

## Supporting Information

### Sugar-based amphiphiles: easily accessible and efficient crude oil spill thickening agents

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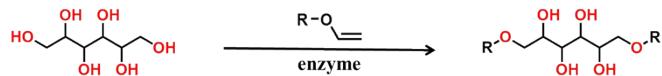
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**Table S1.** Structures of the synthesized amphiphiles with their abbreviations, compound numbers, and both common and systematic names.



a) Mannitol-derived amphiphiles

Abbr.	Comp. #	Structure	Common Name	Systematic Name
M-4	5a		mannitol dibutyrate	(2R,3R,4R,5R)-2,3,4,5-tetrahydroxyhexane-1,6-diyl dibutyrate
M-8	5b		mannitol dicaprylate	(2R,3R,4R,5R)-2,3,4,5-tetrahydroxyhexane-1,6-diyl dioctanoate
M-10	5c		mannitol dicaprate	(2R,3R,4R,5R)-2,3,4,5-tetrahydroxyhexane-1,6-diyl bis(decanoate)
M-12	5d		mannitol dilaurate	(2R,3R,4R,5R)-2,3,4,5-tetrahydroxyhexane-1,6-diyl didodecanoate
M-14	5e		mannitol dimyristate	(2R,3R,4R,5R)-2,3,4,5-tetrahydroxyhexane-1,6-diyl ditetradecanoate

b) Vinyl caprylate-derived amphiphiles

Abbr.	Comp. #	Structure	Common Name	Systematic Name
S-8	6b		sorbitol dicaprylate	(2R,3R,4R,5S)-2,3,4,5-tetrahydroxyhexane-1,6-diyl dioctanoate
G-8*	7b		galactitol dicaprylate	(2R,3S,4R,5S)-2,3,4,5-tetrahydroxyhexane-1,6-diyl dioctanoate
X-8	8b		xylitol dicaprylate	(2R,3r,4S)-2,3,4-trihydroxypentane-1,5-diyl dioctanoate

a) All reactions were carried out at 50 °C for 24 to 48 hrs; b) Abbr. = abbreviation & Comp. = compound;

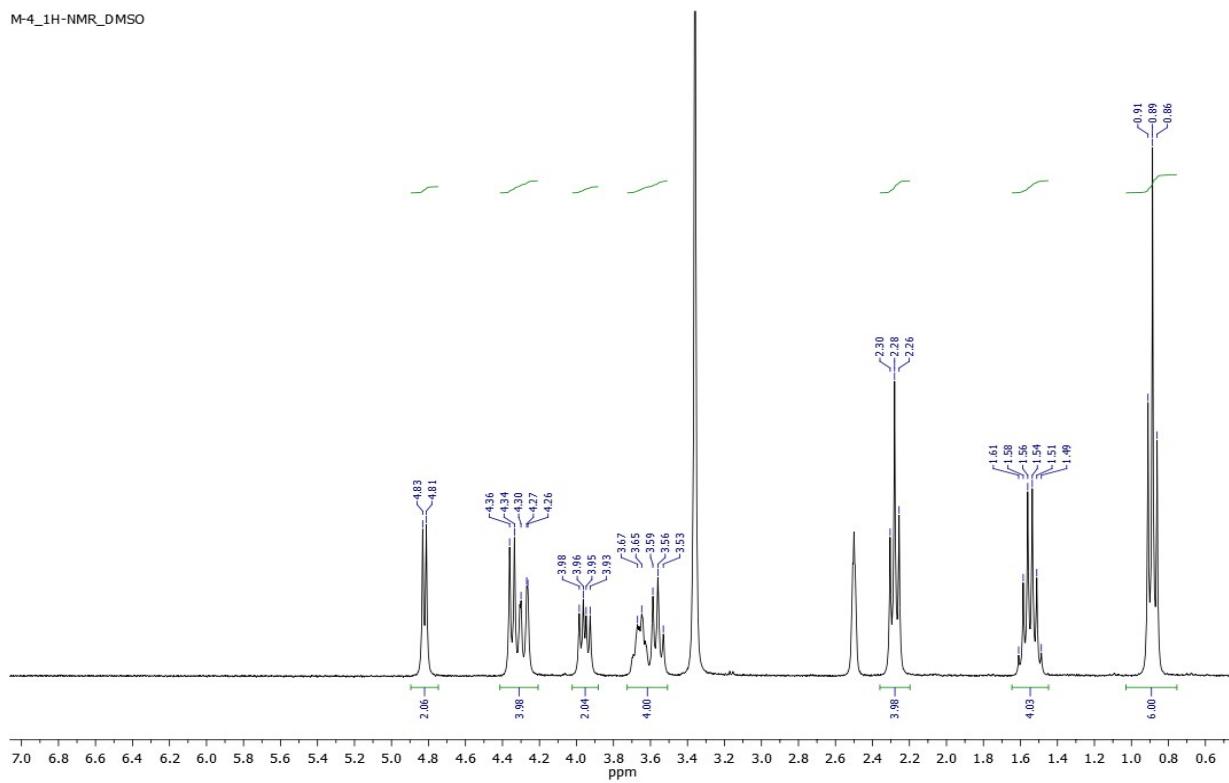
\*G-8 product was never obtained under the enzymatic conditions; hence, the structure shown here is the attempted product rather than the synthesized product.

**Table S2.** Classification and composition of the three used crude oils: South Louisiana Crude Oil (SLCO), Arabian Light Crude Oil (ALCO) and Prudhoe Bay Crude Oil (PBCO).<sup>1,2</sup>

