

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF091118\
 Data File : BF108945.D
 Acq On : 11 Sep 2018 19:44
 Operator : JU/SJ
 Sample : J4724-16
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EP1-SUT-4USTS-S

Quant Time: Sep 12 03:35:51 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF090518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 05 18:03:31 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.74	152	1843	20.00	ng	0.00
21) Naphthalene-d8	8.09	136	320785	20.00	ng	0.07
38) Acenaphthene-d10	9.77	164	254479	20.00	ng	0.00
63) Phenanthrene-d10	11.24	188	402541	20.00	ng	0.00
75) Chrysene-d12	13.87	240	335283	20.00	ng	0.00
86) Perylene-d12	15.28	264	429444	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.35	112	591841	5320.04	ng	0.01
7) Phenol-d6	6.39	99	537467	3754.91	ng	0.02
23) Nitrobenzene-d5	7.37	82	70766	13.81	ng	0.07
41) 2,4,6-Tribromophenol	10.55	330	171943	70.78	ng	0.00
44) 2-Fluorobiphenyl	9.10	172	928726	62.26	ng	0.00
78) Terphenyl-d14	12.83	244	689173	47.93	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.43	88	577	10.792	ng	# 53
6) Aniline	6.54	93	47535	246.914	ng	# 1
9) Benzaldehyde	6.23	77	30243	297.826	ng	# 11
10) Phenol	6.28	94	9666	58.959	ng	# 1
11) bis(2-Chloroethyl)ether	6.54	93	54174	412.494	ng	# 1
15) Benzyl Alcohol	6.87	79	40485	368.555	ng	# 23
17) 2-Methylphenol	6.98	107	1779	17.095	ng	# 1
18) Hexachloroethane	7.23	117	298485	5890.923	ng	# 1
19) n-Nitroso-di-n-propylamine	7.23	70	324825	3264.719	ng	# 87
20) 3+4-Methylphenols	7.17	107	57120	468.577	ng	# 1
22) Acetophenone	7.22	105	1945712	267.258	ng	# 50
24) Nitrobenzene	7.36	77	606236	110.226	ng	# 53
25) Isophorone	7.64	82	129998	12.714	ng	# 1
26) 2-Nitrophenol	7.71	139	35431	16.352	ng	# 1
28) bis(2-Chloroethoxy)methane	7.83	93	14826	2.169	ng	# 1
29) 2,4-Dichlorophenol	7.95	162	47824	10.541	ng	# 34
30) 1,2,4-Trichlorobenzene	8.01	180	10546	2.145	ng	# 46
31) Naphthalene	8.12	128	2170002	150.358	ng	# 95
32) Benzoic acid	7.83	122	9111	7.293	ng	# 1
33) 4-Chloroaniline	8.15	127	31061	4.686	ng	# 1
37) 2-Methylnaphthalene	8.84	142	157956	16.567	ng	# 93
49) Dimethylphthalate	9.48	163	289971	16.291	ng	# 99
53) 2,4-Dinitrophenol	10.05	184	369	9.561	ng	# 1
64) 4,6-Dinitro-2-methylphenol	10.55	198	456	7.115	ng	# 1
66) 4-Bromophenyl-phenylether	10.55	248	10700	2.353	ng	# 1
73) Di-n-butylphthalate	11.81	149	81123	3.769	ng	# 99
79) Butylbenzylphthalate	13.30	149	23926	2.059	ng	# 89
83) Bis(2-ethylhexyl)phthalate	13.87	149	573015	39.246	ng	# 99
84) Di-n-octyl phthalate	14.48	149	54267	2.229	ng	# 30

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

