

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM102218\
 Data File : BM017274.D
 Acq On : 22 Oct 2018 12:05
 Operator : MJ/SJ
 Sample : SSTDC0040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :

Quant Time: Oct 22 13:05:14 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM101718.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 17 14:23:17 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.66	152	270670	20.00	ng	-0.01
21) Naphthalene-d8	10.44	136	1268486	20.00	ng	-0.01
39) Acenaphthene-d10	14.30	164	777547	20.00	ng	-0.01
64) Phenanthrene-d10	17.06	188	1777274	20.00	ng	-0.01
76) Chrysene-d12	21.25	240	1917796	20.00	ng	-0.01
87) Perylene-d12	23.46	264	1707026	20.00	ng	-0.02

System Monitoring Compounds

5) 2-Fluorophenol	5.27	112	1362824	85.16	ng	-0.01
7) Phenol-d6	6.85	99	1921115	88.14	ng	-0.01
23) Nitrobenzene-d5	8.82	82	1944937	89.67	ng	-0.02
42) 2,4,6-Tribromophenol	15.80	330	969656	92.21	ng	-0.01
45) 2-Fluorobiphenyl	12.92	172	4859629	83.16	ng	-0.01
79) Terphenyl-d14	19.70	244	7724005	90.78	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.26	88	310247	37.965	ng	90
3) Pyridine	3.64	79	794105	42.233	ng	# 86
4) n-Nitrosodimethylamine	3.56	42	305356	40.453	ng	# 71
6) Aniline	7.00	93	1193943	43.795	ng	97
8) 2-Chlorophenol	7.23	128	822346	44.418	ng	97
9) Benzaldehyde	6.82	77	565725	40.708	ng	95
10) Phenol	6.87	94	1013628	43.211	ng	99
11) bis(2-Chloroethyl)ether	7.10	93	804949	43.047	ng	99
12) 1,3-Dichlorobenzene	7.55	146	909605	42.146	ng	98
13) 1,4-Dichlorobenzene	7.70	146	934243	42.369	ng	98
14) 1,2-Dichlorobenzene	8.01	146	891475	41.962	ng	99
15) Benzyl Alcohol	7.90	79	777103	48.778	ng	96
16) 2,2'-oxybis(1-Chloropropan	8.19	45	1068546	40.902	ng	99
17) 2-Methylphenol	8.10	107	684195	45.442	ng	98
18) Hexachloroethane	8.72	117	325745	43.167	ng	98
19) n-Nitroso-di-n-propylamine	8.47	70	687115	47.877	ng	98
20) 3+4-Methylphenols	8.43	107	978615	47.522	ng	99
22) Acetophenone	8.48	105	1338690	41.633	ng	# 99
24) Nitrobenzene	8.86	77	987290	43.034	ng	94
25) Isophorone	9.38	82	1617977	45.177	ng	98
26) 2-Nitrophenol	9.56	139	517089	48.140	ng	96
27) 2,4-Dimethylphenol	9.63	122	711721	44.942	ng	99
28) bis(2-Chloroethoxy)methane	9.86	93	1089426	43.218	ng	98
29) 2,4-Dichlorophenol	10.09	162	852862	46.383	ng	99
30) 1,2,4-Trichlorobenzene	10.30	180	928764	42.106	ng	98
31) Naphthalene	10.49	128	2576957	41.455	ng	99
32) Benzoic acid	9.79	122	575342	48.691	ng	93
33) 4-Chloroaniline	10.60	127	1217321	45.342	ng	98
34) Hexachlorobutadiene	10.77	225	574664	41.118	ng	98
35) Caprolactam	11.41	113	318403	47.516	ng	91
36) 4-Chloro-3-methylphenol	11.73	107	970863	46.842	ng	99
37) 2-Methylnaphthalene	12.10	142	1895163	44.552	ng	99
38) 1-Methylnaphthalene	12.33	142	1842140	44.494	ng	99
40) 1,2,4,5-Tetrachlorobenzene	12.47	216	1139901	41.629	ng	99