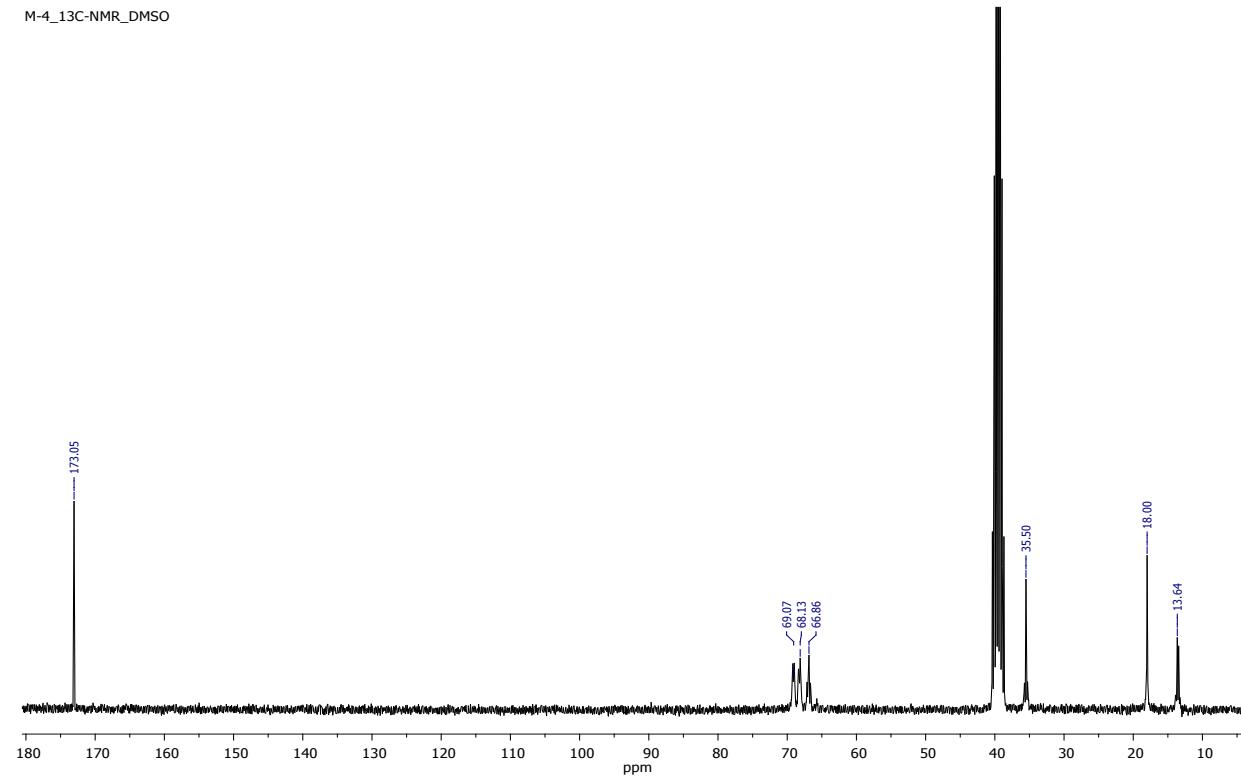
Crude oil:	South Louisiana Crude Oil (SLCO)	Arabian Light Crude Oil (ALCO)	Prudhoe Bay Crude Oil (PBCO)
Physical Properties			
API gravity*	36.61	31.95	26.46
Density (g/cm <sup>3</sup> )	0.840	0.864	0.894
Dynamic Viscosity (cP) (at 15 °C)	7	14	39
Classification	Very Light	Light	Heavy
Chemical Composition			
Sulfur (wt %) (classification)	0.0 (sweet)	2.0 (sour)	0.9 (sour)
Paraffins (% vol)	79	63	27
Naphthenes (% vol)	45	18	36
Aromatics (% vol)	19	19	28
Classification	Paraffinic	Paraffinic	Naphthenic

\* API gravity =  $\frac{141.5}{SG} - 131.5$ , where SG is specific gravity of the crude oil at 15.6 °C (60 °F).

