

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF091421\
 Data File : BF125484.D
 Acq On : 14 Sep 2021 20:41
 Operator : JU/CG
 Sample : M3724-10
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TOTE-COMP

Quant Time: Sep 15 01:06:39 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF081821.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 14 17:17:24 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.875	152	34114	20.000	ng	0.00
21) Naphthalene-d8	8.157	136	142703	20.000	ng	# 0.00
39) Acenaphthene-d10	9.916	164	74441	20.000	ng	0.00
64) Phenanthrene-d10	11.404	188	142980	20.000	ng	0.00
76) Chrysene-d12	14.051	240	114021	20.000	ng	#-0.01
86) Perylene-d12	15.539	264	82857	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.487	112	222643	98.784	ng	-0.02
7) Phenol-d6	6.504	99	255899	87.789	ng	-0.02
23) Nitrobenzene-d5	7.434	82	194099	68.476	ng	-0.02
42) 2,4,6-Tribromophenol	10.704	330	57239	77.029	ng	-0.01
45) 2-Fluorobiphenyl	9.233	172	374940	70.199	ng	0.00
79) Terphenyl-d14	12.986	244	391281	54.669	ng	-0.01
Target Compounds						
10) Phenol	6.522	94	92632	30.345	ng	Qvalue 99
20) 3+4-Methylphenols	7.281	107	131053	54.055	ng	# 73
50) Dimethylphthalate	9.628	163	26319	5.102	ng	# 98

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

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