

## Supporting Information

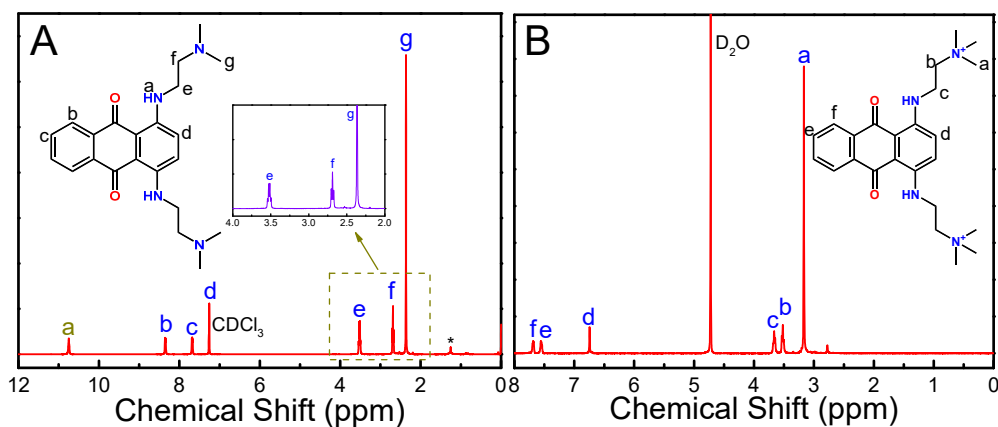
A quaternized anthraquinone derivative for pH-neutral aqueous organic redox flow batteries

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**Fig. S1** (A)  $^1\text{H}$  NMR spectrum of BDEAQ. (B)  $^1\text{H}$  NMR spectrum of BDEAQI<sub>2</sub>.

## Single Mass Analysis

Tolerance = 20.0 PPM / DBE: min = -1.5, max = 50.0

Element prediction: Off

Number of isotope peaks used for i-FIT = 3

Monoisotopic Mass, Even Electron Ions

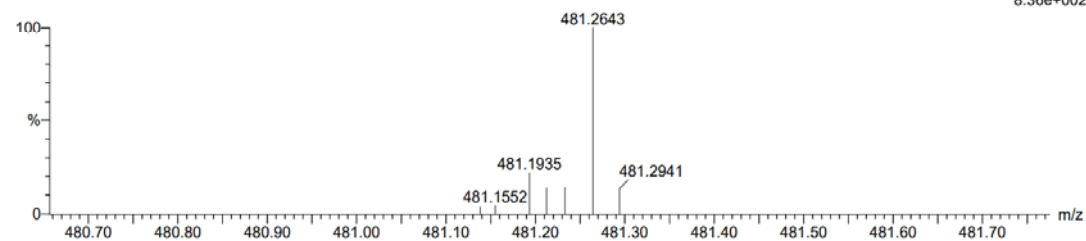
1558 formula(e) evaluated with 1 results within limits (up to 50 closest results for each mass)

Elements Used:

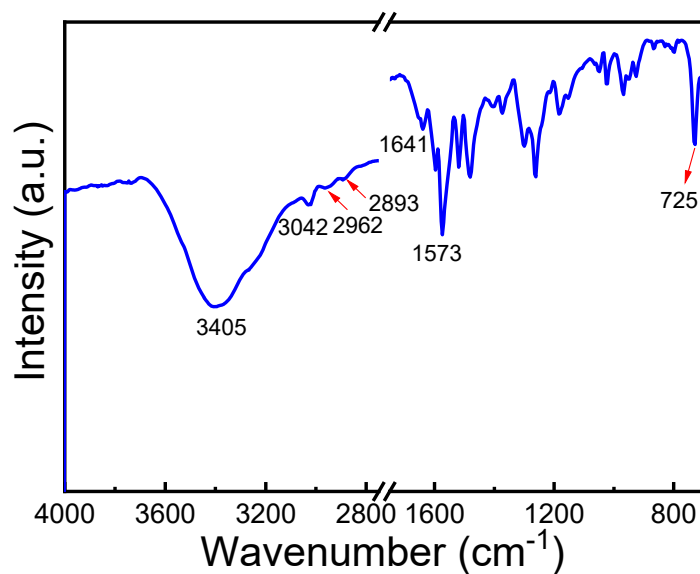
C: 24-24 H: 35-35 N: 0-10 O: 0-41 Na: 0-3 Cl: 1-2

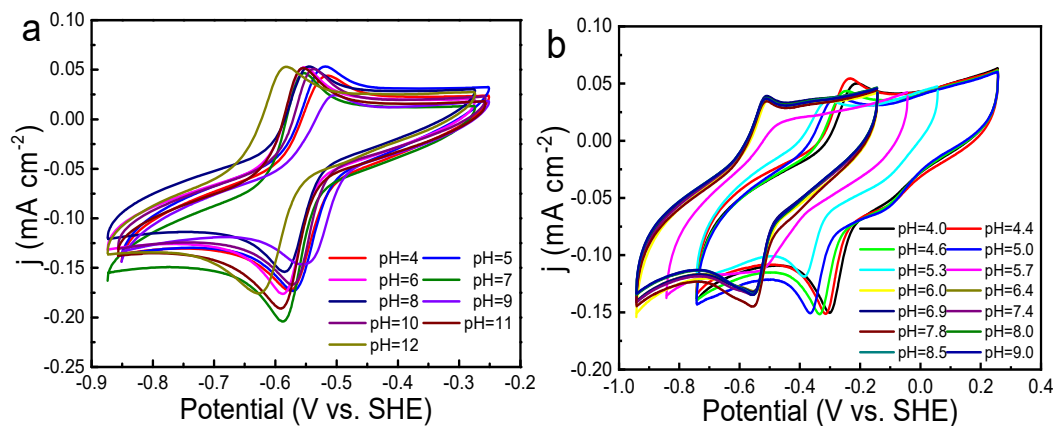
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230216-7-17 (0.093)

1: TOF MS ES+  
8.36e+002Minimum: -1.5  
Maximum: 5.0 20.0 50.0

Mass	Calc. Mass	mDa	PPM	DBE	i-FIT	Norm	Conf (%)	Formula
481.2126	481.2137	-1.1	-2.3	8.5	51.8	n/a	n/a	C24 H35 N4 O2 Cl2

Fig. S2 MS spectrum of BDEAQCl<sub>2</sub> in the form of positive ions.Fig. S3 FTIR spectrum of BDEAQCl<sub>2</sub>.



**Fig. S4** CVs of 1 mM BDEAQC1<sub>2</sub> on a glassy carbon electrode in unbuffered (a) and buffered (b) 1 M NaCl solutions with different pH. The pH buffer is a 0.1 M acetic acid/sodium acetate solution containing 1 M NaCl.