

wwPDB X-ray Structure Validation Summary Report (i)

May 18, 2020 – 06:38 pm BST

PDB ID	:	1Q1L
Title	:	Crystal Structure of Chorismate Synthase
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Deposited on	:	2003-07-21
$\operatorname{Resolution}$:	2.05 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 (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: X-RAY DIFFRACTION

The reported resolution of this entry is 2.05 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	141614	1773 (2.04-2.04)
Ramachandran outliers	138981	1752 (2.04-2.04)
Sidechain outliers	138945	1752 (2.04-2.04)

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 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 <=5%

Note EDS was not executed.

Mol	Chain	Length	Quality of chain					
1	А	401	59%	23%	•	16%		
1	В	401	56%	25%	•	16%		
1	С	401	56%	25%	•	17%		
1	D	401	58%	23%	•	17%		



2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 11091 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.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace	
1	Λ	336	Total	С	Ν	Ο	\mathbf{S}	\mathbf{Se}	0	1	0
	л	000	2635	1674	465	489	1	6	0		0
1	В	336	Total	С	Ν	Ο	S	\mathbf{Se}	0	2	0
			2640	1675	465	493	1	6			0
1	C	222	Total	С	Ν	Ο	S	Se	0	9	0
		000	2622	1662	462	491	1	6		2	0
1	П	224	Total	С	Ν	Ο	S	Se	0	3	0
	034	2634	1670	465	492	1	6	0	0	0	

• Molecule 1 is a protein called Chorismate synthase.

There are 40 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
А	-2	GLY	-	EXPRESSION TAG	UNP 066493
А	-1	SER	-	EXPRESSION TAG	UNP 066493
А	0	HIS	-	EXPRESSION TAG	UNP 066493
А	1	MSE	MET	MODIFIED RESIDUE	UNP 066493
А	257	MSE	MET	MODIFIED RESIDUE	UNP 066493
А	258	MSE	MET	MODIFIED RESIDUE	UNP 066493
A	312	MSE	MET	MODIFIED RESIDUE	UNP 066493
A	320	MSE	MET	MODIFIED RESIDUE	UNP 066493
А	363	MSE	MET	MODIFIED RESIDUE	UNP 066493
А	381	MSE	MET	MODIFIED RESIDUE	UNP 066493
В	-2	GLY	_	EXPRESSION TAG	UNP 066493
В	-1	SER	_	EXPRESSION TAG	UNP 066493
В	0	HIS	-	EXPRESSION TAG	UNP 066493
В	1	MSE	MET	MODIFIED RESIDUE	UNP 066493
В	257	MSE	MET	MODIFIED RESIDUE	UNP 066493
В	258	MSE	MET	MODIFIED RESIDUE	UNP 066493
В	312	MSE	MET	MODIFIED RESIDUE	UNP 066493
В	320	MSE	MET	MODIFIED RESIDUE	UNP 066493
В	363	MSE	MET	MODIFIED RESIDUE	UNP 066493
В	381	MSE	MET	MODIFIED RESIDUE	UNP 066493
С	-2	GLY	-	EXPRESSION TAG	UNP 066493

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Chain	Residue	Modelled	Actual	Comment	Reference
С	-1	SER	-	EXPRESSION TAG	UNP 066493
С	0	HIS	-	EXPRESSION TAG	UNP 066493
С	1	MSE	MET	MODIFIED RESIDUE	UNP 066493
С	257	MSE	MET	MODIFIED RESIDUE	UNP 066493
С	258	MSE	MET	MODIFIED RESIDUE	UNP 066493
С	312	MSE	MET	MODIFIED RESIDUE	UNP 066493
С	320	MSE	MET	MODIFIED RESIDUE	UNP 066493
С	363	MSE	MET	MODIFIED RESIDUE	UNP 066493
С	381	MSE	MET	MODIFIED RESIDUE	UNP 066493
D	-2	GLY	-	EXPRESSION TAG	UNP 066493
D	-1	SER	-	EXPRESSION TAG	UNP 066493
D	0	HIS	-	EXPRESSION TAG	UNP 066493
D	1	MSE	MET	MODIFIED RESIDUE	UNP 066493
D	257	MSE	MET	MODIFIED RESIDUE	UNP 066493
D	258	MSE	MET	MODIFIED RESIDUE	UNP 066493
D	312	MSE	MET	MODIFIED RESIDUE	UNP 066493
D	320	MSE	MET	MODIFIED RESIDUE	UNP 066493
D	363	MSE	MET	MODIFIED RESIDUE	UNP 066493
D	381	MSE	MET	MODIFIED RESIDUE	UNP 066493

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• Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	А	157	Total O 157 157	0	0
2	В	137	Total O 137 137	0	0
2	С	127	Total O 127 127	0	0
2	D	139	Total O 139 139	0	0





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. Residues are colorcoded 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. 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.

