

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA G\DATA\BG091217\
 Data File : BG028681.D
 Acq On : 12 Sep 2017 22:29
 Operator : SJ/JU
 Sample : I5123-04MSD
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 EP-13MSD

Manual Integrations
 APPROVED

Sohil
 9/13/2017 6:00:13 PM

Quant Time: Sep 13 05:05:49 2017
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG091217.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 12 16:57:11 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.33	152	28290	20.00	ng	0.00
21) Naphthalene-d8	11.15	136	118702	20.00	ng	0.00
38) Acenaphthene-d10	14.94	164	91470	20.00	ng	0.00
63) Phenanthrene-d10	17.69	188	251031	20.00	ng	0.00
75) Chrysene-d12	22.03	240	284634	20.00	ng	0.00
86) Perylene-d12	25.53	264	273796	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.90	112	208298	121.49	ng	0.00
7) Phenol-d6	7.49	99	318251	123.29	ng	0.00
23) Nitrobenzene-d5	9.50	82	208306	83.95	ng	0.00
41) 2,4,6-Tribromophenol	16.44	330	187606	124.01	ng	0.00
44) 2-Fluorobiphenyl	13.56	172	520370	75.86	ng	0.00
78) Terphenyl-d14	20.28	244	976051	73.13	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.84	88	22441	32.322	ng	97
3) Pyridine	4.23	79	59595	30.090	ng	95
4) n-Nitrosodimethylamine	4.15	42	36260	40.247	ng	# 80
6) Aniline	7.66	93	87391	29.089	ng	96
8) 2-Chlorophenol	7.89	128	77892	40.754	ng	91
9) Benzaldehyde	7.48	77	10667	6.661	ng	# 86
10) Phenol	7.51	94	128760	47.264	ng	89
11) bis(2-Chloroethyl)ether	7.74	93	75253	38.475	ng	96
12) 1,3-Dichlorobenzene	8.22	146	81148	39.429	ng	92
13) 1,4-Dichlorobenzene	8.36	146	82524	38.995	ng	93
14) 1,2-Dichlorobenzene	8.69	146	81002	38.743	ng	97
15) Benzyl Alcohol	8.57	79	87942	43.641	ng	90
16) 2,2'-oxybis(1-Chloropropan	8.85	45	136153	38.302	ng	98
17) 2-Methylphenol	8.77	107	75896	42.633	ng	93
18) Hexachloroethane	9.42	117	30508	39.333	ng	97
19) n-Nitroso-di-n-propylamine	9.12	70	72599	39.674	ng	89
20) 3+4-Methylphenols	9.10	107	105427	42.340	ng	97
22) Acetophenone	9.15	105	129999	37.456	ng	# 90
24) Nitrobenzene	9.54	77	107789	39.229	ng	95
25) Isophorone	10.06	82	201728	40.179	ng	99
26) 2-Nitrophenol	10.26	139	45245	38.690	ng	# 88
27) 2,4-Dimethylphenol	10.30	122	90848	42.500	ng	98
28) bis(2-Chloroethoxy)methane	10.53	93	108442	39.773	ng	96
29) 2,4-Dichlorophenol	10.80	162	92013	43.263	ng	93
30) 1,2,4-Trichlorobenzene	11.01	180	93815	38.666	ng	94
31) Naphthalene	11.20	128	248765	40.126	ng	98
32) Benzoic acid	10.41	122	4620	8.743	ng	# 50
33) 4-Chloroaniline	11.32	127	62037	23.887	ng	92
34) Hexachlorobutadiene	11.47	225	65741	36.787	ng	97
35) Caprolactam	12.08	113	35671	46.771	ng	# 73
36) 4-Chloro-3-methylphenol	12.42	107	117532	42.463	ng	98
37) 2-Methylnaphthalene	12.78	142	200517	43.321	ng	97
39) 1,2,4,5-Tetrachlorobenzene	13.15	216	127381	33.869	ng	98
40) Hexachlorocyclopentadiene	13.12	237	119794	81.802	ng	96

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.39	196	89443	37.470	ng	97
43) 2,4,5-Trichlorophenol	13.47	196	102881	42.892	ng	91
45) 1,1'-Biphenyl	13.78	154	275450	36.371	ng	99
46) 2-Chloronaphthalene	13.83	162	212397	38.451	ng	99
47) 2-Nitroaniline	14.03	65	87826	42.781	ng	93
48) Acenaphthylene	14.67	152	355941	40.333	ng	97
49) Dimethylphthalate	14.39	163	380886	49.658	ng	96
50) 2,6-Dinitrotoluene	14.52	165	73286	42.940	ng	# 81
51) Acenaphthene	15.01	154	215353	40.171	ng	95
52) 3-Nitroaniline	14.85	138	48037	31.828	ng	93
53) 2,4-Dinitrophenol	15.06	184	14371	22.383	ng	96
54) Dibenzofuran	15.34	168	376526	44.042	ng	94
55) 4-Nitrophenol	15.16	139	86350	78.089	ng	# 81
56) 2,4-Dinitrotoluene	15.30	165	108788	45.332	ng	# 88
57) Fluorene	15.99	166	319473	41.857	ng	97
58) 2,3,4,6-Tetrachlorophenol	15.57	232	97174	45.968	ng	# 94
59) Diethylphthalate	15.74	149	345291	43.806	ng	97
60) 4-Chlorophenyl-phenylether	15.97	204	177672	40.881	ng	93
61) 4-Nitroaniline	16.02	138	66791	41.772	ng	# 86
62) Azobenzene	16.27	77	306069	46.169	ng	90
64) 4,6-Dinitro-2-methylphenol	16.07	198	33974	20.974	ng	95
65) n-Nitrosodiphenylamine	16.19	169	276462	38.418	ng	97
66) 4-Bromophenyl-phenylether	16.87	248	128992	39.935	ng	94
67) Hexachlorobenzene	17.00	284	134081	36.472	ng	# 88
68) Atrazine	17.13	200	102592	33.203	ng	98
69) Pentachlorophenol	17.35	266	133761	80.529	ng	94
70) Phenanthrene	17.74	178	540775	42.125	ng	95
71) Anthracene	17.83	178	549791	42.397	ng	96
72) Carbazole	18.10	167	471695	40.388	ng	# 94
73) Di-n-butylphthalate	18.62	149	576599	40.264	ng	# 93
74) Fluoranthene	19.74	202	656035	41.854	ng	93
76) Benzidine	19.91	184	16243	2.201	ng	# 92
77) Pyrene	20.10	202	676830	41.225	ng	96
79) Butylbenzylphthalate	20.96	149	261549	39.446	ng	97
80) Benzo(a)anthracene	22.01	228	708217	41.336	ng	93
81) 3,3'-Dichlorobenzidine	21.91	252	122502	17.635	ng	# 96
82) Chrysene	22.08	228	667886	41.085	ng	93
83) Bis(2-ethylhexyl)phthalate	21.86	149	368842	39.128	ng	100
84) Di-n-octyl phthalate	23.16	149	647981	39.344	ng	# 89
85) Indeno(1,2,3-cd)pyrene	29.57	276	851790	40.781	ng	# 99
87) Benzo(b)fluoranthene	24.41	252	718799	43.819	ng	99
88) Benzo(k)fluoranthene	24.49	252	716584m	43.733	ng	
89) Benzo(a)pyrene	25.36	252	641433	41.196	ng	96
90) Dibenzo(a,h)anthracene	29.62	278	713257	43.939	ng	98
91) Benzo(g,h,i)perylene	30.83	276	684190	43.196	ng	99

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

