

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG092018\
 Data File : BG036851.D
 Acq On : 20 Sep 2018 20:04
 Operator : JU/SJ
 Sample : SSTDICV040
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 ICVBG092018

Manual Integrations
 APPROVED

Sohil
 9/21/2018 10:49:08 AM

Quant Time: Sep 21 10:39:34 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG092018.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Sep 20 20:07:11 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.05	152	38649	20.00	ng	0.00
21) Naphthalene-d8	10.86	136	196221	20.00	ng	0.00
38) Acenaphthene-d10	14.68	164	136401	20.00	ng	0.00
63) Phenanthrene-d10	17.42	188	348265	20.00	ng	0.00
75) Chrysene-d12	21.69	240	337323	20.00	ng	0.00
86) Perylene-d12	24.90	264	347709	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.62	112	168010	83.37	ng	0.00
7) Phenol-d6	7.21	99	267823	85.90	ng	0.00
23) Nitrobenzene-d5	9.21	82	285556	77.93	ng	0.00
41) 2,4,6-Tribromophenol	16.16	330	132819	81.39	ng	0.00
44) 2-Fluorobiphenyl	13.30	172	712600	77.81	ng	0.00
78) Terphenyl-d14	20.02	244	975950	76.08	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.53	88	39709	43.061	ng	95
3) Pyridine	3.92	79	114152	42.871	ng	96
4) n-Nitrosodimethylamine	3.84	42	48402	40.899	ng	99
6) Aniline	7.37	93	167715	41.454	ng	96
8) 2-Chlorophenol	7.61	128	94761	40.496	ng	94
9) Benzaldehyde	7.19	77	83965	40.753	ng	99
10) Phenol	7.24	94	137198	42.635	ng	97
11) bis(2-Chloroethyl)ether	7.47	93	115088	38.664	ng	97
12) 1,3-Dichlorobenzene	7.94	146	114867	40.040	ng	97
13) 1,4-Dichlorobenzene	8.08	146	119742	40.787	ng	96
14) 1,2-Dichlorobenzene	8.40	146	113185	39.784	ng	99
15) Benzyl Alcohol	8.28	79	107115	42.338	ng	98
16) 2,2'-oxybis(1-Chloropropan	8.57	45	228267	39.944	ng	96
17) 2-Methylphenol	8.49	107	91391	40.772	ng	98
18) Hexachloroethane	9.13	117	43640	40.393	ng	92
19) n-Nitroso-di-n-propylamine	8.85	70	107466	41.431	ng	95
20) 3+4-Methylphenols	8.82	107	132563	42.511	ng	97
22) Acetophenone	8.87	105	179162	38.854	ng	97
24) Nitrobenzene	9.25	77	151811	38.797	ng	94
25) Isophorone	9.78	82	270678	40.428	ng	97
26) 2-Nitrophenol	9.96	139	64035	41.056	ng	95
27) 2,4-Dimethylphenol	10.02	122	96599	39.760	ng	97
28) bis(2-Chloroethoxy)methane	10.26	93	160191	38.775	ng	98
29) 2,4-Dichlorophenol	10.50	162	116771	41.461	ng	96
30) 1,2,4-Trichlorobenzene	10.72	180	136450	38.714	ng	98
31) Naphthalene	10.90	128	360917	38.657	ng	98
32) Benzoic acid	10.16	122	67409m	37.642	ng	
33) 4-Chloroaniline	11.01	127	165907	39.083	ng	98
34) Hexachlorobutadiene	11.19	225	94646	38.830	ng	98
35) Caprolactam	11.79	113	48025	41.430	ng	98
36) 4-Chloro-3-methylphenol	12.13	107	141822	42.202	ng	98
37) 2-Methylnaphthalene	12.50	142	281602	39.602	ng	99
39) 1,2,4,5-Tetrachlorobenzene	12.87	216	186342	39.169	ng	99
40) Hexachlorocyclopentadiene	12.85	237	95786	39.148	ng	97

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.11	196	109058	41.281	ng	97
43) 2,4,5-Trichlorophenol	13.18	196	119555	42.441	ng	98
45) 1,1'-Biphenyl	13.51	154	406402	38.904	ng	98
46) 2-Chloronaphthalene	13.55	162	318347	38.751	ng	99
47) 2-Nitroaniline	13.76	65	112684	39.657	ng	97
48) Acenaphthylene	14.40	152	503593	39.119	ng	99
49) Dimethylphthalate	14.13	163	427918	38.975	ng	100
50) 2,6-Dinitrotoluene	14.25	165	96899	40.450	ng	94
51) Acenaphthene	14.74	154	305348	39.246	ng	98
52) 3-Nitroaniline	14.58	138	101555	40.231	ng	99
53) 2,4-Dinitrophenol	14.79	184	38986	39.056	ng	98
54) Dibenzofuran	15.07	168	514070	39.063	ng	98
55) 4-Nitrophenol	14.89	139	65540	40.642	ng	97
56) 2,4-Dinitrotoluene	15.04	165	135075	40.410	ng	89
57) Fluorene	15.72	166	380991	38.733	ng	98
58) 2,3,4,6-Tetrachlorophenol	15.30	232	113553	39.736	ng	97
59) Diethylphthalate	15.49	149	425744	38.446	ng	99
60) 4-Chlorophenyl-phenylether	15.71	204	244715	39.285	ng	98
61) 4-Nitroaniline	15.75	138	104201	39.244	ng	96
62) Azobenzene	16.00	77	410521	38.582	ng	98
64) 4,6-Dinitro-2-methylphenol	15.80	198	74949	41.163	ng	98
65) n-Nitrosodiphenylamine	15.93	169	375677	39.074	ng	99
66) 4-Bromophenyl-phenylether	16.61	248	157493	39.758	ng	98
67) Hexachlorobenzene	16.73	284	160089	39.239	ng	97
68) Atrazine	16.87	200	153096	39.326	ng	99
69) Pentachlorophenol	17.07	266	104871	40.826	ng	97
70) Phenanthrene	17.46	178	655177	39.039	ng	99
71) Anthracene	17.56	178	649903	39.163	ng	100
72) Carbazole	17.83	167	627528	38.784	ng	98
73) Di-n-butylphthalate	18.38	149	704279	39.195	ng	99
74) Fluoranthene	19.47	202	777984	38.511	ng	100
76) Benzidine	19.65	184	364065	35.703	ng	99
77) Pyrene	19.83	202	793332	39.192	ng	100
79) Butylbenzylphthalate	20.71	149	308518	38.238	ng	95
80) Benzo(a)anthracene	21.67	228	745982	37.884	ng	99
81) 3,3'-Dichlorobenzidine	21.59	252	290611	38.773	ng	99
82) Chrysene	21.73	228	713360	37.495	ng	99
83) Bis(2-ethylhexyl)phthalate	21.57	149	422711	37.503	ng	99
84) Di-n-octyl phthalate	22.78	149	710961	38.310	ng	98
85) Indeno(1,2,3-cd)pyrene	28.55	276	801502	39.231	ng	# 94
87) Benzo(b)fluoranthene	23.88	252	714884	38.579	ng	98
88) Benzo(k)fluoranthene	23.95	252	716293	38.530	ng	98
89) Benzo(a)pyrene	24.75	252	691665	38.609	ng	98
90) Dibenzo(a,h)anthracene	28.62	278	653141	38.901	ng	98
91) Benzo(g,h,i)perylene	29.69	276	657904	39.044	ng	97

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

