

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG050620\
 Data File : BG045249.D
 Acq On : 6 May 2020 12:15
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Sohil
 5/8/2020 5:20:03 PM

Quant Time: May 06 13:03:59 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG041520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 06 12:59:59 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.98	152	75709	20.00	ng	0.00
21) Naphthalene-d8	10.79	136	298964	20.00	ng	0.00
39) Acenaphthene-d10	14.63	164	212300	20.00	ng	0.00
64) Phenanthrene-d10	17.38	188	484382	20.00	ng	0.00
76) Chrysene-d12	21.64	240	395456	20.00	ng	0.00
87) Perylene-d12	24.82	264	463062	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.55	112	345619	75.37	ng	0.00
7) Phenol-d6	7.15	99	501377	73.59	ng	0.00
23) Nitrobenzene-d5	9.14	82	478401	73.02	ng	0.00
42) 2,4,6-Tribromophenol	16.12	330	231310	70.53	ng	0.00
45) 2-Fluorobiphenyl	13.25	172	1102480	76.15	ng	0.00
79) Terphenyl-d14	19.99	244	1622917	89.45	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.44	88	80580	39.539	ng	97
3) Pyridine	3.84	79	195177	31.104	ng	96
4) n-Nitrosodimethylamine	3.76	42	76935	40.023	ng	91
6) Aniline	7.30	93	330883	37.351	ng	100
8) 2-Chlorophenol	7.55	128	195673	39.702	ng	99
9) Benzaldehyde	7.12	77	154643	40.389	ng	98
10) Phenol	7.18	94	267279	39.202	ng	99
11) bis(2-Chloroethyl)ether	7.40	93	194267	35.788	ng	95
12) 1,3-Dichlorobenzene	7.87	146	221651	38.166	ng	98
13) 1,4-Dichlorobenzene	8.02	146	223959	38.701	ng	96
14) 1,2-Dichlorobenzene	8.34	146	212932	38.461	ng	96
15) Benzyl Alcohol	8.22	79	212883	37.267	ng	94
16) 2,2'-oxybis(1-Chloropropan	8.51	45	185841	34.876	ng	97
17) 2-Methylphenol	8.43	107	181767	34.554	ng	96
18) Hexachloroethane	9.07	117	89074	40.945	ng	97
19) n-Nitroso-di-n-propylamine	8.79	70	173402	35.987	ng	97
20) 3+4-Methylphenols	8.76	107	246091	33.423	ng	98
22) Acetophenone	8.80	105	336556	41.628	ng	# 97
24) Nitrobenzene	9.18	77	260238	38.748	ng	96
25) Isophorone	9.71	82	488259	36.389	ng	99
26) 2-Nitrophenol	9.90	139	118888	41.096	ng	97
27) 2,4-Dimethylphenol	9.96	122	186382	40.604	ng	97
28) bis(2-Chloroethoxy)methane	10.19	93	279536	39.651	ng	99
29) 2,4-Dichlorophenol	10.44	162	204357	38.555	ng	95
30) 1,2,4-Trichlorobenzene	10.66	180	235833	39.298	ng	95
31) Naphthalene	10.85	128	636365	38.910	ng	98
32) Benzoic acid	10.11	122	112068	32.582	ng	96
33) 4-Chloroaniline	10.95	127	272031	35.665	ng	98
34) Hexachlorobutadiene	11.14	225	165343	40.186	ng	98
35) Caprolactam	11.72	113	75854	35.958	ng	94
36) 4-Chloro-3-methylphenol	12.08	107	235230	37.779	ng	98
37) 2-Methylnaphthalene	12.46	142	482887	38.040	ng	97
38) 1-Methylnaphthalene	12.67	142	451401	37.667	ng	97
40) 1,2,4,5-Tetrachlorobenzene	12.83	216	281201	41.138	ng	100

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.81	237	185878	43.508	nq	100
43) 2,4,6-Trichlorophenol	13.06	196	187625	38.892	nq	97
44) 2,4,5-Trichlorophenol	13.14	196	197953	36.049	nq	97
46) 1,1'-Biphenyl	13.46	154	637281	39.494	nq	99
47) 2-Chloronaphthalene	13.50	162	472592	38.329	nq	98
48) 2-Nitroaniline	13.70	65	151473	37.339	nq	96
49) Acenaphthylene	14.35	152	788463	39.259	nq	99
50) Dimethylphthalate	14.08	163	629830	36.435	nq	99
51) 2,6-Dinitrotoluene	14.19	165	144959	37.491	nq	93
52) Acenaphthene	14.69	154	505037m	37.167	nq	
53) 3-Nitroaniline	14.52	138	143661	33.877	nq	# 97
54) 2,4-Dinitrophenol	14.73	184	80884	35.180	nq	97
55) Dibenzofuran	15.03	168	752859	38.003	nq	99
56) 4-Nitrophenol	14.85	139	113830	34.620	nq	99
57) 2,4-Dinitrotoluene	14.99	165	208107	36.830	nq	# 95
58) Fluorene	15.67	166	615751	38.052	nq	98
59) 2,3,4,6-Tetrachlorophenol	15.26	232	187740	38.424	nq	98
60) Diethylphthalate	15.45	149	659882	36.614	nq	98
61) 4-Chlorophenyl-phenylether	15.67	204	339753	36.478	nq	99
62) 4-Nitroaniline	15.69	138	157635	35.839	nq	97
63) Azobenzene	15.96	77	625767	37.660	nq	97
65) 4,6-Dinitro-2-methylphenol	15.76	198	133671	41.984	nq	99
66) n-Nitrosodiphenylamine	15.88	169	542935	39.012	nq	100
67) 4-Bromophenyl-phenylether	16.57	248	225494	36.085	nq	98
68) Hexachlorobenzene	16.69	284	249260	38.461	nq	97
69) Atrazine	16.83	200	222686	43.735	nq	98
70) Pentachlorophenol	17.03	266	146143	36.758	nq	99
71) Phenanthrene	17.42	178	966213	38.818	nq	99
72) Anthracene	17.51	178	988723	39.599	nq	100
73) Carbazole	17.78	167	922821	36.420	nq	98
74) Di-n-butylphthalate	18.34	149	1097271	38.969	nq	99
75) Fluoranthene	19.43	202	1150266	38.413	nq	99
77) Benzidine	19.60	184	549374	48.416	nq	96
78) Pyrene	19.79	202	1141053	41.854	nq	98
80) Butylbenzylphthalate	20.68	149	472339	41.797	nq	96
81) Benzo(a)anthracene	21.63	228	1050172	40.973	nq	98
82) 3,3'-Dichlorobenzidine	21.54	252	442439	43.369	nq	98
83) Chrysene	21.69	228	1034829	42.020	nq	99
84) Bis(2-ethylhexyl)phthalate	21.54	149	648383	41.717	nq	99
85) Di-n-octyl phthalate	22.74	149	1091530	40.612	nq	99
86) Indeno(1,2,3-cd)pyrene	28.42	276	1330102	40.680	nq	# 97
88) Benzo(b)fluoranthene	23.82	252	1102434	38.007	nq	98
89) Benzo(k)fluoranthene	23.88	252	1110680	40.264	nq	98
90) Benzo(a)pyrene	24.67	252	1095247	39.993	nq	# 97
91) Dibenzo(a,h)anthracene	28.49	278	1128659	40.900	nq	99
92) Benzo(g,h,i)perylene	29.56	276	1095923	39.612	nq	98

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

