

wwPDB X-ray Structure Validation Summary Report (i)

Sep 3, 2023 – 09:27 AM EDT

PDB ID	:	3RNB
Title	:	Structure of the Toluene/o-Xylene Monooxygenase Hydroxylase
		T201S/F176W Double Mutant
Authors	:	Gucinski, G.; Song, W.J.; Lippard, S.J.; Sazinsky, M.H.
Deposited on	:	2011-04-22
Resolution	:	2.64 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 types of validation reports are described at http://www.wwpdb.org/validation/2017/FAQs#types.

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

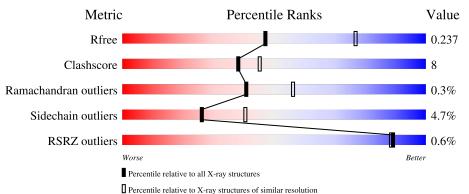
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.35
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: $X\text{-}RAY \, DIFFRACTION$

The reported resolution of this entry is 2.64 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	$egin{array}{c} { m Whole \ archive} \ (\#{ m Entries}) \end{array}$	${f Similar\ resolution}\ (\#{ m Entries,\ resolution\ range}({ m \AA}))$
R_{free}	130704	1426 (2.66-2.62)
Clashscore	141614	1472 (2.66-2.62)
Ramachandran outliers	138981	1446 (2.66-2.62)
Sidechain outliers	138945	1446 (2.66-2.62)
RSRZ outliers	127900	1408 (2.66-2.62)

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 <=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	А	498	80%	18%	••
2	В	330	% 80%	17%	••
3	С	86	62% 35%		•••



 $\mathbf{2}$

Entry composition (i)

There are 7 unique types of molecules in this entry. The entry contains 7526 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 Toluene o-xylene monooxygenase component.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
1	А	491	Total 4024	C 2568	N 676	0 754	S 26	0	1	0

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	176	TRP	PHE	engineered mutation	UNP Q6IV66
А	201	SER	THR	engineered mutation	UNP Q6IV66
А	445	LYS	GLU	engineered mutation	UNP Q6IV66

• Molecule 2 is a protein called Toluene o-xylene monooxygenase component.

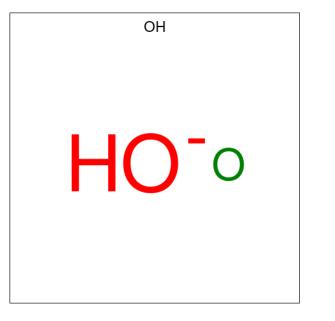
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
2	В	323	Total 2649	C 1680	N 468	0 491	S 10	0	0	0

• Molecule 3 is a protein called Toluene o-xylene monooxygenase component.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
3	С	85	Total 686	C 431	N 122	O 128	${ m S}{ m 5}$	0	0	0

• Molecule 4 is HYDROXIDE ION (three-letter code: OH) (formula: HO).



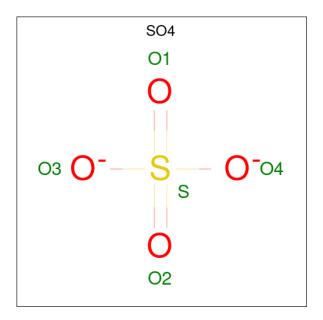


Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	А	1	Total O 1 1	0	0

• Molecule 5 is FE (III) ION (three-letter code: FE) (formula: Fe).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	А	2	TotalFe22	0	0

• Molecule 6 is SULFATE ION (three-letter code: SO4) (formula: O_4S).





Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	А	1	$\begin{array}{ccc} \text{Total} & \text{O} & \text{S} \\ 5 & 4 & 1 \end{array}$	0	0
6	А	1	$\begin{array}{ccc} \text{Total} & \text{O} & \text{S} \\ 5 & 4 & 1 \end{array}$	0	0
6	В	1	$\begin{array}{ccc} \text{Total} & \text{O} & \text{S} \\ 5 & 4 & 1 \end{array}$	0	0

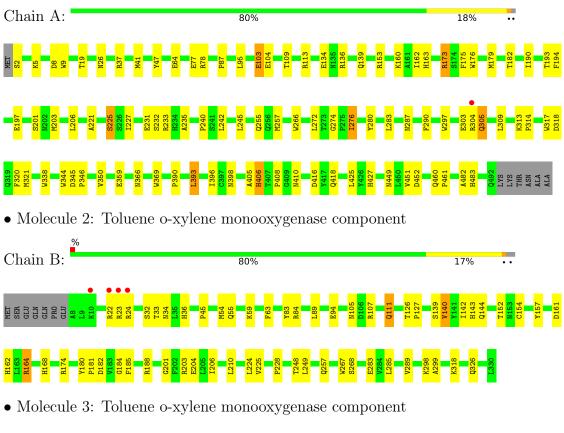
• Molecule 7 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	А	78	Total O 78 78	0	0
7	В	65	$\begin{array}{cc} \text{Total} & \text{O} \\ 65 & 65 \end{array}$	0	0
7	С	6	Total O 6 6	0	0

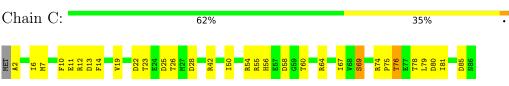


3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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: Toluene o-xylene monooxygenase component





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants	183.16Å 183.16Å 68.08Å	Depositor
a, b, c, α , β , γ	90.00° 90.00° 120.00°	Depositor
Resolution (Å)	44.99 - 2.64	Depositor
Resolution (A)	44.99 - 2.64	EDS
% Data completeness	99.4 (44.99-2.64)	Depositor
(in resolution range)	99.4(44.99-2.64)	EDS
R _{merge}	(Not available)	Depositor
R _{sym}	0.09	Depositor
$< I/\sigma(I) > 1$	8.17 (at 2.65 Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
D D	0.175 , 0.235	Depositor
R, R_{free}	0.182 , 0.237	DCC
R_{free} test set	1916 reflections (4.99%)	wwPDB-VP
Wilson B-factor $(Å^2)$	44.0	Xtriage
Anisotropy	0.161	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.33, 29.9	EDS
L-test for twinning ²	$< L > = 0.51, < L^2 > = 0.34$	Xtriage
Estimated twinning fraction	0.026 for -h,-k,l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7526	wwPDB-VP
Average B, all atoms $(Å^2)$	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 2.80% of the height of the origin peak. No significant pseudotranslation is detected.

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.



¹Intensities estimated from amplitudes.

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: OH, FE, SO4

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 Bor		Chain Bond lengths		Bond angles	
MOI	Unam	RMSZ	# Z > 5	RMSZ	# Z > 5	
1	А	0.92	0/4149	0.87	3/5640~(0.1%)	
2	В	0.97	4/2721~(0.1%)	0.88	2/3699~(0.1%)	
3	С	0.80	0/700	0.90	0/948	
All	All	0.93	4/7570~(0.1%)	0.88	5/10287~(0.0%)	

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Ζ	Observed(A)	Ideal(Å)
2	В	283	GLU	CG-CD	5.87	1.60	1.51
2	В	182	ASP	CB-CG	5.81	1.64	1.51
2	В	157	TYR	CE2-CZ	5.54	1.45	1.38
2	В	140	VAL	CB-CG1	5.30	1.64	1.52

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
1	А	393	LEU	CA-CB-CG	7.17	131.78	115.30
1	А	103	GLU	OE1-CD-OE2	-5.94	116.17	123.30
1	А	173	ARG	NE-CZ-NH2	-5.46	117.57	120.30
2	В	84	ARG	NE-CZ-NH1	5.18	122.89	120.30
2	В	164	ARG	NE-CZ-NH1	-5.09	117.75	120.30

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



Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	А	4024	0	3771	61	0
2	В	2649	0	2548	42	0
3	С	686	0	674	26	0
4	А	1	0	0	1	0
5	А	2	0	0	0	0
6	А	10	0	0	1	0
6	В	5	0	0	1	0
7	А	78	0	0	1	0
7	В	65	0	0	2	0
7	С	6	0	0	0	0
All	All	7526	0	6993	109	0

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.

