

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG010620\
 Data File : BG043943.D
 Acq On : 6 Jan 2020 22:49
 Operator : JU
 Sample : IDOC-W-02-CG
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 IDOC-W-02-CG

Manual Integrations
 APPROVED

mohammad
 1/7/2020 12:30:57 PM

Quant Time: Jan 07 07:08:39 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG123019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 31 13:32:23 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.96	152	108173	20.00	ng	0.00
21) Naphthalene-d8	10.76	136	480952	20.00	ng	0.00
39) Acenaphthene-d10	14.59	164	322988	20.00	ng	0.00
64) Phenanthrene-d10	17.34	188	705809	20.00	ng	0.00
76) Chrysene-d12	21.59	240	619145	20.00	ng	0.00
87) Perylene-d12	24.71	264	696086	20.00	ng	-0.01

System Monitoring Compounds						
5) 2-Fluorophenol	5.54	112	817370	138.74	ng	0.00
7) Phenol-d6	7.13	99	1084369	131.68	ng	0.00
23) Nitrobenzene-d5	9.12	82	686436	76.35	ng	0.00
42) 2,4,6-Tribromophenol	16.08	330	416581	123.25	ng	0.00
45) 2-Fluorobiphenyl	13.22	172	1483772	83.23	ng	0.00
79) Terphenyl-d14	19.95	244	2140052	82.75	ng	0.00

Target Compounds					Qvalue	
2) 1,4-Dioxane	3.43	88	113673	44.252	ng	100
3) Pyridine	3.82	79	281363	37.077	ng	97
4) n-Nitrosodimethylamine	3.74	42	147474	43.650	ng	98
6) Aniline	7.28	93	370866	34.287	ng	97
8) 2-Chlorophenol	7.53	128	352218	50.925	ng	99
9) Benzaldehyde	7.09	77	206398	39.160	ng	99
10) Phenol	7.15	94	437048	51.510	ng	99
11) bis(2-Chloroethyl)ether	7.38	93	322972	44.098	ng	99
12) 1,3-Dichlorobenzene	7.85	146	356260	44.018	ng	97
13) 1,4-Dichlorobenzene	7.99	146	357621	44.782	ng	98
14) 1,2-Dichlorobenzene	8.32	146	337271	44.001	ng	99
15) Benzyl Alcohol	8.19	79	300145	46.181	ng	99
16) 2,2'-oxybis(1-Chloropropan	8.49	45	460834	40.670	ng	99
17) 2-Methylphenol	8.40	107	303265	49.391	ng	97
18) Hexachloroethane	9.04	117	133656	43.013	ng	96
19) n-Nitroso-di-n-propylamine	8.76	70	258196	42.101	ng	97
20) 3+4-Methylphenols	8.73	107	426979	49.848	ng	95
22) Acetophenone	8.77	105	481034	40.231	ng	99
24) Nitrobenzene	9.16	77	375079	42.123	ng	99
25) Isophorone	9.69	82	701298	41.578	ng	99
26) 2-Nitrophenol	9.87	139	192706	45.743	ng	97
27) 2,4-Dimethylphenol	9.94	122	338194	53.330	ng	99
28) bis(2-Chloroethoxy)methane	10.17	93	453702	40.347	ng	99
29) 2,4-Dichlorophenol	10.41	162	349223	46.167	ng	99
30) 1,2,4-Trichlorobenzene	10.63	180	346062	40.802	ng	97
31) Naphthalene	10.81	128	1023703	43.985	ng	100
32) Benzoic acid	10.07	122	137749	29.456	ng	95
33) 4-Chloroaniline	10.91	127	269759	24.301	ng	98
34) Hexachlorobutadiene	11.11	225	200441	38.707	ng	99
35) Caprolactam	11.68	113	130697m	46.413	ng	
36) 4-Chloro-3-methylphenol	12.04	107	372447	46.251	ng	97
37) 2-Methylnaphthalene	12.42	142	773271	44.131	ng	97
38) 1-Methylnaphthalene	12.63	142	731817	44.067	ng	99
40) 1,2,4,5-Tetrachlorobenzene	12.79	216	364919	38.547	ng	99

