



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 4, 2026 – 06:02 PM UTC

PDB ID : 1I0D / pdb_00001i0d
Title : HIGH RESOLUTION STRUCTURE OF THE ZINC/CADMIUM-CONTAINING PHOSPHOTRIESTERASE FROM PSEUDOMONAS DIMINUTA
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Deposited on : 2001-01-29
Resolution : 1.30 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 types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

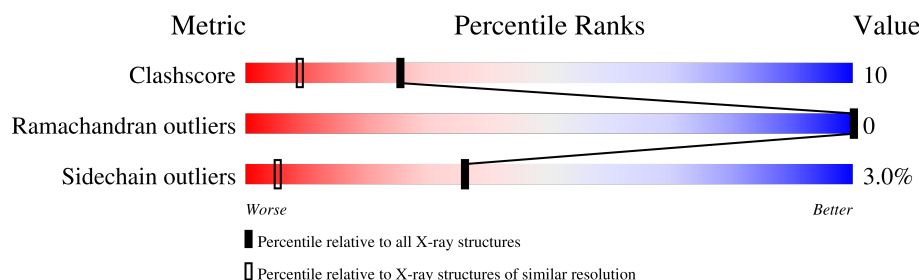
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.30 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	190562	1595 (1.30-1.30)
Ramachandran outliers	187476	1551 (1.30-1.30)
Sidechain outliers	187428	1551 (1.30-1.30)

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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	332	
1	B	332	

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
5	EDO	A	417	-	-	X	-
5	EDO	B	418	-	-	X	-

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 5923 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 PHOSPHOTRIESTERASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	331	Total	C	N	O	S	0	9	0
			2556	1617	453	479	7			
1	B	331	Total	C	N	O	S	0	7	0
			2538	1603	455	473	7			

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

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

- Molecule 3 is CADMIUM ION (CCD ID: CD) (formula: Cd).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Cd	0	0
			1	1		
3	B	1	Total	Cd	0	0
			1	1		

- Molecule 4 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Na	0	0
			1	1		
4	B	1	Total	Na	0	0
			1	1		

- Molecule 5 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



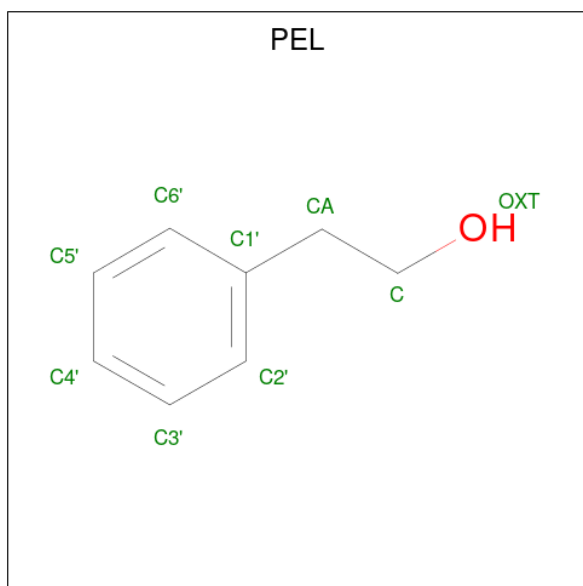
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	1
			5	2	3		
5	A	1	Total	C	O	0	0
			4	2	2		
5	A	1	Total	C	O	0	0
			4	2	2		
5	A	1	Total	C	O	0	0
			4	2	2		
5	A	1	Total	C	O	0	0
			4	2	2		
5	A	1	Total	C	O	0	0
			4	2	2		
5	A	1	Total	C	O	0	0
			4	2	2		
5	B	1	Total	C	O	0	1
			5	2	3		
5	B	1	Total	C	O	0	0
			4	2	2		
5	B	1	Total	C	O	0	0
			4	2	2		
5	B	1	Total	C	O	0	0
			4	2	2		
5	B	1	Total	C	O	0	0
			4	2	2		

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

- Molecule 6 is 2-PHENYL-ETHANOL (CCD ID: PEL) (formula: $C_8H_{10}O$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			9	8	1		
6	B	1	Total	C	O	0	0
			9	8	1		

- Molecule 7 is FORMIC ACID (CCD ID: FMT) (formula: CH_2O_2).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total	C	O	0	0
			3	1	2		
7	B	1	Total	C	O	0	0
			3	1	2		

- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	356	Total	O	0	0
			356	356		
8	B	381	Total	O	0	0
			381	381		

Note EDS was not executed.