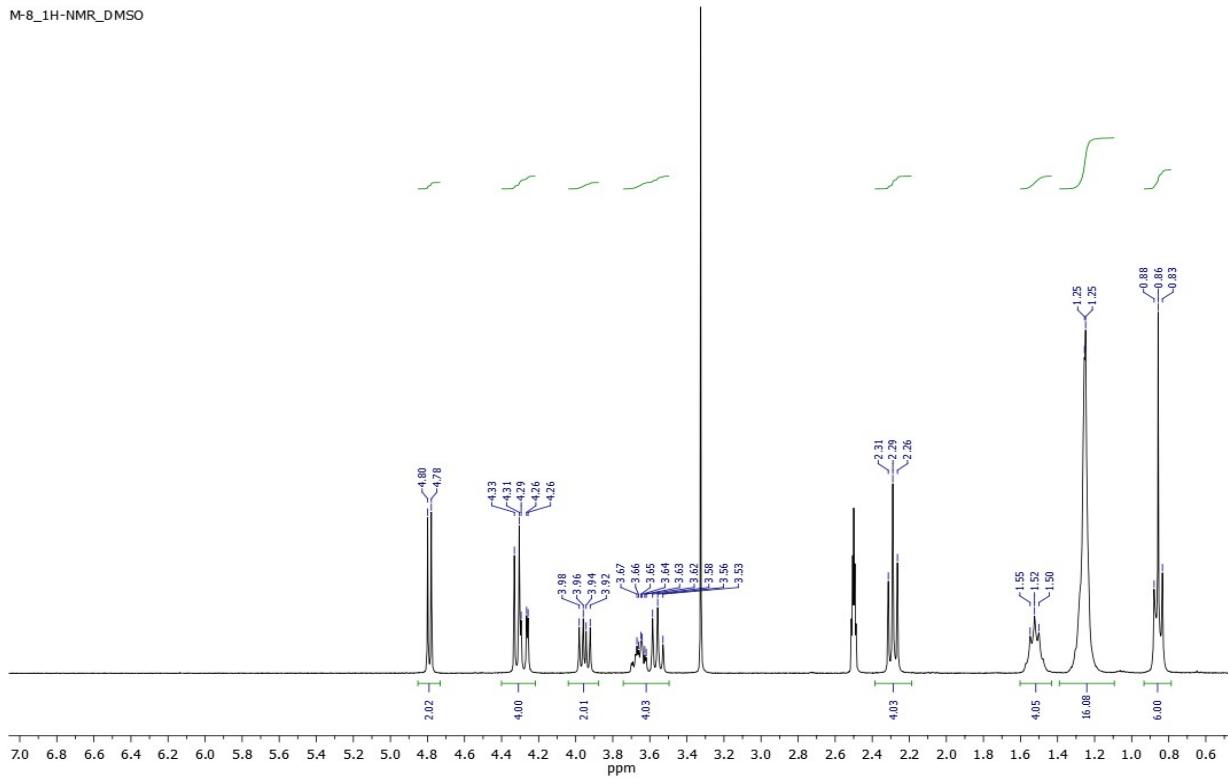
M-4\_1H-NMR\_DMSO

**Fig. S1**  $^1\text{H}$ -NMR of M-4 [(2R,3R,4R,5R)-2,3,4,5-tetrahydroxyhexane-1,6-diyl dibutyrate].

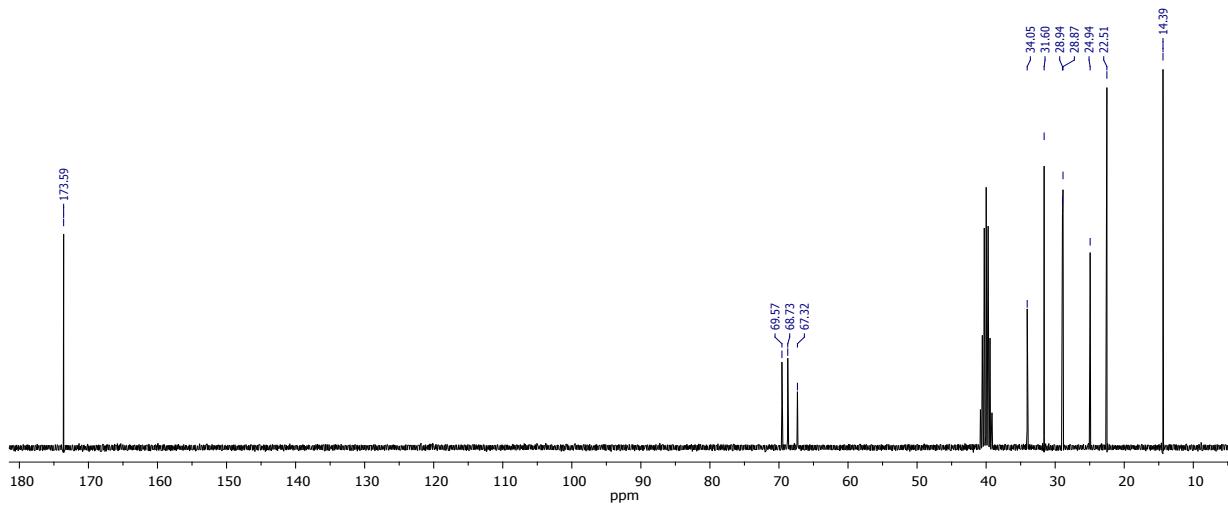
M-4\_13C-NMR\_DMSO

**Fig. S2**  $^{13}\text{C}$ -NMR of M-4 [(2R,3R,4R,5R)-2,3,4,5-tetrahydroxyhexane-1,6-diyl dibutyrate].

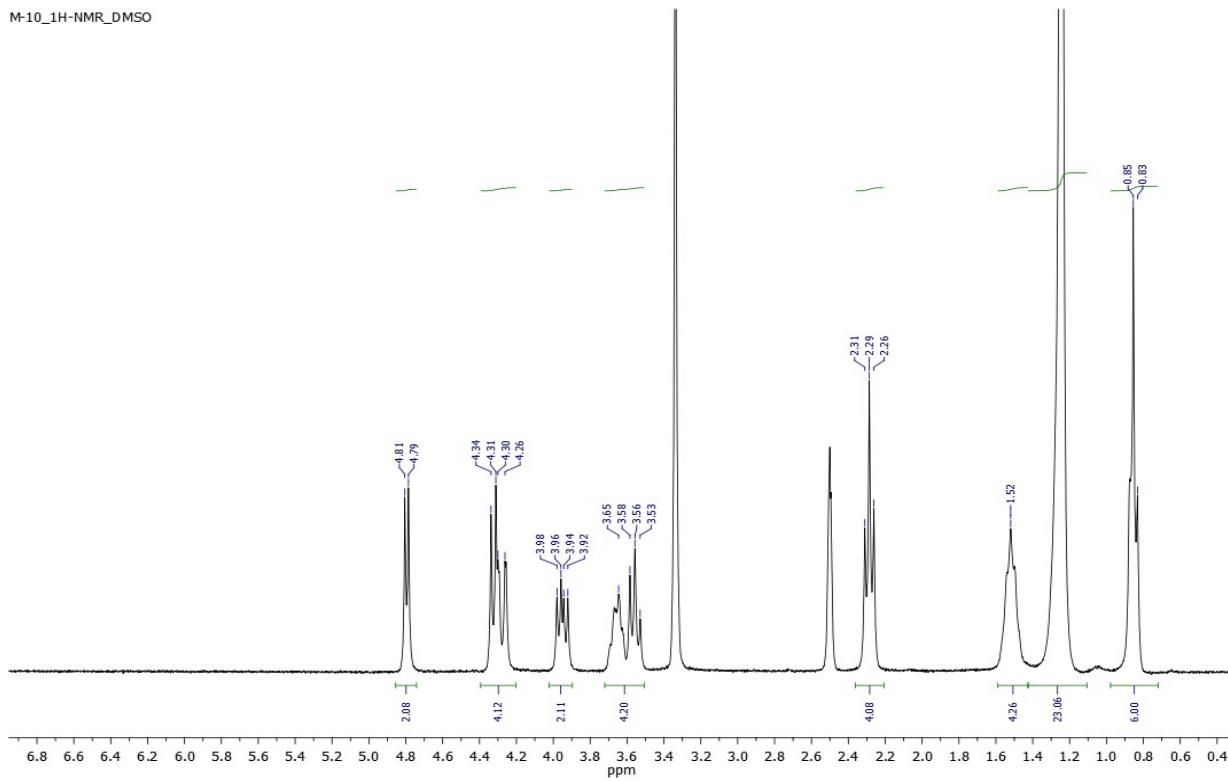
