

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121021\
 Data File : BM033394.D
 Acq On : 10 Dec 2021 15:01
 Operator : CG/JU
 Sample : M4919-03
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 FDR-MW-4-20211202

Quant Time: Dec 11 02:10:44 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM120821.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 09 07:11:07 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.901	152	58288	20.000	ng	-0.01
21) Naphthalene-d8	10.701	136	240346	20.000	ng	0.00
39) Acenaphthene-d10	14.530	164	177829	20.000	ng	0.00
64) Phenanthrene-d10	17.271	188	415452	20.000	ng	0.00
76) Chrysene-d12	21.430	240	387367	20.000	ng	-0.01
86) Perylene-d12	23.747	264	321344	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.490	112	193921	54.579	ng	0.00
7) Phenol-d6	7.078	99	156716	31.634	ng	-0.01
23) Nitrobenzene-d5	9.066	82	522240	92.261	ng	-0.01
42) 2,4,6-Tribromophenol	16.018	330	309506	162.703	ng	0.00
45) 2-Fluorobiphenyl	13.154	172	1058771	82.357	ng	0.00
79) Terphenyl-d14	19.883	244	1491230	71.322	ng	0.00

Target Compounds	Qvalue

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

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