

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA G\DATA\BG090717\
 Data File : BG028599.D
 Acq On : 7 Sep 2017 18:34
 Operator : SJ/JU
 Sample : SSTDICV040
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 ICVBG090717

Quant Time: Sep 07 19:22:30 2017
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG090717.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Sep 07 18:39:50 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.34	152	40418	20.00	ng	0.00
21) Naphthalene-d8	11.16	136	163741	20.00	ng	0.00
38) Acenaphthene-d10	14.95	164	110798	20.00	ng	0.00
63) Phenanthrene-d10	17.70	188	273342	20.00	ng	0.00
75) Chrysene-d12	22.03	240	317938	20.00	ng	0.00
86) Perylene-d12	25.54	264	305156	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.92	112	184490	76.85	ng	0.00
7) Phenol-d6	7.50	99	272501	78.26	ng	0.00
23) Nitrobenzene-d5	9.51	82	283175	79.41	ng	0.00
41) 2,4,6-Tribromophenol	16.44	330	137887	81.88	ng	0.00
44) 2-Fluorobiphenyl	13.57	172	675076	79.71	ng	0.00
78) Terphenyl-d14	20.28	244	1202219	81.47	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.86	88	38603	39.314	ng	95
3) Pyridine	4.25	79	116489	38.205	ng	93
4) n-Nitrosodimethylamine	4.16	42	54434	37.462	ng	# 74
6) Aniline	7.67	93	173416	38.901	ng	95
8) 2-Chlorophenol	7.91	128	104198	38.453	ng	98
9) Benzaldehyde	7.48	77	78731	38.318	ng	96
10) Phenol	7.52	94	147644	39.147	ng	94
11) bis(2-Chloroethyl)ether	7.75	93	105242	37.952	ng	98
12) 1,3-Dichlorobenzene	8.23	146	119680	37.715	ng	90
13) 1,4-Dichlorobenzene	8.38	146	122168	39.046	ng	97
14) 1,2-Dichlorobenzene	8.69	146	117625	38.256	ng	96
15) Benzyl Alcohol	8.57	79	105291	38.148	ng	94
16) 2,2'-oxybis(1-Chloropropan	8.85	45	213837	37.401	ng	97
17) 2-Methylphenol	8.78	107	92835	38.381	ng	95
18) Hexachloroethane	9.42	117	44930	36.552	ng	94
19) n-Nitroso-di-n-propylamine	9.14	70	106608	38.722	ng	91
20) 3+4-Methylphenols	9.10	107	131893	39.383	ng	96
22) Acetophenone	9.16	105	181659	39.935	ng	# 91
24) Nitrobenzene	9.55	77	151816	40.300	ng	94
25) Isophorone	10.07	82	261003	39.750	ng	99
26) 2-Nitrophenol	10.26	139	62778	40.183	ng	93
27) 2,4-Dimethylphenol	10.31	122	111028	39.349	ng	98
28) bis(2-Chloroethoxy)methane	10.54	93	155358	39.647	ng	98
29) 2,4-Dichlorophenol	10.80	162	112100	40.466	ng	94
30) 1,2,4-Trichlorobenzene	11.02	180	135158	39.651	ng	98
31) Naphthalene	11.21	128	349744	39.350	ng	100
32) Benzoic acid	10.46	122	75481	42.458	ng	88
33) 4-Chloroaniline	11.32	127	149343	39.859	ng	96
34) Hexachlorobutadiene	11.48	225	97975	39.779	ng	95
35) Caprolactam	12.09	113	42195	40.240	ng	94
36) 4-Chloro-3-methylphenol	12.42	107	141675	40.167	ng	95
37) 2-Methylnaphthalene	12.79	142	261274	39.868	ng	96
39) 1,2,4,5-Tetrachlorobenzene	13.15	216	173590	40.335	ng	# 96
40) Hexachlorocyclopentadiene	13.12	237	62789	38.616	ng	97

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42) 2,4,6-Trichlorophenol	13.39	196	110542	40.210	ng	93
43) 2,4,5-Trichlorophenol	13.47	196	110830	40.343	ng	94
45) 1,1'-Biphenyl	13.78	154	369145	39.681	ng	97
46) 2-Chloronaphthalene	13.84	162	286315	40.521	ng	99
47) 2-Nitroaniline	14.04	65	113439	39.377	ng	93
48) Acenaphthylene	14.68	152	429737	39.658	ng	97
49) Dimethylphthalate	14.39	163	359046	39.330	ng	98
50) 2,6-Dinitrotoluene	14.52	165	80909	40.455	ng	# 80
51) Acenaphthene	15.02	154	268426	40.449	ng	93
52) 3-Nitroaniline	14.86	138	80498	39.851	ng	99
53) 2,4-Dinitrophenol	15.07	184	45819	39.862	ng	96
54) Dibenzofuran	15.35	168	414856	40.444	ng	92
55) 4-Nitrophenol	15.17	139	58587	40.703	ng	# 80
56) 2,4-Dinitrotoluene	15.31	165	117631	40.810	ng	# 96
57) Fluorene	16.00	166	377365	40.246	ng	99
58) 2,3,4,6-Tetrachlorophenol	15.57	232	98485	40.141	ng	# 95
59) Diethylphthalate	15.74	149	382799	40.020	ng	96
60) 4-Chlorophenyl-phenylether	15.97	204	212803	40.184	ng	92
61) 4-Nitroaniline	16.02	138	88478	40.460	ng	# 84
62) Azobenzene	16.27	77	334456	39.148	ng	94
64) 4,6-Dinitro-2-methylphenol	16.08	198	79947	41.183	ng	91
65) n-Nitrosodiphenylamine	16.19	169	323787	40.113	ng	99
66) 4-Bromophenyl-phenylether	16.87	248	143653	40.473	ng	95
67) Hexachlorobenzene	17.01	284	158103	40.476	ng	# 89
68) Atrazine	17.13	200	114714	39.399	ng	97
69) Pentachlorophenol	17.36	266	73508	40.434	ng	94
70) Phenanthrene	17.74	178	591041	40.661	ng	95
71) Anthracene	17.84	178	597250	40.404	ng	98
72) Carbazole	18.11	167	527805	39.911	ng	94
73) Di-n-butylphthalate	18.62	149	630127	39.238	ng	# 95
74) Fluoranthene	19.74	202	745627	40.179	ng	93
76) Benzidine	19.91	184	249655	38.651	ng	99
77) Pyrene	20.10	202	764139	40.521	ng	95
79) Butylbenzylphthalate	20.97	149	293376	40.219	ng	91
80) Benzo(a)anthracene	22.01	228	762824	40.072	ng	94
81) 3,3'-Dichlorobenzidine	21.91	252	294672	40.287	ng	99
82) Chrysene	22.08	228	710711	39.589	ng	95
83) Bis(2-ethylhexyl)phthalate	21.86	149	411292	39.134	ng	97
84) Di-n-octyl phthalate	23.17	149	713731	39.621	ng	# 91
85) Indeno(1,2,3-cd)pyrene	29.59	276	856816	40.287	ng	# 95
87) Benzo(b)fluoranthene	24.42	252	757560	39.634	ng	97
88) Benzo(k)fluoranthene	24.49	252	744288	40.327	ng	94
89) Benzo(a)pyrene	25.38	252	718624	40.083	ng	95
90) Dibenzo(a,h)anthracene	29.63	278	750477	40.342	ng	99
91) Benzo(g,h,i)perylene	30.85	276	732584	40.152	ng	98

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

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