

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF093021\
 Data File : BF125734.D
 Acq On : 30 Sep 2021 15:34
 Operator : JU/CG
 Sample : M3971-11
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 KEARNY-COMP

Quant Time: Sep 30 16:02:41 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	6.887	152	212019	20.000	ng	0.00
21) Naphthalene-d8	8.169	136	860829	20.000	ng	# 0.00
39) Acenaphthene-d10	9.922	164	462051	20.000	ng	0.00
64) Phenanthrene-d10	11.404	188	794122	20.000	ng	# 0.00
76) Chrysene-d12	14.039	240	567948	20.000	ng	0.00
86) Perylene-d12	15.515	264	438181	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.516	112	1487604	115.599	ng	0.01
7) Phenol-d6	6.510	99	1644121	95.031	ng	0.00
23) Nitrobenzene-d5	7.445	82	1166239	94.497	ng	0.00
42) 2,4,6-Tribromophenol	10.710	330	575237	133.527	ng	0.00
45) 2-Fluorobiphenyl	9.245	172	2240111	87.508	ng	0.00
79) Terphenyl-d14	12.992	244	2462466	75.084	ng	0.00
Target Compounds						
77) Benzidine	12.739	184	1342125	89.954	ng	Qvalue 97

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

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