

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_G\DATA\BG032618\
 Data File : BG033340.D
 Acq On : 26 Mar 2018 21:48
 Operator : SJ/JU
 Sample : J1867-14MSD
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C19013071MSD

Manual Integrations
 APPROVED

Sohil
 3/27/2018 2:57:48 PM

Quant Time: Mar 27 08:19:16 2018
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG031918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 23 15:11:46 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.87	152	41423	20.00	ng	-0.03
21) Naphthalene-d8	11.75	136	168245	20.00	ng	-0.03
38) Acenaphthene-d10	15.48	164	128672	20.00	ng	-0.03
63) Phenanthrene-d10	18.21	188	331345	20.00	ng	-0.03
75) Chrysene-d12	22.56	240	347377	20.00	ng	-0.04
86) Perylene-d12	26.32	264	376180	20.00	ng	-0.06

System Monitoring Compounds

5) 2-Fluorophenol	6.28	112	192108	86.96	ng	-0.03
7) Phenol-d6	7.97	99	184162	48.52	ng	-0.03
23) Nitrobenzene-d5	10.07	82	398986	108.95	ng	-0.03
41) 2,4,6-Tribromophenol	16.96	330	319679	166.12	ng	-0.03
44) 2-Fluorobiphenyl	14.11	172	966644	108.45	ng	-0.03
78) Terphenyl-d14	20.73	244	1612118	99.25	ng	-0.03

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.93	88	21720	19.361	ng	86
3) Pyridine	4.39	79	47312m	15.256	ng	
4) n-Nitrosodimethylamine	4.30	42	27417	21.543	ng	# 90
6) Aniline	8.16	93	76276	17.088	ng	97
8) 2-Chlorophenol	8.41	128	111544	44.168	ng	94
9) Benzaldehyde	7.96	77	72598m	30.097	ng	
10) Phenol	8.01	94	66287	17.688	ng	94
11) bis(2-Chloroethyl)ether	8.25	93	141952	46.407	ng	96
12) 1,3-Dichlorobenzene	8.75	146	135010	40.890	ng	96
13) 1,4-Dichlorobenzene	8.90	146	133870	40.485	ng	93
14) 1,2-Dichlorobenzene	9.23	146	133023	41.218	ng	97
15) Benzyl Alcohol	9.10	79	98540	32.974	ng	89
16) 2,2'-oxybis(1-Chloropropan	9.39	45	217443	43.606	ng	85
17) 2-Methylphenol	9.30	107	90742m	35.545	ng	
18) Hexachloroethane	9.99	117	51066	40.013	ng	90
19) n-Nitroso-di-n-propylamine	9.67	70	128885	46.340	ng	# 96
20) 3+4-Methylphenols	9.64	107	113488	31.222	ng	85
22) Acetophenone	9.71	105	231738	50.276	ng	# 96
24) Nitrobenzene	10.11	77	189379	48.316	ng	91
25) Isophorone	10.63	82	370983	52.197	ng	97
26) 2-Nitrophenol	10.84	139	81771	55.104	ng	# 95
27) 2,4-Dimethylphenol	10.87	122	115288	53.480	ng	89
28) bis(2-Chloroethoxy)methane	11.11	93	197172	55.601	ng	94
29) 2,4-Dichlorophenol	11.38	162	148538	56.028	ng	95
30) 1,2,4-Trichlorobenzene	11.60	180	161732	46.954	ng	98
31) Naphthalene	11.80	128	429424	49.254	ng	98
32) Benzoic acid	10.95	122	30628m	15.284	ng	
33) 4-Chloroaniline	11.89	127	93731	23.996	ng	93
34) Hexachlorobutadiene	12.06	225	119189	45.758	ng	94
35) Caprolactam	12.63	113	24837	23.914	ng	98
36) 4-Chloro-3-methylphenol	12.96	107	171159	49.500	ng	99
37) 2-Methylnaphthalene	13.35	142	343881	50.817	ng	90
39) 1,2,4,5-Tetrachlorobenzene	13.70	216	230714	49.500	ng	97
40) Hexachlorocyclopentadiene	13.68	237	263370	96.391	ng	90

