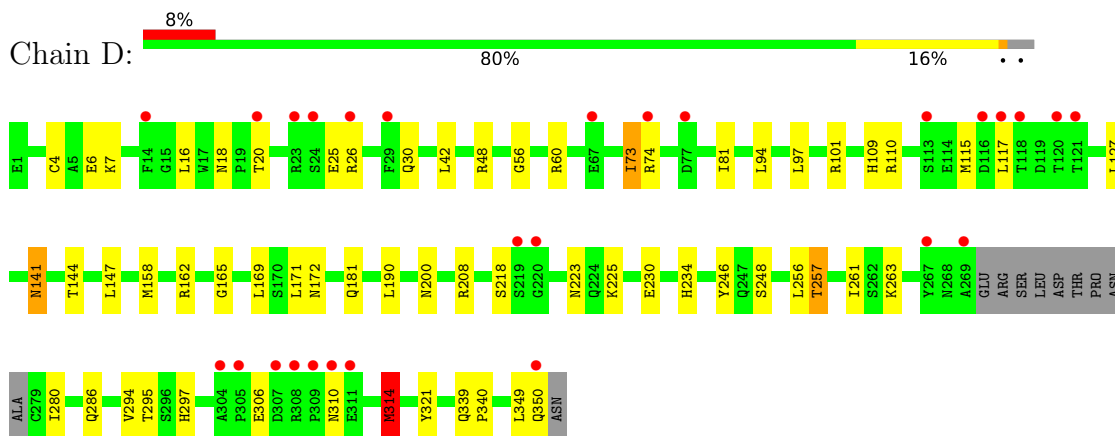
- Molecule 1: EspG protein



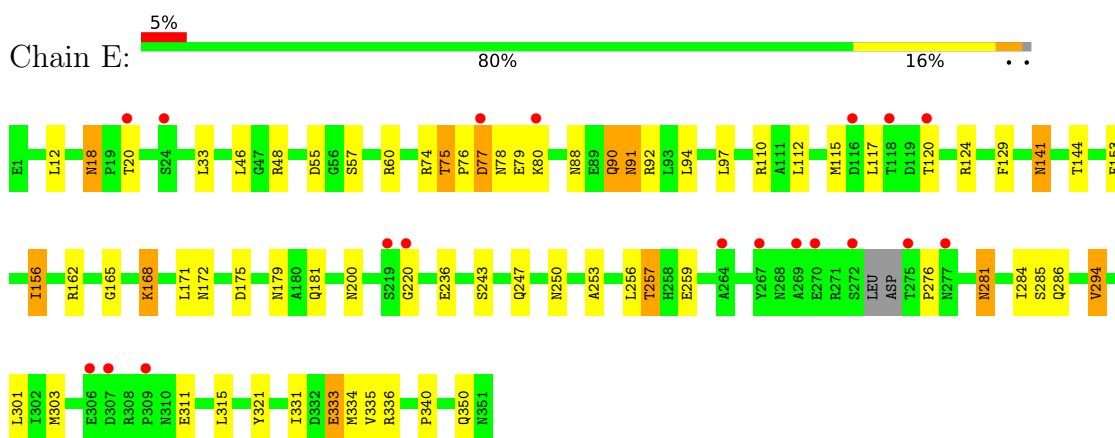
- Molecule 1: EspG protein



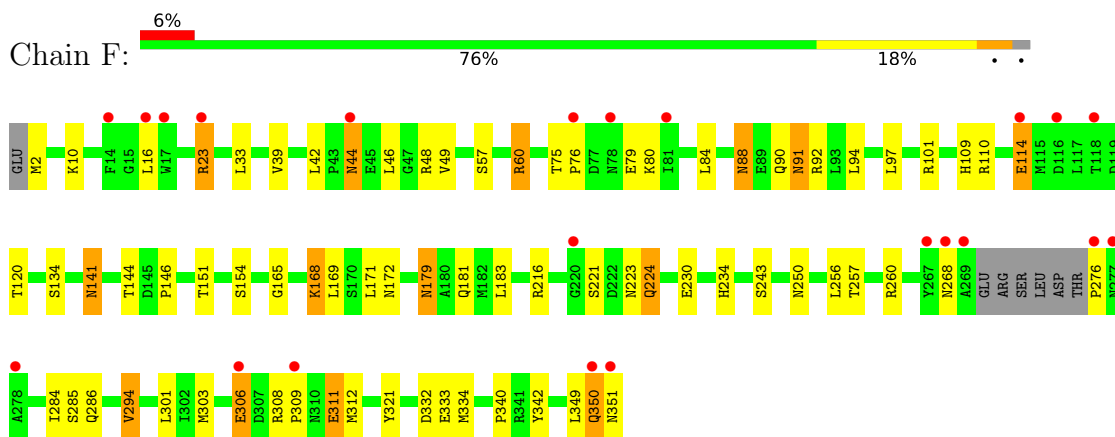
- Molecule 1: EspG protein



