

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111622\
 Data File : BG055652.D
 Acq On : 16 Nov 2022 18:19
 Operator : CG/JU
 Sample : N5607-13MS
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :

BNA_G

ClientSampleId :

SB-40BMS

Manual Integrations

APPROVED

Reviewed By :Christian Giraldo 11/17/2022
 Supervised By :mohammad ahmed 11/17/2022

Quant Time: Nov 17 02:12:00 2022

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG111622.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Thu Nov 17 02:02:46 2022

Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.252	152	40524	20.000	ng	0.00
21) Naphthalene-d8	11.084	136	174513	20.000	ng	0.00
39) Acenaphthene-d10	14.874	164	130512	20.000	ng	0.00
64) Phenanthrene-d10	17.618	188	291894	20.000	ng	0.00
76) Chrysene-d12	21.919	240	282166	20.000	ng	0.00
86) Perylene-d12	25.326	264	303285	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.779	112	290289	129.479	ng	0.00
7) Phenol-d6	7.394	99	390451	130.830	ng	0.00
23) Nitrobenzene-d5	9.427	82	232488	70.414	ng	0.00
42) 2,4,6-Tribromophenol	16.360	330	206731	112.490	ng	0.00
45) 2-Fluorobiphenyl	13.499	172	602204	69.126	ng	0.00
79) Terphenyl-d14	20.203	244	991354	70.272	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.616	88	39692	41.427	ng	97
3) Pyridine	4.028	79	109099m	37.118	ng	
4) n-Nitrosodimethylamine	3.934	42	67351	43.281	ng	98
6) Aniline	7.565	93	84649	22.582	ng	97
8) 2-Chlorophenol	7.812	128	132175	51.571	ng	97
9) Benzaldehyde	7.377	77	75615	37.307	ng	98
10) Phenol	7.418	94	172779	51.705	ng	98
11) bis(2-Chloroethyl)ether	7.659	93	116765	45.596	ng	99
12) 1,3-Dichlorobenzene	8.141	146	135074	44.826	ng	97
13) 1,4-Dichlorobenzene	8.287	146	135568	44.522	ng	99
14) 1,2-Dichlorobenzene	8.611	146	133630	45.797	ng	99
15) Benzyl Alcohol	8.487	79	122763m	50.614	ng	
16) 2,2'-oxybis(1-Chloropr...	8.781	45	210219	45.346	ng	97
17) 2-Methylphenol	8.687	107	122227	51.520	ng	99
18) Hexachloroethane	9.351	117	47126	44.966	ng	98
19) n-Nitroso-di-n-propyla...	9.057	70	104320	45.494	ng	97
20) 3+4-Methylphenols	9.016	107	168539	50.855	ng	96
22) Acetophenone	9.081	105	198279	43.725	ng	# 98
24) Nitrobenzene	9.468	77	149775	43.429	ng	99
25) Isophorone	9.991	82	290676	46.464	ng	99
26) 2-Nitrophenol	10.185	139	75346	47.326	ng	98
27) 2,4-Dimethylphenol	10.232	122	139152	57.429	ng	99
28) bis(2-Chloroethoxy)met...	10.467	93	166316	49.521	ng	98
29) 2,4-Dichlorophenol	10.720	162	142994	49.305	ng	98
30) 1,2,4-Trichlorobenzene	10.943	180	143722	42.655	ng	99
31) Naphthalene	11.137	128	406249	43.063	ng	99
32) Benzoic acid	10.361	122	103190	51.763	ng	96
33) 4-Chloroaniline	11.231	127	70520	18.121	ng	99
34) Hexachlorobutadiene	11.413	225	94259	40.969	ng	97
35) Caprolactam	12.007	113	47732m	47.560	ng	
36) 4-Chloro-3-methylphenol	12.336	107	158752	49.803	ng	99
37) 2-Methylnaphthalene	12.723	142	310099	43.901	ng	99
38) 1-Methylnaphthalene	12.941	142	294937	43.011	ng	99
40) 1,2,4,5-Tetrachloroben...	13.082	216	177151	43.193	ng	98
41) Hexachlorocyclopentadiene	13.058	237	226302	109.477	ng	99
43) 2,4,6-Trichlorophenol	13.317	196	129044	46.125	ng	98

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.387	196	139778	45.464	ng	98
46) 1,1'-Biphenyl	13.711	154	428432	45.048	ng	99
47) 2-Chloronaphthalene	13.763	162	322112	43.574	ng	99
48) 2-Nitroaniline	13.957	65	109985	45.254	ng	98
49) Acenaphthylene	14.604	152	534059	43.872	ng	99
50) Dimethylphthalate	14.322	163	439521	42.246	ng	99
51) 2,6-Dinitrotoluene	14.439	165	95970	45.114	ng	96
52) Acenaphthene	14.938	154	385853m	48.499	ng	
53) 3-Nitroaniline	14.768	138	45602	19.528	ng	97
54) 2,4-Dinitrophenol	14.974	184	119540	97.736	ng	99
55) Dibenzofuran	15.273	168	528923	43.095	ng	100
56) 4-Nitrophenol	15.068	139	163479	89.196	ng	96
57) 2,4-Dinitrotoluene	15.226	165	133138	44.391	ng	97
58) Fluorene	15.920	166	415669	42.404	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.491	232	127691	42.863	ng	99
60) Diethylphthalate	15.673	149	437487	41.592	ng	100
61) 4-Chlorophenyl-phenyle...	15.902	204	232839	43.211	ng	99
62) 4-Nitroaniline	15.931	138	95885	39.279	ng	98
63) Azobenzene	16.196	77	388560	42.312	ng	99
65) 4,6-Dinitro-2-methylph...	15.984	198	77078	47.290	ng	98
66) n-Nitrosodiphenylamine	16.114	169	371641	46.199	ng	99
67) 4-Bromophenyl-phenylether	16.795	248	153017	46.009	ng	99
68) Hexachlorobenzene	16.919	284	166760	45.220	ng	99
69) Atrazine	17.054	200	150287	47.238	ng	99
70) Pentachlorophenol	17.259	266	211513	93.759	ng	98
71) Phenanthrene	17.659	178	677871	43.034	ng	100
72) Anthracene	17.753	178	687423	44.916	ng	99
73) Carbazole	18.011	167	613803	44.611	ng	98
74) Di-n-butylphthalate	18.558	149	734444	42.989	ng	100
75) Fluoranthene	19.657	202	783244	40.872	ng	99
77) Benzidine	19.827	184	198724	30.443	ng	98
78) Pyrene	20.015	202	807255	42.266	ng	99
80) Butylbenzylphthalate	20.884	149	334458	44.732	ng	100
81) Benzo(a)anthracene	21.895	228	833143	42.869	ng	99
82) 3,3'-Dichlorobenzidine	21.795	252	139430	20.417	ng	99
83) Chrysene	21.966	228	783214	42.333	ng	99
84) Bis(2-ethylhexyl)phtha...	21.778	149	483308	44.978	ng	99
85) Di-n-octyl phthalate	23.053	149	828569	45.058	ng	100
87) Indeno(1,2,3-cd)pyrene	29.263	276	1012170	40.738	ng	# 94
88) Benzo(b)fluoranthene	24.239	252	886369	44.876	ng	99
89) Benzo(k)fluoranthene	24.310	252	851769	43.262	ng	99
90) Benzo(a)pyrene	25.162	252	781052	46.081	ng	99
91) Dibenzo(a,h)anthracene	29.328	278	854898	42.106	ng	99
92) Benzo(g,h,i)perylene	30.497	276	839060	41.003	ng	98

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

