

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG033023\
 Data File : BG056924.D
 Acq On : 30 Mar 2023 14:42
 Operator : CG/JU
 Sample : SSTDICCC040
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDICCC040

Quant Time: Mar 30 15:50:48 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG033023.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 30 15:48:01 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.416	152	20823	20.000	ng	0.00
21) Naphthalene-d8	11.259	136	83766	20.000	ng	0.00
39) Acenaphthene-d10	15.031	164	66418	20.000	ng	0.00
64) Phenanthrene-d10	17.774	188	165056	20.000	ng	0.00
76) Chrysene-d12	22.097	240	154024	20.000	ng	0.00
86) Perylene-d12	25.646	264	168601	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.919	112	105556	82.479	ng	0.00
7) Phenol-d6	7.541	99	163184	86.321	ng	0.00
23) Nitrobenzene-d5	9.597	82	165743	101.309	ng	0.00
42) 2,4,6-Tribromophenol	16.511	330	87406	96.120	ng	0.00
45) 2-Fluorobiphenyl	13.662	172	390718	78.773	ng	0.00
79) Terphenyl-d14	20.341	244	720691	76.555	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.728	88	24078	42.914	ng	100
3) Pyridine	4.151	79	77641	43.748	ng	100
4) n-Nitrosodimethylamine	4.057	42	34752	42.009	ng	100
6) Aniline	7.723	93	104127	39.633	ng	100
8) 2-Chlorophenol	7.969	128	52372	41.360	ng	100
9) Benzaldehyde	7.535	77	52037	52.050	ng	100
10) Phenol	7.570	94	78721	40.708	ng	100
11) bis(2-Chloroethyl)ether	7.811	93	56272	38.783	ng	100
12) 1,3-Dichlorobenzene	8.304	146	57627	38.198	ng	100
13) 1,4-Dichlorobenzene	8.451	146	58403	38.440	ng	100
14) 1,2-Dichlorobenzene	8.780	146	56723	38.841	ng	100
15) Benzyl Alcohol	8.645	79	67177	39.686	ng	100
16) 2,2'-oxybis(1-Chloropr...	8.933	45	65625	33.689	ng	100
17) 2-Methylphenol	8.845	107	56491	41.569	ng	100
18) Hexachloroethane	9.520	117	21879	45.460	ng	100
19) n-Nitroso-di-n-propyla...	9.215	70	57150	40.382	ng	100
20) 3+4-Methylphenols	9.174	107	78141	41.858	ng	100
22) Acetophenone	9.244	105	100170	39.137	ng	100
24) Nitrobenzene	9.638	77	83359	48.666	ng	100
25) Isophorone	10.155	82	159336	40.018	ng	100
26) 2-Nitrophenol	10.354	139	33335	49.485	ng	100
27) 2,4-Dimethylphenol	10.396	122	57340	37.249	ng	100
28) bis(2-Chloroethoxy)met...	10.631	93	78110	38.494	ng	100
29) 2,4-Dichlorophenol	10.889	162	61086	40.039	ng	100
30) 1,2,4-Trichlorobenzene	11.118	180	64805	35.657	ng	100
31) Naphthalene	11.312	128	183176	39.456	ng	100
32) Benzoic acid	10.513	122	44685	45.038	ng	100
33) 4-Chloroaniline	11.406	127	84057	39.630	ng	100
34) Hexachlorobutadiene	11.588	225	48856	37.616	ng	100
35) Caprolactam	12.170	113	21233	43.270	ng	100
36) 4-Chloro-3-methylphenol	12.493	107	69133	41.090	ng	100
37) 2-Methylnaphthalene	12.886	142	146387	39.935	ng	100
38) 1-Methylnaphthalene	13.104	142	135503	39.615	ng	100
40) 1,2,4,5-Tetrachloroben...	13.245	216	93291	37.784	ng	100
41) Hexachlorocyclopentadiene	13.221	237	54670	40.824	ng	100
43) 2,4,6-Trichlorophenol	13.474	196	60358	40.291	ng	100

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.544	196	70583	39.313	ng	100
46) 1,1'-Biphenyl	13.867	154	194841	38.726	ng	100
47) 2-Chloronaphthalene	13.920	162	143861	36.133	ng	100
48) 2-Nitroaniline	14.114	65	52818	54.825	ng	100
49) Acenaphthylene	14.760	152	236378	36.539	ng	100
50) Dimethylphthalate	14.467	163	214751	39.113	ng	100
51) 2,6-Dinitrotoluene	14.590	165	46237	47.955	ng	100
52) Acenaphthene	15.101	154	165023	41.354	ng	100
53) 3-Nitroaniline	14.925	138	45149	42.362	ng	100
54) 2,4-Dinitrophenol	15.130	184	27597	43.507	ng	100
55) Dibenzofuran	15.430	168	248066	36.796	ng	100
56) 4-Nitrophenol	15.213	139	33411	40.916	ng	100
57) 2,4-Dinitrotoluene	15.377	165	65438	48.869	ng	100
58) Fluorene	16.076	166	202908	38.066	ng	100
59) 2,3,4,6-Tetrachlorophenol	15.647	232	65756	41.788	ng	100
60) Diethylphthalate	15.818	149	212376	38.346	ng	100
61) 4-Chlorophenyl-phenyle...	16.053	204	118937	36.524	ng	100
62) 4-Nitroaniline	16.082	138	47349	42.628	ng	100
63) Azobenzene	16.346	77	210933	38.054	ng	100
65) 4,6-Dinitro-2-methylph...	16.141	198	41017	48.122	ng	100
66) n-Nitrosodiphenylamine	16.264	169	175668	36.584	ng	100
67) 4-Bromophenyl-phenylether	16.951	248	79847	37.679	ng	100
68) Hexachlorobenzene	17.075	284	80640	37.350	ng	100
69) Atrazine	17.198	200	74004	40.136	ng	100
70) Pentachlorophenol	17.416	266	59048	41.297	ng	100
71) Phenanthrene	17.815	178	333137	37.098	ng	100
72) Anthracene	17.909	178	344118	37.451	ng	100
73) Carbazole	18.167	167	301229	38.030	ng	100
74) Di-n-butylphthalate	18.696	149	350813	39.576	ng	100
75) Fluoranthene	19.801	202	423336	35.728	ng	100
77) Benzidine	19.965	184	171067	46.436	ng	100
78) Pyrene	20.165	202	422871	36.353	ng	100
80) Butylbenzylphthalate	21.022	149	148459	42.501	ng	100
81) Benzo(a)anthracene	22.074	228	426653	37.248	ng	100
82) 3,3'-Dichlorobenzidine	21.974	252	144490	38.880	ng	100
83) Chrysene	22.150	228	397387	37.111	ng	100
84) Bis(2-ethylhexyl)phtha...	21.933	149	217717	41.595	ng	100
85) Di-n-octyl phthalate	23.255	149	364354	41.881	ng	100
87) Indeno(1,2,3-cd)pyrene	29.752	276	501978	38.016	ng	100
88) Benzo(b)fluoranthene	24.512	252	425080	38.592	ng	100
89) Benzo(k)fluoranthene	24.582	252	432594	39.938	ng	100
90) Benzo(a)pyrene	25.481	252	421718	39.831	ng	100
91) Dibenzo(a,h)anthracene	29.822	278	411102	37.591	ng	100
92) Benzo(g,h,i)perylene	31.044	276	402808	37.695	ng	100

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

