

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121021\
 Data File : BM033403.D
 Acq On : 10 Dec 2021 20:26
 Operator : CG/JU
 Sample : M4927-02
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PH1-BOT-4

Quant Time: Dec 11 02:14:41 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	7.907	152	96362	20.000	ng	0.00
21) Naphthalene-d8	10.701	136	365397	20.000	ng	0.00
39) Acenaphthene-d10	14.530	164	199048	20.000	ng	0.00
64) Phenanthrene-d10	17.265	188	363418	20.000	ng	-0.01
76) Chrysene-d12	21.430	240	309072	20.000	ng	#-0.01
86) Perylene-d12	23.747	264	300753	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.490	112	477594	81.309	ng	0.00
7) Phenol-d6	7.078	99	799923	97.670	ng	-0.01
23) Nitrobenzene-d5	9.066	82	563316	65.459	ng	-0.01
42) 2,4,6-Tribromophenol	16.012	330	185715	87.220	ng	-0.01
45) 2-Fluorobiphenyl	13.154	172	1054133	73.255	ng	0.00
79) Terphenyl-d14	19.877	244	1207528	72.383	ng	-0.01
Target Compounds						
10) Phenol	7.101	94	23256	2.845	ng	Qvalue 95

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

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