



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 02:24 pm GMT

PDB ID : 5EAS
Title : 5-EPI-ARISTOLOCHENE SYNTHASE FROM NICOTIANA TABACUM
Authors : Starks, C.M.; Back, K.; Chappell, J.; Noel, J.P.
Deposited on : 1997-06-19
Resolution : 2.25 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
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

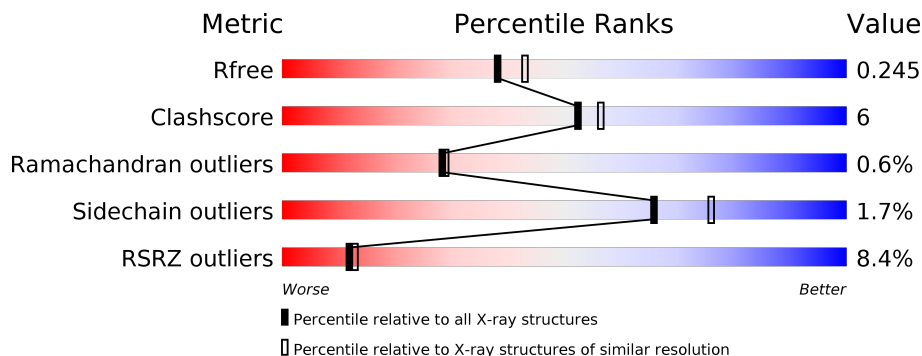
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 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))
R_{free}	100719	1062 (2.26-2.26)
Clashscore	112137	1178 (2.26-2.26)
Ramachandran outliers	110173	1145 (2.26-2.26)
Sidechain outliers	110143	1146 (2.26-2.26)
RSRZ outliers	101464	1066 (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. 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	548	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5369 atoms, of which 916 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 5-EPI-ARISTOLOCHENE SYNTHASE.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	514	Total	C	H	N	O	S	0	0	0
			5092	2671	916	690	797	18			

There are 5 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	44	GLN	LYS	CONFLICT	UNP Q40577
A	55	SER	ASN	CONFLICT	UNP Q40577
A	62	ARG	MET	CONFLICT	UNP Q40577
A	73	ILE	THR	CONFLICT	UNP Q40577
A	89	GLU	ASP	CONFLICT	UNP Q40577

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total	Mg	0	0
			2	2		

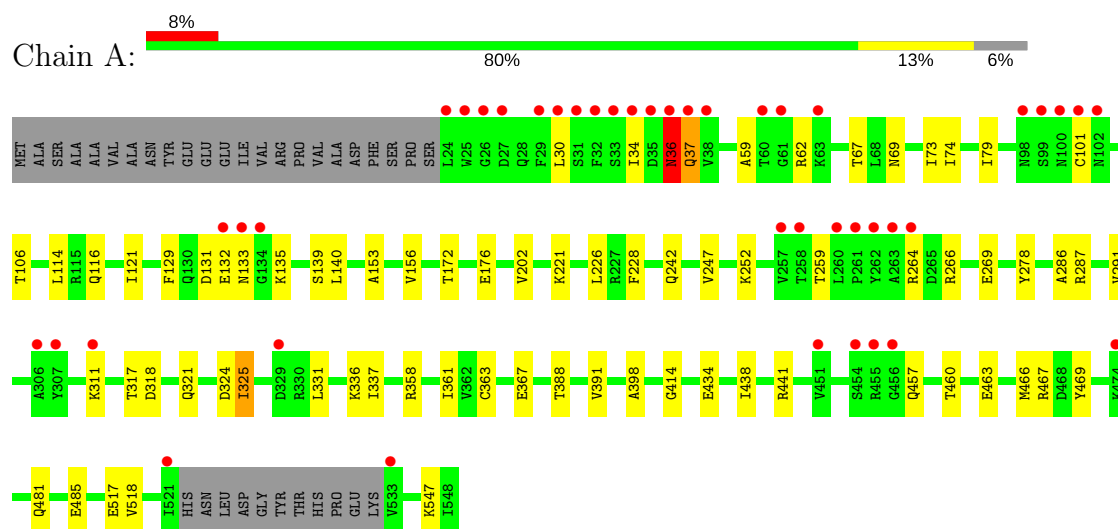
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	275	Total	O	0	0
			275	275		

