

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121821\
 Data File : BM033533.D
 Acq On : 18 Dec 2021 23:41
 Operator : CG/JU
 Sample : M5064-01RE
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SASP2-121-1-(0-5)RE

Quant Time: Dec 20 07:10:21 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	45387	20.000	ng	-0.01
21) Naphthalene-d8	10.742	136	203706	20.000	ng	-0.01
39) Acenaphthene-d10	14.577	164	146407	20.000	ng	-0.01
64) Phenanthrene-d10	17.324	188	311450	20.000	ng	#-0.01
76) Chrysene-d12	21.500	240	273869	20.000	ng	-0.01
86) Perylene-d12	23.912	264	219842	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.496	112	264794	100.670	ng	0.00
7) Phenol-d6	7.101	99	377221	106.865	ng	-0.01
23) Nitrobenzene-d5	9.101	82	288096	60.726	ng	-0.01
42) 2,4,6-Tribromophenol	16.071	330	159019	96.555	ng	0.00
45) 2-Fluorobiphenyl	13.201	172	628929	59.423	ng	-0.01
79) Terphenyl-d14	19.948	244	1006905	57.882	ng	0.00
Target Compounds						
10) Phenol	7.125	94	9209	2.582	ng	# 57

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

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