

Tris-imidazolium and benzimidazolium ionic liquids: A new class of biodegradable surfactants

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Supplementary Information

Structural and characterization data

Tris-((2-chloro-acetayloxy)methyl)ethane (3)

Colourless crystals; Yield: 27.94 g (96%); m.p 42-44°C. Molecular Formula: C₁₁H₁₅Cl₃O₆; Mol. Wt.: 349.59; FTIR (cm⁻¹): 2960, 2852 (C-H)_{Aliph}, 1732 (C=O), 1174, 1150 (O-C), 789 (C-Cl); ¹H-NMR (400 MHz, CDCl₃) δ ppm: 4.08 (s, 6H, CH₂-Cl), 4.04 (s, 6H, CH₂-O), 1.02 (s, 3H, CH₃); ¹³C-NMR (100 MHz, CDCl₃) δ ppm: 167.00 (C=O), 66.98 (CH₂-O), 40.73 (CH₂-Cl), 38.42 (-C-), 16.92 (CH₃); EIMS (m/z): 347.9 (10%)(M), 273.0 (12%), 197.0 (45%), 121.0 (66%), 90.9 (100%).

1-Hexyl-1H-imidazole (6c)

This compound was prepared by using 1-bromohexane (10.10 g, 8.60 mL, 61.2 mmole) to give pale yellow oil in 91% yield (8.48 g). Molecular Formula: C₉H₁₆N₂; Mol. Wt.: 152.24; FTIR (cm⁻¹): 3106 (C-H)_{Ar}, 2955, 2929, 2858 (C-H)_{Aliph}, 1506 (C=N), 1460 (C=C)_{Ar}; ¹H-NMR (400 MHz, CDCl₃) δ ppm: 7.29 (s, H, C-H_{Imidazole}), 6.88 (s, H, C-H_{Imidazole}), 6.75 (s, H, C-H_{Imidazole}), 3.76 (t, J = 7.02 Hz, 2H, α-CH₂), 1.64-1.56 (m, 2H, β-CH₂), 1.13 (bs, 6H, bulk-CH₂), 0.72 (t, J = 6.71 Hz, 3H, ω-CH₃); ¹³C-NMR (100 MHz, CDCl₃) δ ppm: 136.95 (CH_{Imidazole}), 129.19 (CH_{Imidazole}), 118.75 (CH_{Imidazole}), 46.91 (α-CH₂), 31.17 (ω-2), 30.98 (bulk-CH₂), 26.02 (β), 22.35 (ω-1), 13.90 (ω); EIMS (m/z): 152.1 (15%)(M), 137.0 (3%), 125.1 (35%), 109.1 (6%), 96.0 (19%), 84.0 (89%), 69.0 (23%), 49.0 (100%).

1-Octyl-1H-imidazole (6d)

This compound was prepared by using 1-bromo-octane (11.82 g, 10.65 mL, 61.2 mmole) to give pale yellow oil in 94% yield (10.37 g). Molecular Formula: C₁₁H₂₀N₂; Mol. Wt.: 180.29; FTIR (cm⁻¹): 3108 (C-H)_{Ar}, 2958, 2925, 2855 (C-H)_{Aliph}, 1677 (C=N), 1506, 1461 (C=C)_{Ar}; ¹H-NMR (400 MHz, CDCl₃) δ ppm: 7.32 (s, H, C-H_{Imidazole}), 6.91 (s, H, C-H_{Imidazole}), 6.77 (s, H, C-H_{Imidazole}), 3.78 (t, J = 7.07 Hz, 2H, α-CH₂), 1.66-1.59 (m, 2H, β-CH₂), 1.15 (bs, 10H, bulk-CH₂), 0.75 (t, J = 7.07 Hz, 3H, ω-CH₃); ¹³C-NMR (100 MHz, CDCl₃) δ ppm: 137.04 (CH_{Imidazole}), 129.28 (CH_{Imidazole}), 118.77 (CH_{Imidazole}), 47.01 (α-CH₂), 31.72 (ω-2), 30.89, 29.09, 29.02 (bulk-CH₂), 26.52 (β), 22.60 (ω-1), 14.06 (ω); EIMS (m/z): 180.1 (45%)(M), 165.1 (10%), 151.1 (24%), 137.1 (20%), 109.1 (36%), 96.0 (40%), 82.0 (100%), 69.0 (55%), 55.0 (25%), 43.1 (18%).

1-Decyl-1H-imidazole (6e)

This compound was prepared by using 1-bromodecane (13.54 g, 12.65 mL, 61.2 mmole) to give pale yellow oil in 93% yield (11.86g). Molecular Formula: C₁₃H₂₄N₂; Mol. Wt.: 208.34; FTIR (cm⁻¹): 3107 (C-

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H)_{Ar}, 2956, 2925, 2855 (C-H)_{Aliph}, 1678(C=N), 1506, 1461 (C=C)_{Ar}; ¹H-NMR (400 MHz, CDCl₃) δ ppm: 7.38 (s, H, C-H_{Imidazole}), 6.97 (s, H, C-H_{Imidazole}), 6.83 (s, H, C-H_{Imidazole}), 3.85 (t, J = 7.07 Hz, 2H, α-CH₂), 1.73-1.66 (m, 2H, β-CH₂), 1.19 (bs, 14H, bulk-CH₂), 0.81 (t, J = 7.07Hz, 3H, ω-CH₃); ¹³C-NMR (100 MHz, CDCl₃) δ ppm: 137.07 (CH_{Imidazole}), 129.33 (CH_{Imidazole}), 118.79 (CH_{Imidazole}), 47.06 (α-CH₂), 31.88 (ω-2), 31.11, 29.50, 29.46, 29.28, 29.10 (bulk -CH₂), 26.57 (β), 22.69 (ω-1), 14.13(ω); EIMS (m/z): 207.2 (32%)(M⁺), 193.1 (9%), 179.1 (18%), 165.1 (14%), 151.1 (18%), 137.1 (23%), 123.1 (35%), 109.1 (28%), 96.0 (48%), 82.0 (100%), 69.0 (48%), 55.0 (45%), 43.1 (29%).

1-Dodecyl-1H-imidazole (6f)

This compound was prepared by using 1-bromododecane (15.25 g, 14.7 mL, 61.2 mmole) to give pale yellow oil in 96% yield (13.88 g). Molecular Formula: C₁₅H₂₈N₂; Mol. Wt.: 236.40; FTIR (cm⁻¹): 3010 (C-H)_{Ar}, 2955, 2924, 2854 (C-H)_{Aliph}, 1672 (C=N), 1505, 1495, 1452 (C=C)_{Ar}; ¹H-NMR (400 MHz, CDCl₃) δ ppm: 7.29 (s, H, C-H_{Imidazole}), 6.88 (s, H, C-H_{Imidazole}), 6.74 (s, H, C-H_{Imidazole}), 3.76 (t, J = 7.09 Hz, 2H, α-CH₂), 1.64-1.57 (m, 2H, β-CH₂), 1.11 (bs, 18H, bulk-CH₂), 0.74 (t, J = 7.09 Hz, 3H, ω-CH₃); ¹³C-NMR (100 MHz, CDCl₃) δ ppm: 136.26 (CH_{Imidazole}), 128.89 (CH_{Imidazole}), 118.34 (CH_{Imidazole}), 46.60 (α-CH₂), 31.52 (ω-2), 30.70, 29.22 (2), 29.13, 29.04, 28.95, 28.68 (bulk-CH₂), 26.15 (β), 22.29 (ω-1), 13.71 (ω); EIMS (m/z): 235.2 (40%)(M⁺), 221.2 (32%), 207.2 (38%), 193.1 (22%), 179.1 (28%), 165.1 (14%), 151.1 (25%), 137.1 (18%), 123.1 (35%), 109.1 (30%), 96.0 (48%), 82.0 (100%), 69.0 (35%), 55.0 (45%), 43.1 (35%).

