

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG021319\
 Data File : BG039490.D
 Acq On : 13 Feb 2019 15:55
 Operator : JU/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040

Quant Time: Feb 14 02:46:40 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG012919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jan 29 15:41:51 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.17	152	52329	20.00	ng	-0.01
21) Naphthalene-d8	11.00	136	234532	20.00	ng	-0.01
39) Acenaphthene-d10	14.80	164	162889	20.00	ng	-0.01
64) Phenanthrene-d10	17.54	188	391238	20.00	ng	-0.01
76) Chrysene-d12	21.83	240	388093	20.00	ng	-0.01
87) Perylene-d12	25.21	264	399597	20.00	ng	-0.03

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.71	112	249978	81.76	ng	0.00
7) Phenol-d6	7.31	99	381437	84.75	ng	-0.01
23) Nitrobenzene-d5	9.35	82	354957	94.55	ng	-0.01
42) 2,4,6-Tribromophenol	16.28	330	156680	89.33	ng	0.00
45) 2-Fluorobiphenyl	13.42	172	879736	77.48	ng	-0.01
79) Terphenyl-d14	20.13	244	1389719	75.80	ng	-0.02

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	3.61	88	53542	40.034	ng	93
3) Pyridine	4.01	79	149163	42.125	ng	94
4) n-Nitrosodimethylamine	3.93	42	66044	37.295	ng	89
6) Aniline	7.49	93	228588	40.549	ng	99
8) 2-Chlorophenol	7.73	128	142265	42.567	ng	98
9) Benzaldehyde	7.31	77	94046	42.626	ng	93
10) Phenol	7.34	94	195542	41.681	ng	97
11) bis(2-Chloroethyl)ether	7.59	93	148070	39.942	ng	95
12) 1,3-Dichlorobenzene	8.06	146	161067	39.782	ng	97
13) 1,4-Dichlorobenzene	8.21	146	163602	40.541	ng	99
14) 1,2-Dichlorobenzene	8.53	146	159104	40.727	ng	97
15) Benzyl Alcohol	8.41	79	146782	41.543	ng	100
16) 2,2'-oxybis(1-Chloropropan	8.69	45	288938	34.514	ng	99
17) 2-Methylphenol	8.60	107	123854	40.796	ng	96
18) Hexachloroethane	9.26	117	55720	40.134	ng	99
19) n-Nitroso-di-n-propylamine	8.98	70	126856	39.713	ng	96
20) 3+4-Methylphenols	8.93	107	181904	43.510	ng	98
22) Acetophenone	9.00	105	251110	39.859	ng	95
24) Nitrobenzene	9.39	77	182697	45.046	ng	99
25) Isophorone	9.90	82	318908	38.333	ng	98
26) 2-Nitrophenol	10.10	139	90687	54.625	ng	94
27) 2,4-Dimethylphenol	10.14	122	137627	40.895	ng	95
28) bis(2-Chloroethoxy)methane	10.39	93	207141	40.424	ng	98
29) 2,4-Dichlorophenol	10.63	162	164508	44.726	ng	93
30) 1,2,4-Trichlorobenzene	10.85	180	171477	41.526	ng	100
31) Naphthalene	11.05	128	462227	39.727	ng	99
32) Benzoic acid	10.24	122	72986	35.916	ng	96
33) 4-Chloroaniline	11.15	127	223014	41.339	ng	98
34) Hexachlorobutadiene	11.31	225	114776	41.795	ng	94
35) Caprolactam	11.95	113	67267	44.195	ng	# 85
36) 4-Chloro-3-methylphenol	12.25	107	185492	43.607	ng	98
37) 2-Methylnaphthalene	12.64	142	342275	39.719	ng	97
38) 1-Methylnaphthalene	12.85	142	331592	39.918	ng	98
40) 1,2,4,5-Tetrachlorobenzene	12.99	216	215144	41.512	ng	99

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41) Hexachlorocyclopentadiene	12.96	237	103633	36.696	nq	98
43) 2,4,6-Trichlorophenol	13.23	196	140702	44.155	nq	99
44) 2,4,5-Trichlorophenol	13.30	196	154166	46.160	nq	99
46) 1,1'-Biphenyl	13.63	154	519051	38.298	nq	100
47) 2-Chloronaphthalene	13.68	162	404407	39.665	nq	99
48) 2-Nitroaniline	13.89	65	137489	47.127	nq	93
49) Acenaphthylene	14.52	152	602631	39.172	nq	99
50) Dimethylphthalate	14.24	163	519729	39.506	nq	100
51) 2,6-Dinitrotoluene	14.37	165	115155	46.995	nq	94
52) Acenaphthene	14.86	154	393661	39.421	nq	99
53) 3-Nitroaniline	14.71	138	123211	46.630	nq	# 97
54) 2,4-Dinitrophenol	14.91	184	22648	30.433	nq	90
55) Dibenzofuran	15.19	168	634118	39.605	nq	99
56) 4-Nitrophenol	14.99	139	90764	41.188	nq	87
57) 2,4-Dinitrotoluene	15.16	165	161543	50.755	nq	# 89
58) Fluorene	15.84	166	462696	38.766	nq	100
59) 2,3,4,6-Tetrachlorophenol	15.41	232	137214	43.879	nq	97
60) Diethylphthalate	15.60	149	518017	38.084	nq	98
61) 4-Chlorophenyl-phenylether	15.82	204	274708	40.647	nq	97
62) 4-Nitroaniline	15.87	138	125187	45.609	nq	96
63) Azobenzene	16.12	77	484765	36.463	nq	98
65) 4,6-Dinitro-2-methylphenol	15.91	198	59146	40.917	nq	92
66) n-Nitrosodiphenylamine	16.04	169	464423	39.040	nq	97
67) 4-Bromophenyl-phenylether	16.72	248	179706	41.404	nq	94
68) Hexachlorobenzene	16.84	284	183270	42.086	nq	95
69) Atrazine	16.98	200	160112	36.830	nq	97
70) Pentachlorophenol	17.18	266	105082	34.720	nq	99
71) Phenanthrene	17.58	178	778002	38.421	nq	100
72) Anthracene	17.68	178	769187	38.564	nq	99
73) Carbazole	17.95	167	746543	37.505	nq	99
74) Di-n-butylphthalate	18.48	149	875135	37.685	nq	99
75) Fluoranthene	19.58	202	902953	38.418	nq	99
77) Benzidine	19.76	184	369805	29.064	nq	98
78) Pyrene	19.95	202	941142	38.955	nq	98
80) Butylbenzylphthalate	20.81	149	398995	39.150	nq	97
81) Benzo(a)anthracene	21.81	228	896413	39.090	nq	99
82) 3,3'-Dichlorobenzidine	21.72	252	344289	40.490	nq	98
83) Chrysene	21.88	228	861248	39.453	nq	99
84) Bis(2-ethylhexyl)phthalate	21.68	149	560822	38.572	nq	99
85) Di-n-octyl phthalate	22.94	149	948720	39.496	nq	96
86) Indeno(1,2,3-cd)pyrene	29.11	276	975150	41.634	nq	99
88) Benzo(b)fluoranthene	24.13	252	866433	39.759	nq	100
89) Benzo(k)fluoranthene	24.20	252	846798	38.704	nq	98
90) Benzo(a)pyrene	25.05	252	844095	40.219	nq	98
91) Dibenzo(a,h)anthracene	29.18	278	808955	41.158	nq	99
92) Benzo(g,h,i)perylene	30.33	276	792918	40.475	nq	97

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

