WORLDWIDE wwPDB PROTEIN DATA BANK

wwPDB X-ray Structure Validation Summary Report (i)

Apr 30, 2024 – 07:49 pm BST

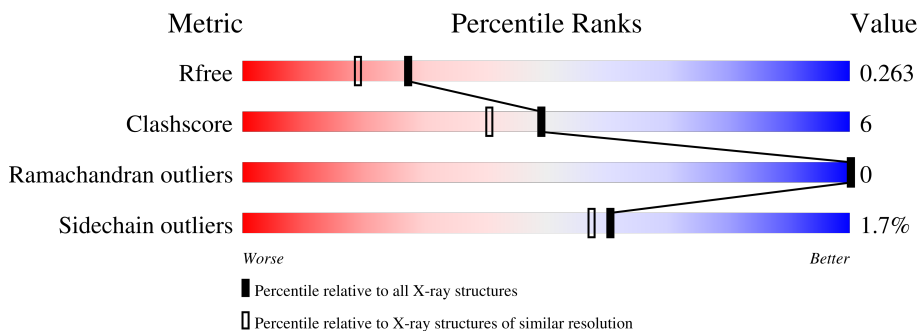|		
PDB ID	:	4A9H
Title	:	N-TERMINAL BROMODOMAIN OF HUMAN BRD2 WITH 1-(2-methyl-1, 2,3,4-tetrahydroquinolin-1-yl)ethan-1-one
Authors	:	Chung, C.W.; Bamborough, P.
Deposited on	:	2011-11-26
Resolution	:	2.05 Å(reported)

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 2.05 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	Percentile Ranks	Value
Rfree		0.263
Clashscore		6
Ramachandran outliers		0
Sidechain outliers		1.7%

Worse   *Better*

▮ Percentile relative to all X-ray structures

▯ Percentile relative to X-ray structures of similar resolution

Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	1692 (2.04-2.04)
Clashscore	141614	1773 (2.04-2.04)
Ramachandran outliers	138981	1752 (2.04-2.04)
Sidechain outliers	138945	1752 (2.04-2.04)

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 <=5%

Mol	Chain	Length	Quality of chain
1	A	154	59% 13% • 27%
1	B	154	63% 6% 31%
1	C	154	64% 5% • 31%

2　Entry composition ⓘ

There are 5 unique types of molecules in this entry. The entry contains 3213 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 BROMODOMAIN-CONTAINING PROTEIN 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	112	Total	C	N	O	S	0	7	0
			996	643	165	177	11			
1	B	107	Total	C	N	O	S	0	6	0
			942	611	158	163	10			
1	C	107	Total	C	N	O	S	0	0	0
			895	582	150	154	9			

There are 60 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	47	GLY	-	expression tag	UNP P25440
A	48	SER	-	expression tag	UNP P25440
A	49	SER	-	expression tag	UNP P25440
A	50	HIS	-	expression tag	UNP P25440
A	51	HIS	-	expression tag	UNP P25440
A	52	HIS	-	expression tag	UNP P25440
A	53	HIS	-	expression tag	UNP P25440
A	54	HIS	-	expression tag	UNP P25440
A	55	HIS	-	expression tag	UNP P25440
A	56	SER	-	expression tag	UNP P25440
A	57	SER	-	expression tag	UNP P25440
A	58	GLY	-	expression tag	UNP P25440
A	59	LEU	-	expression tag	UNP P25440
A	60	VAL	-	expression tag	UNP P25440
A	61	PRO	-	expression tag	UNP P25440
A	62	ARG	-	expression tag	UNP P25440
A	63	GLY	-	expression tag	UNP P25440
A	64	SER	-	expression tag	UNP P25440
A	65	HIS	-	expression tag	UNP P25440
A	66	MET	-	expression tag	UNP P25440
B	47	GLY	-	expression tag	UNP P25440
B	48	SER	-	expression tag	UNP P25440
B	49	SER	-	expression tag	UNP P25440

