

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF091922\
 Data File : BF130324.D
 Acq On : 19 Sep 2022 16:59
 Operator : CG\JU
 Sample : N4648-01DL 2X
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 N48854-1040DL

Quant Time: Sep 20 02:25:38 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	6.663	152	96166	20.000	ng	0.00
21) Naphthalene-d8	7.963	136	306633	20.000	ng	0.02
39) Acenaphthene-d10	9.692	164	162692	20.000	ng	0.00
64) Phenanthrene-d10	11.169	188	228960	20.000	ng	0.00
76) Chrysene-d12	13.798	240	168777	20.000	ng	0.00
86) Perylene-d12	15.186	264	219559	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.251	112	128772	23.226	ng	0.00
7) Phenol-d6	6.304	99	102410	14.762	ng	0.00
23) Nitrobenzene-d5	7.239	82	250145	41.677	ng	0.00
42) 2,4,6-Tribromophenol	10.474	330	69231	44.410	ng	0.00
45) 2-Fluorobiphenyl	9.027	172	418481	37.457	ng	0.00
79) Terphenyl-d14	12.751	244	243033	23.853	ng	0.00
Target Compounds						
						Qvalue
10) Phenol	6.328	94	1617250	198.926	ng	93
17) 2-Methylphenol	6.957	107	3290818	640.962	ng	97
20) 3+4-Methylphenols	7.128	107	4140773	620.844	ng	# 52
27) 2,4-Dimethylphenol	7.663	122	4987419	1296.540	ng	# 82
31) Naphthalene	7.986	128	687169	43.499	ng	99
38) 1-Methylnaphthalene	8.763	142	22160	2.290	ng	81
73) Carbazole	11.404	167	61610	5.503	ng	98

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

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