

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111522\
 Data File : BG055632.D
 Acq On : 15 Nov 2022 21:50
 Operator : CG/JU
 Sample : N5431-08
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 DISSOLVED-20221178-COMP-13-MODIFIED-ELUTRIATE

Quant Time: Nov 16 01:46:13 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	8.255	152	39725	20.000	ng	0.00
21) Naphthalene-d8	11.086	136	161853	20.000	ng	0.00
39) Acenaphthene-d10	14.876	164	117720	20.000	ng	0.00
64) Phenanthrene-d10	17.620	188	306681	20.000	ng	0.00
76) Chrysene-d12	21.915	240	313313	20.000	ng	0.00
86) Perylene-d12	25.329	264	322768	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.781	112	175639	84.215	ng	0.00
7) Phenol-d6	7.391	99	179062	62.155	ng	0.00
23) Nitrobenzene-d5	9.424	82	264443	110.761	ng	0.00
42) 2,4,6-Tribromophenol	16.357	330	232446	202.373	ng	0.00
45) 2-Fluorobiphenyl	13.507	172	602730	77.230	ng	0.00
79) Terphenyl-d14	20.205	244	1227971	78.722	ng	0.00
Target Compounds						
50) Dimethylphthalate	14.312	163	21592	2.414	ng	Qvalue 99

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

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