

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062521\
 Data File : BG049093.D
 Acq On : 25 Jun 2021 16:38
 Operator : CG/JU
 Sample : M2824-02MS
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 ABN-1MS

Quant Time: Jun 25 17:32:44 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG061121.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 24 11:02:36 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.108	152	34699	20.00	ng	0.00
21) Naphthalene-d8	10.928	136	141395	20.00	ng	# 0.00
39) Acenaphthene-d10	14.747	164	99545	20.00	ng	0.00
64) Phenanthrene-d10	17.497	188	227548	20.00	ng	# 0.00
76) Chrysene-d12	21.786	240	221348	20.00	ng	0.00
86) Perylene-d12	25.106	264	238569	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.664	112	10004	4.49	ng	0.00
7) Phenol-d6	7.256	99	59386	17.78	ng	0.00
23) Nitrobenzene-d5	9.277	82	220993	67.39	ng	0.00
42) 2,4,6-Tribromophenol	16.234	330	4243	2.88	ng	0.00
45) 2-Fluorobiphenyl	13.373	172	465328	66.04	ng	0.00
79) Terphenyl-d14	20.100	244	831744	68.82	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.537	88	34527	31.73	ng	# 80
3) Pyridine	3.942	79	85354	28.91	ng	91
4) n-Nitrosodimethylamine	3.854	42	67887	48.02	ng	# 79
6) Aniline	7.421	93	145572	34.60	ng	# 92
8) 2-Chlorophenol	7.668	128	115681	47.08	ng	91
9) Benzaldehyde	7.239	77	86087	49.80	ng	# 85
10) Phenol	7.286	94	153477	44.47	ng	90
11) bis(2-Chloroethyl)ether	7.521	93	103077	40.25	ng	83
12) 1,3-Dichlorobenzene	7.997	146	116207	42.34	ng	97
13) 1,4-Dichlorobenzene	8.143	146	119538	43.69	ng	99
14) 1,2-Dichlorobenzene	8.461	146	115599	43.92	ng	94
15) Benzyl Alcohol	8.343	79	127784	41.76	ng	96
16) 2,2'-oxybis(1-Chloropr...	8.637	45	175548	36.23	ng	84
17) 2-Methylphenol	8.549	107	99127	43.79	ng	94
18) Hexachloroethane	9.195	117	48826	44.48	ng	95
19) n-Nitroso-di-n-propyla...	8.913	70	93853	35.94	ng	96
20) 3+4-Methylphenols	8.878	107	133777	43.23	ng	96
22) Acetophenone	8.931	105	185801	43.19	ng	# 97
24) Nitrobenzene	9.319	77	150067	40.63	ng	95
25) Isophorone	9.841	82	252825	35.39	ng	# 96
26) 2-Nitrophenol	10.029	139	61860	49.37	ng	97
27) 2,4-Dimethylphenol	10.088	122	106328	46.81	ng	94
28) bis(2-Chloroethoxy)met...	10.323	93	137543	41.58	ng	93
29) 2,4-Dichlorophenol	10.570	162	115525	45.33	ng	92
30) 1,2,4-Trichlorobenzene	10.787	180	127505	43.64	ng	96
31) Naphthalene	10.981	128	340292	40.40	ng	99
32) Benzoic acid	10.241	122	73824	49.75	ng	# 73
33) 4-Chloroaniline	11.081	127	83264	23.28	ng	95
34) Hexachlorobutadiene	11.263	225	90804	43.60	ng	93
35) Caprolactam	11.868	113	37663	51.12	ng	# 48
36) 4-Chloro-3-methylphenol	12.203	107	135972	43.65	ng	# 92
37) 2-Methylnaphthalene	12.579	142	257722	40.51	ng	97
38) 1-Methylnaphthalene	12.797	142	238625	39.95	ng	96
40) 1,2,4,5-Tetrachloroben...	12.944	216	151688	47.99	ng	98
41) Hexachlorocyclopentadiene	12.926	237	222192	109.35	ng	96
43) 2,4,6-Trichlorophenol	13.179	196	107789	50.97	ng	97

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.261	196	120003	54.87	ng	91
46) 1,1'-Biphenyl	13.584	154	322469	43.94	ng	98
47) 2-Chloronaphthalene	13.625	162	258120	44.04	ng	99
48) 2-Nitroaniline	13.825	65	101098	47.27	ng	96
49) Acenaphthylene	14.471	152	418357	42.84	ng	95
50) Dimethylphthalate	14.201	163	350522	45.20	ng	100
51) 2,6-Dinitrotoluene	14.319	165	80059	50.70	ng	92
52) Acenaphthene	14.812	154	247305	39.29	ng	94
53) 3-Nitroaniline	14.648	138	53463	33.44	ng	83
54) 2,4-Dinitrophenol	14.853	184	102086	139.56	ng	91
55) Dibenzofuran	15.147	168	404293	41.32	ng	97
56) 4-Nitrophenol	14.959	139	129569	99.41	ng	92
57) 2,4-Dinitrotoluene	15.106	165	113401	58.46	ng	94
58) Fluorene	15.799	166	344305	43.03	ng	98
59) 2,3,4,6-Tetrachlorophenol	15.370	232	108485	49.01	ng	99
60) Diethylphthalate	15.558	149	373072	44.42	ng	99
61) 4-Chlorophenyl-phenyle...	15.782	204	189877	45.38	ng	98
62) 4-Nitroaniline	15.811	138	80075	49.53	ng	89
63) Azobenzene	16.075	77	355446	36.99	ng	90
65) 4,6-Dinitro-2-methylph...	15.864	198	65080	60.31	ng	92
66) n-Nitrosodiphenylamine	15.999	169	299479	42.23	ng	97
67) 4-Bromophenyl-phenylether	16.680	248	134210	47.24	ng	92
68) Hexachlorobenzene	16.804	284	145663	47.30	ng	97
69) Atrazine	16.945	200	128516	56.46	ng	97
70) Pentachlorophenol	17.145	266	190487	102.93	ng	97
71) Phenanthrene	17.538	178	559616	43.71	ng	98
72) Anthracene	17.632	178	572825	44.52	ng	98
73) Carbazole	17.897	167	521504	42.29	ng	99
74) Di-n-butylphthalate	18.455	149	632342	45.72	ng	98
75) Fluoranthene	19.548	202	703806	45.58	ng	99
77) Benzidine	19.724	184	359212	80.56	ng	97
78) Pyrene	19.906	202	698983	43.51	ng	96
80) Butylbenzylphthalate	20.787	149	277728	45.82	ng	94
81) Benzo(a)anthracene	21.769	228	705481	44.39	ng	99
82) 3,3'-Dichlorobenzidine	21.675	252	193132	38.00	ng	99
83) Chrysene	21.833	228	671873	44.09	ng	97
84) Bis(2-ethylhexyl)phtha...	21.663	149	396092	45.84	ng	97
85) Di-n-octyl phthalate	22.914	149	680059	46.41	ng	98
87) Indeno(1,2,3-cd)pyrene	28.913	276	862661	48.05	ng	98
88) Benzo(b)fluoranthene	24.048	252	758096	44.58	ng	# 93
89) Benzo(k)fluoranthene	24.119	252	717357	44.46	ng	# 99
90) Benzo(a)pyrene	24.953	252	728649	46.94	ng	98
91) Dibenzo(a,h)anthracene	28.978	278	722849	48.17	ng	# 96
92) Benzo(g,h,i)perylene	30.094	276	719228	47.53	ng	99

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

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