

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM072318\
 Data File : BM016048.D
 Acq On : 24 Jul 2018 04:49
 Operator : SJ/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDCCC040

Quant Time: Jul 24 05:23:16 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM072318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 23 16:08:38 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.23	152	41778	20.00	ng	0.00
21) Naphthalene-d8	11.05	136	177736	20.00	ng	0.00
38) Acenaphthene-d10	14.85	164	98142	20.00	ng	0.00
63) Phenanthrene-d10	17.58	188	213049	20.00	ng	0.00
75) Chrysene-d12	21.70	240	242257	20.00	ng	0.00
86) Perylene-d12	24.07	264	242654	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.74	112	209556	83.32	ng	0.00
7) Phenol-d6	7.38	99	266902	81.26	ng	0.00
23) Nitrobenzene-d5	9.42	82	327230	85.63	ng	0.00
41) 2,4,6-Tribromophenol	16.34	330	97709	78.50	ng	0.00
44) 2-Fluorobiphenyl	13.47	172	674448	83.62	ng	0.00
78) Terphenyl-d14	20.15	244	971177	83.92	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.52	88	46659	39.596	ng	# 100
3) Pyridine	3.96	79	109720	38.651	ng	# 72
4) n-Nitrosodimethylamine	3.87	42	88728	40.073	ng	# 35
6) Aniline	7.55	93	151024	39.935	ng	# 89
8) 2-Chlorophenol	7.79	128	121251	40.564	ng	96
9) Benzaldehyde	7.37	77	93467	40.503	ng	92
10) Phenol	7.41	94	132714	40.409	ng	97
11) bis(2-Chloroethyl)ether	7.65	93	100976	38.918	ng	87
12) 1,3-Dichlorobenzene	8.11	146	140127	40.050	ng	# 89
13) 1,4-Dichlorobenzene	8.26	146	141059	39.752	ng	# 94
14) 1,2-Dichlorobenzene	8.58	146	137149	39.500	ng	# 93
15) Benzyl Alcohol	8.48	79	107095	40.949	ng	# 85
16) 2,2'-oxybis(1-Chloropropan	8.75	45	155047	39.046	ng	98
17) 2-Methylphenol	8.67	107	91942	40.958	ng	# 82
18) Hexachloroethane	9.30	117	63997	41.722	ng	96
19) n-Nitroso-di-n-propylamine	9.05	70	96471	40.185	ng	# 76
20) 3+4-Methylphenols	9.01	107	124420	38.523	ng	87
22) Acetophenone	9.07	105	177758	39.719	ng	# 84
24) Nitrobenzene	9.46	77	154651	40.206	ng	99
25) Isophorone	9.97	82	210065	40.188	ng	# 94
26) 2-Nitrophenol	10.17	139	62333	38.820	ng	# 50
27) 2,4-Dimethylphenol	10.22	122	88505	39.563	ng	# 81
28) bis(2-Chloroethoxy)methane	10.45	93	135168	39.787	ng	98
29) 2,4-Dichlorophenol	10.70	162	111554	41.965	ng	97
30) 1,2,4-Trichlorobenzene	10.91	180	129771	40.179	ng	97
31) Naphthalene	11.10	128	355173	39.484	ng	99
32) Benzoic acid	10.37	122	62276	36.519	ng	86
33) 4-Chloroaniline	11.22	127	150465	40.990	ng	# 87
34) Hexachlorobutadiene	11.36	225	93326	41.656	ng	97
35) Caprolactam	12.02	113	29259	39.740	ng	90
36) 4-Chloro-3-methylphenol	12.33	107	120462	41.201	ng	88
37) 2-Methylnaphthalene	12.69	142	244606	40.593	ng	# 90
39) 1,2,4,5-Tetrachlorobenzene	13.06	216	135764	40.633	ng	# 100
40) Hexachlorocyclopentadiene	13.02	237	60546	41.049	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.30	196	86457	41.358	nq	99
43) 2,4,5-Trichlorophenol	13.37	196	86945	40.322	nq	# 83
45) 1,1'-Biphenyl	13.69	154	351340	39.668	nq	99
46) 2-Chloronaphthalene	13.74	162	279564	40.582	nq	# 89
47) 2-Nitroaniline	13.96	65	91303	40.089	nq	# 76
48) Acenaphthylene	14.58	152	411249	40.779	nq	100
49) Dimethylphthalate	14.30	163	338800	39.446	nq	99
50) 2,6-Dinitrotoluene	14.45	165	72029	41.680	nq	# 64
51) Acenaphthene	14.92	154	246809	39.793	nq	97
52) 3-Nitroaniline	14.77	138	72712	38.955	nq	# 67
53) 2,4-Dinitrophenol	15.00	184	25330	37.643	nq	# 82
54) Dibenzofuran	15.25	168	411216	40.233	nq	# 83
55) 4-Nitrophenol	15.08	139	51592	38.557	nq	85
56) 2,4-Dinitrotoluene	15.23	165	97839	39.595	nq	# 87
57) Fluorene	15.89	166	306504	40.356	nq	99
58) 2,3,4,6-Tetrachlorophenol	15.47	232	76378	41.209	nq	# 100
59) Diethylphthalate	15.65	149	374613	40.465	nq	95
60) 4-Chlorophenyl-phenylether	15.88	204	170061	40.222	nq	# 83
61) 4-Nitroaniline	15.93	138	70717	39.487	nq	# 10
62) Azobenzene	16.17	77	409750	42.027	nq	76
64) 4,6-Dinitro-2-methylphenol	15.99	198	51841	40.089	nq	85
65) n-Nitrosodiphenylamine	16.10	169	287540	40.987	nq	97
66) 4-Bromophenyl-phenylether	16.77	248	99012	41.042	nq	# 76
67) Hexachlorobenzene	16.89	284	108280	40.758	nq	# 64
68) Atrazine	17.03	200	103223	41.054	nq	97
69) Pentachlorophenol	17.23	266	65446	39.783	nq	98
70) Phenanthrene	17.63	178	473492	39.697	nq	99
71) Anthracene	17.72	178	468904	40.449	nq	98
72) Carbazole	17.99	167	485285	40.408	nq	98
73) Di-n-butylphthalate	18.51	149	618669	41.341	nq	# 96
74) Fluoranthene	19.61	202	539579	40.554	nq	99
76) Benzidine	19.79	184	272724	40.286	nq	99
77) Pyrene	19.97	202	574081	39.543	nq	99
79) Butylbenzylphthalate	20.82	149	301206	40.841	nq	# 79
80) Benzo(a)anthracene	21.69	228	589636	39.518	nq	99
81) 3,3'-Dichlorobenzidine	21.61	252	243938	40.589	nq	97
82) Chrysene	21.74	228	568935	39.750	nq	98
83) Bis(2-ethylhexyl)phthalate	21.57	149	436112	40.802	nq	# 90
84) Di-n-octyl phthalate	22.48	149	722091	40.197	nq	# 89
85) Indeno(1,2,3-cd)pyrene	26.50	276	672986	39.623	nq	# 94
87) Benzo(b)fluoranthene	23.35	252	585843	41.006	nq	# 95
88) Benzo(k)fluoranthene	23.40	252	572683	39.553	nq	98
89) Benzo(a)pyrene	23.97	252	554290	40.357	nq	100
90) Dibenzo(a,h)anthracene	26.50	278	572481	40.231	nq	97
91) Benzo(g,h,i)perylene	27.25	276	535679	39.802	nq	# 93

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

