Research Article

Scale Effects of Physical Modelling of Anaerobic Treatment Plants in the Treatment of Cassava Wastewater

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Abstract: Three Anaerobic Baffled Reactors (ABRs) are used to evaluate the extent of scale effects of cyanide inhibition of cassava wastewater treatment. The reactors (physical models) have aspect ratios of (53:16:30; 4:1:1; 10:3:6). Kinetic analyses of specific growth rate μ_{max} and half saturation constant k_s are evaluated for the respective reactors. For the respective reactors, non-inhibited cassava wastewater treatment process showed Kincannon model yields ($\mu_{max} = 17.24$ day⁻¹; 21.74day⁻¹; 21.28 day⁻¹

and $k_s = 1.00mgCODL^{-1}$; 2.37 $mgCODL^{-1}$; 1.23 $mgCODL^{-1}$) while Monod model yields ($\mu_{max} = 10.87 \text{ day}^{-1}$; 12.82 day⁻¹; 13.70 day⁻¹ and $k_s = 0.87 mgCODL^{-1}$; 1.92 $mgCODL^{-1}$; 2.32 $mgCODL^{-1}$). Coefficient of determination R² is used to verify the respective models to yield values of ($R^2 = 0.839$; 0.960; 0.948) for Kincannon model and ($R^2 = 0.917$; 0.842; 0.830) for Monod model.

Keywords: Anaerobic, Treatment, Cassava Wastewater, Monod model and Kincannon model

1. Introduction

Cassava (Manihot esculenta crantz, also known as manioc or yucca) is one of the leading food and feed plants in the world: it ranks fourth among staple crops with a global production of about 160 milllion tons per year (1). Most of these are grown in three regions, West Africa, and the Congo basin, tropical South America, and South East Asia (2), while in Western countries it is not commonly used, because of the presence of cyanoglucosides (linamarin and lotaustralin).

Nigeria currently is the largest producer of cassava in the whole world with an annual output of over 34 million tons of tuberous root and it plays a dominant role in the rural economy in the southern agroecological zone (3). As a food crop, cassava fits well into the farming system of the small holder farmers in Nigeria because it is available year round, thus providing household food security. Cassava tubers can be kept in the ground prior to harvesting up to two years but once harvested, they begin to deteriorate. To forestall early deterioration and also due to its bulky nature, cassava is usually traded in their processed form. The bulky roots contain much moisture (60-65%), making their transportation from rural areas difficult and expensive. Processing the tuber into a dry form reduces the moisture content and converts into a more durable and stable product with less volume which makes it more transportable (4; 5). Over the years cassava has been transformed into a number of product both for domestic use (depending on local customs and preference) and industrial uses. Cassava in the fresh form contains cyanide which is extremely toxic to humans and animals and therefore needs to be processed to reduce the cyanide content to safe level (6). Traditional cassava processing method involve several steps including peeling, soaking, grinding, steeping in water and left in air to allow fermentation to occur, drying, milling, roasting, steaming, pounding and mixing in cold or hot water. The wastewater from cassava processing or its derivative (garri) ends up with domestic sewage if processed in small quantities while others end up being carried with industrial wastes if processed in large industrial quantities. Lastly, others percolate into the soil depending on the processors. Wastewater from cassava processing, if released directly into the environment before proper treatment, is a source of pollution. This effect on the

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B.C.Okoro (Correspondence) bc1okoro@yahoo.com environment is yet to be addressed properly in developing countries due to inadequate equipment and lack of research materials.

A great number of recent studies have reported many biotechnological approaches to improve the safety and quality of cassava flour (7; 8; 9), and the effect of different processing modalities of the tuberous roots

2. Materials and Method 2.1 The Reactor

on the level of these toxic substances and functional properties has been assessed (10; 11; 12). The objective of this study is to study the scale effect of a physical model of anaerobic treatment plants in the treatment of cassava wastewater as well as to develop a mathematical model describing the microbial kinetics.



Fig. 1: Scheme of the ABR. 1. Feed Tank; 2. Peristaltic Pump; 3. Influent; 4. Sampling Ports; 5. Effluent.

