

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110123\
 Data File : BM042539.D
 Acq On : 01 Nov 2023 14:44
 Operator : MA/JU
 Sample : PB156734TB
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB156734TB

Quant Time: Nov 01 23:42:28 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM103023.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 31 02:55:18 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.922	152	86623	20.000	ng	0.00
21) Naphthalene-d8	10.739	136	311400	20.000	ng	-0.01
39) Acenaphthene-d10	14.569	164	186275	20.000	ng	-0.01
64) Phenanthrene-d10	17.321	188	366846	20.000	ng	0.00
76) Chrysene-d12	21.503	240	308302	20.000	ng	0.00
86) Perylene-d12	23.921	264	348181	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.457	112	767582	143.009	ng	0.00
7) Phenol-d6	7.087	99	959564	130.355	ng	-0.01
23) Nitrobenzene-d5	9.122	82	653663	97.853	ng	-0.01
42) 2,4,6-Tribromophenol	16.068	330	359889	147.982	ng	0.00
45) 2-Fluorobiphenyl	13.192	172	1340178	96.811	ng	0.00
79) Terphenyl-d14	19.939	244	1943695	106.900	ng	0.00

Target Compounds	Qvalue

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

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