Full wwPDB X-ray Structure Validation Report (i)

May 14, 2020 – 09:25 pm BST

PDB ID	:	1BND
Title	:	STRUCTURE OF THE BRAIN-DERIVED NEUROTROPHIC FACTOR(SLASH)NEUROTROPHIN 3 HETERODIMER
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Deposited on	:	1994-12-12
Resolution	:	2.30 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 (i) symbol.

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

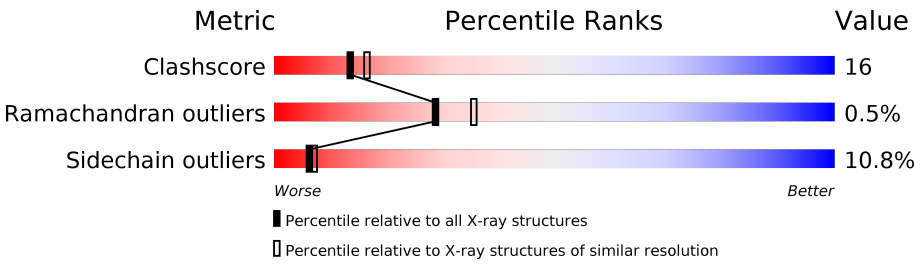
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)

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 >=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 <=5%

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	119	70% 18% • 8%
2	B	119	57% 28% 5% • 9%

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	IPA	A	500	-	-	X	-

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 1937 atoms, of which 141 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 BRAIN DERIVED NEUROTROPHIC FACTOR.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	109	858	535	153	161	9	0	0	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	9	GLN	GLU	CONFLICT	UNP P23560

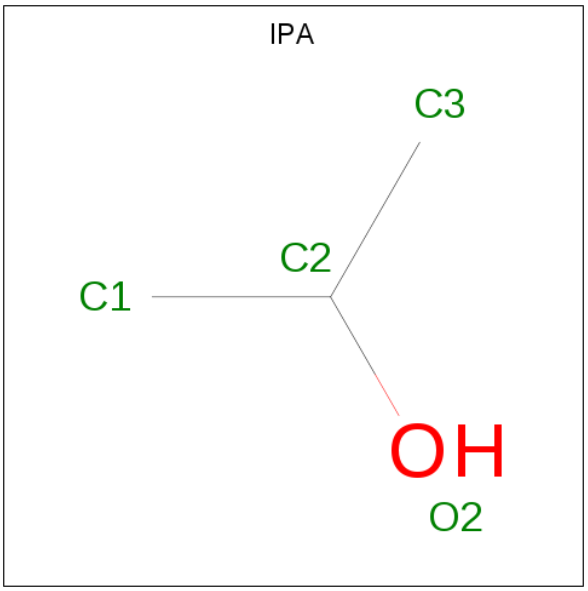
- Molecule 2 is a protein called NEUROTROPHIN 3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	108	864	532	162	164	6	0	0	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	11	VAL	TYR	CONFLICT	UNP P20783
B	44	GLN	GLY	CONFLICT	UNP P20783

- Molecule 3 is ISOPROPYL ALCOHOL (three-letter code: IPA) (formula: C_3H_8O).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	H	O		
3	A	1	5	3	1	1	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	H	O		
4	A	37	111	74	37	0	0
			Total	H	O		
4	B	33	99	66	33	0	0

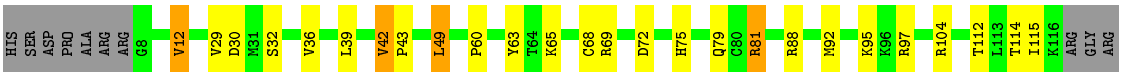
3 Residue-property plots (i)

These plots are drawn for all protein, RNA and DNA 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.