The laboratory scale ABR was constructed from 6mm thick stainless steel, with external dimensions of lengths, widths, depths and working volumes of three different sizes as shown in table 1. Fig. 1 shows a schematic diagram of the reactor. The reactors were divided into different number of equal compartments by vertical baffles with each compartment of each the reactors having downcomer and riser regions created by a further vertical baffle. For each of the reactors, the widths of upcomers were multiples of the widths of downcomer. The lower parts of the downcomer baffles were angled at 45^0 in order to direct the flow

evenly through the upcomer. This produced effective mixing and contact between the wastewater and anaerobic sludge at the base of each riser. Each compartment was equipped with sampling ports that allowed biological solids and liquid samples to be withdrawn. The operating temperature was maintained constant at $35\pm0.5^{\circ}$ C by putting the reactor in a water bath equipped with a temperature regulator. The influent feed was pumped using variable speed peristaltic pump. The outlet was connected to a glass U-tube for level control and to trap solids.

Tuble 1. Dimensions of the Different Reactor	Mouti Dizes.		
Dimensions	Reactor 1	Reactor 2	Reactor 3
Length (cm)	53	40	50
Width (cm)	16	10	15
Depth (cm)	30	10	30
Working Volume (L)	13.57	4	10
Upcomer Width/ Downcomer Width	2.6	4	2.4
No of Compartments	5	4	5
Volume of Each Compartment (L)	2.7	4	2

Table 1: Dimensions of the Different Reactor Model Sizes.

2.2 Start-up of ABR

Start-up without seed sludge was rather difficult and time consuming for suspended growth anaerobic reactors. The following 3 steps were taken: (i) the reactor was filled with cassava wastewater and allowed to rest for 15 days (ii) the sludge bed was allowed through a process of sludge accumulation by settling and sludge improvement and (iii) after 15 days, feeding of the wastewater was resumed at a flow rate of 5.33litres per day and HRT of 6 days with a very low organic loading rate (OLR) of 0.067kgCOD/m³.day. The resumed wastewater feeding is helped the development of sludge bed at the bottom of individual chambers of the ABR. This process of feeding the system followed by two weeks rest is based on the experiment made in Kanpur (India) for the start-up of a UASB plant without inoculum (13).

2.3 Characterization of Wastewater

The cassava wastewater from a cassava processing factory at Imo Polytechnic Umuagwo in eastern Nigeria was used as feed. The supernatant of the wastewater after the simple gravity settling, used in the investigation, had low TSS, as approximately 90% of the solids were removed. The supernatant wastewater was diluted to achieve the COD concentration required for each loading rate with water. In order to achieve pH and alkalinity adjustment, the supernatant was neutralized by NaOH and NaHCO₃. A COD: N: P ratio of 300:5: 1 was kept during operation using NH₄C1 and K₂HPO₄. The micro-nutrient deficiency was added occasionally to correct growth conditions according to (14).

2.4 Procedure for Experiment

The wastewater was collected twice a day from the cassava processing plant, and it was intermittently mixed to feed the reactor with a consistent quality. The wastewater came from processing cassava specie

(bogot) that had no cyanide content (Table 2). The wastewater was fed to the reactor with the help of a variable speed peristaltic pump. The ABR was operated at various hydraulic retention times (HRTs) by varying the flow rate of influent wastewater (Q_{inf}) , thereby varying the organic loading rate (OLR). The wastewater flowed from the downcomer to the upcomer within an individual chamber through the sludge bed formed at the bottom of the individual chambers. After receiving treatment in the particular chamber, wastewater entered the next chamber from the top. Due to the specific design and .positioning of the baffle, the wastewater is evenly distributed in the upcomer and the vertical upflow velocity (V_{up}) could be significantly reduced. The treated effluent was collected from the outlet of the 3rd compartment (C3). The reactor was kept in a temperature controlled chamber maintained at 35 °C.