1-Benzyl-1H-imidazole (6g)

This compound was prepared by using benzyl bromide (10.47 g, 7.28 mL, 61.2 mmole) to give brown viscous syrup crystallized after 5-7 days. Re-crystallization from hexane gave off-white crystals in 86% yield (7.75 g); m.p 70-72°C. Molecular Formula: C₁₀H₁₀N₂; Mol. Wt.: 158.20; FTIR (cm⁻¹): 3113, 3028 (C-H)_{Ar}, 2942 (C-H)_{Aliph}, 1664, 1603, 1585 (C=N), 1505, 1495, 1449 (C=C)_{Ar}; ¹H-NMR (400 MHz, DMSO-d₆) δ ppm: 7.74 (s, H, C-H_{Imidazole}), 7.46-7.17 (m, 6H, 5H_{Ar}, C-H_{Imidazole}), 6.90 (s, H, C-H_{Imidazole}), 5.18 (s, 2H, Ar-CH₂-N); ¹³C-NMR (100 MHz, DMSO-d₆) δ ppm: 137.91 (-C_{Ar}-CH₂-), 137.80 (CH_{Imidazole}), 129.22 (2CH_{Ar}), 128.89 (CH_{Imidazole}), 128.35 (CH_{Ar}), 128.05 (2CH_{Ar}), 120.19 (CH_{Imidazole}), 50.13 (Ar-CH₂-); EIMS (m/z): 158.0 (15%)(M), 131.0 (10%), 118.0 (15%), 104.0 (11%), 91.0 (100%), 77.0 (7.5%), 65.0 (14%).

1-Butyl-1H-benzimidazole (7b)

This compound was prepared by using 1-bromobutane (8.38 g, 6.57 mL, 61.2 mmole) to give yellow oil in 83% yield (8.84 g). Molecular Formula: C₁₁H₁₄N₂; Mol. Wt.: 174.24; FTIR (cm⁻¹): 3055 (C-H)_{Ar}, 2955, 2930, 2855 (C-H)_{Aliph}, 1612 (C=N), 1494, 1450 (C=C)_{Ar}; ¹H-NMR (400 MHz, DMSO-d₆) δ ppm: 8.21 (s, H, C-H_{BImidazole}), 7.64 (d, J= 8.05 Hz, H, C-H_{Ar}), 7.58 (d, J= 8.05 Hz, H, C-H_{Ar}), 7.26-7.16 (m, 2H, CH_{Ar}), 4.23 (t, J= 7.07 Hz, 2H, α-CH₂), 1.72-1.79 (m, 2H, β-CH₂), 1.28-1.18 (m, 2H, (ω-1)), 0.87 (t, J=7.23 Hz, 3H, ω-CH₃); ¹³C-NMR (100 MHz, DMSO-d₆) δ ppm: 144.49 (CH_{BImidazole}), 144.07 (C_{Ar}), 134.35 (C_{Ar}), 122.68 (CH_{Ar}), 121.85 (CH_{Ar}), 119.97 (CH_{Ar}), 110.82 (CH_{Ar}), 44.30 (α-CH₂), 31.98 (ω-2), 19.87 (ω-1),

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13.86 (ω); EIMS (m/z): 174.1 (50%)(M), 159.0 (12%), 145.0 (18%), 131.0 (100%), 118.0 (30%), 104.0 (12%), 90.0 (10%) 77.0 (25%).

1-Hexyl-1H-benzimidazole (7c)

This compound was prepared by using 1-bromohexane (10.10 g, 8.63 mL, 61.2 mmole) to give yellow oil in 88% yield (10.90 g). Molecular Formula: C₁₃H₁₈N₂; Mol. Wt.: 202.30; FTIR (cm⁻¹): 3058 (C-H)_{Ar}, 2954, 2928, 2857 (C-H)_{Aliph}, 1615 (C=N), 1494, 1457 (C=C)_{Ar}; ¹H-NMR (400 MHz, DMSO-d₆) δ ppm: 8.23 (s, H, C-H_{BImidazole}), 7.69 (d, J = 7.56 Hz, H, C-H_{Ar}), 7.54 (d, J = 7.56 Hz, H, C-H_{Ar}), 7.25-7.17 (m, 2H, CH_{Ar}), 4.18 (t, J = 7.07 Hz, 2H, α -CH₂), 1.76-1.69 (m, β -CH₂), 1.18 (bs, 6H, bulk-CH₂), 0.78 (t, J = 6.34 Hz, 3H, ω -CH₃); ¹³C-NMR (100 MHz, DMSO-d₆) δ ppm: 143.86 (CH_{BImidazole}), 143.55 (C_{Ar}), 133.77 (C_{Ar}), 122.05 (CH_{Ar}), 121.22 (CH_{Ar}), 119.40 (CH_{Ar}), 110.17 (CH_{Ar}), 44.02 (α -CH₂), 30.67 (ω -2), 29.32 (bulk-CH₂), 25.73 (β), 21.93 (ω -1), 13.78 (ω); EIMS (m/z): 202.1 (43%)(M), 187.1 (5%), 173.1 (18%), 159.0 (15%), 145.0 (15%), 131.0 (100%), 118.0 (19%), 104.0 (12%), 90.0 (7%) 77.0 (17%).

1-Octyl-1H-benzimidazole (7d)

This compound was prepared by using 1-bromoocetane (11.82 g, 10.65 mL, 61.2 mmole) to give yellow oil in 90% yield (12.69 g). Molecular Formula: C₁₅H₂₂N₂; Mol. Wt.: 230.35; FTIR (cm⁻¹): 3058(C-H)_{Ar}, 2955, 2925, 2854 (C-H)_{Aliph}, 1615 (C=N), 1494, 1458 (C=C)_{Ar}; ¹H-NMR (400 MHz, DMSO-d₆) δ ppm: 8.21 (s, H, C-H_{BImidazole}), 7.64 (d, J = 7.07 Hz, H, C-H_{Ar}), 7.56 (d, J = 7.07 Hz, H, C-H_{Ar}), 7.25-7.16 (m, 2H, CH_{Ar}), 4.20 (t, J = 7.07 Hz, 2H, α -CH₂), 1.79-1.72 (m, β -CH₂), 1.17 (bs, 10H, bulk-CH₂), 0.80 (t, J = 7.07 Hz, 3H, ω -CH₃); ¹³C-NMR (100 MHz, DMSO-d₆) δ ppm: 144.39 (CH_{BImidazole}), 144.06 (C_{Ar}), 134.29 (C_{Ar}), 122.60 (CH_{Ar}), 121.77 (CH_{Ar}), 119.94 (CH_{Ar}), 110.69 (CH_{Ar}), 44.58 (α -CH₂), 31.72 (ω -2), 29.92, 29.14, 29.05 (bulk-CH₂), 26.66 (β), 22.59 (ω -1), 14.32 (ω); EIMS (m/z): 230.1 (40%)(M), 215.1 (9%), 201.1 (12%), 187.1 (24%), 173.1 (39%), 159.0 (25%), 145.0 (39%), 131.0 (100%), 118.0 (47%), 104.0 (17%), 90.0 (9%), 77.0 (22%).

