

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM090220\
 Data File : BM027301.D
 Acq On : 02 Sep 2020 12:16
 Operator : JU/CG
 Sample : SSTDICC020
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICC020

Quant Time: Sep 02 13:48:40 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM090220.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 02 13:40:32 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.57	152	114145	20.00	ng	0.00
21) Naphthalene-d8	10.35	136	448578	20.00	ng	0.00
39) Acenaphthene-d10	14.22	164	328262	20.00	ng	0.00
64) Phenanthrene-d10	16.96	188	795966	20.00	ng	0.00
76) Chrysene-d12	21.17	240	942371	20.00	ng	0.00
86) Perylene-d12	23.34	264	1000103	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.20	112	267736	45.73	ng	0.00
7) Phenol-d6	6.76	99	358605	44.61	ng	0.00
23) Nitrobenzene-d5	8.72	82	403448	37.11	ng	0.00
42) 2,4,6-Tribromophenol	15.71	330	220061	41.40	ng	0.00
45) 2-Fluorobiphenyl	12.83	172	931757	37.29	ng	0.00
79) Terphenyl-d14	19.61	244	1906015	37.78	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.17	88	58451	20.889	ng	99
3) Pyridine	3.56	79	170222	21.671	ng	96
4) n-Nitrosodimethylamine	3.47	42	66237	19.330	ng	89
6) Aniline	6.91	93	219565	24.042	ng	99
8) 2-Chlorophenol	7.15	128	138230	22.686	ng	100
9) Benzaldehyde	6.73	77	144688	26.286	ng	99
10) Phenol	6.79	94	173959	22.705	ng	97
11) bis(2-Chloroethyl)ether	7.01	93	144807	23.122	ng	95
12) 1,3-Dichlorobenzene	7.47	146	169209	19.842	ng	98
13) 1,4-Dichlorobenzene	7.61	146	172208	20.015	ng	99
14) 1,2-Dichlorobenzene	7.92	146	165063	20.227	ng	99
15) Benzyl Alcohol	7.82	79	154120	21.091	ng	99
16) 2,2'-oxybis(1-Chloropropan	8.10	45	110964	22.085	ng	99
17) 2-Methylphenol	8.02	107	119243	22.493	ng	98
18) Hexachloroethane	8.65	117	67363	19.292	ng	96
19) n-Nitroso-di-n-propylamine	8.38	70	147900	23.968	ng	99
20) 3+4-Methylphenols	8.35	107	161203	22.725	ng	99
22) Acetophenone	8.39	105	243421	20.191	ng	# 97
24) Nitrobenzene	8.76	77	213735	19.421	ng	99
25) Isophorone	9.29	82	358572	22.516	ng	99
26) 2-Nitrophenol	9.47	139	75995	22.123	ng	97
27) 2,4-Dimethylphenol	9.54	122	134110	25.145	ng	98
28) bis(2-Chloroethoxy)methane	9.77	93	202267	20.813	ng	100
29) 2,4-Dichlorophenol	10.00	162	151661	20.389	ng	99
30) 1,2,4-Trichlorobenzene	10.22	180	179700	18.403	ng	99
31) Naphthalene	10.40	128	451941	19.878	ng	99
32) Benzoic acid	9.66	122	70543	18.459	ng	99
33) 4-Chloroaniline	10.51	127	197450	21.021	ng	98
34) Hexachlorobutadiene	10.69	225	121137	17.209	ng	99
35) Caprolactam	11.28	113	43065	21.358	ng	92
36) 4-Chloro-3-methylphenol	11.65	107	165957	21.584	ng	98
37) 2-Methylnaphthalene	12.02	142	354849	20.610	ng	98
38) 1-Methylnaphthalene	12.24	142	337326	20.358	ng	98
40) 1,2,4,5-Tetrachlorobenzene	12.40	216	221357	19.147	ng	99

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41) Hexachlorocyclopentadiene	12.38	237	140479	21.984	nq	99
43) 2,4,6-Trichlorophenol	12.64	196	135026	18.916	nq	98
44) 2,4,5-Trichlorophenol	12.71	196	148485	19.616	nq	96
46) 1,1'-Biphenyl	13.04	154	492377	20.013	nq	98
47) 2-Chloronaphthalene	13.08	162	391564	18.666	nq	99
48) 2-Nitroaniline	13.29	65	121570	18.159	nq	95
49) Acenaphthylene	13.93	152	596715	20.455	nq	99
50) Dimethylphthalate	13.69	163	518084	19.440	nq	100
51) 2,6-Dinitrotoluene	13.79	165	109988	20.755	nq	93
52) Acenaphthene	14.28	154	369435	19.949	nq	97
53) 3-Nitroaniline	14.12	138	100469	19.658	nq	97
54) 2,4-Dinitrophenol	14.33	184	55507	20.062	nq	94
55) Dibenzofuran	14.62	168	635323	20.594	nq	100
56) 4-Nitrophenol	14.44	139	83679	20.517	nq	98
57) 2,4-Dinitrotoluene	14.59	165	154143	20.937	nq	97
58) Fluorene	15.27	166	521291	20.468	nq	97
59) 2,3,4,6-Tetrachlorophenol	14.85	232	146688	19.783	nq	97
60) Diethylphthalate	15.06	149	538890	20.382	nq	98
61) 4-Chlorophenyl-phenylether	15.27	204	284221	19.430	nq	98
62) 4-Nitroaniline	15.29	138	104562	18.231	nq	90
63) Azobenzene	15.56	77	572284	21.098	nq	97
65) 4,6-Dinitro-2-methylphenol	15.36	198	86614	22.157	nq	95
66) n-Nitrosodiphenylamine	15.49	169	463154	20.487	nq	99
67) 4-Bromophenyl-phenylether	16.16	248	181806	18.918	nq	95
68) Hexachlorobenzene	16.28	284	213235	18.684	nq	99
69) Atrazine	16.44	200	137773	16.097	nq	98
70) Pentachlorophenol	16.62	266	122290	19.811	nq	99
71) Phenanthrene	17.00	178	875662	19.790	nq	99
72) Anthracene	17.10	178	886151	20.949	nq	99
73) Carbazole	17.37	167	787839	20.249	nq	99
74) Di-n-butylphthalate	17.95	149	940351	20.925	nq	100
75) Fluoranthene	19.03	202	1086328	19.654	nq	100
77) Benzidine	19.22	184	427644	26.162	nq	99
78) Pyrene	19.39	202	1140903	19.896	nq	100
80) Butylbenzylphthalate	20.32	149	434590	20.735	nq	96
81) Benzo(a)anthracene	21.15	228	1185392	19.412	nq	99
82) 3,3'-Dichlorobenzidine	21.09	252	446241	20.969	nq	99
83) Chrysene	21.20	228	1147735	19.352	nq	100
84) Bis(2-ethylhexyl)phthalate	21.11	149	681030	22.831	nq	99
85) Di-n-octyl phthalate	21.97	149	1142784	23.220	nq	100
87) Indeno(1,2,3-cd)pyrene	25.50	276	1336267	17.071	nq	100
88) Benzo(b)fluoranthene	22.69	252	1260186	18.896	nq	100
89) Benzo(k)fluoranthene	22.74	252	1237323	18.950	nq	99
90) Benzo(a)pyrene	23.24	252	1089352	17.893	nq	100
91) Dibenzo(a,h)anthracene	25.51	278	1130554	17.354	nq	99
92) Benzo(g,h,i)perylene	26.16	276	1105328	17.611	nq	98

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

