

# wwPDB X-ray Structure Validation Summary Report (i)

Dec 8, 2023 – 03:11 am GMT

PDB ID : 1W80

Title: Crystal structure of the alpha-adaptin appendage domain, from the AP2 adap-

tor complex, bound to 2 peptides from Synaptojanin170

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Deposited on : 2004-09-14

Resolution : 1.90 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 (1)) were used in the production of this report:

Mol Probity : 4.02b-467

Mogul : 1.8.4, CSD as541be (2020)

Xtriage (Phenix) : 1.13 EDS : 2.36

Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)

 $Refmac \quad : \quad 5.8.0158$ 

CCP4 : 7.0.044 (Gargrove) roteins) : Engh & Huber (2001)

Ideal geometry (proteins) : Engh & Huber (2001) Ideal geometry (DNA, RNA) : Parkinson et al. (1996)

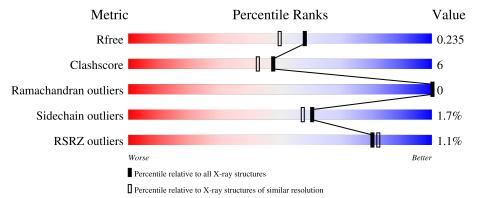
Validation Pipeline (wwPDB-VP) : 2.36

## 1 Overall quality at a glance (i)

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 $(\# \mathrm{Entries})$	$\begin{array}{c} {\rm Similar\ resolution} \\ (\#{\rm Entries},{\rm resolution\ range}(\mathring{\rm A})) \end{array}$
$R_{free}$	130704	6207 (1.90-1.90)
Clashscore	141614	6847 (1.90-1.90)
Ramachandran outliers	138981	6760 (1.90-1.90)
Sidechain outliers	138945	6760 (1.90-1.90)
RSRZ outliers	127900	6082 (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 <=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	250	84%		13% •••				
2	Р	12	58%	8%	33%				
	0		8%						
3	Q	12	58%	8%	33%				



## 2 Entry composition (i)

There are 8 unique types of molecules in this entry. The entry contains 2439 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 ADAPTER-RELATED PROTEIN COMPLEX 2 ALPHA 2 SUBUNIT.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
1	A	248	Total 1995	C 1268	N 341	O 377	S 9	0	4	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	889	GLY	VAL	conflict	UNP P17247
A	890	ALA	LEU	conflict	UNP P17247

• Molecule 2 is a protein called SYNAPTOJANIN 1.

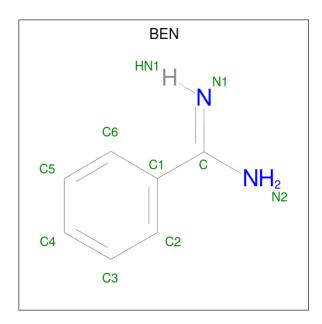
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf	Trace		
2	Р	8	Total 68	C 47		O 11	0	0	0

• Molecule 3 is a protein called SYNAPTOJANIN 1.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf	Trace		
3	Q	8	Total 65	C 41	N 9	O 15	0	0	0

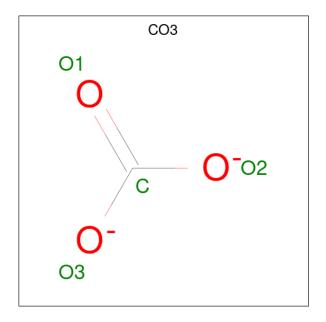
• Molecule 4 is BENZAMIDINE (three-letter code: BEN) (formula: C<sub>7</sub>H<sub>8</sub>N<sub>2</sub>).





Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total 9	C 7	N 2	0	0

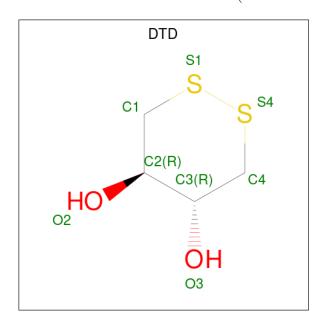
 $\bullet$  Molecule 5 is CARBONATE ION (three-letter code: CO3) (formula: CO3).



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

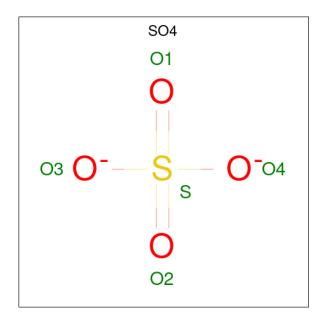


 $\bullet$  Molecule 6 is DITHIANE DIOL (three-letter code: DTD) (formula:  $\mathrm{C_4H_8O_2S_2}).$ 



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	A	1	Total 8	C 4	O 2	S 2	0	0

 $\bullet$  Molecule 7 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



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

• Molecule 8 is water.



