wwPDB X-ray Structure Validation Summary Report (i)

Oct 5, 2023 – 05:21 AM EDT

PDB ID	:	6V0Q
Title	:	Crystal structure of the bromodomain of human BRD7 bound to TG003
Authors	:	Karim, M.R.; Chan, A.; Schonbrunn, E.
Deposited on	:	2019-11-19
Resolution	:	1.69 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

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

MolProbity	:	FAILED
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	FAILED
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)	:	2.35.1

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 1.69 Å.

There are no overall percentile quality scores available for this entry.

MolProbity and EDS failed to run properly - the sequence quality summary graphics cannot be shown.

2 Entry composition (i)

There are 7 unique types of molecules in this entry. The entry contains 4503 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.

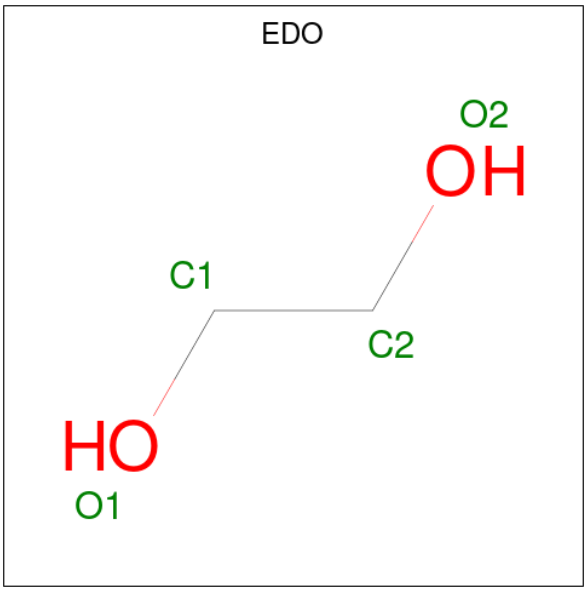
- Molecule 1 is a protein called Bromodomain-containing protein 7.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	119	Total	C	N	O	S	0	3	0
			988	633	161	184	10			
1	B	117	Total	C	N	O	S	0	4	0
			977	628	159	180	10			
1	C	117	Total	C	N	O	S	0	4	0
			977	627	159	180	11			
1	D	117	Total	C	N	O	S	0	5	0
			978	627	159	182	10			

There are 8 discrepancies between the modelled and reference sequences:

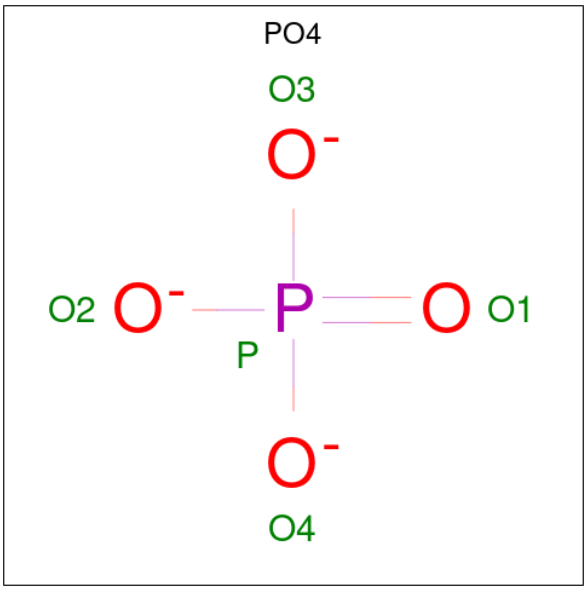
Chain	Residue	Modelled	Actual	Comment	Reference
A	128	GLU	-	expression tag	UNP Q9NPI1
A	129	SER	-	expression tag	UNP Q9NPI1
B	128	GLU	-	expression tag	UNP Q9NPI1
B	129	SER	-	expression tag	UNP Q9NPI1
C	128	GLU	-	expression tag	UNP Q9NPI1
C	129	SER	-	expression tag	UNP Q9NPI1
D	128	GLU	-	expression tag	UNP Q9NPI1
D	129	SER	-	expression tag	UNP Q9NPI1

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



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

- Molecule 3 is PHOSPHATE ION (three-letter code: PO4) (formula: O_4P).