M-8\_1H-NMR\_DMSO

**Fig. S3**  $^1\text{H}$ -NMR of M-8 [(2R,3R,4R,5R)-2,3,4,5-tetrahydroxyhexane-1,6-diyl dioctanoate].

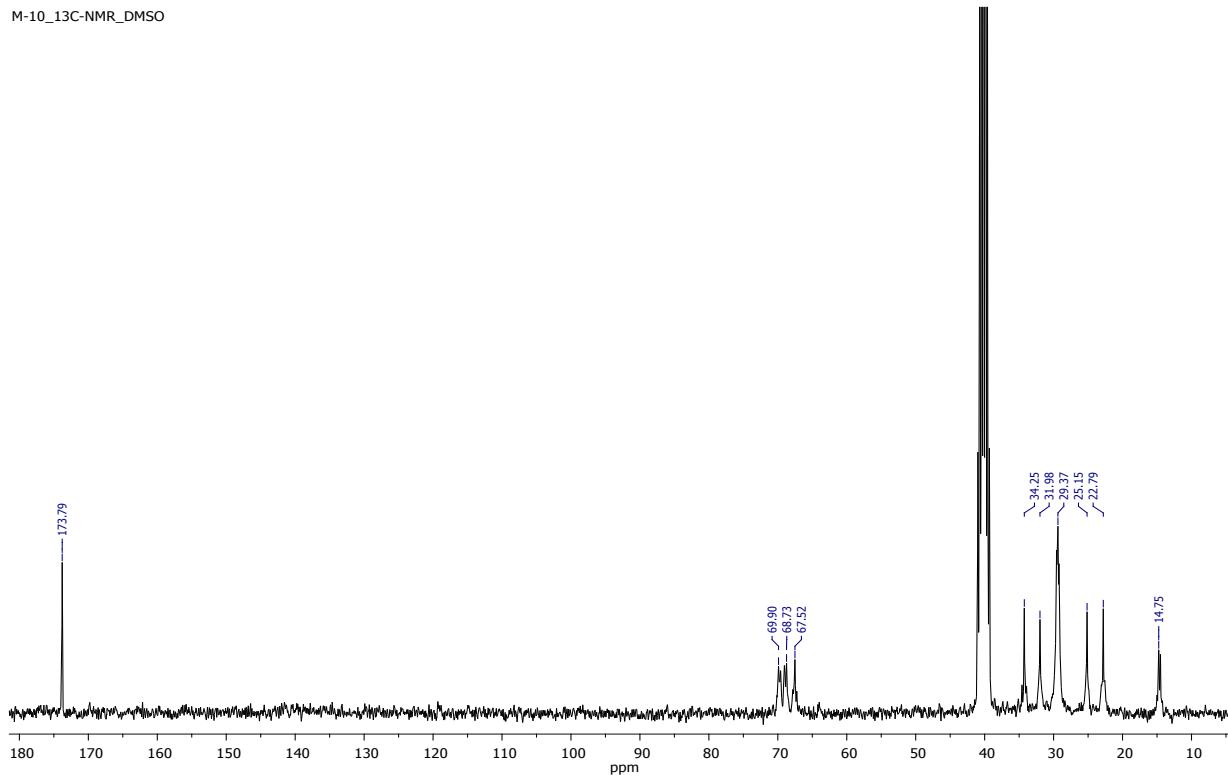
M-8\_13C-NMR\_DMSO

**Fig. S4**  $^{13}\text{C}$ -NMR of M-8 [(2R,3R,4R,5R)-2,3,4,5-tetrahydroxyhexane-1,6-diyl dioctanoate].

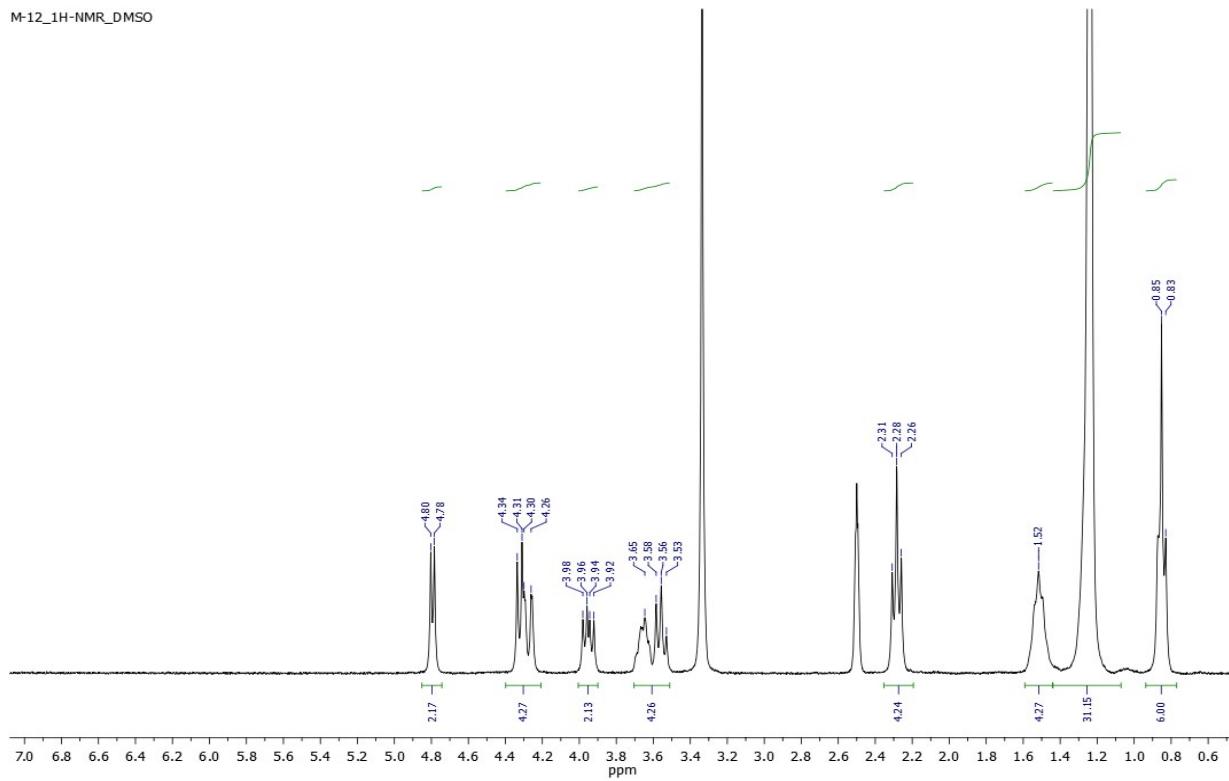
