

Data Path : Z:\HPCHEM1\BNA G\DATA\BG032918\
 Data File : BG033443.D
 Acq On : 30 Mar 2018 7:29
 Operator : SJ/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Sohil
 3/30/2018 6:04:49 PM

Quant Time: Mar 30 08:27:40 2018
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG031918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 28 18:34:58 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.87	152	31942	20.00	ng	0.00
21) Naphthalene-d8	11.75	136	153018	20.00	ng	0.00
38) Acenaphthene-d10	15.49	164	124654	20.00	ng	0.00
63) Phenanthrene-d10	18.21	188	332289	20.00	ng	0.00
75) Chrysene-d12	22.57	240	337141	20.00	ng	0.00
86) Perylene-d12	26.33	264	370392	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	6.30	112	144976	85.11	ng	0.00
7) Phenol-d6	7.98	99	240780	82.26	ng	0.00
23) Nitrobenzene-d5	10.07	82	259773	78.00	ng	0.00
41) 2,4,6-Tribromophenol	16.96	330	151781	81.41	ng	0.00
44) 2-Fluorobiphenyl	14.11	172	646829	74.91	ng	0.00
78) Terphenyl-d14	20.73	244	1176575	74.63	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.93	88	33875	39.158	ng	87
3) Pyridine	4.39	79	100995	42.233	ng	97
4) n-Nitrosodimethylamine	4.30	42	40699	41.471	ng	# 75
6) Aniline	8.17	93	146005	42.419	ng	96
8) 2-Chlorophenol	8.42	128	83082	42.662	ng	96
9) Benzaldehyde	7.98	77	64187	34.508	ng	88
10) Phenol	8.01	94	119399	41.318	ng	90
11) bis(2-Chloroethyl)ether	8.25	93	91471	38.780	ng	93
12) 1,3-Dichlorobenzene	8.75	146	99797	39.197	ng	93
13) 1,4-Dichlorobenzene	8.90	146	104712	41.066	ng	90
14) 1,2-Dichlorobenzene	9.23	146	97314	39.103	ng	97
15) Benzyl Alcohol	9.11	79	99002	42.961	ng	94
16) 2,2'-oxybis(1-Chloropropan	9.39	45	157068	40.848	ng	86
17) 2-Methylphenol	9.31	107	80690	40.989	ng	# 93
18) Hexachloroethane	9.99	117	39710	40.351	ng	88
19) n-Nitroso-di-n-propylamine	9.68	70	93175	43.444	ng	96
20) 3+4-Methylphenols	9.64	107	120128	42.859	ng	93
22) Acetophenone	9.71	105	170888	40.764	ng	# 95
24) Nitrobenzene	10.12	77	136176	38.199	ng	90
25) Isophorone	10.63	82	268691	41.567	ng	98
26) 2-Nitrophenol	10.84	139	57799	42.825	ng	# 89
27) 2,4-Dimethylphenol	10.87	122	79691	40.645	ng	93
28) bis(2-Chloroethoxy)methane	11.11	93	123862	38.404	ng	95
29) 2,4-Dichlorophenol	11.38	162	96895	40.185	ng	96
30) 1,2,4-Trichlorobenzene	11.60	180	118476	37.819	ng	88
31) Naphthalene	11.80	128	307605	38.793	ng	99
32) Benzoic acid	11.06	122	27122m	14.881	ng	
33) 4-Chloroaniline	11.90	127	133878	37.685	ng	# 88
34) Hexachlorobutadiene	12.06	225	87891	37.100	ng	96
35) Caprolactam	12.64	113	43943	46.519	ng	98
36) 4-Chloro-3-methylphenol	12.97	107	125606	39.941	ng	95
37) 2-Methylnaphthalene	13.35	142	239974	38.991	ng	90
39) 1,2,4,5-Tetrachlorobenzene	13.71	216	166364	36.844	ng	98
40) Hexachlorocyclopentadiene	13.68	237	88167	33.309	ng	97

