

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG062620\
 Data File : BG045867.D
 Acq On : 26 Jun 2020 10:43
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040

Quant Time: Jun 26 12:18:03 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG061520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 25 15:44:32 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.92	152	54131	20.00	ng	0.00
21) Naphthalene-d8	10.72	136	212098	20.00	ng	0.00
39) Acenaphthene-d10	14.57	164	163459	20.00	ng	0.00
64) Phenanthrene-d10	17.32	188	406411	20.00	ng	0.00
76) Chrysene-d12	21.57	240	408560	20.00	ng	0.00
86) Perylene-d12	24.69	264	456073	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.51	112	256131	79.93	ng	0.00
7) Phenol-d6	7.13	99	405216	83.13	ng	0.00
23) Nitrobenzene-d5	9.08	82	379807	81.81	ng	0.00
42) 2,4,6-Tribromophenol	16.06	330	207292	84.00	ng	0.00
45) 2-Fluorobiphenyl	13.18	172	892362	80.23	ng	0.00
79) Terphenyl-d14	19.93	244	1602772	79.50	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.36	88	55641	37.911	ng	95
3) Pyridine	3.76	79	177423	40.795	ng	97
4) n-Nitrosodimethylamine	3.68	42	64372	44.095	ng	# 99
6) Aniline	7.24	93	235905	41.104	ng	99
8) 2-Chlorophenol	7.49	128	143189	41.746	ng	96
9) Benzaldehyde	7.05	77	119334	40.761	ng	98
10) Phenol	7.16	94	196831	41.676	ng	97
11) bis(2-Chloroethyl)ether	7.33	93	140988	39.302	ng	97
12) 1,3-Dichlorobenzene	7.80	146	164494	40.242	ng	98
13) 1,4-Dichlorobenzene	7.95	146	166669	40.822	ng	99
14) 1,2-Dichlorobenzene	8.27	146	156209	39.964	ng	95
15) Benzyl Alcohol	8.17	79	162384	43.050	ng	97
16) 2,2'-oxybis(1-Chloropropan	8.44	45	139663	41.309	ng	79
17) 2-Methylphenol	8.40	107	146893	41.793	ng	96
18) Hexachloroethane	8.99	117	62802	40.500	ng	91
19) n-Nitroso-di-n-propylamine	8.72	70	139749	43.610	ng	97
20) 3+4-Methylphenols	8.73	107	198258	42.890	ng	89
22) Acetophenone	8.74	105	234477	40.278	ng	# 98
24) Nitrobenzene	9.12	77	201082	40.620	ng	99
25) Isophorone	9.64	82	386227	42.971	ng	99
26) 2-Nitrophenol	9.83	139	85458	41.437	ng	94
27) 2,4-Dimethylphenol	9.92	122	134426	42.692	ng	98
28) bis(2-Chloroethoxy)methane	10.12	93	194928	41.519	ng	100
29) 2,4-Dichlorophenol	10.40	162	157895	42.981	ng	99
30) 1,2,4-Trichlorobenzene	10.58	180	176339	40.547	ng	98
31) Naphthalene	10.77	128	471045	40.622	ng	99
32) Benzoic acid	10.13	122	89580	44.640	ng	95
33) 4-Chloroaniline	10.89	127	209758	43.196	ng	97
34) Hexachlorobutadiene	11.06	225	127128	41.102	ng	97
35) Caprolactam	11.67	113	59182	49.783	ng	94
36) 4-Chloro-3-methylphenol	12.05	107	184017	43.384	ng	99
37) 2-Methylnaphthalene	12.38	142	363667	42.117	ng	98
38) 1-Methylnaphthalene	12.60	142	345016	42.224	ng	96
40) 1,2,4,5-Tetrachlorobenzene	12.76	216	209142	38.857	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.73	237	90263	35.585	ng	97
43) 2,4,6-Trichlorophenol	13.01	196	142694	39.103	ng	100
44) 2,4,5-Trichlorophenol	13.10	196	163322	39.269	ng	98
46) 1,1'-Biphenyl	13.39	154	467639	39.960	ng	98
47) 2-Chloronaphthalene	13.44	162	373271	39.486	ng	98
48) 2-Nitroaniline	13.65	65	136089	43.278	ng	94
49) Acenaphthylene	14.28	152	609656	40.263	ng	99
50) Dimethylphthalate	14.03	163	526203	41.093	ng	98
51) 2,6-Dinitrotoluene	14.14	165	122657	42.647	ng	98
52) Acenaphthene	14.62	154	370151	40.329	ng	99
53) 3-Nitroaniline	14.48	138	121892	42.430	ng	96
54) 2,4-Dinitrophenol	14.69	184	65131	45.540	ng	# 91
55) Dibenzofuran	14.97	168	616899	41.375	ng	97
56) 4-Nitrophenol	14.85	139	78843	40.377	ng	99
57) 2,4-Dinitrotoluene	14.94	165	177978	43.190	ng	98
58) Fluorene	15.62	166	514374	41.488	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.21	232	150945	41.815	ng	98
60) Diethylphthalate	15.39	149	561515	42.895	ng	98
61) 4-Chlorophenyl-phenylether	15.61	204	299344	41.891	ng	98
62) 4-Nitroaniline	15.65	138	134648	45.669	ng	95
63) Azobenzene	15.90	77	529690	42.141	ng	97
65) 4,6-Dinitro-2-methylphenol	15.71	198	110165	43.862	ng	92
66) n-Nitrosodiphenylamine	15.82	169	456143	39.626	ng	99
67) 4-Bromophenyl-phenylether	16.50	248	210819	39.922	ng	99
68) Hexachlorobenzene	16.63	284	223097	40.742	ng	94
69) Atrazine	16.77	200	162869	36.462	ng	96
70) Pentachlorophenol	16.98	266	88375	37.235	ng	95
71) Phenanthrene	17.36	178	848334	40.473	ng	98
72) Anthracene	17.45	178	844818	40.476	ng	99
73) Carbazole	17.73	167	835241	42.027	ng	99
74) Di-n-butylphthalate	18.28	149	995581	42.308	ng	99
75) Fluoranthene	19.37	202	1092747	42.397	ng	99
77) Benzidine	19.55	184	450628	42.427	ng	99
78) Pyrene	19.73	202	1098180	40.140	ng	100
80) Butylbenzylphthalate	20.62	149	472825	41.289	ng	100
81) Benzo(a)anthracene	21.56	228	1088628	41.098	ng	98
82) 3,3'-Dichlorobenzidine	21.47	252	415547	41.474	ng	98
83) Chrysene	21.62	228	1041064	40.634	ng	99
84) Bis(2-ethylhexyl)phthalate	21.46	149	655844	41.174	ng	98
85) Di-n-octyl phthalate	22.64	149	1136582	41.366	ng	99
87) Indeno(1,2,3-cd)pyrene	28.25	276	1380419	41.751	ng	# 99
88) Benzo(b)fluoranthene	23.71	252	1179860	41.260	ng	98
89) Benzo(k)fluoranthene	23.78	252	1111097	40.598	ng	97
90) Benzo(a)pyrene	24.55	252	1101350	41.231	ng	99
91) Dibenzo(a,h)anthracene	28.31	278	1107119	41.756	ng	99
92) Benzo(g,h,i)perylene	29.36	276	1108552	41.772	ng	98

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

