

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG101022\
 Data File : BG055110.D
 Acq On : 10 Oct 2022 13:01
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040

Quant Time: Oct 11 04:02:40 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG100322.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Oct 08 03:04:59 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.332	152	28677	20.000	ng	0.00
21) Naphthalene-d8	11.158	136	122209	20.000	ng	-0.01
39) Acenaphthene-d10	14.936	164	93196	20.000	ng	0.00
64) Phenanthrene-d10	17.668	188	235025	20.000	ng	0.00
76) Chrysene-d12	21.963	240	229334	20.000	ng	0.00
86) Perylene-d12	25.412	264	248039	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.847	112	119701	92.713	ng	0.00
7) Phenol-d6	7.457	99	170248	88.008	ng	0.00
23) Nitrobenzene-d5	9.501	82	181528	88.758	ng	0.00
42) 2,4,6-Tribromophenol	16.411	330	98845	103.416	ng	0.00
45) 2-Fluorobiphenyl	13.567	172	449473	75.291	ng	0.00
79) Terphenyl-d14	20.248	244	882772	77.801	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.679	88	27875	38.158	ng	99
3) Pyridine	4.096	79	87606	39.881	ng	93
4) n-Nitrosodimethylamine	4.002	42	45607	36.304	ng	# 91
6) Aniline	7.639	93	109891	38.825	ng	96
8) 2-Chlorophenol	7.886	128	69759	46.973	ng	97
9) Benzaldehyde	7.451	77	56839	36.575	ng	94
10) Phenol	7.486	94	95762	43.562	ng	97
11) bis(2-Chloroethyl)ether	7.733	93	68495	38.016	ng	95
12) 1,3-Dichlorobenzene	8.220	146	81823	40.028	ng	96
13) 1,4-Dichlorobenzene	8.367	146	81099	39.203	ng	98
14) 1,2-Dichlorobenzene	8.691	146	78458	39.037	ng	97
15) Benzyl Alcohol	8.555	79	76066	43.778	ng	94
16) 2,2'-oxybis(1-Chloropr...	8.855	45	119922	35.139	ng	99
17) 2-Methylphenol	8.755	107	67698	43.710	ng	96
18) Hexachloroethane	9.431	117	27479	41.892	ng	90
19) n-Nitroso-di-n-propyla...	9.131	70	68707	36.666	ng	# 97
20) 3+4-Methylphenols	9.078	107	95476	44.265	ng	98
22) Acetophenone	9.155	105	120761	38.966	ng	# 97
24) Nitrobenzene	9.542	77	95568	43.035	ng	92
25) Isophorone	10.065	82	178844	41.144	ng	97
26) 2-Nitrophenol	10.259	139	37469	57.787	ng	# 84
27) 2,4-Dimethylphenol	10.300	122	68892	46.233	ng	96
28) bis(2-Chloroethoxy)met...	10.541	93	88985	40.008	ng	98
29) 2,4-Dichlorophenol	10.794	162	78129	43.305	ng	97
30) 1,2,4-Trichlorobenzene	11.017	180	85241	39.938	ng	96
31) Naphthalene	11.211	128	253643	39.325	ng	99
32) Benzoic acid	10.406	122	53727	54.019	ng	92
33) 4-Chloroaniline	11.305	127	109858	41.190	ng	98
34) Hexachlorobutadiene	11.493	225	53712	38.810	ng	97
35) Caprolactam	12.069	113	31365	54.108	ng	# 87
36) 4-Chloro-3-methylphenol	12.392	107	90679	43.840	ng	99
37) 2-Methylnaphthalene	12.792	142	187767	39.503	ng	98
38) 1-Methylnaphthalene	13.009	142	182467	39.466	ng	99
40) 1,2,4,5-Tetrachloroben...	13.150	216	101386	37.259	ng	99
41) Hexachlorocyclopentadiene	13.127	237	50942	43.475	ng	97
43) 2,4,6-Trichlorophenol	13.379	196	74905	45.472	ng	97

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.444	196	82700	44.474	ng	98
46) 1,1'-Biphenyl	13.779	154	243836	38.160	ng	98
47) 2-Chloronaphthalene	13.826	162	197998	38.454	ng	99
48) 2-Nitroaniline	14.014	65	67945	46.991	ng	97
49) Acenaphthylene	14.666	152	326456	38.288	ng	100
50) Dimethylphthalate	14.378	163	290117	45.865	ng	99
51) 2,6-Dinitrotoluene	14.495	165	57739	44.812	ng	90
52) Acenaphthene	15.001	154	210252	39.714	ng	99
53) 3-Nitroaniline	14.830	138	68994	45.897	ng	97
54) 2,4-Dinitrophenol	15.024	184	29514	50.198	ng	# 72
55) Dibenzofuran	15.330	168	329061	38.958	ng	99
56) 4-Nitrophenol	15.107	139	54766	49.221	ng	99
57) 2,4-Dinitrotoluene	15.277	165	84875	47.985	ng	# 87
58) Fluorene	15.976	166	268691	39.752	ng	97
59) 2,3,4,6-Tetrachlorophenol	15.547	232	81261	47.137	ng	99
60) Diethylphthalate	15.729	149	306160	46.908	ng	99
61) 4-Chlorophenyl-phenyle...	15.964	204	142439	38.767	ng	98
62) 4-Nitroaniline	15.982	138	72781	47.829	ng	99
63) Azobenzene	16.252	77	272700	37.097	ng	98
65) 4,6-Dinitro-2-methylph...	16.035	198	43724	49.912	ng	92
66) n-Nitrosodiphenylamine	16.170	169	243755	39.824	ng	96
67) 4-Bromophenyl-phenylether	16.852	248	98711	39.713	ng	97
68) Hexachlorobenzene	16.975	284	112053	38.992	ng	95
69) Atrazine	17.104	200	95658	47.969	ng	98
70) Pentachlorophenol	17.310	266	72968	47.756	ng	97
71) Phenanthrene	17.715	178	478787	38.782	ng	99
72) Anthracene	17.803	178	466783	38.875	ng	99
73) Carbazole	18.062	167	434796	41.255	ng	99
74) Di-n-butylphthalate	18.602	149	526667	47.871	ng	99
75) Fluoranthene	19.701	202	597086	38.768	ng	100
77) Benzidine	19.866	184	200467	50.741	ng	98
78) Pyrene	20.060	202	594370	38.977	ng	99
80) Butylbenzylphthalate	20.929	149	225730	47.350	ng	97
81) Benzo(a)anthracene	21.946	228	572166	39.274	ng	99
82) 3,3'-Dichlorobenzidine	21.846	252	195126	46.073	ng	96
83) Chrysene	22.016	228	535969	38.171	ng	99
84) Bis(2-ethylhexyl)phtha...	21.828	149	326727	45.956	ng	99
85) Di-n-octyl phthalate	23.121	149	556396	46.573	ng	96
87) Indeno(1,2,3-cd)pyrene	29.384	276	716359	40.020	ng	99
88) Benzo(b)fluoranthene	24.308	252	572753	39.190	ng	100
89) Benzo(k)fluoranthene	24.384	252	591074	40.146	ng	100
90) Benzo(a)pyrene	25.248	252	496513	39.639	ng	98
91) Dibenzo(a,h)anthracene	29.449	278	592705	40.761	ng	100
92) Benzo(g,h,i)perylene	30.624	276	581709	39.852	ng	99

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

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