

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG031324\
 Data File : BG060624.D
 Acq On : 13 Mar 2024 12:45
 Operator : MA/JU
 Sample : SSTDICC020
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDICC020

Quant Time: Mar 14 00:34:10 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	8.329	152	90858	20.000	ng	0.00
21) Naphthalene-d8	11.167	136	417680	20.000	ng	0.00
39) Acenaphthene-d10	14.951	164	277017	20.000	ng	0.00
64) Phenanthrene-d10	17.700	188	570637	20.000	ng	0.00
76) Chrysene-d12	22.019	240	516988	20.000	ng	0.00
86) Perylene-d12	25.573	264	577409	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.808	112	242367	41.933	ng	0.00
7) Phenol-d6	7.436	99	354655	42.300	ng	0.00
23) Nitrobenzene-d5	9.522	82	332054	41.375	ng	0.00
42) 2,4,6-Tribromophenol	16.431	330	129431	40.321	ng	0.00
45) 2-Fluorobiphenyl	13.576	172	884429	43.166	ng	0.00
79) Terphenyl-d14	20.274	244	1286764	44.989	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.652	88	52988	21.902	ng	96
3) Pyridine	4.069	79	147746	20.663	ng	97
4) n-Nitrosodimethylamine	3.999	42	73762	21.030	ng	99
6) Aniline	7.647	93	210666	20.563	ng	97
8) 2-Chlorophenol	7.871	128	131254	20.774	ng	98
9) Benzaldehyde	7.465	77	106078	22.602	ng	99
10) Phenol	7.459	94	176508	20.981	ng	98
11) bis(2-Chloroethyl)ether	7.741	93	138435	20.643	ng	99
12) 1,3-Dichlorobenzene	8.206	146	144454	21.202	ng	96
13) 1,4-Dichlorobenzene	8.364	146	144509	20.923	ng	97
14) 1,2-Dichlorobenzene	8.681	146	141643	20.650	ng	97
15) Benzyl Alcohol	8.564	79	135851	20.495	ng	97
16) 2,2'-oxybis(1-Chloropr...	8.834	45	276234m	21.219	ng	
17) 2-Methylphenol	8.740	107	136153	21.077	ng	98
18) Hexachloroethane	9.410	117	50189	21.273	ng	97
19) n-Nitroso-di-n-propyla...	9.128	70	122687	21.110	ng	98
20) 3+4-Methylphenols	9.069	107	180255	20.689	ng	98
22) Acetophenone	9.169	105	233626	20.863	ng	# 98
24) Nitrobenzene	9.563	77	166192	20.713	ng	95
25) Isophorone	10.074	82	331176	20.617	ng	98
26) 2-Nitrophenol	10.274	139	60556	20.000	ng	90
27) 2,4-Dimethylphenol	10.291	122	145979	20.533	ng	98
28) bis(2-Chloroethoxy)met...	10.556	93	201308	20.905	ng	98
29) 2,4-Dichlorophenol	10.791	162	129294	20.263	ng	91
30) 1,2,4-Trichlorobenzene	11.014	180	140341	20.715	ng	96
31) Naphthalene	11.220	128	486119	20.837	ng	99
32) Benzoic acid	10.362	122	70869m	19.223	ng	
33) 4-Chloroaniline	11.337	127	205521	20.891	ng	97
34) Hexachlorobutadiene	11.455	225	96539	20.891	ng	96
35) Caprolactam	12.113	113	44911	20.909	ng	89
36) 4-Chloro-3-methylphenol	12.395	107	162871	20.828	ng	96
37) 2-Methylnaphthalene	12.800	142	341309	21.182	ng	99
38) 1-Methylnaphthalene	13.018	142	318948	21.089	ng	98
40) 1,2,4,5-Tetrachloroben...	13.141	216	176978	20.803	ng	98
41) Hexachlorocyclopentadiene	13.100	237	84157	20.116	ng	96
43) 2,4,6-Trichlorophenol	13.382	196	112415	20.579	ng	99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.447	196	134217	20.770	ng	96
46) 1,1'-Biphenyl	13.793	154	441331	20.959	ng	98
47) 2-Chloronaphthalene	13.840	162	333133	20.963	ng	97
48) 2-Nitroaniline	14.052	65	104049	19.838	ng	95
49) Acenaphthylene	14.680	152	542031	20.938	ng	98
50) Dimethylphthalate	14.392	163	435245	20.950	ng	98
51) 2,6-Dinitrotoluene	14.533	165	79106	20.664	ng	97
52) Acenaphthene	15.015	154	327967	21.200	ng	97
53) 3-Nitroaniline	14.868	138	91804	20.828	ng	92
54) 2,4-Dinitrophenol	15.062	184	27445	19.650	ng	93
55) Dibenzofuran	15.344	168	527279	21.051	ng	99
56) 4-Nitrophenol	15.127	139	62733	19.073	ng	98
57) 2,4-Dinitrotoluene	15.309	165	97187	20.181	ng	97
58) Fluorene	15.996	166	424512	21.183	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.550	232	109191	20.208	ng	# 91
60) Diethylphthalate	15.738	149	448057	20.960	ng	98
61) 4-Chlorophenyl-phenyle...	15.979	204	214833	21.164	ng	97
62) 4-Nitroaniline	16.026	138	93717	20.769	ng	97
63) Azobenzene	16.273	77	469785	21.236	ng	98
65) 4,6-Dinitro-2-methylph...	16.055	198	43448	19.974	ng	96
66) n-Nitrosodiphenylamine	16.196	169	384564	21.534	ng	98
67) 4-Bromophenyl-phenylether	16.878	248	130430	21.196	ng	98
68) Hexachlorobenzene	16.966	284	148201	20.947	ng	98
69) Atrazine	17.125	200	128074	21.441	ng	97
70) Pentachlorophenol	17.318	266	87319	19.042	ng	98
71) Phenanthrene	17.741	178	662913	21.566	ng	99
72) Anthracene	17.835	178	660609	21.255	ng	99
73) Carbazole	18.106	167	580276	21.472	ng	98
74) Di-n-butylphthalate	18.617	149	709848	21.554	ng	99
75) Fluoranthene	19.733	202	732153	21.533	ng	99
77) Benzidine	19.921	184	225396	17.824	ng	99
78) Pyrene	20.098	202	769695	21.334	ng	98
80) Butylbenzylphthalate	20.961	149	280091	20.535	ng	96
81) Benzo(a)anthracene	21.995	228	741202	20.822	ng	100
82) 3,3'-Dichlorobenzidine	21.907	252	245366	20.451	ng	99
83) Chrysene	22.072	228	715983	20.715	ng	97
84) Bis(2-ethylhexyl)phtha...	21.843	149	411197	20.410	ng	99
85) Di-n-octyl phthalate	23.170	149	670011	20.195	ng	100
87) Indeno(1,2,3-cd)pyrene	29.692	276	808189	20.529	ng	# 95
88) Benzo(b)fluoranthene	24.422	252	663426	20.885	ng	99
89) Benzo(k)fluoranthene	24.498	252	696669	20.876	ng	98
90) Benzo(a)pyrene	25.403	252	647273	20.784	ng	99
91) Dibenzo(a,h)anthracene	29.780	278	683427	20.430	ng	98
92) Benzo(g,h,i)perylene	30.997	276	665311	20.552	ng	99

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

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