2.5 Mathematical Model Formulation 2.5.1 Nomenclature

 S_i = Substrate concentration in the influent (mgl⁻¹); S_e = Substrate concentration in the effluent (mgl⁻¹); k_s = Half saturation constant (mgl⁻¹); μ = Specific growth rate of organism (per day); μ_{max} = Maximum specific growth rate of organism (per day); X = concentration of active biomass (mg/L); r_A = Rate of utilization of substrate (mg/l.day)

2.5.2 Monod Model for ABR

The Monod model is described as:

$$r_A = \frac{ds}{dt} = \frac{Q}{V} \left(S_i - S_g \right) = \mu X \tag{1}$$

$$= \frac{\mu_{max}S_e}{K_s + S_e}X \qquad = \qquad \frac{XV}{Q(S_i - S_e)} = \qquad \frac{K_s}{\mu_{max}}\frac{1}{S_e} + \frac{1}{\mu_{max}} \tag{2}$$

Applying experimental results to Equations (2), graph will be plotted. In this, graph XV/ $(S_i\text{-}S_e)$ is plotted against $1/S_e$

Variety	Total Leaves	Free Leaves	Total Roots	Free Roots	Total Cyanide
	Cyanide (µg/g)	Cyanide (µg/g)	Cyanide (µg/g)	Cyanide (µg/g)	Content Ratio in
					Leaves and Roots
Java Brown	490	33(6.7)	185	9(4.9)	2.6
Datu	541	19(3.5)	120	9(7.5)	4.5
Bogot	456	21(4.6)	n.d.	n.d.	n.d.
Lakan	189	13(6.9)	45	0.4(1.0)	4.2

Table 2: Cassava Species of varying Cyanide Concentrations

n.d. = not detected; Source: 15.

Results and Discussion Mathematical Model Calibration Table 3: Computations for Reactor 1

Table	5. Compu	tations for Kea				
S/N	x_i $= SV_i$ $- SV_e$ (mg/ L.d)	$\frac{Q\left(S_i-S_e\right)}{\substack{V\\(\text{mg}/\\\text{L.d})}}$	$\frac{V}{Q (S_i - S_e)}$ (mg/ L.d)	$y = x.V$ $\overline{Q(S_i - S_e)}$ (mg/ L.d)	Se (mg/L	$x = \frac{1}{S_e}$ (L/mg)
1	0.9	7.6	0.1315789	0.1184211	1.3	0.7692308
2	0.9	7.8	0.1282051	0.1153846	1.5	0.6666667
3	1.5	9	0.1111111	0.1666667	1.7	0.5882353
4	1.5	9.5	0.1052632	0.1578947	1.8	0.5555556
5	1.6	10.2	0.0980392	0.1568627	2	0.5
6	1.6	12.8	0.078125	0.125	2.2	0.4545455
7	1.7	13	0.0769231	0.1307692	2.5	0.4
8	1.8	13.5	0.0740741	0.1333333	2.8	0.3571429
9	1.8	14	0.0714286	0.1285714	3	0.3333333
10	1.9	15	0.0666667	0.1266667	3.2	0.3125
11	1.9	16	0.0625	0.11875	3.4	0.2941176
12	2	18.4	0.0543478	0.1086957	3.6	0.2777778
13	2	18.6	0.0537634	0.1075269	3.8	0.2631579
14	2	18.8	0.0531915	0.106383	4.8	0.2083333
15	2.1	20	0.05	0.105	5.6	0.1785714
16	2.1	20.4	0.0490196	0.1029412	5.8	0.1724138
17	2.1	20.6	0.0485437	0.1019417	6.4	0.15625
18	2.2	20.8	0.0480769	0.1057692	6.8	0.1470588
19	2.2	20.8	0.0480769	0.1057692	7.5	0.1333333
20	2.3	22.6	0.0442478	0.1017699	8.4	0.1190476
21	2.3	24.2	0.0413223	0.0950413	8.8	0.1136364
22	2.3	24.8	0.0403226	0.0927419	9.2	0.1086957
23	2.4	26	0.0384615	0.0923077	10.8	0.0925926
24	2.4	26.2	0.0381679	0.0916031	12.6	0.0793651



Fig. 2: A Plot of
$$y = \frac{XV}{Q(s_i - s_e)}$$
 (L.d/mg) Versus $x = \frac{1}{s_e}(L/mg)$ for Reactor 1

From linear regression;