Chain A: 80% 16% .



Chain B: 85% 14%



4 Data and refinement statistics

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

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	128.44Å 90.03Å 68.39Å 90.00° 91.72° 90.00°	Depositor
Resolution (Å)	30.00 – 1.30	Depositor
% Data completeness (in resolution range)	(Not available) (30.00-1.30)	Depositor
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	TNT	Depositor
R, R_{free}	0.191 , 0.224	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	5923	wwPDB-VP
Average B, all atoms (Å ²)	20.0	wwPDB-VP

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: CD, PEL, ZN, NA, FMT, 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	0.98	1/2652 (0.0%)	1.40	18/3605 (0.5%)
1	B	0.95	0/2626	1.33	15/3566 (0.4%)
All	All	0.97	1/5278 (0.0%)	1.36	33/7171 (0.5%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	236	ASP	C-N	-5.38	1.26	1.34

The worst 5 of 33 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	203	ALA	N-CA-C	-8.24	93.33	108.24
1	B	106	ILE	N-CA-C	-8.13	103.45	112.80
1	A	203	ALA	N-CA-C	-7.77	93.74	107.73
1	A	58	ILE	N-CA-C	-7.29	103.05	111.00
1	A	236	ASP	N-CA-CB	7.11	122.51	110.49

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	2556	0	2583	48	0
1	B	2538	0	2560	45	1
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
5	A	37	0	54	14	0
5	B	25	0	36	8	0
6	A	9	0	10	0	0
6	B	9	0	10	0	0
7	A	3	0	0	0	0
7	B	3	0	0	1	0
8	A	356	0	0	7	1
8	B	381	0	0	12	0
All	All	5923	0	5253	101	1

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 10.

The worst 5 of 101 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:B:337:ARG:HH22	1:B:343:GLN:HG3	1.26	0.99
1:A:96:ARG:CG	5:A:417:EDO:H21	2.04	0.87
1:B:96:ARG:HD2	5:B:418:EDO:H12	1.61	0.81
1:A:222[A]:SER:HB2	1:A:363:ARG:HH21	1.45	0.81
1:A:96:ARG:HG3	5:A:417:EDO:H21	1.64	0.80

All (1) 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:343:GLN:OE1	8:A:1157:HOH:O[4_545]	2.01	0.19

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	339/332 (102%)	325 (96%)	14 (4%)	0	100	100
1	B	336/332 (101%)	325 (97%)	11 (3%)	0	100	100
All	All	675/664 (102%)	650 (96%)	25 (4%)	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	276/266 (104%)	262 (95%)	14 (5%)	21	1
1	B	272/266 (102%)	267 (98%)	5 (2%)	51	16
All	All	548/532 (103%)	529 (96%)	19 (4%)	36	4

