



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 4, 2024 – 01:46 PM EST

PDB ID : 1RNB
Title : CRYSTAL STRUCTURE OF A BARNASE-D(*GP*C) COMPLEX AT 1.9
ANGSTROMS RESOLUTION
Authors : Janin, J.; Baudet, S.
Deposited on : 1991-03-19
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

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 : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtrriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.36

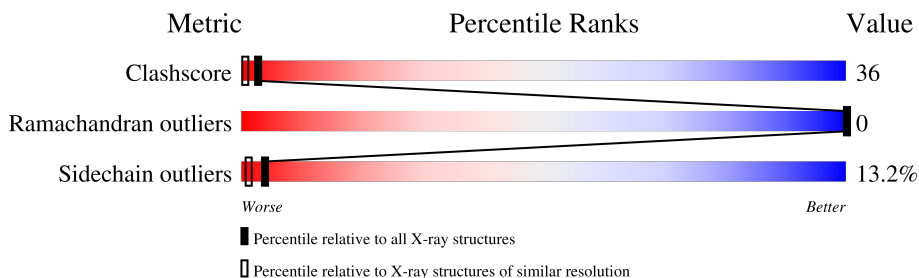
1 Overall quality at a glance

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))
Clashscore	141614	6847 (1.90-1.90)
Ramachandran outliers	138981	6760 (1.90-1.90)
Sidechain outliers	138945	6760 (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 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 $\leq 5\%$

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	C	2	50% 50%
2	A	110	52% 33% 7% 7%

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
3	SO4	C	150	-	-	X	-

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 1008 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 DNA chain called DNA (5'-D(*GP*C)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
1	C	2	38	19	8	10	1	0	0	0

- Molecule 2 is a protein called BARNASE.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
2	A	109	869	550	151	168	0	0	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	22	ASN	ASP	conflict	UNP P00648
A	23	ASP	ASN	conflict	UNP P00648

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	C	1	Total	O	S	0	0
			5	4	1		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	C	8	Total	O	0	0
			8	8		
4	A	88	Total	O	0	0
			88	88		

3 Residue-property plots

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. 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. 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.

Note EDS was not executed.

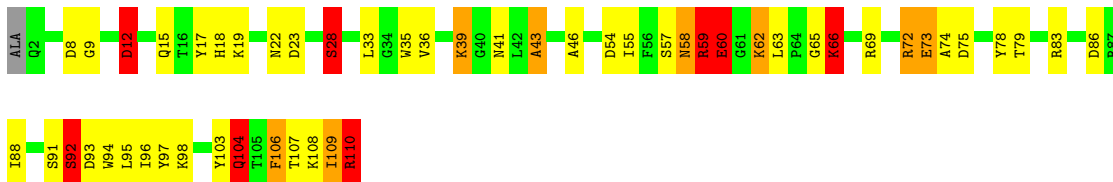
- Molecule 1: DNA (5'-D(*GP*C)-3')

Chain C: 



- Molecule 2: BARNASE

Chain A: 



4 Data and refinement statistics

Xtrriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	57.97Å 57.97Å 85.37Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	(Not available) – 1.90	Depositor
% Data completeness (in resolution range)	(Not available) ((Not available)-1.90)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	PROLSQ	Depositor
R, R_{free}	0.214 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtrriage
Total number of atoms	1008	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP

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: SO4

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	C	3.15	5/42 (11.9%)	3.91	12/63 (19.0%)
2	A	1.11	0/890	2.41	43/1204 (3.6%)
All	All	1.27	5/932 (0.5%)	2.50	55/1267 (4.3%)

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
2	A	1	0

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	112	DC	P-O5'	8.66	1.68	1.59
1	C	111	DG	C5-C4	-6.39	1.33	1.38
1	C	112	DC	N1-C2	-5.88	1.34	1.40
1	C	111	DG	N7-C5	-5.44	1.35	1.39
1	C	112	DC	C4'-O4'	-5.00	1.40	1.45