1-Decyl-1H-benzimidazole (7e)

This compound was prepared by using 1-bromodecane (13.53 g, 12.65 mL, 61.2 mmole) to give a yellow semi-solid in 91% yield (14.38 g). Molecular Formula: C₁₇H₂₆N₂; Mol. Wt.: 258.40; FTIR (cm⁻¹): 3057 (C-H)_{Ar}, 2954, 2923, 2853 (C-H)_{Aliph}, 1615 (C=N), 1494, 1458 (C=C)_{Ar}; ¹H-NMR (400 MHz, DMSO-d₆) δ ppm: 8.20 (s, H, C-H_{BImidazole}), 7.64 (d, J = 7.32 Hz, H, C-H_{Ar}), 7.56 (d, J = 7.32 Hz, H, C-H_{Ar}), 7.25-7.16 (m, 2H, CH_{Ar}), 4.21 (t, J = 7.16 Hz, 2H, α -CH₂), 1.79-1.72 (m, 2H, β -CH₂), 1.18 (bs, 14H, bulk-CH₂), 0.82 (t, J = 7.07 Hz, 3H, ω -CH₃); ¹³C-NMR (100 MHz, DMSO-d₆) δ ppm: 144.34 (CH_{BImidazole}), 143.40 (C_{Ar}), 134.05 (C_{Ar}), 122.93 (CH_{Ar}), 122.13 (CH_{Ar}), 119.73 (CH_{Ar}), 110.82 (CH_{Ar}), 44.63 (α -CH₂), 31.65 (ω -2), 29.69, 29.56, 29.24, 29.02, 28.88 (bulk-CH₂), 26.43 (β), 22.48 (ω -1), 14.30 (ω); EIMS (m/z): 258.2 (60%)(M), 243.2 (16%), 229.1 (32%), 215.1 (27%), 201.1 (30%), 187.1 (36%), 173.1 (53%), 159.1 (39%), 145.0 (50%), 131.0 (100%), 118.0 (70%), 104.0 (21%), 90.0 (10%), 77.0 (25%).

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1-Dodecyl-1H-benzimidazole (7f)

This compound was prepared by using 1-bromododecane (15.25 g, 14.70 mL, 61.2 mmole) to give yellow semi-solid in 94% yield (16.48 g). Molecular Formula: C₁₉H₃₀N₂; Mol. Wt.: 286.45; FTIR (cm⁻¹): 3055 (C-H)_{Ar}, 2958, 2924, 2857 (C-H)_{Aliph}, 1612 (C=N), 1496, 1458 (C=C)_{Ar}; ¹H-NMR (400 MHz, CDCl₃) δ ppm: 7.84 (s, H, C-H_{BImidazole}), 7.82-7.77 (m, H, C-H_{Ar}), 7.37-7.34 (m, H, C-H_{Ar}), 7.29-7.23 (m, 2H, CH_{Ar}), 4.09 (t, J= 7.09 Hz, 2H, α-CH₂), 1.87-1.80 (m, 2H, β-CH₂), 1.23 (bs, 18H, (bulk-CH₂)), 0.87 (t, J= 7.09 Hz, 3H, ω-CH₃). ¹³C-NMR (100 MHz, CDCl₃) δ ppm: 143.75 (C_{Ar}), 142.74 (CH_{BImidazole}), 133.68 (C_{Ar}), 122.57 (CH_{Ar}), 121.78 (CH_{Ar}), 120.17 (CH_{Ar}), 109.49 (CH_{Ar}), 44.89 (α-CH₂), 31.74 (ω-2), 29.62, 29.42 (2), 29.43, 29.25, 29.16, 28.91 (bulk-CH₂), 26.63 (β), 22.51 (ω-1), 13.94 (ω). EIMS (m/z): 286.2 (62%)(M), 271.2 (18%), 257.2 (30%), 243.2 (32%), 229.1 (30%), 215.1 (32%), 201.1 (57%), 187.1 (40%), 173.1 (50%), 159.1 (40%), 145.0 (55%), 131.0 (100%), 118.0 (60%), 104.0 (21%), 90.0 (10%), 77.0 (20%).

1-Benzyl-1H-benzimidazole (7g)

This compound was prepared by using benzyl bromide (10.47 g, 7.28 mL, 61.2 mmole) to give a brown solid re-crystallized from hexane gave off-white crystals in 90% yield (11.47 g); m.p 108-110°C. Molecular Formula: C₁₄H₁₂N₂; Mol. Wt.: 208.26; FTIR (cm⁻¹): 3081, 3032 (C-H)_{Ar}, 2944 (C-H)_{Aliph}, 1666, 1613 (C=N), 1493, 1451, 1442 (C=C)_{Ar}; ¹H-NMR (400 MHz, DMSO-d₆) δ ppm: 8.41 (s, H, C-H_{BImidazole}), 7.67-7.64 (m, H, C-H_{Ar}), 7.51-7.49 (m, H, C-H_{Ar}), 7.34-7.14 (m, 7H, C-H_{Ar}), 5.49 (s, 2H, Ar-CH₂-N); ¹³C-NMR (100 MHz, DMSO-d₆) δ ppm: 144.21 (CH_{BImidazole}), 143.57 (C_{BImidazole}), 136.93 (-C_{Ar}-CH₂-), 133.65 (C_{BImidazole}), 128.67 (2CH_{Ar}), 127.70 (CH_{Ar}), 127.35 (2CH_{Ar}), 122.37 (CH_{BImidazole}), 121.56 (CH_{BImidazole}), 119.48 (CH_{BImidazole}), 110.67 (CH_{BImidazole}), 47.60 (Ar-CH₂-); EIMS (m/z): 208.1 (58%)(M), 103.0 (3%), 91.1 (100%), 77.0 (3%), 65.0 (13%).

Tris-((N-butyl-imidazoliumyl-acetayloxy)methyl)ethane chloride (8b)

This compound was prepared analogously to **8a** using tris-((2-chloro-acetayloxy)methyl)ethane (compound **3**) (1.9 g, 5.43 mmol) and 1-butylimidazole (2.02 g, 2.14 mL, 16.3 mmol) to provide a viscous hygroscopic syrup in 98% yield (3.85 g). Molecular Formula: C₃₂H₅₁Cl₃N₆O₆; Mol. Wt.: 722.14; FTIR (cm⁻¹): 3058 (C-H)_{Ar}, 2959, 2933, 2873 (C-H)_{Aliph}, 1748 (C=O), 1632 (C=N), 1564, 1464 (C=C)_{Ar}, 1199, 1165 (C-O); ¹H-NMR (400 MHz, DMSO-d₆) δ ppm: 9.56 (bt~s, 3H, C-H_{Imidazole}, major), 9.51 (bt~s, 3H, C-H_{Imidazole}, minor), 9.46 (bt~s, 3H, C-H_{Imidazole}, minor), 7.89 (dt, J= 6.83, 1.71 Hz, 6H, C-H_{Imidazole}, major), 7.85 (dt, J= 6.83, 1.71 Hz, 6H, C-H_{Imidazole}, minor), 5.45 (s, 6H, O-CH₂, major), 5.40 (s, 6H, O-CH₂, minor), 4.25 (t, J= 7.07 Hz, 6H, α-CH₂, major), 4.04 (s, 6H, N-CH₂) 3.95 (t, J= 7.07 Hz, 6H, α-CH₂, minor), 1.80-1.73 (m, 6H, β-CH₂, major), 1.69-1.62 (m, 6H, β-CH₂, minor), 1.28-1.14 (m, 6H, (ω-1)), 0.93 (s, 3H, CH₃), 0.88 (t, J= 7.32 Hz, 9H, ω-CH₃); ¹³C-NMR (100 MHz, DMSO-d₆) δ ppm: 166.72 (C=O), 137.44 (CH_{Imidazole}, major), 137.06 (CH_{Imidazole}, minor), 123.95 (CH_{Imidazole}), 122.14 (CH_{Imidazole}), 66.54 (CH₂-O), 49.60 (CH₂-N), 48.78 (α-CH₂), 38.16 (-C-), 32.56 ((ω-2), minor), 31.40 ((ω-2), major), 19.14 ((ω-1), minor), 18.79 ((ω-1), major), 16.38 (CH₃), 13.42 ((ω), minor), 13.32 ((ω), major); HRMS: m/z, [M⁺³-2H]-3Cl⁻ calcd. for C₃₂H₄₉N₆O₆⁵⁺: 613.3714, found: 613.3748.

