

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110123\
 Data File : BM042537.D
 Acq On : 01 Nov 2023 13:31
 Operator : MA/JU
 Sample : PB156808BL
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB156808BL

Quant Time: Nov 01 23:41:42 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	93142	20.000	ng	0.00
21) Naphthalene-d8	10.739	136	338034	20.000	ng	-0.01
39) Acenaphthene-d10	14.568	164	206236	20.000	ng	-0.01
64) Phenanthrene-d10	17.321	188	410602	20.000	ng	0.00
76) Chrysene-d12	21.503	240	308624	20.000	ng	0.00
86) Perylene-d12	23.921	264	327586	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.457	112	814309	141.096	ng	0.00
7) Phenol-d6	7.092	99	1023461	129.305	ng	0.00
23) Nitrobenzene-d5	9.122	82	707467	97.563	ng	-0.01
42) 2,4,6-Tribromophenol	16.068	330	404925	150.385	ng	0.00
45) 2-Fluorobiphenyl	13.192	172	1489838	97.205	ng	0.00
79) Terphenyl-d14	19.939	244	2062705	113.327	ng	0.00

Target Compounds	Qvalue

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

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