

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM031419\
 Data File : BM019164.D
 Acq On : 14 Mar 2019 12:01
 Operator : JU/SJ
 Sample : K1824-01MSRE
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 TP-24MS

Quant Time: Mar 14 13:27:19 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM031119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 11 17:30:46 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.66	152	206273	20.00	ng	0.00
21) Naphthalene-d8	10.44	136	936600	20.00	ng	-0.01
39) Acenaphthene-d10	14.31	164	595205	20.00	ng	-0.01
64) Phenanthrene-d10	17.07	188	1402802	20.00	ng	0.00
76) Chrysene-d12	21.27	240	1478460	20.00	ng	0.00
87) Perylene-d12	23.50	264	1302736	20.00	ng	-0.01

System Monitoring Compounds

5) 2-Fluorophenol	5.26	112	1743118	136.85	ng	0.00
7) Phenol-d6	6.84	99	2688101	144.28	ng	0.00
23) Nitrobenzene-d5	8.83	82	1647926	83.17	ng	0.00
42) 2,4,6-Tribromophenol	15.82	330	1248098	142.02	ng	0.00
45) 2-Fluorobiphenyl	12.92	172	3468242	75.79	ng	-0.01
79) Terphenyl-d14	19.70	244	6203407	83.30	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.22	88	172189	30.389	ng	98
3) Pyridine	3.62	79	273490	17.741	ng	98
4) n-Nitrosodimethylamine	3.53	42	168474	35.155	ng	91
6) Aniline	7.00	93	396067	16.466	ng	99
8) 2-Chlorophenol	7.22	128	649214	42.415	ng	98
9) Benzaldehyde	6.82	77	298751	25.494	ng	98
10) Phenol	6.87	94	1058029	52.695	ng	100
11) bis(2-Chloroethyl)ether	7.10	93	579161	37.802	ng	99
12) 1,3-Dichlorobenzene	7.54	146	583991	36.054	ng	99
13) 1,4-Dichlorobenzene	7.69	146	604421	36.601	ng	99
14) 1,2-Dichlorobenzene	8.00	146	595822	37.085	ng	98
15) Benzyl Alcohol	7.91	79	651019	42.221	ng	99
16) 2,2'-oxybis(1-Chloropropan	8.19	45	581493	37.170	ng	98
17) 2-Methylphenol	8.10	107	581134	44.416	ng	99
18) Hexachloroethane	8.72	117	220921	36.002	ng	95
19) n-Nitroso-di-n-propylamine	8.47	70	597667	39.103	ng	100
20) 3+4-Methylphenols	8.43	107	814276	45.849	ng	99
22) Acetophenone	8.49	105	971418	37.530	ng	# 99
24) Nitrobenzene	8.87	77	824312	40.002	ng	99
25) Isophorone	9.39	82	1539821	40.728	ng	99
26) 2-Nitrophenol	9.57	139	357284	41.829	ng	99
27) 2,4-Dimethylphenol	9.63	122	626590	49.344	ng	99
28) bis(2-Chloroethoxy)methane	9.87	93	913100	40.788	ng	99
29) 2,4-Dichlorophenol	10.10	162	666773	47.411	ng	99
30) 1,2,4-Trichlorobenzene	10.30	180	625440	39.440	ng	99
31) Naphthalene	10.49	128	1949164	39.484	ng	99
32) Benzoic acid	9.77	122	170090	15.816	ng	98
33) 4-Chloroaniline	10.63	127	212940	10.523	ng	99
34) Hexachlorobutadiene	10.75	225	331890	36.470	ng	99
35) Caprolactam	11.45	113	101759	20.312	ng	98
36) 4-Chloro-3-methylphenol	11.75	107	807330	46.524	ng	99
37) 2-Methylnaphthalene	12.11	142	1506523	41.996	ng	99
38) 1-Methylnaphthalene	12.33	142	1439015	40.698	ng	99
40) 1,2,4,5-Tetrachlorobenzene	12.48	216	715073	37.980	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.44	237	755604	88.600	ng	99
43) 2,4,6-Trichlorophenol	12.73	196	564638	46.406	ng	99
44) 2,4,5-Trichlorophenol	12.80	196	603646	46.395	ng	99
46) 1,1'-Biphenyl	13.14	154	2016214	38.788	ng	99
47) 2-Chloronaphthalene	13.18	162	1555370	40.087	ng	100
48) 2-Nitroaniline	13.41	65	575782	44.249	ng	97
49) Acenaphthylene	14.03	152	2610788	39.796	ng	100
50) Dimethylphthalate	13.79	163	2589958	50.866	ng	99
51) 2,6-Dinitrotoluene	13.91	165	478260	43.451	ng	99
52) Acenaphthene	14.37	154	1505096	39.926	ng	99
53) 3-Nitroaniline	14.24	138	243194	20.861	ng	97
54) 2,4-Dinitrophenol	14.46	184	464228	72.625	ng	100
55) Dibenzofuran	14.72	168	2475821	40.725	ng	100
56) 4-Nitrophenol	14.56	139	841891	88.452	ng	98
57) 2,4-Dinitrotoluene	14.70	165	691082	43.254	ng	99
58) Fluorene	15.36	166	2072892	41.366	ng	99
59) 2,3,4,6-Tetrachlorophenol	14.94	232	552936	45.650	ng	99
60) Diethylphthalate	15.15	149	2207977	42.817	ng	99
61) 4-Chlorophenyl-phenylether	15.36	204	1027149	42.204	ng	98
62) 4-Nitroaniline	15.42	138	479323	40.800	ng	99
63) Azobenzene	15.66	77	2305663	42.490	ng	100
65) 4,6-Dinitro-2-methylphenol	15.46	198	348488	37.910	ng	98
66) n-Nitrosodiphenylamine	15.59	169	1833198	41.047	ng	99
67) 4-Bromophenyl-phenylether	16.26	248	619454	40.044	ng	98
68) Hexachlorobenzene	16.36	284	740457	40.261	ng	98
69) Atrazine	16.54	200	656044	43.638	ng	98
70) Pentachlorophenol	16.72	266	923531	83.347	ng	100
71) Phenanthrene	17.11	178	3258157	39.626	ng	100
72) Anthracene	17.20	178	3340588	41.501	ng	99
73) Carbazole	17.48	167	3143730	41.380	ng	99
74) Di-n-butylphthalate	18.03	149	3955132	43.052	ng	100
75) Fluoranthene	19.13	202	3794440	40.217	ng	99
77) Benzidine	19.34	184	343798	8.199	ng	99
78) Pyrene	19.50	202	3868613	41.356	ng	100
80) Butylbenzylphthalate	20.40	149	1882832	46.212	ng	98
81) Benzo(a)anthracene	21.25	228	3638726	39.230	ng	100
82) 3,3'-Dichlorobenzidine	21.19	252	748442	20.946	ng	98
83) Chrysene	21.30	228	3480504	38.787	ng	100
84) Bis(2-ethylhexyl)phthalate	21.17	149	2788953	47.051	ng	# 98
85) Di-n-octyl phthalate	22.05	149	4676045	46.848	ng	100
86) Indeno(1,2,3-cd)pyrene	25.77	276	3482105	36.049	ng	100
88) Benzo(b)fluoranthene	22.83	252	3481634	41.245	ng	100
89) Benzo(k)fluoranthene	22.87	252	3389743	43.658	ng	99
90) Benzo(a)pyrene	23.41	252	3198032	42.157	ng	99
91) Dibenzo(a,h)anthracene	25.77	278	2991573	41.225	ng	99
92) Benzo(g,h,i)perylene	26.46	276	2738954	41.111	ng	100

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

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