

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF091815\
 Data File : BF081708.D
 Acq On : 18 Sep 2015 21:46
 Operator : UM/IZ
 Sample : G3704-13MSD
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-12-9-15-15MSD

Manual Integrations
 APPROVED

MMDadoda
 9/21/2015 2:22:32 PM

Quant Time: Sep 21 10:50:39 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF090115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 15 14:03:59 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.16	152	41276	20.00	ng	-0.06
21) Naphthalene-d8	11.06	136	167974	20.00	ng	-0.06
38) Acenaphthene-d10	15.20	164	78127	20.00	ng	-0.06
63) Phenanthrene-d10	18.70	188	162732	20.00	ng	-0.06
75) Chrysene-d12	23.35	240	140763	20.00	ng	-0.03
86) Perylene-d12	24.78	264	91862	20.00	ng	-0.03

System Monitoring Compounds

5) 2-Fluorophenol	5.32	112	158984	59.76	ng	-0.03
7) Phenol-d6	7.59	99	130205	36.97	ng	-0.06
23) Nitrobenzene-d5	9.48	82	368891	105.96	ng	-0.06
41) 2,4,6-Tribromophenol	17.11	330	115580	166.15	ng	-0.06
44) 2-Fluorobiphenyl	13.72	172	651627	106.88	ng	-0.06
78) Terphenyl-d14	22.07	244	590606	97.67	ng	-0.04

Target Compounds

						Qvalue
2) 1,4-Dioxane	1.57	88	35107	29.38	ng	# 81
3) Pyridine	2.12	79	29575	8.98	ng	# 75
4) n-Nitrosodimethylamine	2.05	42	36523	19.95	ng	# 59
6) Aniline	7.48	93	79359	15.96	ng	# 84
8) 2-Chlorophenol	7.72	128	110675	37.75	ng	# 79
9) Benzaldehyde	7.19	77	56891	25.78	ng	# 73
10) Phenol	7.61	94	52206m	12.78	ng	
11) bis(2-Chloroethyl)ether	7.70	93	139939	47.50	ng	# 80
12) 1,3-Dichlorobenzene	8.01	146	129007	41.19	ng	# 87
13) 1,4-Dichlorobenzene	8.21	146	132552	41.56	ng	# 90
14) 1,2-Dichlorobenzene	8.53	146	129942	45.25	ng	# 91
15) Benzyl Alcohol	8.61	79	80020	27.70	ng	# 68
16) 2,2'-oxybis(1-Chloropropan	8.93	45	298359	52.83	ng	95
17) 2-Methylphenol	8.96	107	71521	30.58	ng	# 79
18) Hexachloroethane	9.26	117	54343	43.80	ng	88
19) n-Nitroso-di-n-propylamine	9.25	70	119815	49.99	ng	# 89
20) 3+4-Methylphenols	9.34	107	80691	25.59	ng	# 80
22) Acetophenone	9.16	105	213267	46.48	ng	# 73
24) Nitrobenzene	9.51	77	177671	49.14	ng	# 61
25) Isophorone	10.10	82	309912	47.86	ng	# 84
26) 2-Nitrophenol	10.24	139	68720	43.61	ng	# 21
27) 2,4-Dimethylphenol	10.52	122	107240	39.40	ng	# 74
28) bis(2-Chloroethoxy)methane	10.71	93	182402	48.76	ng	# 91
29) 2,4-Dichlorophenol	10.86	162	102834	43.57	ng	93
30) 1,2,4-Trichlorobenzene	10.97	180	116333	45.37	ng	# 97
31) Naphthalene	11.12	128	409136	45.67	ng	99
32) Benzoic acid	10.93	122	15458m	10.08	ng	
33) 4-Chloroaniline	11.36	127	62375	17.07	ng	# 82
34) Hexachlorobutadiene	11.46	225	73298	46.97	ng	99
35) Caprolactam	12.29	113	2767	3.82	ng	# 28
36) 4-Chloro-3-methylphenol	12.66	107	116662	44.16	ng	# 76
37) 2-Methylnaphthalene	12.77	142	285718	49.01	ng	# 91
39) 1,2,4,5-Tetrachlorobenzene	13.18	216	122871	49.73	ng	96
40) Hexachlorocyclopentadiene	13.16	237	116100	95.75	ng	97

