

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG050118\
 Data File : BG034080.D
 Acq On : 1 May 2018 15:42
 Operator : SJ/JU
 Sample : SSTDICC060
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDICC060

Manual Integrations
 APPROVED

Sohil
 5/2/2018 4:50:36 PM

Quant Time: May 01 16:28:23 2018
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG050118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue May 01 15:30:49 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.08	152	66456	20.00	ng	0.00
21) Naphthalene-d8	10.90	136	279325	20.00	ng	0.00
38) Acenaphthene-d10	14.72	164	224287	20.00	ng	0.00
63) Phenanthrene-d10	17.48	188	573285	20.00	ng	0.00
75) Chrysene-d12	21.77	240	560160	20.00	ng	0.00
86) Perylene-d12	25.06	264	562759	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.64	112	491530	118.08	ng	0.00
7) Phenol-d6	7.22	99	764261	125.88	ng	0.00
23) Nitrobenzene-d5	9.24	82	897489	182.85	ng	0.00
41) 2,4,6-Tribromophenol	16.21	330	363432	138.89	ng	0.00
44) 2-Fluorobiphenyl	13.35	172	1736115	113.71	ng	0.00
78) Terphenyl-d14	20.10	244	2833920	113.66	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.56	88	98902	56.230	ng	83
3) Pyridine	3.95	79	318189	62.524	ng	86
4) n-Nitrosodimethylamine	3.86	42	179426	77.389	ng	# 79
6) Aniline	7.40	93	514455	63.390	ng	92
8) 2-Chlorophenol	7.64	128	242192	51.927	ng	85
9) Benzaldehyde	7.21	77	175055	52.875	ng	95
10) Phenol	7.25	94	393780	63.313	ng	# 76
11) bis(2-Chloroethyl)ether	7.50	93	300835	63.241	ng	98
12) 1,3-Dichlorobenzene	7.97	146	301143	56.087	ng	# 88
13) 1,4-Dichlorobenzene	8.12	146	306825	56.421	ng	95
14) 1,2-Dichlorobenzene	8.43	146	280976m	53.130	ng	
15) Benzyl Alcohol	8.31	79	423178	79.934	ng	# 83
16) 2,2'-oxybis(1-Chloropropan	8.61	45	407835	43.957	ng	82
17) 2-Methylphenol	8.51	107	264729m	57.437	ng	
18) Hexachloroethane	9.17	117	127993	64.409	ng	94
19) n-Nitroso-di-n-propylamine	8.89	70	337795	65.272	ng	98
20) 3+4-Methylphenols	8.85	107	375962	56.138	ng	# 76
22) Acetophenone	8.90	105	549543	72.285	ng	# 94
24) Nitrobenzene	9.29	77	478109	89.391	ng	# 84
25) Isophorone	9.81	82	865423	78.425	ng	95
26) 2-Nitrophenol	10.00	139	144096	68.185	ng	# 67
27) 2,4-Dimethylphenol	10.06	122	252978	63.999	ng	# 80
28) bis(2-Chloroethoxy)methane	10.30	93	423684	72.648	ng	98
29) 2,4-Dichlorophenol	10.53	162	296804	67.356	ng	97
30) 1,2,4-Trichlorobenzene	10.75	180	360864	73.117	ng	99
31) Naphthalene	10.94	128	875974	60.984	ng	97
32) Benzoic acid	10.21	122	221758	60.715	ng	# 80
33) 4-Chloroaniline	11.05	127	399722	59.205	ng	# 82
34) Hexachlorobutadiene	11.24	225	309575	104.852	ng	95
35) Caprolactam	11.84	113	107566m	57.416	ng	
36) 4-Chloro-3-methylphenol	12.17	107	385192	71.621	ng	93
37) 2-Methylnaphthalene	12.55	142	688382	62.595	ng	# 93
39) 1,2,4,5-Tetrachlorobenzene	12.92	216	504409	67.727	ng	98
40) Hexachlorocyclopentadiene	12.90	237	311040	75.955	ng	91

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.15	196	287337	63.278	nq	99
43) 2,4,5-Trichlorophenol	13.22	196	322902	61.854	nq	94
45) 1,1'-Biphenyl	13.56	154	1009979	55.628	nq	98
46) 2-Chloronaphthalene	13.60	162	771041	56.008	nq	95
47) 2-Nitroaniline	13.79	65	334677	77.861	nq #	85
48) Acenaphthylene	14.44	152	1206158	52.709	nq	99
49) Dimethylphthalate	14.18	163	1091368	55.317	nq	98
50) 2,6-Dinitrotoluene	14.29	165	242170	64.770	nq #	83
51) Acenaphthene	14.79	154	843880	58.704	nq	97
52) 3-Nitroaniline	14.62	138	219055	53.956	nq #	77
53) 2,4-Dinitrophenol	14.82	184	87246	65.249	nq #	76
54) Dibenzofuran	15.12	168	1180010	55.368	nq #	86
55) 4-Nitrophenol	14.91	139	165669	52.030	nq #	42
56) 2,4-Dinitrotoluene	15.08	165	330388	66.451	nq	90
57) Fluorene	15.77	166	1019333	55.271	nq	97
58) 2,3,4,6-Tetrachlorophenol	15.34	232	329638	68.964	nq	97
59) Diethylphthalate	15.55	149	1158159	55.536	nq	98
60) 4-Chlorophenyl-phenylether	15.77	204	616008	63.860	nq	96
61) 4-Nitroaniline	15.79	138	231382	53.677	nq #	39
62) Azobenzene	16.06	77	1237900	66.591	nq	93
64) 4,6-Dinitro-2-methylphenol	15.84	198	170356	67.277	nq	95
65) n-Nitrosodiphenylamine	15.98	169	899904	53.993	nq	99
66) 4-Bromophenyl-phenylether	16.67	248	413481	65.675	nq #	91
67) Hexachlorobenzene	16.78	284	427122	63.614	nq	96
68) Atrazine	16.93	200	307312	47.065	nq	97
69) Pentachlorophenol	17.12	266	300558	65.169	nq	99
70) Phenanthrene	17.52	178	1674719	54.454	nq	98
71) Anthracene	17.61	178	1688240	54.892	nq	100
72) Carbazole	17.88	167	1504475	50.705	nq #	96
73) Di-n-butylphthalate	18.45	149	1814093	49.146	nq #	97
74) Fluoranthene	19.53	202	2002997	53.841	nq	93
76) Benzidine	19.70	184	824036	54.477	nq	98
77) Pyrene	19.89	202	2032071m	55.308	nq	
79) Butylbenzylphthalate	20.79	149	822761	50.993	nq	89
80) Benzo(a)anthracene	21.75	228	2069215	58.581	nq	97
81) 3,3'-Dichlorobenzidine	21.67	252	783562	59.494	nq	96
82) Chrysene	21.82	228	1996279	59.184	nq	97
83) Bis(2-ethylhexyl)phthalate	21.68	149	1121862	49.234	nq	98
84) Di-n-octyl phthalate	22.94	149	1810938	50.071	nq	97
85) Indeno(1,2,3-cd)pyrene	28.85	276	2300157	59.907	nq #	72
87) Benzo(b)fluoranthene	24.02	252	2121170	61.785	nq #	96
88) Benzo(k)fluoranthene	24.09	252	2024828	59.375	nq #	95
89) Benzo(a)pyrene	24.91	252	2008066	61.073	nq #	96
90) Dibenzo(a,h)anthracene	28.93	278	1881086	59.126	nq #	94
91) Benzo(g,h,i)perylene	30.03	276	1872688	59.820	nq #	90

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