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.93	196	152927	56.473	ng	97
43) 2,4,5-Trichlorophenol	14.00	196	173097	54.079	ng	98
45) 1,1'-Biphenyl	14.33	154	478801	48.434	ng	99
46) 2-Chloronaphthalene	14.38	162	376772	49.430	ng	97
47) 2-Nitroaniline	14.57	65	148159	51.895	ng	97
48) Acenaphthylene	15.21	152	601516	47.278	ng	98
49) Dimethylphthalate	14.91	163	566377	50.579	ng	100
50) 2,6-Dinitrotoluene	15.04	165	122743	53.607	ng	97
51) Acenaphthene	15.55	154	456553	55.665	ng	95
52) 3-Nitroaniline	15.38	138	71291	29.699	ng	# 89
53) 2,4-Dinitrophenol	15.58	184	118855	98.305	ng	# 71
54) Dibenzofuran	15.88	168	631055	52.089	ng	92
55) 4-Nitrophenol	15.66	139	76310	45.208	ng	# 87
56) 2,4-Dinitrotoluene	15.82	165	172686	51.222	ng	91
57) Fluorene	16.52	166	534379	51.775	ng	99
58) 2,3,4,6-Tetrachlorophenol	16.09	232	177299m	58.839	ng	
59) Diethylphthalate	16.25	149	540604	49.718	ng	96
60) 4-Chlorophenyl-phenylether	16.50	204	295513	49.808	ng	98
61) 4-Nitroaniline	16.54	138	111065	45.689	ng	# 83
62) Azobenzene	16.79	77	517936	49.172	ng	97
64) 4,6-Dinitro-2-methylphenol	16.58	198	97044	48.958	ng	95
65) n-Nitrosodiphenylamine	16.71	169	462878	48.933	ng	96
66) 4-Bromophenyl-phenylether	17.39	248	198195	50.932	ng	98
67) Hexachlorobenzene	17.52	284	210680	51.300	ng	95
68) Atrazine	17.63	200	198219	51.712	ng	97
69) Pentachlorophenol	17.85	266	271988	96.494	ng	99
70) Phenanthrene	18.26	178	860415	49.861	ng	99
71) Anthracene	18.35	178	879827	50.355	ng	99
72) Carbazole	18.60	167	753803	48.685	ng	97
73) Di-n-butylphthalate	19.10	149	892766	49.538	ng	# 99
74) Fluoranthene	20.21	202	1054654	49.388	ng	96
76) Benzidine	20.37	184	177936	18.888	ng	98
77) Pyrene	20.57	202	1071041	46.229	ng	99
79) Butylbenzylphthalate	21.42	149	417148	48.792	ng	97
80) Benzo(a)anthracene	22.54	228	1086272	49.109	ng	100
81) 3,3'-Dichlorobenzidine	22.42	252	276368	35.111	ng	99
82) Chrysene	22.62	228	1033751	48.280	ng	97
83) Bis(2-ethylhexyl)phthalate	22.36	149	586265	49.744	ng	98
84) Di-n-octyl phthalate	23.76	149	1003431	55.607	ng	97
85) Indeno(1,2,3-cd)pyrene	30.71	276	1267253	54.741	ng	# 88
87) Benzo(b)fluoranthene	25.11	252	1100086	50.617	ng	# 96
88) Benzo(k)fluoranthene	25.20	252	1085137	49.367	ng	# 98
89) Benzo(a)pyrene	26.15	252	1072404	51.381	ng	# 96
90) Dibenzo(a,h)anthracene	30.78	278	1065556	54.199	ng	# 94
91) Benzo(g,h,i)perylene	32.09	276	1058953	54.394	ng	# 94

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

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