

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111522\
 Data File : BG055636.D
 Acq On : 16 Nov 2022 00:35
 Operator : CG/JU
 Sample : N5431-05RE
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 DISSOLVED-20221170-COMP-05-MODIFIED-ELUTRIATERE

Quant Time: Nov 16 01:47:48 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	8.259	152	40182	20.000	ng	0.00
21) Naphthalene-d8	11.085	136	169590	20.000	ng	0.00
39) Acenaphthene-d10	14.875	164	124326	20.000	ng	0.00
64) Phenanthrene-d10	17.619	188	310985	20.000	ng	0.00
76) Chrysene-d12	21.914	240	310433	20.000	ng	0.00
86) Perylene-d12	25.333	264	327780	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.780	112	81750	38.752	ng	0.00
7) Phenol-d6	7.389	99	80592	27.656	ng	0.00
23) Nitrobenzene-d5	9.422	82	141471	56.552	ng	0.00
42) 2,4,6-Tribromophenol	16.355	330	128345	105.803	ng	0.00
45) 2-Fluorobiphenyl	13.506	172	351129	42.601	ng	0.00
79) Terphenyl-d14	20.204	244	681564	44.098	ng	0.00

Target Compounds	Qvalue

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

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