

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061121\
 Data File : BM030401.D
 Acq On : 11 Jun 2021 12:55
 Operator : CG/JU
 Sample : PB136995BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB136995BL

Quant Time: Jun 11 14:11:01 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM060921.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 11 14:10:50 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.839	152	207048	20.000	ng	0.00
21) Naphthalene-d8	10.627	136	774677	20.000	ng	0.00
39) Acenaphthene-d10	14.457	164	445334	20.000	ng	0.00
64) Phenanthrene-d10	17.192	188	794643	20.000	ng	0.00
76) Chrysene-d12	21.356	240	604509	20.000	ng	0.00
86) Perylene-d12	23.638	264	609484	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.422	112	1891572	148.519	ng	0.00
7) Phenol-d6	7.004	99	2353729	127.660	ng	0.00
23) Nitrobenzene-d5	8.992	82	1380909	86.842	ng	0.00
42) 2,4,6-Tribromophenol	15.945	330	671054	119.851	ng	0.00
45) 2-Fluorobiphenyl	13.086	172	3072687	90.736	ng	0.00
79) Terphenyl-d14	19.809	244	3285825	95.472	ng	0.00

Target Compounds	Qvalue

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

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