

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM020719\
 Data File : BM018733.D
 Acq On : 07 Feb 2019 11:59
 Operator : JU/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDCCC040

Quant Time: Feb 08 10:04:06 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM010919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Feb 04 15:10:38 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.72	152	245242	20.00	ng	0.00
21) Naphthalene-d8	10.50	136	958725	20.00	ng	0.00
39) Acenaphthene-d10	14.36	164	525407	20.00	ng	0.00
64) Phenanthrene-d10	17.12	188	1138091	20.00	ng	0.00
76) Chrysene-d12	21.31	240	1148220	20.00	ng	0.00
87) Perylene-d12	23.57	264	1140990	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.30	112	1099807	82.81	ng	0.00
7) Phenol-d6	6.89	99	1413170	83.77	ng	0.00
23) Nitrobenzene-d5	8.88	82	1339969	81.02	ng	0.00
42) 2,4,6-Tribromophenol	15.86	330	565579	84.54	ng	0.00
45) 2-Fluorobiphenyl	12.98	172	3222317	80.03	ng	0.00
79) Terphenyl-d14	19.75	244	4700974	86.31	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.25	88	207868	38.400	ng	98
3) Pyridine	3.65	79	553430	41.007	ng	96
4) n-Nitrosodimethylamine	3.58	42	231140	41.417	ng	# 92
6) Aniline	7.06	93	834664	44.356	ng	98
8) 2-Chlorophenol	7.28	128	645660	41.624	ng	96
9) Benzaldehyde	6.88	77	343108	33.841	ng	99
10) Phenol	6.92	94	715314	41.729	ng	99
11) bis(2-Chloroethyl)ether	7.15	93	542045	40.243	ng	97
12) 1,3-Dichlorobenzene	7.60	146	737054	39.899	ng	98
13) 1,4-Dichlorobenzene	7.75	146	752722	40.202	ng	99
14) 1,2-Dichlorobenzene	8.06	146	718323	40.460	ng	99
15) Benzyl Alcohol	7.96	79	502550	41.247	ng	96
16) 2,2'-oxybis(1-Chloropropan	8.25	45	746106	43.346	ng	96
17) 2-Methylphenol	8.16	107	486845	41.840	ng	99
18) Hexachloroethane	8.78	117	258098	38.661	ng	97
19) n-Nitroso-di-n-propylamine	8.53	70	442985	40.352	ng	94
20) 3+4-Methylphenols	8.49	107	677369	42.170	ng	99
22) Acetophenone	8.55	105	932187	39.306	ng	# 98
24) Nitrobenzene	8.93	77	659161	40.508	ng	97
25) Isophorone	9.45	82	997479	39.830	ng	98
26) 2-Nitrophenol	9.63	139	349392	44.405	ng	96
27) 2,4-Dimethylphenol	9.68	122	482925	40.948	ng	98
28) bis(2-Chloroethoxy)methane	9.93	93	695826	40.221	ng	99
29) 2,4-Dichlorophenol	10.16	162	589363	41.968	ng	99
30) 1,2,4-Trichlorobenzene	10.36	180	689118	39.534	ng	99
31) Naphthalene	10.56	128	1854438	40.163	ng	99
32) Benzoic acid	9.83	122	344821	44.045	ng	99
33) 4-Chloroaniline	10.68	127	810568	44.576	ng	98
34) Hexachlorobutadiene	10.82	225	423668	38.473	ng	99
35) Caprolactam	11.49	113	190976	39.999	ng	89
36) 4-Chloro-3-methylphenol	11.80	107	609010	41.336	ng	98
37) 2-Methylnaphthalene	12.17	142	1288416	39.495	ng	99
38) 1-Methylnaphthalene	12.39	142	1258027	39.855	ng	97
40) 1,2,4,5-Tetrachlorobenzene	12.54	216	743358	40.459	ng	98

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41) Hexachlorocyclopentadiene	12.50	237	343650	34.080	ng	100
43) 2,4,6-Trichlorophenol	12.79	196	482017	43.202	ng	99
44) 2,4,5-Trichlorophenol	12.86	196	494431	42.136	ng	96
46) 1,1'-Biphenyl	13.19	154	1839070	40.088	ng	99
47) 2-Chloronaphthalene	13.24	162	1433206	40.287	ng	99
48) 2-Nitroaniline	13.46	65	372089	42.513	ng	94
49) Acenaphthylene	14.09	152	2113211	40.761	ng	100
50) Dimethylphthalate	13.83	163	1692413	39.221	ng	99
51) 2,6-Dinitrotoluene	13.96	165	377709	41.324	ng	94
52) Acenaphthene	14.43	154	1274922	40.192	ng	98
53) 3-Nitroaniline	14.29	138	397568	43.144	ng	97
54) 2,4-Dinitrophenol	14.50	184	157282	36.127	ng	92
55) Dibenzofuran	14.77	168	2067037	39.438	ng	99
56) 4-Nitrophenol	14.60	139	289660	36.386	ng	98
57) 2,4-Dinitrotoluene	14.75	165	512822	39.654	ng	99
58) Fluorene	15.42	166	1552615	39.766	ng	99
59) 2,3,4,6-Tetrachlorophenol	14.99	232	423213	40.690	ng	97
60) Diethylphthalate	15.20	149	1702133	39.075	ng	99
61) 4-Chlorophenyl-phenylether	15.42	204	882123	39.201	ng	98
62) 4-Nitroaniline	15.46	138	368988	36.711	ng	97
63) Azobenzene	15.70	77	1449389	40.810	ng	98
65) 4,6-Dinitro-2-methylphenol	15.50	198	296671	41.695	ng	91
66) n-Nitrosodiphenylamine	15.63	169	1474352	41.763	ng	99
67) 4-Bromophenyl-phenylether	16.31	248	524070	41.149	ng	98
68) Hexachlorobenzene	16.41	284	581146	40.760	ng	100
69) Atrazine	16.59	200	462233	36.218	ng	99
70) Pentachlorophenol	16.76	266	389505	41.713	ng	97
71) Phenanthrene	17.16	178	2451591	40.499	ng	99
72) Anthracene	17.25	178	2418336	41.544	ng	100
73) Carbazole	17.53	167	2450916	40.982	ng	99
74) Di-n-butylphthalate	18.08	149	2730194	41.677	ng	99
75) Fluoranthene	19.18	202	2750146	39.405	ng	99
77) Benzidine	19.37	184	959406	33.840	ng	100
78) Pyrene	19.55	202	2907185	42.715	ng	99
80) Butylbenzylphthalate	20.45	149	1235818	42.553	ng	98
81) Benzo(a)anthracene	21.29	228	2754659	40.723	ng	100
82) 3,3'-Dichlorobenzidine	21.23	252	1011680	39.551	ng	100
83) Chrysene	21.34	228	2660231	40.726	ng	100
84) Bis(2-ethylhexyl)phthalate	21.22	149	1731077	41.278	ng	100
85) Di-n-octyl phthalate	22.10	149	2889436	40.965	ng	96
86) Indeno(1,2,3-cd)pyrene	25.87	276	2928086	38.875	ng	98
88) Benzo(b)fluoranthene	22.89	252	2608979	40.277	ng	99
89) Benzo(k)fluoranthene	22.93	252	2653973	40.655	ng	99
90) Benzo(a)pyrene	23.47	252	2518415	40.610	ng	98
91) Dibenzo(a,h)anthracene	25.87	278	2440431	38.838	ng	100
92) Benzo(g,h,i)perylene	26.57	276	2434821	39.819	ng	98

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