All (55) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	86	ASP	CB-CG-OD1	15.60	132.34	118.30
2	A	72	ARG	NE-CZ-NH2	-14.71	112.94	120.30
2	A	69	ARG	NE-CZ-NH1	12.48	126.54	120.30
2	A	12	ASP	CB-CG-OD1	-11.48	107.96	118.30
2	A	104	GLN	CB-CG-CD	11.00	140.20	111.60
2	A	12	ASP	CA-CB-CG	-10.16	91.04	113.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	111	DG	O4'-C4'-C3'	-10.13	99.92	106.00
1	C	112	DC	C2-N3-C4	-10.02	114.89	119.90
2	A	59	ARG	CB-CA-C	9.52	129.44	110.40
2	A	109	ILE	CA-CB-CG1	9.13	128.35	111.00
1	C	112	DC	C5-C6-N1	-8.68	116.66	121.00
1	C	111	DG	OP2-P-O3'	8.48	123.85	105.20
2	A	93	ASP	CB-CG-OD1	8.42	125.88	118.30
1	C	112	DC	N1-C2-N3	8.38	125.06	119.20
2	A	28	SER	N-CA-CB	-7.86	98.71	110.50
1	C	112	DC	N1-C2-O2	-7.83	114.20	118.90
2	A	75	ASP	CB-CG-OD1	-7.64	111.42	118.30
2	A	103	TYR	CB-CG-CD1	7.62	125.57	121.00
2	A	78	TYR	CB-CG-CD1	-7.57	116.46	121.00
2	A	73	GLU	OE1-CD-OE2	-7.48	114.32	123.30
2	A	86	ASP	CB-CG-OD2	-7.30	111.73	118.30
2	A	107	THR	CA-CB-CG2	7.29	122.61	112.40
2	A	110	ARG	NE-CZ-NH1	-7.23	116.69	120.30
2	A	95	LEU	CB-CG-CD2	-7.13	98.88	111.00
2	A	110	ARG	CD-NE-CZ	-7.05	113.73	123.60
2	A	59	ARG	NE-CZ-NH1	-6.77	116.92	120.30
2	A	83	ARG	NE-CZ-NH1	-6.70	116.95	120.30
1	C	111	DG	C2-N3-C4	6.41	115.11	111.90
2	A	54	ASP	CB-CG-OD2	-6.34	112.60	118.30
2	A	46	ALA	CB-CA-C	6.29	119.53	110.10
2	A	91	SER	C-N-CA	6.12	137.00	121.70
2	A	73	GLU	CG-CD-OE2	6.07	130.44	118.30
2	A	17	TYR	CB-CG-CD1	-6.01	117.39	121.00
2	A	66	LYS	N-CA-CB	5.85	121.14	110.60
1	C	112	DC	C5'-C4'-O4'	5.84	120.39	109.30
1	C	111	DG	C1'-O4'-C4'	5.82	115.92	110.10
2	A	94	TRP	N-CA-C	5.82	126.70	111.00
2	A	69	ARG	NE-CZ-NH2	-5.79	117.40	120.30
2	A	23	ASP	CB-CG-OD2	5.79	123.52	118.30
1	C	111	DG	O4'-C1'-C2'	-5.77	101.28	105.90
2	A	43	ALA	CA-C-O	-5.74	108.04	120.10
2	A	97	TYR	CB-CG-CD2	5.73	124.44	121.00
2	A	106	PHE	CB-CG-CD1	-5.67	116.83	120.80
2	A	12	ASP	OD1-CG-OD2	5.61	133.97	123.30
1	C	111	DG	N1-C2-N3	-5.57	120.56	123.90
1	C	111	DG	C5-C6-N1	5.53	114.27	111.50
2	A	60	GLU	OE1-CD-OE2	5.42	129.81	123.30
2	A	12	ASP	CA-C-O	5.29	131.21	120.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	41	ASN	N-CA-CB	-5.25	101.14	110.60
2	A	79	THR	CA-CB-OG1	-5.25	97.98	109.00
2	A	63	LEU	CD1-CG-CD2	5.11	125.81	110.50
2	A	58	ASN	C-N-CA	-5.06	109.04	121.70
2	A	92	SER	N-CA-CB	-5.06	102.90	110.50
2	A	95	LEU	CA-CB-CG	5.03	126.86	115.30
2	A	35	TRP	CH2-CZ2-CE2	-5.02	112.38	117.40

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	A	109	ILE	CB

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	C	38	0	24	1	0
2	A	869	0	835	61	0
3	C	5	0	0	1	1
4	A	88	0	0	17	2
4	C	8	0	0	2	1
All	All	1008	0	859	63	2

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 36.

All (63) 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)
2:A:15:GLN:NE2	2:A:110:ARG:HG3	1.83	0.92
2:A:15:GLN:HG3	2:A:110:ARG:NH2	1.86	0.91
2:A:15:GLN:HE22	2:A:110:ARG:HG3	1.36	0.90
2:A:36:VAL:HB	2:A:39:LYS:HG2	1.53	0.90