Data Path : Z:\HPCHEM1\BNA G\DATA\BG032918\
 Data File : BG033443.D
 Acq On : 30 Mar 2018 7:29
 Operator : SJ/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Sohil
 3/30/2018 6:04:49 PM

Quant Time: Mar 30 08:27:40 2018
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG031918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 28 18:34:58 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.93	196	103177	39.329	nq	99
43) 2,4,5-Trichlorophenol	14.01	196	125725	40.545	nq	96
45) 1,1'-Biphenyl	14.33	154	364532	38.064	nq	97
46) 2-Chloronaphthalene	14.38	162	280764	38.022	nq	94
47) 2-Nitroaniline	14.58	65	113369	40.989	nq	92
48) Acenaphthylene	15.22	152	479065	38.867	nq	98
49) Dimethylphthalate	14.91	163	438439	40.416	nq	100
50) 2,6-Dinitrotoluene	15.05	165	91940	41.449	nq	98
51) Acenaphthene	15.55	154	289176	36.394	nq	96
52) 3-Nitroaniline	15.39	138	94591	40.675	nq	# 97
53) 2,4-Dinitrophenol	15.60	184	5269m	11.792	nq	
54) Dibenzofuran	15.88	168	455455	38.806	nq	94
55) 4-Nitrophenol	15.67	139	58883	36.009	nq	# 84
56) 2,4-Dinitrotoluene	15.82	165	131588	40.290	nq	92
57) Fluorene	16.52	166	390464	39.051	nq	99
58) 2,3,4,6-Tetrachlorophenol	16.09	232	119284	40.862	nq	99
59) Diethylphthalate	16.25	149	432199	41.029	nq	96
60) 4-Chlorophenyl-phenylether	16.50	204	223546	38.893	nq	100
61) 4-Nitroaniline	16.54	138	98542	41.844	nq	# 81
62) Azobenzene	16.78	77	398792	39.081	nq	97
64) 4,6-Dinitro-2-methylphenol	16.58	198	35400	17.808	nq	92
65) n-Nitrosodiphenylamine	16.71	169	362980	38.263	nq	97
66) 4-Bromophenyl-phenylether	17.38	248	150424	38.546	nq	95
67) Hexachlorobenzene	17.51	284	157884	38.335	nq	96
68) Atrazine	17.62	200	153172	39.846	nq	98
69) Pentachlorophenol	17.85	266	101371m	35.861	nq	
70) Phenanthrene	18.26	178	662244	38.268	nq	98
71) Anthracene	18.35	178	687146	39.215	nq	99
72) Carbazole	18.61	167	604907	38.958	nq	97
73) Di-n-butylphthalate	19.10	149	712907	39.446	nq	# 98
74) Fluoranthene	20.21	202	813779	38.000	nq	97
76) Benzidine	20.37	184	431436	47.189	nq	99
77) Pyrene	20.57	202	845009	37.580	nq	99
79) Butylbenzylphthalate	21.42	149	324896	39.155	nq	98
80) Benzo(a)anthracene	22.54	228	815237	37.975	nq	99
81) 3,3'-Dichlorobenzidine	22.42	252	318968	41.754	nq	# 97
82) Chrysene	22.62	228	806537	38.812	nq	98
83) Bis(2-ethylhexyl)phthalate	22.36	149	456193	39.883	nq	98
84) Di-n-octyl phthalate	23.76	149	734414	41.935	nq	99
85) Indeno(1,2,3-cd)pyrene	30.71	276	932639	41.510	nq	# 95
87) Benzo(b)fluoranthene	25.12	252	792627	37.040	nq	# 94
88) Benzo(k)fluoranthene	25.20	252	827562	38.237	nq	# 98
89) Benzo(a)pyrene	26.15	252	785580	38.227	nq	# 98
90) Dibenzo(a,h)anthracene	30.79	278	771509	39.855	nq	# 97
91) Benzo(g,h,i)perylene	32.09	276	761394	39.721	nq	# 92

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