M-10\_1H-NMR\_DMSO

**Fig. S5**  $^1\text{H}$ -NMR of M-10 [(2R,3R,4R,5R)-2,3,4,5-tetrahydroxyhexane-1,6-diyl bis(decanoate)].

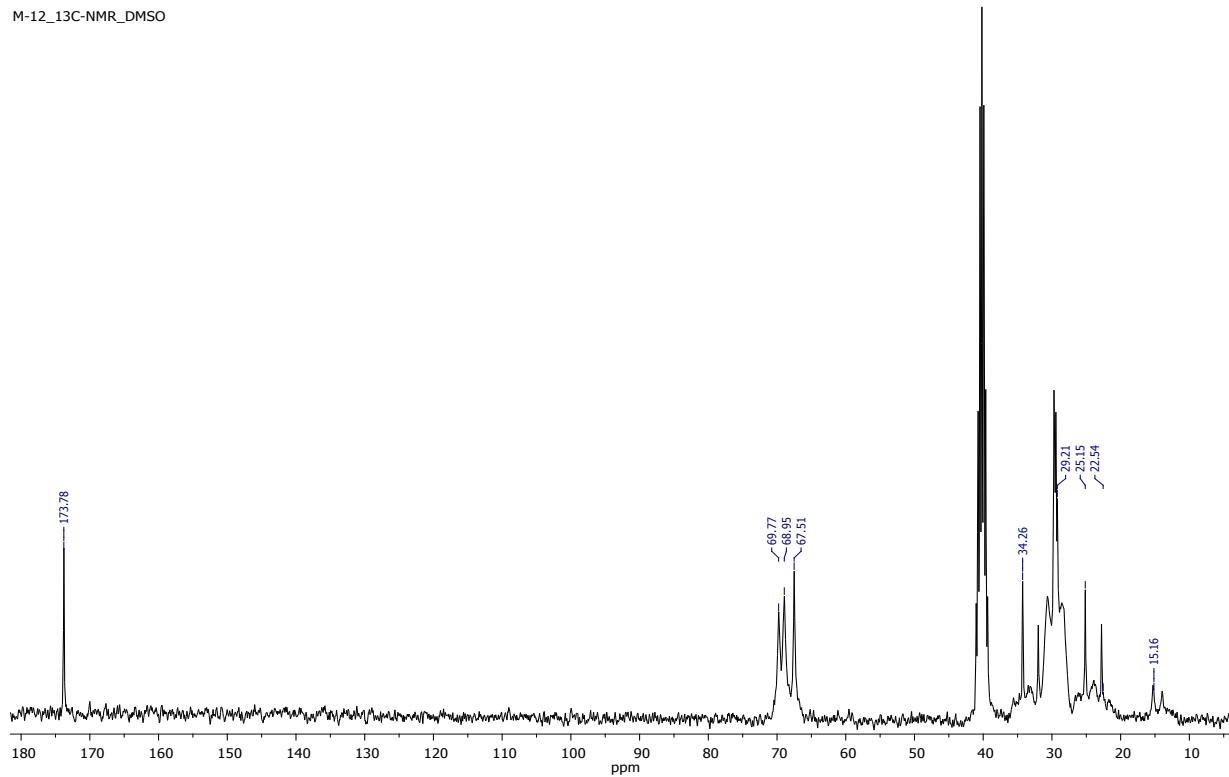
M-10\_13C-NMR\_DMSO

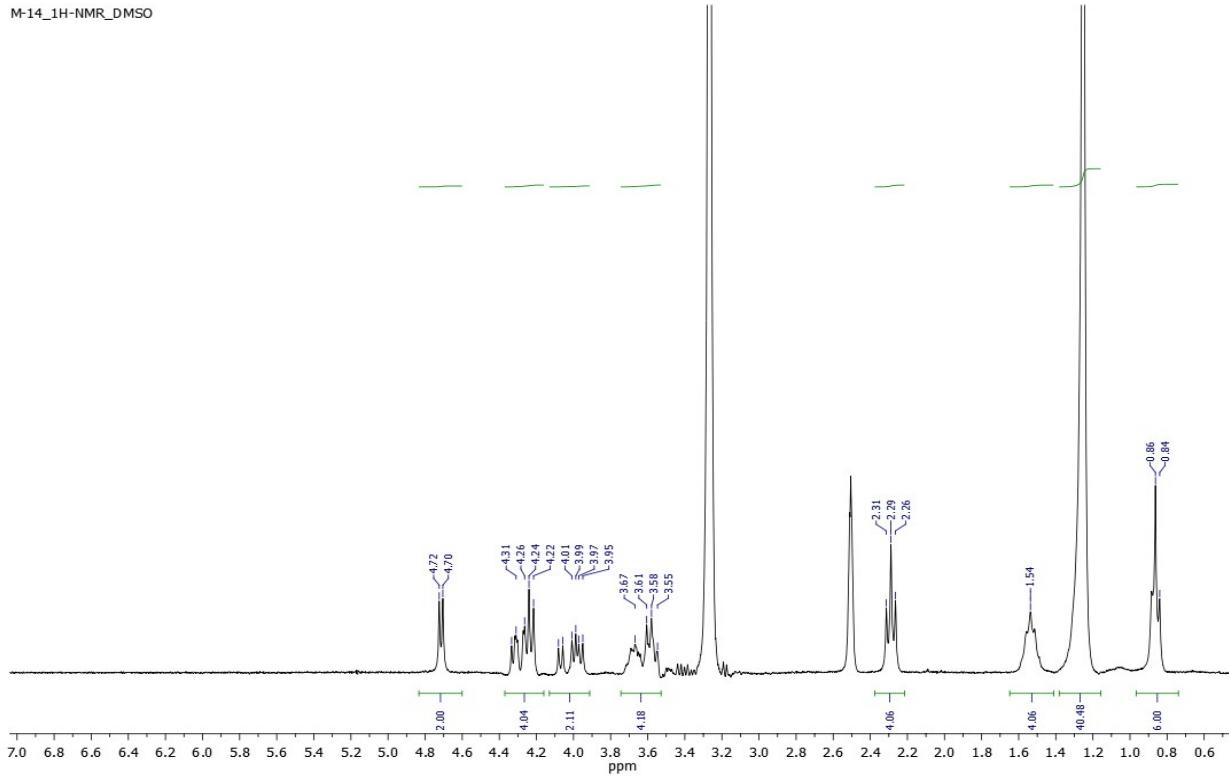
**Fig. S6**  $^{13}\text{C}$ -NMR of M-10 [(2R,3R,4R,5R)-2,3,4,5-tetrahydroxyhexane-1,6-diyl bis(decanoate)].