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 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: 5-EPI-ARISTOLOCHENE SYNTHASE



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	124.73Å 124.73Å 118.94Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.25 23.96 – 2.25	Depositor EDS
% Data completeness (in resolution range)	94.0 (20.00-2.25) 95.3 (23.96-2.25)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.10 (at 2.26Å)	Xtriage
Refinement program	X-PLOR 3.1	Depositor
R, R_{free}	0.240 , 0.200 0.203 , 0.245	Depositor DCC
R_{free} test set	2085 reflections (4.86%)	DCC
Wilson B-factor (Å ²)	31.5	Xtriage
Anisotropy	0.264	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 50.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.012 for -h,-l,-k 0.006 for l,-k,h	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	5369	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.44% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

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: 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	0/4262	0.51	0/5775

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 [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	4176	916	4138	52	0
2	A	2	0	0	0	0
3	A	275	0	0	4	0
All	All	4453	916	4138	52	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.

All (52) 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:133:ASN:ND2	1:A:135:LYS:HD3	1.85	0.92

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:36:ASN:HD22	1:A:36:ASN:N	1.66	0.90
1:A:36:ASN:ND2	1:A:36:ASN:N	2.36	0.74
1:A:36:ASN:HD22	1:A:36:ASN:H	1.38	0.71
1:A:36:ASN:ND2	1:A:36:ASN:H	1.91	0.67
1:A:30:LEU:HG	1:A:252:LYS:HZ3	1.60	0.66
1:A:287:ARG:O	1:A:291:VAL:HG23	1.94	0.66
1:A:358:ARG:O	1:A:361:ILE:HG12	1.95	0.65
1:A:463:GLU:HA	1:A:466:MET:HE3	1.79	0.64
1:A:34:ILE:HD12	1:A:242:GLN:NE2	2.14	0.63
1:A:101:CYS:SG	1:A:106:THR:HB	2.38	0.63
1:A:133:ASN:HD21	1:A:135:LYS:HD3	1.64	0.63
1:A:259:THR:HB	1:A:337:ILE:HD13	1.82	0.62
1:A:59:ALA:HB3	1:A:62:ARG:HD3	1.82	0.60
1:A:317:THR:O	1:A:321:GLN:HG2	2.03	0.58
1:A:34:ILE:HD12	1:A:242:GLN:HE21	1.69	0.58
1:A:363:CYS:O	1:A:367:GLU:HG2	2.05	0.56
1:A:129:PHE:HB3	1:A:140:LEU:HD11	1.87	0.56
1:A:331:LEU:O	1:A:336:LYS:HE3	2.07	0.55
1:A:73:ILE:HG12	3:A:873:HOH:O	2.06	0.54
1:A:202:VAL:HG23	1:A:518:VAL:HG21	1.90	0.53
1:A:324:ASP:O	1:A:325:ILE:HB	2.09	0.53
1:A:131:ASP:C	1:A:133:ASN:H	2.13	0.52
1:A:133:ASN:HD22	1:A:135:LYS:HD3	1.70	0.52
1:A:69:ASN:O	1:A:73:ILE:HG13	2.09	0.51
1:A:30:LEU:CG	1:A:252:LYS:HZ3	2.24	0.51
1:A:153:ALA:O	1:A:156:VAL:HG22	2.12	0.50
1:A:457:GLN:O	1:A:460:THR:HG23	2.11	0.50
1:A:286:ALA:HA	1:A:414:GLY:HA3	1.94	0.50
1:A:37:GLN:HA	1:A:37:GLN:HE21	1.76	0.49
1:A:131:ASP:OD1	1:A:133:ASN:HB3	2.13	0.48
1:A:221:LYS:HD2	1:A:226:LEU:HD22	1.95	0.48
1:A:114:LEU:HD12	1:A:121:ILE:HD12	1.97	0.47
1:A:172:THR:O	1:A:176:GLU:HG3	2.15	0.45
1:A:30:LEU:HG	1:A:252:LYS:NZ	2.29	0.45
1:A:74:ILE:HG23	1:A:79:ILE:HG13	1.99	0.45
1:A:481:GLN:O	1:A:485:GLU:HG3	2.16	0.45
1:A:398:ALA:HB1	1:A:441:ARG:HD3	1.99	0.44
1:A:388:THR:HG21	1:A:467:ARG:HD2	1.99	0.44
1:A:30:LEU:CG	1:A:252:LYS:NZ	2.80	0.44
1:A:101:CYS:SG	1:A:106:THR:CB	3.06	0.43
1:A:391:VAL:HG21	1:A:469:TYR:HE1	1.83	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:266:ARG:HG2	1:A:269:GLU:HB2	2.00	0.43
1:A:30:LEU:HB2	1:A:252:LYS:HZ1	1.84	0.42
1:A:62:ARG:HH21	1:A:67:THR:HA	1.84	0.42
1:A:434:GLU:O	1:A:438:ILE:HG13	2.19	0.42
1:A:517:GLU:HB3	3:A:661:HOH:O	2.19	0.42
1:A:116:GLN:NE2	3:A:618:HOH:O	2.53	0.42
1:A:30:LEU:HA	1:A:252:LYS:HZ2	1.85	0.41
1:A:30:LEU:CB	1:A:252:LYS:NZ	2.83	0.41
1:A:547:LYS:HD2	3:A:827:HOH:O	2.20	0.41
1:A:247:VAL:HG11	1:A:291:VAL:HG11	2.02	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	510/548 (93%)	483 (95%)	24 (5%)	3 (1%)	28 29

