

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121821\
 Data File : BM033526.D
 Acq On : 18 Dec 2021 19:26
 Operator : CG/JU
 Sample : M5115-02
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 MW-2R

Quant Time: Dec 20 07:08:48 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM121621.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 16 17:10:15 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.931	152	29496	20.000	ng	-0.01
21) Naphthalene-d8	10.736	136	137464	20.000	ng	#-0.02
39) Acenaphthene-d10	14.577	164	111114	20.000	ng	#-0.01
64) Phenanthrene-d10	17.324	188	257173	20.000	ng	-0.01
76) Chrysene-d12	21.500	240	240403	20.000	ng	-0.01
86) Perylene-d12	23.912	264	219639	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.496	112	157483	92.129	ng	0.00
7) Phenol-d6	7.101	99	136334	59.431	ng	-0.01
23) Nitrobenzene-d5	9.101	82	335779	104.882	ng	-0.01
42) 2,4,6-Tribromophenol	16.071	330	215207	172.178	ng	0.00
45) 2-Fluorobiphenyl	13.201	172	761340	94.782	ng	-0.01
79) Terphenyl-d14	19.948	244	1302463	85.295	ng	0.00
Target Compounds						
						Qvalue
52) Acenaphthene	14.636	154	15267	2.495	ng	96
58) Fluorene	15.624	166	26350	3.166	ng	99
73) Carbazole	17.736	167	58625	4.745	ng	# 94

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

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