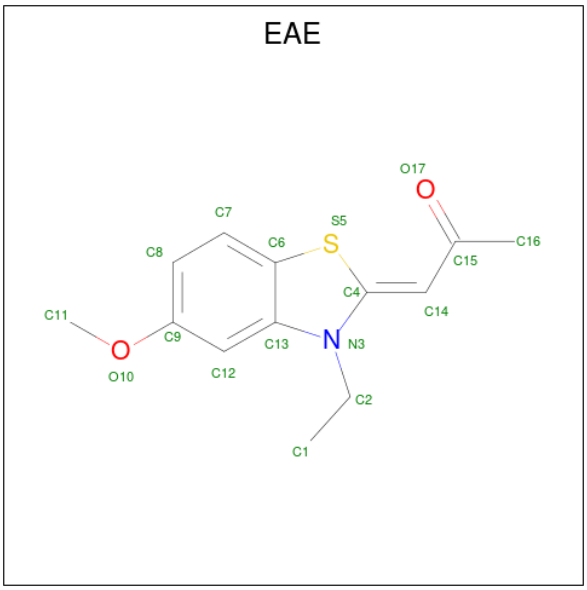


Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total O P 5 4 1	0	0

- Molecule 4 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

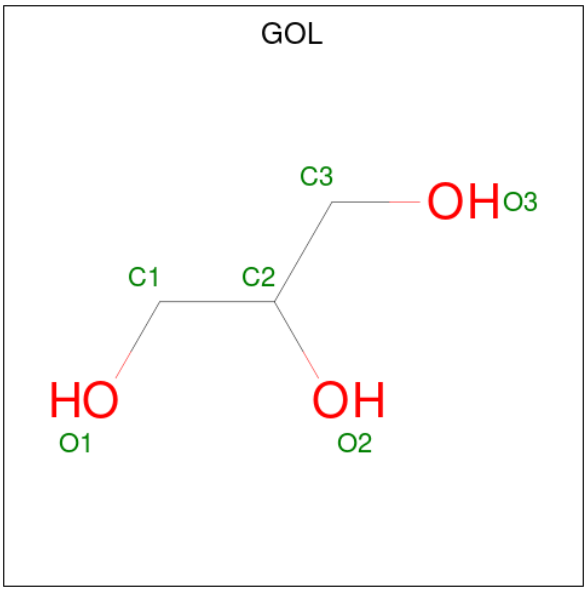
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total Cl 1 1	0	0

- Molecule 5 is (1 {Z})-1-(3-ethyl-5-methoxy-1,3-benzothiazol-2-ylidene)propan-2-one (three-letter code: EAE) (formula: $C_{13}H_{15}NO_2S$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total C N O S 17 13 1 2 1	0	0
5	B	1	Total C N O S 17 13 1 2 1	0	0
5	C	1	Total C N O S 17 13 1 2 1	0	0
5	D	1	Total C N O S 17 13 1 2 1	0	0

- Molecule 6 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	1	Total C O 6 3 3	0	0

- Molecule 7 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	142	Total O 142 142	0	0
7	B	138	Total O 138 138	0	0
7	C	92	Total O 92 92	0	0
7	D	115	Total O 115 115	0	0

MolProbity and EDS failed to run properly - this section is therefore empty.

3 Data and refinement statistics (i)

EDS failed to run properly - this section is therefore incomplete.

