

Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG041119\
 Data File : BG040439.D
 Acq On : 11 Apr 2019 13:28
 Operator : JU/SJ
 Sample : PB118804BS
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB118804BS

Manual Integrations
 APPROVED

Sohil
 4/15/2019 12:22:59 AM

Quant Time: Apr 12 03:50:05 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	8.04	152	54508	20.00	ng	-0.03
21) Naphthalene-d8	10.86	136	237685	20.00	ng	-0.03
39) Acenaphthene-d10	14.67	164	153225	20.00	ng	-0.03
64) Phenanthrene-d10	17.42	188	415620	20.00	ng	-0.02
76) Chrysene-d12	21.70	240	439389	20.00	ng	-0.04
87) Perylene-d12	24.97	264	431005	20.00	ng	-0.06

System Monitoring Compounds

5) 2-Fluorophenol	5.60	112	456257	139.15	ng	-0.02
7) Phenol-d6	7.21	99	631028	139.26	ng	-0.02
23) Nitrobenzene-d5	9.22	82	390986	87.44	ng	-0.02
42) 2,4,6-Tribromophenol	16.17	330	263367	159.04	ng	-0.02
45) 2-Fluorobiphenyl	13.30	172	880593	85.16	ng	-0.02
79) Terphenyl-d14	20.02	244	1590126	81.41	ng	-0.03

