

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG022822\
 Data File : BG052508.D
 Acq On : 28 Feb 2022 19:22
 Operator : CG/JU
 Sample : N1654-01
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 20220091-BERKELEY-3-ELUTRIATE

Quant Time: Mar 01 01:00:18 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG020722.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 08 01:21:20 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.216	152	24671	20.000	ng	0.00
21) Naphthalene-d8	11.054	136	96864	20.000	ng	#-0.01
39) Acenaphthene-d10	14.866	164	69029	20.000	ng	0.00
64) Phenanthrene-d10	17.616	188	172251	20.000	ng	0.00
76) Chrysene-d12	21.927	240	178148	20.000	ng	#-0.01
86) Perylene-d12	25.387	264	175987	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.749	112	133262	94.142	ng	0.00
7) Phenol-d6	7.365	99	134559	61.608	ng	0.00
23) Nitrobenzene-d5	9.391	82	240546	107.929	ng	0.00
42) 2,4,6-Tribromophenol	16.353	330	172304	173.744	ng	0.00
45) 2-Fluorobiphenyl	13.486	172	558728	112.470	ng	0.00
79) Terphenyl-d14	20.206	244	896045	88.819	ng	0.00
Target Compounds						
10) Phenol	7.394	94	7887	3.634	ng	Qvalue 87

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

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