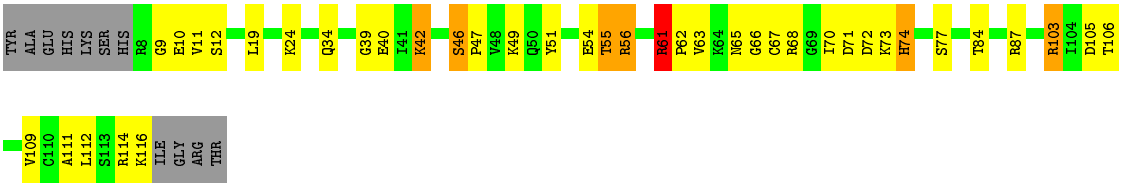
- Molecule 1: BRAIN DERIVED NEUROTROPHIC FACTOR

Chain A: 70% 18% • 8%



- Molecule 2: NEUROTROPHIN 3

Chain B: 57% 28% 5% • 9%



4 Data and refinement statistics (i)

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

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	98.40Å 45.10Å 68.00Å 90.00° 118.30° 90.00°	Depositor
Resolution (Å)	20.00 – 2.30	Depositor
% Data completeness (in resolution range)	(Not available) (20.00-2.30)	Depositor
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR	Depositor
R, R_{free}	0.182 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	1937	wwPDB-VP
Average B, all atoms ($Å^2$)	44.0	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: IPA

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	0.71	0/872	0.90	0/1173
2	B	0.77	0/879	1.05	5/1186 (0.4%)
All	All	0.74	0/1751	0.98	5/2359 (0.2%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	74	HIS	CA-CB-CG	8.41	127.89	113.60
2	B	74	HIS	N-CA-CB	7.52	124.13	110.60
2	B	61	ARG	NE-CZ-NH2	6.92	123.76	120.30
2	B	74	HIS	CB-CA-C	-5.43	99.54	110.40
2	B	63	VAL	O-C-N	-5.36	114.12	122.70

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	858	0	864	31	0
2	B	864	0	851	30	0
3	A	4	1	8	9	0
4	A	37	74	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	33	66	0	1	0
All	All	1796	141	1723	55	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 16.

