

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF092921\
 Data File : BF125714.D
 Acq On : 29 Sep 2021 18:19
 Operator : JU/CG
 Sample : M3964-03
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CC-3

Quant Time: Sep 30 03:27:45 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF092421.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 24 15:38:32 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.886	152	240770	20.000	ng	0.00
21) Naphthalene-d8	8.169	136	973936	20.000	ng	# 0.00
39) Acenaphthene-d10	9.922	164	473749	20.000	ng	0.00
64) Phenanthrene-d10	11.404	188	693316	20.000	ng	# 0.00
76) Chrysene-d12	14.045	240	474916	20.000	ng	0.00
86) Perylene-d12	15.521	264	463371	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.510	112	1014158	69.398	ng	0.00
7) Phenol-d6	6.516	99	1615102	82.206	ng	0.00
23) Nitrobenzene-d5	7.445	82	1048666	75.102	ng	0.00
42) 2,4,6-Tribromophenol	10.704	330	142901	32.352	ng	-0.01
45) 2-Fluorobiphenyl	9.245	172	1994024	73.878	ng	0.00
79) Terphenyl-d14	12.992	244	1628309	59.376	ng	0.00
Target Compounds						
77) Benzidine	12.733	184	107356	8.605	ng	# 95

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

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