M-12\_1H-NMR\_DMSO

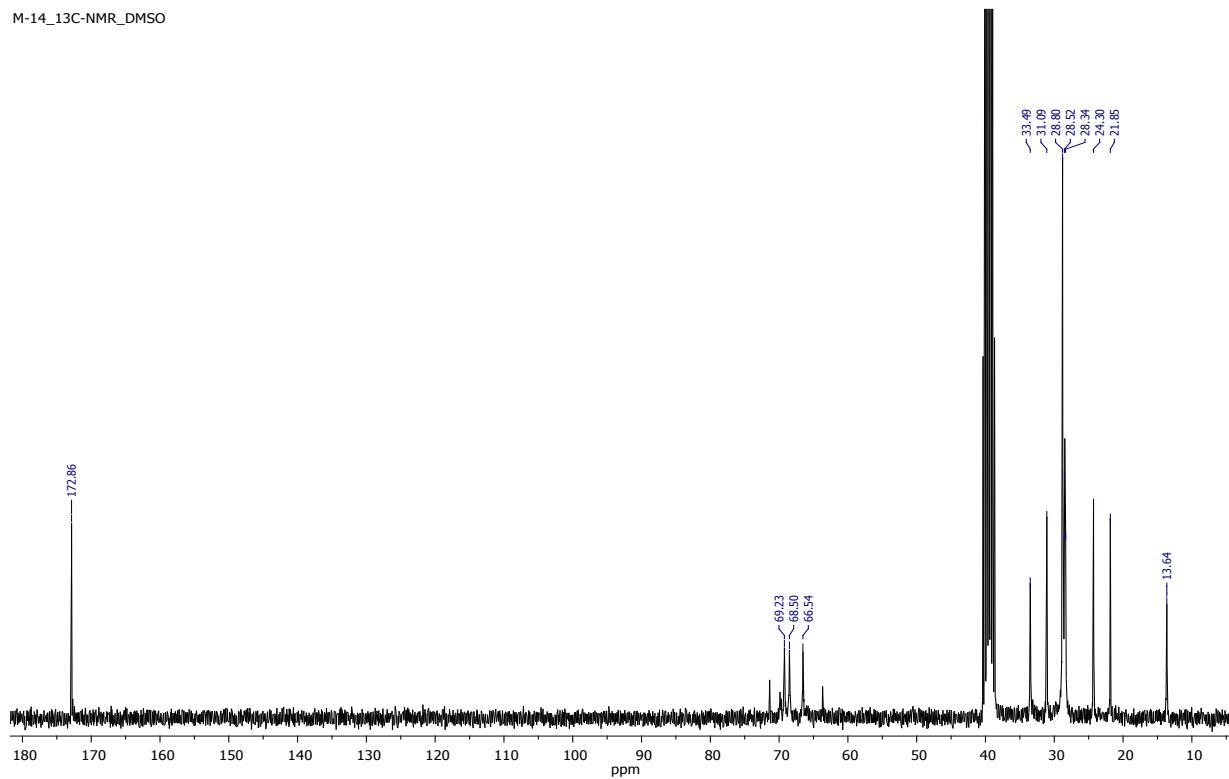
**Fig. S7** <sup>1</sup>H-NMR of M-12 [(2R,3R,4R,5R)-2,3,4,5-tetrahydroxyhexane-1,6-diyl didodecanoate].

