

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG090519\
 Data File : BG042727.D
 Acq On : 5 Sep 2019 19:24
 Operator : HP/JU
 Sample : K4668-03MSD
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SU-04-090319MSD

Manual Integrations
 APPROVED

mohammad
 9/10/2019 9:51:51 AM

Quant Time: Sep 06 07:16:42 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG082219.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 22 14:29:15 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.00	152	42001	20.00	ng	0.00
21) Naphthalene-d8	10.82	136	151153	20.00	ng	0.00
39) Acenaphthene-d10	14.66	164	114416	20.00	ng	0.00
64) Phenanthrene-d10	17.41	188	281809	20.00	ng	0.00
76) Chrysene-d12	21.69	240	319763	20.00	ng	0.00
87) Perylene-d12	24.93	264	388333	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.56	112	203094	97.93	ng	0.00
7) Phenol-d6	7.17	99	240909	86.00	ng	0.00
23) Nitrobenzene-d5	9.17	82	163718	74.85	ng	0.00
42) 2,4,6-Tribromophenol	16.15	330	195391	120.15	ng	0.00
45) 2-Fluorobiphenyl	13.27	172	537199	79.41	ng	0.00
79) Terphenyl-d14	20.02	244	1122730	75.91	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.43	88	24776	28.628	ng	# 51
3) Pyridine	3.84	79	43173	19.454	ng	93
4) n-Nitrosodimethylamine	3.74	42	24908	31.425	ng	89
6) Aniline	7.32	93	46392	12.417	ng	97
8) 2-Chlorophenol	7.56	128	84332	33.554	ng	97
9) Benzaldehyde	7.13	77	41818	29.818	ng	97
10) Phenol	7.19	94	80571	28.289	ng	92
11) bis(2-Chloroethyl)ether	7.41	93	66002	29.507	ng	96
12) 1,3-Dichlorobenzene	7.88	146	96923	30.314	ng	99
13) 1,4-Dichlorobenzene	8.03	146	99521	30.730	ng	97
14) 1,2-Dichlorobenzene	8.36	146	94074	30.332	ng	97
15) Benzyl Alcohol	8.24	79	61651	30.591	ng	99
16) 2,2'-oxybis(1-Chloropropan	8.53	45	80951	26.701	ng	99
17) 2-Methylphenol	8.45	107	64382	30.690	ng	95
18) Hexachloroethane	9.09	117	32757	29.545	ng	94
19) n-Nitroso-di-n-propylamine	8.81	70	51986	28.645	ng	# 94
20) 3+4-Methylphenols	8.78	107	85391	29.416	ng	97
22) Acetophenone	8.83	105	119710	37.029	ng	# 97
24) Nitrobenzene	9.21	77	76802	34.674	ng	98
25) Isophorone	9.74	82	143423	33.225	ng	98
26) 2-Nitrophenol	9.93	139	51697	37.244	ng	96
27) 2,4-Dimethylphenol	9.99	122	78221	40.910	ng	96
28) bis(2-Chloroethoxy)methane	10.22	93	89444	32.875	ng	97
29) 2,4-Dichlorophenol	10.48	162	95833	38.308	ng	98
30) 1,2,4-Trichlorobenzene	10.68	180	111178	36.207	ng	98
31) Naphthalene	10.87	128	253189	36.079	ng	98
32) Benzoic acid	10.15	122	20893	20.663	ng	95
33) 4-Chloroaniline	10.99	127	14806	4.570	ng	97
34) Hexachlorobutadiene	11.16	225	76117	36.885	ng	98
35) Caprolactam	11.76	113	20056m	24.027	ng	
36) 4-Chloro-3-methylphenol	12.12	107	83633	35.217	ng	95
37) 2-Methylnaphthalene	12.48	142	203230	36.067	ng	99
38) 1-Methylnaphthalene	12.70	142	189585	35.421	ng	99
40) 1,2,4,5-Tetrachlorobenzene	12.86	216	141691	38.563	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.83	237	146316	75.809	nq	98
43) 2,4,6-Trichlorophenol	13.09	196	88108	35.476	nq	97
44) 2,4,5-Trichlorophenol	13.17	196	95683	37.177	nq	95
46) 1,1'-Biphenyl	13.49	154	269695	36.465	nq	96
47) 2-Chloronaphthalene	13.53	162	225423	35.348	nq	99
48) 2-Nitroaniline	13.74	65	48008	31.218	nq	98
49) Acenaphthylene	14.38	152	348212	36.293	nq	99
50) Dimethylphthalate	14.11	163	304802	36.921	nq	100
51) 2,6-Dinitrotoluene	14.23	165	68605	35.852	nq	95
52) Acenaphthene	14.72	154	206891	35.512	nq	99
53) 3-Nitroaniline	14.56	138	29743	15.542	nq	100
54) 2,4-Dinitrophenol	14.78	184	84841	66.886	nq	97
55) Dibenzofuran	15.06	168	359726	37.303	nq	99
56) 4-Nitrophenol	14.89	139	85294	62.722	nq	96
57) 2,4-Dinitrotoluene	15.02	165	100467	36.664	nq	99
58) Fluorene	15.71	166	289170	36.683	nq	98
59) 2,3,4,6-Tetrachlorophenol	15.29	232	97968m	38.491	nq	
60) Diethylphthalate	15.47	149	282415	34.995	nq	99
61) 4-Chlorophenyl-phenylether	15.69	204	172962	36.398	nq	97
62) 4-Nitroaniline	15.73	138	69694	34.027	nq	92
63) Azobenzene	15.99	77	182454	32.994	nq	93
65) 4,6-Dinitro-2-methylphenol	15.80	198	68027	38.502	nq	96
66) n-Nitrosodiphenylamine	15.91	169	268889	34.696	nq	99
67) 4-Bromophenyl-phenylether	16.59	248	124953	34.956	nq	98
68) Hexachlorobenzene	16.72	284	132791	34.173	nq	97
69) Atrazine	16.86	200	107764	39.321	nq	97
70) Pentachlorophenol	17.07	266	149613	74.102	nq	97
71) Phenanthrene	17.46	178	510566	36.474	nq	99
72) Anthracene	17.54	178	527654	37.759	nq	99
73) Carbazole	17.81	167	467166	37.510	nq	98
74) Di-n-butylphthalate	18.37	149	537632	36.519	nq	100
75) Fluoranthene	19.47	202	702496	41.121	nq	96
77) Benzidine	19.65	184	49575	5.408	nq	# 95
78) Pyrene	19.82	202	704148	35.918	nq	97
80) Butylbenzylphthalate	20.71	149	250736	32.517	nq	94
81) Benzo(a)anthracene	21.67	228	701525	36.065	nq	98
82) 3,3'-Dichlorobenzidine	21.58	252	165706	21.181	nq	98
83) Chrysene	21.73	228	679184	36.205	nq	99
84) Bis(2-ethylhexyl)phthalate	21.57	149	360796	33.700	nq	98
85) Di-n-octyl phthalate	22.78	149	624733	33.928	nq	99
86) Indeno(1,2,3-cd)pyrene	28.65	276	904853	35.493	nq	# 97
88) Benzo(b)fluoranthene	23.91	252	780058	36.538	nq	# 98
89) Benzo(k)fluoranthene	23.97	252	747109	36.407	nq	# 98
90) Benzo(a)pyrene	24.78	252	763947	37.710	nq	# 99
91) Dibenzo(a,h)anthracene	28.70	278	750993	36.613	nq	98
92) Benzo(g,h,i)perylene	29.80	276	758806	36.988	nq	97

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

