W O R L D W I D E
wwPDB
P R O T E I N   D A T A   B A N K

wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 13, 2017 – 02:06 pm GMT

PDB ID	:	3L0V
Title	:	Crystal structure of catalytic domain of TACE with the first hydantoin inhibitor occupying the S1' pocket
Authors	:	Orth, P.
Deposited on	:	2009-12-10
Resolution	:	1.75 Å(reported)

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

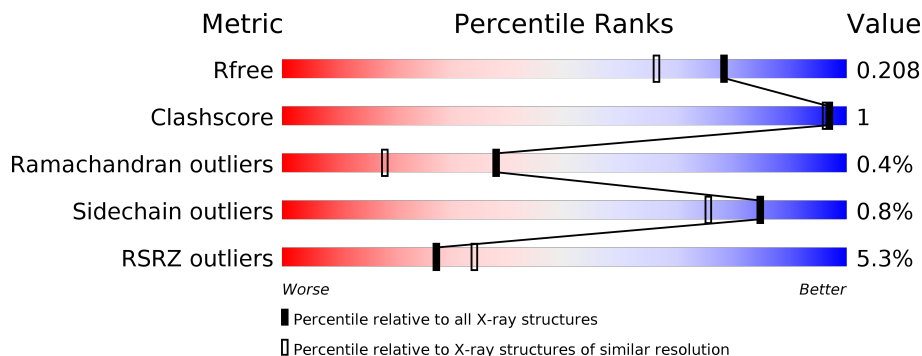|		
MolProbity	:	4.02b-467
Mogul	:	1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix)	:	1.9-1692
EDS	:	trunk28620
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	recalc28949

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 1.75 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   Percentile Ranks   Value

- Rfree — 0.208
- Clashscore — 1
- Ramachandran outliers — 0.4%
- Sidechain outliers — 0.8%
- RSRZ outliers — 5.3%

Worse   Better

▮ Percentile relative to all X-ray structures

▯ Percentile relative to X-ray structures of similar resolution

Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	100719	1762 (1.76-1.76)
Clashscore	112137	1889 (1.76-1.76)
Ramachandran outliers	110173	1868 (1.76-1.76)
Sidechain outliers	110143	1868 (1.76-1.76)
RSRZ outliers	101464	1770 (1.76-1.76)

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 <=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	270	3% / 92% / 6%
1	B	270	7% / 89% / 7%

2 Entry composition ⓘ

There are 4 unique types of molecules in this entry. The entry contains 4288 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 Disintegrin and metalloproteinase domain-containing protein 17.

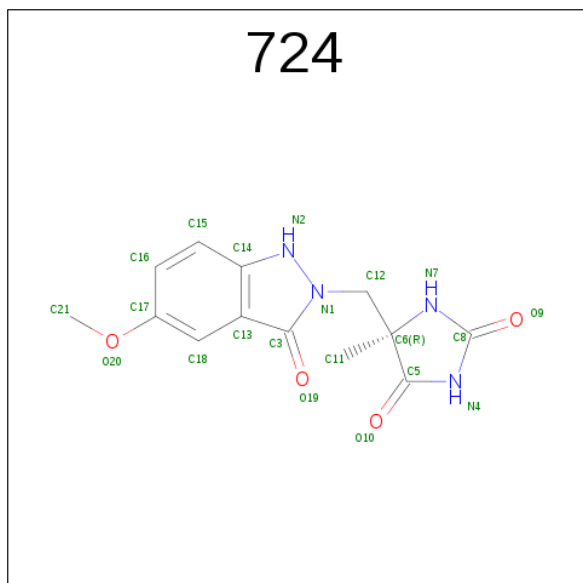| Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	255	Total	C	N	O	S	0	1	0
			1980	1249	329	389	13			
1	B	252	Total	C	N	O	S	0	0	0
			1923	1216	322	373	12			

