

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM120621\
 Data File : BM033306.D
 Acq On : 06 Dec 2021 11:04
 Operator : CG/JU
 Sample : PB141148BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB141148BL

Quant Time: Dec 06 13:02:23 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM112421.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 06 13:01:33 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.925	152	166370	20.000	ng	0.00
21) Naphthalene-d8	10.719	136	663627	20.000	ng	# 0.00
39) Acenaphthene-d10	14.548	164	389556	20.000	ng	0.00
64) Phenanthrene-d10	17.289	188	724819	20.000	ng	# 0.00
76) Chrysene-d12	21.447	240	521303	20.000	ng	# 0.00
86) Perylene-d12	23.777	264	466355	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.513	112	1443898	143.320	ng	0.00
7) Phenol-d6	7.107	99	1898404	140.698	ng	0.00
23) Nitrobenzene-d5	9.089	82	1267426	88.661	ng	0.00
42) 2,4,6-Tribromophenol	16.036	330	451069	106.221	ng	0.00
45) 2-Fluorobiphenyl	13.172	172	2333804	84.490	ng	0.00
79) Terphenyl-d14	19.895	244	2515952	95.434	ng	0.00

Target Compounds	Qvalue

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

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