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Tris-((*N*-octyl-imidazoliumyl-acetayloxy)methyl)ethane chloride (8d)

This compound was prepared analogously to **8a** using tris-((2-chloro-acetayloxy)methyl)ethane (compound **3**) (1.9 g, 5.43 mmol) and 1-octylimidazole (**6d**) (2.94 g, 16.3 mmol) to provide a viscous hygroscopic syrup in 97% yield (4.70 g). Molecular Formula: $C_{44}H_{75}Cl_3N_6O_6$; Mol. Wt.: 890.46; FTIR (cm⁻¹): 3058 (C-H)_{Ar}, 2955, 2925, 2855 (C-H)_{Aliph}, 1749 (C=O), 1668 (C=N), 1564, 1464 (C=C)_{Ar}, 1199, 1167 (C-O); ¹H-NMR (400 MHz, DMSO-*d*₆) δ ppm: 9.55 (bt~s, 3H, C-H_{Imidazole}, major), 9.48 (bt~s, 3H, C-H_{Imidazole}, minor), 9.41 (bt~s, 3H, C-H_{Imidazole}, minor), 7.89 (dt, 6H, *J*=8.86, 1.72 Hz, C-H_{Imidazole}, major), 7.85 (dt, 6H, *J*=8.86, 1.72 Hz, C-H_{Imidazole}, minor), 5.45 (s, 6H, O-CH₂, major), 5.39 (s, 6H, O-CH₂, minor), 5.37 (s, 6H, O-CH₂, minor), 4.24 (t, *J*=7.15 Hz, 6H, α-CH₂), 4.06 (bs, 6H, N-CH₂), 1.82-1.74 (m, 6H, β-CH₂), 1.24 (bs, 30H, bulk-CH₂), 0.95 (s, 3H, CH₃), 0.85 (t, *J*=6.98 Hz, 9H, ω-CH₃); ¹³C-NMR (100 MHz, DMSO-*d*₆) δ ppm: 167.99 (C=O, major), 166.85 (C=O, minor), 137.31 (CH_{Imidazole}, minor), 137.21 (CH_{Imidazole}, major), 123.87 (CH_{Imidazole}), 122.05 (CH_{Imidazole}, minor), 121.91 (CH_{Imidazole}, major), 63.92 (CH₂-O, major), 63.05 (CH₂-O, minor), 49.75 (CH₂-N, major), 49.63 (CH₂-N, minor), 48.94 (α-CH₂, minor), 48.87 (α-CH₂, major), 40.62 (-C-), 31.10 (ω-2), 29.34, 28.44, 28.26 (bulk-CH₂), 25.39 (β), 22.01 (ω-1), 16.57 (CH₃, major), 16.42 (CH₃, minor), 13.89 (ω); HRMS: m/z, [M⁺³-2H]-3Cl⁻ calcd. for $C_{44}H_{73}N_6O_6^{5+}$: 781.5592, found: 781.5608.

Tris-((*N*-decyl-imidazoliumyl-acetayloxy)methyl)ethane chloride (8e)

This compound was prepared analogously to **8a** using tris-((2-chloro-acetayloxy)methyl)ethane (compound **3**) (1.9 g, 5.43 mmol) and 1-decylimidazole (**6e**) (3.39 g, 16.3 mmol) to provide a viscous hygroscopic syrup in 99% yield (5.24 g). Molecular Formula: $C_{50}H_{87}Cl_3N_6O_6$; Mol. Wt.: 974.62; FTIR (cm⁻¹): 3058 (C-H)_{Ar}, 2955, 2929, 2857 (C-H)_{Aliph}, 1749 (C=O), 1666 (C=N), 1564, 1462 (C=C)_{Ar}, 1199, 1165 (C-O); ¹H-NMR (400 MHz, DMSO-*d*₆) δ ppm: 9.64 (bt~s, 3H, C-H_{Imidazole}, major), 9.57 (bt~s, 3H, C-H_{Imidazole}, minor), 9.51 (bt~s, 3H, C-H_{Imidazole}, minor), 7.94 (t, *J*=1.81 Hz, 6H, C-H_{Imidazole}, major), 7.89 (t, *J*=1.81 Hz, 6H, C-H_{Imidazole}, major), 7.84 (t, *J*=1.81 Hz, 6H, C-H_{Imidazole}, minor), 7.81 (t, *J*=1.81 Hz, 6H, C-H_{Imidazole}, minor), 5.49 (s, 6H, O-CH₂, major), 5.45 (s, 6H, O-CH₂, minor), 5.40 (s, 6H, O-CH₂, minor), 4.25 (t, *J*=7.25 Hz, 6H, α-CH₂, major), 4.05 (s, 6H, N-CH₂), 4.01 (t, 6H, *J*=7.25 Hz, α-CH₂, minor), 1.82-1.75 (m, 6H, β-CH₂, major), 1.73-1.67 (m, 6H, β-CH₂, minor), 1.24 (bs, 42H, bulk-CH₂), 0.95 (s, 3H, CH₃), 0.84 (t, 9H, *J*=6.80 Hz, ω-CH₃); ¹³C-NMR (100 MHz, DMSO-*d*₆) δ ppm: 166.72 (C=O), 137.41 (CH_{Imidazole}, major), 136.42 (CH_{Imidazole}, minor), 123.91 (CH_{Imidazole}), 122.04(CH_{Imidazole}), 66.48 (CH₂-O), 49.52 (CH₂-N), 48.92 (α-CH₂), 38.08 (-C-), 31.15 (ω-2), 30.16 (minor), 29.40 (major), 28.53, 28.49, 28.37, 28.31 (bulk-CH₂), 25.76 (β, minor), 25.47 (β, major), 22.05 (ω-1), 16.39 (CH₃), 13.93 (ω); HRMS: m/z, [M⁺³-2H]-3Cl⁻ calcd. for $C_{50}H_{85}N_6O_6^{5+}$: 865.6531, found: 865.6518.

Tris-((*N*-hexyl-benzimidazoliumyl-acetayloxy)methyl)ethane chloride (9c)

This compound was prepared analogously to **9b** using tris-((2-chloro-acetayloxy)methyl)ethane (compound **3**) (1.9 g, 5.43 mmol) and 1-hexyl-benzimidazole (**7c**) (3.30 g, 16.3 mmol) to provide a viscous hygroscopic syrup in 99% yield (5.15 g). Molecular Formula: $C_{50}H_{69}Cl_3N_6O_6$; Mol. Wt.: 956.48; FTIR (cm⁻¹): 3020 (C-H)_{Ar}, 2954, 2929, 2859 (C-H)_{Aliph}, 1748 (C=O), 1617(C=N), 1563, 1486, 1464 (C=C)_{Ar}, 1196, 1160 (C-O); ¹H-NMR (400 MHz, DMSO-*d*₆) δ ppm: 10.39 (s, 3H, C-H_{BImidazole}, major),