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:108:LYS:HE3	2:A:110:ARG:OXT	1.72	0.89
2:A:62:LYS:HB3	4:A:313:HOH:O	1.73	0.88
2:A:92:SER:HB3	4:A:315:HOH:O	1.74	0.87
2:A:62:LYS:CB	4:A:313:HOH:O	2.23	0.87
2:A:15:GLN:HA	2:A:110:ARG:HH21	1.41	0.84
2:A:57:SER:O	2:A:59:ARG:NH2	2.13	0.82
2:A:36:VAL:HB	2:A:39:LYS:CG	2.12	0.79
2:A:8:ASP:OD1	4:A:329:HOH:O	2.01	0.77
2:A:36:VAL:HG11	2:A:39:LYS:HD3	1.68	0.74
2:A:15:GLN:NE2	2:A:110:ARG:CG	2.51	0.73
2:A:15:GLN:HE22	2:A:110:ARG:CG	2.02	0.73
2:A:15:GLN:HA	2:A:110:ARG:NH2	2.03	0.72
2:A:18:HIS:HE1	4:A:260:HOH:O	1.73	0.71
2:A:96:ILE:N	2:A:110:ARG:O	2.18	0.70
2:A:15:GLN:CG	2:A:110:ARG:NH2	2.54	0.70
2:A:36:VAL:CG1	2:A:39:LYS:HD3	2.22	0.69
2:A:96:ILE:HD12	2:A:110:ARG:HE	1.59	0.68
1:C:111:DG:H2'	1:C:111:DG:N3	2.09	0.68
3:C:150:SO4:O4	4:C:257:HOH:O	2.14	0.65
2:A:65:GLY:O	2:A:66:LYS:HD3	1.97	0.65
2:A:15:GLN:NE2	2:A:110:ARG:NE	2.44	0.65
2:A:60:GLU:HB2	2:A:62:LYS:HD2	1.80	0.64
2:A:15:GLN:HG3	2:A:110:ARG:CZ	2.27	0.63
2:A:55:ILE:O	4:A:305:HOH:O	2.15	0.63
2:A:9:GLY:O	2:A:12:ASP:HB3	2.00	0.62
2:A:15:GLN:CA	2:A:110:ARG:HH21	2.11	0.61
2:A:28:SER:HB2	4:A:208:HOH:O	2.01	0.60
2:A:8:ASP:O	2:A:12:ASP:HB2	2.03	0.59
2:A:12:ASP:HB3	4:A:211:HOH:O	2.02	0.59
2:A:15:GLN:CG	2:A:110:ARG:CZ	2.80	0.58
2:A:15:GLN:NE2	2:A:110:ARG:CD	2.67	0.58
2:A:57:SER:O	2:A:58:ASN:CB	2.50	0.56
2:A:98:LYS:O	2:A:106:PHE:HA	2.06	0.55
2:A:15:GLN:CB	2:A:110:ARG:NH2	2.70	0.55
2:A:57:SER:O	2:A:58:ASN:HB3	2.07	0.54
2:A:43:ALA:HA	4:A:328:HOH:O	2.06	0.54
2:A:9:GLY:O	2:A:12:ASP:CB	2.56	0.53
2:A:109:ILE:HG22	2:A:110:ARG:HB2	1.90	0.53
2:A:72:ARG:NE	4:A:212:HOH:O	2.12	0.53
2:A:55:ILE:HD11	2:A:72:ARG:NH2	2.25	0.52
2:A:15:GLN:CA	2:A:110:ARG:NH2	2.71	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:55:ILE:HD12	4:A:212:HOH:O	2.12	0.50
2:A:22:ASN:HB2	4:A:230:HOH:O	2.12	0.49
2:A:36:VAL:O	2:A:39:LYS:HG2	2.13	0.48
2:A:73:GLU:HA	2:A:88:ILE:O	2.14	0.48
2:A:18:HIS:CE1	4:A:260:HOH:O	2.58	0.48
2:A:15:GLN:CD	2:A:110:ARG:CZ	2.84	0.47
2:A:15:GLN:CB	2:A:110:ARG:HH21	2.29	0.46
4:C:257:HOH:O	2:A:104:GLN:NE2	2.46	0.45
2:A:108:LYS:NZ	4:A:250:HOH:O	2.50	0.45
2:A:55:ILE:CD1	4:A:212:HOH:O	2.64	0.45
2:A:59:ARG:HD3	2:A:59:ARG:HA	1.80	0.45
2:A:62:LYS:HB2	4:A:313:HOH:O	2.04	0.44
2:A:62:LYS:H	2:A:62:LYS:HG2	1.36	0.43
2:A:55:ILE:HD11	2:A:72:ARG:CZ	2.49	0.42
2:A:62:LYS:HB2	2:A:106:PHE:CE2	2.56	0.41
2:A:92:SER:CB	4:A:315:HOH:O	2.48	0.41
2:A:74:ALA:HB3	2:A:88:ILE:HG22	2.03	0.40
2:A:58:ASN:OD1	2:A:62:LYS:HG2	2.20	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:C:240:HOH:O	4:A:263:HOH:O[5_666]	2.07	0.13
3:C:150:SO4:O4	4:A:273:HOH:O[5_666]	2.18	0.02

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
2	A	107/110 (97%)	103 (96%)	4 (4%)	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
2	A	91/92 (99%)	79 (87%)	12 (13%)	4 1

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	A	12	ASP
2	A	19	LYS
2	A	28	SER
2	A	33	LEU
2	A	39	LYS
2	A	59	ARG
2	A	60	GLU
2	A	62	LYS
2	A	66	LYS
2	A	92	SER
2	A	104	GLN
2	A	110	ARG

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

Mol	Chain	Res	Type
2	A	15	GLN
2	A	18	HIS

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 monosaccharides 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	$\# Z > 2$
3	SO4	C	150	-	4,4,4	0.70	0	6,6,6	0.69	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 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	150	SO4	1	1

5.7 Other polymers [i](#)

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

6.1 Protein, DNA and RNA chains

EDS was not executed - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.