



Full wwPDB X-ray Structure Validation Report ⓘ

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PDB ID : 3ECM
Title : Crystal structure of the unliganded PDE8A catalytic domain
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Deposited on : 2008-09-01
Resolution : 1.90 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
<http://wwpdb.org/validation/2016/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 : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

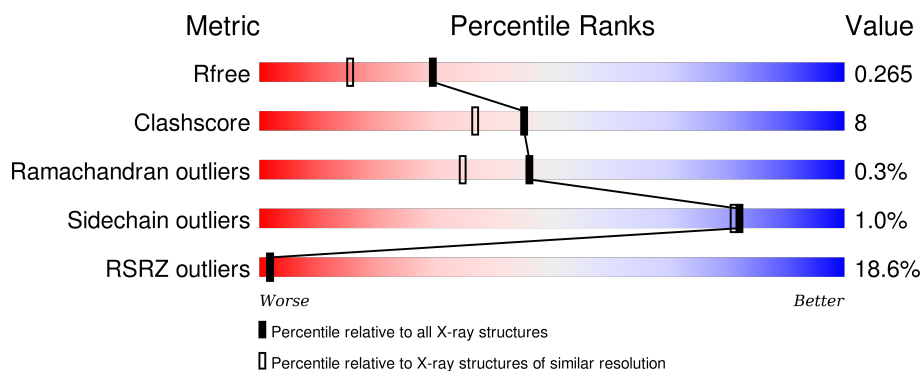
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION


The reported resolution of this entry is 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 (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	91344	4755 (1.90-1.90)
Clashscore	102246	5398 (1.90-1.90)
Ramachandran outliers	100387	5338 (1.90-1.90)
Sidechain outliers	100360	5339 (1.90-1.90)
RSRZ outliers	91569	4766 (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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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 $\leq 5\%$. 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	338	

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	MG	A	2	-	-	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 2814 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 High affinity cAMP-specific and IBMX-insensitive 3',5'-cyclic phosphodiesterase 8A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	338	Total	C	N	O	S	0	0	0
			2744	1734	464	527	19			

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Zn	0	0
			1	1		

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Mg	0	0
			1	1		

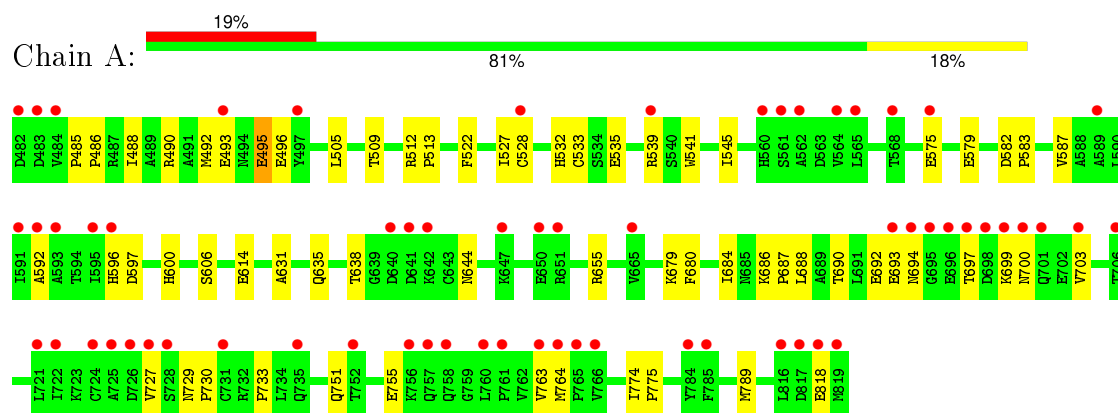
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	68	Total	O	0	0
			68	68		

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 errors 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: High affinity cAMP-specific and IBMX-insensitive 3',5'-cyclic phosphodiesterase 8A