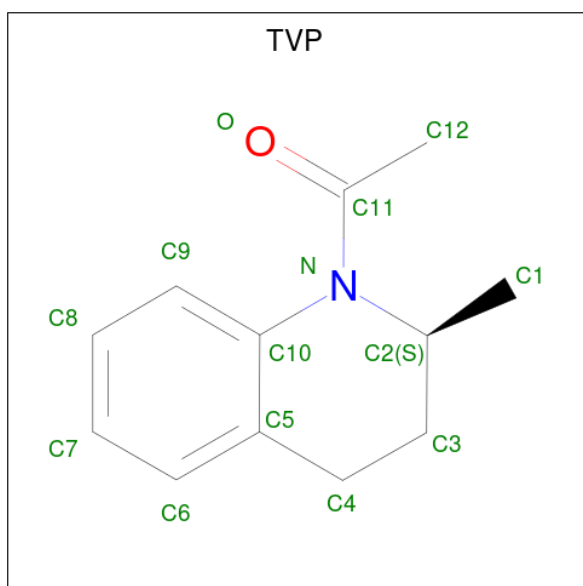*Continued on next page...*

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Chain	Residue	Modelled	Actual	Comment	Reference
B	50	HIS	-	expression tag	UNP P25440
B	51	HIS	-	expression tag	UNP P25440
B	52	HIS	-	expression tag	UNP P25440
B	53	HIS	-	expression tag	UNP P25440
B	54	HIS	-	expression tag	UNP P25440
B	55	HIS	-	expression tag	UNP P25440
B	56	SER	-	expression tag	UNP P25440
B	57	SER	-	expression tag	UNP P25440
B	58	GLY	-	expression tag	UNP P25440
B	59	LEU	-	expression tag	UNP P25440
B	60	VAL	-	expression tag	UNP P25440
B	61	PRO	-	expression tag	UNP P25440
B	62	ARG	-	expression tag	UNP P25440
B	63	GLY	-	expression tag	UNP P25440
B	64	SER	-	expression tag	UNP P25440
B	65	HIS	-	expression tag	UNP P25440
B	66	MET	-	expression tag	UNP P25440
C	47	GLY	-	expression tag	UNP P25440
C	48	SER	-	expression tag	UNP P25440
C	49	SER	-	expression tag	UNP P25440
C	50	HIS	-	expression tag	UNP P25440
C	51	HIS	-	expression tag	UNP P25440
C	52	HIS	-	expression tag	UNP P25440
C	53	HIS	-	expression tag	UNP P25440
C	54	HIS	-	expression tag	UNP P25440
C	55	HIS	-	expression tag	UNP P25440
C	56	SER	-	expression tag	UNP P25440
C	57	SER	-	expression tag	UNP P25440
C	58	GLY	-	expression tag	UNP P25440
C	59	LEU	-	expression tag	UNP P25440
C	60	VAL	-	expression tag	UNP P25440
C	61	PRO	-	expression tag	UNP P25440
C	62	ARG	-	expression tag	UNP P25440
C	63	GLY	-	expression tag	UNP P25440
C	64	SER	-	expression tag	UNP P25440
C	65	HIS	-	expression tag	UNP P25440
C	66	MET	-	expression tag	UNP P25440

- Molecule 2 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: $C_2H_6O_2$).

EDO

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

- Molecule 3 is (2S)-1-ACETYL-2-METHYL-1,2,3,4-TETRAHYDROQUINOLINE (three-letter code: TVP) (formula: $C_{12}H_{15}NO$).
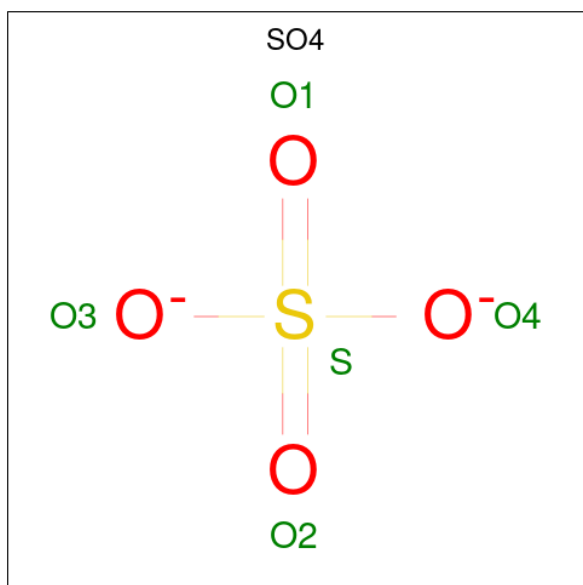

TVP

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total 14	C 12	N 1	O 1	0	0

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	B	1	Total	C	N	O	0	0
			14	12	1	1		
3	C	1	Total	C	N	O	0	0
			14	12	1	1		

- Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O_4S).

