

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF091421\
 Data File : BF125482.D
 Acq On : 14 Sep 2021 19:36
 Operator : JU/CG
 Sample : M3724-02
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 31190

Quant Time: Sep 15 01:06:14 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	140407	20.000	ng	0.00
21) Naphthalene-d8	8.157	136	550234	20.000	ng	# 0.00
39) Acenaphthene-d10	9.916	164	300469	20.000	ng	0.00
64) Phenanthrene-d10	11.404	188	535935	20.000	ng	0.00
76) Chrysene-d12	14.051	240	245511	20.000	ng	#-0.01
86) Perylene-d12	15.539	264	284573	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.516	112	891309	96.084	ng	0.01
7) Phenol-d6	6.516	99	1039242	86.623	ng	0.00
23) Nitrobenzene-d5	7.439	82	848207	77.607	ng	-0.01
42) 2,4,6-Tribromophenol	10.710	330	373050	124.377	ng	0.00
45) 2-Fluorobiphenyl	9.233	172	1503904	69.759	ng	0.00
79) Terphenyl-d14	12.992	244	1442804	93.622	ng	0.00
Target Compounds						Qvalue
20) 3+4-Methylphenols	7.292	107	25998	2.605	ng	# 74
32) Benzoic acid	7.963	122	170217	30.125	ng	87
50) Dimethylphthalate	9.627	163	53265	2.558	ng	# 95

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

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