There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	266	ALA	SER	ENGINEERED	UNP P78536
A	353	GLY	VAL	ENGINEERED	UNP P78536
A	452	GLN	ASN	ENGINEERED	UNP P78536
A	477	GLY	-	EXPRESSION TAG	UNP P78536
A	478	SER	-	EXPRESSION TAG	UNP P78536
A	479	HIS	-	EXPRESSION TAG	UNP P78536
A	480	HIS	-	EXPRESSION TAG	UNP P78536
A	481	HIS	-	EXPRESSION TAG	UNP P78536
A	482	HIS	-	EXPRESSION TAG	UNP P78536
A	483	HIS	-	EXPRESSION TAG	UNP P78536
A	484	HIS	-	EXPRESSION TAG	UNP P78536
B	266	ALA	SER	ENGINEERED	UNP P78536
B	353	GLY	VAL	ENGINEERED	UNP P78536
B	452	GLN	ASN	ENGINEERED	UNP P78536
B	477	GLY	-	EXPRESSION TAG	UNP P78536
B	478	SER	-	EXPRESSION TAG	UNP P78536
B	479	HIS	-	EXPRESSION TAG	UNP P78536
B	480	HIS	-	EXPRESSION TAG	UNP P78536
B	481	HIS	-	EXPRESSION TAG	UNP P78536
B	482	HIS	-	EXPRESSION TAG	UNP P78536
B	483	HIS	-	EXPRESSION TAG	UNP P78536
B	484	HIS	-	EXPRESSION TAG	UNP P78536

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

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

- Molecule 3 is (5R)-5-[(5-METHOXY-3-OXO-1,3-DIHYDRO-2H-INDAZOL-2-YL)METHYL]-5-METHYLIMIDAZOLIDINE-2,4-DIONE (three-letter code: 724) (formula: $C_{13}H_{14}N_4O_4$).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total 21	C 13	N 4	O 4	0	0
3	B	1	Total 21	C 13	N 4	O 4	0	0

- Molecule 4 is water.

| Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	183	Total 183	O 183	0	0
4	B	158	Total 158	O 158	0	0

3 Residue-property plots ⓘ

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 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: Disintegrin and metalloproteinase domain-containing protein 17
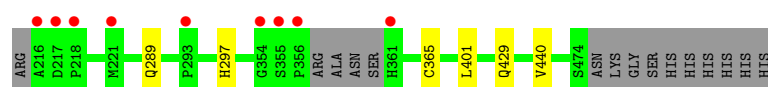
Chain A:

• Molecule 1: Disintegrin and metalloproteinase domain-containing protein 17

Chain B:
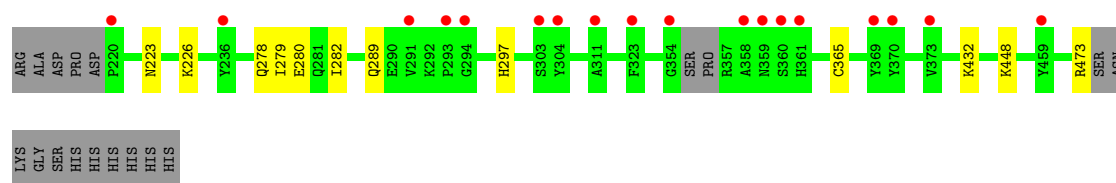
4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	72.04Å 76.84Å 103.80Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.89 – 1.75	Depositor
	46.89 – 1.75	EDS
% Data completeness (in resolution range)	(Not available) (46.89-1.75)	Depositor
	99.5 (46.89-1.75)	EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.46 (at 1.75Å)	Xtriage
Refinement program	BUSTER 2.9.3	Depositor
R, R_{free}	0.189 , 0.208	Depositor
	0.189 , 0.208	DCC
R_{free} test set	1142 reflections (1.99%)	DCC
Wilson B-factor (Å^2)	25.9	Xtriage
Anisotropy	0.254	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.35 , 45.8	EDS
L-test for twinning[2]	$< \lvert L \rvert > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	4288	wwPDB-VP
Average B, all atoms (Å^2)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.07% of the height of the origin peak. No significant pseudotranslation is detected.*

---

[1]Intensities estimated from amplitudes.
[2]Theoretical values of $< \lvert L \rvert >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 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, 724

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	0.50	0/2028	0.61	0/2744
1	B	0.50	0/1966	0.61	0/2663
All	All	0.50	0/3994	0.61	0/5407

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	1980	0	1855	2	0
1	B	1923	0	1776	6	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	21	0	13	0	0
3	B	21	0	13	0	0
4	A	183	0	0	0	0
4	B	158	0	0	2	0
All	All	4288	0	3657	8	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 1.

