

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG091715\
 Data File : BG018721.D
 Acq On : 17 Sep 2015 00:43
 Operator : TP
 Sample : G3651-07MSD
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 CPS033031MSD

Manual Integrations
 APPROVED

MMDadoda
 9/17/2015 3:48:52 PM

Quant Time: Sep 17 06:01:10 2015
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG091115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 11 10:44:16 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.00	152	53385	20.00	ng	0.00
21) Naphthalene-d8	10.80	136	228785	20.00	ng	0.00
38) Acenaphthene-d10	14.63	164	157092	20.00	ng	0.00
63) Phenanthrene-d10	17.37	188	413175	20.00	ng	0.00
75) Chrysene-d12	21.63	240	422725	20.00	ng	0.00
86) Perylene-d12	24.81	264	437303	20.00	ng	-0.01

System Monitoring Compounds

5) 2-Fluorophenol	5.58	112	112873	36.52	ng	0.00
7) Phenol-d6	7.16	99	69850	15.76	ng	0.00
23) Nitrobenzene-d5	9.16	82	255117	62.40	ng	0.00
41) 2,4,6-Tribromophenol	16.12	330	30022	14.19	ng	0.00
44) 2-Fluorobiphenyl	13.25	172	700816	61.93	ng	0.00
78) Terphenyl-d14	19.98	244	1008564	62.65	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.47	88	14969	11.59	ng	# 100
3) Pyridine	3.87	79	44685	11.16	ng	94
4) n-Nitrosodimethylamine	3.78	42	17154	12.30	ng	# 76
6) Aniline	7.32	93	144652	23.69	ng	99
8) 2-Chlorophenol	7.57	128	82337	23.07	ng	94
9) Benzaldehyde	7.14	77	46339	19.42	ng	92
10) Phenol	7.19	94	30822	6.58	ng	97
11) bis(2-Chloroethyl)ether	7.42	93	110602	30.49	ng	98
12) 1,3-Dichlorobenzene	7.89	146	107488	25.89	ng	95
13) 1,4-Dichlorobenzene	8.04	146	109528	26.31	ng	95
14) 1,2-Dichlorobenzene	8.36	146	108617	27.32	ng	98
15) Benzyl Alcohol	8.23	79	64968	20.34	ng	99
16) 2,2'-oxybis(1-Chloropropan	8.54	45	127155	30.87	ng	97
17) 2-Methylphenol	8.44	107	40808	12.55	ng	97
18) Hexachloroethane	9.08	117	35849	24.08	ng	96
19) n-Nitroso-di-n-propylamine	8.80	70	89442	28.66	ng	96
20) 3+4-Methylphenols	8.77	107	47315	10.36	ng	97
22) Acetophenone	8.82	105	178410	30.78	ng	# 98
24) Nitrobenzene	9.20	77	130246	30.00	ng	98
25) Isophorone	9.72	82	261321	29.72	ng	100
26) 2-Nitrophenol	9.93	139	18282	9.03	ng	93
27) 2,4-Dimethylphenol	9.97	122	58865	16.00	ng	96
28) bis(2-Chloroethoxy)methane	10.21	93	161733	32.03	ng	98
29) 2,4-Dichlorophenol	10.46	162	86146	23.18	ng	98
30) 1,2,4-Trichlorobenzene	10.67	180	118618	28.80	ng	95
31) Naphthalene	10.85	128	360242	29.40	ng	99
33) 4-Chloroaniline	10.96	127	190030	34.28	ng	97
34) Hexachlorobutadiene	11.14	225	65465	26.96	ng	99
35) Caprolactam	11.70	113	12856	8.04	ng	93
36) 4-Chloro-3-methylphenol	12.09	107	58723	14.16	ng	90
37) 2-Methylnaphthalene	12.46	142	278904	31.10	ng	99
39) 1,2,4,5-Tetrachlorobenzene	12.83	216	150637	30.23	ng	99
42) 2,4,6-Trichlorophenol	13.08	196	25191	7.30	ng	99
43) 2,4,5-Trichlorophenol	13.15	196	72345m	19.11	ng	

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
45) 1,1'-Biphenyl	13.46	154	387093	30.99	ng	100
46) 2-Chloronaphthalene	13.51	162	298077	30.97	ng	99
47) 2-Nitroaniline	13.71	65	93078	32.14	ng	88
48) Acenaphthylene	14.35	152	531106	31.22	ng	98
49) Dimethylphthalate	14.08	163	584162	44.91	ng	100
50) 2,6-Dinitrotoluene	14.20	165	95251	33.70	ng	98
51) Acenaphthene	14.69	154	304727	30.79	ng	97
52) 3-Nitroaniline	14.53	138	108544	32.14	ng	90
54) Dibenzofuran	15.02	168	522935	33.99	ng	99
56) 2,4-Dinitrotoluene	14.98	165	81508	20.34	ng	# 93
57) Fluorene	15.68	166	427342	32.27	ng	97
59) Diethylphthalate	15.44	149	451225	33.26	ng	98
60) 4-Chlorophenyl-phenylether	15.66	204	210242	33.04	ng	95
61) 4-Nitroaniline	15.69	138	103102	28.31	ng	96
62) Azobenzene	15.96	77	386907	32.27	ng	97
65) n-Nitrosodiphenylamine	15.88	169	374152	31.27	ng	96
66) 4-Bromophenyl-phenylether	16.56	248	141342	33.64	ng	97
67) Hexachlorobenzene	16.68	284	154839	33.91	ng	95
68) Atrazine	16.82	200	135828	28.55	ng	98
70) Phenanthrene	17.41	178	701586	32.28	ng	99
71) Anthracene	17.50	178	716011	32.54	ng	98
72) Carbazole	17.77	167	660993	31.03	ng	100
73) Di-n-butylphthalate	18.33	149	791000	31.49	ng	99
74) Fluoranthene	19.42	202	838825	31.17	ng	97
76) Benzidine	19.59	184	244405	19.25	ng	97
77) Pyrene	19.78	202	859150	33.07	ng	98
79) Butylbenzylphthalate	20.66	149	335883	31.80	ng	97
80) Benzo(a)anthracene	21.61	228	808435	33.22	ng	100
81) 3,3'-Dichlorobenzidine	21.52	252	205719	22.25	ng	99
82) Chrysene	21.67	228	733963	32.03	ng	99
83) Bis(2-ethylhexyl)phthalate	21.52	149	512134	33.61	ng	98
84) Di-n-octyl phthalate	22.71	149	858277	33.31	ng	100
85) Indeno(1,2,3-cd)pyrene	28.44	276	949067	32.61	ng	# 100
87) Benzo(b)fluoranthene	23.81	252	837780	33.04	ng	98
88) Benzo(k)fluoranthene	23.87	252	801980	31.57	ng	98
89) Benzo(a)pyrene	24.66	252	800810	33.19	ng	97
90) Dibenzo(a,h)anthracene	28.50	278	770741	32.44	ng	98
91) Benzo(g,h,i)perylene	29.58	276	773522	31.97	ng	97

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

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