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

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	133	Total	O	0	0
			133	133		
5	B	102	Total	O	0	0
			102	102		
5	C	85	Total	O	0	0
			85	85		

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.
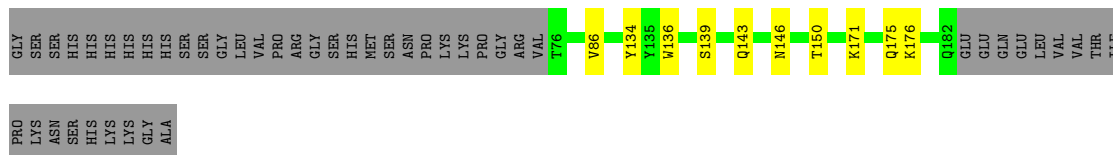
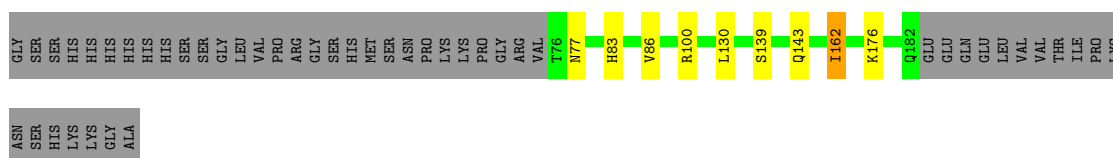• Molecule 1: BROMODOMAIN-CONTAINING PROTEIN 2

Chain A:
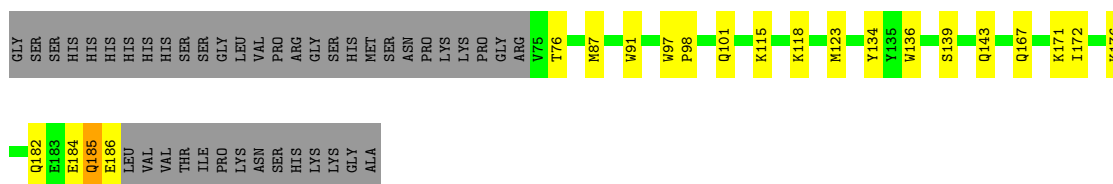
• Molecule 1: BROMODOMAIN-CONTAINING PROTEIN 2

Chain B:

• Molecule 1: BROMODOMAIN-CONTAINING PROTEIN 2

Chain C:

4 Data and refinement statistics ⓘ

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	114.23Å 55.53Å 67.97Å 90.00° 94.85° 90.00°	Depositor
Resolution (Å)	67.88 – 2.05 56.91 – 2.04	Depositor EDS
% Data completeness (in resolution range)	99.9 (67.88-2.05) 98.0 (56.91-2.04)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	3.56 (at 2.05Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.159 , 0.205 0.235 , 0.263	Depositor DCC
R_{free} test set	1342 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å^2)	30.1	Xtriage
Anisotropy	0.518	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.37 , 60.8	EDS
L-test for twinning[2]	$< |L| > = 0.49, < L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.93	EDS
Total number of atoms	3213	wwPDB-VP
Average B, all atoms (Å^2)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.95% of the height of the origin peak. No significant pseudotranslation is detected.*

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[1]Intensities estimated from amplitudes.
[2]Theoretical values of $< |L| >$, $< L^2 >$ 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: TVP, SO4, EDO

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	0.57	1/1021 (0.1%)	0.59	0/1380
1	B	0.54	1/967 (0.1%)	0.56	0/1308
1	C	0.54	0/920	0.55	0/1245
All	All	0.55	2/2908 (0.1%)	0.57	0/3933

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	136	TRP	CD2-CE2	5.12	1.47	1.41
1	A	136	TRP	CD2-CE2	5.05	1.47	1.41

There are no bond angle outliers.

