

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG083119\
 Data File : BG042623.D
 Acq On : 31 Aug 2019 7:31
 Operator : HP/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040

Quant Time: Aug 31 13:28:44 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG082219.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 22 14:29:15 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.99	152	33577	20.00	ng	-0.01
21) Naphthalene-d8	10.81	136	138265	20.00	ng	-0.01
39) Acenaphthene-d10	14.66	164	105624	20.00	ng	0.00
64) Phenanthrene-d10	17.41	188	240557	20.00	ng	-0.01
76) Chrysene-d12	21.68	240	250094	20.00	ng	-0.01
87) Perylene-d12	24.92	264	302476	20.00	ng	-0.02

System Monitoring Compounds

5) 2-Fluorophenol	5.55	112	130990	79.01	ng	-0.01
7) Phenol-d6	7.16	99	172723	77.13	ng	-0.01
23) Nitrobenzene-d5	9.16	82	157487	78.71	ng	-0.01
42) 2,4,6-Tribromophenol	16.15	330	121486	80.92	ng	-0.01
45) 2-Fluorobiphenyl	13.28	172	525048	84.07	ng	0.00
79) Terphenyl-d14	20.02	244	981782	84.87	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.42	88	26608	38.458	ng	98
3) Pyridine	3.82	79	68509	38.616	ng	99
4) n-Nitrosodimethylamine	3.73	42	25961	40.971	ng	85
6) Aniline	7.31	93	114132	38.211	ng	99
8) 2-Chlorophenol	7.56	128	80527	40.078	ng	98
9) Benzaldehyde	7.12	77	43237	38.565	ng	99
10) Phenol	7.19	94	86587	38.028	ng	96
11) bis(2-Chloroethyl)ether	7.41	93	65122	36.418	ng	99
12) 1,3-Dichlorobenzene	7.89	146	102722	40.189	ng	98
13) 1,4-Dichlorobenzene	8.03	146	106100	40.980	ng	98
14) 1,2-Dichlorobenzene	8.35	146	100430	40.506	ng	96
15) Benzyl Alcohol	8.24	79	63667	39.517	ng	97
16) 2,2'-oxybis(1-Chloropropan	8.53	45	88695	36.595	ng	98
17) 2-Methylphenol	8.45	107	65351	38.967	ng	94
18) Hexachloroethane	9.09	117	34916	39.393	ng	90
19) n-Nitroso-di-n-propylamine	8.81	70	54625	37.651	ng	98
20) 3+4-Methylphenols	8.78	107	87256	37.600	ng	96
22) Acetophenone	8.82	105	118109	39.939	ng	# 99
24) Nitrobenzene	9.20	77	79686	39.329	ng	96
25) Isophorone	9.73	82	152956	38.736	ng	98
26) 2-Nitrophenol	9.93	139	51283	40.390	ng	97
27) 2,4-Dimethylphenol	9.99	122	69033	39.470	ng	99
28) bis(2-Chloroethoxy)methane	10.22	93	95769	38.481	ng	98
29) 2,4-Dichlorophenol	10.47	162	94096	41.120	ng	99
30) 1,2,4-Trichlorobenzene	10.68	180	117011	41.659	ng	99
31) Naphthalene	10.87	128	257986	40.189	ng	99
32) Benzoic acid	10.14	122	38609	33.409	ng	97
33) 4-Chloroaniline	10.98	127	116332	39.257	ng	98
34) Hexachlorobutadiene	11.16	225	82327	43.613	ng	99
35) Caprolactam	11.75	113	28055	36.742	ng	96
36) 4-Chloro-3-methylphenol	12.11	107	85025	39.141	ng	99
37) 2-Methylnaphthalene	12.48	142	210499	40.839	ng	99
38) 1-Methylnaphthalene	12.70	142	198560	40.556	ng	97
40) 1,2,4,5-Tetrachlorobenzene	12.85	216	142815	42.104	ng	99

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41) Hexachlorocyclopentadiene	12.83	237	61204	34.350	nq	98
43) 2,4,6-Trichlorophenol	13.09	196	92009	40.130	nq	98
44) 2,4,5-Trichlorophenol	13.17	196	91164	38.370	nq	97
46) 1,1'-Biphenyl	13.49	154	272482	39.909	nq	98
47) 2-Chloronaphthalene	13.53	162	236048	40.095	nq	99
48) 2-Nitroaniline	13.73	65	52947	37.296	nq	98
49) Acenaphthylene	14.38	152	351369	39.671	nq	100
50) Dimethylphthalate	14.10	163	303765	39.858	nq	99
51) 2,6-Dinitrotoluene	14.23	165	68870	38.986	nq	91
52) Acenaphthene	14.72	154	209923	39.032	nq	98
53) 3-Nitroaniline	14.56	138	65886	37.293	nq	98
54) 2,4-Dinitrophenol	14.78	184	37391	35.204	nq	97
55) Dibenzofuran	15.06	168	359273	40.357	nq	99
56) 4-Nitrophenol	14.89	139	44908	35.773	nq	93
57) 2,4-Dinitrotoluene	15.02	165	99172	39.204	nq	97
58) Fluorene	15.71	166	291710	40.085	nq	98
59) 2,3,4,6-Tetrachlorophenol	15.29	232	95547m	40.665	nq	
60) Diethylphthalate	15.47	149	295283	39.635	nq	100
61) 4-Chlorophenyl-phenylether	15.70	204	179836	40.995	nq	96
62) 4-Nitroaniline	15.73	138	72704	38.451	nq	96
63) Azobenzene	15.99	77	187384	36.706	nq	95
65) 4,6-Dinitro-2-methylphenol	15.80	198	60308	39.987	nq	96
66) n-Nitrosodiphenylamine	15.91	169	267089	40.373	nq	98
67) 4-Bromophenyl-phenylether	16.59	248	125831	41.238	nq	98
68) Hexachlorobenzene	16.72	284	137429	41.432	nq	98
69) Atrazine	16.86	200	84886	36.285	nq	97
70) Pentachlorophenol	17.07	266	63526	36.860	nq	96
71) Phenanthrene	17.45	178	496415	41.545	nq	99
72) Anthracene	17.54	178	496888	41.655	nq	99
73) Carbazole	17.81	167	431507	40.589	nq	99
74) Di-n-butylphthalate	18.36	149	518940	41.294	nq	99
75) Fluoranthene	19.46	202	642510	44.060	nq	97
77) Benzidine	19.64	184	270539	37.730	nq	98
78) Pyrene	19.83	202	643508	41.968	nq	97
80) Butylbenzylphthalate	20.70	149	236110	39.150	nq	98
81) Benzo(a)anthracene	21.67	228	635772	41.789	nq	98
82) 3,3'-Dichlorobenzidine	21.58	252	241187	39.418	nq	# 99
83) Chrysene	21.73	228	605464	41.266	nq	99
84) Bis(2-ethylhexyl)phthalate	21.57	149	333096	39.780	nq	98
85) Di-n-octyl phthalate	22.78	149	567389	39.397	nq	99
86) Indeno(1,2,3-cd)pyrene	28.62	276	792501	39.746	nq	# 97
88) Benzo(b)fluoranthene	23.90	252	688150m	41.383	nq	
89) Benzo(k)fluoranthene	23.96	252	664307	41.561	nq	# 98
90) Benzo(a)pyrene	24.77	252	650963	41.254	nq	98
91) Dibenzo(a,h)anthracene	28.69	278	644312	40.328	nq	98
92) Benzo(g,h,i)perylene	29.79	276	642864	40.231	nq	96

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

