

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG041023\
 Data File : BG057038.D
 Acq On : 10 Apr 2023 9:55
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040

Quant Time: Apr 11 04:10:53 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG033023.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Apr 08 04:45:47 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.405	152	14723	20.000	ng	0.00
21) Naphthalene-d8	11.248	136	62397	20.000	ng	# 0.00
39) Acenaphthene-d10	15.026	164	50185	20.000	ng	0.00
64) Phenanthrene-d10	17.763	188	125177	20.000	ng	0.00
76) Chrysene-d12	22.081	240	103400	20.000	ng	0.00
86) Perylene-d12	25.623	264	107530	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.914	112	73791	80.214	ng	0.00
7) Phenol-d6	7.536	99	113928	82.804	ng	0.00
23) Nitrobenzene-d5	9.586	82	116186	76.190	ng	0.00
42) 2,4,6-Tribromophenol	16.506	330	66679	83.170	ng	0.00
45) 2-Fluorobiphenyl	13.651	172	287216	77.606	ng	0.00
79) Terphenyl-d14	20.330	244	531514	86.983	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.717	88	14702	36.457	ng	98
3) Pyridine	4.140	79	50822	39.128	ng	97
4) n-Nitrosodimethylamine	4.046	42	22103	37.013	ng	99
6) Aniline	7.718	93	70617	40.148	ng	99
8) 2-Chlorophenol	7.965	128	38456	41.969	ng	94
9) Benzaldehyde	7.530	77	31566	36.578	ng	94
10) Phenol	7.565	94	55721	40.872	ng	97
11) bis(2-Chloroethyl)ether	7.806	93	39140	40.304	ng	97
12) 1,3-Dichlorobenzene	8.299	146	40725	40.223	ng	98
13) 1,4-Dichlorobenzene	8.440	146	41692	40.800	ng	97
14) 1,2-Dichlorobenzene	8.769	146	40428	40.951	ng	91
15) Benzyl Alcohol	8.640	79	45150	39.493	ng	96
16) 2,2'-oxybis(1-Chloropr...	8.922	45	46061	39.356	ng	99
17) 2-Methylphenol	8.840	107	40207	41.338	ng	89
18) Hexachloroethane	9.510	117	15362	41.131	ng	93
19) n-Nitroso-di-n-propyla...	9.210	70	40411	41.063	ng	96
20) 3+4-Methylphenols	9.169	107	56150	41.285	ng	98
22) Acetophenone	9.233	105	71394	38.452	ng	96
24) Nitrobenzene	9.627	77	60079	38.535	ng	97
25) Isophorone	10.150	82	111400	38.362	ng	# 94
26) 2-Nitrophenol	10.344	139	24666	40.988	ng	# 85
27) 2,4-Dimethylphenol	10.385	122	41296	38.866	ng	94
28) bis(2-Chloroethoxy)met...	10.626	93	56263	38.701	ng	99
29) 2,4-Dichlorophenol	10.884	162	43184	37.792	ng	100
30) 1,2,4-Trichlorobenzene	11.101	180	47538	39.154	ng	98
31) Naphthalene	11.301	128	132809	39.049	ng	99
32) Benzoic acid	10.514	122	27230	37.309	ng	92
33) 4-Chloroaniline	11.395	127	62426	40.640	ng	96
34) Hexachlorobutadiene	11.577	225	35248	38.648	ng	96
35) Caprolactam	12.165	113	15625	39.650	ng	91
36) 4-Chloro-3-methylphenol	12.488	107	49817	38.984	ng	95
37) 2-Methylnaphthalene	12.876	142	105553	39.104	ng	98
38) 1-Methylnaphthalene	13.093	142	99527	39.267	ng	99
40) 1,2,4,5-Tetrachloroben...	13.234	216	67278	38.659	ng	99
41) Hexachlorocyclopentadiene	13.210	237	37727	39.346	ng	97
43) 2,4,6-Trichlorophenol	13.463	196	44181	39.784	ng	95

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.539	196	51258	38.962	ng	98
46) 1,1'-Biphenyl	13.862	154	141754	38.586	ng	99
47) 2-Chloronaphthalene	13.909	162	108428	39.907	ng	99
48) 2-Nitroaniline	14.103	65	38081	39.197	ng	94
49) Acenaphthylene	14.749	152	176429	39.521	ng	99
50) Dimethylphthalate	14.462	163	158119	39.987	ng	# 99
51) 2,6-Dinitrotoluene	14.585	165	34968	40.930	ng	92
52) Acenaphthene	15.090	154	121674	39.520	ng	98
53) 3-Nitroaniline	14.920	138	35049	41.446	ng	92
54) 2,4-Dinitrophenol	15.120	184	20838	43.333	ng	92
55) Dibenzofuran	15.419	168	183167	39.638	ng	99
56) 4-Nitrophenol	15.208	139	25523	40.815	ng	# 90
57) 2,4-Dinitrotoluene	15.372	165	49418	41.012	ng	97
58) Fluorene	16.065	166	152169	39.795	ng	98
59) 2,3,4,6-Tetrachlorophenol	15.637	232	48017	40.086	ng	# 98
60) Diethylphthalate	15.807	149	161922	40.847	ng	100
61) 4-Chlorophenyl-phenyle...	16.042	204	89635	39.925	ng	99
62) 4-Nitroaniline	16.077	138	37952	43.316	ng	86
63) Azobenzene	16.336	77	152203	38.527	ng	97
65) 4,6-Dinitro-2-methylph...	16.130	198	32029	42.248	ng	97
66) n-Nitrosodiphenylamine	16.253	169	133373	39.691	ng	98
67) 4-Bromophenyl-phenylether	16.941	248	58443	37.734	ng	96
68) Hexachlorobenzene	17.070	284	63201	41.274	ng	95
69) Atrazine	17.187	200	53518	38.368	ng	98
70) Pentachlorophenol	17.411	266	41305	37.599	ng	96
71) Phenanthrene	17.804	178	248493	38.936	ng	99
72) Anthracene	17.898	178	255392	39.036	ng	100
73) Carbazole	18.157	167	216969	38.037	ng	95
74) Di-n-butylphthalate	18.679	149	260203	39.393	ng	99
75) Fluoranthene	19.790	202	303832	37.555	ng	99
77) Benzidine	19.954	184	99180	35.092	ng	98
78) Pyrene	20.154	202	302085	42.240	ng	96
80) Butylbenzylphthalate	21.006	149	107747	43.781	ng	95
81) Benzo(a)anthracene	22.057	228	289294	40.431	ng	99
82) 3,3'-Dichlorobenzidine	21.957	252	97992	40.557	ng	100
83) Chrysene	22.134	228	273503	40.824	ng	99
84) Bis(2-ethylhexyl)phtha...	21.910	149	148968	41.802	ng	98
85) Di-n-octyl phthalate	23.226	149	250808	41.764	ng	99
87) Indeno(1,2,3-cd)pyrene	29.717	276	320954	40.387	ng	# 97
88) Benzo(b)fluoranthene	24.489	252	281021	41.793	ng	99
89) Benzo(k)fluoranthene	24.566	252	282060	42.017	ng	99
90) Benzo(a)pyrene	25.459	252	267906	40.545	ng	100
91) Dibenzo(a,h)anthracene	29.776	278	273540	41.875	ng	98
92) Benzo(g,h,i)perylene	31.004	276	261801	40.930	ng	98

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

