

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF091521\
 Data File : BF125511.D
 Acq On : 15 Sep 2021 18:55
 Operator : JU/CG
 Sample : M3724-10RE
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TOTE-COMP

Quant Time: Sep 16 03:06:45 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF091521.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 15 15:08:32 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.869	152	314655	20.000	ng	0.00
21) Naphthalene-d8	8.151	136	1228148	20.000	ng	# 0.00
39) Acenaphthene-d10	9.910	164	663632	20.000	ng	0.00
64) Phenanthrene-d10	11.398	188	1217321	20.000	ng	0.00
76) Chrysene-d12	14.045	240	912140	20.000	ng	# 0.00
86) Perylene-d12	15.533	264	735203	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.510	112	992700	50.128	ng	0.02
7) Phenol-d6	6.522	99	1191355	45.264	ng	0.01
23) Nitrobenzene-d5	7.439	82	958362	38.528	ng	0.00
42) 2,4,6-Tribromophenol	10.704	330	278698	34.481	ng	0.00
45) 2-Fluorobiphenyl	9.227	172	1649253	36.467	ng	0.00
79) Terphenyl-d14	12.986	244	1611092	28.641	ng	0.00
Target Compounds						
10) Phenol	6.534	94	448178	15.027	ng	Qvalue 94
20) 3+4-Methylphenols	7.286	107	627545	33.757	ng	# 75
50) Dimethylphthalate	9.622	163	136976	2.895	ng	# 97

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

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