

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG082318\
 Data File : BG036355.D
 Acq On : 22 Aug 2018 23:10
 Operator : JU/SJ
 Sample : SSTDICC025
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDICC025

Quant Time: Aug 23 01:16:48 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG082318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 23 01:15:05 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.07	152	48898	20.00	ng	0.00
21) Naphthalene-d8	10.88	136	209504	20.00	ng	0.00
38) Acenaphthene-d10	14.70	164	132295	20.00	ng	0.00
63) Phenanthrene-d10	17.44	188	335813	20.00	ng	0.00
75) Chrysene-d12	21.71	240	356777	20.00	ng	0.00
86) Perylene-d12	24.94	264	377829	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.63	112	149821	50.57	ng	0.00
7) Phenol-d6	7.22	99	216293	49.25	ng	0.00
23) Nitrobenzene-d5	9.23	82	189867	38.96	ng	0.00
41) 2,4,6-Tribromophenol	16.18	330	79082	45.35	ng	0.00
44) 2-Fluorobiphenyl	13.33	172	467112	47.39	ng	0.00
78) Terphenyl-d14	20.05	244	794173	49.22	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.55	88	33535	22.397	ng	99
3) Pyridine	3.94	79	99508	25.367	ng	99
4) n-Nitrosodimethylamine	3.85	42	43818	25.766	ng	98
6) Aniline	7.39	93	135981	23.945	ng	95
8) 2-Chlorophenol	7.63	128	80350	26.395	ng	97
9) Benzaldehyde	7.21	77	75958	28.004	ng	97
10) Phenol	7.24	94	110984	24.908	ng	97
11) bis(2-Chloroethyl)ether	7.49	93	88231	25.233	ng	97
12) 1,3-Dichlorobenzene	7.96	146	91428	25.103	ng	99
13) 1,4-Dichlorobenzene	8.11	146	93342	25.252	ng	99
14) 1,2-Dichlorobenzene	8.43	146	86217	24.301	ng	96
15) Benzyl Alcohol	8.30	79	85947	21.856	ng	99
16) 2,2'-oxybis(1-Chloropropan	8.60	45	178930	50.590	ng	99
17) 2-Methylphenol	8.50	107	72850	24.196	ng	96
18) Hexachloroethane	9.16	117	32730	23.254	ng	99
19) n-Nitroso-di-n-propylamine	8.88	70	78526	21.537	ng	96
20) 3+4-Methylphenols	8.83	107	101949	24.455	ng	99
22) Acetophenone	8.90	105	136224	21.274	ng	# 97
24) Nitrobenzene	9.28	77	103376	20.847	ng	99
25) Isophorone	9.79	82	189895	20.844	ng	98
26) 2-Nitrophenol	9.99	139	37474	21.173	ng	99
27) 2,4-Dimethylphenol	10.04	122	72751	23.966	ng	98
28) bis(2-Chloroethoxy)methane	10.28	93	113223	22.269	ng	97
29) 2,4-Dichlorophenol	10.52	162	82637	23.985	ng	94
30) 1,2,4-Trichlorobenzene	10.74	180	95555	22.717	ng	98
31) Naphthalene	10.93	128	260149	23.800	ng	98
32) Benzoic acid	10.14	122	45022	21.802	ng	99
33) 4-Chloroaniline	11.03	127	118096	24.631	ng	96
34) Hexachlorobutadiene	11.22	225	65248	20.180	ng	97
35) Caprolactam	11.79	113	34008	22.939	ng	96
36) 4-Chloro-3-methylphenol	12.14	107	95039	20.905	ng	99
37) 2-Methylnaphthalene	12.53	142	192937	23.819	ng	99
39) 1,2,4,5-Tetrachlorobenzene	12.89	216	115508	23.121	ng	98
40) Hexachlorocyclopentadiene	12.88	237	58488	22.440	ng	93

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42) 2,4,6-Trichlorophenol	13.13	196	71119	24.382	ng	97
43) 2,4,5-Trichlorophenol	13.20	196	77282	23.555	ng	98
45) 1,1'-Biphenyl	13.54	154	274205	26.369	ng	97
46) 2-Chloronaphthalene	13.58	162	209195	25.814	ng	99
47) 2-Nitroaniline	13.77	65	66373	23.470	ng	96
48) Acenaphthylene	14.42	152	328761	24.496	ng	99
49) Dimethylphthalate	14.15	163	272966	23.530	ng	97
50) 2,6-Dinitrotoluene	14.27	165	54794	23.251	ng	97
51) Acenaphthene	14.76	154	205753	24.456	ng	99
52) 3-Nitroaniline	14.59	138	61686	25.249	ng	98
53) 2,4-Dinitrophenol	14.79	184	18401	18.062	ng	# 77
54) Dibenzofuran	15.09	168	330843	24.849	ng	97
55) 4-Nitrophenol	14.88	139	48178	27.016	ng	92
56) 2,4-Dinitrotoluene	15.05	165	70126	21.077	ng	99
57) Fluorene	15.75	166	249444	22.603	ng	99
58) 2,3,4,6-Tetrachlorophenol	15.31	232	73214	23.452	ng	97
59) Diethylphthalate	15.52	149	279060	23.476	ng	98
60) 4-Chlorophenyl-phenylether	15.74	204	151770	23.309	ng	98
61) 4-Nitroaniline	15.75	138	66621	26.110	ng	94
62) Azobenzene	16.03	77	282153	23.993	ng	98
64) 4,6-Dinitro-2-methylphenol	15.81	198	32650	17.990	ng	93
65) n-Nitrosodiphenylamine	15.95	169	245333	25.495	ng	96
66) 4-Bromophenyl-phenylether	16.63	248	94376	24.152	ng	93
67) Hexachlorobenzene	16.75	284	100222	24.862	ng	95
68) Atrazine	16.90	200	99681	25.610	ng	98
69) Pentachlorophenol	17.09	266	70390	27.897	ng	96
70) Phenanthrene	17.48	178	433108	25.652	ng	98
71) Anthracene	17.57	178	430154	25.589	ng	99
72) Carbazole	17.84	167	438351	27.230	ng	98
73) Di-n-butylphthalate	18.41	149	498022	26.375	ng	100
74) Fluoranthene	19.49	202	544511	24.236	ng	99
76) Benzidine	19.67	184	272835	28.222	ng	98
77) Pyrene	19.85	202	553043	24.402	ng	99
79) Butylbenzylphthalate	20.74	149	220168	26.783	ng	98
80) Benzo(a)anthracene	21.69	228	542752	24.392	ng	98
81) 3,3'-Dichlorobenzidine	21.60	252	220822	27.856	ng	99
82) Chrysene	21.76	228	514116	24.808	ng	99
83) Bis(2-ethylhexyl)phthalate	21.62	149	312313	27.885	ng	98
84) Di-n-octyl phthalate	22.85	149	526792	28.056	ng	99
85) Indeno(1,2,3-cd)pyrene	28.60	276	591990	25.669	ng	99
87) Benzo(b)fluoranthene	23.91	252	519184	22.689	ng	97
88) Benzo(k)fluoranthene	23.98	252	510132	22.933	ng	100
89) Benzo(a)pyrene	24.78	252	501223	23.564	ng	100
90) Dibenzo(a,h)anthracene	28.68	278	482089	23.143	ng	97
91) Benzo(g,h,i)perylene	29.75	276	481794	23.561	ng	98

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

