



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 09:20 AM UTC

PDB ID : 1GQE / pdb_00001gqe
Title : Polypeptide Chain Release Factor 2 (RF2) from Escherichia coli
Authors : Vestergaard, B.; Kjeldgaard, M.
Deposited on : 2001-11-22
Resolution : 1.81 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

<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

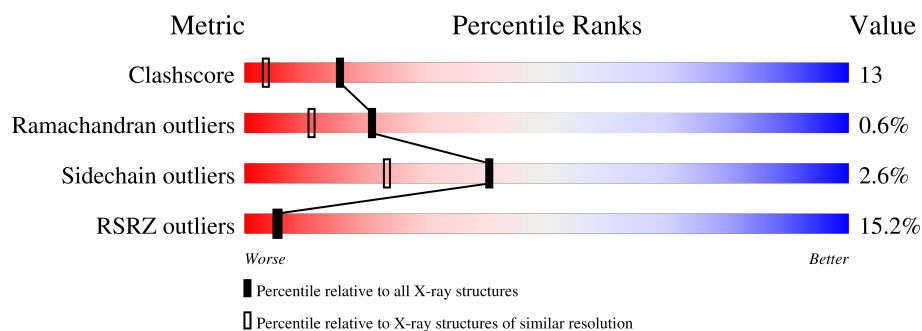
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.81 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))
Clashscore	190562	1148 (1.82-1.82)
Ramachandran outliers	187476	1140 (1.82-1.82)
Sidechain outliers	187428	1140 (1.82-1.82)
RSRZ outliers	180081	1112 (1.82-1.82)

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	365	 14% 76% 21% ..

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3152 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 RELEASE FACTOR 2.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	362	Total	C	N	O	S	Se	75	0	0
			2866	1762	504	590	2	8			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	246	ALA	THR	engineered mutation	UNP P07012
A	298	VAL	LEU	conflict	UNP P07012

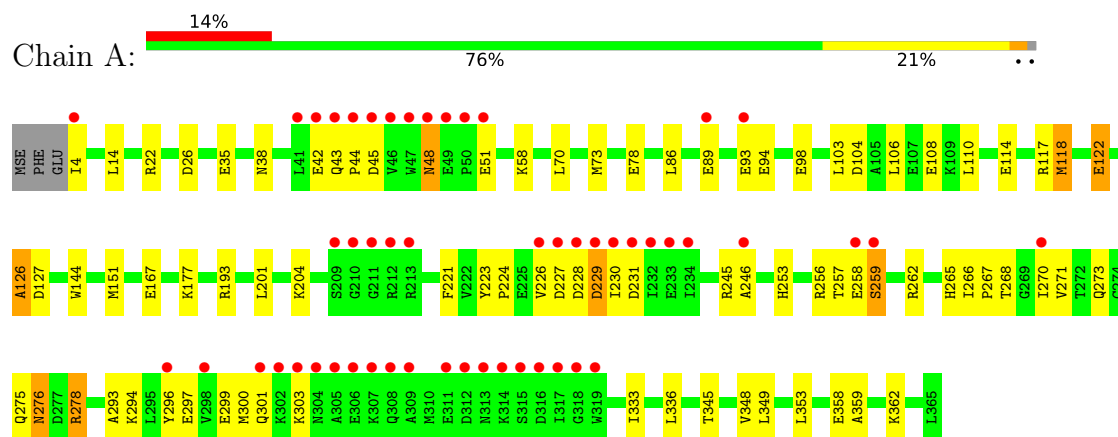
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	286	Total	O	0	0
			286	286		

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: RELEASE FACTOR 2



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	57.39Å 49.89Å 63.14Å 90.00° 107.01° 90.00°	Depositor
Resolution (Å)	18.18 – 1.81 18.18 – 1.81	Depositor EDS
% Data completeness (in resolution range)	96.8 (18.18-1.81) 95.5 (18.18-1.81)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.84 (at 1.81Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.222 , 0.247 0.228 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	17.0	Xtriage
Anisotropy	0.176	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 49.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	3152	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.21% 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	0.34	0/2898	0.85	10/3891 (0.3%)

There are no bond length outliers.

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	257	THR	N-CA-C	-10.12	100.68	113.23
1	A	127	ASP	N-CA-C	-7.58	100.80	110.19
1	A	256	ARG	N-CA-C	-6.37	104.41	111.36
1	A	126	ALA	N-CA-C	5.65	117.99	110.53
1	A	193	ARG	N-CA-C	5.47	118.75	111.75
1	A	204	LYS	N-CA-C	-5.46	101.08	109.76
1	A	276	ASN	N-CA-C	5.41	119.73	113.18
1	A	229	ASP	N-CA-C	-5.35	106.89	113.41
1	A	177	LYS	N-CA-C	-5.23	106.05	112.90
1	A	48	ASN	N-CA-C	-5.02	106.41	112.54

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	2866	0	2765	73	0
2	A	286	0	0	9	0
All	All	3152	0	2765	73	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 13.

