

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG042922\
 Data File : BG053377.D
 Acq On : 29 Apr 2022 14:55
 Operator : CG/JU
 Sample : PB144441BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB144441BS

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 05/02/2022
 Supervised By : Jagrut Upadhyay 05/02/2022

Quant Time: May 02 01:02:44 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	8.105	152	29659	20.000	ng	0.00
21) Naphthalene-d8	10.913	136	116023	20.000	ng	0.00
39) Acenaphthene-d10	14.726	164	88016	20.000	ng	0.00
64) Phenanthrene-d10	17.475	188	210776	20.000	ng	0.00
76) Chrysene-d12	21.757	240	218592	20.000	ng	0.00
86) Perylene-d12	25.053	264	228945	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.673	112	228445	132.033	ng	0.00
7) Phenol-d6	7.271	99	315641	118.299	ng	0.00
23) Nitrobenzene-d5	9.263	82	218244	76.377	ng	-0.01
42) 2,4,6-Tribromophenol	16.218	330	153179	109.390	ng	0.00
45) 2-Fluorobiphenyl	13.351	172	474338	74.312	ng	0.00
79) Terphenyl-d14	20.077	244	975391	75.361	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.547	88	26873	32.523	ng	# 89
3) Pyridine	3.952	79	67733	31.081	ng	93
4) n-Nitrosodimethylamine	3.864	42	45871	42.505	ng	# 94
6) Aniline	7.424	93	112500	36.382	ng	95
8) 2-Chlorophenol	7.671	128	80139	42.903	ng	94
9) Benzaldehyde	7.236	77	61503m	38.562	ng	
10) Phenol	7.295	94	112307	42.286	ng	90
11) bis(2-Chloroethyl)ether	7.512	93	77478	40.469	ng	95
12) 1,3-Dichlorobenzene	7.994	146	88086m	40.582	ng	
13) 1,4-Dichlorobenzene	8.140	146	89287	40.350	ng	96
14) 1,2-Dichlorobenzene	8.458	146	84759	39.721	ng	96
15) Benzyl Alcohol	8.340	79	90302m	38.076	ng	
16) 2,2'-oxybis(1-Chloropr...	8.622	45	125639	39.307	ng	97
17) 2-Methylphenol	8.546	107	73600	40.952	ng	93
18) Hexachloroethane	9.192	117	33920	40.877	ng	96
19) n-Nitroso-di-n-propyla...	8.904	70	72972	37.357	ng	# 97
20) 3+4-Methylphenols	8.875	107	99668	39.284	ng	92
22) Acetophenone	8.922	105	126302	39.239	ng	94
24) Nitrobenzene	9.310	77	111922	40.269	ng	91
25) Isophorone	9.832	82	195643	40.282	ng	98
26) 2-Nitrophenol	10.020	139	44571	38.350	ng	# 82
27) 2,4-Dimethylphenol	10.079	122	77122	46.333	ng	92
28) bis(2-Chloroethoxy)met...	10.308	93	103545	37.948	ng	97
29) 2,4-Dichlorophenol	10.561	162	84652	40.454	ng	99
30) 1,2,4-Trichlorobenzene	10.778	180	91843	38.780	ng	96
31) Naphthalene	10.966	128	242678	39.205	ng	99
32) Benzoic acid	10.244	122	37924m	35.991	ng	
33) 4-Chloroaniline	11.072	127	83089	31.345	ng	95
34) Hexachlorobutadiene	11.248	225	66961	38.062	ng	98
35) Caprolactam	11.836	113	27541m	37.340	ng	
36) 4-Chloro-3-methylphenol	12.188	107	96315	39.536	ng	94
37) 2-Methylnaphthalene	12.564	142	175049	38.217	ng	95
38) 1-Methylnaphthalene	12.781	142	167283	37.416	ng	93
40) 1,2,4,5-Tetrachloroben...	12.928	216	111754	37.384	ng	99
41) Hexachlorocyclopentadiene	12.911	237	122829	79.596	ng	94
43) 2,4,6-Trichlorophenol	13.169	196	76490	37.112	ng	96

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.245	196	80596	36.696	ng	98
46) 1,1'-Biphenyl	13.563	154	239358	37.763	ng	99
47) 2-Chloronaphthalene	13.610	162	185263	36.636	ng	97
48) 2-Nitroaniline	13.803	65	75749	39.472	ng	95
49) Acenaphthylene	14.450	152	320041	40.429	ng	99
50) Dimethylphthalate	14.174	163	261264	37.112	ng	100
51) 2,6-Dinitrotoluene	14.297	165	56984	38.800	ng	# 81
52) Acenaphthene	14.790	154	222754m	41.495	ng	
53) 3-Nitroaniline	14.632	138	50176	32.310	ng	98
54) 2,4-Dinitrophenol	14.843	184	66449	71.109	ng	93
55) Dibenzofuran	15.125	168	306797	36.498	ng	98
56) 4-Nitrophenol	14.943	139	89254	75.359	ng	# 80
57) 2,4-Dinitrotoluene	15.084	165	83204	39.793	ng	88
58) Fluorene	15.771	166	255497	37.633	ng	97
59) 2,3,4,6-Tetrachlorophenol	15.354	232	77113	36.154	ng	99
60) Diethylphthalate	15.531	149	274513	37.217	ng	99
61) 4-Chlorophenyl-phenyle...	15.760	204	142665	37.070	ng	98
62) 4-Nitroaniline	15.789	138	61437	37.491	ng	# 80
63) Azobenzene	16.053	77	269466	36.235	ng	97
65) 4,6-Dinitro-2-methylph...	15.859	198	51281	34.873	ng	95
66) n-Nitrosodiphenylamine	15.971	169	228790	37.747	ng	99
67) 4-Bromophenyl-phenylether	16.653	248	101809	37.696	ng	95
68) Hexachlorobenzene	16.788	284	112016	38.298	ng	96
69) Atrazine	16.917	200	103124	43.556	ng	99
70) Pentachlorophenol	17.128	266	118118	73.894	ng	92
71) Phenanthrene	17.516	178	441051	38.309	ng	97
72) Anthracene	17.610	178	448529	39.325	ng	99
73) Carbazole	17.874	167	414334	37.371	ng	98
74) Di-n-butylphthalate	18.421	149	489234	38.907	ng	99
75) Fluoranthene	19.525	202	586341	39.731	ng	99
77) Benzidine	19.695	184	286791	45.984	ng	98
78) Pyrene	19.883	202	582096	37.390	ng	98
80) Butylbenzylphthalate	20.759	149	218325	37.616	ng	97
81) Benzo(a)anthracene	21.740	228	581974	38.939	ng	99
82) 3,3'-Dichlorobenzidine	21.646	252	172522	31.387	ng	99
83) Chrysene	21.804	228	528480	38.100	ng	99
84) Bis(2-ethylhexyl)phtha...	21.628	149	312333	39.811	ng	96
85) Di-n-octyl phthalate	22.862	149	527597	40.210	ng	97
87) Indeno(1,2,3-cd)pyrene	28.830	276	656722	38.609	ng	# 96
88) Benzo(b)fluoranthene	24.007	252	610366	42.254	ng	97
89) Benzo(k)fluoranthene	24.078	252	565299	39.819	ng	98
90) Benzo(a)pyrene	24.900	252	581626	47.264	ng	99
91) Dibenzo(a,h)anthracene	28.889	278	567511	40.538	ng	98
92) Benzo(g,h,i)perylene	30.011	276	565531	40.985	ng	98

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

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