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	996	0	987	17	1
1	B	942	0	943	5	2
1	C	895	0	895	5	0
2	A	8	0	12	0	0
3	A	14	0	15	2	0
3	B	14	0	15	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	C	14	0	15	3	0
4	A	5	0	0	0	0
4	C	5	0	0	0	0
5	A	133	0	0	4	1
5	B	102	0	0	0	0
5	C	85	0	0	1	0
All	All	3213	0	2882	34	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 6.

The worst 5 of 34 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:76[B]:THR:HG21	1:A:184:GLU:HG3	1.30	1.08
1:A:118:LYS:O	5:A:2070:HOH:O	1.84	0.95
1:A:91:TRP:HA	1:A:123[B]:MET:HE1	1.51	0.92
1:A:76[B]:THR:HG21	1:A:184:GLU:CG	2.01	0.89
1:A:87[B]:MET:HE3	1:A:91:TRP:HB2	1.58	0.86

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)
1:B:134:TYR:OH	5:A:2062:HOH:O[4_555]	1.89	0.31
1:A:115:LYS:NZ	1:B:134:TYR:O[4_545]	2.14	0.06

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	117/154 (76%)	117 (100%)	0	0	100	100
1	B	111/154 (72%)	111 (100%)	0	0	100	100
1	C	105/154 (68%)	104 (99%)	1 (1%)	0	100	100
All	All	333/462 (72%)	332 (100%)	1 (0%)	0	100	100

There are no Ramachandran outliers to report.

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	110/139 (79%)	107 (97%)	3 (3%)	44	38
1	B	104/139 (75%)	104 (100%)	0	100	100
1	C	98/139 (70%)	96 (98%)	2 (2%)	55	50
All	All	312/417 (75%)	307 (98%)	5 (2%)	60	59

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	101	GLN
1	A	182	GLN
1	A	185	GLN
1	C	77	ASN
1	C	162	ILE

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

Mol	Chain	Res	Type
1	B	133	ASN
1	B	175	GLN
1	C	80	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no monosaccharides in this entry.

5.6 Ligand geometry ⓘ

7 ligands are 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	$\#|Z| > 2$
4	SO4	C	1184	-	4,4,4	0.22	0	6,6,6	0.37	0
3	TVP	C	1183	-	15,15,15	0.50	0	21,21,21	0.92	0
4	SO4	A	1190	-	4,4,4	0.36	0	6,6,6	0.15	0
2	EDO	A	1187	-	3,3,3	0.24	0	2,2,2	0.54	0
3	TVP	A	1189	-	15,15,15	0.43	0	21,21,21	0.99	1 (4%)
3	TVP	B	1183	-	15,15,15	0.34	0	21,21,21	0.89	1 (4%)
2	EDO	A	1188	-	3,3,3	0.32	0	2,2,2	0.55	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	TVP	C	1183	-	-	0/4/17/17	0/2/2/2
2	EDO	A	1187	-	-	0/1/1/1	-
3	TVP	A	1189	-	-	0/4/17/17	0/2/2/2

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	TVP	B	1183	-	-	0/4/17/17	0/2/2/2
2	EDO	A	1188	-	-	1/1/1/1	-

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	A	1189	TVP	C3-C4-C5	-2.90	107.37	112.87
3	B	1183	TVP	C3-C4-C5	-2.06	108.96	112.87

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1188	EDO	O1-C1-C2-O2

There are no ring outliers.

3 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	1183	TVP	3	0
3	A	1189	TVP	2	0
3	B	1183	TVP	3	0

5.7 Other polymers ⓘ

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 ⓘ

Unable to reproduce the depositors R factor - this section is therefore empty.

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

Unable to reproduce the depositors R factor - this section is therefore empty.

6.3   Carbohydrates ⓘ

Unable to reproduce the depositors R factor - this section is therefore empty.

6.4   Ligands ⓘ

Unable to reproduce the depositors R factor - this section is therefore empty.

6.5   Other polymers ⓘ

Unable to reproduce the depositors R factor - this section is therefore empty.