

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG042722\
 Data File : BG053342.D
 Acq On : 27 Apr 2022 11:35
 Operator : CG/JU
 Sample : PB144386BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :

BNA_G

ClientSampleId :

PB144386BS

Manual Integrations

APPROVED

Reviewed By : Christian Giraldo 04/28/2022
 Supervised By : Jagrut Upadhyay 04/28/2022

Quant Time: Apr 28 06:40:13 2022

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG040822.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Sun Apr 24 01:13:30 2022

Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.101	152	24541	20.000	ng	-0.01
21) Naphthalene-d8	10.914	136	100382	20.000	ng	0.00
39) Acenaphthene-d10	14.727	164	77455	20.000	ng	0.00
64) Phenanthrene-d10	17.476	188	194816	20.000	ng	0.00
76) Chrysene-d12	21.759	240	193278	20.000	ng	0.00
86) Perylene-d12	25.054	264	208583	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.669	112	205135	143.286	ng	0.00
7) Phenol-d6	7.266	99	288023	130.460	ng	0.00
23) Nitrobenzene-d5	9.264	82	199159	80.558	ng	-0.01
42) 2,4,6-Tribromophenol	16.219	330	152024	123.368	ng	0.00
45) 2-Fluorobiphenyl	13.352	172	454338	80.884	ng	0.00
79) Terphenyl-d14	20.079	244	965167	84.338	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.548	88	23961	35.046	ng	# 92
3) Pyridine	3.947	79	62577	34.703	ng	90
4) n-Nitrosodimethylamine	3.859	42	40873	45.773	ng	97
6) Aniline	7.425	93	106915	41.786	ng	98
8) 2-Chlorophenol	7.672	128	78443	50.753	ng	95
9) Benzaldehyde	7.237	77	55424m	41.998	ng	
10) Phenol	7.296	94	107287	48.821	ng	91
11) bis(2-Chloroethyl)ether	7.513	93	72948	46.049	ng	93
12) 1,3-Dichlorobenzene	7.995	146	78471m	43.692	ng	
13) 1,4-Dichlorobenzene	8.142	146	80604	44.022	ng	94
14) 1,2-Dichlorobenzene	8.459	146	78238	44.311	ng	99
15) Benzyl Alcohol	8.341	79	87939m	44.813	ng	
16) 2,2'-oxybis(1-Chloropr...	8.623	45	116100	43.898	ng	99
17) 2-Methylphenol	8.547	107	72009	48.422	ng	96
18) Hexachloroethane	9.187	117	30527	44.460	ng	96
19) n-Nitroso-di-n-propyla...	8.905	70	70816	43.813	ng	# 93
20) 3+4-Methylphenols	8.876	107	97598	46.491	ng	91
22) Acetophenone	8.923	105	123355	44.294	ng	94
24) Nitrobenzene	9.311	77	107915	44.877	ng	92
25) Isophorone	9.828	82	196753	46.823	ng	97
26) 2-Nitrophenol	10.022	139	45727	45.475	ng	95
27) 2,4-Dimethylphenol	10.080	122	78006	54.167	ng	93
28) bis(2-Chloroethoxy)met...	10.309	93	104240	44.155	ng	97
29) 2,4-Dichlorophenol	10.562	162	84848	46.866	ng	99
30) 1,2,4-Trichlorobenzene	10.779	180	88376	43.130	ng	97
31) Naphthalene	10.967	128	237381	44.324	ng	99
32) Benzoic acid	10.239	122	36981	40.103	ng	91
33) 4-Chloroaniline	11.073	127	73156	31.898	ng	97
34) Hexachlorobutadiene	11.249	225	62830	41.279	ng	96
35) Caprolactam	11.837	113	30022m	47.046	ng	
36) 4-Chloro-3-methylphenol	12.189	107	98369	46.671	ng	93
37) 2-Methylnaphthalene	12.565	142	174089m	43.930	ng	
38) 1-Methylnaphthalene	12.783	142	169204	43.742	ng	98
40) 1,2,4,5-Tetrachloroben...	12.935	216	112886	42.912	ng	97
41) Hexachlorocyclopentadiene	12.912	237	124407	91.611	ng	92
43) 2,4,6-Trichlorophenol	13.170	196	76789	42.337	ng	95

