

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF100416\
 Data File : BF090369.D
 Acq On : 5 Oct 2016 3:45
 Operator : UM/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040EC

Manual Integrations
 APPROVED

sohil
 10/5/2016 7:27:09 PM

Quant Time: Oct 05 04:50:36 2016
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF100416.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 05 03:16:32 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.01	152	152653	20.00	ng	0.00
21) Naphthalene-d8	8.30	136	590003	20.00	ng	0.00
38) Acenaphthene-d10	10.06	164	272460	20.00	ng	0.01
63) Phenanthrene-d10	11.55	188	463077	20.00	ng	0.00
75) Chrysene-d12	14.20	240	484668	20.00	ng	0.01
86) Perylene-d12	15.71	264	365799	20.00	ng	0.01

System Monitoring Compounds

5) 2-Fluorophenol	5.62	112	937485	96.86	ng	0.01
7) Phenol-d6	6.63	99	1072851	88.68	ng	0.01
23) Nitrobenzene-d5	7.57	82	885321	87.47	ng	0.00
41) 2,4,6-Tribromophenol	10.85	330	195847	82.76	ng	0.01
44) 2-Fluorobiphenyl	9.38	172	1405646	84.11	ng	0.00
78) Terphenyl-d14	13.14	244	1352821	64.98	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.70	88	236417	50.05	ng	# 93
3) Pyridine	3.49	79	621407	49.62	ng	# 84
4) n-Nitrosodimethylamine	3.42	42	243648	53.71	ng	# 60
6) Aniline	6.67	93	707754	42.58	ng	99
8) 2-Chlorophenol	6.79	128	453975	41.74	ng	95
9) Benzaldehyde	6.55	77	261170	46.51	ng	90
10) Phenol	6.65	94	655365	44.54	ng	94
11) bis(2-Chloroethyl)ether	6.74	93	419330	40.13	ng	# 82
12) 1,3-Dichlorobenzene	6.95	146	472627m	41.84	ng	
13) 1,4-Dichlorobenzene	7.03	146	444840	38.78	ng	95
14) 1,2-Dichlorobenzene	7.18	146	407743	37.39	ng	96
15) Benzyl Alcohol	7.15	79	405674	47.57	ng	98
16) 2,2'-oxybis(1-Chloropropan	7.29	45	758274	47.48	ng	97
17) 2-Methylphenol	7.26	107	362646	43.05	ng	97
18) Hexachloroethane	7.53	117	110964	26.74	ng	# 82
19) n-Nitroso-di-n-propylamine	7.42	70	380664	49.27	ng	# 83
20) 3+4-Methylphenols	7.41	107	443841	41.44	ng	96
22) Acetophenone	7.42	105	605232	44.31	ng	# 87
24) Nitrobenzene	7.59	77	533470	49.36	ng	96
25) Isophorone	7.83	82	898189	46.64	ng	99
26) 2-Nitrophenol	7.91	139	190020	45.57	ng	93
27) 2,4-Dimethylphenol	7.95	122	390944	41.55	ng	99
28) bis(2-Chloroethoxy)methane	8.04	93	487076	40.74	ng	# 96
29) 2,4-Dichlorophenol	8.15	162	306711	39.82	ng	95
30) 1,2,4-Trichlorobenzene	8.25	180	299721	34.74	ng	94
31) Naphthalene	8.33	128	1142173	40.50	ng	97
32) Benzoic acid	8.04	122	275995	45.89	ng	97
33) 4-Chloroaniline	8.37	127	457131	40.58	ng	# 89
34) Hexachlorobutadiene	8.44	225	159043	31.32	ng	99
35) Caprolactam	8.71	113	92662	46.00	ng	90
36) 4-Chloro-3-methylphenol	8.84	107	351609	42.51	ng	94
37) 2-Methylnaphthalene	9.01	142	636149	35.95	ng	100
39) 1,2,4,5-Tetrachlorobenzene	9.18	216	291668	37.92	ng	99
40) Hexachlorocyclopentadiene	9.17	237	29295	11.18	ng	97

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.29	196	177423	36.67	ng	97
43) 2,4,5-Trichlorophenol	9.33	196	187885	38.60	ng	# 88
45) 1,1'-Biphenyl	9.48	154	919757	43.60	ng	97
46) 2-Chloronaphthalene	9.50	162	612417	38.68	ng	99
47) 2-Nitroaniline	9.59	65	267138	61.81	ng	97
48) Acenaphthylene	9.93	152	996961	41.62	ng	96
49) Dimethylphthalate	9.78	163	737515	45.73	ng	97
50) 2,6-Dinitrotoluene	9.83	165	149588	42.40	ng	97
51) Acenaphthene	10.10	154	600904	41.23	ng	99
52) 3-Nitroaniline	10.01	138	205563	51.44	ng	# 91
53) 2,4-Dinitrophenol	10.11	184	17332	19.94	ng	# 83
54) Dibenzofuran	10.27	168	819863	42.54	ng	99
55) 4-Nitrophenol	10.15	139	172188	54.11	ng	98
56) 2,4-Dinitrotoluene	10.23	165	184374	43.42	ng	# 21
57) Fluorene	10.61	166	732247	45.42	ng	100
58) 2,3,4,6-Tetrachlorophenol	10.38	232	144968	42.23	ng	# 86
59) Diethylphthalate	10.47	149	707857	42.94	ng	# 90
60) 4-Chlorophenyl-phenylether	10.60	204	309039	40.42	ng	92
61) 4-Nitroaniline	10.61	138	192574	53.57	ng	93
62) Azobenzene	10.76	77	865296	50.36	ng	90
64) 4,6-Dinitro-2-methylphenol	10.65	198	33216	20.90	ng	# 61
65) n-Nitrosodiphenylamine	10.71	169	573408	36.31	ng	98
66) 4-Bromophenyl-phenylether	11.09	248	176265	35.22	ng	# 89
67) Hexachlorobenzene	11.16	284	194590	33.08	ng	# 87
68) Atrazine	11.24	200	135091	35.73	ng	97
69) Pentachlorophenol	11.35	266	122182	37.06	ng	98
70) Phenanthrene	11.57	178	923931	37.19	ng	97
71) Anthracene	11.63	178	1059516	42.72	ng	97
72) Carbazole	11.78	167	980446	43.01	ng	98
73) Di-n-butylphthalate	12.11	149	1182247	43.39	ng	98
74) Fluoranthene	12.77	202	1041286	44.19	ng	90
76) Benzidine	12.89	184	564744	42.37	ng	99
77) Pyrene	13.00	202	1135355	36.59	ng	98
79) Butylbenzylphthalate	13.62	149	614298	50.23	ng	# 87
80) Benzo(a)anthracene	14.19	228	1086227	39.97	ng	98
81) 3,3'-Dichlorobenzidine	14.16	252	413773	44.20	ng	# 98
82) Chrysene	14.22	228	904772	35.53	ng	98
83) Bis(2-ethylhexyl)phthalate	14.18	149	884804	45.22	ng	# 96
84) Di-n-octyl phthalate	14.80	149	1488559	52.37	ng	95
85) Indeno(1,2,3-cd)pyrene	17.23	276	745004	27.22	ng	# 94
87) Benzo(b)fluoranthene	15.26	252	931075	44.26	ng	# 96
88) Benzo(k)fluoranthene	15.30	252	942905m	46.93	ng	
89) Benzo(a)pyrene	15.64	252	810678	42.25	ng	# 94
90) Dibenzo(a,h)anthracene	17.25	278	627559	33.35	ng	96
91) Benzo(g,h,i)perylene	17.69	276	568852	30.87	ng	95

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

