

Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG030819\
 Data File : BG039817.D
 Acq On : 8 Mar 2019 12:23
 Operator : JU/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040

Quant Time: Mar 08 13:57:31 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG030419.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 04 14:38:15 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.15	152	40823	20.00	ng	-0.02
21) Naphthalene-d8	10.97	136	196749	20.00	ng	-0.02
39) Acenaphthene-d10	14.78	164	132727	20.00	ng	-0.01
64) Phenanthrene-d10	17.52	188	316537	20.00	ng	-0.01
76) Chrysene-d12	21.81	240	307201	20.00	ng	-0.02
87) Perylene-d12	25.18	264	316384	20.00	ng	-0.02

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.69	112	194879	81.44	ng	-0.02
7) Phenol-d6	7.30	99	300366	84.18	ng	-0.02
23) Nitrobenzene-d5	9.33	82	293586	80.70	ng	-0.01
42) 2,4,6-Tribromophenol	16.27	330	123983	80.14	ng	-0.01
45) 2-Fluorobiphenyl	13.41	172	713329	76.52	ng	-0.01
79) Terphenyl-d14	20.11	244	1083522	77.66	ng	-0.01

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	3.59	88	41394	37.470	ng	95
3) Pyridine	3.99	79	122663	41.474	ng	94
4) n-Nitrosodimethylamine	3.91	42	57862	41.240	ng	# 78
6) Aniline	7.48	93	185720	39.730	ng	96
8) 2-Chlorophenol	7.71	128	118197	40.715	ng	97
9) Benzaldehyde	7.30	77	92335	40.119	ng	98
10) Phenol	7.33	94	162018	41.917	ng	99
11) bis(2-Chloroethyl)ether	7.57	93	122083	40.511	ng	95
12) 1,3-Dichlorobenzene	8.04	146	127848	38.939	ng	98
13) 1,4-Dichlorobenzene	8.19	146	132487	39.616	ng	99
14) 1,2-Dichlorobenzene	8.51	146	125582	38.865	ng	99
15) Benzyl Alcohol	8.39	79	117239	41.423	ng	98
16) 2,2'-oxybis(1-Chloropropan	8.67	45	237675	39.433	ng	98
17) 2-Methylphenol	8.59	107	102730	41.682	ng	96
18) Hexachloroethane	9.24	117	47616	39.681	ng	100
19) n-Nitroso-di-n-propylamine	8.96	70	102299	39.834	ng	94
20) 3+4-Methylphenols	8.91	107	145369	42.457	ng	95
22) Acetophenone	8.98	105	200444	37.958	ng	# 94
24) Nitrobenzene	9.38	77	149887	39.275	ng	96
25) Isophorone	9.89	82	292040	38.313	ng	99
26) 2-Nitrophenol	10.08	139	75607	41.033	ng	97
27) 2,4-Dimethylphenol	10.12	122	116188	40.544	ng	98
28) bis(2-Chloroethoxy)methane	10.37	93	173925	37.547	ng	96
29) 2,4-Dichlorophenol	10.62	162	132539	41.078	ng	97
30) 1,2,4-Trichlorobenzene	10.83	180	138430	38.128	ng	97
31) Naphthalene	11.03	128	402323	38.622	ng	98
32) Benzoic acid	10.25	122	69307	36.980	ng	98
33) 4-Chloroaniline	11.14	127	175796	39.798	ng	96
34) Hexachlorobutadiene	11.29	225	92700	37.364	ng	91
35) Caprolactam	11.92	113	47618	38.635	ng	96
36) 4-Chloro-3-methylphenol	12.24	107	147701	39.934	ng	96
37) 2-Methylnaphthalene	12.62	142	303446	38.963	ng	99
38) 1-Methylnaphthalene	12.84	142	294019	38.848	ng	100
40) 1,2,4,5-Tetrachlorobenzene	12.98	216	172324	38.324	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.95	237	81077	33.647	ng	100
43) 2,4,6-Trichlorophenol	13.22	196	108928	41.379	ng	96
44) 2,4,5-Trichlorophenol	13.29	196	119222	40.179	ng	97
46) 1,1'-Biphenyl	13.62	154	424051	38.845	ng	95
47) 2-Chloronaphthalene	13.66	162	316375	38.524	ng	98
48) 2-Nitroaniline	13.87	65	110549	41.887	ng	98
49) Acenaphthylene	14.51	152	525212	38.958	ng	99
50) Dimethylphthalate	14.23	163	425601	38.348	ng	99
51) 2,6-Dinitrotoluene	14.36	165	97247	41.332	ng	96
52) Acenaphthene	14.84	154	314113	38.711	ng	99
53) 3-Nitroaniline	14.69	138	92971	39.310	ng	# 97
54) 2,4-Dinitrophenol	14.89	184	46266	34.389	ng	96
55) Dibenzofuran	15.17	168	514989	38.683	ng	99
56) 4-Nitrophenol	14.98	139	77993	36.863	ng	86
57) 2,4-Dinitrotoluene	15.14	165	137808	41.684	ng	# 86
58) Fluorene	15.83	166	410243	39.198	ng	100
59) 2,3,4,6-Tetrachlorophenol	15.40	232	117600	40.581	ng	98
60) Diethylphthalate	15.58	149	426538	38.823	ng	98
61) 4-Chlorophenyl-phenylether	15.81	204	216418	38.751	ng	98
62) 4-Nitroaniline	15.86	138	101145	39.448	ng	96
63) Azobenzene	16.10	77	391358	39.296	ng	98
65) 4,6-Dinitro-2-methylphenol	15.90	198	64268	34.339	ng	98
66) n-Nitrosodiphenylamine	16.03	169	364652	39.793	ng	99
67) 4-Bromophenyl-phenylether	16.71	248	138918	39.208	ng	95
68) Hexachlorobenzene	16.82	284	144844	39.438	ng	96
69) Atrazine	16.97	200	141310	39.182	ng	98
70) Pentachlorophenol	17.17	266	83492	35.996	ng	95
71) Phenanthrene	17.57	178	674624	38.693	ng	100
72) Anthracene	17.66	178	667471	39.285	ng	98
73) Carbazole	17.93	167	597961	39.039	ng	97
74) Di-n-butylphthalate	18.46	149	715569	39.706	ng	99
75) Fluoranthene	19.57	202	782902	37.995	ng	99
77) Benzidine	19.75	184	283599	29.092	ng	97
78) Pyrene	19.93	202	786920	39.421	ng	99
80) Butylbenzylphthalate	20.80	149	320578	41.451	ng	95
81) Benzo(a)anthracene	21.79	228	756168	39.275	ng	99
82) 3,3'-Dichlorobenzidine	21.71	252	269683	38.366	ng	99
83) Chrysene	21.86	228	726627	38.929	ng	100
84) Bis(2-ethylhexyl)phthalate	21.65	149	447089	41.138	ng	99
85) Di-n-octyl phthalate	22.91	149	763909	42.452	ng	95
86) Indeno(1,2,3-cd)pyrene	29.05	276	826204	39.018	ng	98
88) Benzo(b)fluoranthene	24.10	252	743439	39.405	ng	99
89) Benzo(k)fluoranthene	24.17	252	692438	39.428	ng	99
90) Benzo(a)pyrene	25.03	252	698311	40.055	ng	98
91) Dibenzo(a,h)anthracene	29.13	278	668030	39.086	ng	99
92) Benzo(g,h,i)perylene	30.29	276	655704	38.199	ng	98

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

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