

Data Path : Z:\HPCHEM1\BNA F\DATA\BF042315\
 Data File : BF078757.D
 Acq On : 23 Apr 2015 19:35
 Operator : TP/IZ
 Sample : G1991-03MS
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-75DMS

Quant Time: Apr 24 11:20:42 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF040115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 24 00:49:54 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	5.76	152	26187	20.00	ng	0.01
21) Naphthalene-d8	7.92	136	108520	20.00	ng	0.01
38) Acenaphthene-d10	11.10	164	55575	20.00	ng	0.00
63) Phenanthrene-d10	13.83	188	108362	20.00	ng	0.00
75) Chrysene-d12	17.71	240	118327	20.00	ng	0.00
86) Perylene-d12	19.42	264	120515	20.00	ng	-0.08

System Monitoring Compounds						
5) 2-Fluorophenol	3.77	112	198183	124.78	ng	0.00
7) Phenol-d6	5.27	99	262952	132.11	ng	0.01
23) Nitrobenzene-d5	6.70	82	150981	84.33	ng	0.00
41) 2,4,6-Tribromophenol	12.57	330	65279	141.63	ng	0.00
44) 2-Fluorobiphenyl	9.93	172	287448	73.93	ng	0.00
78) Terphenyl-d14	16.38	244	317005	60.54	ng	0.00

Target Compounds						Qvalue
2) 1,4-Dioxane	1.34	88	23365	32.53	ng	99
3) Pyridine	1.83	79	59429	31.59	ng	98
4) n-Nitrosodimethylamine	1.78	42	30816	42.55	ng	92
6) Aniline	5.24	93	61597	21.82	ng	# 86
8) 2-Chlorophenol	5.42	128	79646	42.41	ng	98
9) Benzaldehyde	5.06	77	8643	6.55	ng	95
10) Phenol	5.28	94	99292	41.63	ng	90
11) bis(2-Chloroethyl)ether	5.39	93	68717	40.07	ng	99
12) 1,3-Dichlorobenzene	5.66	146	79360	38.47	ng	# 93
13) 1,4-Dichlorobenzene	5.78	146	81663	39.33	ng	96
14) 1,2-Dichlorobenzene	6.01	146	77549	39.83	ng	99
15) Benzyl Alcohol	6.03	79	66335	47.42	ng	95
16) 2,2'-oxybis(1-Chloropropan	6.27	45	92030	40.23	ng	64
17) 2-Methylphenol	6.26	107	61794	42.38	ng	95
18) Hexachloroethane	6.56	117	26990	38.54	ng	# 80
19) n-Nitroso-di-n-propylamine	6.49	70	54827	41.75	ng	92
20) 3+4-Methylphenols	6.54	107	81995	42.15	ng	96
22) Acetophenone	6.46	105	110028	43.52	ng	# 92
24) Nitrobenzene	6.73	77	79850	42.66	ng	89
25) Isophorone	7.15	82	143146	42.13	ng	# 92
26) 2-Nitrophenol	7.28	139	38175	42.80	ng	95
27) 2,4-Dimethylphenol	7.44	122	69491	41.90	ng	97
28) bis(2-Chloroethoxy)methane	7.60	93	89087	40.89	ng	99
29) 2,4-Dichlorophenol	7.72	162	66158	43.85	ng	96
30) 1,2,4-Trichlorobenzene	7.84	180	66477	38.43	ng	98
31) Naphthalene	7.95	128	221657	39.24	ng	99
32) Benzoic acid	7.63	122	7342m	14.07	ng	
33) 4-Chloroaniline	8.11	127	57554	24.56	ng	100
34) Hexachlorobutadiene	8.20	225	36810	38.75	ng	96
35) Caprolactam	8.74	113	20555	45.64	ng	# 85
36) 4-Chloro-3-methylphenol	9.07	107	68778	45.18	ng	94
37) 2-Methylnaphthalene	9.21	142	152880	39.87	ng	98
39) 1,2,4,5-Tetrachlorobenzene	9.52	216	65432	38.00	ng	99
40) Hexachlorocyclopentadiene	9.51	237	28665	42.19	ng	97

