

Data Path : Z:\HPCHEM1\BNA F\DATA\BF100717\
 Data File : BF099220.D
 Acq On : 8 Oct 2017 4:47
 Operator : SJ/JU
 Sample : I5714-03
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TR-05-100517

Quant Time: Oct 08 06:06:42 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF092317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sun Oct 08 03:54:40 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	136724	20.00	ng	0.00
21) Naphthalene-d8	8.15	136	508557	20.00	ng	0.00
38) Acenaphthene-d10	9.90	164	193500	20.00	ng	0.00
63) Phenanthrene-d10	11.39	188	297915	20.00	ng	0.00
75) Chrysene-d12	14.03	240	239966	20.00	ng	0.00
86) Perylene-d12	15.50	264	182947	20.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.50	112	956186	111.33	ng	0.02
7) Phenol-d6	6.50	99	1032438	97.44	ng	0.00
23) Nitrobenzene-d5	7.43	82	694717	87.61	ng	0.00
41) 2,4,6-Tribromophenol	10.69	330	188672	119.16	ng	0.00
44) 2-Fluorobiphenyl	9.22	172	1192515	102.88	ng	0.00
78) Terphenyl-d14	12.97	244	867198	82.19	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
9) Benzaldehyde	6.64	77	19705	3.206	ng	83
15) Benzyl Alcohol	7.00	79	113002	14.538	ng	97
17) 2-Methylphenol	7.00	107	60809	7.641	ng	# 22
19) n-Nitroso-di-n-propylamine	7.43	70	88369	13.064	ng	# 75
20) 3+4-Methylphenols	7.00	107	60808	6.040	ng	# 1
25) Isophorone	7.43	82	694708	42.372	ng	# 64
50) 2,6-Dinitrotoluene	9.90	165	24374	7.666	ng	# 25
56) 2,4-Dinitrotoluene	9.90	165	24374	5.907	ng	# 22
66) 4-Bromophenyl-phenylether	10.69	248	12528	4.296	ng	# 1

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

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