

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG031020\
 Data File : BG044714.D
 Acq On : 10 Mar 2020 14:25
 Operator : CG/JU
 Sample : SSTDICC080
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDICC080

Manual Integrations
 APPROVED

mohammad
 3/11/2020 4:56:55 PM

Quant Time: Mar 10 15:48:07 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG031020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 10 14:03:59 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.05	152	104042	20.00	ng	0.00
21) Naphthalene-d8	10.87	136	388742	20.00	ng	0.00
39) Acenaphthene-d10	14.69	164	258600	20.00	ng	0.00
64) Phenanthrene-d10	17.44	188	566467	20.00	ng	0.00
76) Chrysene-d12	21.71	240	480303	20.00	ng	0.00
87) Perylene-d12	24.93	264	581306	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.62	112	1017684	146.90	ng	0.00
7) Phenol-d6	7.21	99	1469958	140.38	ng	0.00
23) Nitrobenzene-d5	9.22	82	1433224	167.86	ng	0.00
42) 2,4,6-Tribromophenol	16.18	330	548633	154.45	ng	0.00
45) 2-Fluorobiphenyl	13.32	172	2666672	143.65	ng	0.00
79) Terphenyl-d14	20.06	244	3502665	130.17	ng	0.00

Target Compounds					Qvalue
2) 1,4-Dioxane	3.53	88	246728	74.301 ng	98
3) Pyridine	3.93	79	750036m	72.849 ng	
4) n-Nitrosodimethylamine	3.83	42	296723	68.132 ng	95
6) Aniline	7.38	93	923867	70.428 ng	98
8) 2-Chlorophenol	7.62	128	514870	73.329 ng	99
9) Benzaldehyde	7.19	77	352917	57.420 ng	97
10) Phenol	7.24	94	735435	71.296 ng	97
11) bis(2-Chloroethyl)ether	7.48	93	576773	72.770 ng	97
12) 1,3-Dichlorobenzene	7.95	146	622449	74.542 ng	97
13) 1,4-Dichlorobenzene	8.09	146	613873	73.584 ng	99
14) 1,2-Dichlorobenzene	8.42	146	581025	73.679 ng	97
15) Benzyl Alcohol	8.29	79	604622	66.929 ng	98
16) 2,2'-oxybis(1-Chloropropan	8.59	45	1019482	68.081 ng	99
17) 2-Methylphenol	8.49	107	496328	70.223 ng	93
18) Hexachloroethane	9.15	117	249155	74.464 ng	98
19) n-Nitroso-di-n-propylamine	8.88	70	517380	66.137 ng	97
20) 3+4-Methylphenols	8.83	107	686990	69.772 ng	99
22) Acetophenone	8.88	105	862450	75.997 ng	# 97
24) Nitrobenzene	9.26	77	755675	83.478 ng	98
25) Isophorone	9.79	82	1360228	72.243 ng	98
26) 2-Nitrophenol	9.97	139	268551	90.720 ng	96
27) 2,4-Dimethylphenol	10.03	122	455558	77.119 ng	98
28) bis(2-Chloroethoxy)methane	10.27	93	800810	75.547 ng	99
29) 2,4-Dichlorophenol	10.51	162	533969	79.445 ng	97
30) 1,2,4-Trichlorobenzene	10.73	180	623765	79.894 ng	97
31) Naphthalene	10.92	128	1668689	76.489 ng	98
32) Benzoic acid	10.21	122	399531	96.440 ng	96
33) 4-Chloroaniline	11.02	127	762839	75.918 ng	98
34) Hexachlorobutadiene	11.22	225	415653	77.923 ng	98
35) Caprolactam	11.81	113	208452m	74.295 ng	
36) 4-Chloro-3-methylphenol	12.14	107	625796	73.520 ng	97