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.77	196	45716	42.10	nq	97
43) 2,4,5-Trichlorophenol	9.85	196	47393	41.77	nq	# 92
45) 1,1'-Biphenyl	10.09	154	178427	39.25	nq	98
46) 2-Chloronaphthalene	10.09	162	137867	38.55	nq	97
47) 2-Nitroaniline	10.33	65	40873	46.19	nq	89
48) Acenaphthylene	10.83	152	227509	39.39	nq	100
49) Dimethylphthalate	10.74	163	190788	45.70	nq	98
50) 2,6-Dinitrotoluene	10.83	165	36821	42.86	nq	# 53
51) Acenaphthene	11.15	154	130428	40.11	nq	100
52) 3-Nitroaniline	11.11	138	35988	36.84	nq	99
53) 2,4-Dinitrophenol	11.34	184	6039	29.42	nq	89
54) Dibenzofuran	11.48	168	206869	41.65	nq	98
55) 4-Nitrophenol	11.54	139	52711	73.52	nq	85
56) 2,4-Dinitrotoluene	11.57	165	50006	44.95	nq	# 80
57) Fluorene	12.11	166	168996	41.98	nq	100
58) 2,3,4,6-Tetrachlorophenol	11.76	232	35415	42.10	nq	97
59) Diethylphthalate	12.06	149	169941	42.21	nq	97
60) 4-Chlorophenyl-phenylether	12.18	204	75415	40.72	nq	# 88
61) 4-Nitroaniline	12.23	138	35911	36.37	nq	95
62) Azobenzene	12.47	77	156891	42.70	nq	93
64) 4,6-Dinitro-2-methylphenol	12.32	198	11432	26.02	nq	# 65
65) n-Nitrosodiphenylamine	12.41	169	145217	38.74	nq	99
66) 4-Bromophenyl-phenylether	13.07	248	46307	40.35	nq	93
67) Hexachlorobenzene	13.11	284	46921	37.31	nq	# 87
68) Atrazine	13.50	200	55681	51.29	nq	99
69) Pentachlorophenol	13.53	266	35357	65.01	nq	95
70) Phenanthrene	13.87	178	251628	40.14	nq	100
71) Anthracene	13.97	178	249354	40.37	nq	99
72) Carbazole	14.31	167	248380	42.91	nq	99
73) Di-n-butylphthalate	14.99	149	274621	41.88	nq	100
74) Fluoranthene	15.77	202	275912	41.74	nq	95
76) Benzidine	16.03	184	164210	36.04	nq	98
77) Pyrene	16.08	202	285668	38.46	nq	98
79) Butylbenzylphthalate	17.07	149	124941	44.13	nq	99
80) Benzo(a)anthracene	17.70	228	265254	37.68	nq	98
81) 3,3'-Dichlorobenzidine	17.71	252	89210	37.83	nq	100
82) Chrysene	17.75	228	256323	38.75	nq	99
83) Bis(2-ethylhexyl)phthalate	17.85	149	182549	41.11	nq	99
84) Di-n-octyl phthalate	18.63	149	329184	45.15	nq	# 100
85) Indeno(1,2,3-cd)pyrene	20.57	276	300625m	40.05	nq	
87) Benzo(b)fluoranthene	18.99	252	276416	37.11	nq	99
88) Benzo(k)fluoranthene	19.03	252	255330m	38.58	nq	
89) Benzo(a)pyrene	19.35	252	259495	39.92	nq	# 95
90) Dibenzo(a,h)anthracene	20.59	278	252600m	38.81	nq	
91) Benzo(g,h,i)perylene	20.88	276	249765m	38.28	nq	

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

