

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061021\
 Data File : BM030383.D
 Acq On : 10 Jun 2021 20:29
 Operator : CG/JU
 Sample : M2613-24
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 18-B12(126-130)

Quant Time: Jun 11 07:53:48 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	7.851	152	3448	20.000	ng	0.00
21) Naphthalene-d8	10.639	136	11232	20.000	ng	# 0.00
39) Acenaphthene-d10	14.468	164	7295	20.000	ng	0.00
64) Phenanthrene-d10	17.204	188	17961	20.000	ng	# 0.00
76) Chrysene-d12	21.368	240	20765	20.000	ng	# 0.00
86) Perylene-d12	23.650	264	17668	20.000	ng	# 0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.428	112	2217	10.453	ng	0.00
7) Phenol-d6	7.022	99	21013	68.437	ng	0.01
23) Nitrobenzene-d5	8.998	82	7944	34.456	ng	-0.01
42) 2,4,6-Tribromophenol	0.000	330	0	0.000	ng	
45) 2-Fluorobiphenyl	13.092	172	38190	68.845	ng	0.00
79) Terphenyl-d14	19.815	244	102502	86.703	ng	0.00
Target Compounds						
10) Phenol	7.045	94	5443	17.548	ng	Qvalue 98

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

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