All (73) 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:246:ALA:H	1:A:258:GLU:HG2	1.23	1.00
1:A:333:ILE:CD1	1:A:348:VAL:HG21	1.95	0.97
1:A:73:MSE:HE1	1:A:106:LEU:HB3	1.55	0.88
1:A:151:MSE:HE3	1:A:348:VAL:C	2.05	0.82
1:A:333:ILE:HD13	1:A:348:VAL:HG21	1.61	0.82
1:A:345:THR:O	1:A:348:VAL:HG22	1.84	0.78
1:A:268:THR:OG1	1:A:270:ILE:HG22	1.87	0.75
1:A:122:GLU:H	1:A:122:GLU:CD	1.95	0.74
1:A:300:MSE:HE3	1:A:303:LYS:HB3	1.72	0.72
1:A:151:MSE:HE2	1:A:353:LEU:HD21	1.73	0.70
1:A:38:ASN:O	1:A:42:GLU:HG3	1.92	0.69
1:A:245:ARG:HA	1:A:258:GLU:CB	2.21	0.69
1:A:151:MSE:HE3	1:A:348:VAL:O	1.92	0.69
1:A:333:ILE:HD12	1:A:348:VAL:HG21	1.75	0.68
1:A:89:GLU:HG3	2:A:2086:HOH:O	1.93	0.66
1:A:44:PRO:HA	2:A:2045:HOH:O	1.94	0.66
1:A:278:ARG:HH11	1:A:278:ARG:HB3	1.61	0.66
1:A:258:GLU:O	1:A:259:SER:HB2	1.95	0.65
1:A:348:VAL:HG23	1:A:349:LEU:N	2.12	0.65
1:A:73:MSE:CE	1:A:106:LEU:HB3	2.26	0.63
1:A:94:GLU:O	1:A:98:GLU:HG3	1.98	0.63
1:A:86:LEU:C	1:A:86:LEU:HD23	2.25	0.61
1:A:227:ASP:O	1:A:228:ASP:HB2	2.00	0.61
1:A:151:MSE:HE2	1:A:353:LEU:CD2	2.31	0.60
1:A:270:ILE:HD11	1:A:294:LYS:HD3	1.83	0.60
1:A:270:ILE:HD12	2:A:2217:HOH:O	2.00	0.60
1:A:266:ILE:HB	1:A:267:PRO:HD3	1.84	0.60
1:A:245:ARG:HA	1:A:258:GLU:HB3	1.83	0.60
1:A:297:GLU:O	1:A:301:GLN:HG3	2.03	0.58
1:A:245:ARG:HA	1:A:258:GLU:HB2	1.85	0.58
1:A:270:ILE:HD11	1:A:294:LYS:CD	2.34	0.57
1:A:51:GLU:HB3	1:A:228:ASP:OD2	2.05	0.56
1:A:221:PHE:CZ	1:A:297:GLU:HG2	2.40	0.56
1:A:246:ALA:N	1:A:258:GLU:HG2	2.07	0.56
1:A:258:GLU:O	1:A:259:SER:CB	2.54	0.55
1:A:246:ALA:H	1:A:258:GLU:CG	2.10	0.55
1:A:299:GLU:HG2	2:A:2236:HOH:O	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:14:LEU:HD12	1:A:103:LEU:HD11	1.93	0.51
1:A:265:HIS:HB3	1:A:270:ILE:HG22	1.94	0.50
1:A:118:MSE:N	1:A:118:MSE:HE3	2.27	0.49
1:A:348:VAL:CG2	1:A:349:LEU:N	2.76	0.49
1:A:221:PHE:HZ	1:A:297:GLU:HG2	1.78	0.49
1:A:45:ASP:HA	1:A:48:ASN:ND2	2.28	0.48
1:A:270:ILE:HD12	1:A:271:VAL:H	1.77	0.48
1:A:271:VAL:HG21	2:A:2046:HOH:O	2.13	0.48
1:A:275:GLN:NE2	2:A:2220:HOH:O	2.43	0.48
1:A:110:LEU:O	1:A:114:GLU:HG3	2.13	0.48
1:A:151:MSE:CE	1:A:353:LEU:HD21	2.42	0.48
1:A:359:ALA:HA	1:A:362:LYS:HE3	1.95	0.48
1:A:262:ARG:HE	1:A:273:GLN:NE2	2.12	0.47
1:A:167:GLU:HG3	1:A:273:GLN:HB2	1.97	0.47
1:A:278:ARG:HB3	1:A:278:ARG:NH1	2.29	0.46
1:A:78:GLU:HG3	2:A:2079:HOH:O	2.16	0.45
1:A:224:PRO:O	1:A:293:ALA:HB1	2.17	0.45
1:A:104:ASP:O	1:A:108:GLU:HG3	2.17	0.45
1:A:336:LEU:N	1:A:336:LEU:HD12	2.32	0.45
1:A:151:MSE:CE	1:A:348:VAL:O	2.62	0.44
1:A:227:ASP:O	1:A:228:ASP:CB	2.64	0.44
1:A:265:HIS:HB3	1:A:270:ILE:CG2	2.47	0.44
1:A:230:ILE:HD13	1:A:296:TYR:CD1	2.52	0.44
1:A:144:TRP:CE2	1:A:201:LEU:HB2	2.53	0.44
1:A:358:GLU:O	1:A:362:LYS:HE2	2.19	0.43
1:A:22:ARG:HG2	1:A:70:LEU:HD13	1.99	0.43
1:A:270:ILE:HD11	1:A:294:LYS:NZ	2.33	0.43
1:A:43:GLN:HA	1:A:44:PRO:HD3	1.87	0.42
1:A:226:VAL:HB	1:A:230:ILE:HD12	2.01	0.42
1:A:117:ARG:C	1:A:118:MSE:HE3	2.44	0.42
1:A:35:GLU:HG3	2:A:2039:HOH:O	2.19	0.42
1:A:58:LYS:HE2	1:A:126:ALA:HA	2.01	0.42
1:A:4:ILE:HD12	1:A:4:ILE:N	2.36	0.41
1:A:278:ARG:HH11	1:A:278:ARG:CB	2.31	0.41
1:A:299:GLU:O	1:A:303:LYS:HB2	2.21	0.41
1:A:231:ASP:HB2	2:A:2199:HOH:O	2.20	0.41