All (55) 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:95:LYS:HD3	1:A:97:ARG:HH12	1.33	0.93
1:A:104:ARG:NE	3:A:500:IPA:H33	1.85	0.92
1:A:29:VAL:HG23	3:A:500:IPA:H32	1.55	0.88
2:B:42:LYS:HA	2:B:47:PRO:HB3	1.57	0.84
2:B:62:PRO:HD2	2:B:66:GLY:O	1.76	0.83
2:B:42:LYS:HD2	2:B:42:LYS:H	1.45	0.82
2:B:61:ARG:O	2:B:61:ARG:HG2	1.79	0.80
1:A:69:ARG:HG2	2:B:112:LEU:HD12	1.67	0.74
1:A:30:ASP:O	3:A:500:IPA:H31	1.93	0.69
1:A:104:ARG:CD	3:A:500:IPA:H33	2.21	0.69
1:A:29:VAL:HG21	3:A:500:IPA:H13	1.77	0.65
2:B:65:ASN:OD1	2:B:72:ASP:HB3	1.96	0.65
1:A:69:ARG:O	2:B:112:LEU:HD11	1.99	0.63
1:A:115:ILE:HD12	1:A:115:ILE:H	1.65	0.62
2:B:103:ARG:HG2	2:B:103:ARG:HH11	1.67	0.60
1:A:29:VAL:CG2	3:A:500:IPA:H32	2.32	0.56
1:A:88:ARG:NH2	2:B:51:TYR:CE1	2.75	0.55
2:B:67:CYS:SG	2:B:77:SER:HB2	2.47	0.54
1:A:30:ASP:O	3:A:500:IPA:C3	2.54	0.54
1:A:79:GLN:HE21	1:A:81:ARG:HH12	1.57	0.53
1:A:29:VAL:CG2	3:A:500:IPA:H13	2.39	0.52
1:A:81:ARG:HD3	1:A:112:THR:HB	1.90	0.52
2:B:56:ARG:H	2:B:56:ARG:HD2	1.75	0.52
2:B:65:ASN:ND2	2:B:72:ASP:HB2	2.25	0.51
2:B:103:ARG:NH1	2:B:103:ARG:HG2	2.26	0.50
1:A:63:TYR:N	1:A:63:TYR:CD1	2.80	0.49
2:B:65:ASN:OD1	2:B:72:ASP:CB	2.61	0.49
1:A:60:PRO:O	1:A:65:LYS:HD2	2.13	0.48
2:B:116:LYS:HB3	2:B:116:LYS:NZ	2.28	0.48
1:A:95:LYS:HD3	1:A:97:ARG:NH1	2.16	0.47
2:B:54:GLU:HG2	2:B:105:ASP:HB2	1.97	0.47
4:A:364:HOH:O	2:B:55:THR:CG2	2.63	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:61:ARG:CG	2:B:61:ARG:O	2.58	0.47
1:A:12:VAL:HG23	2:B:112:LEU:HG	1.97	0.46
2:B:65:ASN:HD21	2:B:72:ASP:HB2	1.81	0.46
2:B:24:LYS:HA	4:B:367:HOH:O	2.15	0.46
1:A:92:MET:HA	1:A:97:ARG:O	2.16	0.46
2:B:70:ILE:HG22	2:B:71:ASP:N	2.32	0.45
1:A:42:VAL:HA	1:A:43:PRO:HD3	1.83	0.45
1:A:88:ARG:NH1	4:A:318:HOH:O	2.48	0.45
2:B:71:ASP:OD1	2:B:73:LYS:HB2	2.16	0.45
1:A:114:THR:HG22	2:B:10:GLU:HG2	2.00	0.44
1:A:115:ILE:HD12	2:B:9:GLY:O	2.18	0.44
1:A:104:ARG:HD2	3:A:500:IPA:H33	1.98	0.43
2:B:11:VAL:HG11	2:B:68:ARG:HD2	2.00	0.43
2:B:84:THR:HB	2:B:106:THR:HG21	2.01	0.43
2:B:77:SER:HA	2:B:111:ALA:O	2.20	0.42
1:A:79:GLN:HG2	1:A:81:ARG:HH11	1.84	0.42
1:A:63:TYR:N	1:A:63:TYR:HD1	2.17	0.42
2:B:40:GLU:CD	2:B:47:PRO:HB2	2.40	0.42
2:B:39:GLY:O	2:B:49:LYS:HG3	2.20	0.42
1:A:81:ARG:HD3	1:A:112:THR:CB	2.50	0.42
1:A:72:ASP:OD1	1:A:75:HIS:CD2	2.74	0.41
1:A:79:GLN:HE21	1:A:81:ARG:NH1	2.18	0.41
1:A:49:LEU:HD23	1:A:49:LEU:HA	1.89	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	107/119 (90%)	97 (91%)	10 (9%)	0	100 100
2	B	106/119 (89%)	98 (92%)	7 (7%)	1 (1%)	17 20

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	213/238 (90%)	195 (92%)	17 (8%)	1 (0%)	29 35

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	46	SER

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	97/105 (92%)	89 (92%)	8 (8%)	11 14
2	B	98/107 (92%)	85 (87%)	13 (13%)	4 4
All	All	195/212 (92%)	174 (89%)	21 (11%)	6 7

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	12	VAL
1	A	32	SER
1	A	36	VAL
1	A	39	LEU
1	A	42	VAL
1	A	49	LEU
1	A	68	CYS
1	A	81	ARG
2	B	12	SER
2	B	19	LEU
2	B	34	GLN
2	B	42	LYS
2	B	46	SER
2	B	55	THR
2	B	56	ARG
2	B	61	ARG
2	B	74	HIS

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Mol	Chain	Res	Type
2	B	87	ARG
2	B	103	ARG
2	B	109	VAL
2	B	114	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	75	HIS
1	A	79	GLN
2	B	76	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 carbohydrates 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	IPA	A	500	-	3,3,3	0.40	0	3,3,3	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 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	500	IPA	9	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)

EDS was not executed - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains (i)

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates (i)

EDS was not executed - this section is therefore empty.

6.4 Ligands (i)

EDS was not executed - this section is therefore empty.

6.5 Other polymers (i)

EDS was not executed - this section is therefore empty.