

Data Path : Z:\HPCHEM1\BNA F\DATA\BF011316\
 Data File : BF084484.D
 Acq On : 14 Jan 2016 11:59
 Operator : SJ/UM
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040EC

Quant Time: Jan 14 16:35:21 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF123115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jan 14 14:15:52 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.83	152	280274	20.00	ng	0.00
21) Naphthalene-d8	8.12	136	1138599	20.00	ng	0.00
38) Acenaphthene-d10	9.87	164	469234	20.00	ng	0.00
63) Phenanthrene-d10	11.36	188	748727	20.00	ng	0.00
75) Chrysene-d12	14.00	240	584000	20.00	ng	0.00
86) Perylene-d12	15.41	264	582931	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.44	112	1432853	82.69	ng	0.00
7) Phenol-d6	6.48	99	1683592	77.36	ng	0.00
23) Nitrobenzene-d5	7.40	82	1462296	71.48	ng	0.00
41) 2,4,6-Tribromophenol	10.67	330	333266	70.35	ng	0.00
44) 2-Fluorobiphenyl	9.20	172	2298964	86.28	ng	0.00
78) Terphenyl-d14	12.95	244	1734558	66.30	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.40	88	352052	42.89	ng	# 83
3) Pyridine	3.12	79	995318	43.49	ng	# 80
4) n-Nitrosodimethylamine	3.07	42	416472	35.45	ng	# 92
6) Aniline	6.50	93	1194676	37.53	ng	95
8) 2-Chlorophenol	6.62	128	827446	42.09	ng	94
9) Benzaldehyde	6.38	77	522025	36.84	ng	88
10) Phenol	6.49	94	893796	36.12	ng	77
11) bis(2-Chloroethyl)ether	6.58	93	810516	42.29	ng	95
12) 1,3-Dichlorobenzene	6.77	146	829882	40.92	ng	# 94
13) 1,4-Dichlorobenzene	6.85	146	864562	42.51	ng	95
14) 1,2-Dichlorobenzene	7.00	146	725311	37.24	ng	94
15) Benzyl Alcohol	6.98	79	599892	32.77	ng	93
16) 2,2'-oxybis(1-Chloropropan	7.12	45	1202577	34.45	ng	86
17) 2-Methylphenol	7.10	107	680148	41.28	ng	98
18) Hexachloroethane	7.34	117	290204	35.68	ng	91
19) n-Nitroso-di-n-propylamine	7.25	70	490623	33.30	ng	# 74
20) 3+4-Methylphenols	7.25	107	701627	36.23	ng	# 55
22) Acetophenone	7.25	105	1076345	39.31	ng	# 93
24) Nitrobenzene	7.42	77	895756	43.17	ng	97
25) Isophorone	7.66	82	1517092	38.77	ng	97
26) 2-Nitrophenol	7.74	139	376067	39.82	ng	# 82
27) 2,4-Dimethylphenol	7.78	122	696751	36.45	ng	95
28) bis(2-Chloroethoxy)methane	7.88	93	963635	40.32	ng	# 96
29) 2,4-Dichlorophenol	7.98	162	625500	39.28	ng	99
30) 1,2,4-Trichlorobenzene	8.06	180	652503	39.69	ng	96
31) Naphthalene	8.14	128	2023145	39.16	ng	98
32) Benzoic acid	7.89	122	231419	22.62	ng	97
33) 4-Chloroaniline	8.20	127	933020	39.40	ng	# 88
34) Hexachlorobutadiene	8.26	225	360667	38.17	ng	99
35) Caprolactam	8.57	113	182898	34.07	ng	# 52
36) 4-Chloro-3-methylphenol	8.68	107	621099	34.82	ng	94
37) 2-Methylnaphthalene	8.84	142	1383646	37.31	ng	96
39) 1,2,4,5-Tetrachlorobenzene	9.00	216	611906	45.36	ng	100
40) Hexachlorocyclopentadiene	8.99	237	102149	12.54	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.12	196	393260	38.97	nq	95
43) 2,4,5-Trichlorophenol	9.16	196	403380	41.69	nq	96
45) 1,1'-Biphenyl	9.30	154	1540986	41.32	nq	99
46) 2-Chloronaphthalene	9.32	162	1134045	38.71	nq	95
47) 2-Nitroaniline	9.42	65	443815	45.90	nq	# 74
48) Acenaphthylene	9.73	152	1755915	40.57	nq	97
49) Dimethylphthalate	9.61	163	1497242	44.63	nq	# 91
50) 2,6-Dinitrotoluene	9.66	165	314241	42.71	nq	# 73
51) Acenaphthene	9.90	154	1135847	39.13	nq	96
52) 3-Nitroaniline	9.84	138	417686	45.82	nq	# 79
53) 2,4-Dinitrophenol	9.94	184	31443	13.14	nq	86
54) Dibenzofuran	10.08	168	1530012	40.39	nq	92
55) 4-Nitrophenol	10.00	139	213736	32.30	nq	# 71
56) 2,4-Dinitrotoluene	10.06	165	385339	41.38	nq	98
57) Fluorene	10.42	166	1088294	37.01	nq	99
58) 2,3,4,6-Tetrachlorophenol	10.20	232	308042	37.29	nq	87
59) Diethylphthalate	10.30	149	1373988	41.54	nq	95
60) 4-Chlorophenyl-phenylether	10.42	204	579978	47.77	nq	92
61) 4-Nitroaniline	10.44	138	304550	37.47	nq	86
62) Azobenzene	10.58	77	1447482	39.90	nq	97
64) 4,6-Dinitro-2-methylphenol	10.46	198	60750	16.42	nq	# 37
65) n-Nitrosodiphenylamine	10.53	169	1007058	42.07	nq	98
66) 4-Bromophenyl-phenylether	10.91	248	360757	44.77	nq	95
67) Hexachlorobenzene	10.97	284	345521	38.09	nq	# 87
68) Atrazine	11.06	200	296023	37.79	nq	99
69) Pentachlorophenol	11.16	266	144440	30.71	nq	96
70) Phenanthrene	11.38	178	1520979	38.84	nq	97
71) Anthracene	11.44	178	1678099	43.71	nq	99
72) Carbazole	11.58	167	1392226	37.09	nq	100
73) Di-n-butylphthalate	11.93	149	1919987	49.30	nq	98
74) Fluoranthene	12.57	202	1446044	35.34	nq	99
76) Benzidine	12.69	184	735338	35.12	nq	99
77) Pyrene	12.80	202	1484029	32.38	nq	99
79) Butylbenzylphthalate	13.41	149	651181	32.84	nq	# 83
80) Benzo(a)anthracene	13.99	228	1243500	36.26	nq	99
81) 3,3'-Dichlorobenzidine	13.95	252	541437	45.24	nq	# 98
82) Chrysene	14.02	228	1187587	36.04	nq	97
83) Bis(2-ethylhexyl)phthalate	13.99	149	832229	33.22	nq	# 94
84) Di-n-octyl phthalate	14.59	149	1529785	36.17	nq	# 88
85) Indeno(1,2,3-cd)pyrene	16.80	276	1260380	48.25	nq	97
87) Benzo(b)fluoranthene	15.01	252	1495410m	39.22	nq	
88) Benzo(k)fluoranthene	15.04	252	1074590m	34.46	nq	
89) Benzo(a)pyrene	15.36	252	1273031	39.36	nq	# 96
90) Dibenzo(a,h)anthracene	16.82	278	1044967	37.55	nq	96
91) Benzo(g,h,i)perylene	17.22	276	1010297	35.05	nq	99

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