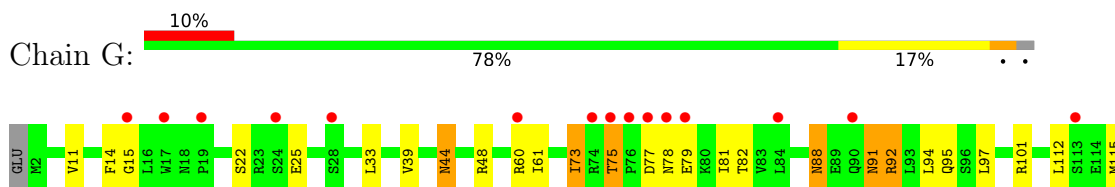
- Molecule 1: EspG protein

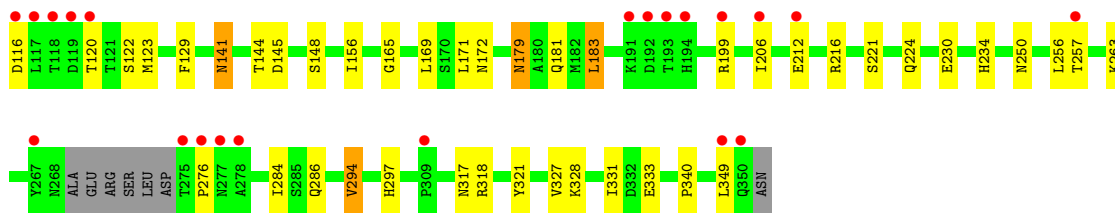


- Molecule 1: EspG protein

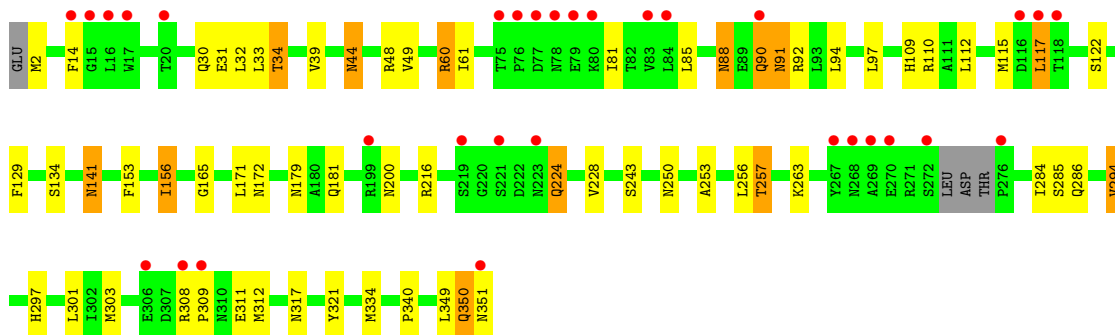
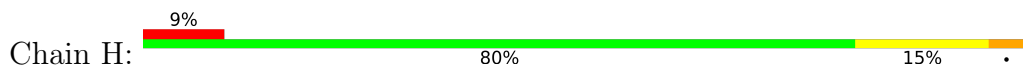


