

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061021\
 Data File : BM030374.D
 Acq On : 10 Jun 2021 14:56
 Operator : CG/JU
 Sample : M2613-15
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 18-B12(90-94)

Quant Time: Jun 10 15:51:45 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	7.845	152	194045	20.000	ng	0.00
21) Naphthalene-d8	10.633	136	763166	20.000	ng	# 0.00
39) Acenaphthene-d10	14.468	164	529542	20.000	ng	0.00
64) Phenanthrene-d10	17.204	188	1291349	20.000	ng	0.00
76) Chrysene-d12	21.362	240	1428406	20.000	ng	0.00
86) Perylene-d12	23.638	264	1281466	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.428	112	1434870	120.210	ng	0.00
7) Phenol-d6	7.010	99	1726371	99.908	ng	0.00
23) Nitrobenzene-d5	8.998	82	1190918	76.024	ng	-0.01
42) 2,4,6-Tribromophenol	15.951	330	932748	140.098	ng	0.00
45) 2-Fluorobiphenyl	13.092	172	2973752	73.850	ng	0.00
79) Terphenyl-d14	19.815	244	6422639	78.976	ng	0.00
Target Compounds						
32) Benzoic acid	9.881	122	23548	7.327	ng	Qvalue 96
84) Bis(2-ethylhexyl)phtha...	21.274	149	13157	4.452	ng	95

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

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