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 8.

The worst 5 of 109 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:369:TRP:CH2	1:A:406:HIS:O	2.12	1.01
3:C:75:PRO:O	3:C:76:THR:HB	1.61	0.98
1:A:427:HIS:HE1	3:C:76:THR:HG23	1.29	0.97
1:A:427:HIS:HE1	3:C:76:THR:CG2	1.82	0.92
1:A:8:ASP:O	2:B:174:ARG:HD2	1.71	0.90

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	А	490/498~(98%)	460 (94%)	28~(6%)	2~(0%)	34	48



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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Perce	ntiles
2	В	321/330~(97%)	315~(98%)	6~(2%)	0	100	100
3	С	83/86~(96%)	75~(90%)	7 (8%)	1 (1%)	13	18
All	All	894/914 (98%)	850 (95%)	41 (5%)	3~(0%)	41	56

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All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	А	406	HIS
3	С	69	SER
1	А	408	PRO

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	А	417/422~(99%)	397~(95%)	20~(5%)	25 39
2	В	275/282~(98%)	264 (96%)	11 (4%)	31 47
3	С	77/79~(98%)	72 (94%)	5 (6%)	17 26
All	All	769/783~(98%)	733~(95%)	36~(5%)	26 40

 $5~{\rm of}~36$ residues with a non-rotameric side chain are listed below:

Mol	Chain	Res	Type
2	В	206	ILE
3	С	85	ASP
2	В	248	THR
3	С	50	ILE
1	А	280	TYR

Sometimes side chains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 19 such side chains are listed below:

Mol	Chain	Res	Type
2	В	153	ASN

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Mol	Chain	Res	Type
2	В	326	GLN
3	С	56	HIS
2	В	168	HIS
1	А	483	HIS

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 monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 6 ligands modelled in this entry, 1 is modelled with single atom and 2 are monoatomic - leaving 3 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		be Chain	Res	Link	Bond lengths			Bond angles		
	Mol Type Chain Res		Counts	RMSZ	# Z >2	Counts	RMSZ	# Z >2		
6	SO4	А	503	-	4,4,4	0.10	0	$6,\!6,\!6$	0.23	0
6	SO4	А	502	-	4,4,4	0.17	0	$6,\!6,\!6$	0.77	0
6	SO4	В	331	-	4,4,4	0.18	0	$6,\!6,\!6$	0.25	0

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.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	А	502	SO4	1	0
6	В	331	SO4	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)

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, 95^{th} 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	$OWAB(Å^2)$	$\mathbf{Q}{<}0.9$
1	А	491/498~(98%)	-0.29	1 (0%) 95 96	29, 45, 60, 77	0
2	В	323/330~(97%)	-0.40	4 (1%) 79 77	31, 41, 60, 86	0
3	С	85/86~(98%)	-0.43	0 100 100	40, 53, 65, 74	0
All	All	899/914 (98%)	-0.34	5 (0%) 89 88	29, 45, 61, 86	0

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	В	22	ARG	3.4
2	В	23	ARG	3.2
1	А	304	ARG	2.5
2	В	10	LYS	2.4
2	В	24	ARG	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 monosaccharides 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. The B-factors column lists the minimum, median, 95^{th} 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	B-factors(Å ²)	Q<0.9
6	SO4	А	503	5/5	0.91	0.16	49,51,52,52	5
6	SO4	А	502	5/5	0.92	0.19	42,42,46,46	5
6	SO4	В	331	5/5	0.96	0.11	47,48,49,49	5
5	FE	А	501	1/1	0.99	0.17	43,43,43,43	0
4	OH	А	499	1/1	1.00	0.16	49,49,49,49	0
5	FE	А	500	1/1	1.00	0.17	38,38,38,38	0

6.5 Other polymers (i)

There are no such residues in this entry.