- Molecule 1: EspG protein

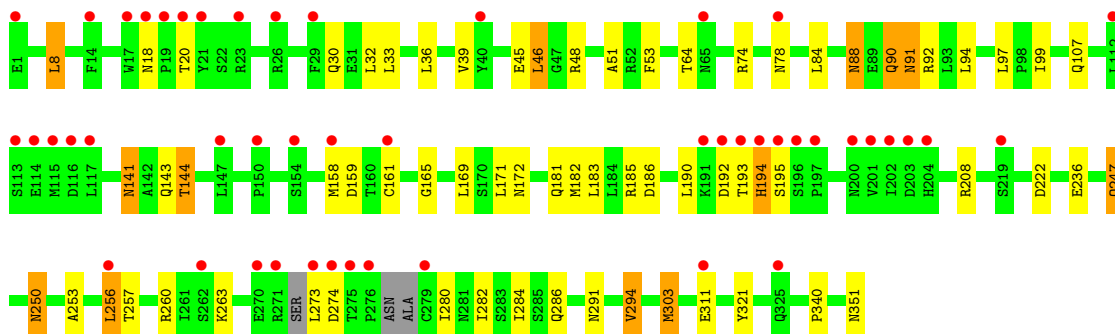
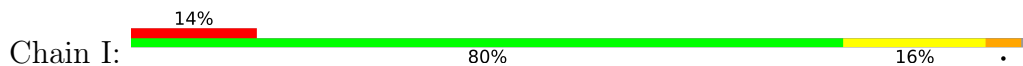




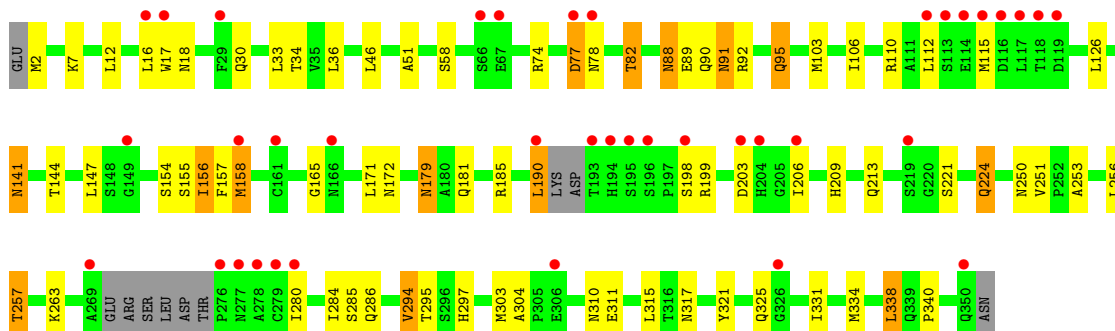
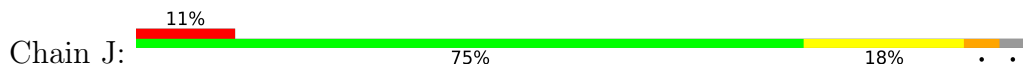
● Molecule 1: EspG protein



