

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA G\DATA\BG022118\
 Data File : BG032758.D
 Acq On : 21 Feb 2018 12:48
 Operator : SJ/JU
 Sample : SSTDICCC040
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDICCC040

Manual Integrations
 APPROVED

Sohil
 2/22/2018 3:13:48 PM

Quant Time: Feb 21 13:59:51 2018
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG022118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 21 13:32:56 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.94	152	62950	20.00	ng	0.00
21) Naphthalene-d8	11.82	136	286130	20.00	ng	0.00
38) Acenaphthene-d10	15.55	164	158630	20.00	ng	0.00
63) Phenanthrene-d10	18.28	188	348454	20.00	ng	0.00
75) Chrysene-d12	22.64	240	314294	20.00	ng	0.00
86) Perylene-d12	26.46	264	334374	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	6.34	112	323836	103.21	ng	0.00
7) Phenol-d6	8.03	99	446020	103.96	ng	0.00
23) Nitrobenzene-d5	10.13	82	299918	66.62	ng	0.00
41) 2,4,6-Tribromophenol	17.01	330	136467	55.87	ng	0.00
44) 2-Fluorobiphenyl	14.18	172	882710	69.59	ng	0.00
78) Terphenyl-d14	20.80	244	1104476	80.09	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.98	88	68069	51.862	ng	88
3) Pyridine	4.44	79	195867	58.180	ng	95
4) n-Nitrosodimethylamine	4.34	42	76980	48.663	ng	# 95
6) Aniline	8.22	93	297221	58.616	ng	97
8) 2-Chlorophenol	8.48	128	175234	48.569	ng	94
9) Benzaldehyde	8.03	77	123111	48.555	ng	95
10) Phenol	8.05	94	221986	53.201	ng	92
11) bis(2-Chloroethyl)ether	8.31	93	181065	54.128	ng	94
12) 1,3-Dichlorobenzene	8.82	146	202419	39.780	ng	97
13) 1,4-Dichlorobenzene	8.97	146	204428	40.664	ng	95
14) 1,2-Dichlorobenzene	9.31	146	193764	38.215	ng	96
15) Benzyl Alcohol	9.17	79	161160	50.117	ng	90
16) 2,2'-oxybis(1-Chloropropan	9.47	45	309570	85.988	ng	97
17) 2-Methylphenol	9.36	107	164431	55.645	ng	95
18) Hexachloroethane	10.06	117	69614	37.924	ng	# 86
19) n-Nitroso-di-n-propylamine	9.75	70	153374	53.177	ng	# 98
20) 3+4-Methylphenols	9.69	107	226162	50.639	ng	94
22) Acetophenone	9.78	105	289823	47.626	ng	# 93
24) Nitrobenzene	10.18	77	170337	36.136	ng	92
25) Isophorone	10.70	82	397927	46.359	ng	# 94
26) 2-Nitrophenol	10.90	139	56110	22.256	ng	91
27) 2,4-Dimethylphenol	10.93	122	170971	51.617	ng	90
28) bis(2-Chloroethoxy)methane	11.18	93	233184	53.142	ng	96
29) 2,4-Dichlorophenol	11.44	162	158597	32.714	ng	93
30) 1,2,4-Trichlorobenzene	11.67	180	185946	27.359	ng	97
31) Naphthalene	11.87	128	594753	42.808	ng	99
32) Benzoic acid	11.05	122	92913	38.489	ng	# 82
33) 4-Chloroaniline	11.96	127	266083	48.267	ng	95
34) Hexachlorobutadiene	12.14	225	104503	18.761	ng	97
35) Caprolactam	12.70	113	65298	47.431	ng	# 82
36) 4-Chloro-3-methylphenol	13.01	107	187672	43.037	ng	91
37) 2-Methylnaphthalene	13.42	142	430350	38.187	ng	94
39) 1,2,4,5-Tetrachlorobenzene	13.77	216	187716	25.055	ng	# 96
40) Hexachlorocyclopentadiene	13.75	237	97644	20.594	ng	93

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.99	196	119597	31.982	nq	99
43) 2,4,5-Trichlorophenol	14.06	196	139276	29.887	nq	97
45) 1,1'-Biphenyl	14.39	154	554520	42.484	nq	95
46) 2-Chloronaphthalene	14.45	162	416237	40.300	nq	98
47) 2-Nitroaniline	14.62	65	89400	39.792	nq	# 88
48) Acenaphthylene	15.28	152	683856	45.004	nq	98
49) Dimethylphthalate	14.98	163	539811	39.368	nq	99
50) 2,6-Dinitrotoluene	15.10	165	78615	27.839	nq	97
51) Acenaphthene	15.61	154	424070	40.866	nq	98
52) 3-Nitroaniline	15.43	138	99017	39.706	nq	95
53) 2,4-Dinitrophenol	15.62	184	21396	19.459	nq	# 1
54) Dibenzofuran	15.94	168	599264	38.814	nq	96
55) 4-Nitrophenol	15.69	139	75410	46.560	nq	# 81
56) 2,4-Dinitrotoluene	15.87	165	89825	23.085	nq	91
57) Fluorene	16.59	166	493762	38.888	nq	98
58) 2,3,4,6-Tetrachlorophenol	16.14	232	108781m	24.793	nq	
59) Diethylphthalate	16.31	149	569174	42.661	nq	97
60) 4-Chlorophenyl-phenylether	16.56	204	240909	28.737	nq	94
61) 4-Nitroaniline	16.58	138	110086	42.817	nq	89
62) Azobenzene	16.86	77	507867	59.248	nq	96
64) 4,6-Dinitro-2-methylphenol	16.63	198	35611	15.210	nq	95
65) n-Nitrosodiphenylamine	16.77	169	454759	44.499	nq	98
66) 4-Bromophenyl-phenylether	17.45	248	144847	29.046	nq	# 90
67) Hexachlorobenzene	17.58	284	156181	29.159	nq	# 94
68) Atrazine	17.69	200	150677	34.375	nq	95
69) Pentachlorophenol	17.91	266	100906	30.773	nq	98
70) Phenanthrene	18.32	178	774384	41.370	nq	99
71) Anthracene	18.41	178	780120	40.732	nq	99
72) Carbazole	18.66	167	770592	47.189	nq	98
73) Di-n-butylphthalate	19.17	149	960946	52.264	nq	98
74) Fluoranthene	20.27	202	855829	34.248	nq	94
76) Benzidine	20.42	184	417017	56.347	nq	98
77) Pyrene	20.63	202	879233	47.022	nq	97
79) Butylbenzylphthalate	21.50	149	411298	66.774	nq	91
80) Benzo(a)anthracene	22.62	228	805493	40.933	nq	99
81) 3,3'-Dichlorobenzidine	22.50	252	306781	41.744	nq	# 96
82) Chrysene	22.69	228	768585	39.610	nq	97
83) Bis(2-ethylhexyl)phthalate	22.47	149	613672	69.089	nq	# 95
84) Di-n-octyl phthalate	23.92	149	957972	68.742	nq	98
85) Indeno(1,2,3-cd)pyrene	30.92	276	923828	40.191	nq	97
87) Benzo(b)fluoranthene	25.24	252	808430	40.570	nq	# 96
88) Benzo(k)fluoranthene	25.32	252	784606	38.045	nq	# 96
89) Benzo(a)pyrene	26.28	252	759069	39.101	nq	# 96
90) Dibenzo(a,h)anthracene	31.02	278	773381	40.016	nq	# 94
91) Benzo(g,h,i)perylene	32.32	276	758631	40.606	nq	# 91

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

