

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110322\
 Data File : BG055448.D
 Acq On : 3 Nov 2022 22:14
 Operator : CG/JU
 Sample : N5306-07
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 DISSOLVED-20221167-COMP-02-MODIFIED-ELUTRIATE

Quant Time: Nov 04 05:32:10 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG110122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 04 05:28:43 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.236	152	26716	20.000	ng	0.00
21) Naphthalene-d8	11.062	136	112568	20.000	ng	0.00
39) Acenaphthene-d10	14.851	164	82525	20.000	ng	0.00
64) Phenanthrene-d10	17.583	188	199435	20.000	ng	0.00
76) Chrysene-d12	21.855	240	203539	20.000	ng	0.00
86) Perylene-d12	25.227	264	220349	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.762	112	116311	78.184	ng	0.00
7) Phenol-d6	7.384	99	136848	66.334	ng	0.00
23) Nitrobenzene-d5	9.411	82	149441	70.752	ng	0.00
42) 2,4,6-Tribromophenol	16.332	330	143431	127.419	ng	0.00
45) 2-Fluorobiphenyl	13.476	172	403632	72.517	ng	0.00
79) Terphenyl-d14	20.157	244	769159	73.483	ng	0.00
Target Compounds						
10) Phenol	7.413	94	7766	3.326	ng	Qvalue 97

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

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