 $y = mx + c \equiv y = 0.078x + 0.092$ (3)

$$\frac{1}{\mu_{max}} = c = intercept; \ i.e. \ \mu_{max} = \frac{1}{c} = \frac{1}{0.092} = 10.87/day \quad (4)$$

and;
$$\frac{K_s}{\mu_{max}} = m = slope; \ i.e. K_s = \mu_{max} * m$$
 (5)

$$\therefore K_s = 10.87 * 0.078 = 0.87 mgCOD/L$$
(6)

Substituting μ_{max} and k_s into equation 1 gives;

$$r_A = \frac{ds}{dt} = \frac{Q}{V}(S_i - S_e) = \frac{\mu_{max} * S_e}{K_s + S_e} = \frac{\mathbf{10.87} * S_e}{\mathbf{0.87} + S_e}$$
(7)

S/N	x_i $= SV_i$ $- SV_e$ (mg/ L.d)	$\frac{Q (S_i - S_e)}{V \atop (mg/ \\ L.d)}$	$\frac{V}{\substack{Q \ (S_i - S_e) \\ (mg/\\ L.d)}}$	$y = x.V$ $\overline{Q(S_i - S_e)}$ (mg/ L.d)	S _e (mg/L	$x = \frac{1}{S_e}$ (L/ mg)
1	1	6.4	0.15625	0.15625	1.5	0.6666667
2	1	6.5	0.1538462	0.1538462	1.5	0.6666667
3	1.5	6.8	0.1470588	0.2205882	1.6	0.625
4	1.5	7	0.1428571	0.2142857	1.6	0.625
5	1.6	9.8	0.1020408	0.1632653	1.8	0.5555556
6	1.6	12	0.0833333	0.1333333	2.2	0.4545455
7	1.8	14.6	0.0684932	0.1232877	2.5	0.4
8	1.8	14.8	0.0675676	0.1216216	2.6	0.3846154
9	1.8	15	0.0666667	0.12	2.8	0.3571429
10	2	15.6	0.0641026	0.1282051	3	0.3333333
11	2	17	0.0588235	0.1176471	3.2	0.3125
12	2.2	19	0.0526316	0.1157895	3.3	0.3030303
13	2.2	20.5	0.0487805	0.1073171	3.4	0.2941176
14	2.3	18.8	0.0531915	0.1223404	3.6	0.2777778
15	2.3	20	0.05	0.115	4.2	0.2380952
16	2.4	20.4	0.0490196	0.1176471	4.8	0.2083333
17	2.4	20.6	0.0485437	0.1165049	5.8	0.1724138
18	2.4	20.8	0.0480769	0.1153846	6.6	0.1515152
19	2.4	20.8	0.0480769	0.1153846	7.8	0.1282051
20	2.3	22.6	0.0442478	0.1017699	8	0.125
21	2.3	25	0.04	0.092	8.4	0.1190476
22	2.3	26.4	0.0378788	0.0871212	10.6	0.0943396
23	2.4	27.5	0.0363636	0.0872727	11.6	0.0862069
24	2.6	28	0.0357143	0.0928571	16.6	0.060241

 Table 4: Computations for Reactor 2



Fig. 3: A Plot of $y = \frac{XV}{Q(S_i - S_e)}$ (/L.d/mg) Versus $x = \frac{1}{s_e}$ (L/mg) for Reactor 2 From linear regression equation;

 $y = mx + c \equiv y = 0.150 x + 0.078$ (8)

By comparison with equation 2;

$$\frac{1}{\mu_{max}} = c = intercept; \ i. e. \ \mu_{max} = \frac{1}{c} = \frac{1}{0.078} = 12.82/day \tag{9}$$

and;
$$\frac{K_s}{M_s} = m = slope; \ i. e. \ K_s = \mu_{max} * m \tag{10}$$

$$\frac{\kappa_s}{\mu_{max}} = m = slope; \ i.e. K_s = \mu_{max} * m \tag{10}$$

$$\therefore K_s = 12.82 * 0.150 = 1.92mgCOD/L$$
(11)

Substituting μ_{max} and k_s into equation 1 gives;

$$r_A = \frac{ds}{dt} = \frac{Q}{V}(S_i - S_e) = \frac{\mu_{max} * S_e}{K_s + S_e} = \frac{\mathbf{12.82} * S_e}{\mathbf{1.92} + S_e}$$
(12)

