



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 01:27 PM UTC

PDB ID : 1G19 / pdb_00001g19
Title : STRUCTURE OF RECA PROTEIN
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Deposited on : 2000-10-11
Resolution : 3.00 Å(reported)

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

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<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

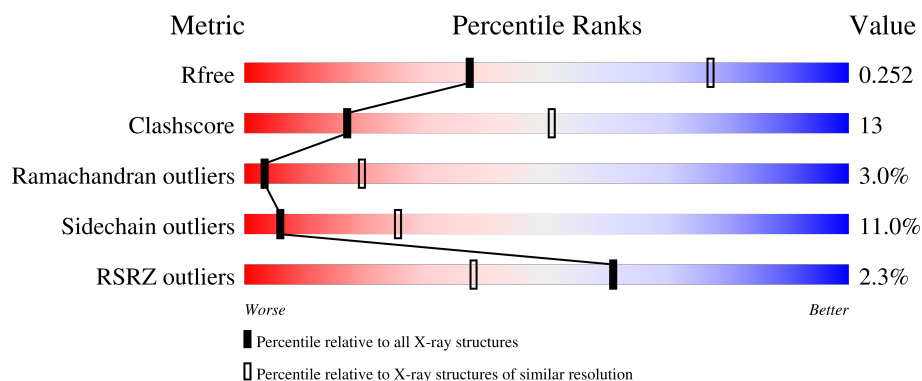
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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	2672 (3.00-3.00)
Clashscore	190562	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)
RSRZ outliers	180081	2671 (3.00-3.00)

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	350	2% 59% 23% 6% 11%

2 Entry composition [i](#)

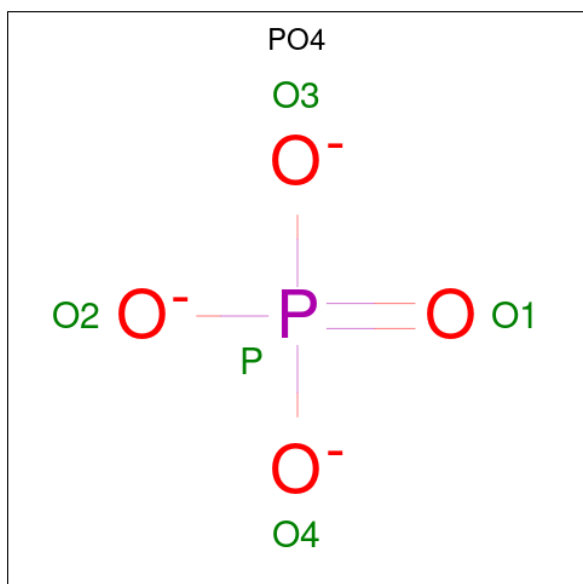
There are 3 unique types of molecules in this entry. The entry contains 2346 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 RECA PROTEIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	311	Total	C	N	O	S	0	0	0
			2276	1423	400	446	7			

- Molecule 2 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	P	0	0
			5	4	1		

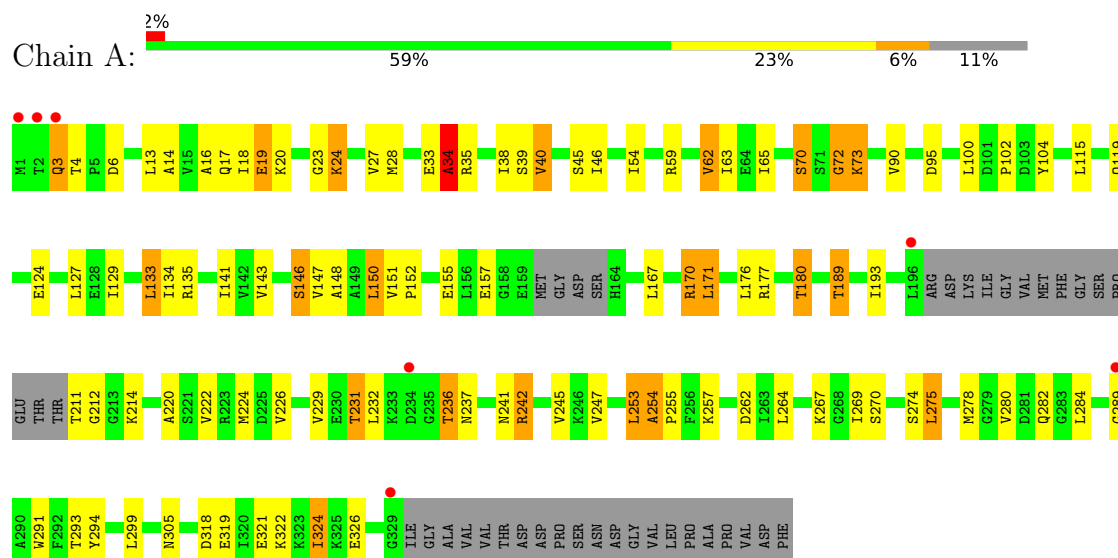
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	65	Total	O	0	0
			65	65		