37) 2-Methylnaphthalene	12.52	142	1248766	75.852 ng	97
38) 1-Methylnaphthalene	12.74	142	1174705	75.479 ng	96
40) 1,2,4,5-Tetrachlorobenzene	12.89	216	669357	79.733 ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.88	237	399308	81.341	nq	98
43) 2,4,6-Trichlorophenol	13.12	196	462821	79.524	nq	96
44) 2,4,5-Trichlorophenol	13.19	196	477577	80.016	nq	98
46) 1,1'-Biphenyl	13.53	154	1584790	77.051	nq	95
47) 2-Chloronaphthalene	13.57	162	1231613	78.501	nq	98
48) 2-Nitroaniline	13.76	65	482398	91.167	nq	98
49) Acenaphthylene	14.42	152	1893934	75.348	nq	97
50) Dimethylphthalate	14.15	163	1575903	73.302	nq	99
51) 2,6-Dinitrotoluene	14.26	165	352591	91.272	nq	98
52) Acenaphthene	14.76	154	1282851	78.931	nq	98
53) 3-Nitroaniline	14.59	138	392817	88.235	nq	96
54) 2,4-Dinitrophenol	14.79	184	177134	98.757	nq	# 74
55) Dibenzofuran	15.09	168	1793677	73.652	nq	95
56) 4-Nitrophenol	14.89	139	312763	90.292	nq	97
57) 2,4-Dinitrotoluene	15.04	165	492354	95.384	nq	95
58) Fluorene	15.74	166	1465041	72.733	nq	100
59) 2,3,4,6-Tetrachlorophenol	15.31	232	444167m	78.191	nq	
60) Diethylphthalate	15.52	149	1658719	72.165	nq	99
61) 4-Chlorophenyl-phenylether	15.74	204	810882	74.190	nq	98
62) 4-Nitroaniline	15.76	138	415685	84.772	nq	98
63) Azobenzene	16.03	77	1706848	70.960	nq	96
65) 4,6-Dinitro-2-methylphenol	15.81	198	230788	99.726	nq	98
66) n-Nitrosodiphenylamine	15.95	169	1324421	76.469	nq	98
67) 4-Bromophenyl-phenylether	16.63	248	566166	79.248	nq	98
68) Hexachlorobenzene	16.75	284	608838	78.924	nq	97
69) Atrazine	16.90	200	457897	68.412	nq	100
70) Pentachlorophenol	17.09	266	362556	82.825	nq	97
71) Phenanthrene	17.48	178	2309506	74.354	nq	95
72) Anthracene	17.58	178	2259107	73.844	nq	97
73) Carbazole	17.84	167	2177070	74.306	nq	96
74) Di-n-butylphthalate	18.41	149	2624361	71.122	nq	96
75) Fluoranthene	19.49	202	2672312	71.186	nq	96
77) Benzidine	19.66	184	1227460	64.020	nq	96
78) Pyrene	19.85	202	2652266m	72.397	nq	
80) Butylbenzylphthalate	20.74	149	1310345	80.133	nq	93
81) Benzo(a)anthracene	21.70	228	2666667	74.035	nq	94
82) 3,3'-Dichlorobenzidine	21.61	252	1073128	74.829	nq	97
83) Chrysene	21.77	228	2534842	73.937	nq	94
84) Bis(2-ethylhexyl)phthalate	21.63	149	1723964	75.600	nq	# 96
85) Di-n-octyl phthalate	22.85	149	2914302	74.420	nq	97
86) Indeno(1,2,3-cd)pyrene	28.61	276	3564957	78.734	nq	99
88) Benzo(b)fluoranthene	23.92	252	3017101	76.260	nq	98
89) Benzo(k)fluoranthene	23.99	252	2779728	74.956	nq	95
90) Benzo(a)pyrene	24.79	252	2836416	76.976	nq	98
91) Dibenzo(a,h)anthracene	28.69	278	2806382	76.378	nq	99
92) Benzo(g,h,i)perylene	29.77	276	2860451	77.720	nq	99

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

