

# Supporting Information

## **Synthesis of short-chain-length and medium-chain-length polyhydroxyalkanoate blends from activated sludge by manipulating octanoic acid and nonanoic acid as carbon sources**

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**Table S1**  $^{13}\text{C}$  NMR chemical shifts of 3-hydroxyalkanoate (ppm).

Carbon site	scl-PHAs		mcl-PHAs			
	3HB	3HV	3HHx	3HHp	3HO	3HN
C1	169.1	169.36	169.36	169.36	169.36	169.36
C2	40.74	–	39.05	39.05	39.05	39.05
C3	67.56	71.83	70.54	70.79	70.79	70.79
C4	19.72	–	35.85	33.44	33.71	33.71
C5		9.26	18.27	27.11	24.66	24.96
C6			13.75	22.37	31.47	28.98
C7				13.87	22.44	31.64
C8					13.92	22.51
C9						13.99

Note: The symbols of “–” indicate that no signal peaks responding to the C2 and C4 in  $^{13}\text{C}$  NMR spectrum are detected because very low content of 3HV has been extracted from active sludge.

1 **Table S2** GC-MS analysis of methyl esterified samples PHAs produced from active sludge under aerobic condition at 30 °C.

Peak	Retention time (min)	m/z	Molecular weight	Methyl esterified PHA monomer structure	Similarity index (%)
P1	2.725	25, 29, 43, 59, 74, 87, 103, 119, 133, 149, 179	118	Methyl ester of 3HB	95
P2	3.833	26, 31, 43, 59, 71, 85, 103, 114, 131	132	Methyl ester of 3HV	91
P3	5.183	26, 41, 43, 61, 71, 97, 103, 113, 128, 145	146	Methyl ester of 3HHx	84
P4	6.752	27, 41, 43, 61, 74, 85, 103, 113, 127, 142, 159	160	Methyl ester of 3HHp	88
P5	8.283	27, 41, 43, 61, 74, 96, 103, 125, 141, 156, 173	174	Methyl ester of 3HO	89
P6	9.725	27, 41, 43, 61, 74, 96, 103, 113, 139, 155, 170, 187	188	Methyl ester of 3HN	90

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