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.49	88	53669	34.810	ng	# 95
3) Pyridine	3.89	79	132367	31.887	ng	97
4) n-Nitrosodimethylamine	3.82	42	72364	37.686	ng	# 90
6) Aniline	7.37	93	182879	31.063	ng	98
8) 2-Chlorophenol	7.61	128	163244	42.531	ng	96
9) Benzaldehyde	7.19	77	74051	23.824	ng	93
10) Phenol	7.23	94	215102	43.733	ng	99
11) bis(2-Chloroethyl)ether	7.46	93	157516	39.985	ng	97
12) 1,3-Dichlorobenzene	7.93	146	159779	38.295	ng	95
13) 1,4-Dichlorobenzene	8.08	146	162373	39.073	ng	100
14) 1,2-Dichlorobenzene	8.40	146	156951	39.495	ng	96
15) Benzyl Alcohol	8.29	79	142927	39.165	ng	98
16) 2,2'-oxybis(1-Chloropropan	8.57	45	297461	39.344	ng	98
17) 2-Methylphenol	8.48	107	145862	42.857	ng	99
18) Hexachloroethane	9.12	117	59639	37.729	ng	97
19) n-Nitroso-di-n-propylamine	8.85	70	123006	37.861	ng	97
20) 3+4-Methylphenols	8.81	107	200837	43.559	ng	98
22) Acetophenone	8.88	105	242032	40.822	ng	# 98
24) Nitrobenzene	9.26	77	186001	40.168	ng	98
25) Isophorone	9.77	82	361572	39.298	ng	99
26) 2-Nitrophenol	9.97	139	90865	41.413	ng	98
27) 2,4-Dimethylphenol	10.01	122	167779	48.140	ng	98
28) bis(2-Chloroethoxy)methane	10.26	93	215876	39.658	ng	99
29) 2,4-Dichlorophenol	10.50	162	166144	43.919	ng	99
30) 1,2,4-Trichlorobenzene	10.71	180	170333	41.006	ng	97
31) Naphthalene	10.91	128	497600	40.844	ng	99
32) Benzoic acid	10.17	122	66993	31.814	ng	93
33) 4-Chloroaniline	11.03	127	145305	27.961	ng	97
34) Hexachlorobutadiene	11.17	225	101959	38.380	ng	94
35) Caprolactam	11.81	113	64480m	42.951	ng	
36) 4-Chloro-3-methylphenol	12.14	107	190567	43.818	ng	98
37) 2-Methylnaphthalene	12.51	142	372258	41.314	ng	98
38) 1-Methylnaphthalene	12.73	142	359165	41.107	ng	98
40) 1,2,4,5-Tetrachlorobenzene	12.87	216	190367	39.090	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.84	237	213513	79.848	ng	96
43) 2,4,6-Trichlorophenol	13.11	196	137795	43.886	ng	97
44) 2,4,5-Trichlorophenol	13.19	196	150236	43.898	ng	96
46) 1,1'-Biphenyl	13.51	154	484258	39.999	ng	98
47) 2-Chloronaphthalene	13.56	162	378684	40.777	ng	100
48) 2-Nitroaniline	13.77	65	135012	43.101	ng	98
49) Acenaphthylene	14.40	152	622718	39.549	ng	100
50) Dimethylphthalate	14.13	163	503015	41.946	ng	99
51) 2,6-Dinitrotoluene	14.26	165	117391	43.597	ng	96
52) Acenaphthene	14.74	154	370713	41.089	ng	97
53) 3-Nitroaniline	14.60	138	97774	34.304	ng	92
54) 2,4-Dinitrophenol	14.80	184	128348	89.628	ng	94
55) Dibenzofuran	15.07	168	619364	42.722	ng	98
56) 4-Nitrophenol	14.90	139	208403	90.595	ng	97
57) 2,4-Dinitrotoluene	15.04	165	174756	46.708	ng	96
58) Fluorene	15.72	166	490289	43.014	ng	97
59) 2,3,4,6-Tetrachlorophenol	15.30	232	149777	48.228	ng	98
60) Diethylphthalate	15.48	149	546761	44.556	ng	100
61) 4-Chlorophenyl-phenylether	15.71	204	262575	43.097	ng	100
62) 4-Nitroaniline	15.76	138	141792	46.356	ng	95
63) Azobenzene	16.00	77	502578	43.419	ng	100
65) 4,6-Dinitro-2-methylphenol	15.80	198	96789	41.451	ng	98
66) n-Nitrosodiphenylamine	15.93	169	476808	39.204	ng	98
67) 4-Bromophenyl-phenylether	16.60	248	176614	37.761	ng	97
68) Hexachlorobenzene	16.72	284	184102	39.148	ng	97
69) Atrazine	16.87	200	181868	40.199	ng	98
70) Pentachlorophenol	17.07	266	211893	77.936	ng	95
71) Phenanthrene	17.47	178	894660	40.954	ng	99
72) Anthracene	17.56	178	920476	42.096	ng	98
73) Carbazole	17.83	167	886780	42.853	ng	99
74) Di-n-butylphthalate	18.36	149	1070001	43.622	ng	99
75) Fluoranthene	19.47	202	1133033	43.800	ng	99
77) Benzidine	19.66	184	421237	26.548	ng	99
78) Pyrene	19.83	202	1116047	38.625	ng	99
80) Butylbenzylphthalate	20.70	149	520318	42.082	ng	96
81) Benzo(a)anthracene	21.68	228	1044523	38.595	ng	98
82) 3,3'-Dichlorobenzidine	21.59	252	307659	29.884	ng	97
83) Chrysene	21.74	228	1039207	39.954	ng	100
84) Bis(2-ethylhexyl)phthalate	21.54	149	756214	42.785	ng	99
85) Di-n-octyl phthalate	22.76	149	1287950	42.770	ng	100
86) Indeno(1,2,3-cd)pyrene	28.74	276	1274939	40.077	ng	100
88) Benzo(b)fluoranthene	23.92	252	1132104	42.903	ng	99
89) Benzo(k)fluoranthene	23.99	252	1089717	44.906	ng	99
90) Benzo(a)pyrene	24.82	252	1107608	44.736	ng	99
91) Dibenzo(a,h)anthracene	28.80	278	1041778	45.305	ng	98
92) Benzo(g,h,i)perylene	29.93	276	1063267	44.993	ng	98

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

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