

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF091922\
 Data File : BF130325.D
 Acq On : 19 Sep 2022 17:34
 Operator : CG\JU
 Sample : N4648-01DL2 10X
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 N48854-1040DL2

Quant Time: Sep 20 02:26:12 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Sep 19 15:06:14 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.657	152	106269	20.000	ng	0.00
21) Naphthalene-d8	7.945	136	391268	20.000	ng	0.00
39) Acenaphthene-d10	9.692	164	210509	20.000	ng	0.00
64) Phenanthrene-d10	11.168	188	348088	20.000	ng	0.00
76) Chrysene-d12	13.792	240	177860	20.000	ng	-0.01
86) Perylene-d12	15.180	264	200840	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.245	112	26988	4.405	ng	-0.01
7) Phenol-d6	6.286	99	23061	3.008	ng	-0.01
23) Nitrobenzene-d5	7.222	82	58077	7.583	ng	-0.01
42) 2,4,6-Tribromophenol	10.469	330	18652	9.247	ng	-0.01
45) 2-Fluorobiphenyl	9.016	172	104580	7.234	ng	0.00
79) Terphenyl-d14	12.751	244	85099	7.926	ng	0.00
Target Compounds						
						Qvalue
10) Phenol	6.304	94	375596	41.807	ng	94
17) 2-Methylphenol	6.928	107	826563	145.687	ng	99
20) 3+4-Methylphenols	7.086	107	1059357	143.734	ng	# 53
27) 2,4-Dimethylphenol	7.628	122	1287662	262.335	ng	# 88
31) Naphthalene	7.969	128	191124	9.481	ng	99

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

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