

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM022322\
 Data File : BM034050.D
 Acq On : 23 Feb 2022 13:34
 Operator : CG/JU
 Sample : N1627-02
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 20220096-BERKELEY-8-ELUTRIATE

Quant Time: Feb 23 14:14:44 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM022222.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 22 17:55:56 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.019	152	217274	20.000	ng	0.00
21) Naphthalene-d8	10.836	136	787331	20.000	ng	# 0.00
39) Acenaphthene-d10	14.654	164	500935	20.000	ng	0.00
64) Phenanthrene-d10	17.395	188	1045535	20.000	ng	# 0.00
76) Chrysene-d12	21.565	240	1101234	20.000	ng	0.00
86) Perylene-d12	24.024	264	1169031	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.566	112	619017	50.787	ng	0.00
7) Phenol-d6	7.166	99	548527	34.272	ng	0.00
23) Nitrobenzene-d5	9.178	82	932387	63.908	ng	0.00
42) 2,4,6-Tribromophenol	16.136	330	619648	101.920	ng	0.00
45) 2-Fluorobiphenyl	13.283	172	2341277	62.269	ng	0.00
79) Terphenyl-d14	20.012	244	3169309	50.115	ng	0.00
Target Compounds						
10) Phenol	7.195	94	52470	3.333	ng	Qvalue 92
74) Di-n-butylphthalate	18.365	149	248142	4.105	ng	# 94

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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