

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF071615\
 Data File : BF080511.D
 Acq On : 16 Jul 2015 15:26
 Operator : TP/UM
 Sample : SSTDICV040
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 ICVBF071615

Quant Time: Jul 16 17:18:12 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF071615.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 16 17:11:01 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.07	152	17969	20.00	ng	0.00
21) Naphthalene-d8	8.36	136	78827	20.00	ng	0.00
38) Acenaphthene-d10	10.12	164	34534	20.00	ng	0.00
63) Phenanthrene-d10	11.61	188	64611	20.00	ng	0.00
75) Chrysene-d12	14.43	240	53199	20.00	ng	0.00
86) Perylene-d12	16.16	264	49712	20.00	ng	0.01

System Monitoring Compounds

5) 2-Fluorophenol	5.70	112	92268	87.54	ng	0.00
7) Phenol-d6	6.72	99	122402	85.51	ng	0.00
23) Nitrobenzene-d5	7.64	82	110879	82.57	ng	0.00
41) 2,4,6-Tribromophenol	10.91	330	24465	91.05	ng	0.00
44) 2-Fluorobiphenyl	9.44	172	217559	85.73	ng	0.00
78) Terphenyl-d14	13.23	244	222019	86.53	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.76	88	21827	41.93	ng	# 81
3) Pyridine	3.71	79	45789	38.71	ng	87
4) n-Nitrosodimethylamine	3.55	42	29802	45.23	ng	81
6) Aniline	6.75	93	77494	41.99	ng	# 90
8) 2-Chlorophenol	6.86	128	50783	42.95	ng	83
9) Benzaldehyde	6.62	77	42666	43.91	ng	# 68
10) Phenol	6.73	94	66274	43.04	ng	# 63
11) bis(2-Chloroethyl)ether	6.81	93	45768	40.83	ng	89
12) 1,3-Dichlorobenzene	7.01	146	61086	41.30	ng	# 92
13) 1,4-Dichlorobenzene	7.09	146	62182m	40.66	ng	
14) 1,2-Dichlorobenzene	7.25	146	54083	40.31	ng	97
15) Benzyl Alcohol	7.22	79	41926	41.96	ng	# 69
16) 2,2'-oxybis(1-Chloropropan	7.34	45	84983	42.27	ng	63
17) 2-Methylphenol	7.33	107	40150	42.42	ng	# 78
18) Hexachloroethane	7.60	117	18495	40.43	ng	# 64
19) n-Nitroso-di-n-propylamine	7.48	70	35410	40.76	ng	# 81
20) 3+4-Methylphenols	7.49	107	54618	42.71	ng	# 79
22) Acetophenone	7.48	105	73213	41.97	ng	# 78
24) Nitrobenzene	7.66	77	53927	39.78	ng	# 75
25) Isophorone	7.89	82	96222	40.57	ng	# 91
26) 2-Nitrophenol	7.98	139	23251	39.52	ng	# 81
27) 2,4-Dimethylphenol	8.02	122	45505	40.04	ng	# 80
28) bis(2-Chloroethoxy)methane	8.10	93	58798	42.21	ng	# 92
29) 2,4-Dichlorophenol	8.22	162	43284	41.79	ng	97
30) 1,2,4-Trichlorobenzene	8.30	180	44858	37.72	ng	97
31) Naphthalene	8.38	128	154278	39.85	ng	100
32) Benzoic acid	8.10	122	22076	37.74	ng	# 59
33) 4-Chloroaniline	8.44	127	59789	39.73	ng	# 89
34) Hexachlorobutadiene	8.50	225	28450	41.32	ng	98
35) Caprolactam	8.78	113	10004	40.87	ng	94
36) 4-Chloro-3-methylphenol	8.92	107	43977	40.20	ng	# 80
37) 2-Methylnaphthalene	9.07	142	95504	41.28	ng	# 90
39) 1,2,4,5-Tetrachlorobenzene	9.24	216	46318	41.82	ng	98
40) Hexachlorocyclopentadiene	9.23	237	22376	41.17	ng	96