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: RECA PROTEIN



4 Data and refinement statistics

Property	Value	Source
Space group	P 61	Depositor
Cell constants a, b, c, α , β , γ	108.07Å 108.07Å 72.78Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	15.00 – 3.00 15.00 – 3.00	Depositor EDS
% Data completeness (in resolution range)	91.6 (15.00-3.00) 90.8 (15.00-3.00)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.99 (at 3.00Å)	Xtriage
Refinement program	X-PLOR 3.851	Depositor
R, R_{free}	0.217 , 0.270 0.204 , 0.252	Depositor DCC
R_{free} test set	704 reflections (7.86%)	wwPDB-VP
Wilson B-factor (Å ²)	41.8	Xtriage
Anisotropy	0.133	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.24 , 63.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.065 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	2346	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 3.60% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: PO4

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	0.85	0/2300	1.19	15/3103 (0.5%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2

There are no bond length outliers.

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	212	GLY	N-CA-C	-8.23	104.20	114.16
1	A	95	ASP	N-CA-C	7.11	120.32	108.73
1	A	119	GLN	CA-C-N	7.03	127.44	119.92
1	A	119	GLN	C-N-CA	7.03	127.44	119.92
1	A	254	ALA	CA-C-N	6.27	127.68	119.84
1	A	254	ALA	C-N-CA	6.27	127.68	119.84
1	A	133	LEU	N-CA-C	-6.06	104.79	111.82
1	A	24	LYS	N-CA-C	6.00	117.50	111.07
1	A	319	GLU	N-CA-C	-5.86	104.26	111.40
1	A	151	VAL	CA-C-N	5.71	125.69	120.03
1	A	151	VAL	C-N-CA	5.71	125.69	120.03
1	A	72	GLY	CA-C-O	-5.69	110.67	120.57
1	A	253	LEU	CA-CB-CG	-5.69	96.40	116.30
1	A	135	ARG	N-CA-C	5.46	119.56	113.01
1	A	34	ALA	N-CA-C	5.15	121.77	110.80

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	104	TYR	Sidechain
1	A	72	GLY	Mainchain