Table 5	. Computa	tions for Reactor	3			
S/N	x_i $= SV_i$ $- SV_e$ (mg/ L.d)	$\frac{Q (S_i - S_e)}{V \atop (mg/ \\ L.d)}$	$\frac{V}{\substack{Q \ (S_i - S_e) \\ (mg/\\ L.d)}}$	$y = x_i \cdot V$ $\overline{Q(S_i - S_e)}$ (mg/ L.d)	S _e (mg/L	$x = \frac{1}{S_e}$ (L/mg)
1	12	5.8	0 1724138	0 2068966	12	0 8333333
2	1.2	6	0.1666667	0.2166667	1.2	0.7142857
3	1.3	6	0.1666667	0.2166667	1.4	0.7142857
4	1.3	6.8	0.1470588	0.1911765	1.4	0.7142857
5	1.4	7.4	0.1351351	0.1891892	1.8	0.5555556
6	1.4	8.6	0.1162791	0.1627907	2	0.5
7	1.5	14	0.0714286	0.1071429	2	0.5
8	1.5	14.2	0.0704225	0.1056338	2.6	0.3846154
9	1.5	14.6	0.0684932	0.1027397	2.8	0.3571429
10	1.8	15.6	0.0641026	0.1153846	4	0.25
11	1.8	16	0.0625	0.1125	4	0.25
12	2	16	0.0625	0.125	4	0.25
13	2.1	18	0.0555556	0.1166667	4	0.25
14	2.2	18.6	0.0537634	0.1182796	4.6	0.2173913
15	2.2	20	0.05	0.11	4.8	0.2083333
16	2.2	20.2	0.049505	0.1089109	4.8	0.2083333
17	2.3	20.4	0.0490196	0.1127451	6	0.1666667
18	2.3	20.6	0.0485437	0.1116505	6.6	0.1515152
19	2.4	20.8	0.0480769	0.1153846	6.8	0.1470588
20	2.3	22	0.0454545	0.1045455	8.6	0.1162791
21	2.3	24	0.0416667	0.0958333	8.8	0.1136364
22	2.3	28.6	0.034965	0.0804196	10.2	0.0980392
23	2.4	30.6	0.0326797	0.0784314	12.4	0.0806452
24	2.8	34.6	0.0289017	0.0809249	14	0.0714286

Table 5: Computations for Reactor 3

From linear regression equation;

$$y = mx + c \equiv y = 0.169x + 0.073$$
 (13)

By comparison with equation 1;

$$\frac{1}{\mu_{max}} = c = intercept; \ i. e. \ \mu_{max} = \frac{1}{c} = \frac{1}{0.073} = \mathbf{13.70} day^{-1}$$
(14)

and;
$$\frac{K_s}{\mu_{max}} = m; \quad K_s = \mu_{max} * m$$
 (15)

$$\therefore K_s = 13.70 * 0.169 = 2.32mg/L$$
(16)



Fig. 4: A Plot of $y = \frac{XV}{Q(S_i - S_e)}$ (/L.d/mg) Versus $x = \frac{1}{S_e}$ (L/mg) for Reactor 3

Substituting μ_{\max} and k_s into equation 2 gives; $r_A = \frac{ds}{dt} = \frac{Q}{V} (S_i - S_e) = \frac{\mu_{max} * S_e}{K_s + S_e} = \frac{\mathbf{13.70} * S_e}{\mathbf{2.32} + S_e}$ (17)

Observed y	Simulated y'
x_i . V	x_i . V
$\overline{\boldsymbol{O}\left(\boldsymbol{S}_{i}-\boldsymbol{S}_{a}\right)}$	$\overline{\boldsymbol{0}\left(\boldsymbol{S}_{i}-\boldsymbol{S}_{e}\right)}$
(mg/L.d)	(mg/L.d)
(8,,	(8,)
0.1184211	0.11456
0.1153846	0.114208
0.1666667	0.137882
0.1578947	0.135333
0.1568627	0.131
0.125	0.127455
0.1307692	0.1232
0.1333333	0.119857
0.1285714	0.125214
0.1266667	0.116375
0.11875	0.114941
0.1086957	0.113667
0.1075269	0.112526
0.106383	0.10825
0.105	0.105929
0.1029412	0.105448
0.1019417	0.104188
0.1057692	0.103471
0.1057692	0.1024
0.1017699	0.101286
0.0950413	0.100864
0.0927419	0.100478
0.0923077	0.099222
0.0916031	0.09819

3.2 Model Verification Table 6: Simulation for Reactor 1

Fig. 5: A Scatter plot of Observed y versus Simulated y' for Reactor 1



The coefficient of determination, R^2 for reactor 1 in Monod Model yielded 0.916, suggesting a satisfactory fitting of the developed model.

