

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG060719\
 Data File : BG041114.D
 Acq On : 8 Jun 2019 1:35
 Operator : HP/JU
 Sample : PB120393BSD
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB120393BSD

Manual Integrations
 APPROVED

mohammad
 6/11/2019 8:54:34 AM

Quant Time: Jun 08 18:08:41 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG052919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Jun 08 17:33:39 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.11	152	39332	20.00	ng	0.00
21) Naphthalene-d8	10.93	136	162154	20.00	ng	0.00
39) Acenaphthene-d10	14.74	164	125570	20.00	ng	0.00
64) Phenanthrene-d10	17.48	188	310283	20.00	ng	0.00
76) Chrysene-d12	21.76	240	306085	20.00	ng	0.00
87) Perylene-d12	25.09	264	321726	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.64	112	309060	141.69	ng	0.00
7) Phenol-d6	7.25	99	441488	131.06	ng	0.00
23) Nitrobenzene-d5	9.29	82	299458	81.98	ng	0.00
42) 2,4,6-Tribromophenol	16.23	330	239829	128.65	ng	0.00
45) 2-Fluorobiphenyl	13.36	172	726254	83.29	ng	0.00
79) Terphenyl-d14	20.08	244	1268384	87.95	ng	0.00

Target Compounds				Qvalue		
2) 1,4-Dioxane	3.50	88	40130	40.544	ng	94
3) Pyridine	3.90	79	106986	36.529	ng	96
4) n-Nitrosodimethylamine	3.83	42	58347	42.925	ng	99
6) Aniline	7.43	93	136566	30.372	ng	99
8) 2-Chlorophenol	7.66	128	127337	48.968	ng	99
9) Benzaldehyde	7.25	77	37712	18.767	ng	93
10) Phenol	7.28	94	175202	48.507	ng	98
11) bis(2-Chloroethyl)ether	7.52	93	112360	42.882	ng	94
12) 1,3-Dichlorobenzene	7.99	146	125369	40.365	ng	92
13) 1,4-Dichlorobenzene	8.14	146	130383	41.473	ng	97
14) 1,2-Dichlorobenzene	8.46	146	130811	42.854	ng	98
15) Benzyl Alcohol	8.35	79	142279	45.659	ng	98
16) 2,2'-oxybis(1-Chloropropan	8.63	45	173378	40.493	ng	98
17) 2-Methylphenol	8.54	107	125624	48.037	ng	97
18) Hexachloroethane	9.19	117	49598	40.933	ng	94
19) n-Nitroso-di-n-propylamine	8.91	70	110586	42.658	ng	97
20) 3+4-Methylphenols	8.87	107	168193	46.430	ng	96
22) Acetophenone	8.94	105	204531	44.965	ng	# 97
24) Nitrobenzene	9.33	77	170861	45.901	ng	97
25) Isophorone	9.84	82	317374	45.657	ng	99
26) 2-Nitrophenol	10.04	139	77485	50.494	ng	96
27) 2,4-Dimethylphenol	10.08	122	139298	54.963	ng	96
28) bis(2-Chloroethoxy)methane	10.33	93	167905	44.754	ng	96
29) 2,4-Dichlorophenol	10.57	162	156268	48.555	ng	97
30) 1,2,4-Trichlorobenzene	10.78	180	162341	44.341	ng	97
31) Naphthalene	10.98	128	399785	45.920	ng	98
32) Benzoic acid	10.23	122	87385	44.962	ng	95
33) 4-Chloroaniline	11.09	127	86696	21.462	ng	98
34) Hexachlorobutadiene	11.24	225	117294	41.870	ng	97
35) Caprolactam	11.88	113	48676m	45.800	ng	
36) 4-Chloro-3-methylphenol	12.19	107	177465	48.282	ng	94
37) 2-Methylnaphthalene	12.57	142	310362	46.286	ng	99
38) 1-Methylnaphthalene	12.79	142	294627	46.628	ng	99
40) 1,2,4,5-Tetrachlorobenzene	12.93	216	226043	44.662	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.90	237	296837	106.111	ng	96
43) 2,4,6-Trichlorophenol	13.17	196	149261	45.588	ng	96
44) 2,4,5-Trichlorophenol	13.25	196	160988	46.623	ng	98
46) 1,1'-Biphenyl	13.58	154	424585	45.679	ng	99
47) 2-Chloronaphthalene	13.62	162	357776	44.837	ng	99
48) 2-Nitroaniline	13.83	65	126629	44.235	ng	98
49) Acenaphthylene	14.46	152	554912	46.205	ng	99
50) Dimethylphthalate	14.19	163	476587	44.836	ng	98
51) 2,6-Dinitrotoluene	14.32	165	105928	46.220	ng	98
52) Acenaphthene	14.80	154	323661	44.487	ng	98
53) 3-Nitroaniline	14.65	138	70164	30.616	ng	100
54) 2,4-Dinitrophenol	14.86	184	141070	104.524	ng	99
55) Dibenzofuran	15.14	168	561834	45.894	ng	99
56) 4-Nitrophenol	14.94	139	156767	99.729	ng	98
57) 2,4-Dinitrotoluene	15.10	165	148404	45.841	ng	# 99
58) Fluorene	15.78	166	442921	45.431	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.36	232	163859	48.287	ng	99
60) Diethylphthalate	15.54	149	478666	44.126	ng	99
61) 4-Chlorophenyl-phenylether	15.77	204	281812	44.472	ng	96
62) 4-Nitroaniline	15.81	138	101874	41.989	ng	99
63) Azobenzene	16.06	77	442592	44.890	ng	98
65) 4,6-Dinitro-2-methylphenol	15.86	198	92221	54.129	ng	95
66) n-Nitrosodiphenylamine	15.98	169	410118	47.019	ng	100
67) 4-Bromophenyl-phenylether	16.66	248	181402	43.321	ng	93
68) Hexachlorobenzene	16.78	284	188535	42.283	ng	99
69) Atrazine	16.92	200	170226	50.578	ng	99
70) Pentachlorophenol	17.12	266	231368	95.965	ng	97
71) Phenanthrene	17.52	178	750217	46.418	ng	98
72) Anthracene	17.62	178	764847	47.982	ng	99
73) Carbazole	17.89	167	659410	48.212	ng	99
74) Di-n-butylphthalate	18.42	149	771403	45.378	ng	100
75) Fluoranthene	19.53	202	958304	48.004	ng	99
77) Benzidine	19.71	184	381302	52.221	ng	99
78) Pyrene	19.89	202	959969	48.246	ng	98
80) Butylbenzylphthalate	20.76	149	356826	46.082	ng	99
81) Benzo(a)anthracene	21.74	228	923162	45.309	ng	98
82) 3,3'-Dichlorobenzidine	21.65	252	212377	29.635	ng	99
83) Chrysene	21.81	228	909104	46.626	ng	99
84) Bis(2-ethylhexyl)phthalate	21.61	149	502664	46.115	ng	99
85) Di-n-octyl phthalate	22.85	149	840751	46.312	ng	100
86) Indeno(1,2,3-cd)pyrene	28.92	276	1104691	46.251	ng	100
88) Benzo(b)fluoranthene	24.02	252	949639	48.665	ng	98
89) Benzo(k)fluoranthene	24.09	252	908888	47.068	ng	98
90) Benzo(a)pyrene	24.93	252	917745	49.295	ng	98
91) Dibenzo(a,h)anthracene	28.98	278	903598	48.573	ng	99
92) Benzo(g,h,i)perylene	30.11	276	910314	50.368	ng	98

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

