

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG031324\
 Data File : BG060625.D
 Acq On : 13 Mar 2024 13:26
 Operator : MA/JU
 Sample : SSTDICCC040
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDICCC040

Quant Time: Mar 14 00:36:20 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG031324.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 14 00:26:58 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.324	152	106013	20.000	ng	0.00
21) Naphthalene-d8	11.173	136	497769	20.000	ng	0.00
39) Acenaphthene-d10	14.951	164	328973	20.000	ng	0.00
64) Phenanthrene-d10	17.701	188	696137	20.000	ng	0.00
76) Chrysene-d12	22.025	240	617246	20.000	ng	0.00
86) Perylene-d12	25.574	264	701669	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.809	112	550683	81.657	ng	0.00
7) Phenol-d6	7.437	99	812780	83.083	ng	0.00
23) Nitrobenzene-d5	9.528	82	792109	82.818	ng	0.00
42) 2,4,6-Tribromophenol	16.438	330	323020	84.735	ng	0.00
45) 2-Fluorobiphenyl	13.577	172	1935969	79.566	ng	0.00
79) Terphenyl-d14	20.280	244	2705895	79.239	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.653	88	110507	39.147	ng	100
3) Pyridine	4.070	79	336570	40.342	ng	100
4) n-Nitrosodimethylamine	4.000	42	164505	40.196	ng	100
6) Aniline	7.648	93	493519	41.285	ng	100
8) 2-Chlorophenol	7.871	128	303349	41.149	ng	100
9) Benzaldehyde	7.472	77	229773	41.958	ng	100
10) Phenol	7.466	94	409643	41.733	ng	100
11) bis(2-Chloroethyl)ether	7.742	93	321460	41.082	ng	100
12) 1,3-Dichlorobenzene	8.206	146	318132	40.017	ng	100
13) 1,4-Dichlorobenzene	8.365	146	330113	40.964	ng	100
14) 1,2-Dichlorobenzene	8.682	146	322553	40.301	ng	100
15) Benzyl Alcohol	8.565	79	328405	42.462	ng	100
16) 2,2'-oxybis(1-Chloropr...	8.847	45	622027	40.950	ng	100
17) 2-Methylphenol	8.741	107	317128	42.074	ng	100
18) Hexachloroethane	9.411	117	111624	40.549	ng	100
19) n-Nitroso-di-n-propyla...	9.135	70	284355	41.934	ng	100
20) 3+4-Methylphenols	9.076	107	428947	42.196	ng	100
22) Acetophenone	9.170	105	545271	40.858	ng	# 100
24) Nitrobenzene	9.569	77	398820	41.708	ng	100
25) Isophorone	10.075	82	784387	40.975	ng	100
26) 2-Nitrophenol	10.275	139	151954	39.639	ng	100
27) 2,4-Dimethylphenol	10.298	122	335920	39.647	ng	100
28) bis(2-Chloroethoxy)met...	10.562	93	464038	40.435	ng	100
29) 2,4-Dichlorophenol	10.792	162	315062	41.432	ng	100
30) 1,2,4-Trichlorobenzene	11.015	180	320069	39.642	ng	100
31) Naphthalene	11.226	128	1108429	39.867	ng	100
32) Benzoic acid	10.416	122	216291	39.620	ng	100
33) 4-Chloroaniline	11.338	127	479022	40.859	ng	100
34) Hexachlorobutadiene	11.456	225	220492	40.037	ng	100
35) Caprolactam	12.137	113	105104	41.059	ng	100
36) 4-Chloro-3-methylphenol	12.401	107	383623	41.166	ng	100
37) 2-Methylnaphthalene	12.801	142	769502	40.072	ng	100
38) 1-Methylnaphthalene	13.018	142	728241	40.404	ng	100
40) 1,2,4,5-Tetrachloroben...	13.142	216	403057	39.896	ng	100
41) Hexachlorocyclopentadiene	13.101	237	207004	41.666	ng	100
43) 2,4,6-Trichlorophenol	13.383	196	266406	41.067	ng	100

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.453	196	318571	41.513	ng	100
46) 1,1'-Biphenyl	13.794	154	1000255	40.001	ng	100
47) 2-Chloronaphthalene	13.847	162	769343	40.767	ng	100
48) 2-Nitroaniline	14.058	65	270577	43.441	ng	100
49) Acenaphthylene	14.687	152	1237694	40.259	ng	100
50) Dimethylphthalate	14.399	163	1001529	40.593	ng	100
51) 2,6-Dinitrotoluene	14.534	165	194157	42.708	ng	100
52) Acenaphthene	15.016	154	732296	39.861	ng	100
53) 3-Nitroaniline	14.875	138	218189	41.683	ng	100
54) 2,4-Dinitrophenol	15.069	184	80681	38.544	ng	100
55) Dibenzofuran	15.351	168	1186967	39.904	ng	100
56) 4-Nitrophenol	15.134	139	161680	41.393	ng	100
57) 2,4-Dinitrotoluene	15.310	165	245584	42.941	ng	100
58) Fluorene	15.997	166	956224	40.180	ng	100
59) 2,3,4,6-Tetrachlorophenol	15.551	232	267240m	41.648	ng	
60) Diethylphthalate	15.745	149	1022382	40.274	ng	100
61) 4-Chlorophenyl-phenyle...	15.980	204	479922	39.812	ng	100
62) 4-Nitroaniline	16.038	138	228536	42.647	ng	100
63) Azobenzene	16.273	77	1066263	40.586	ng	100
65) 4,6-Dinitro-2-methylph...	16.062	198	121169	39.470	ng	100
66) n-Nitrosodiphenylamine	16.197	169	878417	40.319	ng	100
67) 4-Bromophenyl-phenylether	16.879	248	296170	39.453	ng	100
68) Hexachlorobenzene	16.967	284	343833	39.837	ng	100
69) Atrazine	17.131	200	298049	40.902	ng	100
70) Pentachlorophenol	17.319	266	236904	42.349	ng	100
71) Phenanthrene	17.742	178	1491147	39.765	ng	100
72) Anthracene	17.836	178	1520851	40.112	ng	100
73) Carbazole	18.107	167	1327862	40.278	ng	100
74) Di-n-butylphthalate	18.618	149	1668496	41.530	ng	100
75) Fluoranthene	19.734	202	1687334	40.678	ng	100
77) Benzidine	19.922	184	673112	44.583	ng	100
78) Pyrene	20.098	202	1745868	40.530	ng	100
80) Butylbenzylphthalate	20.962	149	687362	42.209	ng	100
81) Benzo(a)anthracene	22.002	228	1731050	40.731	ng	100
82) 3,3'-Dichlorobenzidine	21.908	252	609950	42.580	ng	100
83) Chrysene	22.072	228	1661759	40.269	ng	100
84) Bis(2-ethylhexyl)phtha...	21.843	149	1019512	42.384	ng	100
85) Di-n-octyl phthalate	23.177	149	1697272	42.849	ng	100
87) Indeno(1,2,3-cd)pyrene	29.711	276	1958915	40.946	ng	100
88) Benzo(b)fluoranthene	24.429	252	1595256	41.326	ng	100
89) Benzo(k)fluoranthene	24.505	252	1633488	40.280	ng	100
90) Benzo(a)pyrene	25.410	252	1544804	40.819	ng	100
91) Dibenzo(a,h)anthracene	29.805	278	1658306	40.793	ng	100
92) Benzo(g,h,i)perylene	31.015	276	1619419	41.166	ng	100

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

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