M-12\_13C-NMR\_DMSO

**Fig. S8** <sup>13</sup>C-NMR of M-12 [(2R,3R,4R,5R)-2,3,4,5-tetrahydroxyhexane-1,6-diyl didodecanoate].

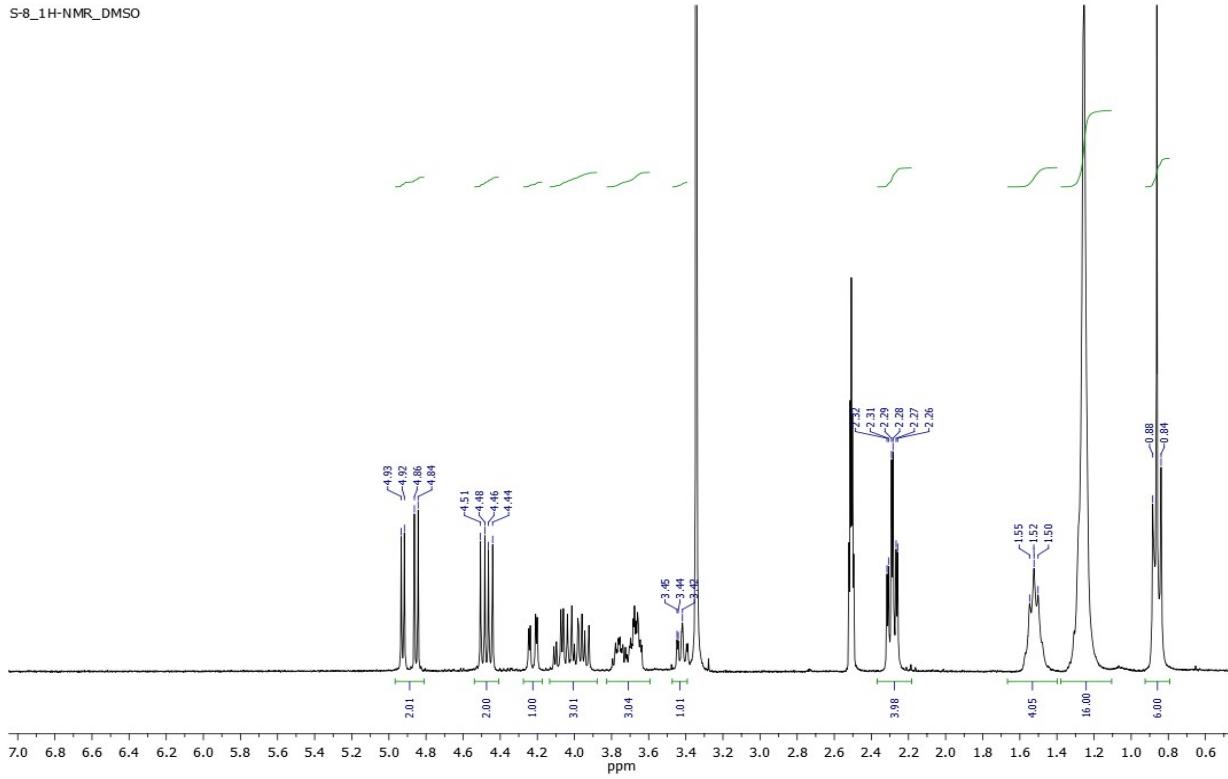


**Fig. S9**  $^1\text{H}$ -NMR of M-14 [(2R,3R,4R,5R)-2,3,4,5-tetrahydroxyhexane-1,6-diyl ditetradecanoate].

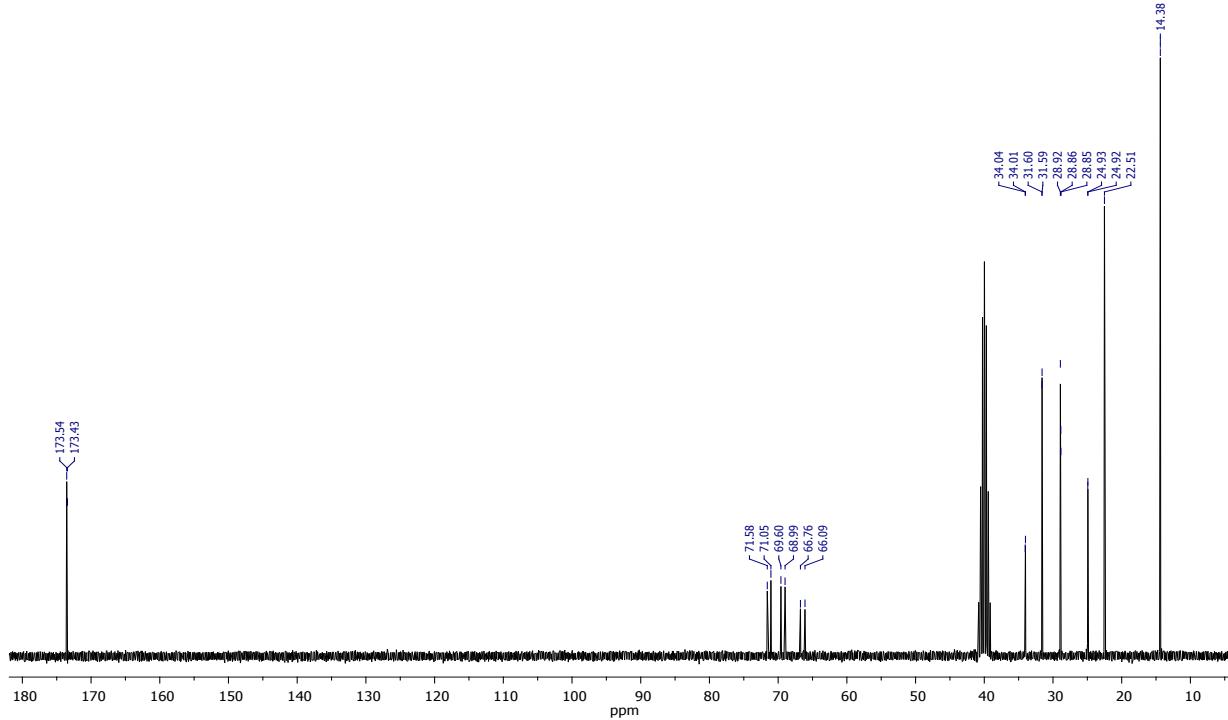


**Fig. S10.**  $^{13}\text{C}$ -NMR of M-14 [(2R,3R,4R,5R)-2,3,4,5-tetrahydroxyhexane-1,6-diyl ditetradecanoate].

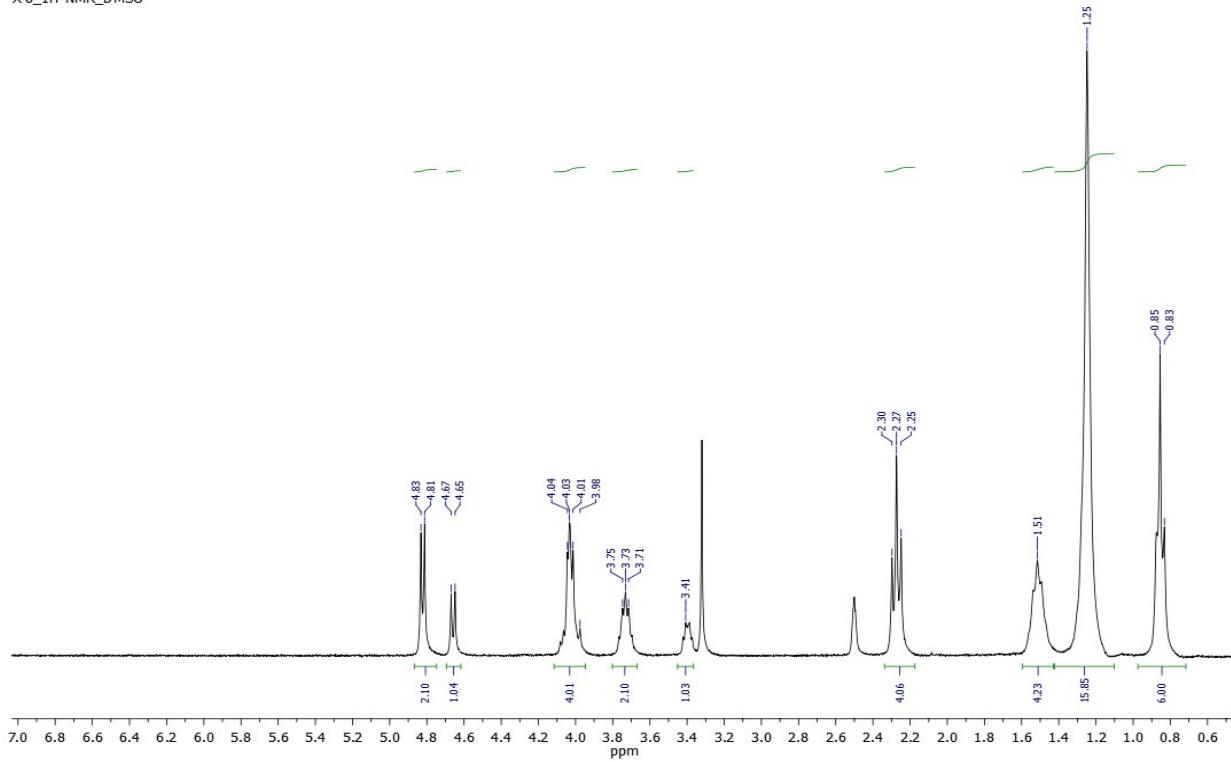
S-8\_1H-NMR\_DMSO

**Fig. S11**  $^1\text{H}$ -NMR of S-8 [(2R,3R,4R,5S)-2,3,4,5-tetrahydroxyhexane-1,6-diyl dioctanoate].

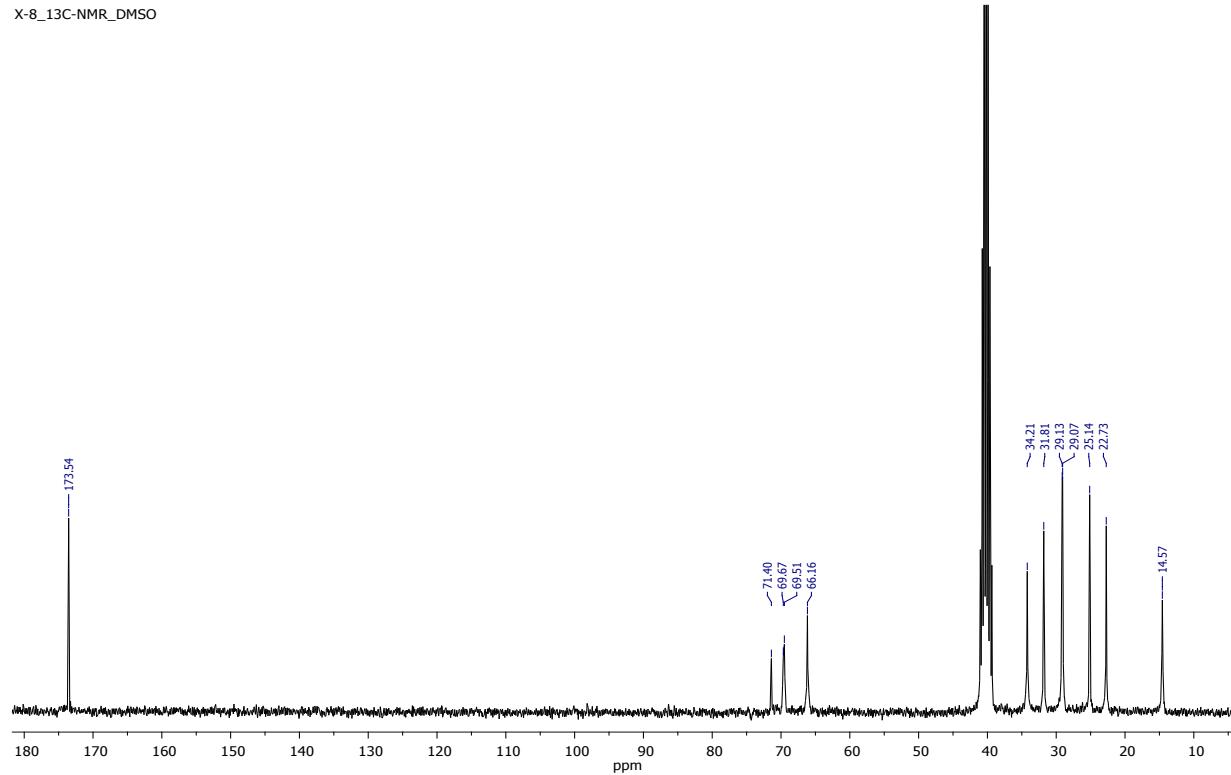
S-8\_13C-NMR\_DMSO