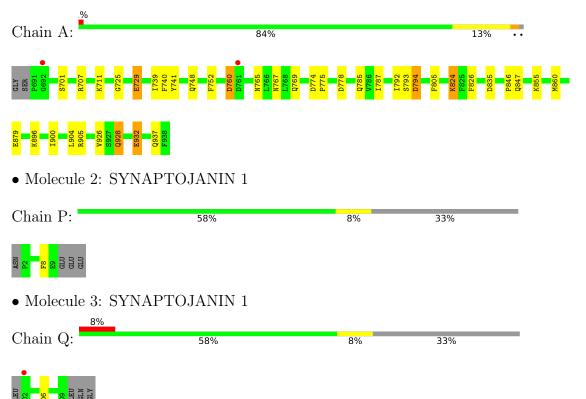
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	A	263	Total O 263 263	0	0
8	Р	7	Total O 7 7	0	0
8	Q	7	Total O 7 7	0	0



## 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: ADAPTER-RELATED PROTEIN COMPLEX 2 ALPHA 2 SUBUNIT





## 4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants	146.60Å 67.32Å 39.72Å	Donositor
a, b, c, $\alpha$ , $\beta$ , $\gamma$	90.00° 94.53° 90.00°	Depositor
Resolution (Å)	72.55 - 1.90	Depositor
Resolution (A)	39.60 - 1.90	EDS
% Data completeness	95.0 (72.55-1.90)	Depositor
(in resolution range)	95.0 (39.60-1.90)	EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	0.09	Depositor
$< I/\sigma(I) > 1$	1.39 (at 1.89Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
D D.	0.177 , 0.224	Depositor
$R, R_{free}$	0.188 , 0.235	DCC
$R_{free}$ test set	1467 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	27.3	Xtriage
Anisotropy	0.069	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$ , $B_{sol}(Å^2)$	0.34, 54.9	EDS
L-test for twinning <sup>2</sup>	$ < L >=0.49, < L^2>=0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	2439	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 7.19% of the height of the origin peak. No significant pseudotranslation is detected.

<sup>&</sup>lt;sup>2</sup>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.



<sup>&</sup>lt;sup>1</sup>Intensities estimated from amplitudes.

# 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: CO3, BEN, SO4, DTD

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	Bo	nd lengths	Bond angles		
IVIOI	Chain	RMSZ	# Z  > 5	RMSZ	# Z  > 5	
1	A	1.37	11/2034~(0.5%)	1.12	6/2754 (0.2%)	
2	Р	1.34	0/71	0.89	0/95	
3	Q	1.39	0/66	0.99	0/86	
All	All	1.37	$11/2171 \ (0.5\%)$	1.11	$6/2935 \ (0.2\%)$	

The worst 5 of 11 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	$\mathbf{Z}$	Observed(A)	Ideal(Å)
1	A	729	GLU	CB-CG	-7.38	1.38	1.52
1	A	879	GLU	CG-CD	7.07	1.62	1.51
1	A	896	LYS	CE-NZ	6.04	1.64	1.49
1	A	701	SER	CB-OG	5.82	1.49	1.42
1	A	826	PHE	CD1-CE1	-5.82	1.27	1.39

The worst 5 of 6 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^o)$	$\operatorname{Ideal}({}^o)$
1	A	794	ASP	CB-CG-OD1	-7.42	111.63	118.30
1	A	860	MET	CG-SD-CE	5.88	109.60	100.20
1	A	904	LEU	CA-CB-CG	5.71	128.44	115.30
1	A	794	ASP	CB-CG-OD2	5.47	123.22	118.30
1	A	824	LYS	CD-CE-NZ	-5.35	99.40	111.70

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	1995	0	1988	26	0
2	Р	68	0	64	1	0
3	Q	65	0	50	0	0
4	A	9	0	7	0	0
5	A	12	0	0	0	0
6	A	8	0	8	3	0
7	A	5	0	0	0	0
8	A	263	0	0	9	0
8	Р	7	0	0	0	0
8	Q	7	0	0	0	0
All	All	2439	0	2117	26	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 6.