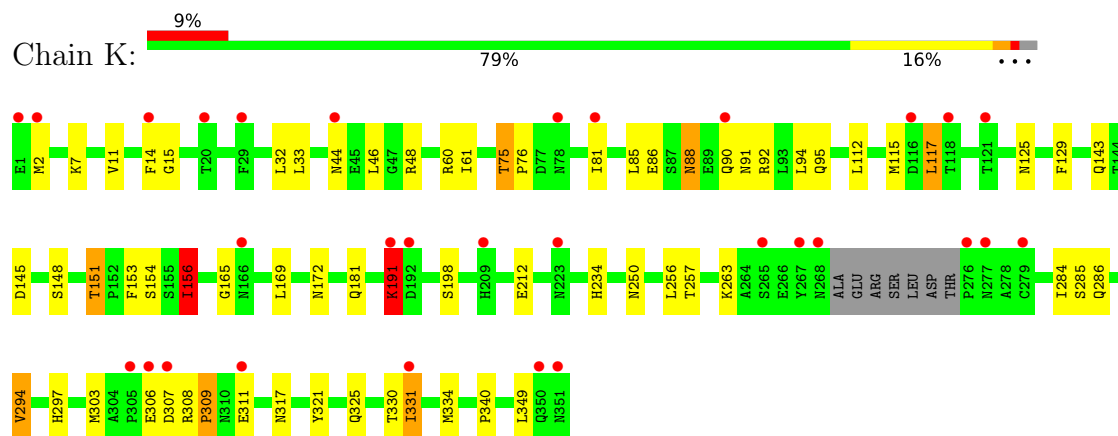
● Molecule 1: EspG protein



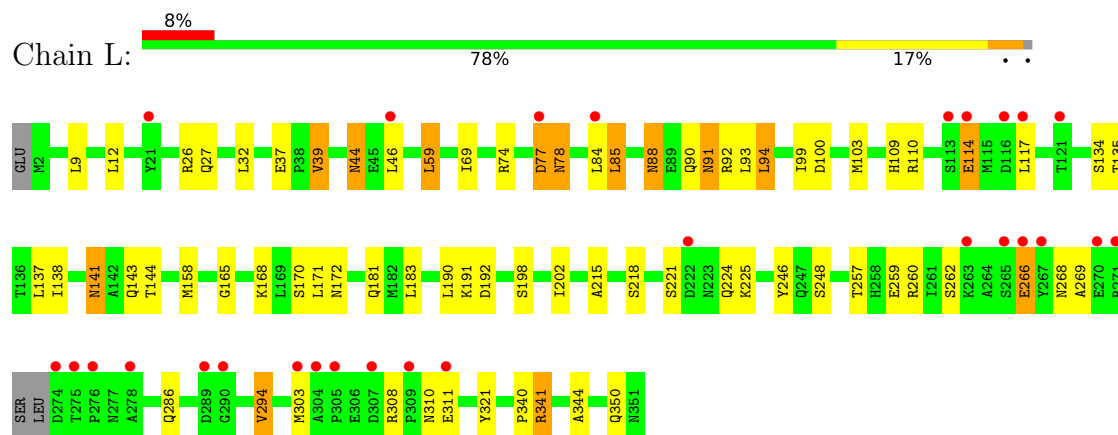
● Molecule 1: EspG protein



- Molecule 1: EspG protein



- Molecule 1: EspG protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	94.61Å 167.74Å 361.69Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.15 50.00 – 2.15	Depositor EDS
% Data completeness (in resolution range)	99.8 (50.00-2.15) 99.8 (50.00-2.15)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.10 (at 2.16Å)	Xtrriage
Refinement program	REFMAC 5.5.0072	Depositor
R, R_{free}	0.200 , 0.239 0.214 , 0.249	Depositor DCC
R_{free} test set	15668 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	29.7	Xtrriage
Anisotropy	0.631	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 41.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	34644	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 14.67% 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

Bond lengths and bond angles in the following residue types are not validated in this section: ACY, FMT, MG

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	0.55	0/2767	0.80	0/3758
1	B	0.54	0/2767	0.81	0/3758
1	C	0.53	0/2718	0.83	3/3689 (0.1%)
1	D	0.57	2/2682 (0.1%)	0.82	0/3641
1	E	0.51	0/2743	0.81	1/3724 (0.0%)
1	F	0.54	0/2705	0.85	4/3672 (0.1%)
1	G	0.50	0/2674	0.83	2/3633 (0.1%)
1	H	0.51	0/2725	0.81	2/3699 (0.1%)
1	I	0.51	0/2703	0.86	1/3673 (0.0%)
1	J	0.49	1/2669 (0.0%)	0.83	4/3623 (0.1%)
1	K	0.50	1/2686 (0.0%)	0.84	3/3649 (0.1%)
1	L	0.55	1/2720 (0.0%)	0.85	4/3695 (0.1%)
All	All	0.52	5/32559 (0.0%)	0.83	24/44214 (0.1%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	314	MET	SD-CE	-8.00	1.59	1.79
1	L	294	VAL	C-O	-6.48	1.17	1.24
1	J	251	VAL	CA-CB	6.01	1.57	1.54
1	D	314	MET	C-O	-5.56	1.17	1.24
1	K	14	PHE	C-O	-5.28	1.17	1.24

The worst 5 of 24 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	80	LYS	N-CA-C	10.37	125.35	108.34
1	K	14	PHE	N-CA-C	8.67	121.98	111.40
1	G	14	PHE	N-CA-C	7.76	123.22	112.04
1	J	156	ILE	N-CA-C	7.14	117.28	110.42
1	F	350	GLN	N-CA-C	6.91	120.25	110.23

