

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG121418\
 Data File : BG038588.D
 Acq On : 15 Dec 2018 4:20
 Operator : JU/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Sohil
 12/17/2018 4:11:57 PM

Quant Time: Dec 15 05:50:19 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG121218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 12 15:26:27 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.06	152	30539	20.00	ng	0.00
21) Naphthalene-d8	10.88	136	128344	20.00	ng	0.00
39) Acenaphthene-d10	14.70	164	81925	20.00	ng	0.00
64) Phenanthrene-d10	17.45	188	197187	20.00	ng	0.00
76) Chrysene-d12	21.72	240	193943	20.00	ng	0.00
87) Perylene-d12	25.02	264	208612	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.61	112	142635	82.84	ng	0.00
7) Phenol-d6	7.21	99	196749	83.24	ng	0.00
23) Nitrobenzene-d5	9.24	82	206307	82.12	ng	0.00
42) 2,4,6-Tribromophenol	16.19	330	90893	80.50	ng	0.00
45) 2-Fluorobiphenyl	13.32	172	471504	78.95	ng	0.00
79) Terphenyl-d14	20.04	244	718233	80.60	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.49	88	31048	38.745	ng	96
3) Pyridine	3.90	79	86234	39.085	ng	# 89
4) n-Nitrosodimethylamine	3.82	42	46598	43.104	ng	# 85
6) Aniline	7.39	93	124225	41.169	ng	98
8) 2-Chlorophenol	7.62	128	81899	41.103	ng	97
9) Benzaldehyde	7.20	77	59875	39.048	ng	96
10) Phenol	7.24	94	103305	41.138	ng	97
11) bis(2-Chloroethyl)ether	7.48	93	81976	41.002	ng	95
12) 1,3-Dichlorobenzene	7.95	146	96045	40.767	ng	100
13) 1,4-Dichlorobenzene	8.09	146	95909	39.749	ng	99
14) 1,2-Dichlorobenzene	8.41	146	90775	39.906	ng	99
15) Benzyl Alcohol	8.30	79	79327	42.996	ng	99
16) 2,2'-oxybis(1-Chloropropan	8.58	45	184466	40.277	ng	99
17) 2-Methylphenol	8.50	107	70825	42.096	ng	96
18) Hexachloroethane	9.14	117	35825	40.632	ng	92
19) n-Nitroso-di-n-propylamine	8.86	70	68122	41.034	ng	# 98
20) 3+4-Methylphenols	8.83	107	100142	43.098	ng	96
22) Acetophenone	8.89	105	125934	39.284	ng	98
24) Nitrobenzene	9.28	77	102664	40.395	ng	97
25) Isophorone	9.79	82	171683	38.035	ng	# 98
26) 2-Nitrophenol	9.99	139	53055	40.787	ng	96
27) 2,4-Dimethylphenol	10.03	122	77129	41.373	ng	97
28) bis(2-Chloroethoxy)methane	10.27	93	103131	40.487	ng	96
29) 2,4-Dichlorophenol	10.52	162	88609	42.725	ng	98
30) 1,2,4-Trichlorobenzene	10.74	180	102377	40.867	ng	98
31) Naphthalene	10.93	128	249793	39.502	ng	99
32) Benzoic acid	10.19	122	41209m	37.596	ng	
33) 4-Chloroaniline	11.04	127	112652	41.033	ng	98
34) Hexachlorobutadiene	11.19	225	69276	40.869	ng	98
35) Caprolactam	11.84	113	32105	37.406	ng	96
36) 4-Chloro-3-methylphenol	12.15	107	93607	39.552	ng	96
37) 2-Methylnaphthalene	12.53	142	183264	40.623	ng	96
38) 1-Methylnaphthalene	12.75	142	174144	39.780	ng	99
40) 1,2,4,5-Tetrachlorobenzene	12.89	216	124267	40.579	ng	99