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	360/365 (99%)	347 (96%)	11 (3%)	2 (1%)	21	11

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	259	SER
1	A	253	HIS

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	307/301 (102%)	299 (97%)	8 (3%)	40	23

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	26	ASP
1	A	93	GLU
1	A	118	MSE
1	A	122	GLU
1	A	223	TYR
1	A	229	ASP
1	A	276	ASN
1	A	278	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	12	GLN
1	A	38	ASN
1	A	48	ASN
1	A	72	GLN
1	A	112	GLN
1	A	273	GLN
1	A	275	GLN
1	A	276	ASN
1	A	280	GLN
1	A	301	GLN
1	A	304	ASN
1	A	355	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

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

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, 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	343/365 (93%)	0.63	52 (15%) 5 5	10, 19, 62, 108	0

All (52) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	312	ASP	8.8
1	A	313	ASN	8.6
1	A	319	TRP	8.5
1	A	315	SER	7.6
1	A	318	GLY	7.1
1	A	314	LYS	6.7
1	A	317	ILE	6.7
1	A	311	GLU	6.7
1	A	309	ALA	6.5
1	A	44	PRO	5.4
1	A	305	ALA	4.8
1	A	258	GLU	4.5
1	A	232	ILE	4.4
1	A	49	GLU	4.4
1	A	46	VAL	4.3
1	A	307	LYS	4.2
1	A	209	SER	4.1
1	A	211	GLY	3.9
1	A	296	TYR	3.8
1	A	228	ASP	3.7
1	A	230	ILE	3.3
1	A	43	GLN	3.3
1	A	303	LYS	3.2
1	A	47	TRP	3.2
1	A	302	LYS	3.2
1	A	301	GLN	3.1
1	A	231	ASP	3.0

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Mol	Chain	Res	Type	RSRZ
1	A	212	ARG	3.0
1	A	233	GLU	3.0
1	A	316	ASP	2.9
1	A	234	ILE	2.8
1	A	229	ASP	2.8
1	A	246	ALA	2.8
1	A	298	VAL	2.8
1	A	4	ILE	2.7
1	A	306	GLU	2.7
1	A	48	ASN	2.6
1	A	308	GLN	2.6
1	A	45	ASP	2.6
1	A	93	GLU	2.5
1	A	259	SER	2.5
1	A	41	LEU	2.5
1	A	227	ASP	2.5
1	A	210	GLY	2.4
1	A	89	GLU	2.2
1	A	50	PRO	2.1
1	A	213	ARG	2.1
1	A	51	GLU	2.1
1	A	226	VAL	2.1
1	A	42	GLU	2.1
1	A	304	ASN	2.1
1	A	270	ILE	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 oligosaccharides 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.