5 of 19 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	36	ARG
1	B	303	LEU
1	B	344	GLU
1	B	299	SER
1	A	303	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	38	ASN
1	A	206	GLN
1	A	212	GLN
1	B	38	ASN
1	B	212	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 27 ligands modelled in this entry, 6 are monoatomic - leaving 21 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	$\# Z > 2$
5	EDO	A	406[B]	-	3,3,3	0.32	0	2,2,2	0.61	0
5	EDO	A	410	-	3,3,3	0.35	0	2,2,2	0.69	0
5	EDO	B	418	-	3,3,3	0.41	0	2,2,2	0.27	0
5	EDO	A	409	-	3,3,3	0.35	0	2,2,2	0.37	0
5	EDO	B	412	-	3,3,3	0.39	0	2,2,2	0.50	0
5	EDO	A	406[A]	-	3,3,3	0.33	0	2,2,2	0.72	0
5	EDO	B	413	-	3,3,3	0.45	0	2,2,2	0.49	0
5	EDO	A	416	-	3,3,3	0.59	0	2,2,2	0.29	0
7	FMT	A	369	3,1,2	2,2,2	0.33	0	1,1,1	0.05	0
5	EDO	A	408	-	3,3,3	0.47	0	2,2,2	0.23	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	EDO	B	407	-	3,3,3	0.25	0	2,2,2	0.57	0
5	EDO	A	417	-	3,3,3	0.48	0	2,2,2	0.18	0
6	PEL	A	415	-	9,9,9	1.15	0	10,10,10	0.69	0
5	EDO	A	419	-	3,3,3	0.43	0	2,2,2	0.32	0
7	FMT	B	369	3,1,2	2,2,2	0.00	0	1,1,1	0.05	0
5	EDO	B	405[B]	-	3,3,3	0.23	0	2,2,2	0.69	0
6	PEL	B	414	-	9,9,9	1.54	1 (11%)	10,10,10	0.81	0
5	EDO	A	421	-	3,3,3	0.49	0	2,2,2	0.41	0
5	EDO	B	405[A]	-	3,3,3	0.34	0	2,2,2	0.75	0
5	EDO	A	411	-	3,3,3	0.43	0	2,2,2	0.30	0
5	EDO	B	420	-	3,3,3	0.31	0	2,2,2	0.56	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
5	EDO	A	406[B]	-	-	0/1/1/1	-
5	EDO	A	410	-	-	1/1/1/1	-
5	EDO	B	418	-	-	1/1/1/1	-
5	EDO	A	409	-	-	1/1/1/1	-
5	EDO	B	412	-	-	1/1/1/1	-
5	EDO	A	406[A]	-	-	0/1/1/1	-
5	EDO	B	413	-	-	1/1/1/1	-
5	EDO	A	416	-	-	1/1/1/1	-
5	EDO	A	408	-	-	0/1/1/1	-
5	EDO	B	407	-	-	0/1/1/1	-
5	EDO	A	417	-	-	1/1/1/1	-
6	PEL	A	415	-	-	1/3/3/3	0/1/1/1
5	EDO	A	419	-	-	1/1/1/1	-
5	EDO	B	405[B]	-	-	0/1/1/1	-
6	PEL	B	414	-	-	2/3/3/3	0/1/1/1
5	EDO	A	421	-	-	1/1/1/1	-
5	EDO	B	405[A]	-	-	0/1/1/1	-
5	EDO	A	411	-	-	0/1/1/1	-
5	EDO	B	420	-	-	1/1/1/1	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	B	414	PEL	C4'-C3'	2.19	1.43	1.38

There are no bond angle outliers.

There are no chirality outliers.

5 of 13 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	A	415	PEL	OXT-C-CA-C1'
6	B	414	PEL	OXT-C-CA-C1'
5	A	410	EDO	O1-C1-C2-O2
5	A	419	EDO	O1-C1-C2-O2
5	B	412	EDO	O1-C1-C2-O2

There are no ring outliers.

11 monomers are involved in 23 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	410	EDO	1	0
5	B	418	EDO	5	0
5	A	409	EDO	1	0
5	B	412	EDO	2	0
5	A	416	EDO	3	0
5	A	408	EDO	2	0
5	A	417	EDO	7	0
7	B	369	FMT	1	0
5	B	405[A]	EDO	1	0
5	A	411	EDO	1	0
5	B	420	EDO	1	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](#)

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates [i](#)

EDS was not executed - this section is therefore empty.

6.4 Ligands [i](#)

EDS was not executed - this section is therefore empty.

6.5 Other polymers [i](#)

EDS was not executed - this section is therefore empty.