



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 29, 2017 – 07:30 AM EDT

PDB ID : 5UV0
Title : Crystal Structure of (+)-Limonene Synthase from Citrus sinensis
Authors : Prem Kumar, R.; Oprian, D.D.
Deposited on : 2017-02-17
Resolution : 2.30 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 : rb-20029077
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) : rb-20029077

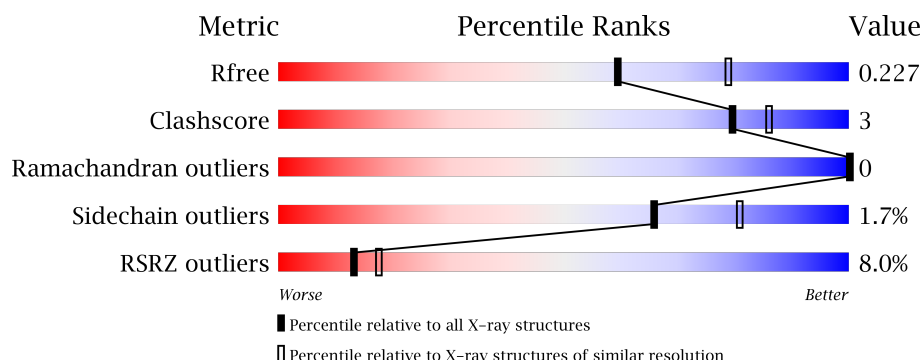
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION


The reported resolution of this entry is 2.30 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))
R_{free}	100719	4130 (2.30-2.30)
Clashscore	112137	4751 (2.30-2.30)
Ramachandran outliers	110173	4705 (2.30-2.30)
Sidechain outliers	110143	4704 (2.30-2.30)
RSRZ outliers	101464	4156 (2.30-2.30)

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 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	607	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4489 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 (+)-limonene synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	522	Total	C	N	O	S	0	0	0
			4315	2793	712	792	18			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	245	ALA	VAL	conflict	UNP A0A1C9J6A7

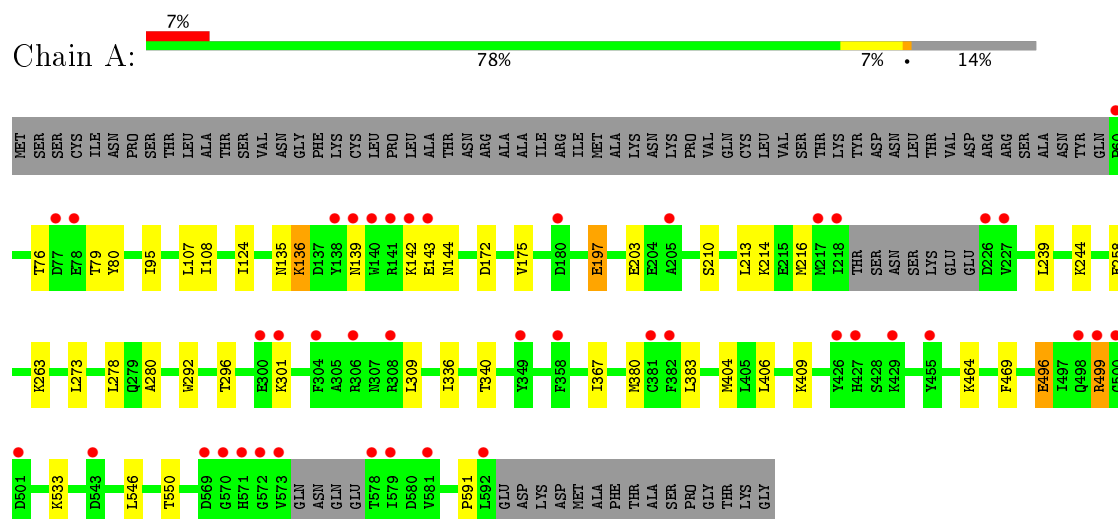
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	174	Total	O	0	0
			174	174		

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: (+)-limonene synthase



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	85.79 Å 85.79 Å 216.35 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.99 – 2.30 19.99 – 2.30	Depositor EDS
% Data completeness (in resolution range)	99.8 (19.99-2.30) 99.8 (19.99-2.30)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.90 (at 2.30 Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.190 , 0.229 0.185 , 0.227	Depositor DCC
R_{free} test set	1775 reflections (4.84%)	DCC
Wilson B-factor (Å ²)	49.7	Xtriage
Anisotropy	0.071	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 40.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	4489	wwPDB-VP
Average B, all atoms (Å ²)	56.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.86% 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](#)

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	0.39	0/4425	0.52	0/5987

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	4315	0	4243	27	0
2	A	174	0	0	2	0
All	All	4489	0	4243	27	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 3.

