

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG073018\
 Data File : BG035987.D
 Acq On : 30 Jul 2018 23:28
 Operator : JU/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Sohil
 7/31/2018 12:20:58 PM

Quant Time: Jul 31 01:43:57 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG071218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 25 11:04:14 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.02	152	34008	20.00	ng	0.00
21) Naphthalene-d8	10.82	136	143607	20.00	ng	0.00
38) Acenaphthene-d10	14.65	164	106288	20.00	ng	0.00
63) Phenanthrene-d10	17.39	188	281763	20.00	ng	0.00
75) Chrysene-d12	21.65	240	293526	20.00	ng	0.00
86) Perylene-d12	24.85	264	292463	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.61	112	147999	72.25	ng	0.00
7) Phenol-d6	7.19	99	228571	74.79	ng	0.00
23) Nitrobenzene-d5	9.18	82	258638	75.24	ng	0.00
41) 2,4,6-Tribromophenol	16.14	330	109300	78.02	ng	0.00
44) 2-Fluorobiphenyl	13.27	172	608899	76.75	ng	0.00
78) Terphenyl-d14	20.00	244	1037624	77.92	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.52	88	35088	33.656	ng	# 70
3) Pyridine	3.91	79	95772	35.188	ng	# 70
4) n-Nitrosodimethylamine	3.83	42	37869	32.404	ng	# 71
6) Aniline	7.35	93	143201	36.151	ng	99
8) 2-Chlorophenol	7.59	128	80893	38.829	ng	94
9) Benzaldehyde	7.17	77	62759	33.468	ng	95
10) Phenol	7.22	94	116125	37.610	ng	96
11) bis(2-Chloroethyl)ether	7.44	93	91353	37.779	ng	85
12) 1,3-Dichlorobenzene	7.91	146	96222	38.247	ng	95
13) 1,4-Dichlorobenzene	8.06	146	96804	37.871	ng	98
14) 1,2-Dichlorobenzene	8.38	146	94740	38.581	ng	96
15) Benzyl Alcohol	8.26	79	101612	36.613	ng	90
16) 2,2'-oxybis(1-Chloropropan	8.54	45	72241	35.635	ng	78
17) 2-Methylphenol	8.47	107	79941	38.080	ng	# 92
18) Hexachloroethane	9.10	117	37407	38.274	ng	90
19) n-Nitroso-di-n-propylamine	8.83	70	91201	35.438	ng	# 87
20) 3+4-Methylphenols	8.80	107	112426	38.604	ng	86
22) Acetophenone	8.84	105	165351	37.026	ng	# 93
24) Nitrobenzene	9.22	77	125625	36.337	ng	96
25) Isophorone	9.75	82	230653	36.144	ng	97
26) 2-Nitrophenol	9.94	139	51188	41.690	ng	# 91
27) 2,4-Dimethylphenol	9.99	122	80838	38.895	ng	92
28) bis(2-Chloroethoxy)methane	10.23	93	132853	37.782	ng	99
29) 2,4-Dichlorophenol	10.48	162	96287	40.585	ng	96
30) 1,2,4-Trichlorobenzene	10.69	180	117139	40.265	ng	92
31) Naphthalene	10.88	128	285705	38.193	ng	98
32) Benzoic acid	10.17	122	24367m	17.416	ng	
33) 4-Chloroaniline	10.99	127	122444	37.555	ng	94
34) Hexachlorobutadiene	11.16	225	91030	40.127	ng	98
35) Caprolactam	11.77	113	35410	34.779	ng	# 61
36) 4-Chloro-3-methylphenol	12.11	107	120298	37.828	ng	100
37) 2-Methylnaphthalene	12.48	142	220741	39.608	ng	92
39) 1,2,4,5-Tetrachlorobenzene	12.84	216	151948	37.876	ng	97
40) Hexachlorocyclopentadiene	12.82	237	69806	33.179	ng	89

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.08	196	92468	39.414	nq	95
43) 2,4,5-Trichlorophenol	13.16	196	103116	39.290	nq	91
45) 1,1'-Biphenyl	13.48	154	317500	38.529	nq	96
46) 2-Chloronaphthalene	13.52	162	243808	38.037	nq	97
47) 2-Nitroaniline	13.73	65	83361	36.787	nq	93
48) Acenaphthylene	14.37	152	414117	38.569	nq	98
49) Dimethylphthalate	14.10	163	349938	37.578	nq	99
50) 2,6-Dinitrotoluene	14.22	165	78166	41.219	nq	94
51) Acenaphthene	14.71	154	257832	38.548	nq	98
52) 3-Nitroaniline	14.55	138	72975	37.651	nq #	97
53) 2,4-Dinitrophenol	14.77	184	17151	22.723	nq #	73
54) Dibenzofuran	15.05	168	411951	38.727	nq	95
55) 4-Nitrophenol	14.88	139	40591	29.407	nq #	79
56) 2,4-Dinitrotoluene	15.01	165	113682	41.786	nq #	86
57) Fluorene	15.69	166	344584	38.650	nq	99
58) 2,3,4,6-Tetrachlorophenol	15.28	232	93867	37.303	nq	95
59) Diethylphthalate	15.46	149	354279	36.998	nq	97
60) 4-Chlorophenyl-phenylether	15.68	204	204158	38.961	nq	96
61) 4-Nitroaniline	15.72	138	75306	37.143	nq #	78
62) Azobenzene	15.97	77	338187	35.805	nq	96
64) 4,6-Dinitro-2-methylphenol	15.77	198	57945	36.944	nq	88
65) n-Nitrosodiphenylamine	15.90	169	314725	39.242	nq	100
66) 4-Bromophenyl-phenylether	16.58	248	134442	41.127	nq	95
67) Hexachlorobenzene	16.70	284	138323	41.090	nq	96
68) Atrazine	16.84	200	116972	35.424	nq	95
69) Pentachlorophenol	17.04	266	58537	28.548	nq	98
70) Phenanthrene	17.43	178	543223	38.637	nq	97
71) Anthracene	17.53	178	545729	38.982	nq	99
72) Carbazole	17.80	167	494320	37.181	nq	98
73) Di-n-butylphthalate	18.35	149	584672	37.214	nq	98
74) Fluoranthene	19.44	202	699178	36.919	nq	97
76) Benzidine	19.62	184	191467	24.811	nq	98
77) Pyrene	19.80	202	716374	38.598	nq	99
79) Butylbenzylphthalate	20.68	149	262348	39.527	nq	98
80) Benzo(a)anthracene	21.63	228	699181	38.224	nq	99
81) 3,3'-Dichlorobenzidine	21.55	252	248387	38.945	nq	98
82) Chrysene	21.70	228	657548	38.792	nq	98
83) Bis(2-ethylhexyl)phthalate	21.53	149	359825	39.878	nq	99
84) Di-n-octyl phthalate	22.73	149	623862	41.293	nq #	93
85) Indeno(1,2,3-cd)pyrene	28.49	276	729059	38.849	nq #	89
87) Benzo(b)fluoranthene	23.83	252	659382	37.094	nq #	97
88) Benzo(k)fluoranthene	23.90	252	677544	38.988	nq #	98
89) Benzo(a)pyrene	24.70	252	633290	38.295	nq #	95
90) Dibenzo(a,h)anthracene	28.55	278	609592	37.547	nq #	97
91) Benzo(g,h,i)perylene	29.63	276	590731	37.183	nq #	92

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