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.52	196	78115	45.81	ng	98
43) 2,4,5-Trichlorophenol	13.64	196	81544	46.87	ng #	83
45) 1,1'-Biphenyl	13.91	154	348908	48.54	ng	96
46) 2-Chloronaphthalene	13.90	162	258155	49.73	ng #	87
47) 2-Nitroaniline	14.25	65	110026	52.01	ng #	59
48) Acenaphthylene	14.86	152	443007	48.11	ng	97
49) Dimethylphthalate	14.80	163	337564	55.61	ng #	97
50) 2,6-Dinitrotoluene	14.89	165	63690	46.23	ng #	21
51) Acenaphthene	15.28	154	253424	48.38	ng	90
52) 3-Nitroaniline	15.25	138	36098	23.02	ng #	47
53) 2,4-Dinitrophenol	15.55	184	61174	86.40	ng #	24
54) Dibenzofuran	15.71	168	398079	52.04	ng #	85
55) 4-Nitrophenol	15.90	139	19183m	19.47	ng	
56) 2,4-Dinitrotoluene	15.85	165	91925	52.40	ng #	67
57) Fluorene	16.52	166	312183	53.78	ng	96
58) 2,3,4,6-Tetrachlorophenol	16.07	232	70329	52.95	ng	93
59) Diethylphthalate	16.49	149	347855	54.48	ng	95
60) 4-Chlorophenyl-phenylether	16.61	204	148743	54.98	ng #	86
61) 4-Nitroaniline	16.72	138	61504	38.81	ng #	13
62) Azobenzene	16.97	77	394049	55.51	ng	81
64) 4,6-Dinitro-2-methylphenol	16.81	198	43332	47.61	ng #	74
65) n-Nitrosodiphenylamine	16.93	169	266797	45.06	ng	92
66) 4-Bromophenyl-phenylether	17.75	248	83735	50.89	ng #	91
67) Hexachlorobenzene	17.80	284	84866	49.33	ng #	82
68) Atrazine	18.35	200	95471	55.71	ng #	78
69) Pentachlorophenol	18.35	266	65593	78.50	ng	98
70) Phenanthrene	18.77	178	446931	49.40	ng	97
71) Anthracene	18.88	178	452901	48.73	ng	98
72) Carbazole	19.36	167	418843	48.02	ng	96
73) Di-n-butylphthalate	20.41	149	560667	54.44	ng #	98
74) Fluoranthene	21.39	202	488465	49.88	ng	89
76) Benzidine	21.71	184	86104	14.25	ng #	97
77) Pyrene	21.74	202	505642	48.49	ng	99
79) Butylbenzylphthalate	22.77	149	229205	50.54	ng #	77
80) Benzo(a)anthracene	23.34	228	407841	45.50	ng	96
81) 3,3'-Dichlorobenzidine	23.35	252	82401	27.25	ng #	93
82) Chrysene	23.39	228	392185	48.81	ng	95
83) Bis(2-ethylhexyl)phthalate	23.47	149	333041	55.65	ng #	93
84) Di-n-octyl phthalate	24.12	149	510480	51.81	ng	96
85) Indeno(1,2,3-cd)pyrene	25.82	276	266621	45.22	ng #	84
87) Benzo(b)fluoranthene	24.45	252	336731m	51.34	ng	
88) Benzo(k)fluoranthene	24.47	252	272484m	50.07	ng	
89) Benzo(a)pyrene	24.73	252	289323	52.06	ng #	88
90) Dibenzo(a,h)anthracene	25.83	278	224932	52.38	ng #	79
91) Benzo(g,h,i)perylene	26.12	276	216193	52.75	ng #	76

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