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG121418\
 Data File : BG038588.D
 Acq On : 15 Dec 2018 4:20
 Operator : JU/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Sohil
 12/17/2018 4:11:57 PM

Quant Time: Dec 15 05:50:19 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG121218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 12 15:26:27 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.85	237	65316	38.822	nq	96
43) 2,4,6-Trichlorophenol	13.13	196	77949	41.398	nq	97
44) 2,4,5-Trichlorophenol	13.21	196	81081	40.671	nq	97
46) 1,1'-Biphenyl	13.53	154	274433	39.958	nq	99
47) 2-Chloronaphthalene	13.58	162	210822	39.739	nq	96
48) 2-Nitroaniline	13.79	65	71335	42.215	nq	99
49) Acenaphthylene	14.42	152	315528	39.784	nq	98
50) Dimethylphthalate	14.15	163	267470	39.979	nq	99
51) 2,6-Dinitrotoluene	14.28	165	61204	40.369	nq	96
52) Acenaphthene	14.76	154	185699	39.157	nq	96
53) 3-Nitroaniline	14.62	138	63648	41.182	nq	99
54) 2,4-Dinitrophenol	14.82	184	28457	34.683	nq	93
55) Dibenzofuran	15.09	168	321523	39.551	nq	99
56) 4-Nitrophenol	14.92	139	46478	39.641	nq	97
57) 2,4-Dinitrotoluene	15.06	165	85186	39.892	nq	100
58) Fluorene	15.74	166	236820	39.320	nq	99
59) 2,3,4,6-Tetrachlorophenol	15.32	232	71465	39.844	nq	99
60) Diethylphthalate	15.50	149	283240	38.565	nq	99
61) 4-Chlorophenyl-phenylether	15.73	204	150140	39.954	nq	98
62) 4-Nitroaniline	15.78	138	65180	40.965	nq	98
63) Azobenzene	16.02	77	240726	39.235	nq	98
65) 4,6-Dinitro-2-methylphenol	15.83	198	56171	39.933	nq	96
66) n-Nitrosodiphenylamine	15.95	169	233514	40.304	nq	98
67) 4-Bromophenyl-phenylether	16.62	248	97414	40.451	nq	98
68) Hexachlorobenzene	16.74	284	100727	40.349	nq	97
69) Atrazine	16.89	200	88100	40.974	nq	98
70) Pentachlorophenol	17.09	266	60354	40.874	nq	95
71) Phenanthrene	17.49	178	406190	39.724	nq	99
72) Anthracene	17.58	178	406670	40.286	nq	100
73) Carbazole	17.85	167	402503	40.317	nq	97
74) Di-n-butylphthalate	18.38	149	464198	40.522	nq	100
75) Fluoranthene	19.49	202	500499	39.560	nq	97
77) Benzidine	19.67	184	253789	44.247	nq	98
78) Pyrene	19.86	202	502322	39.206	nq	100
80) Butylbenzylphthalate	20.73	149	207074	41.368	nq	98
81) Benzo(a)anthracene	21.70	228	480402	40.650	nq	99
82) 3,3'-Dichlorobenzidine	21.62	252	195192	41.645	nq	98
83) Chrysene	21.77	228	454978	39.970	nq	99
84) Bis(2-ethylhexyl)phthalate	21.57	149	295061	41.561	nq	99
85) Di-n-octyl phthalate	22.80	149	490743	41.275	nq	99
86) Indeno(1,2,3-cd)pyrene	28.80	276	542696	40.553	nq	# 86
88) Benzo(b)fluoranthene	23.96	252	470450	41.074	nq	99
89) Benzo(k)fluoranthene	24.03	252	463977	40.680	nq	98
90) Benzo(a)pyrene	24.86	252	458520	41.496	nq	99
91) Dibenzo(a,h)anthracene	28.86	278	455371	41.390	nq	99
92) Benzo(g,h,i)perylene	30.00	276	442609	40.822	nq	98

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

