



wwPDB X-ray Structure Validation Summary Report ⓘ

May 26, 2020 – 10:33 am BST

PDB ID : 1JEN
Title : HUMAN S-ADENOSYLMETHIONINE DECARBOXYLASE
Authors : Ekstrom, J.L.; Mathews, I.I.; Stanley, B.A.; Pegg, A.E.; Ealick, S.E.
Deposited on : 1999-02-23
Resolution : 2.25 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 ⓘ symbol.

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)
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

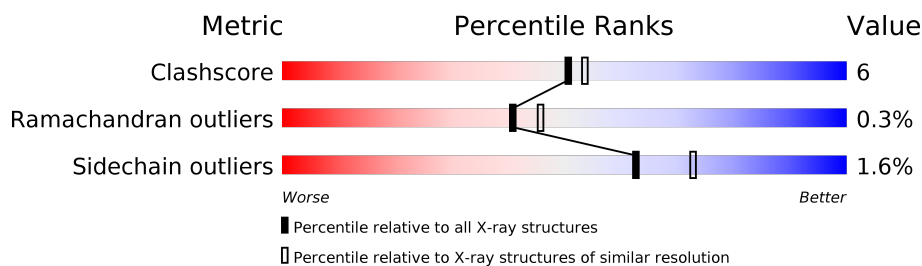
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.25 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	1487 (2.26-2.26)
Ramachandran outliers	138981	1449 (2.26-2.26)
Sidechain outliers	138945	1450 (2.26-2.26)

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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	B	67	
1	D	67	
2	A	267	
2	C	267	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5333 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 PROTEIN (S-ADENOSYLMETHIONINE DECARBOXY-LASE (BETA CHAIN)).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	57	Total	C	N	O	S	0	0	0
			464	296	76	91	1			
1	D	54	Total	C	N	O	S	0	0	0
			438	282	70	85	1			

- Molecule 2 is a protein called PROTEIN (S-ADENOSYLMETHIONINE DECARBOXY-LASE (ALPHA CHAIN)).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	A	249	Total	C	N	O	S	0	0	0
			2013	1296	325	375	17			
2	C	248	Total	C	N	O	S	0	0	0
			2004	1291	323	373	17			

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	45	Total	O	0	0
			45	45		
3	A	158	Total	O	0	0
			158	158		
3	D	46	Total	O	0	0
			46	46		
3	C	165	Total	O	0	0
			165	165		

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.

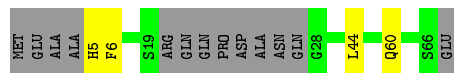
• Molecule 1: PROTEIN (S-ADENOSYLMETHIONINE DECARBOXYLASE (BETA CHAIN))

Chain B: 




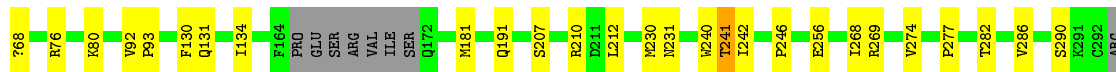
• Molecule 1: PROTEIN (S-ADENOSYLMETHIONINE DECARBOXYLASE (BETA CHAIN))

Chain D: 




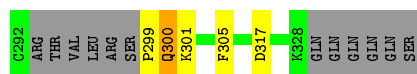
• Molecule 2: PROTEIN (S-ADENOSYLMETHIONINE DECARBOXYLASE (ALPHA CHAIN))

Chain A: 



• Molecule 2: PROTEIN (S-ADENOSYLMETHIONINE DECARBOXYLASE (ALPHA CHAIN))

Chain C: 



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	74.60 Å 55.80 Å 90.10 Å 90.00° 109.60° 90.00°	Depositor
Resolution (Å)	20.00 – 2.25	Depositor
% Data completeness (in resolution range)	98.9 (20.00-2.25)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	0.05	Depositor
Refinement program	X-PLOR 3.8	Depositor
R, R_{free}	0.177 , 0.229	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	5333	wwPDB-VP
Average B, all atoms (Å ²)	20.0	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: PYR

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	B	0.51	0/472	0.66	0/634
1	D	0.44	0/446	0.68	0/601
2	A	0.51	0/2058	0.69	1/2773 (0.0%)
2	C	0.52	0/2049	0.68	0/2761
All	All	0.51	0/5025	0.68	1/6769 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	246	PRO	N-CA-C	5.04	125.20	112.10

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	B	464	0	455	8	0
1	D	438	0	431	6	0
2	A	2013	0	1960	28	0
2	C	2004	0	1952	25	0
3	A	158	0	0	4	0
3	B	45	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	C	165	0	0	2	0
3	D	46	0	0	1	0
All	All	5333	0	4798	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 6.

The worst 5 of 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)
2:A:131:GLN:HE22	2:C:153:ASN:HD21	1.23	0.86
2:A:68:PYR:H33	2:A:80:LYS:O	1.79	0.82
2:A:131:GLN:HE22	2:C:153:ASN:ND2	1.78	0.80
2:C:68:PYR:H33	2:C:80:LYS:O	1.82	0.79
2:A:300:GLN:HA	2:A:300:GLN:HE21	1.50	0.77

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	B	53/67 (79%)	52 (98%)	0	1 (2%)	8	4
1	D	50/67 (75%)	50 (100%)	0	0	100	100
2	A	242/267 (91%)	232 (96%)	10 (4%)	0	100	100
2	C	241/267 (90%)	232 (96%)	8 (3%)	1 (0%)	34	37
All	All	586/668 (88%)	566 (97%)	18 (3%)	2 (0%)	41	46

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	5	HIS
2	C	300	GLN

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	B	52/60 (87%)	51 (98%)	1 (2%)	57	66
1	D	50/60 (83%)	49 (98%)	1 (2%)	55	64
2	A	225/243 (93%)	221 (98%)	4 (2%)	59	68
2	C	224/243 (92%)	221 (99%)	3 (1%)	69	79
All	All	551/606 (91%)	542 (98%)	9 (2%)	62	73

5 of 9 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	A	317	ASP
2	C	317	ASP
2	C	241	THR
2	A	300	GLN
1	D	44	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 10 such sidechains are listed below:

Mol	Chain	Res	Type
2	A	311	GLN
1	D	5	HIS
2	C	131	GLN
2	A	300	GLN
1	D	60	GLN

5.3.3 RNA ⓘ

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](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
2	A	1
2	C	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	A	68:PYR	C2	69:SER	N	2.36
1	C	68:PYR	C2	69:SER	N	2.35

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.