

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080218\
 Data File : BG036070.D
 Acq On : 2 Aug 2018 22:42
 Operator : JU/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Sohil
 8/3/2018 12:58:38 PM

Quant Time: Aug 03 09:06:19 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG071218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 31 12:30:06 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.02	152	32206	20.00	ng	0.00
21) Naphthalene-d8	10.82	136	135845	20.00	ng	0.00
38) Acenaphthene-d10	14.64	164	103749	20.00	ng	0.00
63) Phenanthrene-d10	17.39	188	290046	20.00	ng	0.00
75) Chrysene-d12	21.65	240	323092	20.00	ng	0.00
86) Perylene-d12	24.85	264	322883	20.00	ng	-0.01

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.60	112	146785	75.67	ng	0.00
7) Phenol-d6	7.19	99	227237	78.51	ng	0.00
23) Nitrobenzene-d5	9.19	82	249401	76.70	ng	0.00
41) 2,4,6-Tribromophenol	16.14	330	117859	86.19	ng	0.00
44) 2-Fluorobiphenyl	13.27	172	602175	77.76	ng	0.00
78) Terphenyl-d14	19.99	244	1114816	76.06	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	3.52	88	34475	34.918	ng	# 69
3) Pyridine	3.92	79	91309	35.426	ng	# 72
4) n-Nitrosodimethylamine	3.83	42	39575	35.759	ng	# 82
6) Aniline	7.35	93	139581	37.208	ng	96
8) 2-Chlorophenol	7.59	128	80884	40.997	ng	93
9) Benzaldehyde	7.16	77	64305	36.211	ng	97
10) Phenol	7.22	94	115308	39.435	ng	92
11) bis(2-Chloroethyl)ether	7.44	93	88675	38.724	ng	85
12) 1,3-Dichlorobenzene	7.91	146	95669	40.155	ng	96
13) 1,4-Dichlorobenzene	8.06	146	98026	40.494	ng	95
14) 1,2-Dichlorobenzene	8.38	146	93501	40.207	ng	93
15) Benzyl Alcohol	8.26	79	97377	37.050	ng	92
16) 2,2'-oxybis(1-Chloropropan	8.55	45	67636	35.230	ng	76
17) 2-Methylphenol	8.46	107	78109	39.289	ng	# 92
18) Hexachloroethane	9.10	117	36053	38.952	ng	95
19) n-Nitroso-di-n-propylamine	8.83	70	89584	36.757	ng	# 84
20) 3+4-Methylphenols	8.79	107	112697	40.862	ng	95
22) Acetophenone	8.84	105	159362	37.724	ng	# 92
24) Nitrobenzene	9.23	77	123163	37.661	ng	95
25) Isophorone	9.74	82	224539	37.197	ng	99
26) 2-Nitrophenol	9.94	139	50878	43.805	ng	# 85
27) 2,4-Dimethylphenol	10.00	122	77126	39.229	ng	90
28) bis(2-Chloroethoxy)methane	10.23	93	124005	37.280	ng	97
29) 2,4-Dichlorophenol	10.48	162	92637	41.277	ng	96
30) 1,2,4-Trichlorobenzene	10.68	180	111113	40.376	ng	97
31) Naphthalene	10.87	128	279624	39.516	ng	100
32) Benzoic acid	10.17	122	34562m	26.114	ng	
33) 4-Chloroaniline	10.98	127	118359	38.376	ng	# 87
34) Hexachlorobutadiene	11.15	225	87051	40.565	ng	96
35) Caprolactam	11.78	113	36453	37.850	ng	# 67
36) 4-Chloro-3-methylphenol	12.11	107	122131	40.599	ng	91
37) 2-Methylnaphthalene	12.48	142	216157	41.001	ng	96
39) 1,2,4,5-Tetrachlorobenzene	12.84	216	151167	38.604	ng	99
40) Hexachlorocyclopentadiene	12.82	237	69592	33.887	ng	91

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.08	196	91857	40.112	nq	96
43) 2,4,5-Trichlorophenol	13.16	196	98403	38.411	nq	96
45) 1,1'-Biphenyl	13.48	154	310623	38.617	nq	99
46) 2-Chloronaphthalene	13.52	162	243613	38.937	nq	99
47) 2-Nitroaniline	13.73	65	84376	38.146	nq	88
48) Acenaphthylene	14.37	152	409316	39.055	nq	97
49) Dimethylphthalate	14.10	163	355464	39.105	nq	99
50) 2,6-Dinitrotoluene	14.22	165	77939	42.105	nq	96
51) Acenaphthene	14.71	154	253205	38.783	nq	99
52) 3-Nitroaniline	14.55	138	76098	40.223	nq	# 95
53) 2,4-Dinitrophenol	14.77	184	27370m	32.970	nq	
54) Dibenzofuran	15.04	168	409646	39.453	nq	93
55) 4-Nitrophenol	14.87	139	44091m	32.724	nq	
56) 2,4-Dinitrotoluene	15.01	165	112648	42.419	nq	89
57) Fluorene	15.69	166	348425	40.037	nq	95
58) 2,3,4,6-Tetrachlorophenol	15.27	232	99803	40.632	nq	# 94
59) Diethylphthalate	15.46	149	362283	38.760	nq	96
60) 4-Chlorophenyl-phenylether	15.68	204	203076	39.703	nq	97
61) 4-Nitroaniline	15.72	138	80069	40.459	nq	# 75
62) Azobenzene	15.98	77	345644	37.490	nq	94
64) 4,6-Dinitro-2-methylphenol	15.78	198	60610	37.539	nq	83
65) n-Nitrosodiphenylamine	15.90	169	324690	39.329	nq	98
66) 4-Bromophenyl-phenylether	16.58	248	137306	40.804	nq	95
67) Hexachlorobenzene	16.70	284	141925	40.956	nq	95
68) Atrazine	16.85	200	118876	34.972	nq	98
69) Pentachlorophenol	17.05	266	69510m	32.932	nq	
70) Phenanthrene	17.43	178	563142	38.910	nq	100
71) Anthracene	17.52	178	562197	39.012	nq	98
72) Carbazole	17.79	167	524214	38.303	nq	98
73) Di-n-butylphthalate	18.34	149	613517	37.935	nq	# 98
74) Fluoranthene	19.44	202	747418	38.340	nq	96
76) Benzidine	19.62	184	338156	39.810	nq	98
77) Pyrene	19.80	202	755170	36.965	nq	97
79) Butylbenzylphthalate	20.68	149	290773	39.801	nq	# 94
80) Benzo(a)anthracene	21.63	228	771065	38.296	nq	99
81) 3,3'-Dichlorobenzidine	21.55	252	282338	40.217	nq	98
82) Chrysene	21.70	228	714651	38.303	nq	96
83) Bis(2-ethylhexyl)phthalate	21.53	149	402252	40.500	nq	99
84) Di-n-octyl phthalate	22.73	149	682016	41.011	nq	# 94
85) Indeno(1,2,3-cd)pyrene	28.48	276	816654	39.534	nq	# 89
87) Benzo(b)fluoranthene	23.84	252	732854	37.343	nq	# 96
88) Benzo(k)fluoranthene	23.90	252	736282	38.376	nq	# 96
89) Benzo(a)pyrene	24.70	252	702193	38.461	nq	# 95
90) Dibenzo(a,h)anthracene	28.54	278	685211	38.229	nq	# 94
91) Benzo(g,h,i)perylene	29.63	276	677436	38.623	nq	# 93

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

