

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120522\
 Data File : BG055887.D
 Acq On : 5 Dec 2022 16:26
 Operator : CG/JU
 Sample : N5870-04MS
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 ML-4MS

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 12/06/2022
 Supervised By : Jagrut Upadhyay 12/06/2022

Quant Time: Dec 06 04:46:45 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG111622.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 06 04:41:17 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.172	152	27213	20.000	ng	0.00
21) Naphthalene-d8	10.998	136	124201	20.000	ng	0.00
39) Acenaphthene-d10	14.805	164	99744	20.000	ng	0.00
64) Phenanthrene-d10	17.549	188	239610	20.000	ng	0.00
76) Chrysene-d12	21.832	240	221562	20.000	ng	0.00
86) Perylene-d12	25.181	264	235402	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.704	112	194589	129.248	ng	0.00
7) Phenol-d6	7.331	99	265092	132.273	ng	0.00
23) Nitrobenzene-d5	9.347	82	165445	70.407	ng	0.00
42) 2,4,6-Tribromophenol	16.291	330	150965	107.485	ng	0.00
45) 2-Fluorobiphenyl	13.424	172	421076	63.244	ng	0.00
79) Terphenyl-d14	20.134	244	731222	66.010	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.506	88	24411	37.941	ng	95
3) Pyridine	3.929	79	67796	34.348	ng	92
4) n-Nitrosodimethylamine	3.841	42	48646	46.551	ng	# 91
6) Aniline	7.484	93	77644	30.845	ng	96
8) 2-Chlorophenol	7.737	128	82765	48.088	ng	95
9) Benzaldehyde	7.302	77	40231m	29.558	ng	
10) Phenol	7.355	94	110682	49.324	ng	95
11) bis(2-Chloroethyl)ether	7.578	93	72451	42.130	ng	96
12) 1,3-Dichlorobenzene	8.060	146	85078	42.044	ng	97
13) 1,4-Dichlorobenzene	8.207	146	86321	42.215	ng	97
14) 1,2-Dichlorobenzene	8.530	146	84367	43.057	ng	98
15) Benzyl Alcohol	8.412	79	83030m	50.977	ng	
16) 2,2'-oxybis(1-Chloropr...	8.700	45	145035	46.588	ng	96
17) 2-Methylphenol	8.618	107	79135	49.672	ng	98
18) Hexachloroethane	9.264	117	29748	42.269	ng	97
19) n-Nitroso-di-n-propyla...	8.976	70	69982	45.447	ng	93
20) 3+4-Methylphenols	8.947	107	109408	49.161	ng	99
22) Acetophenone	9.000	105	130418	40.410	ng	94
24) Nitrobenzene	9.394	77	100689	41.023	ng	98
25) Isophorone	9.911	82	196316	44.093	ng	98
26) 2-Nitrophenol	10.105	139	48840	43.104	ng	94
27) 2,4-Dimethylphenol	10.157	122	91708m	53.180	ng	
28) bis(2-Chloroethoxy)met...	10.387	93	110044	46.039	ng	99
29) 2,4-Dichlorophenol	10.651	162	95562	46.298	ng	97
30) 1,2,4-Trichlorobenzene	10.857	180	91798	38.281	ng	97
31) Naphthalene	11.051	128	265257	39.507	ng	99
32) Benzoic acid	10.322	122	51013	35.955	ng	93
33) 4-Chloroaniline	11.162	127	60757	21.937	ng	99
34) Hexachlorobutadiene	11.327	225	59456	36.311	ng	98
35) Caprolactam	11.938	113	33348m	46.688	ng	
36) 4-Chloro-3-methylphenol	12.278	107	108795	47.957	ng	97
37) 2-Methylnaphthalene	12.643	142	207995	41.374	ng	99
38) 1-Methylnaphthalene	12.860	142	196544	40.273	ng	97
40) 1,2,4,5-Tetrachloroben...	13.007	216	119571	38.147	ng	99
41) Hexachlorocyclopentadiene	12.984	237	88283	55.882	ng	96
43) 2,4,6-Trichlorophenol	13.248	196	85729	40.095	ng	97

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44) 2,4,5-Trichlorophenol	13.330	196	96203	40.943	ng	99
46) 1,1'-Biphenyl	13.642	154	291980	40.171	ng	99
47) 2-Chloronaphthalene	13.689	162	220595	39.046	ng	99
48) 2-Nitroaniline	13.888	65	81510	43.883	ng	98
49) Acenaphthylene	14.529	152	373728	40.172	ng	99
50) Dimethylphthalate	14.247	163	316442	39.798	ng	99
51) 2,6-Dinitrotoluene	14.376	165	67509	41.524	ng	91
52) Acenaphthene	14.870	154	265124m	43.604	ng	
53) 3-Nitroaniline	14.711	138	41133	23.047	ng	98
54) 2,4-Dinitrophenol	14.928	184	71823	76.836	ng	96
55) Dibenzofuran	15.199	168	372915	39.757	ng	99
56) 4-Nitrophenol	15.028	139	101734	72.629	ng	90
57) 2,4-Dinitrotoluene	15.169	165	96776	42.220	ng	98
58) Fluorene	15.851	166	300628	40.128	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.428	232	85453	37.533	ng	99
60) Diethylphthalate	15.598	149	319251	39.713	ng	100
61) 4-Chlorophenyl-phenyle...	15.833	204	164150	39.860	ng	98
62) 4-Nitroaniline	15.874	138	71869	38.523	ng	95
63) Azobenzene	16.127	77	295261	42.070	ng	97
65) 4,6-Dinitro-2-methylph...	15.933	198	54467	40.709	ng	98
66) n-Nitrosodiphenylamine	16.045	169	272735	41.302	ng	99
67) 4-Bromophenyl-phenylether	16.726	248	108658	39.800	ng	98
68) Hexachlorobenzene	16.855	284	119816	39.580	ng	98
69) Atrazine	16.985	200	112451	43.058	ng	98
70) Pentachlorophenol	17.202	266	112662	60.838	ng	99
71) Phenanthrene	17.590	178	504675	39.030	ng	100
72) Anthracene	17.678	178	517297	41.175	ng	99
73) Carbazole	17.948	167	455672	40.345	ng	99
74) Di-n-butylphthalate	18.477	149	551196	39.303	ng	99
75) Fluoranthene	19.588	202	583067	37.065	ng	99
77) Benzidine	19.758	184	233347	45.525	ng	98
78) Pyrene	19.952	202	592934	39.537	ng	99
80) Butylbenzylphthalate	20.810	149	233719	39.809	ng	98
81) Benzo(a)anthracene	21.808	228	591064	38.732	ng	100
82) 3,3'-Dichlorobenzidine	21.714	252	97422	18.168	ng	97
83) Chrysene	21.879	228	557840	38.398	ng	99
84) Bis(2-ethylhexyl)phtha...	21.679	149	342959	40.647	ng	100
85) Di-n-octyl phthalate	22.925	149	584496	40.479	ng	96
87) Indeno(1,2,3-cd)pyrene	29.047	276	707109	36.667	ng	98
88) Benzo(b)fluoranthene	24.112	252	620075	40.447	ng	100
89) Benzo(k)fluoranthene	24.182	252	601346	39.351	ng	100
90) Benzo(a)pyrene	25.022	252	548000	41.655	ng	100
91) Dibenzo(a,h)anthracene	29.100	278	593094	37.635	ng	99
92) Benzo(g,h,i)perylene	30.263	276	586923	36.952	ng	99

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

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