

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM060419\
 Data File : BM020709.D
 Acq On : 04 Jun 2019 18:48
 Operator : HP/JU
 Sample : PB120137BS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB120137BS

Manual Integrations
 APPROVED

mohammad
 6/5/2019 3:41:32 PM

Quant Time: Jun 05 03:51:10 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\8270-BM053119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri May 31 15:48:49 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.83	152	316143	20.00	ng	0.00
21) Naphthalene-d8	10.62	136	1199603	20.00	ng	-0.01
39) Acenaphthene-d10	14.46	164	615611	20.00	ng	-0.01
64) Phenanthrene-d10	17.20	188	1195680	20.00	ng	-0.01
76) Chrysene-d12	21.38	240	1009918	20.00	ng	-0.02
87) Perylene-d12	23.67	264	1180340	20.00	ng	-0.02

System Monitoring Compounds

5) 2-Fluorophenol	5.41	112	2435584	135.52	ng	0.00
7) Phenol-d6	6.99	99	3155460	119.24	ng	0.00
23) Nitrobenzene-d5	8.99	82	1924039	83.10	ng	-0.01
42) 2,4,6-Tribromophenol	15.95	330	1039197	136.25	ng	-0.01
45) 2-Fluorobiphenyl	13.09	172	4154020	89.71	ng	0.00
79) Terphenyl-d14	19.83	244	4717759	78.14	ng	-0.02

