

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF090421\
 Data File : BF125368.D
 Acq On : 4 Sep 2021 14:34
 Operator : JU/CG
 Sample : M3646-13
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 1992

Quant Time: Sep 06 08:27:54 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF081821.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 03 11:56:34 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.898	152	127125	20.000	ng	0.00
21) Naphthalene-d8	8.180	136	467584	20.000	ng	# 0.00
39) Acenaphthene-d10	9.939	164	214811	20.000	ng	0.00
64) Phenanthrene-d10	11.427	188	351356	20.000	ng	0.00
76) Chrysene-d12	14.062	240	286380	20.000	ng	# 0.00
86) Perylene-d12	15.556	264	327226	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.522	112	320583	38.170	ng	0.00
7) Phenol-d6	6.528	99	248803	22.905	ng	0.00
23) Nitrobenzene-d5	7.463	82	722141	77.752	ng	0.00
42) 2,4,6-Tribromophenol	10.727	330	179970	83.930	ng	0.00
45) 2-Fluorobiphenyl	9.257	172	1127639	73.163	ng	0.00
79) Terphenyl-d14	13.010	244	1099821	61.181	ng	0.00
Target Compounds						
						Qvalue
3) Pyridine	3.510	79	74555	6.814	ng	# 91
10) Phenol	6.545	94	42119	3.703	ng	89
20) 3+4-Methylphenols	7.310	107	794516	87.941	ng	# 74
27) 2,4-Dimethylphenol	7.851	122	19471	2.752	ng	# 93
31) Naphthalene	8.198	128	56541	2.361	ng	# 91
37) 2-Methylnaphthalene	8.892	142	35039	2.214	ng	95
38) 1-Methylnaphthalene	8.992	142	39153	2.650	ng	98
71) Phenanthrene	11.451	178	39968	2.064	ng	# 92

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

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