

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG050319\
 Data File : BG040731.D
 Acq On : 3 May 2019 19:08
 Operator : JU/SJ
 Sample : K2635-02MSD
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 AU-05-050219-AMSD

Quant Time: May 03 23:17:51 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG042319.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 24 05:35:33 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.15	152	37061	20.00	ng	-0.01
21) Naphthalene-d8	10.97	136	152084	20.00	ng	-0.01
39) Acenaphthene-d10	14.77	164	117239	20.00	ng	-0.02
64) Phenanthrene-d10	17.52	188	328790	20.00	ng	-0.02
76) Chrysene-d12	21.81	240	332800	20.00	ng	-0.02
87) Perylene-d12	25.16	264	358690	20.00	ng	-0.04

System Monitoring Compounds

5) 2-Fluorophenol	5.68	112	349349	166.16	ng	0.00
7) Phenol-d6	7.29	99	480142	148.32	ng	0.00
23) Nitrobenzene-d5	9.32	82	389746	118.41	ng	-0.02
42) 2,4,6-Tribromophenol	16.26	330	325275	201.85	ng	-0.01
45) 2-Fluorobiphenyl	13.40	172	825510	101.69	ng	-0.02
79) Terphenyl-d14	20.12	244	1756077	116.90	ng	-0.02

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.55	88	45247	47.322	ng	# 21
3) Pyridine	3.96	79	122392	44.162	ng	98
4) n-Nitrosodimethylamine	3.88	42	75177	48.904	ng	96
6) Aniline	7.47	93	137634	32.077	ng	98
8) 2-Chlorophenol	7.71	128	137690	53.635	ng	97
9) Benzaldehyde	7.28	77	91428	50.527	ng	93
10) Phenol	7.31	94	167957	48.267	ng	97
11) bis(2-Chloroethyl)ether	7.57	93	129456	49.611	ng	98
12) 1,3-Dichlorobenzene	8.03	146	108862	38.851	ng	97
13) 1,4-Dichlorobenzene	8.18	146	115155	40.540	ng	97
14) 1,2-Dichlorobenzene	8.50	146	111560	40.725	ng	92
15) Benzyl Alcohol	8.38	79	172726	51.280	ng	97
16) 2,2'-oxybis(1-Chloropropan	8.67	45	227994	43.885	ng	96
17) 2-Methylphenol	8.58	107	129833	52.722	ng	95
18) Hexachloroethane	9.23	117	39957	34.292	ng	94
19) n-Nitroso-di-n-propylamine	8.95	70	135037	46.341	ng	99
20) 3+4-Methylphenols	8.91	107	182079	53.075	ng	97
22) Acetophenone	8.98	105	241803	57.176	ng	# 95
24) Nitrobenzene	9.37	77	207250	59.016	ng	97
25) Isophorone	9.89	82	383490	52.306	ng	97
26) 2-Nitrophenol	10.07	139	81005	64.350	ng	97
27) 2,4-Dimethylphenol	10.12	122	154519	65.422	ng	99
28) bis(2-Chloroethoxy)methane	10.36	93	192621	52.365	ng	99
29) 2,4-Dichlorophenol	10.61	162	170537	63.511	ng	95
30) 1,2,4-Trichlorobenzene	10.83	180	142881	47.984	ng	97
31) Naphthalene	11.02	128	398646	49.339	ng	98
32) Benzoic acid	10.23	122	88656	54.831	ng	98
33) 4-Chloroaniline	11.13	127	40279	11.244	ng	93
34) Hexachlorobutadiene	11.28	225	96840	44.052	ng	98
35) Caprolactam	11.91	113	49027	46.247	ng	94
36) 4-Chloro-3-methylphenol	12.22	107	204922	60.496	ng	98
37) 2-Methylnaphthalene	12.61	142	325034	53.804	ng	98
38) 1-Methylnaphthalene	12.83	142	310981	52.666	ng	98
40) 1,2,4,5-Tetrachlorobenzene	12.97	216	227785	52.155	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.94	237	202078	78.411	ng	97
43) 2,4,6-Trichlorophenol	13.21	196	162139	61.433	ng	99
44) 2,4,5-Trichlorophenol	13.28	196	181714	63.202	ng	95
46) 1,1'-Biphenyl	13.61	154	461113	50.658	ng	98
47) 2-Chloronaphthalene	13.66	162	374774	52.610	ng	99
48) 2-Nitroaniline	13.86	65	165116	65.116	ng	97
49) Acenaphthylene	14.50	152	624261	51.652	ng	99
50) Dimethylphthalate	14.22	163	555637	56.645	ng	99
51) 2,6-Dinitrotoluene	14.35	165	125309	65.454	ng	92
52) Acenaphthene	14.84	154	367328	52.189	ng	96
53) 3-Nitroaniline	14.68	138	77593	39.557	ng	# 98
54) 2,4-Dinitrophenol	14.88	184	82249	76.511	ng	# 83
55) Dibenzofuran	15.17	168	641401	55.636	ng	99
56) 4-Nitrophenol	14.97	139	186657	109.740	ng	91
57) 2,4-Dinitrotoluene	15.13	165	184469	70.911	ng	95
58) Fluorene	15.82	166	532322	57.128	ng	98
59) 2,3,4,6-Tetrachlorophenol	15.39	232	200041	69.127	ng	96
60) Diethylphthalate	15.58	149	591786	57.180	ng	100
61) 4-Chlorophenyl-phenylether	15.81	204	314919	58.110	ng	96
62) 4-Nitroaniline	15.85	138	131347	59.860	ng	95
63) Azobenzene	16.10	77	574832	53.932	ng	96
65) 4,6-Dinitro-2-methylphenol	15.89	198	79612	48.973	ng	89
66) n-Nitrosodiphenylamine	16.03	169	493588	51.026	ng	98
67) 4-Bromophenyl-phenylether	16.70	248	214947	52.474	ng	96
68) Hexachlorobenzene	16.81	284	224243	52.943	ng	95
69) Atrazine	16.96	200	225774	58.910	ng	96
70) Pentachlorophenol	17.16	266	251853	101.615	ng	99
71) Phenanthrene	17.57	178	960392	54.230	ng	98
72) Anthracene	17.65	178	983683	55.260	ng	100
73) Carbazole	17.92	167	868642	53.560	ng	99
74) Di-n-butylphthalate	18.46	149	1050160	53.648	ng	99
75) Fluoranthene	19.56	202	1254705	56.854	ng	100
77) Benzidine	19.74	184	173650	19.057	ng	97
78) Pyrene	19.93	202	1255892	60.210	ng	98
80) Butylbenzylphthalate	20.80	149	485140	59.675	ng	92
81) Benzo(a)anthracene	21.78	228	1207580	57.413	ng	99
82) 3,3'-Dichlorobenzidine	21.70	252	258133	34.432	ng	99
83) Chrysene	21.85	228	1183784	59.180	ng	98
84) Bis(2-ethylhexyl)phthalate	21.66	149	657209	56.003	ng	96
85) Di-n-octyl phthalate	22.92	149	1126199	56.983	ng	98
86) Indeno(1,2,3-cd)pyrene	29.02	276	1285804	53.166	ng	# 89
88) Benzo(b)fluoranthene	24.09	252	1227049	55.981	ng	97
89) Benzo(k)fluoranthene	24.16	252	1198129	58.530	ng	98
90) Benzo(a)pyrene	25.00	252	1184591	57.094	ng	98
91) Dibenzo(a,h)anthracene	29.09	278	1062089	50.584	ng	# 98
92) Benzo(g,h,i)perylene	30.24	276	1045747	50.044	ng	97

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

