

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM043021\
 Data File : BM029710.D
 Acq On : 29 Apr 2021 15:15
 Operator : CG/JU
 Sample : SSTDICC050
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICC050

Quant Time: Apr 29 15:52:18 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM043021.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 29 15:25:42 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.598	152	69410	20.00	ng	0.00
21) Naphthalene-d8	10.375	136	252732	20.00	ng	0.00
39) Acenaphthene-d10	14.245	164	158057	20.00	ng	0.00
64) Phenanthrene-d10	16.992	188	376804	20.00	ng	0.00
76) Chrysene-d12	21.180	240	414560	20.00	ng	0.00
86) Perylene-d12	23.350	264	430123	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.193	112	336049	97.06	ng	0.00
7) Phenol-d6	6.781	99	459710	88.90	ng	0.00
23) Nitrobenzene-d5	8.751	82	439478	89.56	ng	0.00
42) 2,4,6-Tribromophenol	15.739	330	260792	119.31	ng	0.00
45) 2-Fluorobiphenyl	12.863	172	1223834	105.09	ng	0.00
79) Terphenyl-d14	19.627	244	2355266	107.71	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.122	88	61267	43.90	ng	95
3) Pyridine	3.528	79	126934	35.68	ng	96
4) n-Nitrosodimethylamine	3.440	42	63591	36.43	ng	85
6) Aniline	6.934	93	246328	43.33	ng	96
8) 2-Chlorophenol	7.163	128	204250	47.45	ng	93
9) Benzaldehyde	6.745	77	87549	29.78	ng	97
10) Phenol	6.810	94	218841	43.67	ng	95
11) bis(2-Chloroethyl)ether	7.034	93	171647	44.90	ng	96
12) 1,3-Dichlorobenzene	7.487	146	239114	48.20	ng	97
13) 1,4-Dichlorobenzene	7.634	146	250790	49.37	ng	97
14) 1,2-Dichlorobenzene	7.945	146	243200	48.03	ng	98
15) Benzyl Alcohol	7.845	79	161395	42.96	ng	95
16) 2,2'-oxybis(1-Chloropr...	8.122	45	182383	36.19	ng	96
17) 2-Methylphenol	8.045	107	156064	44.69	ng	96
18) Hexachloroethane	8.663	117	89258	48.66	ng	88
19) n-Nitroso-di-n-propyla...	8.410	70	147365	39.07	ng	90
20) 3+4-Methylphenols	8.375	107	209425	43.60	ng	96
22) Acetophenone	8.422	105	284514	46.96	ng	95
24) Nitrobenzene	8.792	77	218071	43.94	ng	94
25) Isophorone	9.322	82	395959	42.94	ng	96
26) 2-Nitrophenol	9.498	139	117424	54.76	ng	92
27) 2,4-Dimethylphenol	9.569	122	160056	48.32	ng	98
28) bis(2-Chloroethoxy)met...	9.804	93	215918	47.39	ng	99
29) 2,4-Dichlorophenol	10.033	162	227759	52.13	ng	96
30) 1,2,4-Trichlorobenzene	10.245	180	270391	53.73	ng	98
31) Naphthalene	10.428	128	636900	49.01	ng	99
32) Benzoic acid	9.739	122	124706	48.86	ng	93
33) 4-Chloroaniline	10.545	127	250238	48.83	ng	96
34) Hexachlorobutadiene	10.710	225	187334	58.47	ng	97
35) Caprolactam	11.339	113	56043	42.36	ng	86
36) 4-Chloro-3-methylphenol	11.680	107	194825	45.63	ng	96
37) 2-Methylnaphthalene	12.045	142	492036	47.30	ng	98
38) 1-Methylnaphthalene	12.269	142	466393	47.02	ng	100
40) 1,2,4,5-Tetrachloroben...	12.422	216	314409	66.34	ng	# 97
41) Hexachlorocyclopentadiene	12.398	237	164046	62.02	ng	98
43) 2,4,6-Trichlorophenol	12.669	196	203912	62.58	ng	98

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.739	196	204925	56.97	ng	96
46) 1,1'-Biphenyl	13.074	154	593719	51.39	ng	98
47) 2-Chloronaphthalene	13.116	162	481309	52.06	ng	98
48) 2-Nitroaniline	13.327	65	105028	45.18	ng	92
49) Acenaphthylene	13.969	152	723714	51.05	ng	100
50) Dimethylphthalate	13.716	163	601804	50.38	ng	99
51) 2,6-Dinitrotoluene	13.833	165	139716	53.13	ng	# 82
52) Acenaphthene	14.310	154	451323	50.82	ng	98
53) 3-Nitroaniline	14.163	138	113665	51.77	ng	95
54) 2,4-Dinitrophenol	14.374	184	84198	58.19	ng	# 81
55) Dibenzofuran	14.645	168	766058	51.09	ng	96
56) 4-Nitrophenol	14.480	139	94128	48.39	ng	92
57) 2,4-Dinitrotoluene	14.627	165	188386	53.05	ng	# 86
58) Fluorene	15.298	166	632275	50.22	ng	98
59) 2,3,4,6-Tetrachlorophenol	14.880	232	189617	59.24	ng	# 89
60) Diethylphthalate	15.086	149	597909	50.99	ng	97
61) 4-Chlorophenyl-phenyle...	15.298	204	352329	55.30	ng	95
62) 4-Nitroaniline	15.327	138	121157	52.75	ng	88
63) Azobenzene	15.592	77	498282	44.76	ng	90
65) 4,6-Dinitro-2-methylph...	15.392	198	130618	52.15	ng	86
66) n-Nitrosodiphenylamine	15.515	169	547459	47.00	ng	99
67) 4-Bromophenyl-phenylether	16.192	248	219184	55.71	ng	92
68) Hexachlorobenzene	16.304	284	246613	55.13	ng	# 91
69) Atrazine	16.468	200	219844	52.90	ng	96
70) Pentachlorophenol	16.645	266	156311	53.66	ng	98
71) Phenanthrene	17.033	178	1056134	48.68	ng	100
72) Anthracene	17.127	178	1063989	49.71	ng	99
73) Carbazole	17.398	167	968836	48.31	ng	99
74) Di-n-butylphthalate	17.968	149	1159894	53.27	ng	99
75) Fluoranthene	19.051	202	1312892	49.90	ng	99
77) Benzidine	19.245	184	401920	45.79	ng	99
78) Pyrene	19.415	202	1365123	51.18	ng	99
80) Butylbenzylphthalate	20.327	149	524132	52.49	ng	88
81) Benzo(a)anthracene	21.162	228	1356472	49.46	ng	99
82) 3,3'-Dichlorobenzidine	21.103	252	449913	51.35	ng	# 99
83) Chrysene	21.215	228	1363389	50.59	ng	99
84) Bis(2-ethylhexyl)phtha...	21.103	149	857845	53.61	ng	99
85) Di-n-octyl phthalate	21.962	149	1423398	53.01	ng	# 95
87) Indeno(1,2,3-cd)pyrene	25.544	276	1618538	48.67	ng	98
88) Benzo(b)fluoranthene	22.703	252	1427445	50.89	ng	98
89) Benzo(k)fluoranthene	22.750	252	1383610	49.22	ng	99
90) Benzo(a)pyrene	23.262	252	1309666	49.73	ng	99
91) Dibenzo(a,h)anthracene	25.550	278	1399148	48.79	ng	98
92) Benzo(g,h,i)perylene	26.209	276	1305344	49.23	ng	98

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