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41) Hexachlorocyclopentadiene	12.45	237	609544	46.046	ng	99
43) 2,4,6-Trichlorophenol	12.72	196	750029	47.185	ng	98
44) 2,4,5-Trichlorophenol	12.80	196	782331	45.897	ng	99
46) 1,1'-Biphenyl	13.13	154	2911838	41.615	ng	99
47) 2-Chloronaphthalene	13.17	162	2130152	41.382	ng	99
48) 2-Nitroaniline	13.39	65	643810	44.712	ng	97
49) Acenaphthylene	14.03	152	3234307	43.347	ng	100
50) Dimethylphthalate	13.77	163	2652495	44.217	ng	99
51) 2,6-Dinitrotoluene	13.89	165	602859	45.660	ng	100
52) Acenaphthene	14.37	154	1972537	42.105	ng	98
53) 3-Nitroaniline	14.22	138	652055	45.599	ng	95
54) 2,4-Dinitrophenol	14.43	184	307005	43.786	ng	99
55) Dibenzofuran	14.70	168	3232949	41.498	ng	99
56) 4-Nitrophenol	14.53	139	495040	40.462	ng	98
57) 2,4-Dinitrotoluene	14.68	165	832816	46.931	ng	# 98
58) Fluorene	15.36	166	2459569	42.653	ng	99
59) 2,3,4,6-Tetrachlorophenol	14.93	232	718697	46.256	ng	99
60) Diethylphthalate	15.14	149	2706417	45.493	ng	100
61) 4-Chlorophenyl-phenylether	15.36	204	1405733	42.761	ng	97
62) 4-Nitroaniline	15.39	138	673245	42.764	ng	99
63) Azobenzene	15.65	77	2485009	42.975	ng	99
65) 4,6-Dinitro-2-methylphenol	15.44	198	458937	41.348	ng	97
66) n-Nitrosodiphenylamine	15.57	169	2390694	44.461	ng	99
67) 4-Bromophenyl-phenylether	16.25	248	892996	45.760	ng	96
68) Hexachlorobenzene	16.36	284	987462	43.562	ng	98
69) Atrazine	16.53	200	884276	45.949	ng	99
70) Pentachlorophenol	16.70	266	692391	44.589	ng	100
71) Phenanthrene	17.10	178	4003638	41.752	ng	100
72) Anthracene	17.19	178	4009615	44.007	ng	99
73) Carbazole	17.46	167	4036510	43.370	ng	100
74) Di-n-butylphthalate	18.03	149	4675197	47.626	ng	99
75) Fluoranthene	19.12	202	4687565	43.443	ng	99
77) Benzidine	19.32	184	2426355	49.896	ng	98
78) Pyrene	19.48	202	4917587	45.879	ng	100
80) Butylbenzylphthalate	20.40	149	2150179	52.011	ng	97
81) Benzo(a)anthracene	21.23	228	4640086	42.995	ng	100
82) 3,3'-Dichlorobenzidine	21.17	252	1806042	44.901	ng	99
83) Chrysene	21.29	228	4401643	41.206	ng	99
84) Bis(2-ethylhexyl)phthalate	21.17	149	3104065	49.548	ng	99
85) Di-n-octyl phthalate	22.05	149	4996908	47.350	ng	99
86) Indeno(1,2,3-cd)pyrene	25.67	276	3840029m	32.140	ng	
88) Benzo(b)fluoranthene	22.80	252	4142015	44.381	ng	100
89) Benzo(k)fluoranthene	22.84	252	4113879	43.929	ng	99
90) Benzo(a)pyrene	23.36	252	3781620	42.675	ng	100
91) Dibenzo(a,h)anthracene	25.68	278	3242044	36.846	ng	99
92) Benzo(g,h,i)perylene	26.34	276	3088711	35.418	ng	99

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

