

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111222\
 Data File : BM037497.D
 Acq On : 12 Nov 2022 18:27
 Operator : CG/JU
 Sample : PB148287BL
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB148287BL

Quant Time: Nov 14 03:13:04 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM111222.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 14 01:42:00 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.798	152	272535	20.000	ng	0.00
21) Naphthalene-d8	10.598	136	1163462	20.000	ng	0.00
39) Acenaphthene-d10	14.445	164	794680	20.000	ng	0.00
64) Phenanthrene-d10	17.192	188	1721231	20.000	ng	0.00
76) Chrysene-d12	21.374	240	1820180	20.000	ng	0.00
86) Perylene-d12	23.715	264	1832685	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.375	112	2175086	150.716	ng	0.00
7) Phenol-d6	6.987	99	2866070	138.031	ng	0.00
23) Nitrobenzene-d5	8.969	82	1704567	76.135	ng	0.00
42) 2,4,6-Tribromophenol	15.939	330	1514077	137.792	ng	0.00
45) 2-Fluorobiphenyl	13.069	172	4324747	75.633	ng	0.00
79) Terphenyl-d14	19.821	244	8605094	94.417	ng	0.00

Target Compounds	Qvalue

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

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