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.36	196	30838	46.52	ng	99
43) 2,4,5-Trichlorophenol	9.40	196	30503	43.23	ng	96
45) 1,1'-Biphenyl	9.54	154	115069	40.77	ng	95
46) 2-Chloronaphthalene	9.56	162	90752	42.93	ng	# 92
47) 2-Nitroaniline	9.66	65	24588	45.13	ng	# 56
48) Acenaphthylene	9.98	152	149442	42.20	ng	98
49) Dimethylphthalate	9.84	163	98703	41.67	ng	# 96
50) 2,6-Dinitrotoluene	9.89	165	19119	41.25	ng	# 23
51) Acenaphthene	10.16	154	87944	42.03	ng	88
52) 3-Nitroaniline	10.08	138	22603	44.15	ng	# 42
53) 2,4-Dinitrophenol	10.18	184	8527	39.04	ng	# 60
54) Dibenzofuran	10.33	168	121888	40.29	ng	94
55) 4-Nitrophenol	10.25	139	16038	43.49	ng	# 29
56) 2,4-Dinitrotoluene	10.30	165	28012	41.71	ng	# 93
57) Fluorene	10.67	166	98545	41.49	ng	94
58) 2,3,4,6-Tetrachlorophenol	10.45	232	21748	42.24	ng	100
59) Diethylphthalate	10.53	149	92736	41.51	ng	97
60) 4-Chlorophenyl-phenylether	10.66	204	43151	40.60	ng	94
61) 4-Nitroaniline	10.69	138	23292	45.95	ng	# 46
62) Azobenzene	10.81	77	101808	43.38	ng	# 80
64) 4,6-Dinitro-2-methylphenol	10.72	198	11840	37.60	ng	# 69
65) n-Nitrosodiphenylamine	10.77	169	89961	40.94	ng	92
66) 4-Bromophenyl-phenylether	11.15	248	24436	38.83	ng	# 81
67) Hexachlorobenzene	11.22	284	29042	41.74	ng	# 83
68) Atrazine	11.29	200	22068	37.09	ng	83
69) Pentachlorophenol	11.41	266	12001	35.65	ng	96
70) Phenanthrene	11.63	178	150411	41.34	ng	97
71) Anthracene	11.69	178	140356	40.68	ng	99
72) Carbazole	11.85	167	118255	39.08	ng	97
73) Di-n-butylphthalate	12.16	149	146961	41.44	ng	# 98
74) Fluoranthene	12.84	202	136435	40.04	ng	95
76) Benzidine	12.98	184	71709	40.93	ng	# 97
77) Pyrene	13.09	202	144030	41.52	ng	97
79) Butylbenzylphthalate	13.75	149	48191	38.80	ng	# 66
80) Benzo(a)anthracene	14.42	228	123435	41.35	ng	98
81) 3,3'-Dichlorobenzidine	14.39	252	38782	39.73	ng	# 93
82) Chrysene	14.47	228	122942	42.34	ng	96
83) Bis(2-ethylhexyl)phthalate	14.39	149	75797	41.74	ng	# 87
84) Di-n-octyl phthalate	15.14	149	112290	38.40	ng	99
85) Indeno(1,2,3-cd)pyrene	17.75	276	162989	45.83	ng	# 93
87) Benzo(b)fluoranthene	15.68	252	129433	44.99	ng	# 90
88) Benzo(k)fluoranthene	15.71	252	103391m	35.35	ng	
89) Benzo(a)pyrene	16.08	252	112905	41.97	ng	# 93
90) Dibenzo(a,h)anthracene	17.76	278	131481	44.40	ng	# 85
91) Benzo(g,h,i)perylene	18.23	276	137446	44.69	ng	# 79

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