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10.32 (s, 3H, C-H_{BImidazole}, minor), 10.26 (s, 3H, C-H_{BImidazole}, minor), 8.15-8.09 (m, 6H, CH_{Ar}), 7.70-7.61 (m, 6H, CH_{Ar}, major), 7.31-7.24 (m, 6H, CH_{Ar}, minor), 5.85 (s, 6H, O-CH₂, major), 5.80 (s, 6H, O-CH₂, minor), 4.57 (t, *J*=7.07, 6H, α -CH₂, major), 4.29 (t, *J*=7.07, 6H, α -CH₂, minor), 4.03 (s, 6H, N-CH₂, minor), 3.99 (s, 6H, N-CH₂, major), 1.92-1.84 (m, 6H, β -CH₂), 1.26 (bs, 18H, bulk-CH₂), 0.86 (s, 3H, CH₃), 0.82 (t, 9H, overlap, ω -CH₃); ¹³C-NMR (100 MHz, DMSO-d₆) δ ppm: 167.13 (C=O, minor), 166.41 (C=O, major), 143.37 (CH_{BImidazole}), 131.50 (C_{Ar}), 130.61 (C_{Ar}), 126.74 (CH_{Ar}), 126.66 (CH_{Ar}), 114.10 (CH_{Ar}), 113.75 (CH_{Ar}), 66.29 (CH₂-O), 47.55 (CH₂-N), 46.80 (α -CH₂), 38.13 (-C-), 30.56 (ω -2), 28.49 (bulk-CH₂), 25.64 (β , minor), 25.32 (β , major), 21.87 (ω -1), 16.16 (CH₃, major), 16.09 (CH₃, minor), 13.79 (ω); HRMS: m/z, [M⁺³-2H]-3Cl⁻ calcd. for C₅₀H₆₇N₆O₆⁵⁺: 847.5122, found: 847.5162.

Tris-((N-octyl-benzimidazoliumyl-acetayloxy)methyl)ethane chloride (9d)

This compound was prepared analogously to **9b** using tris-((2-chloro-acetayloxy)methyl)ethane (compound **3**) (1.9 g, 5.43 mmol) and 1-octyl-benzimidazole (**7d**) (3.75 g, 16.3 mmol) to provide a viscous hygroscopic syrup in 98% yield (5.54g). Molecular Formula: C₅₆H₈₁Cl₃N₆O₆; Mol. Wt.: 1040.64; FTIR (cm⁻¹): 3134 (C-H)_{Ar}, 2955, 2925, 2855 (C-H)_{Aliph}, 1749 (C=O), 1618 (C=N), 1562, 1486 1462 (C=C)_{Ar}, 1199 (C-O); ¹H-NMR (400 MHz, DMSO-d₆) δ ppm: 10.39 (s, 3H, C-H_{BImidazole}, major), 10.32 (s, 3H, C-H_{BImidazole}, minor), 10.26 (s, 3H, C-H_{BImidazole}, minor), 8.16-8.10 (m, 6H, CH_{Ar}), 7.70-7.61 (m, 6H, CH_{Ar}, major), 7.29-7.19 (m, 6H, CH_{Ar}, minor), 5.85 (s, 6H, O-CH₂, major), 5.80 (s, 6H, O-CH₂, minor), 4.57 (t, *J*=7.07, 6H, α -CH₂, major), 4.25 (t, *J*=7.07, 6H, α -CH₂, minor), 4.04 (s, 6H, N-CH₂, minor), 3.99 (s, 6H, N-CH₂, major), 1.92- 1.85 (m, 6H, β -CH₂), 1.20 (bs, 30H, bulk-CH₂), 0.86 (s, 3H, CH₃), 0.82 (t, 9H, *J*=6.80, ω -CH₃); ¹³C-NMR (100 MHz, DMSO-d₆) δ ppm: 167.83 (C=O, major), 166.51 (C=O, minor), 143.22 (CH_{BImidazole}), 131.56 (C_{Ar}), 130.62 (C_{Ar}), 126.79 (CH_{Ar}), 126.57 (CH_{Ar}), 113.86 (CH_{Ar}), 113.71 (CH_{Ar}), 64.02 (CH₂-O, major), 63.13 (CH₂-O, minor), 47.54 (CH₂-N), 46.77 (α -CH₂), 40.58 (-C-), 31.08 (ω -2), 28.42 (2), 28.31 (bulk-CH₂), 25.61 (β), 21.98 (ω -1), 16.55 (CH₃, major), 16.32 (CH₃, minor), 13.88 (ω); HRMS: m/z, [M⁺³-2H]-3Cl⁻ calcd. for C₅₆H₇₉N₆O₆⁵⁺: 931.6061, found: 931.6144.

Tris-((N-benzyl-benzimidazoliumyl-acetayloxy)methyl)ethane chloride (9g)

This compound was prepared analogously to **9b** using tris-((2-chloro-acetayloxy)methyl)ethane (compound **3**) (1.9 g, 5.43 mmol) and 1-benzyl-benzimidazole (**7g**) (3.39 g, 16.3 mmol) to provide a pale yellow hygroscopic semi-solid in 96% yield (5.08g). Molecular Formula: C₅₃H₅₁Cl₃N₆O₆; Mol. Wt.: 974.37 ; FTIR (cm⁻¹): 3120 (C-H)_{Ar}, 2974 (C-H)_{Aliph}, 1750 (C=O), 1615 (C=N), 1562, 1486, 1455 (C=C)_{Ar}, 1190, 1165 (C-O); ¹H-NMR (400 MHz, DMSO-d₆) δ ppm: 10.30 (s, 3H, C-H_{BImidazole}, major), 10.23 (s, 3H, C-H_{BImidazole}, minor), 10.17 (s, 3H, C-H_{BImidazole}, minor), 8.05 (d, *J*=8.31 Hz, 3H, C-H_{Ar}), 7.99 (d, *J*=8.31 Hz, 3H, C-H_{Ar}), 7.63-7.51 (m, 12H, C-H_{Ar}), 7.41-7.31 (m, 9H, C-H_{Ar}), 5.89 (s, 6H, CH₂-Ar), 5.78 (s, 6H, O-CH₂, major), 5.74 (s, 6H, O-CH₂, minor), 5.67 (s, 6H, O-CH₂, minor), 4.03 (s, 6H, N-CH₂, major), 3.97 (s, 6H, N-CH₂, minor), 0.88 (s, 3H, CH₃, minor), 0.85 (s, 3H, minor), 0.81 (s, 3H, CH₃, major); ¹³C-NMR (100 MHz, DMSO-d₆) δ ppm: 167.19 (C=O, minor), 166.45 (C=O, major), 143.97 (CH_{BImidazole}, minor), 143.66 (CH_{BImidazole}, major), 133.93 (-C_{Ar}-CH₂-), 131.65 (C_{BImidazole}), 130.34 (C_{BImidazole}), 129.00 (2 \times CH_{Ar}), 128.77 (CH_{Ar}, major), 128.70 (CH_{Ar}, minor), 128.35 (2 \times CH_{Ar}), 126.90 (CH_{BImidazole}), 126.84 (CH_{BImidazole}), 114.20 (CH_{BImidazole}), 113.96 (CH_{BImidazole}), 66.36 (CH₂-O, major),

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66.19 (CH₂-O, minor), 49.93 (CH₂-N), 47.84 (Ar-CH₂-, minor), 47.20 (Ar-CH₂-, major), 38.30 (-C-), 16.21 (CH₃, minor), 16.13 (CH₃, major); HRMS: m/z, [M⁺³-2H]-3Cl⁻ calcd. for C₅₃H₄₉N₆O₆⁵⁺: 865.3714, found: 865.3746.

