

Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG040319\
 Data File : BG040328.D
 Acq On : 3 Apr 2019 16:23
 Operator : JU/SJ
 Sample : PB118462BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB118462BS

Manual Integrations
 APPROVED

sohil
 4/4/2019 4:46:29 AM

Quant Time: Apr 04 01:53:04 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG032719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 28 05:47:00 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.05	152	48463	20.00	ng	-0.02
21) Naphthalene-d8	10.86	136	190854	20.00	ng	-0.02
39) Acenaphthene-d10	14.68	164	119977	20.00	ng	-0.02
64) Phenanthrene-d10	17.43	188	324377	20.00	ng	-0.02
76) Chrysene-d12	21.70	240	380770	20.00	ng	-0.03
87) Perylene-d12	24.99	264	372397	20.00	ng	-0.04

System Monitoring Compounds

5) 2-Fluorophenol	5.61	112	382837	131.32	ng	-0.01
7) Phenol-d6	7.20	99	506070	125.61	ng	-0.02
23) Nitrobenzene-d5	9.22	82	308152	85.82	ng	-0.02
42) 2,4,6-Tribromophenol	16.18	330	209501	161.57	ng	-0.01
45) 2-Fluorobiphenyl	13.30	172	706415	87.24	ng	-0.02
79) Terphenyl-d14	20.03	244	1375327	81.26	ng	-0.02

