

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG122118\
 Data File : BG038791.D
 Acq On : 21 Dec 2018 18:09
 Operator : JU/SJ
 Sample : PB115909BSD
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB115909BSD

Manual Integrations
 APPROVED

Sohil
 12/26/2018 3:17:28 PM

Quant Time: Dec 22 01:48:23 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG121218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 12 15:26:27 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.06	152	22194	20.00	ng	0.00
21) Naphthalene-d8	10.88	136	90509	20.00	ng	0.00
39) Acenaphthene-d10	14.70	164	65323	20.00	ng	0.00
64) Phenanthrene-d10	17.44	188	171849	20.00	ng	0.00
76) Chrysene-d12	21.73	240	173722	20.00	ng	0.00
87) Perylene-d12	25.03	264	181040	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.61	112	145974	116.66	ng	0.00
7) Phenol-d6	7.22	99	196514	114.40	ng	0.01
23) Nitrobenzene-d5	9.23	82	102266	57.72	ng	0.00
42) 2,4,6-Tribromophenol	16.19	330	116851	129.80	ng	0.00
45) 2-Fluorobiphenyl	13.32	172	277969	58.38	ng	0.00
79) Terphenyl-d14	20.04	244	612391	76.72	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.49	88	22135	38.008	ng	98
3) Pyridine	3.89	79	48070	29.980	ng	96
4) n-Nitrosodimethylamine	3.82	42	30593	38.940	ng	# 92
6) Aniline	7.38	93	62862	28.666	ng	99
8) 2-Chlorophenol	7.62	128	61224	42.280	ng	98
9) Benzaldehyde	7.20	77	18117	16.258	ng	92
10) Phenol	7.24	94	78466	42.995	ng	97
11) bis(2-Chloroethyl)ether	7.48	93	50290	34.611	ng	97
12) 1,3-Dichlorobenzene	7.94	146	64322	37.568	ng	98
13) 1,4-Dichlorobenzene	8.09	146	67184	38.313	ng	94
14) 1,2-Dichlorobenzene	8.41	146	63219	38.242	ng	96
15) Benzyl Alcohol	8.30	79	54512	40.656	ng	95
16) 2,2'-oxybis(1-Chloropropan	8.58	45	105550	31.712	ng	96
17) 2-Methylphenol	8.50	107	52982	43.331	ng	98
18) Hexachloroethane	9.13	117	22185	34.623	ng	90
19) n-Nitroso-di-n-propylamine	8.86	70	40901	33.901	ng	98
20) 3+4-Methylphenols	8.83	107	72379	42.862	ng	98
22) Acetophenone	8.89	105	88490	39.142	ng	# 96
24) Nitrobenzene	9.28	77	65774	36.699	ng	98
25) Isophorone	9.79	82	119159	37.434	ng	99
26) 2-Nitrophenol	9.99	139	35959	39.200	ng	93
27) 2,4-Dimethylphenol	10.03	122	63786	48.519	ng	97
28) bis(2-Chloroethoxy)methane	10.27	93	66820	37.197	ng	99
29) 2,4-Dichlorophenol	10.52	162	69260	47.356	ng	98
30) 1,2,4-Trichlorobenzene	10.73	180	72559	41.072	ng	96
31) Naphthalene	10.93	128	188760	42.328	ng	99
32) Benzoic acid	10.19	122	33818m	42.087	ng	
33) 4-Chloroaniline	11.04	127	50878	26.279	ng	98
34) Hexachlorobutadiene	11.19	225	47963	40.124	ng	97
35) Caprolactam	11.83	113	23557	38.920	ng	91
36) 4-Chloro-3-methylphenol	12.15	107	72816	43.628	ng	91
37) 2-Methylnaphthalene	12.53	142	143932	45.241	ng	98
38) 1-Methylnaphthalene	12.75	142	138274	44.790	ng	96
40) 1,2,4,5-Tetrachlorobenzene	12.89	216	91231	37.363	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.85	237	100676	75.048	ng	96
43) 2,4,6-Trichlorophenol	13.13	196	65467	43.606	ng	98
44) 2,4,5-Trichlorophenol	13.21	196	71052	44.699	ng	99
46) 1,1'-Biphenyl	13.53	154	195187	35.643	ng	98
47) 2-Chloronaphthalene	13.58	162	162888	38.507	ng	98
48) 2-Nitroaniline	13.79	65	51732	38.395	ng	92
49) Acenaphthylene	14.42	152	265407	41.970	ng	99
50) Dimethylphthalate	14.15	163	211150	39.582	ng	100
51) 2,6-Dinitrotoluene	14.28	165	48521	40.137	ng	91
52) Acenaphthene	14.76	154	151472	40.057	ng	96
53) 3-Nitroaniline	14.62	138	39535	32.082	ng	# 92
54) 2,4-Dinitrophenol	14.82	184	48627	68.407	ng	95
55) Dibenzofuran	15.09	168	262738	40.534	ng	96
56) 4-Nitrophenol	14.92	139	73892	79.040	ng	98
57) 2,4-Dinitrotoluene	15.07	165	71426	41.949	ng	92
58) Fluorene	15.74	166	216492	45.081	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.32	232	68405	47.831	ng	97
60) Diethylphthalate	15.50	149	219742	37.524	ng	99
61) 4-Chlorophenyl-phenylether	15.73	204	122937	41.029	ng	97
62) 4-Nitroaniline	15.78	138	52241	41.177	ng	94
63) Azobenzene	16.02	77	184865	37.788	ng	95
65) 4,6-Dinitro-2-methylphenol	15.83	198	42800	34.914	ng	93
66) n-Nitrosodiphenylamine	15.95	169	199006	39.413	ng	98
67) 4-Bromophenyl-phenylether	16.63	248	84801	40.405	ng	96
68) Hexachlorobenzene	16.74	284	88887	40.856	ng	94
69) Atrazine	16.89	200	81094	43.277	ng	98
70) Pentachlorophenol	17.09	266	90244	70.128	ng	97
71) Phenanthrene	17.49	178	375850	42.176	ng	99
72) Anthracene	17.58	178	384935	43.755	ng	99
73) Carbazole	17.85	167	337777	38.822	ng	98
74) Di-n-butylphthalate	18.38	149	386058	38.670	ng	99
75) Fluoranthene	19.49	202	462003	41.901	ng	97
77) Benzidine	19.68	184	204884	39.879	ng	99
78) Pyrene	19.86	202	463416	40.379	ng	99
80) Butylbenzylphthalate	20.73	149	172030	38.367	ng	97
81) Benzo(a)anthracene	21.71	228	452687	42.764	ng	99
82) 3,3'-Dichlorobenzidine	21.62	252	131207	31.252	ng	98
83) Chrysene	21.78	228	422509	41.438	ng	98
84) Bis(2-ethylhexyl)phthalate	21.57	149	238677	37.533	ng	97
85) Di-n-octyl phthalate	22.80	149	407417	38.255	ng	99
86) Indeno(1,2,3-cd)pyrene	28.82	276	500215	41.729	ng	# 76
88) Benzo(b)fluoranthene	23.98	252	443827	44.651	ng	98
89) Benzo(k)fluoranthene	24.04	252	433691	43.816	ng	98
90) Benzo(a)pyrene	24.87	252	436390	45.508	ng	# 97
91) Dibenzo(a,h)anthracene	28.88	278	415262	43.493	ng	98
92) Benzo(g,h,i)perylene	30.02	276	407301	43.286	ng	98

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