Tris-((N-butyl-imidazoliumyl-acetayloxy)methyl)ethane bis(trifluoromethylsulfonyl)amide (10b)

This compound was prepared analogously to **10a** using tris-((N-butyl-imidazoliumyl-acetayloxy)methyl)-ethane chloride **8b** (0.72 g, 1.0 mmole) and Lithium bis-(trifluoromethanesulphonyl)imide LiNTf₂ (1.0 g, 3.5 mmol) to provide a clear viscous hygroscopic liquid at room temperature in 84% yield (1.22 g). Molecular Formula: C₃₈H₅₁F₁₈N₉O₁₈S₆; Mol. Wt.: 1456.22; FTIR (cm⁻¹): 3055 (C-H)_{Ar}, 2960, 2935, 2872 (C-H)_{Aliph}, 1743 (C=O), 1642 (C=N), 1560, 1466 (C=C)_{Ar}, 1351, 1223 (C-F), 1360, 1152 (O=S=O), 1195, 1163 (C-O); ¹H-NMR (400 MHz, CD₃OD) δ ppm: 8.97 (s, 3H, C-H_{Imidazole}, minor), 8.94 (s, 3H, C-H_{Imidazole}, major), 8.80 (s, 3H, C-H_{Imidazole}, minor), 7.70 (s, 3H, C-H_{Imidazole}), 7.60 (s, 3H, C-H_{Imidazole}, major), 7.50 (s, 3H, C-H_{Imidazole}, minor), 5.20 (s, 6H, O-CH₂, major), 5.19 (s, 6H, O-CH₂, minor), 4.26 (t, J= 7.28 Hz, 6H, α-CH₂), 4.19 (s, 6H, N-CH₂, major), 4.16 (s, 6H, N-CH₂, minor), 4.15 (s, 6H, N-CH₂, minor), 1.80-1.73 (m, 6H, β-CH₂, major), 1.92-1.83 (m, 6H, β-CH₂), 1.42-1.33 (m, 6H, (ω-1), 1.08 (s, 3H, CH₃), 0.98 (t, J= 7.28 Hz, 9H, ω-CH₃); ¹³C-NMR (100 MHz, CD₃OD) δ ppm: 167.90 (C=O, minor), 167.77 (C=O, major), 138.67 (CH_{Imidazole}), 126.10, 122.91, 119.72, 116.54 (q, J=320, CF₃), 125.37 (CH_{Imidazole}), 123.69 (CH_{Imidazole}), 68.83 (CH₂-O, minor), 68.30 (CH₂-O, minor), 68.10 (CH₂-O, major), 51.03 (CH₂-N), 50.87 (α-CH₂), 40.10 (-C-), 33.39 ((ω-2), minor), 33.15 ((ω-2), major), 20.61 ((ω-1), minor), 20.52 ((ω-1), major), 17.04 (CH₃, major), 17.00 (CH₃, minor), 13.83 (ω). ¹⁹F (336, MHz) δ ppm: -80.02; HRMS: m/z, [M⁺³-2H]-3NTF₂ calcd. for C₃₂H₄₉N₆O₆⁵⁺: 613.3714, found: 613.3692; m/z, [NTF₂] calcd. for C₂F₆NO₄S₂⁻: 279.9173, found: 279.9175.

Tris-((N-hexyl-imidazoliumyl-acetayloxy)methyl)ethane bis(trifluoromethylsulfonyl)amide (10c)

This compound was prepared analogously to **10a** using tris-((N-hexyl-imidazoliumyl-acetayloxy)methyl)-ethane chloride **8c** (0.81 g, 1.0 mmole) and Lithium bis-(trifluoromethanesulphonyl)imide LiNTf₂ (1.0 g, 3.5 mmol) to provide a clear viscous hygroscopic liquid at room temperature in 85% yield (1.32 g). Molecular Formula: C₄₄H₆₃F₁₈N₉O₁₈S₆; Mol. Wt.: 1540.38; FTIR (cm⁻¹): 3062 (C-H)_{Ar}, 2950, 2931, 2845 (C-H)_{Aliph}, 1749 (C=O), 1644 (C=N), 1564, 1460(C=C)_{Ar}, 1347, 1220 (C-F), 1366, 1153 (O=S=O), 1185, 1185 (C-O); ¹H-NMR (400 MHz, CD₃OD) δ ppm: 8.98 (bt~s, 3H, C-H_{Imidazole}, minor), 8.96 (bt~s, 3H, C-H_{Imidazole}, minor), 8.93 (bt~s, 3H, C-H_{Imidazole}, major), 7.69 (t, J= 1.81 Hz, 3H, C-H_{Imidazole}, major), 7.63 (t, J = 1.81 Hz, 3H, C-H_{Imidazole}, minor), 7.61 (t, 3H, J= 1.81 Hz, C-H_{Imidazole}, minor), 7.59 (t, J= 1.81 Hz, 3H, C-H_{Imidazole}, major), 5.19 (s, 6H, O-CH₂), 4.24 (t, J= 7.25 Hz, 6H, α-CH₂), 4.19 (s, 6H, N-CH₂, major), 4.14 (s, 6H, N-CH₂, minor), 1.93-1.86 (m, 6H, β-CH₂), 1.34 (bs, 18H, bulk-CH₂), 1.08 (s, 3H, CH₃), 0.90 (t, J = 6.34 Hz, 9H, ω-CH₃); ¹³C-NMR (100 MHz, CD₃OD) δ ppm: 167.77 (C=O), 138.65 (CH_{Imidazole}), 126.15, 122.93, 119.71, 116.49 (q, J=320, CF₃), 125.37 (CH_{Imidazole}), 123.69 (CH_{Imidazole}), 68.23 (CH₂-O, minor), 68.15 (CH₂-O, major), 51.31 (CH₂-N), 50.89 (α-CH₂), 40.11(-C-), 32.35 (ω-2), 31.14 (bulk-CH₂), 26.99 (β-CH₂), 23.58 (ω-1), 17.11 (CH₃, minor), 17.07 (CH₃, major), 14.38 (ω). ¹⁹F (336, MHz) δ ppm: -80.22;

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HRMS: m/z, [M⁺³-2H]-3NTF₂ calcd. for C₃₈H₆₁N₆O₆⁵⁺: 697.4653, found: 697.4623; m/z, [NTF₂] calcd. for C₂F₆NO₄S₂⁻: 279.9173, found: 279.9177.

Tris-((N-decyl-imidazoliumyl-acetayloxy)methyl)ethane bis(trifluoromethylsulfonyl)amide (10e)

