

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG111119\
 Data File : BG043585.D
 Acq On : 11 Nov 2019 16:07
 Operator : JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040

Quant Time: Nov 12 00:46:20 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG102119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 04 15:22:15 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.00	152	38417	20.00	ng	0.00
21) Naphthalene-d8	10.81	136	149812	20.00	ng	0.00
39) Acenaphthene-d10	14.63	164	102343	20.00	ng	0.00
64) Phenanthrene-d10	17.38	188	246745	20.00	ng	0.00
76) Chrysene-d12	21.65	240	263822	20.00	ng	0.00
87) Perylene-d12	24.83	264	310871	20.00	ng	-0.02

System Monitoring Compounds

5) 2-Fluorophenol	5.60	112	180584	86.14	ng	0.00
7) Phenol-d6	7.20	99	250061	82.90	ng	0.00
23) Nitrobenzene-d5	9.17	82	238874	82.20	ng	0.00
42) 2,4,6-Tribromophenol	16.13	330	156588	80.40	ng	0.00
45) 2-Fluorobiphenyl	13.26	172	622467	84.04	ng	0.00
79) Terphenyl-d14	19.99	244	1160524	82.53	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.47	88	33867	41.454	ng	92
3) Pyridine	3.87	79	91740	44.515	ng	94
4) n-Nitrosodimethylamine	3.79	42	46540	44.753	ng	# 89
6) Aniline	7.33	93	143426	41.277	ng	98
8) 2-Chlorophenol	7.58	128	93869	40.561	ng	98
9) Benzaldehyde	7.14	77	60705	40.952	ng	86
10) Phenol	7.23	94	112323	40.751	ng	97
11) bis(2-Chloroethyl)ether	7.42	93	84010	39.422	ng	94
12) 1,3-Dichlorobenzene	7.89	146	120160	41.440	ng	97
13) 1,4-Dichlorobenzene	8.04	146	119702	40.803	ng	96
14) 1,2-Dichlorobenzene	8.36	146	114342	39.918	ng	97
15) Benzyl Alcohol	8.25	79	89333	42.170	ng	99
16) 2,2'-oxybis(1-Chloropropan	8.53	45	121164	39.102	ng	96
17) 2-Methylphenol	8.47	107	81631	40.625	ng	94
18) Hexachloroethane	9.08	117	43582	41.176	ng	93
19) n-Nitroso-di-n-propylamine	8.80	70	73427	37.672	ng	# 90
20) 3+4-Methylphenols	8.80	107	111906	40.614	ng	98
22) Acetophenone	8.83	105	155858	40.625	ng	# 92
24) Nitrobenzene	9.21	77	112301	40.568	ng	# 96
25) Isophorone	9.73	82	206318	38.834	ng	98
26) 2-Nitrophenol	9.92	139	56109	41.984	ng	91
27) 2,4-Dimethylphenol	10.00	122	79278	41.388	ng	95
28) bis(2-Chloroethoxy)methane	10.21	93	119895	40.889	ng	98
29) 2,4-Dichlorophenol	10.48	162	101535	41.371	ng	94
30) 1,2,4-Trichlorobenzene	10.67	180	124932	40.389	ng	96
31) Naphthalene	10.85	128	308080	40.514	ng	98
32) Benzoic acid	10.18	122	53444	38.701	ng	93
33) 4-Chloroaniline	10.97	127	130732	41.940	ng	97
34) Hexachlorobutadiene	11.14	225	92586	40.491	ng	97
35) Caprolactam	11.75	113	31896	37.735	ng	# 89
36) 4-Chloro-3-methylphenol	12.12	107	106014	39.601	ng	97
37) 2-Methylnaphthalene	12.46	142	228808	39.215	ng	97
38) 1-Methylnaphthalene	12.68	142	219910	39.612	ng	95
40) 1,2,4,5-Tetrachlorobenzene	12.83	216	163857	43.262	ng	99

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41) Hexachlorocyclopentadiene	12.81	237	150941	75.864	nq	97
43) 2,4,6-Trichlorophenol	13.08	196	91822	41.166	nq	96
44) 2,4,5-Trichlorophenol	13.17	196	91839	40.726	nq	96
46) 1,1'-Biphenyl	13.47	154	326828	41.978	nq	96
47) 2-Chloronaphthalene	13.51	162	247363	41.757	nq	99
48) 2-Nitroaniline	13.72	65	76297	43.093	nq	96
49) Acenaphthylene	14.36	152	383947	41.644	nq	99
50) Dimethylphthalate	14.09	163	326755	41.220	nq	99
51) 2,6-Dinitrotoluene	14.21	165	71702	41.635	nq	96
52) Acenaphthene	14.70	154	231533	41.180	nq	97
53) 3-Nitroaniline	14.55	138	70991	42.696	nq	98
54) 2,4-Dinitrophenol	14.76	184	55101	53.856	nq	95
55) Dibenzofuran	15.03	168	377616	41.415	nq	98
56) 4-Nitrophenol	14.90	139	45617	40.940	nq	94
57) 2,4-Dinitrotoluene	15.00	165	101724	41.332	nq	97
58) Fluorene	15.68	166	316497	41.387	nq	97
59) 2,3,4,6-Tetrachlorophenol	15.27	232	91523	38.200	nq	93
60) Diethylphthalate	15.45	149	321557	40.371	nq	99
61) 4-Chlorophenyl-phenylether	15.67	204	181446	40.672	nq	96
62) 4-Nitroaniline	15.72	138	75411	43.916	nq	96
63) Azobenzene	15.97	77	266304	41.311	nq	96
65) 4,6-Dinitro-2-methylphenol	15.77	198	56880	39.171	nq	94
66) n-Nitrosodiphenylamine	15.89	169	277106	39.778	nq	98
67) 4-Bromophenyl-phenylether	16.57	248	132420	39.878	nq	98
68) Hexachlorobenzene	16.69	284	149490	39.219	nq	97
69) Atrazine	16.84	200	82291	33.996	nq	99
70) Pentachlorophenol	17.04	266	71232	41.013	nq	99
71) Phenanthrene	17.42	178	502828	39.796	nq	100
72) Anthracene	17.52	178	503877	39.858	nq	99
73) Carbazole	17.79	167	456284	39.857	nq	97
74) Di-n-butylphthalate	18.33	149	562843	40.520	nq	99
75) Fluoranthene	19.43	202	658563	39.980	nq	99
77) Benzidine	19.61	184	243884	45.934	nq	98
78) Pyrene	19.80	202	658058	40.297	nq	100
80) Butylbenzylphthalate	20.68	149	260629	42.126	nq	98
81) Benzo(a)anthracene	21.63	228	682870	41.322	nq	99
82) 3,3'-Dichlorobenzidine	21.55	252	276816	43.308	nq	100
83) Chrysene	21.69	228	669745	40.790	nq	98
84) Bis(2-ethylhexyl)phthalate	21.53	149	371920	42.319	nq	97
85) Di-n-octyl phthalate	22.72	149	640149	42.933	nq	98
86) Indeno(1,2,3-cd)pyrene	28.48	276	874699	42.439	nq	100
88) Benzo(b)fluoranthene	23.83	252	717882	40.773	nq	98
89) Benzo(k)fluoranthene	23.89	252	744075	41.979	nq	98
90) Benzo(a)pyrene	24.69	252	701401	41.717	nq	99
91) Dibenzo(a,h)anthracene	28.54	278	730692	42.489	nq	99
92) Benzo(g,h,i)perylene	29.61	276	709532	41.827	nq	98

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