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	2719	0	2674	27	0
1	B	2719	0	2674	36	0
1	C	2671	0	2629	36	0
1	D	2636	0	2592	36	0
1	E	2696	0	2646	42	0
1	F	2658	0	2617	50	0
1	G	2627	0	2574	46	0
1	H	2678	0	2629	46	0
1	I	2658	0	2587	53	0
1	J	2624	0	2576	50	0
1	K	2639	0	2588	37	0
1	L	2673	0	2621	52	0
2	A	2	0	0	0	0
2	B	1	0	0	0	0
2	C	2	0	0	0	0
2	D	3	0	0	0	0
2	E	1	0	0	0	0
2	F	2	0	0	0	0
2	H	1	0	0	0	0
2	I	2	0	0	0	0
2	J	2	0	0	0	0
2	L	3	0	0	0	0
3	A	6	0	2	0	0
3	B	3	0	1	0	0
3	D	3	0	1	0	0
3	L	6	0	2	1	0
4	J	4	0	3	0	0
5	A	281	0	0	3	0
5	B	287	0	0	0	0
5	C	251	0	0	3	0
5	D	227	0	0	3	0
5	E	243	0	0	1	0
5	F	240	0	0	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	G	164	0	0	2	0
5	H	196	0	0	3	0
5	I	153	0	0	2	0
5	J	152	0	0	1	0
5	K	184	0	0	0	0
5	L	227	0	0	3	0
All	All	34644	0	31416	498	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 8.

The worst 5 of 498 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:J:172:ASN:HD21	1:J:286:GLN:H	1.10	0.99
1:F:23:ARG:HG2	1:F:23:ARG:HH11	1.31	0.95
1:J:156:ILE:HG21	1:J:334:MET:HE1	1.44	0.95
1:D:172:ASN:HD21	1:D:286:GLN:H	1.11	0.94
1:E:172:ASN:HD21	1:E:286:GLN:H	1.16	0.92

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	349/351 (99%)	343 (98%)	6 (2%)	0	100	100
1	B	349/351 (99%)	344 (99%)	4 (1%)	1 (0%)	36	34
1	C	342/351 (97%)	333 (97%)	9 (3%)	0	100	100
1	D	337/351 (96%)	325 (96%)	12 (4%)	0	100	100
1	E	345/351 (98%)	338 (98%)	6 (2%)	1 (0%)	36	34

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	F	340/351 (97%)	334 (98%)	6 (2%)	0	100	100
1	G	339/351 (97%)	333 (98%)	6 (2%)	0	100	100
1	H	343/351 (98%)	337 (98%)	6 (2%)	0	100	100
1	I	342/351 (97%)	340 (99%)	2 (1%)	0	100	100
1	J	335/351 (95%)	327 (98%)	8 (2%)	0	100	100
1	K	340/351 (97%)	330 (97%)	9 (3%)	1 (0%)	36	34
1	L	344/351 (98%)	333 (97%)	11 (3%)	0	100	100
All	All	4105/4212 (98%)	4017 (98%)	85 (2%)	3 (0%)	48	50

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	77	ASP
1	E	220	GLY
1	K	309	PRO

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	312/312 (100%)	289 (93%)	23 (7%)	13	8
1	B	312/312 (100%)	297 (95%)	15 (5%)	23	20
1	C	304/312 (97%)	281 (92%)	23 (8%)	12	8
1	D	302/312 (97%)	280 (93%)	22 (7%)	13	8
1	E	307/312 (98%)	282 (92%)	25 (8%)	11	7
1	F	304/312 (97%)	269 (88%)	35 (12%)	5	2
1	G	295/312 (95%)	266 (90%)	29 (10%)	7	4
1	H	304/312 (97%)	278 (91%)	26 (9%)	10	5
1	I	299/312 (96%)	270 (90%)	29 (10%)	8	4
1	J	298/312 (96%)	270 (91%)	28 (9%)	8	4

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	K	299/312 (96%)	275 (92%)	24 (8%)	11 7
1	L	303/312 (97%)	275 (91%)	28 (9%)	8 4
All	All	3639/3744 (97%)	3332 (92%)	307 (8%)	10 6

5 of 307 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	J	89	GLU
1	L	84	LEU
1	J	158	MET
1	K	81	ILE
1	L	191	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 172 such sidechains are listed below:

Mol	Chain	Res	Type
1	H	317	ASN
1	K	95	GLN
1	I	88	ASN
1	I	351	ASN
1	K	250	ASN

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 26 ligands modelled in this entry, 19 are monoatomic - leaving 7 for Mogul analysis.