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	59.97Å 36.78Å 118.96Å 90.00° 101.77° 90.00°	Depositor
Resolution (Å)	37.68 – 1.69	Depositor
% Data completeness (in resolution range)	98.5 (37.68-1.69)	Depositor
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	6.10 (at 1.69Å)	Xtriage
Refinement program	PHENIX 1.12	Depositor
R, R_{free}	0.181 , 0.218	Depositor
Wilson B-factor (Å^2)	19.7	Xtriage
Anisotropy	0.242	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	0.006 for h,-k,-h-l	Xtriage
Total number of atoms	4503	wwPDB-VP
Average B, all atoms (Å^2)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 34.86 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 6.3822e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

[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.

4 Model quality

4.1 Standard geometry

MolProbity failed to run properly - this section is therefore empty.

4.2 Too-close contacts

MolProbity failed to run properly - this section is therefore empty.

4.3 Torsion angles

4.3.1 Protein backbone

MolProbity failed to run properly - this section is therefore empty.

4.3.2 Protein sidechains

MolProbity failed to run properly - this section is therefore empty.

4.3.3 RNA

MolProbity failed to run properly - this section is therefore empty.

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

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

4.5 Carbohydrates

There are no monosaccharides in this entry.

4.6 Ligand geometry

Of 11 ligands modelled in this entry, 1 is monoatomic - leaving 10 for Mogul analysis.

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	EDO	C	301	-	3,3,3	0.44	0	2,2,2	0.37	0
5	EAE	B	301	-	18,18,18	1.10	0	24,25,25	1.91	8 (33%)
3	PO4	A	303	-	4,4,4	0.85	0	6,6,6	0.46	0
5	EAE	C	302	-	18,18,18	1.15	1 (5%)	24,25,25	2.48	11 (45%)
6	GOL	A	306	-	5,5,5	1.03	0	5,5,5	1.31	0
2	EDO	A	302	-	3,3,3	0.31	0	2,2,2	0.65	0
5	EAE	A	305	-	18,18,18	1.07	0	24,25,25	2.18	11 (45%)
5	EAE	D	302	-	18,18,18	1.14	0	24,25,25	2.80	12 (50%)
2	EDO	A	301	-	3,3,3	0.46	0	2,2,2	0.37	0
2	EDO	D	301	-	3,3,3	0.47	0	2,2,2	0.30	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
2	EDO	C	301	-	-	0/1/1/1	-
5	EAE	B	301	-	-	0/8/8/8	0/2/2/2
5	EAE	C	302	-	-	3/8/8/8	0/2/2/2
6	GOL	A	306	-	-	2/4/4/4	-
2	EDO	A	302	-	-	1/1/1/1	-
5	EAE	A	305	-	-	1/8/8/8	0/2/2/2
5	EAE	D	302	-	-	4/8/8/8	0/2/2/2
2	EDO	A	301	-	-	0/1/1/1	-
2	EDO	D	301	-	-	1/1/1/1	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	C	302	EAE	C12-C13	2.13	1.43	1.39

The worst 5 of 42 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
5	D	302	EAE	C6-S5-C4	5.89	96.30	91.24
5	C	302	EAE	C14-C4-N3	4.58	129.15	124.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
5	C	302	EAE	C6-S5-C4	4.53	95.13	91.24
5	D	302	EAE	C14-C4-N3	4.39	128.95	124.19
5	C	302	EAE	C13-C6-S5	-4.23	105.63	110.87

There are no chirality outliers.

5 of 12 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	C	302	EAE	C1-C2-N3-C4
5	D	302	EAE	C1-C2-N3-C4
6	A	306	GOL	O1-C1-C2-O2
5	C	302	EAE	C12-C9-O10-C11
5	C	302	EAE	C8-C9-O10-C11

There are no ring outliers.

No monomer is involved in short contacts.

4.7 Other polymers (i)

There are no such residues in this entry.

4.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

5 Fit of model and data

5.1 Protein, DNA and RNA chains

EDS failed to run properly - this section is therefore empty.

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

EDS failed to run properly - this section is therefore empty.

5.3 Carbohydrates

EDS failed to run properly - this section is therefore empty.

5.4 Ligands

EDS failed to run properly - this section is therefore empty.

5.5 Other polymers

EDS failed to run properly - this section is therefore empty.