

Data Path : Z:\HPCHEM1\BNA_F\Data\BF111617\
 Data File : BF100625.D
 Acq On : 16 Nov 2017 12:49
 Operator : SJ/JU
 Sample : I6469-01
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SV30666L

Quant Time: Nov 16 13:59:56 2017
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF110817.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 16 11:34:03 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.88	152	127477	20.00	ng	-0.02
21) Naphthalene-d8	8.16	136	500332	20.00	ng	-0.02
38) Acenaphthene-d10	9.92	164	229415	20.00	ng	-0.02
63) Phenanthrene-d10	11.40	188	426936	20.00	ng	-0.02
75) Chrysene-d12	14.04	240	266110	20.00	ng	-0.03
86) Perylene-d12	15.52	264	237798	20.00	ng	-0.03

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.49	112	420565	52.88	ng	-0.02
7) Phenol-d6	6.50	99	321760	32.81	ng	-0.02
23) Nitrobenzene-d5	7.45	82	611737	80.83	ng	-0.02
41) 2,4,6-Tribromophenol	10.71	330	253875	108.60	ng	-0.02
44) 2-Fluorobiphenyl	9.24	172	1120434	74.37	ng	-0.02
78) Terphenyl-d14	12.99	244	894314	77.22	ng	-0.02

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
9) Benzaldehyde	6.65	77	14147	2.020	ng	86
19) n-Nitroso-di-n-propylamine	7.45	70	79846	11.741	ng	# 80
25) Isophorone	7.45	82	611281	38.438	ng	# 64
32) Benzoic acid	7.92	122	2608	9.755	ng	# 60
47) 2-Nitroaniline	9.46	65	250	2.839	ng	# 19
50) 2,6-Dinitrotoluene	9.92	165	31294	9.026	ng	# 26
56) 2,4-Dinitrotoluene	9.92	165	31294	7.155	ng	# 23
64) 4,6-Dinitro-2-methylphenol	10.70	198	526	8.770	ng	# 1
66) 4-Bromophenyl-phenylether	10.71	248	15810	3.454	ng	# 1
84) Di-n-octyl phthalate	14.63	149	67860	3.882	ng	# 74

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

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