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	76.26Å 132.66Å 101.52Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 1.90 30.49 – 1.90	Depositor EDS
% Data completeness (in resolution range)	(Not available) (30.00-1.90) 86.8 (30.49-1.90)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.31 (at 1.91Å)	Xtriage
Refinement program	CNS1.2	Depositor
R, R_{free}	0.236 , 0.265 0.235 , 0.265	Depositor DCC
R_{free} test set	3532 reflections (9.96%)	DCC
Wilson B-factor (Å ²)	32.2	Xtriage
Anisotropy	0.811	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 50.5	EDS
Estimated twinning fraction	0.016 for 1/2*h-1/2*k,-3/2*h-1/2*k,-l 0.023 for 1/2*h+1/2*k,3/2*h-1/2*k,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 36323 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	2814	wwPDB-VP
Average B, all atoms (Å ²)	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.89% of the height of the origin peak. No significant pseudotranslation is detected.*

¹ Intensities estimated from amplitudes.

² Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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: ZN, MG

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	0.33	0/2809	0.52	0/3811

There are no bond length outliers.

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	2744	0	2631	42	0
2	A	1	0	0	0	0
3	A	1	0	0	0	0
4	A	68	0	0	2	0
All	All	2814	0	2631	42	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 8.

All (42) 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:694:ASN:HB2	1:A:700:ASN:ND2	2.00	0.76
1:A:680:PHE:CD1	1:A:684:ILE:HD12	2.25	0.71
1:A:697:THR:HG22	1:A:699:LYS:H	1.60	0.67
1:A:492:MET:O	1:A:495:GLU:HG3	1.96	0.66
1:A:763:VAL:HG23	1:A:764:MET:HG3	1.80	0.63
1:A:485:PRO:HG2	1:A:488:ILE:HD12	1.81	0.63
1:A:528:CYS:HB3	1:A:533:CYS:O	2.00	0.61
1:A:686:LYS:HB2	1:A:687:PRO:HD3	1.84	0.60
1:A:690:THR:O	1:A:693:GLU:HG2	2.03	0.59
1:A:575:GLU:O	1:A:579:GLU:HG2	2.02	0.59
1:A:606:SER:HB3	1:A:763:VAL:HG11	1.84	0.58
1:A:490:ARG:HA	1:A:493:GLU:HG3	1.85	0.58
1:A:697:THR:HB	1:A:700:ASN:HB2	1.88	0.56
1:A:688:LEU:O	1:A:692:GLU:HG3	2.05	0.56
1:A:583:PRO:O	1:A:587:VAL:HG23	2.06	0.55
1:A:697:THR:HB	1:A:700:ASN:CB	2.36	0.55
1:A:699:LYS:O	1:A:703:VAL:HG23	2.06	0.54
1:A:694:ASN:HB2	1:A:700:ASN:HD22	1.72	0.54
1:A:751:GLN:O	1:A:755:GLU:HG3	2.07	0.54
1:A:486:PRO:O	1:A:490:ARG:HG3	2.08	0.53
1:A:729:ASN:HB2	1:A:730:PRO:HD3	1.91	0.53
1:A:614:GLU:HG3	4:A:856:HOH:O	2.10	0.52
1:A:631:ALA:O	1:A:635:GLN:HG3	2.11	0.50
1:A:582:ASP:HB3	1:A:583:PRO:HD2	1.94	0.49
1:A:522:PHE:CG	1:A:527:ILE:HD12	2.48	0.48
1:A:528:CYS:O	1:A:532:HIS:N	2.45	0.48
1:A:592:ALA:O	1:A:596:HIS:HB3	2.15	0.46
1:A:774:ILE:N	1:A:775:PRO:HD2	2.30	0.46
1:A:541:TRP:O	1:A:545:ILE:HG12	2.16	0.45
1:A:505:LEU:O	1:A:509:THR:HG23	2.17	0.45
1:A:512:ARG:N	1:A:513:PRO:CD	2.80	0.44
1:A:485:PRO:CG	1:A:488:ILE:HD12	2.47	0.44
1:A:535:GLU:O	1:A:539:ARG:HG3	2.17	0.44
1:A:727:VAL:HG21	1:A:789:MET:HE1	1.99	0.44
1:A:727:VAL:CG2	1:A:789:MET:HE1	2.47	0.43
1:A:486:PRO:HB2	1:A:490:ARG:NH1	2.34	0.42
1:A:755:GLU:OE2	1:A:763:VAL:HG22	2.20	0.42
1:A:655:ARG:HD3	1:A:655:ARG:HA	1.87	0.42
1:A:597:ASP:O	1:A:600:HIS:HB2	2.20	0.41
1:A:638:THR:HA	1:A:644:ASN:ND2	2.34	0.41
1:A:575:GLU:OE1	1:A:575:GLU:HA	2.20	0.41
1:A:679:LYS:HG3	4:A:849:HOH:O	2.20	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	336/338 (99%)	314 (94%)	21 (6%)	1 (0%)	46 35

