

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG092118\
 Data File : BG036861.D
 Acq On : 21 Sep 2018 16:42
 Operator : JU/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040EC

Manual Integrations
 APPROVED

Sohil
 9/24/2018 9:29:41 AM

Quant Time: Sep 21 17:51:18 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG092018.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Sep 20 20:07:11 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.05	152	48658	20.00	ng	0.00
21) Naphthalene-d8	10.86	136	224593	20.00	ng	0.00
38) Acenaphthene-d10	14.67	164	153053	20.00	ng	0.00
63) Phenanthrene-d10	17.42	188	385448	20.00	ng	0.00
75) Chrysene-d12	21.69	240	402945	20.00	ng	0.00
86) Perylene-d12	24.90	264	423429	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.62	112	187135	73.76	ng	0.00
7) Phenol-d6	7.21	99	274464	69.92	ng	0.00
23) Nitrobenzene-d5	9.21	82	284203	67.76	ng	0.00
41) 2,4,6-Tribromophenol	16.17	330	125831	68.72	ng	0.00
44) 2-Fluorobiphenyl	13.30	172	673127	65.50	ng	0.00
78) Terphenyl-d14	20.03	244	970755	63.35	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.53	88	44350	38.200	ng	98
3) Pyridine	3.93	79	124543	37.152	ng	97
4) n-Nitrosodimethylamine	3.84	42	53780	36.096	ng	# 95
6) Aniline	7.38	93	168510	33.083	ng	99
8) 2-Chlorophenol	7.61	128	101984	34.618	ng	96
9) Benzaldehyde	7.19	77	88638	34.172	ng	94
10) Phenol	7.24	94	141215	34.857	ng	97
11) bis(2-Chloroethyl)ether	7.46	93	117390	31.325	ng	99
12) 1,3-Dichlorobenzene	7.94	146	123002	34.056	ng	98
13) 1,4-Dichlorobenzene	8.08	146	123965	33.539	ng	99
14) 1,2-Dichlorobenzene	8.41	146	118790	33.165	ng	95
15) Benzyl Alcohol	8.29	79	104048	32.666	ng	97
16) 2,2'-oxybis(1-Chloropropan	8.58	45	234832	32.640	ng	98
17) 2-Methylphenol	8.49	107	91209	32.321	ng	95
18) Hexachloroethane	9.13	117	45648	33.561	ng	91
19) n-Nitroso-di-n-propylamine	8.85	70	104734	32.072	ng	95
20) 3+4-Methylphenols	8.82	107	131742	33.557	ng	97
22) Acetophenone	8.87	105	167853	31.803	ng	96
24) Nitrobenzene	9.26	77	147840	33.009	ng	98
25) Isophorone	9.77	82	256470	33.467	ng	99
26) 2-Nitrophenol	9.97	139	64131	35.923	ng	94
27) 2,4-Dimethylphenol	10.02	122	92271	33.181	ng	99
28) bis(2-Chloroethoxy)methane	10.26	93	157579	33.324	ng	99
29) 2,4-Dichlorophenol	10.50	162	115715	35.896	ng	91
30) 1,2,4-Trichlorobenzene	10.71	180	137656	34.122	ng	99
31) Naphthalene	10.91	128	352197	32.958	ng	99
32) Benzoic acid	10.16	122	65952m	33.021	ng	
33) 4-Chloroaniline	11.01	127	161886	33.318	ng	98
34) Hexachlorobutadiene	11.19	225	95497	34.230	ng	97
35) Caprolactam	11.78	113	43519	32.800	ng	98
36) 4-Chloro-3-methylphenol	12.13	107	131436	34.171	ng	98
37) 2-Methylnaphthalene	12.51	142	265689	32.644	ng	98
39) 1,2,4,5-Tetrachlorobenzene	12.87	216	177989	33.342	ng	99
40) Hexachlorocyclopentadiene	12.85	237	94502	34.421	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.11	196	102706	34.647	ng	99
43) 2,4,5-Trichlorophenol	13.18	196	111458	35.262	ng	98
45) 1,1'-Biphenyl	13.51	154	380366	32.450	ng	98
46) 2-Chloronaphthalene	13.55	162	302191	32.782	ng	99
47) 2-Nitroaniline	13.76	65	105090	32.960	ng	99
48) Acenaphthylene	14.40	152	470611	32.580	ng	98
49) Dimethylphthalate	14.13	163	396282	32.166	ng	100
50) 2,6-Dinitrotoluene	14.25	165	88693	32.997	ng	97
51) Acenaphthene	14.74	154	278971	31.955	ng	98
52) 3-Nitroaniline	14.59	138	91477	32.296	ng	98
53) 2,4-Dinitrophenol	14.79	184	35992	33.489	ng	89
54) Dibenzofuran	15.07	168	476273	32.253	ng	99
55) 4-Nitrophenol	14.89	139	62823	34.719	ng	99
56) 2,4-Dinitrotoluene	15.03	165	124590	33.218	ng	91
57) Fluorene	15.72	166	356424	32.293	ng	97
58) 2,3,4,6-Tetrachlorophenol	15.30	232	112798	35.178	ng	93
59) Diethylphthalate	15.48	149	403474	32.471	ng	98
60) 4-Chlorophenyl-phenylether	15.71	204	228067	32.629	ng	97
61) 4-Nitroaniline	15.75	138	99291	33.327	ng	96
62) Azobenzene	16.01	77	392521	32.876	ng	98
64) 4,6-Dinitro-2-methylphenol	15.80	198	70996	35.231	ng	96
65) n-Nitrosodiphenylamine	15.93	169	349836	32.876	ng	97
66) 4-Bromophenyl-phenylether	16.61	248	145719	33.237	ng	98
67) Hexachlorobenzene	16.72	284	152963	33.875	ng	98
68) Atrazine	16.87	200	146321	33.960	ng	98
69) Pentachlorophenol	17.07	266	95335	33.534	ng	98
70) Phenanthrene	17.47	178	624535	33.623	ng	99
71) Anthracene	17.55	178	619010	33.703	ng	99
72) Carbazole	17.82	167	610562	34.095	ng	97
73) Di-n-butylphthalate	18.38	149	686266	34.508	ng	99
74) Fluoranthene	19.47	202	766456	34.280	ng	100
76) Benzidine	19.65	184	411766	33.804	ng	99
77) Pyrene	19.83	202	785077	32.468	ng	100
79) Butylbenzylphthalate	20.71	149	312213	32.394	ng	94
80) Benzo(a)anthracene	21.67	228	764152	32.487	ng	98
81) 3,3'-Dichlorobenzidine	21.58	252	299718	33.476	ng	99
82) Chrysene	21.74	228	735353	32.356	ng	99
83) Bis(2-ethylhexyl)phthalate	21.57	149	429264	31.882	ng	100
84) Di-n-octyl phthalate	22.78	149	721057	32.527	ng	98
85) Indeno(1,2,3-cd)pyrene	28.56	276	841889	34.497	ng	# 94
87) Benzo(b)fluoranthene	23.88	252	756661	33.532	ng	97
88) Benzo(k)fluoranthene	23.95	252	722595	31.918	ng	99
89) Benzo(a)pyrene	24.74	252	721362	33.066	ng	99
90) Dibenzo(a,h)anthracene	28.62	278	683358	33.422	ng	98
91) Benzo(g,h,i)perylene	29.70	276	684272	33.347	ng	97

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

