

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG070620\
 Data File : BG045945.D
 Acq On : 6 Jul 2020 18:20
 Operator : CG/JU
 Sample : SSTDICC010
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDICC010

Manual Integrations
 APPROVED

mohammad
 7/7/2020 12:03:22 PM

Quant Time: Jul 07 01:23:39 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG070620.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 07 00:57:59 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.01	152	128581	20.00	ng	0.00
21) Naphthalene-d8	10.81	136	521798	20.00	ng	0.00
39) Acenaphthene-d10	14.63	164	314743	20.00	ng	0.00
64) Phenanthrene-d10	17.37	188	651102	20.00	ng	0.00
76) Chrysene-d12	21.63	240	635331	20.00	ng	0.00
86) Perylene-d12	24.79	264	674769	20.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.58	112	165353	21.72	ng	0.00
7) Phenol-d6	7.16	99	239555	20.69	ng	0.00
23) Nitrobenzene-d5	9.16	82	165398	14.48	ng	0.00
42) 2,4,6-Tribromophenol	16.11	330	52710	11.09	ng	0.00
45) 2-Fluorobiphenyl	13.26	172	426585	19.92	ng	0.00
79) Terphenyl-d14	19.99	244	621714	19.83	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	3.52	88	35220	10.103	ng	# 87
3) Pyridine	3.90	79	110255	10.672	ng	94
4) n-Nitrosodimethylamine	3.82	42	35267	10.170	ng	# 92
6) Aniline	7.33	93	151967	11.147	ng	98
8) 2-Chlorophenol	7.57	128	87606	10.752	ng	99
9) Benzaldehyde	7.14	77	82512	11.865	ng	95
10) Phenol	7.19	94	118960	10.604	ng	96
11) bis(2-Chloroethyl)ether	7.43	93	99719	11.703	ng	93
12) 1,3-Dichlorobenzene	7.90	146	103119	10.620	ng	98
13) 1,4-Dichlorobenzene	8.05	146	99690	10.279	ng	98
14) 1,2-Dichlorobenzene	8.36	146	98779	10.639	ng	98
15) Benzyl Alcohol	8.24	79	89396	9.977	ng	97
16) 2,2'-oxybis(1-Chloropropan	8.54	45	127484	15.874	ng	97
17) 2-Methylphenol	8.44	107	90984	10.898	ng	93
18) Hexachloroethane	9.10	117	37865	10.280	ng	99
19) n-Nitroso-di-n-propylamine	8.81	70	79576	10.454	ng	99
20) 3+4-Methylphenols	8.77	107	120654	10.988	ng	95
22) Acetophenone	8.82	105	137517	9.602	ng	98
24) Nitrobenzene	9.20	77	90298	7.414	ng	97
25) Isophorone	9.73	82	219909	9.945	ng	99
26) 2-Nitrophenol	9.91	139	26860	5.294	ng	# 92
27) 2,4-Dimethylphenol	9.97	122	80160	10.348	ng	96
28) bis(2-Chloroethoxy)methane	10.21	93	123239	10.670	ng	99
29) 2,4-Dichlorophenol	10.45	162	77517	8.577	ng	98
30) 1,2,4-Trichlorobenzene	10.67	180	91261	8.530	ng	99
31) Naphthalene	10.85	128	290859	10.196	ng	97
32) Benzoic acid	10.04	122	25013m	5.647	ng	
33) 4-Chloroaniline	10.95	127	123420	10.331	ng	97
34) Hexachlorobutadiene	11.16	225	51265	6.737	ng	97
35) Caprolactam	11.69	113	29540	10.100	ng	# 85
36) 4-Chloro-3-methylphenol	12.07	107	89169	8.545	ng	99
37) 2-Methylnaphthalene	12.46	142	201571	9.489	ng	98
38) 1-Methylnaphthalene	12.68	142	187325	9.319	ng	99
40) 1,2,4,5-Tetrachlorobenzene	12.83	216	87664	8.459	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.82	237	50366	10.709	nq	98
43) 2,4,6-Trichlorophenol	13.06	196	56621	8.058	nq	97
44) 2,4,5-Trichlorophenol	13.13	196	61957	7.737	nq	94
46) 1,1'-Biphenyl	13.47	154	238919	10.603	nq	96
47) 2-Chloronaphthalene	13.51	162	190707	10.477	nq	98
48) 2-Nitroaniline	13.69	65	33781	5.579	nq	96
49) Acenaphthylene	14.35	152	311425	10.681	nq	96
50) Dimethylphthalate	14.08	163	238052	9.655	nq	98
51) 2,6-Dinitrotoluene	14.19	165	28055	5.066	nq	98
52) Acenaphthene	14.70	154	185424	10.492	nq	96
53) 3-Nitroaniline	14.51	138	35879	6.486	nq #	93
54) 2,4-Dinitrophenol	14.72	184	8868	3.709	nq #	51
55) Dibenzofuran	15.02	168	288327	10.043	nq	98
56) 4-Nitrophenol	14.81	139	28616	8.029	nq	94
57) 2,4-Dinitrotoluene	14.98	165	34206	4.311	nq #	95
58) Fluorene	15.68	166	234965	9.842	nq	97
59) 2,3,4,6-Tetrachlorophenol	15.25	232	47396m	6.819	nq	
60) Diethylphthalate	15.45	149	248246	9.849	nq	97
61) 4-Chlorophenyl-phenylether	15.67	204	113870	8.276	nq	99
62) 4-Nitroaniline	15.68	138	39158	6.898	nq	90
63) Azobenzene	15.96	77	269857	11.150	nq	96
65) 4,6-Dinitro-2-methylphenol	15.74	198	12311	3.060	nq	90
66) n-Nitrosodiphenylamine	15.88	169	202963	11.006	nq	97
67) 4-Bromophenyl-phenylether	16.56	248	69946	8.268	nq	97
68) Hexachlorobenzene	16.69	284	72512	8.266	nq	96
69) Atrazine	16.83	200	67296	9.404	nq	98
70) Pentachlorophenol	17.02	266	35429	9.838	nq	98
71) Phenanthrene	17.42	178	356612	10.620	nq	98
72) Anthracene	17.50	178	365955	10.944	nq	97
73) Carbazole	17.77	167	362722	11.392	nq	97
74) Di-n-butylphthalate	18.35	149	457411	12.133	nq	96
75) Fluoranthene	19.42	202	425356	10.301	nq	95
77) Benzidine	19.59	184	199723	12.092	nq	99
78) Pyrene	19.78	202	435646	10.240	nq	94
80) Butylbenzylphthalate	20.68	149	199862	11.223	nq	95
81) Benzo(a)anthracene	21.61	228	430492	10.451	nq	94
82) 3,3'-Dichlorobenzidine	21.53	252	157120	10.084	nq	99
83) Chrysene	21.67	228	403397	10.125	nq	97
84) Bis(2-ethylhexyl)phthalate	21.56	149	299411	12.088	nq	97
85) Di-n-octyl phthalate	22.76	149	502858	11.769	nq	99
87) Indeno(1,2,3-cd)pyrene	28.36	276	460858	9.421	nq	99
88) Benzo(b)fluoranthene	23.79	252	421901	9.972	nq	99
89) Benzo(k)fluoranthene	23.85	252	399452	9.865	nq	98
90) Benzo(a)pyrene	24.64	252	391842	9.915	nq	99
91) Dibenzo(a,h)anthracene	28.44	278	382550	9.752	nq	97
92) Benzo(g,h,i)perylene	29.47	276	374163	9.530	nq	99

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

