

Data Path : Z:\HPCHEM1\BNA G\DATA\BG051618\
 Data File : BG034487.D
 Acq On : 16 May 2018 14:17
 Operator : SJ/JU
 Sample : SSTDICC2.5
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDICC2.5

Manual Integrations
 APPROVED

Sohil
 5/17/2018 5:04:08 PM

Quant Time: May 16 19:34:04 2018
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG051618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 16 19:55:14 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.04	152	35957	20.00	ng	0.00
21) Naphthalene-d8	10.86	136	151289	20.00	ng	0.00
38) Acenaphthene-d10	14.69	164	118315	20.00	ng	0.00
63) Phenanthrene-d10	17.45	188	339543	20.00	ng	0.00
75) Chrysene-d12	21.73	240	390258	20.00	ng	0.00
86) Perylene-d12	25.00	264	390203	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.62	112	10732	4.82	ng	0.00
7) Phenol-d6	7.21	99	17545	5.06	ng	0.00
23) Nitrobenzene-d5	9.21	82	19279	4.89	ng	0.00
41) 2,4,6-Tribromophenol	16.18	330	10074	6.15	ng	0.00
44) 2-Fluorobiphenyl	13.31	172	46676	5.75	ng	0.00
78) Terphenyl-d14	20.06	244	91855	5.16	ng	0.00

Target Compounds

						Qvalue
4) n-Nitrosodimethylamine	3.84	42	3357	2.100	ng	# 71
8) 2-Chlorophenol	7.61	128	5973	2.690	ng	# 81
10) Phenol	7.23	94	8603	2.462	ng	# 67
11) bis(2-Chloroethyl)ether	7.46	93	6551	2.381	ng	# 81
12) 1,3-Dichlorobenzene	7.93	146	6844m	2.436	ng	
13) 1,4-Dichlorobenzene	8.09	146	7139m	2.558	ng	
14) 1,2-Dichlorobenzene	8.40	146	7565m	2.854	ng	
15) Benzyl Alcohol	8.28	79	7631	2.065	ng	# 79
16) 2,2'-oxybis(1-Chloropropan	8.57	45	6523	1.723	ng	# 84
17) 2-Methylphenol	8.50	107	6999	3.044	ng	# 73
18) Hexachloroethane	9.14	117	3002	2.529	ng	# 95
19) n-Nitroso-di-n-propylamine	8.84	70	6738	2.165	ng	# 81
20) 3+4-Methylphenols	8.82	107	9057	2.659	ng	# 72
22) Acetophenone	8.86	105	12063	2.341	ng	# 87
24) Nitrobenzene	9.26	77	9171	2.125	ng	# 90
25) Isophorone	9.77	82	18687	2.323	ng	# 88
26) 2-Nitrophenol	9.97	139	3583	2.999	ng	# 94
27) 2,4-Dimethylphenol	10.03	122	6167	2.662	ng	# 97
28) bis(2-Chloroethoxy)methane	10.26	93	11120	2.814	ng	# 91
29) 2,4-Dichlorophenol	10.51	162	5619	2.029	ng	# 90
30) 1,2,4-Trichlorobenzene	10.73	180	8712	2.613	ng	# 91
31) Naphthalene	10.91	128	20654	2.597	ng	# 88
33) 4-Chloroaniline	11.01	127	8089	2.256	ng	# 87
34) Hexachlorobutadiene	11.20	225	6822	2.418	ng	# 94
35) Caprolactam	11.76	113	2352	2.392	ng	# 59
36) 4-Chloro-3-methylphenol	12.15	107	8028	2.236	ng	# 71
37) 2-Methylnaphthalene	12.52	142	17047	2.644	ng	# 91
39) 1,2,4,5-Tetrachlorobenzene	12.89	216	10809	2.361	ng	# 89
40) Hexachlorocyclopentadiene	12.86	237	7181	2.647	ng	# 82
42) 2,4,6-Trichlorophenol	13.13	196	5902	2.219	ng	# 91
43) 2,4,5-Trichlorophenol	13.21	196	6795	2.311	ng	# 69
45) 1,1'-Biphenyl	13.53	154	22962	2.484	ng	# 95
46) 2-Chloronaphthalene	13.56	162	18429	2.634	ng	# 88
47) 2-Nitroaniline	13.76	65	7194	2.571	ng	# 69

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
48) Acenaphthylene	14.41	152	29738	2.662	nq	# 89
49) Dimethylphthalate	14.13	163	25783	2.513	nq	# 98
50) 2,6-Dinitrotoluene	14.26	165	5038	2.557	nq	93
51) Acenaphthene	14.76	154	18579	2.447	nq	# 88
52) 3-Nitroaniline	14.59	138	4492	2.262	nq	# 70
54) Dibenzofuran	15.09	168	32676	2.968	nq	# 62
56) 2,4-Dinitrotoluene	15.05	165	7270	2.717	nq	# 79
57) Fluorene	15.74	166	23595	2.459	nq	# 81
58) 2,3,4,6-Tetrachlorophenol	15.31	232	7983	2.624	nq	# 97
59) Diethylphthalate	15.50	149	26875	2.475	nq	# 85
60) 4-Chlorophenyl-phenylether	15.73	204	15221	2.667	nq	# 82
61) 4-Nitroaniline	15.76	138	6145	2.917	nq	# 69
62) Azobenzene	16.03	77	29471	2.513	nq	94
65) n-Nitrosodiphenylamine	15.94	169	23334	2.520	nq	97
66) 4-Bromophenyl-phenylether	16.63	248	8926	2.117	nq	# 79
67) Hexachlorobenzene	16.75	284	10669	2.493	nq	# 77
68) Atrazine	16.90	200	11446	3.019	nq	99
69) Pentachlorophenol	17.10	266	5347	1.793	nq	# 68
70) Phenanthrene	17.48	178	47141	2.702	nq	98
71) Anthracene	17.58	178	48493m	2.742	nq	
72) Carbazole	17.85	167	37793	2.335	nq	# 89
73) Di-n-butylphthalate	18.41	149	48270	2.477	nq	# 92
74) Fluoranthene	19.50	202	57366	2.602	nq	93
77) Pyrene	19.86	202	59444	2.431	nq	96
79) Butylbenzylphthalate	20.75	149	21794	2.283	nq	# 82
80) Benzo(a)anthracene	21.71	228	66627	2.664	nq	95
81) 3,3'-Dichlorobenzidine	21.62	252	23150	2.448	nq	# 93
82) Chrysene	21.78	228	58117	2.428	nq	96
83) Bis(2-ethylhexyl)phthalate	21.62	149	30922	2.354	nq	# 97
84) Di-n-octyl phthalate	22.85	149	49689	2.368	nq	# 98
85) Indeno(1,2,3-cd)pyrene	28.72	276	64965m	2.466	nq	
87) Benzo(b)fluoranthene	23.96	252	63156	2.562	nq	# 89
88) Benzo(k)fluoranthene	24.03	252	58281	2.474	nq	# 94
89) Benzo(a)pyrene	24.84	252	56799	2.463	nq	# 93
90) Dibenzo(a,h)anthracene	28.79	278	56902	2.618	nq	# 93
91) Benzo(a,h,i)perylene	29.88	276	58706	2.740	nq	# 90

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

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