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.247	196	84067	43.496	ng	99
46) 1,1'-Biphenyl	13.564	154	238485	42.756	ng	98
47) 2-Chloronaphthalene	13.611	162	188225	42.297	ng	97
48) 2-Nitroaniline	13.805	65	75583	44.755	ng	94
49) Acenaphthylene	14.451	152	329019	47.231	ng	98
50) Dimethylphthalate	14.175	163	274216	44.263	ng	98
51) 2,6-Dinitrotoluene	14.292	165	58616	45.354	ng	# 82
52) Acenaphthene	14.792	154	232763m	49.271	ng	
53) 3-Nitroaniline	14.627	138	48718	35.649	ng	96
54) 2,4-Dinitrophenol	14.839	184	69930	85.037	ng	# 88
55) Dibenzofuran	15.126	168	314751	42.550	ng	98
56) 4-Nitrophenol	14.944	139	96560	92.643	ng	85
57) 2,4-Dinitrotoluene	15.085	165	87014	47.290	ng	87
58) Fluorene	15.773	166	263099	44.037	ng	97
59) 2,3,4,6-Tetrachlorophenol	15.356	232	79754	42.491	ng	97
60) Diethylphthalate	15.532	149	282118	43.463	ng	100
61) 4-Chlorophenyl-phenyle...	15.761	204	143621	42.406	ng	96
62) 4-Nitroaniline	15.790	138	63435	43.989	ng	# 80
63) Azobenzene	16.055	77	284281	43.439	ng	99
65) 4,6-Dinitro-2-methylph...	15.855	198	53010	39.002	ng	96
66) n-Nitrosodiphenylamine	15.972	169	242494	43.285	ng	99
67) 4-Bromophenyl-phenylether	16.654	248	103840	41.598	ng	94
68) Hexachlorobenzene	16.783	284	115184	42.608	ng	93
69) Atrazine	16.918	200	106616	48.720	ng	100
70) Pentachlorophenol	17.130	266	123656	83.696	ng	91
71) Phenanthrene	17.517	178	458625	43.099	ng	98
72) Anthracene	17.611	178	467598	44.355	ng	99
73) Carbazole	17.876	167	434852	42.434	ng	99
74) Di-n-butylphthalate	18.422	149	505301	43.477	ng	99
75) Fluoranthene	19.520	202	607615	44.546	ng	99
77) Benzidine	19.697	184	315409	57.196	ng	99
78) Pyrene	19.885	202	606729	44.077	ng	99
80) Butylbenzylphthalate	20.760	149	228631	44.551	ng	99
81) Benzo(a)anthracene	21.741	228	598002	45.252	ng	97
82) 3,3'-Dichlorobenzidine	21.647	252	162333	33.402	ng	100
83) Chrysene	21.806	228	546753	44.580	ng	99
84) Bis(2-ethylhexyl)phtha...	21.629	149	319557	46.067	ng	100
85) Di-n-octyl phthalate	22.863	149	541619	46.685	ng	97
87) Indeno(1,2,3-cd)pyrene	28.826	276	674349	43.516	ng	98
88) Benzo(b)fluoranthene	24.009	252	642734	48.839	ng	98
89) Benzo(k)fluoranthene	24.079	252	582736	45.054	ng	99
90) Benzo(a)pyrene	24.901	252	603823	53.858	ng	98
91) Dibenzo(a,h)anthracene	28.890	278	579121	45.405	ng	98
92) Benzo(g,h,i)perylene	30.018	276	578270	45.999	ng	97

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

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