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.49	88	54822	39.993	ng	99
3) Pyridine	3.90	79	134452	36.429	ng	96
4) n-Nitrosodimethylamine	3.81	42	60802	35.614	ng	# 97
6) Aniline	7.37	93	138209	26.404	ng	96
8) 2-Chlorophenol	7.61	128	130362	38.201	ng	98
9) Benzaldehyde	7.19	77	88595	32.058	ng	96
10) Phenol	7.23	94	171808	39.288	ng	95
11) bis(2-Chloroethyl)ether	7.47	93	128233	36.612	ng	96
12) 1,3-Dichlorobenzene	7.94	146	144642	38.991	ng	96
13) 1,4-Dichlorobenzene	8.08	146	144411	39.086	ng	99
14) 1,2-Dichlorobenzene	8.40	146	133433	37.765	ng	97
15) Benzyl Alcohol	8.29	79	115833	35.700	ng	96
16) 2,2'-oxybis(1-Chloropropan	8.57	45	238003	35.406	ng	99
17) 2-Methylphenol	8.49	107	115800	38.268	ng	97
18) Hexachloroethane	9.12	117	51020	36.303	ng	99
19) n-Nitroso-di-n-propylamine	8.85	70	100739	34.875	ng	98
20) 3+4-Methylphenols	8.82	107	159345	38.871	ng	96
22) Acetophenone	8.88	105	195469	41.058	ng	# 99
24) Nitrobenzene	9.27	77	147024	39.542	ng	97
25) Isophorone	9.78	82	285754	38.678	ng	99
26) 2-Nitrophenol	9.98	139	72161	40.958	ng	98
27) 2,4-Dimethylphenol	10.02	122	128018	45.744	ng	93
28) bis(2-Chloroethoxy)methane	10.26	93	171732	39.290	ng	96
29) 2,4-Dichlorophenol	10.51	162	133602	43.982	ng	97
30) 1,2,4-Trichlorobenzene	10.72	180	136051	40.790	ng	99
31) Naphthalene	10.92	128	401364	41.028	ng	98
32) Benzoic acid	10.12	122	25942	15.342	ng	91
33) 4-Chloroaniline	11.03	127	102775	24.630	ng	98
34) Hexachlorobutadiene	11.18	225	84576	39.648	ng	94
35) Caprolactam	11.80	113	51546m	42.760	ng	
36) 4-Chloro-3-methylphenol	12.14	107	155064	44.404	ng	95
37) 2-Methylnaphthalene	12.52	142	300764	41.570	ng	99
38) 1-Methylnaphthalene	12.73	142	284156	40.502	ng	99
40) 1,2,4,5-Tetrachlorobenzene	12.87	216	153373	40.221	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.84	237	183461	87.622	ng	95
43) 2,4,6-Trichlorophenol	13.12	196	112601	45.800	ng	98
44) 2,4,5-Trichlorophenol	13.19	196	129042	48.154	ng	98
46) 1,1'-Biphenyl	13.51	154	382172	40.315	ng	99
47) 2-Chloronaphthalene	13.57	162	299028	41.123	ng	99
48) 2-Nitroaniline	13.78	65	104509	42.609	ng	97
49) Acenaphthylene	14.41	152	506589	41.090	ng	99
50) Dimethylphthalate	14.13	163	407809	43.431	ng	99
51) 2,6-Dinitrotoluene	14.27	165	91427	43.364	ng	97
52) Acenaphthene	14.75	154	291597	41.276	ng	96
53) 3-Nitroaniline	14.60	138	68822	30.837	ng	97
54) 2,4-Dinitrophenol	14.81	184	79341	70.760	ng	93
55) Dibenzofuran	15.08	168	481888	42.450	ng	98
56) 4-Nitrophenol	14.90	139	178088	98.870	ng	98
57) 2,4-Dinitrotoluene	15.05	165	134951	46.065	ng	97
58) Fluorene	15.73	166	387914	43.464	ng	96
59) 2,3,4,6-Tetrachlorophenol	15.31	232	125740	51.708	ng	99
60) Diethylphthalate	15.48	149	443371	46.143	ng	99
61) 4-Chlorophenyl-phenylether	15.71	204	208163	43.634	ng	98
62) 4-Nitroaniline	15.76	138	113994	47.595	ng	94
63) Azobenzene	16.01	77	385482	42.532	ng	99
65) 4,6-Dinitro-2-methylphenol	15.81	198	69150	37.944	ng	99
66) n-Nitrosodiphenylamine	15.93	169	372947	39.290	ng	98
67) 4-Bromophenyl-phenylether	16.61	248	137428	37.648	ng	98
68) Hexachlorobenzene	16.72	284	144507	39.372	ng	96
69) Atrazine	16.87	200	142742	40.425	ng	98
70) Pentachlorophenol	17.07	266	159678	75.251	ng	96
71) Phenanthrene	17.47	178	710246	41.657	ng	99
72) Anthracene	17.56	178	723938	42.421	ng	99
73) Carbazole	17.84	167	719758	44.565	ng	99
74) Di-n-butylphthalate	18.37	149	874235	45.666	ng	100
75) Fluoranthene	19.48	202	941461	46.632	ng	99
77) Benzidine	19.66	184	444577	32.333	ng	98
78) Pyrene	19.84	202	945909	37.776	ng	99
80) Butylbenzylphthalate	20.71	149	443392	41.381	ng	99
81) Benzo(a)anthracene	21.68	228	901215	38.426	ng	98
82) 3,3'-Dichlorobenzidine	21.60	252	203178	22.773	ng	100
83) Chrysene	21.75	228	891658	39.559	ng	99
84) Bis(2-ethylhexyl)phthalate	21.55	149	627822	40.990	ng	99
85) Di-n-octyl phthalate	22.77	149	1069608	40.988	ng	99
86) Indeno(1,2,3-cd)pyrene	28.75	276	1089025	39.503	ng	# 92
88) Benzo(b)fluoranthene	23.94	252	968477	42.478	ng	99
89) Benzo(k)fluoranthene	24.01	252	908929	43.351	ng	98
90) Benzo(a)pyrene	24.83	252	936455	43.776	ng	99
91) Dibenzo(a,h)anthracene	28.81	278	893763	44.985	ng	99
92) Benzo(g,h,i)perylene	29.94	276	917928	44.956	ng	97

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

