

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG050118\
 Data File : BG034111.D
 Acq On : 2 May 2018 11:50
 Operator : SJ/JU
 Sample : J2607-02
 Misc : 8PPM MDL
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 NJD001-74

Manual Integrations
 APPROVED

Sohil
 5/2/2018 4:51:05 PM

Quant Time: May 02 14:16:24 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 19:08:03 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.07	152	66051	20.00	ng	0.00
21) Naphthalene-d8	10.89	136	255348	20.00	ng	0.00
38) Acenaphthene-d10	14.72	164	196000	20.00	ng	0.00
63) Phenanthrene-d10	17.47	188	554819	20.00	ng	0.00
75) Chrysene-d12	21.76	240	670831	20.00	ng	0.00
86) Perylene-d12	25.05	264	656265	20.00	ng	-0.02

System Monitoring Compounds

5) 2-Fluorophenol	5.64	112	586529	143.47	ng	0.00
7) Phenol-d6	7.22	99	797651	125.33	ng	0.00
23) Nitrobenzene-d5	9.24	82	687254	103.27	ng	0.00
41) 2,4,6-Tribromophenol	16.21	330	442168	162.92	ng	0.00
44) 2-Fluorobiphenyl	13.34	172	1398632	103.98	ng	0.00
78) Terphenyl-d14	20.09	244	2870292	93.73	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.56	88	128	0.075	ng	# 1
3) Pyridine	3.95	79	112	0.021	ng	# 15
4) n-Nitrosodimethylamine	3.86	42	207	0.070	ng	# 1
6) Aniline	7.39	93	7617	0.880	ng	# 84
8) 2-Chlorophenol	7.64	128	176	0.043	ng	# 1
10) Phenol	7.24	94	351	0.055	ng	# 1
11) bis(2-Chloroethyl)ether	7.39	93	7617	1.507	ng	# 26
12) 1,3-Dichlorobenzene	7.93	146	113	0.022	ng	# 72
13) 1,4-Dichlorobenzene	7.93	146	113	0.022	ng	# 89
14) 1,2-Dichlorobenzene	7.93	146	113	0.023	ng	# 92
15) Benzyl Alcohol	8.39	79	13430	1.978	ng	# 43
16) 2,2'-oxybis(1-Chloropropan	8.62	45	356	0.051	ng	# 75
17) 2-Methylphenol	8.48	107	85	0.020	ng	# 13
18) Hexachloroethane	9.13	117	130	0.060	ng	# 1
20) 3+4-Methylphenols	8.83	107	89	0.014	ng	# 12
22) Acetophenone	8.90	105	657	0.076	ng	# 68
24) Nitrobenzene	9.23	77	3083	0.423	ng	# 25
25) Isophorone	9.81	82	92	0.007	ng	# 69
26) 2-Nitrophenol	10.02	139	179	0.089	ng	# 29
27) 2,4-Dimethylphenol	10.16	122	81	0.021	ng	# 1
28) bis(2-Chloroethoxy)methane	10.27	93	115	0.017	ng	# 51
29) 2,4-Dichlorophenol	10.49	162	262	0.056	ng	# 23
30) 1,2,4-Trichlorobenzene	10.76	180	74	0.013	ng	# 12
31) Naphthalene	10.94	128	795	0.059	ng	# 68
33) 4-Chloroaniline	11.04	127	2368	0.391	ng	# 76
34) Hexachlorobutadiene	11.38	225	68	0.014	ng	# 20
35) Caprolactam	11.77	113	71	0.043	ng	# 20
36) 4-Chloro-3-methylphenol	12.14	107	177	0.029	ng	# 21
37) 2-Methylnaphthalene	12.55	142	162	0.015	ng	# 17
39) 1,2,4,5-Tetrachlorobenzene	12.91	216	124	0.016	ng	# 72
40) Hexachlorocyclopentadiene	12.48	237	98	0.022	ng	# 87
42) 2,4,6-Trichlorophenol	13.07	196	111	0.025	ng	# 11
43) 2,4,5-Trichlorophenol	13.21	196	213	0.044	ng	# 17
45) 1,1'-Biphenyl	13.55	154	3416	0.223	ng	# 69

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
46) 2-Chloronaphthalene	13.65	162	67	0.006	nq	# 1
47) 2-Nitroaniline	13.78	65	187	0.040	nq	# 9
48) Acenaphthylene	14.52	152	62	0.003	nq	# 59
49) Dimethylphthalate	14.17	163	174	0.010	nq	# 48
51) Acenaphthene	14.77	154	532	0.042	nq	# 5
52) 3-Nitroaniline	14.64	138	56	0.017	nq	# 1
53) 2,4-Dinitrophenol	14.90	184	104	0.081	nq	# 1
54) Dibenzofuran	15.11	168	316	0.017	nq	# 46
55) 4-Nitrophenol	14.90	139	54	0.022	nq	# 1
57) Fluorene	15.77	166	503	0.032	nq	# 7
58) 2,3,4,6-Tetrachlorophenol	15.35	232	127	0.025	nq	# 33
59) Diethylphthalate	15.54	149	588	0.033	nq	# 1
60) 4-Chlorophenyl-phenylether	15.71	204	97	0.010	nq	# 30
61) 4-Nitroaniline	15.78	138	88	0.025	nq	# 12
62) Azobenzene	16.04	77	213	0.011	nq	# 48
64) 4,6-Dinitro-2-methylphenol	15.78	198	206	0.079	nq	# 68
65) n-Nitrosodiphenylamine	16.20	169	19329	1.278	nq	# 42
67) Hexachlorobenzene	16.92	284	55	0.008	nq	# 41
68) Atrazine	16.96	200	75	0.925	nq	# 61
69) Pentachlorophenol	17.12	266	55	0.011	nq	# 19
70) Phenanthrene	17.51	178	1060	0.037	nq	# 60
71) Anthracene	17.60	178	141	0.005	nq	# 59
72) Carbazole	17.88	167	73	0.003	nq	# 1
73) Di-n-butylphthalate	18.44	149	888	0.028	nq	# 21
74) Fluoranthene	19.49	202	126	0.003	nq	# 65
76) Benzidine	19.70	184	1102115m	63.780	nq	
77) Pyrene	20.09	202	15223	0.362	nq	# 53
79) Butylbenzylphthalate	20.78	149	162	0.010	nq	# 20
80) Benzo(a)anthracene	21.76	228	1682	0.039	nq	# 20
81) 3,3'-Dichlorobenzidine	21.65	252	1069	0.066	nq	# 61
82) Chrysene	21.82	228	55	0.001	nq	# 49
83) Bis(2-ethylhexyl)phthalate	21.69	149	262	0.012	nq	# 61
84) Di-n-octyl phthalate	22.91	149	143	0.004	nq	# 1
85) Indeno(1,2,3-cd)pyrene	28.79	276	71	0.002	nq	# 1
87) Benzo(b)fluoranthene	24.00	252	77	0.002	nq	# 59
88) Benzo(k)fluoranthene	24.11	252	243	0.006	nq	# 1
89) Benzo(a)pyrene	24.89	252	83	0.002	nq	# 9
90) Dibenzo(a,h)anthracene	28.92	278	60	0.002	nq	# 1
91) Benzo(g,h,i)perylene	29.99	276	63	0.002	nq	# 56

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