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	$\# Z > 2$
3	FMT	A	403	-	2,2,2	0.75	0	1,1,1	0.24	0
3	FMT	B	402	-	2,2,2	0.72	0	1,1,1	0.23	0
4	ACY	J	403	-	3,3,3	0.83	0	3,3,3	0.73	0
3	FMT	L	404	-	2,2,2	0.77	0	1,1,1	0.28	0
3	FMT	D	404	-	2,2,2	0.75	0	1,1,1	0.21	0
3	FMT	A	404	-	2,2,2	0.65	0	1,1,1	0.25	0
3	FMT	L	405	-	2,2,2	0.83	0	1,1,1	0.25	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.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	L	405	FMT	1	0

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, 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	351/351 (100%)	0.11	12 (3%) 48 52	15, 27, 50, 63	0
1	B	351/351 (100%)	0.12	24 (6%) 23 26	15, 28, 52, 90	0
1	C	346/351 (98%)	0.34	25 (7%) 21 25	19, 31, 62, 88	0
1	D	341/351 (97%)	0.38	27 (7%) 18 20	18, 32, 60, 71	0
1	E	349/351 (99%)	0.37	19 (5%) 31 35	22, 33, 61, 86	0
1	F	344/351 (98%)	0.33	22 (6%) 25 28	20, 31, 58, 75	0
1	G	343/351 (97%)	0.62	36 (10%) 11 13	24, 41, 70, 85	0
1	H	347/351 (98%)	0.59	31 (8%) 15 17	22, 38, 69, 94	0
1	I	348/351 (99%)	0.75	48 (13%) 6 7	24, 43, 74, 113	0
1	J	341/351 (97%)	0.74	38 (11%) 10 11	26, 41, 63, 80	0
1	K	344/351 (98%)	0.64	30 (8%) 16 17	25, 42, 69, 84	0
1	L	348/351 (99%)	0.37	28 (8%) 18 20	17, 31, 58, 81	0
All	All	4153/4212 (98%)	0.45	340 (8%) 17 19	15, 35, 65, 113	0

The worst 5 of 340 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	79	GLU	6.3
1	E	272	SER	6.1
1	D	309	PRO	6.0
1	I	273	LEU	5.9
1	D	118	THR	5.4

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 oligosaccharides 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, 95th 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 ²)	Q<0.9
2	MG	D	402	1/1	0.76	0.14	44,44,44,44	0
4	ACY	J	403	4/4	0.76	0.21	57,57,57,57	0
2	MG	L	402	1/1	0.79	0.23	60,60,60,60	0
2	MG	I	402	1/1	0.83	0.12	64,64,64,64	0
3	FMT	L	405	3/3	0.85	0.19	35,35,35,35	0
2	MG	L	403	1/1	0.86	0.13	49,49,49,49	0
2	MG	E	401	1/1	0.86	0.08	47,47,47,47	0
2	MG	H	401	1/1	0.86	0.14	63,63,63,63	0
3	FMT	B	402	3/3	0.88	0.13	44,44,44,44	0
2	MG	L	401	1/1	0.90	0.08	34,34,34,34	0
2	MG	A	402	1/1	0.90	0.07	31,31,31,31	0
2	MG	J	401	1/1	0.90	0.09	46,46,46,46	0
3	FMT	A	403	3/3	0.91	0.11	39,39,39,39	0
2	MG	D	403	1/1	0.91	0.08	54,54,54,54	0
2	MG	C	401	1/1	0.92	0.06	39,39,39,39	0
2	MG	C	402	1/1	0.92	0.13	59,59,59,59	0
3	FMT	A	404	3/3	0.92	0.10	34,34,34,35	0
2	MG	I	401	1/1	0.94	0.05	34,34,34,34	0
2	MG	F	402	1/1	0.94	0.08	46,46,46,46	0
2	MG	D	401	1/1	0.94	0.06	34,34,34,34	0
2	MG	J	402	1/1	0.94	0.11	39,39,39,39	0
2	MG	A	401	1/1	0.95	0.07	27,27,27,27	0
3	FMT	D	404	3/3	0.97	0.06	21,21,21,21	0
2	MG	B	401	1/1	0.97	0.04	35,35,35,35	0
2	MG	F	401	1/1	0.97	0.13	30,30,30,30	0
3	FMT	L	404	3/3	0.98	0.04	21,21,21,21	0

6.5 Other polymers [i](#)

There are no such residues in this entry.