

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF041318\
 Data File : BF104498.D
 Acq On : 14 Apr 2018 00:02
 Operator : JU/SJ
 Sample : PB108229BL
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :

Quant Time: Apr 14 05:18:21 2018
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 13 17:38:31 2018
 Response via : Initial Calibration

Internal Standards	R.T.	OIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.81	152	132878	20.00	ng	0.00
21) Naphthalene-d8	8.09	136	515697	20.00	ng	0.00
38) Acenaphthene-d10	9.85	164	215874	20.00	ng	0.00
63) Phenanthrene-d10	11.34	188	322691	20.00	ng	0.00
75) Chrysene-d12	13.98	240	248928	20.00	ng	-0.01
86) Perylene-d12	15.45	264	183831	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.43	112	556903	64.08	ng	0.00
7) Phenol-d6	6.45	99	673327	63.06	ng	0.00
23) Nitrobenzene-d5	7.37	82	407951	51.66	ng	0.00
41) 2,4,6-Tribromophenol	10.64	330	117248	52.36	ng	0.00
44) 2-Fluorobiphenyl	9.17	172	747561	54.71	ng	0.00
78) Terphenyl-d14	12.93	244	538768	49.34	ng	0.00

Target Compounds

						Qvalue
19) n-Nitroso-di-n-propylamine	7.37	70	52934	7.587	ng	# 72
25) Isophorone	7.37	82	408015	24.431	ng	# 64
49) Dimethylphthalate	9.56	163	44358	2.606	ng	# 99
50) 2,6-Dinitrotoluene	9.85	165	27023	8.579	ng	# 26
56) 2,4-Dinitrotoluene	9.85	165	27023	6.540	ng	# 24
64) 4,6-Dinitro-2-methylphenol	10.64	198	128	7.416	ng	# 1
66) 4-Bromophenyl-phenylether	10.64	248	7272	2.119	ng	# 1
76) Benzidine	12.93	184	6628	Below Cal		98

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

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