Observed y	Simulated y'
$\frac{x_i \cdot v}{O(S - S)}$	$\frac{x_i \cdot v}{O(S - S)}$
$Q(S_i - S_e)$ (mg/L d)	$(\mathbf{x}_i - \mathbf{x}_e)$
(IIIg/1.u)	(ing/L.u)
0.15625	0.178
0.1538462	0.178
0.2205882	0.17175
0.2142857	0.17175
0.1632653	0.161333
0.1333333	0.146182
0.1232877	0.121367
0.1216216	0.135692
0.12	0.131571
0.1282051	0.128
0.1176471	0.124875
0.1157895	0.123455
0.1073171	0.122118
0.1223404	0.119667
0.115	0.113714
0.1176471	0.10925
0.1165049	0.103862
0.1153846	0.100727
0.1153846	0.097231
0.1017699	0.09675
0.092	0.095857
0.0871212	0.092151
0.0872727	0.090931
0.0928571	0.087036

Table 7: Simulation for Reactor 2





The coefficient of determination, R^2 for reactor 2 in Monod Model yielded 0.842, suggesting a satisfactory fitting of the developed model.

Observed v	Simulated v'
x_i . V	x_i . V
$\overline{\boldsymbol{0}\left(\boldsymbol{S}_{i}-\boldsymbol{S}_{i}\right)}$	$\overline{\boldsymbol{0}\left(\boldsymbol{S}_{i}-\boldsymbol{S}_{i}\right)}$
(mg/L.d)	(mg/L.d)
((111 g , 2200)
0.2068966	0.213833
0.2166667	0.193714
0.2166667	0.193714
0.1911765	0.193714
0.1891892	0.166889
0.1627907	0.1575
0.1071429	0.1575
0.1056338	0.138
0.1027397	0.133357
0.1153846	0.11525
0.1125	0.11525
0.125	0.11525
0.1166667	0.11525
0.1182796	0.109739
0.11	0.108208
0.1089109	0.108208
0.1127451	0.101167
0.1116505	0.098606
0.1153846	0.097853
0.1045455	0.092651
0.0958333	0.092205
0.0804196	0.089569
0.0784314	0.086629
0.0809249	0.085071





4. Conclusions

Coefficient of determination \mathbb{R}^2 was used to verify the respective models. The reactors with aspect ratios of (53:16:30; 4:1:1; 10:3:6) yielded specific growth rate μ_{max} and half saturation constant k_s for noninhibited cassava wastewater treatment process as follows:

Monod model yielded:

Reactor $1-(\mu_{max}) = 10.87$ $day^{-1};$ $k_s = 0.87 mgCODL^{-1};$ $R^2 = 0.917)$ Reactor $2-(\mu_{max}) = 12.82$ $day^{-1};$ $k_s = 0.917$

1.92 $mgCODL^{-1}$; $R^2 = 0.842$) Reactor 3-($\mu_{max} = 13.70$ day⁻¹; $k_s = 2.32$ $mgCODL^{-1}$; $R^2 = 0.830$). The evaluation showed that the kinetic parameters developed vary within acceptable limits for any size of reactor (physical model) considered, thus paving way for application in large treatment plants.

The coefficient of determination, R^2 for reactor 3 in Monod Model yielded 0.830, suggesting a satisfactory fitting of the developed model.

However, investigations of the hydrodynamics to date have not taken into account other factors, which include biogas mixing effects, viscosity changes due to extra- cellular polymer production, biomass particle size and the rate of solid particles/biomass within the reactor. It is suggested that further studies should thus be carried out.

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