

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG010919\
 Data File : BG039015.D
 Acq On : 9 Jan 2019 11:00
 Operator : JU/SJ
 Sample : SSTDICC025
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDICC025

Quant Time: Jan 09 12:56:46 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG010919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jan 09 12:51:00 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.03	152	20963	20.00	ng	0.00
21) Naphthalene-d8	10.85	136	89014	20.00	ng	0.00
39) Acenaphthene-d10	14.67	164	66662	20.00	ng	0.00
64) Phenanthrene-d10	17.42	188	171462	20.00	ng	0.00
76) Chrysene-d12	21.70	240	176729	20.00	ng	0.00
87) Perylene-d12	24.98	264	194971	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.58	112	57343	48.52	ng	0.00
7) Phenol-d6	7.19	99	82741	50.99	ng	0.00
23) Nitrobenzene-d5	9.21	82	81918	47.01	ng	0.00
42) 2,4,6-Tribromophenol	16.16	330	55839	60.78	ng	0.00
45) 2-Fluorobiphenyl	13.29	172	240003	49.39	ng	0.00
79) Terphenyl-d14	20.02	244	495278	60.99	ng	0.00

Target Compounds				Qvalue		
2) 1,4-Dioxane	3.46	88	10914	19.841	ng	99
3) Pyridine	3.87	79	31024	20.485	ng	94
4) n-Nitrosodimethylamine	3.79	42	13091	17.641	ng	96
6) Aniline	7.36	93	50791	24.522	ng	99
8) 2-Chlorophenol	7.59	128	33348	24.382	ng	93
9) Benzaldehyde	7.17	77	24955	23.709	ng	95
10) Phenol	7.22	94	41798	24.248	ng	99
11) bis(2-Chloroethyl)ether	7.45	93	31937	23.271	ng	97
12) 1,3-Dichlorobenzene	7.92	146	40210	24.864	ng	96
13) 1,4-Dichlorobenzene	8.07	146	40618	24.524	ng	97
14) 1,2-Dichlorobenzene	8.39	146	39508	25.302	ng	91
15) Benzyl Alcohol	8.27	79	32227	25.447	ng	99
16) 2,2'-oxybis(1-Chloropropan	8.56	45	61748	19.641	ng	97
17) 2-Methylphenol	8.47	107	29191	25.276	ng	100
18) Hexachloroethane	9.11	117	13621	22.506	ng	93
19) n-Nitroso-di-n-propylamine	8.83	70	28621	25.115	ng	# 94
20) 3+4-Methylphenols	8.80	107	42005	26.336	ng	97
22) Acetophenone	8.86	105	59542	26.780	ng	# 98
24) Nitrobenzene	9.25	77	42025	23.842	ng	99
25) Isophorone	9.77	82	71769	22.925	ng	99
26) 2-Nitrophenol	9.96	139	22569	25.016	ng	93
27) 2,4-Dimethylphenol	10.01	122	32825	25.388	ng	93
28) bis(2-Chloroethoxy)methane	10.25	93	44086	24.954	ng	99
29) 2,4-Dichlorophenol	10.49	162	39996	27.806	ng	96
30) 1,2,4-Trichlorobenzene	10.71	180	47592	27.392	ng	94
31) Naphthalene	10.90	128	112310	25.608	ng	99
32) Benzoic acid	10.15	122	15086	21.976	ng	88
33) 4-Chloroaniline	11.02	127	53165	27.921	ng	99
34) Hexachlorobutadiene	11.16	225	32748	27.856	ng	96
35) Caprolactam	11.79	113	16890	28.374	ng	99
36) 4-Chloro-3-methylphenol	12.13	107	42508	25.897	ng	95
37) 2-Methylnaphthalene	12.50	142	84426	26.983	ng	99
38) 1-Methylnaphthalene	12.72	142	82092	27.038	ng	99
40) 1,2,4,5-Tetrachlorobenzene	12.86	216	62243	24.979	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.83	237	28601	20.892	nq	95
43) 2,4,6-Trichlorophenol	13.11	196	39581	25.834	nq	98
44) 2,4,5-Trichlorophenol	13.19	196	42415	26.147	nq	97
46) 1,1'-Biphenyl	13.50	154	129842	23.234	nq	98
47) 2-Chloronaphthalene	13.56	162	103941	24.078	nq	99
48) 2-Nitroaniline	13.77	65	31075	22.600	nq	95
49) Acenaphthylene	14.40	152	161565	25.036	nq	98
50) Dimethylphthalate	14.12	163	142961	26.261	nq	99
51) 2,6-Dinitrotoluene	14.25	165	32003	25.941	nq	97
52) Acenaphthene	14.74	154	95443	24.733	nq	97
53) 3-Nitroaniline	14.59	138	32554	25.886	nq	95
54) 2,4-Dinitrophenol	14.80	184	15192	23.583	nq	96
55) Dibenzofuran	15.07	168	170502	25.776	nq	97
56) 4-Nitrophenol	14.90	139	22339	23.415	nq	99
57) 2,4-Dinitrotoluene	15.04	165	46034	26.493	nq	# 96
58) Fluorene	15.72	166	128283	26.176	nq	98
59) 2,3,4,6-Tetrachlorophenol	15.29	232	39286	26.918	nq	# 99
60) Diethylphthalate	15.48	149	146497	24.514	nq	99
61) 4-Chlorophenyl-phenylether	15.71	204	81171	26.546	nq	99
62) 4-Nitroaniline	15.76	138	35425	27.362	nq	95
63) Azobenzene	16.00	77	115030	23.041	nq	98
65) 4,6-Dinitro-2-methylphenol	15.81	198	31023	25.364	nq	92
66) n-Nitrosodiphenylamine	15.92	169	128641	25.534	nq	98
67) 4-Bromophenyl-phenylether	16.60	248	55296	26.406	nq	94
68) Hexachlorobenzene	16.72	284	59788	27.543	nq	97
69) Atrazine	16.87	200	55542	29.707	nq	94
70) Pentachlorophenol	17.07	266	29494	22.971	nq	95
71) Phenanthrene	17.46	178	229161	25.773	nq	98
72) Anthracene	17.56	178	227835	25.956	nq	98
73) Carbazole	17.83	167	227253	26.178	nq	99
74) Di-n-butylphthalate	18.36	149	248548	24.952	nq	98
75) Fluoranthene	19.47	202	286503	26.043	nq	99
77) Benzidine	19.66	184	141769	27.125	nq	99
78) Pyrene	19.84	202	293357	25.126	nq	98
80) Butylbenzylphthalate	20.71	149	107295	23.523	nq	99
81) Benzo(a)anthracene	21.68	228	275169	25.552	nq	97
82) 3,3'-Dichlorobenzidine	21.60	252	113935	26.676	nq	98
83) Chrysene	21.75	228	260370	25.102	nq	99
84) Bis(2-ethylhexyl)phthalate	21.55	149	151522	23.422	nq	95
85) Di-n-octyl phthalate	22.77	149	257009	23.722	nq	100
86) Indeno(1,2,3-cd)pyrene	28.74	276	309389	25.371	nq	# 96
88) Benzo(b)fluoranthene	23.93	252	270348	25.255	nq	99
89) Benzo(k)fluoranthene	24.00	252	263915	24.758	nq	99
90) Benzo(a)pyrene	24.82	252	261638	25.335	nq	99
91) Dibenzo(a,h)anthracene	28.80	278	257622	25.054	nq	96
92) Benzo(g,h,i)perylene	29.92	276	258928	25.552	nq	97

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

