

Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG052920\
 Data File : BG045535.D
 Acq On : 30 May 2020 00:41
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

mohammad
 6/1/2020 12:22:47 PM

Quant Time: May 30 06:36:57 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG051520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri May 29 11:17:45 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.95	152	64891	20.00	ng	0.00
21) Naphthalene-d8	10.76	136	274063	20.00	ng	0.00
39) Acenaphthene-d10	14.60	164	207783	20.00	ng	0.00
64) Phenanthrene-d10	17.35	188	488729	20.00	ng	0.00
76) Chrysene-d12	21.61	240	459253	20.00	ng	0.00
87) Perylene-d12	24.76	264	516154	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.55	112	284929	85.81	ng	0.00
7) Phenol-d6	7.16	99	422235	89.34	ng	-0.01
23) Nitrobenzene-d5	9.12	82	426008	84.76	ng	0.00
42) 2,4,6-Tribromophenol	16.10	330	219703	82.68	ng	0.00
45) 2-Fluorobiphenyl	13.22	172	988211	80.67	ng	0.00
79) Terphenyl-d14	19.96	244	1644398	83.51	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.40	88	65575	39.83	ng	98
3) Pyridine	3.80	79	193484	42.19	ng	95
4) n-Nitrosodimethylamine	3.72	42	64669	43.10	ng	95
6) Aniline	7.28	93	268019	43.44	ng	98
8) 2-Chlorophenol	7.53	128	158046	43.40	ng	98
9) Benzaldehyde	7.08	77	127589	41.44	ng	96
10) Phenol	7.18	94	213978	43.93	ng	96
11) bis(2-Chloroethyl)ether	7.37	93	166881	43.93	ng	98
12) 1,3-Dichlorobenzene	7.84	146	188420	43.06	ng	98
13) 1,4-Dichlorobenzene	7.99	146	189841	43.96	ng	98
14) 1,2-Dichlorobenzene	8.31	146	181891	43.79	ng	96
15) Benzyl Alcohol	8.20	79	176227	45.14	ng	98
16) 2,2'-oxybis(1-Chloropropan	8.48	45	157209	43.59	ng	97
17) 2-Methylphenol	8.42	107	163327	44.80	ng	96
18) Hexachloroethane	9.04	117	73316	44.33	ng	97
19) n-Nitroso-di-n-propylamine	8.76	70	148187	45.34	ng	98
20) 3+4-Methylphenols	8.75	107	222193	44.76	ng	92
22) Acetophenone	8.78	105	271623	41.82	ng	# 94
24) Nitrobenzene	9.16	77	226504	42.07	ng	98
25) Isophorone	9.68	82	421796	42.62	ng	99
26) 2-Nitrophenol	9.87	139	100129	43.47	ng	94
27) 2,4-Dimethylphenol	9.95	122	148111	43.35	ng	96
28) bis(2-Chloroethoxy)methane	10.17	93	223306	43.02	ng	97
29) 2,4-Dichlorophenol	10.43	162	176448	43.74	ng	94
30) 1,2,4-Trichlorobenzene	10.63	180	207582	43.54	ng	97
31) Naphthalene	10.81	128	541495	42.40	ng	100
32) Benzoic acid	10.13	122	102681	44.99	ng	99
33) 4-Chloroaniline	10.93	127	234037	43.84	ng	97
34) Hexachlorobutadiene	11.10	225	143234	42.51	ng	95
35) Caprolactam	11.69	113	60674	40.97	ng	93
36) 4-Chloro-3-methylphenol	12.08	107	198651	43.70	ng	99
37) 2-Methylnaphthalene	12.42	142	412573	43.34	ng	99
38) 1-Methylnaphthalene	12.64	142	391606m	43.55	ng	
40) 1,2,4,5-Tetrachlorobenzene	12.79	216	235904	40.65	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.78	237	111803	33.38	ng	99
43) 2,4,6-Trichlorophenol	13.04	196	162953	40.42	ng	96
44) 2,4,5-Trichlorophenol	13.12	196	188523	40.92	ng	94
46) 1,1'-Biphenyl	13.43	154	515430	40.37	ng	99
47) 2-Chloronaphthalene	13.48	162	418836	40.40	ng	99
48) 2-Nitroaniline	13.68	65	135631	39.77	ng	91
49) Acenaphthylene	14.32	152	677821	40.70	ng	99
50) Dimethylphthalate	14.06	163	568770	39.90	ng	100
51) 2,6-Dinitrotoluene	14.18	165	128813	40.05	ng	94
52) Acenaphthene	14.66	154	457715m	41.06	ng	
53) 3-Nitroaniline	14.51	138	133194	40.74	ng	99
54) 2,4-Dinitrophenol	14.72	184	73775	38.08	ng	98
55) Dibenzofuran	15.00	168	675830	40.83	ng	97
56) 4-Nitrophenol	14.86	139	95207	38.46	ng	91
57) 2,4-Dinitrotoluene	14.96	165	187614	40.28	ng	98
58) Fluorene	15.65	166	565567	41.59	ng	97
59) 2,3,4,6-Tetrachlorophenol	15.24	232	166165	41.00	ng	97
60) Diethylphthalate	15.42	149	594376	40.06	ng	98
61) 4-Chlorophenyl-phenylether	15.64	204	322473	41.44	ng	100
62) 4-Nitroaniline	15.67	138	134381	39.37	ng	98
63) Azobenzene	15.94	77	560564	40.52	ng	97
65) 4,6-Dinitro-2-methylphenol	15.73	198	115950	40.74	ng	96
66) n-Nitrosodiphenylamine	15.86	169	490925	41.84	ng	99
67) 4-Bromophenyl-phenylether	16.54	248	231786	43.80	ng	96
68) Hexachlorobenzene	16.66	284	233799	42.24	ng	95
69) Atrazine	16.81	200	189395	39.19	ng	98
70) Pentachlorophenol	17.01	266	117574	40.91	ng	99
71) Phenanthrene	17.39	178	895890	41.99	ng	99
72) Anthracene	17.48	178	892038	41.63	ng	97
73) Carbazole	17.75	167	852636	40.82	ng	99
74) Di-n-butylphthalate	18.31	149	1007493	40.19	ng	98
75) Fluoranthene	19.40	202	1101980	40.61	ng	100
77) Benzidine	19.58	184	500426	38.80	ng	100
78) Pyrene	19.76	202	1094774	41.82	ng	99
80) Butylbenzylphthalate	20.65	149	457009	40.44	ng	98
81) Benzo(a)anthracene	21.59	228	1073555	41.96	ng	99
82) 3,3'-Dichlorobenzidine	21.51	252	414436	41.30	ng	99
83) Chrysene	21.65	228	1036429	42.13	ng	99
84) Bis(2-ethylhexyl)phthalate	21.50	149	646288	41.61	ng	99
85) Di-n-octyl phthalate	22.69	149	1110226	41.62	ng	100
86) Indeno(1,2,3-cd)pyrene	28.34	276	1359791	42.27	ng	# 97
88) Benzo(b)fluoranthene	23.76	252	1142399	41.27	ng	99
89) Benzo(k)fluoranthene	23.83	252	1116813	42.94	ng	99
90) Benzo(a)pyrene	24.61	252	1089758	42.27	ng	97
91) Dibenzo(a,h)anthracene	28.39	278	1082400	41.78	ng	100
92) Benzo(g,h,i)perylene	29.45	276	1076986	41.57	ng	99

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

