

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG052919\
 Data File : BG040941.D
 Acq On : 29 May 2019 17:26
 Operator : HP/JU
 Sample : SSTDICC080
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDICC080

Quant Time: May 29 18:04:09 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG052919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 29 16:09:47 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.13	152	58440	20.00	ng	0.00
21) Naphthalene-d8	10.95	136	243059	20.00	ng	0.00
39) Acenaphthene-d10	14.76	164	183583	20.00	ng	0.00
64) Phenanthrene-d10	17.50	188	437043	20.00	ng	0.00
76) Chrysene-d12	21.79	240	419985	20.00	ng	0.00
87) Perylene-d12	25.13	264	460736	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.66	112	501222	151.19	ng	0.00
7) Phenol-d6	7.27	99	759292	148.75	ng	0.00
23) Nitrobenzene-d5	9.31	82	872221	165.81	ng	0.00
42) 2,4,6-Tribromophenol	16.25	330	432107	171.24	ng	0.00
45) 2-Fluorobiphenyl	13.38	172	1915780	150.71	ng	0.00
79) Terphenyl-d14	20.10	244	2790214	147.18	ng	0.00

Target Compounds				Qvalue		
2) 1,4-Dioxane	3.53	88	104669	69.422	ng	96
3) Pyridine	3.94	79	338674	77.497	ng	98
4) n-Nitrosodimethylamine	3.86	42	162924	67.213	ng	# 99
6) Aniline	7.46	93	514728	76.078	ng	98
8) 2-Chlorophenol	7.69	128	295232	72.932	ng	97
9) Benzaldehyde	7.27	77	200102	63.224	ng	91
10) Phenol	7.30	94	405561	73.911	ng	98
11) bis(2-Chloroethyl)ether	7.55	93	304167	73.922	ng	98
12) 1,3-Dichlorobenzene	8.01	146	359408	81.344	ng	97
13) 1,4-Dichlorobenzene	8.17	146	360451	80.475	ng	98
14) 1,2-Dichlorobenzene	8.48	146	343882	79.610	ng	97
15) Benzyl Alcohol	8.37	79	363585	68.455	ng	95
16) 2,2'-oxybis(1-Chloropropan	8.65	45	490745	59.905	ng	97
17) 2-Methylphenol	8.57	107	299600	77.153	ng	95
18) Hexachloroethane	9.22	117	140769	76.615	ng	98
19) n-Nitroso-di-n-propylamine	8.94	70	299006	65.072	ng	98
20) 3+4-Methylphenols	8.90	107	412280	76.213	ng	98
22) Acetophenone	8.97	105	544751	80.598	ng	# 97
24) Nitrobenzene	9.35	77	447034	79.650	ng	98
25) Isophorone	9.87	82	841337	71.803	ng	98
26) 2-Nitrophenol	10.06	139	195843	97.346	ng	96
27) 2,4-Dimethylphenol	10.11	122	306575	81.218	ng	94
28) bis(2-Chloroethoxy)methane	10.35	93	454256	77.270	ng	99
29) 2,4-Dichlorophenol	10.59	162	384838	89.677	ng	94
30) 1,2,4-Trichlorobenzene	10.81	180	445753	93.668	ng	98
31) Naphthalene	11.00	128	1031897	79.912	ng	98
32) Benzoic acid	10.28	122	246002	95.198	ng	98
33) 4-Chloroaniline	11.12	127	482942	84.352	ng	96
34) Hexachlorobutadiene	11.27	225	341304	97.145	ng	97
35) Caprolactam	11.93	113	127851	75.461	ng	90
36) 4-Chloro-3-methylphenol	12.22	107	434776	80.311	ng	100
37) 2-Methylnaphthalene	12.60	142	780593	80.851	ng	98
38) 1-Methylnaphthalene	12.82	142	742054	78.633	ng	# 99
40) 1,2,4,5-Tetrachlorobenzene	12.96	216	586077	85.697	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.93	237	318216	78.854	nq	97
43) 2,4,6-Trichlorophenol	13.20	196	376871	91.190	nq	96
44) 2,4,5-Trichlorophenol	13.27	196	391061	86.862	nq	98
46) 1,1'-Biphenyl	13.60	154	1058918	74.292	nq	98
47) 2-Chloronaphthalene	13.65	162	910183	81.596	nq	97
48) 2-Nitroaniline	13.85	65	334793	84.317	nq	99
49) Acenaphthylene	14.49	152	1322490	69.880	nq	97
50) Dimethylphthalate	14.21	163	1179017	76.759	nq	99
51) 2,6-Dinitrotoluene	14.34	165	257608	85.932	nq	98
52) Acenaphthene	14.83	154	830902	75.390	nq	98
53) 3-Nitroaniline	14.68	138	260079	84.672	nq	94
54) 2,4-Dinitrophenol	14.88	184	135890	91.973	nq	97
55) Dibenzofuran	15.16	168	1340208	74.241	nq	96
56) 4-Nitrophenol	14.96	139	191092	71.747	nq	98
57) 2,4-Dinitrotoluene	15.12	165	372631	91.476	nq	99
58) Fluorene	15.80	166	1072916	73.533	nq	100
59) 2,3,4,6-Tetrachlorophenol	15.38	232	382078	84.317	nq	98
60) Diethylphthalate	15.56	149	1193303	73.633	nq	95
61) 4-Chlorophenyl-phenylether	15.79	204	720001	84.844	nq	95
62) 4-Nitroaniline	15.85	138	273394	79.569	nq	96
63) Azobenzene	16.09	77	1089475	65.278	nq	95
65) 4,6-Dinitro-2-methylphenol	15.88	198	204139	101.492	nq	95
66) n-Nitrosodiphenylamine	16.01	169	976719	75.960	nq	98
67) 4-Bromophenyl-phenylether	16.69	248	479087	87.988	nq	97
68) Hexachlorobenzene	16.80	284	508264	90.276	nq	99
69) Atrazine	16.94	200	328058	64.396	nq	97
70) Pentachlorophenol	17.14	266	268972	81.641	nq	97
71) Phenanthrene	17.55	178	1719902	73.062	nq	94
72) Anthracene	17.64	178	1715021	72.480	nq	96
73) Carbazole	17.91	167	1482501	68.769	nq	97
74) Di-n-butylphthalate	18.44	149	1826099	70.181	nq	# 95
75) Fluoranthene	19.55	202	2086580	71.129	nq	94
77) Benzidine	19.73	184	705363	61.339	nq	97
78) Pyrene	19.91	202	2063889	78.407	nq	94
80) Butylbenzylphthalate	20.78	149	853758	83.217	nq	97
81) Benzo(a)anthracene	21.77	228	2146665	80.874	nq	95
82) 3,3'-Dichlorobenzidine	21.68	252	785396	83.016	nq	97
83) Chrysene	21.84	228	2051369	81.263	nq	93
84) Bis(2-ethylhexyl)phthalate	21.64	149	1200457	81.060	nq	97
85) Di-n-octyl phthalate	22.88	149	1992369	79.883	nq	99
86) Indeno(1,2,3-cd)pyrene	28.99	276	2663170	87.258	nq	# 98
88) Benzo(b)fluoranthene	24.07	252	2217083	78.746	nq	96
89) Benzo(k)fluoranthene	24.14	252	2127561	80.915	nq	96
90) Benzo(a)pyrene	24.98	252	2125637	79.758	nq	98
91) Dibenzo(a,h)anthracene	29.06	278	2163325	80.213	nq	99
92) Benzo(g,h,i)perylene	30.21	276	2070010	77.120	nq	96

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

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