

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG031720\
 Data File : BG044836.D
 Acq On : 17 Mar 2020 17:12
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

mohammad
 3/18/2020 4:00:17 PM

Quant Time: Mar 17 18:14:41 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG031020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 16 17:37:15 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.05	152	82487	20.00	ng	0.00
21) Naphthalene-d8	10.85	136	334477	20.00	ng	0.00
39) Acenaphthene-d10	14.68	164	249379	20.00	ng	0.00
64) Phenanthrene-d10	17.43	188	580734	20.00	ng	0.00
76) Chrysene-d12	21.69	240	473659	20.00	ng	0.00
87) Perylene-d12	24.90	264	558430	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.61	112	386586	72.96	ng	0.00
7) Phenol-d6	7.20	99	569918	74.03	ng	0.00
23) Nitrobenzene-d5	9.20	82	560364	73.52	ng	0.00
42) 2,4,6-Tribromophenol	16.16	330	281904	86.33	ng	0.00
45) 2-Fluorobiphenyl	13.30	172	1246948	71.60	ng	0.00
79) Terphenyl-d14	20.04	244	1957767	77.32	ng	0.00

Target Compounds

					Qvalue
2) 1,4-Dioxane	3.52	88	83963m	32.905	ng
3) Pyridine	3.91	79	252436	33.418	ng
4) n-Nitrosodimethylamine	3.83	42	95775	33.416	ng
6) Aniline	7.37	93	350109	36.546	ng
8) 2-Chlorophenol	7.61	128	201827	38.155	ng
9) Benzaldehyde	7.17	77	161381	32.649	ng
10) Phenol	7.23	94	278129	36.490	ng
11) bis(2-Chloroethyl)ether	7.46	93	214784	35.309	ng
12) 1,3-Dichlorobenzene	7.94	146	238779	36.645	ng
13) 1,4-Dichlorobenzene	8.08	146	233657	36.271	ng
14) 1,2-Dichlorobenzene	8.40	146	224644	36.769	ng
15) Benzyl Alcohol	8.28	79	231485	37.474	ng
16) 2,2'-oxybis(1-Chloropropan	8.58	45	345840	32.216	ng
17) 2-Methylphenol	8.48	107	193783	37.533	ng
18) Hexachloroethane	9.14	117	94414	37.227	ng
19) n-Nitroso-di-n-propylamine	8.85	70	195730	36.005	ng
20) 3+4-Methylphenols	8.81	107	269795	37.907	ng
22) Acetophenone	8.87	105	343348	36.234	ng
24) Nitrobenzene	9.24	77	282594	35.731	ng
25) Isophorone	9.77	82	538163	36.175	ng
26) 2-Nitrophenol	9.96	139	118978	43.868	ng
27) 2,4-Dimethylphenol	10.02	122	180027	37.161	ng
28) bis(2-Chloroethoxy)methane	10.25	93	316249	35.622	ng
29) 2,4-Dichlorophenol	10.50	162	219549	38.800	ng
30) 1,2,4-Trichlorobenzene	10.72	180	252423	37.308	ng
31) Naphthalene	10.91	128	672687	36.306	ng
32) Benzoic acid	10.15	122	156592	42.803	ng
33) 4-Chloroaniline	11.01	127	318446	38.148	ng
34) Hexachlorobutadiene	11.20	225	171001	38.432	ng
35) Caprolactam	11.77	113	94690	44.198	ng
36) 4-Chloro-3-methylphenol	12.12	107	266496	39.489	ng
37) 2-Methylnaphthalene	12.51	142	518944	37.442	ng
38) 1-Methylnaphthalene	12.73	142	489145	37.539	ng
40) 1,2,4,5-Tetrachlorobenzene	12.88	216	297575	36.233	ng

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.87	237	158763	35.372	nq	97
43) 2,4,6-Trichlorophenol	13.11	196	210018	38.532	nq	96
44) 2,4,5-Trichlorophenol	13.18	196	220977	39.094	nq	97
46) 1,1'-Biphenyl	13.51	154	699492	35.100	nq	98
47) 2-Chloronaphthalene	13.56	162	526914	34.612	nq	98
48) 2-Nitroaniline	13.75	65	205949	38.642	nq	97
49) Acenaphthylene	14.40	152	848663	35.583	nq	99
50) Dimethylphthalate	14.13	163	738709	37.192	nq	99
51) 2,6-Dinitrotoluene	14.24	165	161572	40.357	nq	97
52) Acenaphthene	14.74	154	580093	37.139	nq	98
53) 3-Nitroaniline	14.57	138	184064	39.989	nq	96
54) 2,4-Dinitrophenol	14.77	184	93907	48.540	nq	# 85
55) Dibenzofuran	15.08	168	827879	36.550	nq	99
56) 4-Nitrophenol	14.87	139	142012	40.427	nq	98
57) 2,4-Dinitrotoluene	15.03	165	234570	42.721	nq	99
58) Fluorene	15.73	166	688479	36.950	nq	99
59) 2,3,4,6-Tetrachlorophenol	15.30	232	214544m	41.227	nq	
60) Diethylphthalate	15.50	149	782617	37.779	nq	99
61) 4-Chlorophenyl-phenylether	15.72	204	378177	37.694	nq	97
62) 4-Nitroaniline	15.73	138	193686	39.462	nq	100
63) Azobenzene	16.01	77	741468	34.467	nq	99
65) 4,6-Dinitro-2-methylphenol	15.80	198	127965	47.014	nq	98
66) n-Nitrosodiphenylamine	15.93	169	625493	35.775	nq	100
67) 4-Bromophenyl-phenylether	16.62	248	269248	37.087	nq	97
68) Hexachlorobenzene	16.74	284	293941	37.320	nq	97
69) Atrazine	16.88	200	239823	37.439	nq	99
70) Pentachlorophenol	17.07	266	179379	41.392	nq	98
71) Phenanthrene	17.47	178	1134128	36.545	nq	100
72) Anthracene	17.56	178	1115606	36.457	nq	99
73) Carbazole	17.82	167	1063249	36.172	nq	99
74) Di-n-butylphthalate	18.39	149	1324277	37.061	nq	99
75) Fluoranthene	19.47	202	1361876	36.857	nq	99
77) Benzidine	19.65	184	602612	39.190	nq	99
78) Pyrene	19.84	202	1351990	38.905	nq	99
80) Butylbenzylphthalate	20.72	149	572985	37.063	nq	98
81) Benzo(a)anthracene	21.67	228	1277201	37.447	nq	99
82) 3,3'-Dichlorobenzidine	21.59	252	500335	37.919	nq	99
83) Chrysene	21.74	228	1203156	36.929	nq	98
84) Bis(2-ethylhexyl)phthalate	21.60	149	774491	36.299	nq	98
85) Di-n-octyl phthalate	22.81	149	1327732	37.187	nq	99
86) Indeno(1,2,3-cd)pyrene	28.56	276	1535915	35.959	nq	# 97
88) Benzo(b)fluoranthene	23.89	252	1349477	36.605	nq	99
89) Benzo(k)fluoranthene	23.95	252	1252839	35.910	nq	98
90) Benzo(a)pyrene	24.75	252	1248348	36.188	nq	97
91) Dibenzo(a,h)anthracene	28.62	278	1245779	36.606	nq	98
92) Benzo(g,h,i)perylene	29.70	276	1245308	36.218	nq	97

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