This compound was prepared analogously to **10a** using tris-((N-decyl-imidazoliumyl-acetayloxy)methyl)-ethane chloride **8e** (0.97 g, 1.0 mmole) and Lithium bis-(trifluoromethanesulphonyl)imide LiNTf₂ (1.0 g, 3.5 mmol) to provide a clear viscous hygroscopic liquid at room temperature in 94% yield (1.6 g). Molecular Formula: C₅₆H₈₇F₁₈N₉O₁₈S₆; Mol. Wt.: 1708.70; FTIR (cm⁻¹): 3058 (C-H)_{Ar}, 2955, 2929, 2857 (C-H)_{Aliph}, 1749 (C=O), 1666 (C=N), 1564, 1462 (C=C)_{Ar}, 1340, 1210 (C-F), 1360, 1123 (O=S=O), 1199, 1165 (C-O); ¹H-NMR (400 MHz, CD₃OD) δ ppm: 9.00 (bt~s, 3H, C-H_{Imidazole}, minor), 8.98 (bt~s, 3H, C-H_{Imidazole}, minor), 8.95 (bt~s, 3H, C-H_{Imidazole}, major), 7.70 (t, J=1.81, 3H, C-H_{Imidazole}, major), 7.65 (t, 3H, C-H_{Imidazole}, minor), 7.61 (t, J=1.81, 3H, C-H_{Imidazole}, major), 7.56 (t, 3H, J=1.81, C-H_{Imidazole}, minor), 5.31 (s, 6H, O-CH₂, minor), 5.21 (s, 6H, O-CH₂, major), 4.25 (t, J= 7.25 Hz, 6H, α-CH₂), 4.21 (s, 6H, N-CH₂), 4.19 (s, 6H, N-CH₂), 1.94-1.87 (m, 6H, β-CH₂), 1.33 (bs, 42H, bulk-CH₂), 1.09 (s, 3H, CH₃, major), 1.07 (s, 3H, CH₃, minor), 0.90 (t, 9H, J =6.80 Hz, ω-CH₃); ¹³C-NMR (100 MHz, CD₃OD) δ ppm: 167.80 (C=O), 138.67 (CH_{Imidazole}), 126.17, 122.94, 119.71, 116.49 (q, J=321, CF₃), 125.38 (CH_{Imidazole}), 123.70 (CH_{Imidazole}), 68.17 (CH₂-O), 51.32 (CH₂-N), 50.88 (α-CH₂), 40.12 (-C-), 33.00 (ω-2), 31.20, 30.30 (2), 30.15 (2) (bulk-CH₂), 27.33 (β), 23.80 (ω-1), 17.07 (CH₃), 14.53 (ω). ¹⁹F (336, MHz) δ ppm: -80.52; HRMS: m/z, [M⁺³-2H]-3NTF₂ calcd. for C₅₀H₈₅N₆O₆⁵⁺ 865.6531 found: 865.6512; m/z, [NTF₂] calcd. for C₂F₆NO₄S₂⁻: 279.9173, found: 279.9148.

Tris-((N-butyl-benzimidazoliumyl-acetayloxy)methyl)ethane bis(trifluoromethylsulfonyl)-amide (11b)

This compound was prepared analogously to **10a** using tris-((N-butyl-benzimidazoliumyl-acetayloxy)-methyl) ethane chloride **9b** (0.87 g, 1.0 mmole) and Lithium bis-(trifluoromethanesulphonyl)imide LiNTf₂ (1.0 g, 3.5 mmol) to provide a clear viscous hygroscopic liquid at room temperature in 84% yield (1.35 g). Molecular Formula: C₅₀H₅₇F₁₈N₉O₁₈S₆; Mol. Wt.: 1606.39; FTIR (cm⁻¹): 3022 (C-H)_{Ar}, 2950, 2942, 2863 (C-H)_{Aliph}, 1755(C=O), 1618(C=N), 1564, 1485, 1469 (C=C)_{Ar}, 1374, 1233 (C-F), 1354, 1169 (O=S=O), 1187, 1160 (C-O); ¹H-NMR (400 MHz, DMSO-d₆) δ ppm: 9.94 (s, 3H, C-H_{BImidazole}, minor), 9.89 (s, 3H, C-H_{BImidazole}, minor), 9.82 (s, 3H, C-H_{BImidazole}), 8.10-8.04 (m, 6H, CH_{Ar}), 7.68-7.59 (m, 6H, CH_{Ar}, major), 7.27-7.22 (m, 6H, CH_{Ar}, minor), 5.65 (s, 6H, O-CH₂, minor), 5.60 (s, 6H, O-CH₂, major), 4.47 (t, J=7.07, 6H, α-CH₂, major), 4.27 (t, J=7.07, 6H, α-CH₂, minor), 4.01 (s, 6H, N-CH₂, major), 3.98 (s, 6H, N-CH₂, minor), 1.87-1.79 (m, 6H, β-CH₂, major), 1.66-1.59 (m, 6H, β-CH₂, minor), 1.28-1.18 (m, 6H, (ω-1)), 0.93 (s, 3H, CH₃), 0.87 (t, 9H, J =7.21 Hz, ω-CH₃); ¹³C-NMR (100 MHz, DMSO-d₆) δ ppm: 166.76 (C=O), 144.03 (CH_{BImidazole}, major), 143.867 (CH_{BImidazole}, minor), 132.45 (C_{Ar}), 128.90 (C_{Ar}), 125.30, 122.09, 118.88, 115.67 (q, J=321, CF₃), 125.02 (CH_{Ar}), 124.14 (CH_{Ar}), 112.85 (CH_{Ar}), 110.17 (CH_{Ar}), 68.03 (CH₂-O), 46.55 (CH₂-N), 44.03 (α-CH₂, minor), 43.78 (α-CH₂, major), 38.31 (-C-), 32.17 ((ω-2), minor), 31.90 ((ω-2), major), 18.55 ((ω-1), major), 17.95 ((ω-1), minor), 15.74 (CH₃), 13.23 ((ω), minor), 13.12 ((ω), major). ¹⁹F (336, MHz) δ ppm: -80.12; HRMS: m/z, [M⁺³-2H]-3NTF₂ calcd. for

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$C_{44}H_{55}N_6O_6^{5+}$: 763.4183, found: 763.4202; m/z, [NTF₂] calcd. for C₂F₆NO₄S₂⁻: 279.9173, found: 279.9209.

Tris-((N-decyl-benzimidazoliumyl-acetayloxy)methyl)ethane bis(trifluoromethylsulfonyl)-amide (11e)

This compound was prepared analogously to **10a** using tris-((N-decyl-benzimidazoliumyl-acetayloxy)-methyl) ethane chloride **9e** (1.12 g, 1.0 mmole) and Lithium bis-(trifluoromethanesulphonyl)imide LiNTf₂ (1.0 g, 3.5 mmol) to provide a clear viscous hygroscopic liquid at room temperature in 90% yield (1.67 g). Molecular Formula: C₆₈H₉₃F₁₈N₉O₁₈S₆; Mol. Wt.: 1858.88; FTIR (cm⁻¹): 3130 (C-H)_{Ar}, 2949, 2924, 2848 (C-H)_{Aliph}, 1749 (C=O), 1620 (C=N), 1566 1485, 1457 (C=C)_{Ar}, 1367, 1223 (C-F), 1360, 1165 (O=S=O), 1218, 1199 (C-O); ¹H-NMR (400 MHz, CD₃OD) δ ppm: 9.54 (s, 3H, C-H_{BImidazole}, minor), 9.48 (s, 3H, C-H_{BImidazole}, minor), 9.42 (s, 3H, C-H_{BImidazole}, major), 8.02-7.82 (m, 6H, CH_{Ar}), 7.73-7.55 (m, 6H, CH_{Ar}), 5.51 (s, 6H, O-CH₂, minor), 5.50 (s, 6H, O-CH₂, minor), 5.48 (s, 6H, O-CH₂, major), 4.57 (t, J=7.20, 6H, α-CH₂, major), 4.50 (t, J=7.20, 6H, α-CH₂, minor), 4.21 (s, 6H, N-CH₂, major), 4.19 (s, 6H, N-CH₂, minor), 4.17 (s, 6H, N-CH₂, minor), 2.04- 1.94 (m, 6H, β-CH₂), 1.27 (bs, 42H, bulk-CH₂), 1.04 (s, 3H, CH₃, major), 1.02 (s, 3H, CH₃, minor), 0.88 (t, 9H, J=6.78, ω-CH₃); ¹³C-NMR (100 MHz, CD₃OD) δ ppm: 167.66 (C=O, major), 167.57 (C=O, minor), 144.05 (CH_{BImidazole}), 133.31 (C_{Ar}), 132.57 (C_{Ar}), 128.78 (CH_{Ar}), 128.60 (CH_{Ar}), 126.17, 122.95, 119.73, 116.51 (q, J=320, CF₃), 114.79 (2) (CH_{Ar}), 68.07 (CH₂-O, minor), 67.85 (CH₂-O, major), 46.69 (CH₂-N), 41.89 (α-CH₂, minor), 41.70 (α-CH₂, major), 40.50 (-C-), 33.17 (ω-2), 30.72, 30.65, 30.54, 30.25 (2) (bulk-CH₂), 27.58 (β), 23.85 (ω-1), 17.07 (CH₃, minor), 16.97 (CH₃, major), 14.58 (ω). ¹⁹F (336, MHz) δ ppm: -80.06; HRMS: m/z, [M⁺³-2H]-3NTF₂ calcd. for C₆₂H₉₁N₆O₆⁵⁺ : 1015.7000, found: 1015.6977; m/z, [NTF₂] calcd. for C₂F₆NO₄S₂⁻: 279.9173, found: 279.9175.

