

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM022322\
 Data File : BM034054.D
 Acq On : 23 Feb 2022 16:01
 Operator : CG/JU
 Sample : N1632-02
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 20220092B-BERKELEY-4B-ELUTRIATE

Quant Time: Feb 23 16:58:01 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	212696	20.000	ng	0.00
21) Naphthalene-d8	10.836	136	794946	20.000	ng	# 0.00
39) Acenaphthene-d10	14.654	164	513124	20.000	ng	0.00
64) Phenanthrene-d10	17.395	188	1091380	20.000	ng	# 0.00
76) Chrysene-d12	21.565	240	1191724	20.000	ng	0.00
86) Perylene-d12	24.024	264	1278724	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.566	112	793212	66.480	ng	0.00
7) Phenol-d6	7.166	99	679001	43.337	ng	0.00
23) Nitrobenzene-d5	9.178	82	1343176	91.182	ng	0.00
42) 2,4,6-Tribromophenol	16.136	330	1023012	164.269	ng	0.00
45) 2-Fluorobiphenyl	13.283	172	3509609	91.125	ng	0.00
79) Terphenyl-d14	20.012	244	4987671	72.879	ng	0.00
Target Compounds						
10) Phenol	7.190	94	54803	3.556	ng	Qvalue 94

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

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