

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120122\
 Data File : BG055860.D
 Acq On : 1 Dec 2022 18:00
 Operator : CG/JU
 Sample : N5823-01MSD
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 OE-2-113022MSD

Quant Time: Dec 02 04:10:49 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG111622.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 02 04:07:17 2022
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By : Christian Giraldo 12/02/2022
 Supervised By : Jagrut Upadhyay 12/02/2022

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.209	152	40565	20.000	ng	0.00
21) Naphthalene-d8	11.035	136	163763	20.000	ng	0.00
39) Acenaphthene-d10	14.837	164	109958	20.000	ng	0.00
64) Phenanthrene-d10	17.580	188	243388	20.000	ng	0.00
76) Chrysene-d12	21.870	240	227981	20.000	ng	0.00
86) Perylene-d12	25.254	264	251727	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.747	112	279936	124.735	ng	0.00
7) Phenol-d6	7.363	99	357136	119.546	ng	0.00
23) Nitrobenzene-d5	9.384	82	225064	72.640	ng	0.00
42) 2,4,6-Tribromophenol	16.323	330	180630	116.660	ng	0.00
45) 2-Fluorobiphenyl	13.462	172	546585	74.470	ng	0.00
79) Terphenyl-d14	20.160	244	875264	76.789	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.562	88	34625	36.102	ng	93
3) Pyridine	3.973	79	100735	34.238	ng	96
4) n-Nitrosodimethylamine	3.891	42	64196	41.211	ng	98
6) Aniline	7.522	93	99812	26.600	ng	97
8) 2-Chlorophenol	7.774	128	120349	46.909	ng	100
9) Benzaldehyde	7.334	77	61682m	30.402	ng	
10) Phenol	7.392	94	152609	45.623	ng	99
11) bis(2-Chloroethyl)ether	7.616	93	108003	42.132	ng	100
12) 1,3-Dichlorobenzene	8.097	146	134392	44.554	ng	97
13) 1,4-Dichlorobenzene	8.244	146	135340	44.402	ng	100
14) 1,2-Dichlorobenzene	8.567	146	132049	45.209	ng	98
15) Benzyl Alcohol	8.444	79	109293m	45.015	ng	
16) 2,2'-oxybis(1-Chloropr...	8.732	45	193826	41.768	ng	99
17) 2-Methylphenol	8.656	107	106746	44.949	ng	99
18) Hexachloroethane	9.302	117	48701	46.422	ng	97
19) n-Nitroso-di-n-propyla...	9.014	70	90737	39.531	ng	99
20) 3+4-Methylphenols	8.985	107	145064	43.727	ng	98
22) Acetophenone	9.038	105	182024	42.775	ng	# 98
24) Nitrobenzene	9.425	77	137852	42.595	ng	99
25) Isophorone	9.948	82	253471	43.177	ng	99
26) 2-Nitrophenol	10.142	139	70942	47.485	ng	95
27) 2,4-Dimethylphenol	10.195	122	118779m	52.238	ng	
28) bis(2-Chloroethoxy)met...	10.424	93	144534	45.861	ng	98
29) 2,4-Dichlorophenol	10.683	162	125129	45.977	ng	99
30) 1,2,4-Trichlorobenzene	10.900	180	141045	44.609	ng	98
31) Naphthalene	11.088	128	384664	43.451	ng	99
32) Benzoic acid	10.348	122	63254	33.813	ng	97
33) 4-Chloroaniline	11.194	127	84706	23.195	ng	99
34) Hexachlorobutadiene	11.364	225	95637	44.297	ng	97
35) Caprolactam	11.975	113	37436m	39.750	ng	
36) 4-Chloro-3-methylphenol	12.310	107	130387	43.590	ng	99
37) 2-Methylnaphthalene	12.680	142	291556	43.985	ng	99
38) 1-Methylnaphthalene	12.898	142	275379	42.795	ng	100
40) 1,2,4,5-Tetrachloroben...	13.045	216	167245	48.400	ng	100
41) Hexachlorocyclopentadiene	13.015	237	76326	43.826	ng	98
43) 2,4,6-Trichlorophenol	13.280	196	108036	45.834	ng	97

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.362	196	115792	44.702	ng	97
46) 1,1'-Biphenyl	13.673	154	363414	45.354	ng	99
47) 2-Chloronaphthalene	13.720	162	277076	44.488	ng	98
48) 2-Nitroaniline	13.920	65	88580	43.259	ng	96
49) Acenaphthylene	14.560	152	460780	44.928	ng	100
50) Dimethylphthalate	14.278	163	353480	40.327	ng	99
51) 2,6-Dinitrotoluene	14.408	165	78847	43.993	ng	98
52) Acenaphthene	14.901	154	332434m	49.596	ng	
53) 3-Nitroaniline	14.737	138	61359	31.187	ng	97
54) 2,4-Dinitrophenol	14.954	184	47259	45.861	ng	97
55) Dibenzofuran	15.230	168	478158	46.242	ng	98
56) 4-Nitrophenol	15.054	139	122422	79.280	ng	95
57) 2,4-Dinitrotoluene	15.195	165	109041	43.152	ng	97
58) Fluorene	15.882	166	429502	52.005	ng	98
59) 2,3,4,6-Tetrachlorophenol	15.459	232	104322m	41.564	ng	
60) Diethylphthalate	15.630	149	348726	39.350	ng	100
61) 4-Chlorophenyl-phenyle...	15.859	204	191905	42.271	ng	97
62) 4-Nitroaniline	15.900	138	76145	37.024	ng	95
63) Azobenzene	16.153	77	324213	41.905	ng	98
65) 4,6-Dinitro-2-methylph...	15.959	198	39024	28.714	ng	98
66) n-Nitrosodiphenylamine	16.076	169	304764	45.436	ng	98
67) 4-Bromophenyl-phenylether	16.758	248	124841	45.018	ng	96
68) Hexachlorobenzene	16.881	284	137603	44.750	ng	98
69) Atrazine	17.016	200	125418	47.278	ng	99
70) Pentachlorophenol	17.234	266	146442	77.852	ng	100
71) Phenanthrene	17.622	178	1289794	98.201	ng	98
72) Anthracene	17.710	178	724042	56.737	ng	99
73) Carbazole	17.980	167	588437	51.291	ng	99
74) Di-n-butylphthalate	18.509	149	576456	40.466	ng	100
75) Fluoranthene	19.625	202	1509945	94.497	ng	98
77) Benzidine	19.790	184	238682	45.254	ng	100
78) Pyrene	19.983	202	1437355	93.144	ng	96
80) Butylbenzylphthalate	20.841	149	256902	42.525	ng	96
81) Benzo(a)anthracene	21.852	228	1051935	66.992	ng	98
82) 3,3'-Dichlorobenzidine	21.752	252	163050	29.551	ng	97
83) Chrysene	21.922	228	954383	63.844	ng	99
84) Bis(2-ethylhexyl)phtha...	21.711	149	374048	43.083	ng	99
85) Di-n-octyl phthalate	22.968	149	636358	42.830	ng	100
87) Indeno(1,2,3-cd)pyrene	29.167	276	923124	44.764	ng	# 86
88) Benzo(b)fluoranthene	24.179	252	1078411	65.782	ng	99
89) Benzo(k)fluoranthene	24.243	252	834501	51.067	ng	98
90) Benzo(a)pyrene	25.095	252	930575	66.148	ng	99
91) Dibenzo(a,h)anthracene	29.220	278	652120	38.697	ng	99
92) Benzo(g,h,i)perylene	30.389	276	746987	43.980	ng	99

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

Manual Integrations

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