The worst 5 of 26 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic	$\operatorname{Clash}$	
Atom-1	Atom-2	${\rm distance}({\rm \AA})$	$overlap (\AA)$	
1:A:729:GLU:HG3	6:A:1943:DTD:H11	1.70	0.73	
1:A:900:ILE:HD13	1:A:926:VAL:HG21	1.74	0.69	
1:A:835:ASP:OD2	8:A:2154:HOH:O	2.11	0.69	
1:A:729:GLU:OE1	6:A:1943:DTD:O2	2.14	0.66	
1:A:928[B]:GLN:NE2	8:A:2238:HOH:O	2.18	0.65	

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 nu	imber of residues.
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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Perce	ntiles
1	A	$250/250 \; (100\%)$	246 (98%)	4 (2%)	0	100	100
2	Р	6/12 (50%)	6 (100%)	0	0	100	100
3	Q	6/12 (50%)	6 (100%)	0	0	100	100
All	All	262/274 (96%)	258 (98%)	4 (2%)	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	Analysed Rotameric Outlie		Percentiles
1	A	224/221 (101%)	221 (99%)	3 (1%)	69 68
2	Р	7/11 (64%)	7 (100%)	0	100 100
3	Q	7/10 (70%)	6 (86%)	1 (14%)	3 1
All	All	238/242 (98%)	234 (98%)	4 (2%)	60 57

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	760	ASP
1	A	794	ASP
1	A	932	GLU
3	Q	6	ASP

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

Mol	Chain	Res	Type
1	A	765	ASN
1	A	785	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 monosaccharides in this entry.

#### 5.6 Ligand geometry (i)

6 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	В	ond leng	$_{ m gths}$	В	ond ang	gles
MIOI	Type	Chain	nes	Lilik	Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
7	SO4	A	1944	-	4,4,4	0.24	0	6,6,6	0.55	0
5	CO3	A	1941	-	2,3,3	0.57	0	2,3,3	1.50	1 (50%)
6	DTD	A	1943	-	6,8,8	0.77	0	6,10,10	1.23	0
5	CO3	A	1940	-	2,3,3	0.98	0	2,3,3	1.82	1 (50%)
5	CO3	A	1942	-	2,3,3	0.31	0	2,3,3	0.49	0
4	BEN	A	1939	-	9,9,9	2.03	2 (22%)	7,11,11	1.19	1 (14%)

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.

$\mathbf{Mol}$	Type	Chain	$\operatorname{Res}$	Link	Chirals	Torsions	Rings
6	DTD	A	1943	-	-	-	0/0/1/1
4	BEN	A	1939	-	-	2/4/4/4	0/1/1/1



#### All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	$\mathbf{Z}$	$\operatorname{Observed}(\text{\AA})$	$\operatorname{Ideal}( ext{\AA})$
4	A	1939	BEN	C1-C	4.47	1.55	1.47
4	A	1939	BEN	C2-C1	2.79	1.44	1.39

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	$\mathbf{Z}$	$\mathbf{Observed}(^o)$	$\operatorname{Ideal}({}^{o})$
4	A	1939	BEN	C1-C-N2	-2.70	113.97	118.05
5	A	1940	CO3	O3-C-O1	-2.26	113.68	119.55
5	A	1941	CO3	O2-C-O1	2.05	124.86	119.55

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	1939	BEN	N2-C-C1-C2
4	A	1939	BEN	N2-C-C1-C6

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	$\operatorname{Res}$	Type	Clashes	Symm-Clashes
6	A	1943	DTD	3	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,  $95^{th}$  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>	$\# \mathrm{RSRZ}{>}2$	$\mathrm{OWAB}(\mathrm{\AA}^2)$	Q < 0.9
1	A	248/250 (99%)	-0.25	2 (0%) 86 87	20, 30, 46, 66	0
2	P	8/12 (66%)	0.12	0 100 100	33, 33, 57, 58	0
3	Q	8/12 (66%)	0.30	1 (12%) 3 4	34, 45, 48, 53	0
All	All	264/274 (96%)	-0.22	3 (1%) 80 82	20, 31, 47, 66	0

#### All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	761	ASP	2.9
1	A	692	GLY	2.9
3	Q	2	ASP	2.9

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

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

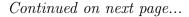
### 6.3 Carbohydrates (i)

There are no monosaccharides 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,  $95^{th}$  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	${f B-factors}({f \AA}^2)$	Q<0.9
6	DTD	A	1943	8/8	0.82	0.27	22,23,30,35	8





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Mol	Type	Chain	Res	Atoms	RSCC	RSR	${f B-factors}({f \AA}^2)$	Q < 0.9
5	CO3	A	1942	4/4	0.91	0.14	59,62,62,63	0
7	SO4	A	1944	5/5	0.92	0.11	70,73,74,76	0
5	CO3	A	1941	4/4	0.93	0.11	64,64,64,64	0
4	BEN	A	1939	9/9	0.93	0.10	26,28,31,32	0
5	CO3	A	1940	4/4	0.97	0.07	35,35,36,39	0

## 6.5 Other polymers (i)

There are no such residues in this entry.

