

Data Path : Z:\HPCHEM1\BNA G\DATA\BG021418\
 Data File : BG032700.D
 Acq On : 15 Feb 2018 4:38
 Operator : SJ/JU
 Sample : J1530-41MS
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 WC-01MS

Manual Integrations
 APPROVED

Sohil
 2/15/2018 11:07:59 AM

Quant Time: Feb 15 06:38:03 2018
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG021218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Feb 12 18:08:01 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.66	152	28110	20.00	ng	-0.02
21) Naphthalene-d8	11.53	136	112433	20.00	ng	-0.01
38) Acenaphthene-d10	15.30	164	95282	20.00	ng	0.00
63) Phenanthrene-d10	18.03	188	240317	20.00	ng	-0.01
75) Chrysene-d12	22.36	240	307500	20.00	ng	-0.01
86) Perylene-d12	25.99	264	362549	20.00	ng	-0.03

System Monitoring Compounds

5) 2-Fluorophenol	6.10	112	159797	114.06	ng	-0.01
7) Phenol-d6	7.78	99	178816	93.34	ng	-0.01
23) Nitrobenzene-d5	9.85	82	166937	94.36	ng	-0.02
41) 2,4,6-Tribromophenol	16.78	330	236712	161.34	ng	-0.01
44) 2-Fluorobiphenyl	13.93	172	610385	80.11	ng	0.00
78) Terphenyl-d14	20.58	244	1358818	100.72	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.77	88	14539	24.807	ng	# 76
3) Pyridine	4.23	79	28090	18.685	ng	# 80
4) n-Nitrosodimethylamine	4.13	42	25091	35.520	ng	# 75
6) Aniline	7.96	93	24272	10.720	ng	# 93
8) 2-Chlorophenol	8.21	128	76958	47.768	ng	# 80
9) Benzaldehyde	7.76	77	8962	7.916	ng	# 94
10) Phenol	7.81	94	67765	36.370	ng	# 87
11) bis(2-Chloroethyl)ether	8.05	93	54297	36.350	ng	# 78
12) 1,3-Dichlorobenzene	8.54	146	91980	40.480	ng	# 94
13) 1,4-Dichlorobenzene	8.69	146	84617	37.693	ng	# 97
14) 1,2-Dichlorobenzene	9.03	146	76326	33.710	ng	# 96
15) Benzyl Alcohol	8.90	79	67320	46.882	ng	# 92
16) 2,2'-oxybis(1-Chloropropan	9.18	45	47528	29.564	ng	# 71
17) 2-Methylphenol	9.10	107	52571	39.840	ng	# 90
18) Hexachloroethane	9.77	117	34239	41.771	ng	# 80
19) n-Nitroso-di-n-propylamine	9.47	70	49104	38.126	ng	# 91
20) 3+4-Methylphenols	9.44	107	78105	39.163	ng	# 93
22) Acetophenone	9.50	105	111031	46.433	ng	# 93
24) Nitrobenzene	9.90	77	81628	44.069	ng	# 90
25) Isophorone	10.42	82	118764	35.212	ng	# 80
26) 2-Nitrophenol	10.62	139	48429m	48.885	ng	# 95
27) 2,4-Dimethylphenol	10.67	122	79500	61.081	ng	# 97
28) bis(2-Chloroethoxy)methane	10.90	93	89237	51.755	ng	# 97
29) 2,4-Dichlorophenol	11.17	162	105164	55.205	ng	# 90
30) 1,2,4-Trichlorobenzene	11.39	180	122037	45.695	ng	# 98
31) Naphthalene	11.58	128	341977	62.640	ng	# 99
32) Benzoic acid	10.80	122	30066	30.578	ng	# 81
33) 4-Chloroaniline	11.69	127	23177	10.699	ng	# 87
34) Hexachlorobutadiene	11.86	225	95837	43.786	ng	# 96
35) Caprolactam	12.45	113	20851	38.544	ng	# 45
36) 4-Chloro-3-methylphenol	12.77	107	95899	55.966	ng	# 87
37) 2-Methylnaphthalene	13.16	142	237054	53.532	ng	# 93
39) 1,2,4,5-Tetrachlorobenzene	13.51	216	183044	40.674	ng	# 97
40) Hexachlorocyclopentadiene	13.49	237	226544	79.547	ng	# 93

