

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG040623\
 Data File : BG057015.D
 Acq On : 6 Apr 2023 16:54
 Operator : CG/JU
 Sample : 02191-02MS
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 72-11982MS

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 04/07/2023
 Supervised By : Jagrut Upadhyay 04/07/2023

Quant Time: Apr 07 01:37:57 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG033023.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 07 01:33:41 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.408	152	16086	20.000	ng	0.00
21) Naphthalene-d8	11.251	136	62929	20.000	ng	# 0.00
39) Acenaphthene-d10	15.022	164	44734	20.000	ng	0.00
64) Phenanthrene-d10	17.766	188	104155	20.000	ng	0.00
76) Chrysene-d12	22.083	240	88993	20.000	ng	0.00
86) Perylene-d12	25.620	264	96874	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.923	112	96853	96.362	ng	0.00
7) Phenol-d6	7.544	99	143706	95.597	ng	0.00
23) Nitrobenzene-d5	9.588	82	91458	59.468	ng	0.00
42) 2,4,6-Tribromophenol	16.503	330	70635	98.840	ng	0.00
45) 2-Fluorobiphenyl	13.648	172	215392	65.291	ng	0.00
79) Terphenyl-d14	20.327	244	332892	63.298	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.726	88	17238	39.124	ng	95
3) Pyridine	4.149	79	50614	35.666	ng	97
4) n-Nitrosodimethylamine	4.055	42	27841	42.671	ng	99
6) Aniline	7.720	93	68844	35.824	ng	97
8) 2-Chlorophenol	7.967	128	52844	52.785	ng	91
9) Benzaldehyde	7.532	77	37078	39.325	ng	94
10) Phenol	7.574	94	73916	49.624	ng	98
11) bis(2-Chloroethyl)ether	7.809	93	49104	46.280	ng	96
12) 1,3-Dichlorobenzene	8.302	146	58637	53.008	ng	94
13) 1,4-Dichlorobenzene	8.449	146	59311	53.124	ng	98
14) 1,2-Dichlorobenzene	8.772	146	55535	51.486	ng	96
15) Benzyl Alcohol	8.643	79	57747	46.232	ng	92
16) 2,2'-oxybis(1-Chloropr...	8.931	45	58001m	45.359	ng	
17) 2-Methylphenol	8.842	107	50048	47.096	ng	93
18) Hexachloroethane	9.512	117	21150	51.830	ng	# 82
19) n-Nitroso-di-n-propyla...	9.212	70	46067	42.844	ng	99
20) 3+4-Methylphenols	9.171	107	67508	45.430	ng	97
22) Acetophenone	9.242	105	86230	46.050	ng	# 97
24) Nitrobenzene	9.630	77	71097	45.216	ng	99
25) Isophorone	10.152	82	122773	41.921	ng	96
26) 2-Nitrophenol	10.346	139	29964	49.371	ng	92
27) 2,4-Dimethylphenol	10.393	122	53092	49.546	ng	96
28) bis(2-Chloroethoxy)met...	10.628	93	63686	43.436	ng	94
29) 2,4-Dichlorophenol	10.887	162	57002	49.463	ng	97
30) 1,2,4-Trichlorobenzene	11.110	180	61874	50.531	ng	97
31) Naphthalene	11.304	128	165056	48.120	ng	99
32) Benzoic acid	10.534	122	35383m	48.070	ng	
33) 4-Chloroaniline	11.398	127	43520	28.093	ng	97
34) Hexachlorobutadiene	11.574	225	45834	49.830	ng	98
35) Caprolactam	12.173	113	17095m	43.014	ng	
36) 4-Chloro-3-methylphenol	12.490	107	60241	46.742	ng	100
37) 2-Methylnaphthalene	12.878	142	121719	44.712	ng	100
38) 1-Methylnaphthalene	13.095	142	113209	44.287	ng	100
40) 1,2,4,5-Tetrachloroben...	13.236	216	82847	53.407	ng	98
41) Hexachlorocyclopentadiene	13.213	237	99140	115.994	ng	94
43) 2,4,6-Trichlorophenol	13.466	196	51596	52.122	ng	97

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.542	196	56303	48.011	ng	98
46) 1,1'-Biphenyl	13.859	154	170989	52.216	ng	97
47) 2-Chloronaphthalene	13.912	162	128360	52.999	ng	97
48) 2-Nitroaniline	14.106	65	41510	47.932	ng	98
49) Acenaphthylene	14.752	152	201427	50.619	ng	99
50) Dimethylphthalate	14.458	163	162883	46.212	ng	# 98
51) 2,6-Dinitrotoluene	14.588	165	36212	47.550	ng	96
52) Acenaphthene	15.087	154	157458m	57.374	ng	
53) 3-Nitroaniline	14.917	138	28457	37.751	ng	93
54) 2,4-Dinitrophenol	15.128	184	49917	116.452	ng	93
55) Dibenzofuran	15.422	168	201797	48.991	ng	96
56) 4-Nitrophenol	15.216	139	58313	104.615	ng	89
57) 2,4-Dinitrotoluene	15.369	165	52213	48.612	ng	100
58) Fluorene	16.062	166	167087	49.021	ng	98
59) 2,3,4,6-Tetrachlorophenol	15.639	232	52926	49.568	ng	# 99
60) Diethylphthalate	15.809	149	164167	46.460	ng	99
61) 4-Chlorophenyl-phenyle...	16.044	204	96349	48.146	ng	95
62) 4-Nitroaniline	16.080	138	35808	45.849	ng	88
63) Azobenzene	16.338	77	166200	47.196	ng	98
65) 4,6-Dinitro-2-methylph...	16.133	198	35630	56.483	ng	98
66) n-Nitrosodiphenylamine	16.256	169	138616	49.577	ng	99
67) 4-Bromophenyl-phenylether	16.937	248	62052	48.150	ng	98
68) Hexachlorobenzene	17.067	284	68909	54.084	ng	99
69) Atrazine	17.190	200	60612	52.224	ng	97
70) Pentachlorophenol	17.407	266	88678	97.015	ng	94
71) Phenanthrene	17.807	178	362368	68.239	ng	99
72) Anthracene	17.895	178	271277	49.833	ng	99
73) Carbazole	18.159	167	235771	49.675	ng	97
74) Di-n-butylphthalate	18.682	149	257167	46.791	ng	100
75) Fluoranthene	19.792	202	473401	70.325	ng	100
77) Benzidine	19.957	184	80677	33.167	ng	98
78) Pyrene	20.156	202	474099	77.024	ng	98
80) Butylbenzylphthalate	21.008	149	103128	48.689	ng	95
81) Benzo(a)anthracene	22.060	228	381211	61.902	ng	99
82) 3,3'-Dichlorobenzidine	21.954	252	69713	33.523	ng	100
83) Chrysene	22.136	228	365034	63.307	ng	99
84) Bis(2-ethylhexyl)phtha...	21.913	149	149980	48.899	ng	98
85) Di-n-octyl phthalate	23.223	149	246361	47.665	ng	99
87) Indeno(1,2,3-cd)pyrene	29.738	276	414034	57.830	ng	# 97
88) Benzo(b)fluoranthene	24.498	252	417912	68.987	ng	99
89) Benzo(k)fluoranthene	24.568	252	340483	56.300	ng	99
90) Benzo(a)pyrene	25.461	252	347152	58.318	ng	99
91) Dibenzo(a,h)anthracene	29.791	278	306321	52.052	ng	97
92) Benzo(g,h,i)perylene	31.018	276	342090	59.365	ng	97

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