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	325	ILE
1	A	36	ASN
1	A	132	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	461/489 (94%)	453 (98%)	8 (2%)	66 76

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	36	ASN
1	A	37	GLN
1	A	139	SER
1	A	228	PHE
1	A	264	ARG
1	A	278	TYR
1	A	311	LYS
1	A	318	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	28	GLN
1	A	36	ASN
1	A	37	GLN
1	A	116	GLN
1	A	133	ASN
1	A	149	ASN
1	A	155	HIS
1	A	242	GLN
1	A	540	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 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

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	514/548 (93%)	0.12	43 (8%) 12 13	17, 37, 100, 131	0

All (43) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	24	LEU	7.0
1	A	101	CYS	6.9
1	A	257	VAL	6.7
1	A	264	ARG	6.2
1	A	307	TYR	5.4
1	A	33	SER	5.4
1	A	29	PHE	5.4
1	A	30	LEU	5.2
1	A	100	ASN	5.0
1	A	34	ILE	4.7
1	A	521	ILE	4.7
1	A	133	ASN	4.3
1	A	35	ASP	4.3
1	A	27	ASP	4.1
1	A	37	GLN	4.0
1	A	31	SER	4.0
1	A	132	GLU	4.0
1	A	98	ASN	3.7
1	A	26	GLY	3.7
1	A	102	ASN	3.6
1	A	25	TRP	3.5
1	A	261	PRO	3.3
1	A	263	ALA	3.3
1	A	36	ASN	3.1
1	A	63	LYS	3.0
1	A	60	THR	3.0
1	A	329	ASP	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	258	THR	2.9
1	A	455	ARG	2.9
1	A	311	LYS	2.8
1	A	32	PHE	2.7
1	A	306	ALA	2.7
1	A	454	SER	2.7
1	A	260	LEU	2.7
1	A	451	VAL	2.7
1	A	533	VAL	2.7
1	A	456	GLY	2.6
1	A	38	VAL	2.5
1	A	262	TYR	2.5
1	A	474	LYS	2.3
1	A	99	SER	2.3
1	A	61	GLY	2.2
1	A	134	GLY	2.2

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 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
2	MG	A	851	1/1	0.92	0.11	-0.35	61,61,61,61	0
2	MG	A	852	1/1	0.75	0.22	-	62,62,62,62	0

6.5 Other polymers [i](#)

There are no such residues in this entry.