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.74	196	107510	47.864	nq	97
43) 2,4,5-Trichlorophenol	13.82	196	120295	42.976	nq #	88
45) 1,1'-Biphenyl	14.14	154	314953	40.173	nq	94
46) 2-Chloronaphthalene	14.19	162	255424	41.172	nq	98
47) 2-Nitroaniline	14.38	65	66480	49.264	nq #	82
48) Acenaphthylene	15.02	152	415834	45.560	nq	99
49) Dimethylphthalate	14.73	163	379660	46.097	nq #	98
50) 2,6-Dinitrotoluene	14.86	165	88043	51.905	nq #	78
51) Acenaphthene	15.36	154	318434	51.087	nq	98
52) 3-Nitroaniline	15.20	138	26359	17.597	nq #	83
53) 2,4-Dinitrophenol	15.39	184	70825	77.770	nq #	63
54) Dibenzofuran	15.69	168	471336	50.825	nq	96
55) 4-Nitrophenol	15.49	139	92291	82.864	nq #	79
56) 2,4-Dinitrotoluene	15.64	165	130060	55.649	nq #	71
57) Fluorene	16.33	166	363586	47.674	nq	100
58) 2,3,4,6-Tetrachlorophenol	15.91	232	139844	53.064	nq #	86
59) Diethylphthalate	16.08	149	386056	48.173	nq	96
60) 4-Chlorophenyl-phenylether	16.31	204	235560	46.781	nq	97
61) 4-Nitroaniline	16.35	138	59328	38.417	nq #	75
62) Azobenzene	16.61	77	240807	46.770	nq	84
64) 4,6-Dinitro-2-methylphenol	16.40	198	65693	38.672	nq	74
65) n-Nitrosodiphenylamine	16.53	169	342159	48.547	nq	99
66) 4-Bromophenyl-phenylether	17.20	248	163875	47.648	nq	98
67) Hexachlorobenzene	17.33	284	169626	45.920	nq #	91
68) Atrazine	17.46	200	153391	50.741	nq	93
69) Pentachlorophenol	17.67	266	221471	97.934	nq	99
70) Phenanthrene	18.07	178	625708	48.469	nq	99
71) Anthracene	18.17	178	658425	49.847	nq	99
72) Carbazole	18.43	167	570562	50.662	nq	98
73) Di-n-butylphthalate	18.94	149	585997	46.212	nq	99
74) Fluoranthene	20.04	202	869568	50.456	nq	95
76) Benzidine	20.21	184	88585	12.234	nq	98
77) Pyrene	20.40	202	863948	47.226	nq	98
79) Butylbenzylphthalate	21.26	149	310695	51.556	nq #	78
80) Benzo(a)anthracene	22.33	228	931739	48.394	nq	99
81) 3,3'-Dichlorobenzidine	22.23	252	176009	24.479	nq #	97
82) Chrysene	22.41	228	893911	47.087	nq	98
83) Bis(2-ethylhexyl)phthalate	22.18	149	373641	42.995	nq #	99
84) Di-n-octyl phthalate	23.53	149	663732	48.680	nq #	89
85) Indeno(1,2,3-cd)pyrene	30.20	276	1277050	56.785	nq #	84
87) Benzo(b)fluoranthene	24.83	252	933699	43.215	nq #	96
88) Benzo(k)fluoranthene	24.91	252	986788	44.130	nq #	96
89) Benzo(a)pyrene	25.82	252	1075099	51.077	nq #	95
90) Dibenzo(a,h)anthracene	30.28	278	1073260	51.217	nq #	95
91) Benzo(g,h,i)perylene	31.53	276	1057903	52.225	nq #	92

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