Tris-((N-dodecyl-benzimidazoliumyl-acetayloxy)methyl)ethane bis(trifluoromethylsulfonyl)-amide (11f)

This compound was prepared analogously to **10a** using tris-((N-dodecyl-benzimidazoliumyl-acetayloxy)-methyl) ethane chloride **9f** (1.21 g, 1.0 mmole) and Lithium bis-(trifluoromethanesulphonyl)imide LiNTf₂ (1.0 g, 3.5 mmol) to provide a clear viscous hygroscopic liquid at room temperature in 96% yield (1.87 g). Molecular Formula: C₇₄H₁₀₅F₁₈N₉O₁₈S₆; Mol. Wt.: 1943.04 ; FTIR (cm⁻¹): 3127 (C-H)_{Ar}, 2965 (C-H)_{Aliph}, 1750 (C=O), 1622 (C=N), 1565, 1485, 1448 (C=C)_{Ar}, 1362, 1222 (C-F), 1358, 1167 (O=S=O), 1210, 1170 (C-O); ¹H-NMR (400 MHz, CD₃OD) δ ppm: 9.72 (s, 3H, C-H_{BImidazole}, minor), 9.69 (s, 3H, C-H_{BImidazole}, minor), 9.65 (s, 3H, C-H_{BImidazole}, major), 8.00-7.91 (m, 6H, C-H_{BImidazole}), 7.71-7.54 (m, 6H, C-H_{Ar}, major), 7.40-7.31 (m, 6H, C-H_{Ar}, minor), 5.60 (s, 6H, O-CH₂, major), 5.57 (s, 6H, O-CH₂, minor), 4.53 (t, J=7.07, 6H, α-CH₂, major), 4.34 (t, J=7.07, 6H, α-CH₂, minor), 4.19 (s, 6H, N-CH₂, major), 4.16 (s, 6H, N-CH₂, minor), 1.98-1.92 (m, 6H, β-CH₂), 1.24 (bs, 54H, bulk-CH₂), 1.03 (s, 3H, CH₃, minor), 1.00 (s, 3H, CH₃, minor), 0.97 (s, 3H, CH₃, major), 0.86 (t, J=7.25, 9H, ω-CH₃); ¹³C-NMR (100 MHz, CD₃OD) δ ppm: 166.89 (C=O), 143.75 (CH_{BImidazole}, major), 143.51(CH_{BImidazole}, minor), 133.10 (C_{BImidazole}), 130.92 (C_{BImidazole}), 128.08 (CH_{Ar}), 127.12 (CH_{Ar}), 126.36, 123.14, 119.92, 116.71 (q, J=320, CF₃), 114.81 (CH_{Ar}), 114.76 (CH_{Ar}), 68.37 (CH₂-O, minor), 68.24 (CH₂-O, major), 45.70 (CH₂-N), 40.18 (α-CH₂), 39.34 (-C-), 31.65 (ω-2), 29.66 (2), 29.32, 29.22, 29.03, 28.93, 28.86 (bulk-CH₂), 25.82 (β),

Supplementary Information: Tris-imidazolium and benzimidazolium ionic liquids: A new class of biodegradable surfactants

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22.53 (ω -1), 16.23 (CH₃, minor), 16.03 (CH₃, major), 14.19 (ω). ¹⁹F (336, MHz) δ ppm: -80.20; HRMS: m/z, [M^{+3-2H}]–3NTF₂ calcd. for C₆₈H₁₀₃N₆O₆⁵⁺: 1099.7939, found: 1099.7977; m/z, [NTF₂] calcd. for C₂F₆NO₄S₂⁻: 279.9173, found: 279.9225.

Tris-((N-benzyl-benzimidazoliumyl-acetayloxy)methyl)ethane bis(trifluoromethylsulfonyl)-amide (11g)

This compound was prepared analogously to **10a** using tris-(*N*-benzyl-benzimidazoliumyl-acetayloxy)-methyl)ethane chloride **9g** (0.97 g, 1.0 mmole) and Lithium bis-(trifluoromethanesulphonyl)imide LiNTf₂ (1.0 g, 3.5 mmol) to provide a clear viscous hygroscopic liquid at room temperature in 96% yield (1.63 g). Molecular Formula: C₅₉H₅₁F₁₈N₉O₁₈S₆; Mol. Wt.: 1708.45; FTIR (cm⁻¹): 3127 (C-H)_{Ar}, 2965 (C-H)_{Aliph}, 1750 (C=O), 1622 (C=N), 1565, 1485, 1448 (C=C)_{Ar}, 1362, 1222 (C-F), 1358, 1167 (O=S=O), 1210, 1170 (C-O); ¹H-NMR (400 MHz, DMSO-*d*₆) δ ppm: 10.44 (s, 3H, C-H_{BImidazole}, major), 10.30 (s, 3H, C-H_{BImidazole}, minor), 10.13 (s, 3H, C-H_{BImidazole}, minor), 8.12-8.00 (m, 6H, C-H_{Ar}), 7.61-7.53 (m, 12H, C-H_{Ar}), 7.39-7.22 (m, 9H, C-H_{Ar}), 5.92 (s, 6H, CH₂-Ar), 5.84 (s, 6H, O-CH₂, major), 5.79 (s, 6H, O-CH₂, minor), 4.04 (s, 6H, N-CH₂, major), 3.99 (s, 6H, N-CH₂, minor), 0.91 (s, 3H, CH₃, minor), 0.86 (s, 3H, CH₃, minor), 0.82 (s, 3H, CH₃, major); ¹³C-NMR (100 MHz, DMSO-*d*₆) δ ppm: 168.01 (C=O, major), 166.57 (C=O, minor), 143.48 (CH_{BImidazole}, major), 142.82 (CH_{BImidazole}, minor), 133.82 (-C_{Ar}-CH₂-), 131.79 (C_{BImidazole}), 130.48 (C_{BImidazole}), 129.10 (2×CH_{Ar}), 128.87 (CH_{Ar}), 128.31 (2×CH_{Ar}), 126.98 (CH_{BImidazole}), 126.81 (CH_{BImidazole}), 124.37, 121.13, 117.90, 114.66 (q, *J*=322, CF₃), 114.06 (CH_{BImidazole}), 113.89 (CH_{BImidazole}), 63.33 (CH₂-O, major), 62.67 (CH₂-O, minor), 50.00 (CH₂-N), 47.72 (Ar-CH₂-, major), 47.61 (Ar-CH₂-, minor), 40.63 (-C-), 16.33 (CH₃, major), 16.08 (CH₃, minor). ¹⁹F (336, MHz) δ ppm: -79.95; HRMS: m/z, [M^{+3-2H}]–3NTF₂ calcd. for C₅₃H₄₉N₆O₆⁵⁺: 865.3714, found: 865.3750; m/z, [NTF₂] calcd. for C₂F₆NO₄S₂⁻: 279.9173, found: 279.9214.