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.35	88	308004	41.859	ng	97
3) Pyridine	3.74	79	852589	36.418	ng	98
4) n-Nitrosodimethylamine	3.66	42	390991	38.743	ng	96
6) Aniline	7.16	93	1079809	28.200	ng	98
8) 2-Chlorophenol	7.39	128	957770	42.838	ng	98
9) Benzaldehyde	6.97	77	292510	15.307	ng	97
10) Phenol	7.02	94	1151243	40.368	ng	98
11) bis(2-Chloroethyl)ether	7.26	93	863005	37.971	ng	97
12) 1,3-Dichlorobenzene	7.71	146	1009347	39.280	ng	99
13) 1,4-Dichlorobenzene	7.86	146	1028824	39.384	ng	98
14) 1,2-Dichlorobenzene	8.17	146	990724	38.972	ng	99
15) Benzyl Alcohol	8.06	79	834706	35.654	ng	97
16) 2,2'-oxybis(1-Chloropropan	8.36	45	1197554	34.129	ng	99
17) 2-Methylphenol	8.26	107	759175	37.857	ng	99
18) Hexachloroethane	8.90	117	377916	37.786	ng	94
19) n-Nitroso-di-n-propylamine	8.63	70	705441	33.162	ng	96
20) 3+4-Methylphenols	8.59	107	1006024	36.872	ng	97
22) Acetophenone	8.65	105	1317410	43.230	ng	# 96
24) Nitrobenzene	9.03	77	1034360	43.761	ng	98
25) Isophorone	9.55	82	1828689	39.706	ng	100
26) 2-Nitrophenol	9.73	139	448297	50.890	ng	99
27) 2,4-Dimethylphenol	9.79	122	811795	49.179	ng	98
28) bis(2-Chloroethoxy)methane	10.03	93	1117571	40.133	ng	99
29) 2,4-Dichlorophenol	10.26	162	833620	45.646	ng	98
30) 1,2,4-Trichlorobenzene	10.47	180	934598	44.506	ng	100
31) Naphthalene	10.67	128	2704478	43.841	ng	100
32) Benzoic acid	9.92	122	462277m	44.868	ng	
33) 4-Chloroaniline	10.78	127	395903	13.790	ng	97
34) Hexachlorobutadiene	10.94	225	543790	43.158	ng	98
35) Caprolactam	11.57	113	231965m	36.562	ng	
36) 4-Chloro-3-methylphenol	11.90	107	848568	39.964	ng	99
37) 2-Methylnaphthalene	12.28	142	1928089	42.168	ng	99
38) 1-Methylnaphthalene	12.50	142	1837344	42.201	ng	100
40) 1,2,4,5-Tetrachlorobenzene	12.64	216	983957	48.773	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.62	237	1249446	103.554	ng	100
43) 2,4,6-Trichlorophenol	12.89	196	616899	47.506	ng	98
44) 2,4,5-Trichlorophenol	12.96	196	648372	48.363	ng	97
46) 1,1'-Biphenyl	13.29	154	2373703	46.169	ng	98
47) 2-Chloronaphthalene	13.33	162	1838837	44.314	ng	100
48) 2-Nitroaniline	13.54	65	509684	39.134	ng	90
49) Acenaphthylene	14.19	152	2879495	44.474	ng	99
50) Dimethylphthalate	13.92	163	2166621	41.195	ng	99
51) 2,6-Dinitrotoluene	14.04	165	466384	43.671	ng	89
52) Acenaphthene	14.53	154	1673115	43.363	ng	99
53) 3-Nitroaniline	14.37	138	273999	22.987	ng	96
54) 2,4-Dinitrophenol	14.57	184	380212	82.507	ng	95
55) Dibenzofuran	14.86	168	2586086	43.154	ng	98
56) 4-Nitrophenol	14.67	139	741806	84.450	ng	91
57) 2,4-Dinitrotoluene	14.83	165	604738	42.292	ng	95
58) Fluorene	15.51	166	2063693	41.876	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.08	232	544653	46.623	ng	98
60) Diethylphthalate	15.29	149	2130065	39.268	ng	99
61) 4-Chlorophenyl-phenylether	15.50	204	1067883	40.926	ng	98
62) 4-Nitroaniline	15.53	138	441376	34.914	ng	96
63) Azobenzene	15.80	77	2011627	38.682	ng	97
65) 4,6-Dinitro-2-methylphenol	15.59	198	234254	43.965	ng	97
66) n-Nitrosodiphenylamine	15.72	169	1760835	46.003	ng	100
67) 4-Bromophenyl-phenylether	16.40	248	620580	44.082	ng	96
68) Hexachlorobenzene	16.50	284	710173	43.201	ng	96
69) Atrazine	16.67	200	551510	48.277	ng	100
70) Pentachlorophenol	16.85	266	810298	103.323	ng	100
71) Phenanthrene	17.25	178	2936681	43.330	ng	100
72) Anthracene	17.34	178	3015806	44.458	ng	100
73) Carbazole	17.61	167	2589592	42.068	ng	99
74) Di-n-butylphthalate	18.17	149	3158738	39.529	ng	100
75) Fluoranthene	19.26	202	3103192	41.172	ng	99
77) Benzidine	19.45	184	1434237	46.868	ng	99
78) Pyrene	19.62	202	3151858	40.669	ng	100
80) Butylbenzylphthalate	20.53	149	1260148	38.476	ng	93
81) Benzo(a)anthracene	21.37	228	2911939	41.404	ng	100
82) 3,3'-Dichlorobenzidine	21.30	252	863267	33.520	ng	100
83) Chrysene	21.42	228	2879060	43.066	ng	100
84) Bis(2-ethylhexyl)phthalate	21.30	149	1773266	37.253	ng	100
85) Di-n-octyl phthalate	22.19	149	2952076	39.063	ng	97
86) Indeno(1,2,3-cd)pyrene	25.99	276	4077042	59.612	ng	99
88) Benzo(b)fluoranthene	22.98	252	3236198	41.482	ng	100
89) Benzo(k)fluoranthene	23.03	252	3119736	40.921	ng	99
90) Benzo(a)pyrene	23.57	252	3162803	44.696	ng	100
91) Dibenzo(a,h)anthracene	26.01	278	3437795	50.076	ng	98
92) Benzo(g,h,i)perylene	26.71	276	3324301	52.132	ng	99

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