The worst 5 of 8 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:432:LYS:HE3	4:B:554:HOH:O	1.75	0.84
1:B:432:LYS:CE	4:B:554:HOH:O	2.48	0.49
1:A:401:LEU:HD13	1:A:440:VAL:HG23	1.99	0.44
1:B:278:GLN:HE21	1:B:473:ARG:HH11	1.66	0.42
1:A:289:GLN:HG2	1:A:297:HIS:CG	2.55	0.42

There are no symmetry-related clashes.

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	252/270 (93%)	247 (98%)	4 (2%)	1 (0%)	38	18
1	B	248/270 (92%)	243 (98%)	4 (2%)	1 (0%)	38	18
All	All	500/540 (93%)	490 (98%)	8 (2%)	2 (0%)	38	18

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	365	CYS
1	B	365	CYS

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	207/230 (90%)	206 (100%)	1 (0%)	91	86
1	B	193/230 (84%)	191 (99%)	2 (1%)	80	67
All	All	400/460 (87%)	397 (99%)	3 (1%)	85	75

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

Mol	Chain	Res	Type
1	A	429	GLN
1	B	223	ASN
1	B	448	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	223	ASN
1	B	278	GLN
1	B	359	ASN

5.3.3   RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5   Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6   Ligand geometry ⓘ

Of 4 ligands modelled in this entry, 2 are monoatomic - leaving 2 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	$\#|Z| > 2$
3	724	A	485	2	18,23,23	2.01	4 (22%)	23,35,35	2.00	7 (30%)
3	724	B	485	2	18,23,23	1.76	4 (22%)	23,35,35	2.17	8 (34%)

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
3	724	A	485	2	-	0/2/22/22	0/3/3/3
3	724	B	485	2	-	0/2/22/22	0/3/3/3

The worst 5 of 8 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	485	724	C18-C13	-3.10	1.35	1.41
3	B	485	724	C15-C14	-2.28	1.37	1.41
3	B	485	724	C8-N7	2.64	1.40	1.35
3	A	485	724	C8-N7	2.83	1.40	1.35
3	A	485	724	C16-C17	2.94	1.44	1.38

The worst 5 of 15 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	A	485	724	O20-C17-C18	-4.99	111.42	124.49
3	B	485	724	C17-C18-C13	-3.92	114.77	120.54
3	A	485	724	C16-C15-C14	-3.90	116.06	120.84
3	B	485	724	C5-N4-C8	-3.68	107.99	111.70
3	B	485	724	O20-C17-C18	-3.41	115.55	124.49

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 ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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	255/270 (94%)	0.34	9 (3%)	44	51	17, 27, 48, 69	0
1	B	252/270 (93%)	0.41	18 (7%)	17	22	18, 29, 53, 66	0
All	All	507/540 (93%)	0.38	27 (5%)	27	34	17, 28, 52, 69	0

The worst 5 of 27 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	216	ALA	5.7
1	A	356	PRO	5.7
1	B	361	HIS	5.3
1	B	358	ALA	4.4
1	B	369	TYR	4.3

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 ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

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(Å²)	Q<0.9
3	724	A	485	21/21	0.92	0.12	0.44	24,32,34,34	0
3	724	B	485	21/21	0.95	0.10	-0.28	18,20,22,22	0
2	ZN	B	2	1/1	1.00	0.09	-	22,22,22,22	0
2	ZN	A	1	1/1	0.99	0.10	-	27,27,27,27	0

6.5 Other polymers ⓘ

There are no such residues in this entry.