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	818	GLU

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	303/303 (100%)	300 (99%)	3 (1%)	82 81

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	495	GLU
1	A	496	GLU
1	A	733	PRO

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	694	ASN
1	A	700	ASN
1	A	705	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](#)

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

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.

No monomer is involved in short contacts.

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 ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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, 95th 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>2	OWAB(Å ²)	Q<0.9
1	A	338/338 (100%)	0.98	63 (18%) 2 2	29, 43, 79, 95	0

All (63) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	819	MET	8.9
1	A	758	GLN	5.8
1	A	760	LEU	5.2
1	A	699	LYS	5.1
1	A	818	GLU	4.9
1	A	697	THR	4.8
1	A	564	VAL	4.7
1	A	698	ASP	4.7
1	A	727	VAL	4.5
1	A	700	ASN	4.3
1	A	693	GLU	4.2
1	A	651	ARG	4.1
1	A	765	PRO	4.1
1	A	725	ALA	3.8
1	A	593	ALA	3.7
1	A	756	LYS	3.7
1	A	483	ASP	3.7
1	A	493	GLU	3.6
1	A	817	ASP	3.6
1	A	695	GLY	3.5
1	A	694	ASN	3.5
1	A	647	LYS	3.5
1	A	642	LYS	3.5
1	A	592	ALA	3.5
1	A	763	VAL	3.4
1	A	696	GLU	3.4
1	A	724	CYS	3.3

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Mol	Chain	Res	Type	RSRZ
1	A	757	GLN	3.3
1	A	721	LEU	3.1
1	A	484	VAL	3.1
1	A	722	ILE	2.9
1	A	595	ILE	2.9
1	A	816	LEU	2.8
1	A	761	PRO	2.8
1	A	726	ASP	2.6
1	A	596	HIS	2.6
1	A	482	ASP	2.6
1	A	640	ASP	2.6
1	A	784	TYR	2.6
1	A	589	ALA	2.4
1	A	497	TYR	2.4
1	A	703	VAL	2.3
1	A	528	CYS	2.3
1	A	565	LEU	2.3
1	A	562	ALA	2.3
1	A	641	ASP	2.3
1	A	735	GLN	2.3
1	A	764	MET	2.3
1	A	560	HIS	2.2
1	A	665	VAL	2.2
1	A	568	THR	2.2
1	A	706	THR	2.2
1	A	785	PHE	2.2
1	A	591	ILE	2.2
1	A	575	GLU	2.1
1	A	561	SER	2.1
1	A	650	GLU	2.1
1	A	728	SER	2.1
1	A	731	CYS	2.0
1	A	752	THR	2.0
1	A	701	GLN	2.0
1	A	539	ARG	2.0
1	A	766	VAL	2.0

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

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

6.3 Carbohydrates [i](#)

There are no carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th 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	LLDF	B-factors(\AA^2)	Q<0.9
3	MG	A	2	1/1	0.98	0.35	3.25	36,36,36,36	0
2	ZN	A	1	1/1	0.99	0.11	-2.55	34,34,34,34	0

6.5 Other polymers [i](#)

There are no such residues in this entry.