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	2276	0	2277	58	0
2	A	5	0	0	1	0
3	A	65	0	0	0	0
All	All	2346	0	2277	58	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 (58) 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:253:LEU:HG	1:A:254:ALA:N	1.76	1.01
1:A:253:LEU:HG	1:A:254:ALA:H	1.26	1.00
1:A:226:VAL:HG22	1:A:245:VAL:HG22	1.62	0.80
1:A:264:LEU:HD21	1:A:270:SER:HB2	1.66	0.77
1:A:278:MET:HE3	1:A:324:ILE:HD11	1.64	0.77
1:A:62:VAL:HG13	1:A:220:ALA:HA	1.68	0.74
1:A:152:PRO:HG2	1:A:167:LEU:HB3	1.75	0.69
1:A:148:ALA:HA	1:A:211:THR:HG21	1.75	0.67
1:A:264:LEU:CD2	1:A:270:SER:HB2	2.27	0.64
1:A:278:MET:HB3	1:A:324:ILE:HD12	1.81	0.62
1:A:129:ILE:O	1:A:133:LEU:HB2	2.00	0.61
1:A:275:LEU:HD11	1:A:321:GLU:HB2	1.83	0.60
1:A:65:ILE:HG12	1:A:224:MET:HE3	1.86	0.58
1:A:40:VAL:HG21	1:A:54:ILE:O	2.03	0.57
1:A:278:MET:HB3	1:A:324:ILE:CD1	2.34	0.57
1:A:124:GLU:HG2	1:A:171:LEU:HD13	1.88	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:176:LEU:O	1:A:180:THR:HB	2.05	0.56
1:A:264:LEU:HB2	1:A:267:LYS:HB2	1.88	0.56
1:A:152:PRO:HG2	1:A:167:LEU:CB	2.36	0.55
1:A:33:GLU:O	1:A:34:ALA:O	2.25	0.54
1:A:148:ALA:CA	1:A:211:THR:HG21	2.37	0.54
1:A:170:ARG:HD3	1:A:171:LEU:N	2.21	0.54
1:A:28:MET:HE1	1:A:254:ALA:N	2.25	0.52
1:A:247:VAL:O	1:A:247:VAL:HG13	2.10	0.52
1:A:231:THR:HA	1:A:241:ASN:HD22	1.76	0.51
1:A:17:GLN:HA	1:A:20:LYS:HD3	1.93	0.50
1:A:253:LEU:O	1:A:254:ALA:C	2.54	0.50
1:A:289:GLY:C	1:A:291:TRP:H	2.21	0.49
1:A:322:LYS:O	1:A:326:GLU:HG3	2.12	0.49
1:A:28:MET:HE1	1:A:253:LEU:HA	1.93	0.49
1:A:73:LYS:NZ	1:A:193:ILE:HG23	2.28	0.49
1:A:141:ILE:HD12	1:A:189:THR:O	2.13	0.49
1:A:127:LEU:HD12	1:A:171:LEU:HD11	1.95	0.48
1:A:177:ARG:O	1:A:180:THR:HG22	2.13	0.48
1:A:147:VAL:HA	1:A:150:LEU:HD22	1.96	0.48
1:A:236:THR:HG23	1:A:237:ASN:H	1.78	0.48
1:A:284:LEU:O	1:A:294:TYR:HD1	1.97	0.47
1:A:19:GLU:HA	1:A:23:GLY:O	2.16	0.46
1:A:146:SER:O	1:A:150:LEU:HD13	2.15	0.46
1:A:229:VAL:HB	1:A:242:ARG:O	2.16	0.45
1:A:155:GLU:HB2	1:A:167:LEU:HD13	1.99	0.45
1:A:247:VAL:HG12	1:A:257:LYS:O	2.16	0.45
1:A:100:LEU:O	1:A:102:PRO:HD3	2.17	0.44
1:A:14:ALA:O	1:A:18:ILE:HG13	2.18	0.44
1:A:39:SER:HB2	1:A:59:ARG:HD2	1.99	0.44
1:A:70:SER:HA	2:A:401:PO4:O2	2.18	0.44
1:A:28:MET:HE1	1:A:254:ALA:H	1.83	0.43
1:A:19:GLU:HG3	1:A:24:LYS:HA	2.01	0.43
1:A:254:ALA:HA	1:A:255:PRO:HD3	1.88	0.43
1:A:236:THR:HG23	1:A:237:ASN:N	2.33	0.43
1:A:262:ASP:O	1:A:269:ILE:HA	2.20	0.42
1:A:3:GLN:HB2	1:A:6:ASP:OD1	2.20	0.42
1:A:148:ALA:CB	1:A:211:THR:HG21	2.50	0.41
1:A:63:ILE:HG13	1:A:222:VAL:HB	2.02	0.41
1:A:13:LEU:O	1:A:16:ALA:HB3	2.21	0.41
1:A:141:ILE:HD12	1:A:189:THR:HG23	2.02	0.41
1:A:46:ILE:HG22	1:A:274:SER:OG	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:280:VAL:C	1:A:282:GLN:H	2.29	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	305/350 (87%)	270 (88%)	26 (8%)	9 (3%)	3 19

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	3	GLN
1	A	34	ALA
1	A	35	ARG
1	A	73	LYS
1	A	170	ARG
1	A	4	THR
1	A	70	SER
1	A	214	LYS
1	A	236	THR

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	228/274 (83%)	203 (89%)	25 (11%)	6 25

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	19	GLU
1	A	27	VAL
1	A	38	ILE
1	A	40	VAL
1	A	45	SER
1	A	62	VAL
1	A	90	VAL
1	A	115	LEU
1	A	134	ILE
1	A	143	VAL
1	A	146	SER
1	A	150	LEU
1	A	157	GLU
1	A	171	LEU
1	A	180	THR
1	A	189	THR
1	A	231	THR
1	A	232	LEU
1	A	242	ARG
1	A	275	LEU
1	A	293	THR
1	A	299	LEU
1	A	305	ASN
1	A	318	ASP
1	A	324	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	17	GLN
1	A	185	ASN
1	A	237	ASN
1	A	241	ASN
1	A	258	GLN
1	A	282	GLN
1	A	301	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](#)

1 ligand is modelled in this entry.

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	$\# Z > 2$
2	PO4	A	401	-	4,4,4	1.68	2 (50%)	6,6,6	0.45	0

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	401	PO4	P-O2	-2.42	1.47	1.54
2	A	401	PO4	P-O4	-2.05	1.48	1.54

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
2	A	401	PO4	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	311/350 (88%)	-0.20	7 (2%) 61 38	2, 18, 69, 96	0

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	329	GLY	5.4
1	A	234	ASP	3.2
1	A	289	GLY	2.3
1	A	3	GLN	2.3
1	A	1	MET	2.2
1	A	196	LEU	2.1
1	A	2	THR	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](#)

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	PO4	A	401	5/5	0.79	0.22	54,54,54,54	0

6.5 Other polymers [i](#)

There are no such residues in this entry.