

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061021\
 Data File : BM030393.D
 Acq On : 11 Jun 2021 06:11
 Operator : CG/JU
 Sample : M2612-32
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 18-B12(50-54)

Quant Time: Jun 11 07:55:40 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM060921.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 08 20:51:08 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.840	152	149476	20.000	ng	-0.01
21) Naphthalene-d8	10.628	136	639965	20.000	ng	#-0.01
39) Acenaphthene-d10	14.463	164	443763	20.000	ng	-0.01
64) Phenanthrene-d10	17.198	188	1043638	20.000	ng	-0.01
76) Chrysene-d12	21.356	240	1141648	20.000	ng	-0.01
86) Perylene-d12	23.633	264	1042232	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.428	112	1510554	164.283	ng	0.00
7) Phenol-d6	7.004	99	1845874	138.676	ng	0.00
23) Nitrobenzene-d5	8.992	82	1191524	90.705	ng	-0.02
42) 2,4,6-Tribromophenol	15.951	330	782222	140.200	ng	0.00
45) 2-Fluorobiphenyl	13.086	172	2984533	88.445	ng	-0.01
79) Terphenyl-d14	19.815	244	6185109	95.159	ng	0.00
Target Compounds						
32) Benzoic acid	9.875	122	46475	10.492	ng	Qvalue 95

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

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