Note EDS was not executed.





• Molecule 1: Chorismate synthase





F388 F398



4 Data and refinement statistics (i)

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

Property	Value	Source	
Space group	C 1 2 1	Depositor	
Cell constants	153.43Å 95.96 Å 113.46 Å	Depositor	
a, b, c, α , β , γ	90.00° 93.04° 90.00°	Depositor	
Resolution (Å)	24.63 - 2.05	Depositor	
% Data completeness	81 2 (24 63-2 05)	Depositor	
(in resolution range)	01.2 (24.05 2.09)	Depositor	
R_{merge}	0.07	Depositor	
R _{sym}	(Not available)	Depositor	
Refinement program	CNS 1.1	Depositor	
R, R_{free}	0.209 , 0.246	Depositor	
Estimated twinning fraction	No twinning to report.	Xtriage	
Total number of atoms	11091	wwPDB-VP	
Average B, all atoms $(Å^2)$	45.0	wwPDB-VP	



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).

Mal	Chain	Bo	ond lengths	Bond angles		
	Cham	RMSZ	# Z > 5	RMSZ	# Z > 5	
1	А	0.75	10/2675~(0.4%)	0.56	0/3579	
1	В	1.15	16/2680~(0.6%)	0.52	0/3588	
1	С	0.96	10/2661~(0.4%)	0.54	0/3560	
1	D	1.11	24/2673~(0.9%)	0.53	0/3577	
All	All	1.01	60/10689~(0.6%)	0.54	0/14304	

The worst 5 of 60 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms		Observed(Å)	Ideal(Å)
1	D	266[A]	VAL	CB-CG1	17.56	1.89	1.52
1	D	266[B]	VAL	CB-CG1	17.56	1.89	1.52
1	С	266[A]	VAL	CB-CG1	17.54	1.89	1.52
1	С	266[B]	VAL	CB-CG1	17.54	1.89	1.52
1	В	35[A]	GLU	CB-CG	17.36	1.85	1.52

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	А	2635	0	2653	104	0
1	В	2640	0	2648	115	0
1	С	2622	0	2624	111	0
1	D	2634	0	2639	127	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes	
2	А	157	0	0	1	0	
2	В	137	0	0	4	0	
2	С	127	0	0	6	0	
2	D	139	0	0	5	0	
All	All	11091	0	10564	407	0	

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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 19.

The worst 5 of 407 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:266[A]:VAL:CA	1:B:266[A]:VAL:CB	1.82	1.56	
1:B:266[B]:VAL:CA	1:B:266[B]:VAL:CB	1.82	1.56	
1:C:266[A]:VAL:CA	1:C:266[A]:VAL:CB	1.82	1.55	
1:C:266[B]:VAL:CA	1:C:266[B]:VAL:CB	1.82	1.55	
1:A:159[B]:LYS:CG	1:A:159[B]:LYS:CB	1.84	1.54	

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	Percer	ntiles
1	A	329/401~(82%)	313~(95%)	12 (4%)	4 (1%)	13	5
1	В	330/401~(82%)	312~(94%)	13~(4%)	5(2%)	10	3
1	С	327/401~(82%)	316~(97%)	10~(3%)	1 (0%)	41	31
1	D	329/401~(82%)	320~(97%)	8 (2%)	1 (0%)	41	31
All	All	1315/1604~(82%)	1261 (96%)	43 (3%)	11 (1%)	19	10

5 of 11 Ramachandran outliers are listed below:



Mol	Chain	Res	Type
1	А	324	PRO
1	В	84	ARG
1	В	324	PRO
1	А	292	LYS
1	В	135	SER

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	Perce	ntiles
1	А	274/322~(85%)	264~(96%)	10 (4%)	35	28
1	В	275/322~(85%)	265~(96%)	10 (4%)	35	28
1	С	273/322~(85%)	262~(96%)	11 (4%)	31	24
1	D	274/322~(85%)	263~(96%)	11 (4%)	31	24
All	All	1096/1288~(85%)	1054 (96%)	42 (4%)	34	26

5 of 42 residues with a non-rotameric sidechain are listed below:

Mol	Chain	\mathbf{Res}	Type
1	В	373	LEU
1	С	153	LEU
1	D	250[A]	ASP
1	В	376	LEU
1	С	6	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 30 such sidechains are listed below:

Mol	Chain	Res	Type
1	В	191	HIS
1	С	45	GLN
1	D	299	ASN
1	В	255	GLN
1	С	83	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)

There are no ligands in this entry.

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.