All (27) 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:591:PRO:O	2:A:701:HOH:O	2.11	0.69
1:A:301:LYS:HG3	1:A:380:MET:SD	2.37	0.64
1:A:108:ILE:HD13	1:A:124:ILE:HG23	1.87	0.56
1:A:406:LEU:HD23	1:A:409:LYS:HE3	1.88	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:258:GLU:O	1:A:263:LYS:NZ	2.27	0.53
1:A:76:THR:HG22	1:A:280:ALA:HB1	1.91	0.52
1:A:135:ASN:OD1	1:A:136:LYS:HE2	2.11	0.51
1:A:197:GLU:HB2	2:A:717:HOH:O	2.10	0.51
1:A:142:LYS:NZ	1:A:144:ASN:HD22	2.11	0.49
1:A:95:ILE:HG12	1:A:107:LEU:HD21	1.94	0.48
1:A:142:LYS:HE3	1:A:142:LYS:HB3	1.64	0.47
1:A:546:LEU:HD22	1:A:550:THR:HG21	1.98	0.46
1:A:203:GLU:OE1	1:A:533:LYS:NZ	2.44	0.45
1:A:336:ILE:O	1:A:340:THR:HG23	2.17	0.45
1:A:80:TYR:HB3	1:A:273:LEU:HD11	1.99	0.44
1:A:292:TRP:O	1:A:296:THR:HG23	2.19	0.43
1:A:309:LEU:HD12	1:A:309:LEU:HA	1.86	0.42
1:A:367:ILE:HD13	1:A:383:LEU:HD13	2.00	0.42
1:A:210:SER:O	1:A:214:LYS:HG2	2.19	0.42
1:A:499:ARG:HA	1:A:499:ARG:HD3	1.93	0.41
1:A:142:LYS:O	1:A:143:GLU:HG2	2.20	0.41
1:A:172:ASP:HB3	1:A:175:VAL:HG12	2.01	0.41
1:A:469:PHE:CD2	1:A:546:LEU:HG	2.56	0.41
1:A:213:LEU:HA	1:A:216:MET:HG3	2.03	0.41
1:A:239:LEU:HD23	1:A:244:LYS:HE2	2.02	0.41
1:A:278:LEU:HD12	1:A:278:LEU:HA	1.95	0.40
1:A:496:GLU:HA	1:A:499:ARG:HB2	2.04	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	516/607 (85%)	501 (97%)	15 (3%)	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	467/541 (86%)	459 (98%)	8 (2%)	66 81

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

Mol	Chain	Res	Type
1	A	79	THR
1	A	136	LYS
1	A	139	ASN
1	A	197	GLU
1	A	404	MET
1	A	464	LYS
1	A	496	GLU
1	A	499	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	144	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

There are no ligands in this entry.

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	522/607 (85%)	0.22	42 (8%) 13 17	35, 54, 84, 103	0

All (42) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	578	THR	8.3
1	A	569	ASP	7.4
1	A	218	ILE	6.6
1	A	579	ILE	5.3
1	A	60	PRO	4.7
1	A	499	ARG	4.4
1	A	570	GLY	4.1
1	A	217	MET	4.1
1	A	139	ASN	3.9
1	A	306	ARG	3.7
1	A	138	TYR	3.6
1	A	300	GLU	3.4
1	A	501	ASP	3.4
1	A	142	LYS	3.3
1	A	500	GLY	3.2
1	A	227	VAL	3.0
1	A	572	GLY	3.0
1	A	349	TYR	2.9
1	A	301	LYS	2.9
1	A	141	ARG	2.8
1	A	358	PHE	2.8
1	A	78	GLU	2.8
1	A	571	HIS	2.7
1	A	573	VAL	2.7
1	A	498	GLN	2.7
1	A	381	CYS	2.6
1	A	455	TYR	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	140	TRP	2.5
1	A	382	PHE	2.5
1	A	205	ALA	2.4
1	A	426	TYR	2.4
1	A	308	ARG	2.3
1	A	592	LEU	2.3
1	A	226	ASP	2.3
1	A	143	GLU	2.3
1	A	427	HIS	2.3
1	A	180	ASP	2.3
1	A	581	VAL	2.2
1	A	77	ASP	2.2
1	A	429	LYS	2.2
1	A	304	PHE	2.1
1	A	543	ASP	2.1

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 ⓘ

There are no ligands in this entry.

6.5 Other polymers ⓘ

There are no such residues in this entry.