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41) Hexachlorocyclopentadiene	12.78	237	482924	98.396	nq	99
43) 2,4,6-Trichlorophenol	13.02	196	278285	42.722	nq	99
44) 2,4,5-Trichlorophenol	13.10	196	301468	44.872	nq	100
46) 1,1'-Biphenyl	13.43	154	968037	41.802	nq	99
47) 2-Chloronaphthalene	13.47	162	763134	40.781	nq	98
48) 2-Nitroaniline	13.66	65	244626	40.639	nq	99
49) Acenaphthylene	14.32	152	1200582	44.638	nq	98
50) Dimethylphthalate	14.05	163	972168	42.391	nq	99
51) 2,6-Dinitrotoluene	14.16	165	225914	43.583	nq	99
52) Acenaphthene	14.66	154	850265	45.079	nq	98
53) 3-Nitroaniline	14.49	138	183622	31.195	nq	96
54) 2,4-Dinitrophenol	14.69	184	200376	76.161	nq	97
55) Dibenzofuran	14.99	168	1148527	44.233	nq	99
56) 4-Nitrophenol	14.80	139	437887	97.077	nq	98
57) 2,4-Dinitrotoluene	14.94	165	319622	45.316	nq	99
58) Fluorene	15.64	166	907222	44.082	nq	99
59) 2,3,4,6-Tetrachlorophenol	15.22	232	265425	45.945	nq	97
60) Diethylphthalate	15.41	149	998819	43.544	nq	99
61) 4-Chlorophenyl-phenylether	15.63	204	471483	42.191	nq	99
62) 4-Nitroaniline	15.65	138	244760	42.107	nq	94
63) Azobenzene	15.93	77	935283	43.967	nq	100
65) 4,6-Dinitro-2-methylphenol	15.71	198	163899	45.338	nq	98
66) n-Nitrosodiphenylamine	15.85	169	863998	41.568	nq	99
67) 4-Bromophenyl-phenylether	16.52	248	311828	39.282	nq	99
68) Hexachlorobenzene	16.65	284	331362	38.739	nq	97
69) Atrazine	16.79	200	335922	49.720	nq	99
70) Pentachlorophenol	16.99	266	374908	79.510	nq	100
71) Phenanthrene	17.38	178	1459689	43.117	nq	99
72) Anthracene	17.47	178	1510271	45.140	nq	99
73) Carbazole	17.73	167	1370742	44.353	nq	99
74) Di-n-butylphthalate	18.30	149	1677972	42.525	nq	99
75) Fluoranthene	19.39	202	1694985	43.779	nq	100
77) Benzidine	19.56	184	855058	54.943	nq	99
78) Pyrene	19.74	202	1703077	42.141	nq	99
80) Butylbenzylphthalate	20.64	149	805566	41.868	nq	98
81) Benzo(a)anthracene	21.57	228	1656380	43.145	nq	99
82) 3,3'-Dichlorobenzidine	21.48	252	508342	33.515	nq	99
83) Chrysene	21.63	228	1559884	42.761	nq	99
84) Bis(2-ethylhexyl)phthalate	21.50	149	1120301	42.198	nq	100
85) Di-n-octyl phthalate	22.68	149	1941461	43.499	nq	99
86) Indeno(1,2,3-cd)pyrene	28.25	276	2023759	40.822	nq	# 88
88) Benzo(b)fluoranthene	23.72	252	1762207	42.823	nq	99
89) Benzo(k)fluoranthene	23.79	252	1698697	43.410	nq	99
90) Benzo(a)pyrene	24.56	252	1643140	42.306	nq	99
91) Dibenzo(a,h)anthracene	28.32	278	1673528	42.904	nq	99
92) Benzo(g,h,i)perylene	29.36	276	1666452	43.010	nq	98

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

