

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121021\
 Data File : BM033409.D
 Acq On : 11 Dec 2021 00:02
 Operator : CG/JU
 Sample : M4927-12
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PH1-BOT-13

Quant Time: Dec 11 02:17:18 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	73670	20.000	ng	-0.01
21) Naphthalene-d8	10.695	136	373048	20.000	ng	-0.01
39) Acenaphthene-d10	14.530	164	225160	20.000	ng	# 0.00
64) Phenanthrene-d10	17.265	188	461789	20.000	ng	-0.01
76) Chrysene-d12	21.430	240	401785	20.000	ng	-0.01
86) Perylene-d12	23.747	264	355806	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.490	112	163431	36.394	ng	0.00
7) Phenol-d6	7.078	99	474043	75.708	ng	-0.01
23) Nitrobenzene-d5	9.066	82	628833	71.574	ng	-0.01
42) 2,4,6-Tribromophenol	16.019	330	227430	94.425	ng	0.00
45) 2-Fluorobiphenyl	13.154	172	1333708	81.935	ng	0.00
79) Terphenyl-d14	19.877	244	1798565	82.934	ng	-0.01

Target Compounds	Qvalue

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

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