

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG062619\
 Data File : BG041477.D
 Acq On : 26 Jun 2019 21:20
 Operator : HP/JU
 Sample : SSTDICC100
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDICC100

Quant Time: Jun 27 06:52:20 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG062619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 27 05:39:03 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.03	152	24119	20.00	ng	0.00
21) Naphthalene-d8	10.85	136	95403	20.00	ng	0.00
39) Acenaphthene-d10	14.68	164	69170	20.00	ng	0.00
64) Phenanthrene-d10	17.43	188	186491	20.00	ng	0.00
76) Chrysene-d12	21.71	240	198765	20.00	ng	0.00
87) Perylene-d12	24.99	264	221015	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.57	112	251594	194.56	ng	0.00
7) Phenol-d6	7.19	99	360884	195.97	ng	0.00
23) Nitrobenzene-d5	9.22	82	414015	184.38	ng	0.00
42) 2,4,6-Tribromophenol	16.17	330	220537	198.26	ng	0.00
45) 2-Fluorobiphenyl	13.30	172	961767	181.07	ng	0.00
79) Terphenyl-d14	20.03	244	1571077	155.54	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.41	88	55595	94.532	ng	97
3) Pyridine	3.82	79	147799	96.493	ng	96
4) n-Nitrosodimethylamine	3.75	42	82129	94.025	ng	# 94
6) Aniline	7.36	93	232889	96.271	ng	97
8) 2-Chlorophenol	7.59	128	139376	92.137	ng	95
10) Phenol	7.22	94	184156	95.195	ng	97
11) bis(2-Chloroethyl)ether	7.45	93	139096	88.588	ng	99
12) 1,3-Dichlorobenzene	7.92	146	180242	93.243	ng	98
13) 1,4-Dichlorobenzene	8.07	146	180922	92.176	ng	99
14) 1,2-Dichlorobenzene	8.39	146	173537	92.033	ng	97
15) Benzyl Alcohol	8.28	79	154072	103.306	ng	97
16) 2,2'-oxybis(1-Chloropropan	8.56	45	197660	91.770	ng	99
17) 2-Methylphenol	8.48	107	132165	94.522	ng	98
18) Hexachloroethane	9.11	117	71842	93.116	ng	95
19) n-Nitroso-di-n-propylamine	8.85	70	132581	92.844	ng	99
20) 3+4-Methylphenols	8.81	107	180305	95.150	ng	97
22) Acetophenone	8.87	105	260696	93.096	ng	# 99
24) Nitrobenzene	9.26	77	206589	91.699	ng	96
25) Isophorone	9.77	82	378801	92.837	ng	99
26) 2-Nitrophenol	9.97	139	90056	96.203	ng	95
27) 2,4-Dimethylphenol	10.02	122	123896	91.520	ng	98
28) bis(2-Chloroethoxy)methane	10.26	93	199686	93.668	ng	99
29) 2,4-Dichlorophenol	10.51	162	170256	94.994	ng	96
30) 1,2,4-Trichlorobenzene	10.71	180	212462	94.855	ng	92
31) Naphthalene	10.91	128	478380	92.488	ng	98
33) 4-Chloroaniline	11.03	127	208859	93.985	ng	96
34) Hexachlorobutadiene	11.17	225	166956	94.863	ng	97
35) Caprolactam	11.84	113	54698	98.684	ng	96
36) 4-Chloro-3-methylphenol	12.14	107	186979	96.707	ng	99
37) 2-Methylnaphthalene	12.51	142	359922	93.200	ng	97
38) 1-Methylnaphthalene	12.73	142	339367	93.078	ng	98
40) 1,2,4,5-Tetrachlorobenzene	12.87	216	273524	91.920	ng	100
41) Hexachlorocyclopentadiene	12.83	237	163403	101.604	ng	94
43) 2,4,6-Trichlorophenol	13.12	196	166527	93.955	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.19	196	165741	94.969	ng	97
46) 1,1'-Biphenyl	13.51	154	498680	93.302	ng	95
47) 2-Chloronaphthalene	13.56	162	425655	91.770	ng	98
48) 2-Nitroaniline	13.78	65	148262	94.613	ng	90
49) Acenaphthylene	14.41	152	631403	91.115	ng	98
50) Dimethylphthalate	14.14	163	572841	92.291	ng	98
51) 2,6-Dinitrotoluene	14.27	165	123009	93.472	ng	93
52) Acenaphthene	14.74	154	376862	92.583	ng	99
53) 3-Nitroaniline	14.61	138	121450	96.120	ng	96
55) Dibenzofuran	15.08	168	648008	91.053	ng	97
56) 4-Nitrophenol	14.91	139	85371	103.254	ng	97
57) 2,4-Dinitrotoluene	15.06	165	179699	95.208	ng	99
58) Fluorene	15.73	166	524544	92.242	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.30	232	172876	93.363	ng	99
60) Diethylphthalate	15.48	149	581847	91.655	ng	99
61) 4-Chlorophenyl-phenylether	15.71	204	341823	93.187	ng	99
62) 4-Nitroaniline	15.78	138	121287	97.706	ng	99
63) Azobenzene	16.01	77	490749	91.631	ng	99
65) 4,6-Dinitro-2-methylphenol	15.82	198	116840	99.285	ng	100
66) n-Nitrosodiphenylamine	15.93	169	464587	91.412	ng	99
67) 4-Bromophenyl-phenylether	16.61	248	232225	93.477	ng	98
68) Hexachlorobenzene	16.72	284	248648	92.797	ng	99
70) Pentachlorophenol	17.07	266	120943	100.709	ng	98
71) Phenanthrene	17.47	178	908276	90.206	ng	98
72) Anthracene	17.56	178	907329	90.148	ng	98
73) Carbazole	17.84	167	777478	90.218	ng	98
74) Di-n-butylphthalate	18.36	149	973890	89.963	ng	98
75) Fluoranthene	19.48	202	1192553	89.002	ng	96
77) Benzidine	19.67	184	357454	89.824	ng	99
78) Pyrene	19.84	202	1181845	88.134	ng	96
80) Butylbenzylphthalate	20.71	149	473183	92.913	ng	98
81) Benzo(a)anthracene	21.68	228	1223653	90.393	ng	96
82) 3,3'-Dichlorobenzidine	21.60	252	439680	93.996	ng	96
83) Chrysene	21.75	228	1162184	89.087	ng	98
84) Bis(2-ethylhexyl)phthalate	21.55	149	660251	92.755	ng	99
85) Di-n-octyl phthalate	22.77	149	1131399	94.317	ng	100
86) Indeno(1,2,3-cd)pyrene	28.78	276	1521154	97.021	ng	# 98
88) Benzo(b)fluoranthene	23.94	252	1262281	91.621	ng	98
89) Benzo(k)fluoranthene	24.02	252	1227192	90.297	ng	99
90) Benzo(a)pyrene	24.84	252	1218488	93.486	ng	98
91) Dibenzo(a,h)anthracene	28.85	278	1218819	93.459	ng	98
92) Benzo(g,h,i)perylene	29.98	276	1220281	94.041	ng	100

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

