

Data Path : Z:\HPCHEM1\BNA G\DATA\BG041018\
 Data File : BG033710.D
 Acq On : 10 Apr 2018 12:58
 Operator : SJ/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Sohil
 4/11/2018 5:18:15 PM

Quant Time: Apr 10 17:03:40 2018
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG031918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Apr 10 16:51:36 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.86	152	33841	20.00	ng	0.00
21) Naphthalene-d8	11.74	136	139354	20.00	ng	0.00
38) Acenaphthene-d10	15.47	164	103443	20.00	ng	0.00
63) Phenanthrene-d10	18.21	188	283129	20.00	ng	0.00
75) Chrysene-d12	22.55	240	292530	20.00	ng	0.00
86) Perylene-d12	26.31	264	319019	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	6.28	112	151450	83.92	ng	0.00
7) Phenol-d6	7.97	99	234239	75.54	ng	0.00
23) Nitrobenzene-d5	10.06	82	242865	80.07	ng	0.00
41) 2,4,6-Tribromophenol	16.95	330	116955	75.60	ng	0.00
44) 2-Fluorobiphenyl	14.10	172	555173	77.48	ng	0.00
78) Terphenyl-d14	20.73	244	1028809	75.21	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.92	88	39275	42.853	ng	88
3) Pyridine	4.39	79	107275	42.341	ng	98
4) n-Nitrosodimethylamine	4.30	42	45134	43.409	ng	# 87
6) Aniline	8.15	93	145465	39.890	ng	# 96
8) 2-Chlorophenol	8.40	128	81037	39.277	ng	98
9) Benzaldehyde	7.96	77	64284m	32.621	ng	
10) Phenol	8.00	94	116196	37.953	ng	93
11) bis(2-Chloroethyl)ether	8.23	93	87481	35.007	ng	89
12) 1,3-Dichlorobenzene	8.74	146	105269	39.026	ng	92
13) 1,4-Dichlorobenzene	8.89	146	100522	37.211	ng	99
14) 1,2-Dichlorobenzene	9.22	146	101278	38.413	ng	96
15) Benzyl Alcohol	9.09	79	96057	39.344	ng	90
16) 2,2'-oxybis(1-Chloropropan	9.37	45	168693	41.409	ng	88
17) 2-Methylphenol	9.29	107	77248m	37.038	ng	
18) Hexachloroethane	9.97	117	40430	38.777	ng	98
19) n-Nitroso-di-n-propylamine	9.66	70	90612	39.878	ng	99
20) 3+4-Methylphenols	9.63	107	110865	37.334	ng	89
22) Acetophenone	9.70	105	157359	41.217	ng	# 98
24) Nitrobenzene	10.10	77	127742	39.347	ng	92
25) Isophorone	10.61	82	245490	41.702	ng	99
26) 2-Nitrophenol	10.83	139	51060	41.542	ng	# 92
27) 2,4-Dimethylphenol	10.86	122	75701	42.396	ng	98
28) bis(2-Chloroethoxy)methane	11.10	93	115525	39.331	ng	99
29) 2,4-Dichlorophenol	11.37	162	89562	40.786	ng	94
30) 1,2,4-Trichlorobenzene	11.59	180	109221	38.283	ng	98
31) Naphthalene	11.79	128	286043	39.611	ng	98
32) Benzoic acid	11.00	122	62711m	37.781	ng	
33) 4-Chloroaniline	11.89	127	123533	38.182	ng	# 89
34) Hexachlorobutadiene	12.05	225	82391	38.188	ng	90
35) Caprolactam	12.62	113	40373	46.931	ng	97
36) 4-Chloro-3-methylphenol	12.95	107	110486	38.578	ng	97
37) 2-Methylnaphthalene	13.34	142	210412	37.540	ng	90
39) 1,2,4,5-Tetrachlorobenzene	13.69	216	139635	37.266	ng	99
40) Hexachlorocyclopentadiene	13.67	237	72273	32.903	ng	92

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.92	196	84073	38.619	nq	99
43) 2,4,5-Trichlorophenol	14.00	196	92067	35.779	nq	98
45) 1,1'-Biphenyl	14.32	154	316627	39.841	nq	98
46) 2-Chloronaphthalene	14.37	162	239008	39.004	nq	95
47) 2-Nitroaniline	14.56	65	99360	43.291	nq	88
48) Acenaphthylene	15.20	152	409552	40.041	nq	98
49) Dimethylphthalate	14.90	163	370482	41.154	nq	99
50) 2,6-Dinitrotoluene	15.03	165	76667	41.650	nq	96
51) Acenaphthene	15.54	154	264394	40.098	nq	96
52) 3-Nitroaniline	15.37	138	77778	40.303	nq	# 97
53) 2,4-Dinitrophenol	15.57	184	26803	33.074	nq	87
54) Dibenzofuran	15.87	168	380641	39.082	nq	94
55) 4-Nitrophenol	15.66	139	52562	38.734	nq	# 79
56) 2,4-Dinitrotoluene	15.81	165	111692	41.210	nq	95
57) Fluorene	16.51	166	332367	40.057	nq	96
58) 2,3,4,6-Tetrachlorophenol	16.08	232	95810	39.551	nq	95
59) Diethylphthalate	16.24	149	374952	42.893	nq	98
60) 4-Chlorophenyl-phenylether	16.48	204	182758	38.316	nq	99
61) 4-Nitroaniline	16.52	138	85596	43.800	nq	89
62) Azobenzene	16.78	77	347402	41.026	nq	97
64) 4,6-Dinitro-2-methylphenol	16.57	198	69185	40.847	nq	95
65) n-Nitrosodiphenylamine	16.70	169	307860	38.088	nq	97
66) 4-Bromophenyl-phenylether	17.38	248	120667	36.290	nq	93
67) Hexachlorobenzene	17.51	284	127998	36.475	nq	94
68) Atrazine	17.61	200	133329	40.707	nq	95
69) Pentachlorophenol	17.84	266	79596	33.047	nq	96
70) Phenanthrene	18.25	178	565005	38.317	nq	99
71) Anthracene	18.34	178	566827	37.966	nq	100
72) Carbazole	18.60	167	535050	40.442	nq	97
73) Di-n-butylphthalate	19.09	149	653950	42.466	nq	# 99
74) Fluoranthene	20.20	202	716858	39.286	nq	98
76) Benzidine	20.36	184	332149	41.869	nq	99
77) Pyrene	20.56	202	734409	37.643	nq	99
79) Butylbenzylphthalate	21.41	149	294617	40.921	nq	99
80) Benzo(a)anthracene	22.53	228	709778	38.105	nq	98
81) 3,3'-Dichlorobenzidine	22.42	252	265441	40.046	nq	# 97
82) Chrysene	22.61	228	689529	38.241	nq	97
83) Bis(2-ethylhexyl)phthalate	22.35	149	425174	42.840	nq	97
84) Di-n-octyl phthalate	23.74	149	681588	44.853	nq	99
85) Indeno(1,2,3-cd)pyrene	30.68	276	786314	40.334	nq	# 90
87) Benzo(b)fluoranthene	25.10	252	677323	36.749	nq	# 96
88) Benzo(k)fluoranthene	25.18	252	706061	37.877	nq	99
89) Benzo(a)pyrene	26.13	252	670424	37.877	nq	99
90) Dibenzo(a,h)anthracene	30.75	278	655431	39.311	nq	98
91) Benzo(g,h,i)perylene	32.05	276	637699	38.625	nq	98

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