**Fig. S12**  $^{13}\text{C}$ -NMR of S-8 [(2R,3R,4R,5S)-2,3,4,5-tetrahydroxyhexane-1,6-diyl dioctanoate]

X-8\_1H-NMR\_DMSO

**Fig. S13**  $^1\text{H}$ -NMR of X-8 [(2R,3r,4S)-2,3,4-trihydroxypentane-1,5-diyl dioctanoate].

X-8\_13C-NMR\_DMSO

**Fig. S14**  $^{13}\text{C}$ -NMR of X-8 [(2R,3r,4S)-2,3,4-trihydroxypentane-1,5-diyl dioctanoate].

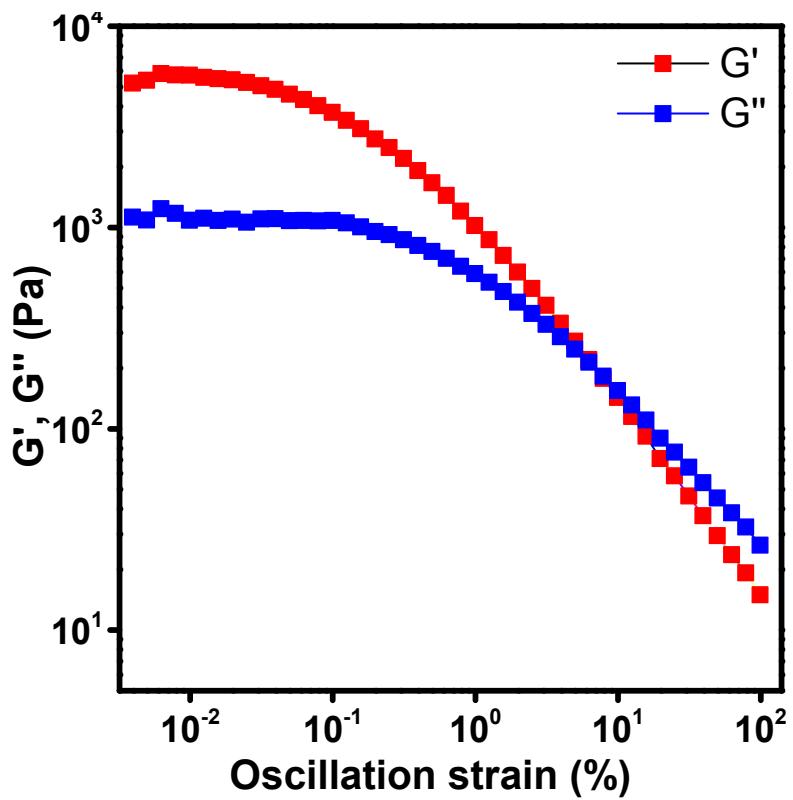


Fig. S15 Rheology data of 5% M-8 in crude oil (SLCO): strain sweep.

## References

1. E. L. Holder, R. N. Conmy and A. D. Venosa, *J. Environ. Prot. (Irvine, Calif.)*, 2015, **06**, 628-639.
2. The American Petroleum Institute Petroleum HPV Testing Group, US EPA: *HIGH PRODUCTION VOLUME (HPV) CHEMICAL CHALLENGE PROGRAM*, 2011, **Consortium Registration # 1100997**, 1-108.