

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111522\
 Data File : BG055631.D
 Acq On : 15 Nov 2022 21:09
 Operator : CG/JU
 Sample : N5431-07
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 DISSOLVED-20221177-COMP-12-MODIFIED-ELUTRIATE

Quant Time: Nov 16 01:45:47 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG110922.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 16 01:36:26 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.256	152	42864	20.000	ng	0.00
21) Naphthalene-d8	11.082	136	174569	20.000	ng	0.00
39) Acenaphthene-d10	14.878	164	127559	20.000	ng	0.00
64) Phenanthrene-d10	17.615	188	327247	20.000	ng	0.00
76) Chrysene-d12	21.916	240	321823	20.000	ng	0.00
86) Perylene-d12	25.330	264	338133	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.782	112	158565	70.461	ng	0.00
7) Phenol-d6	7.386	99	159717	51.380	ng	-0.01
23) Nitrobenzene-d5	9.425	82	249490	96.886	ng	0.00
42) 2,4,6-Tribromophenol	16.358	330	224260	180.187	ng	0.00
45) 2-Fluorobiphenyl	13.503	172	581151	68.722	ng	0.00
79) Terphenyl-d14	20.207	244	1256419	78.415	ng	0.00
Target Compounds						
50) Dimethylphthalate	14.313	163	21886	2.258	ng	Qvalue 99

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

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