

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG071718\
 Data File : BG035700.D
 Acq On : 18 Jul 2018 3:10
 Operator : JU/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Sohil
 7/18/2018 3:27:52 PM

Quant Time: Jul 18 04:47:57 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG071218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 12 14:43:32 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.04	152	57690	20.00	ng	-0.02
21) Naphthalene-d8	10.84	136	236945	20.00	ng	-0.02
38) Acenaphthene-d10	14.66	164	162888	20.00	ng	-0.02
63) Phenanthrene-d10	17.41	188	420818	20.00	ng	-0.02
75) Chrysene-d12	21.68	240	445388	20.00	ng	0.00
86) Perylene-d12	24.88	264	442084	20.00	ng	-0.03

System Monitoring Compounds

5) 2-Fluorophenol	5.62	112	278073	80.03	ng	0.00
7) Phenol-d6	7.21	99	414152	79.88	ng	0.00
23) Nitrobenzene-d5	9.21	82	429591	75.74	ng	-0.02
41) 2,4,6-Tribromophenol	16.16	330	171277	79.78	ng	0.00
44) 2-Fluorobiphenyl	13.29	172	963282	79.23	ng	-0.02
78) Terphenyl-d14	20.01	244	1531748	75.81	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.54	88	63207	35.739	ng	# 68
3) Pyridine	3.94	79	181370	39.283	ng	# 76
4) n-Nitrosodimethylamine	3.85	42	71430	36.031	ng	# 72
6) Aniline	7.37	93	253697	37.754	ng	100
8) 2-Chlorophenol	7.61	128	144596	40.915	ng	99
9) Benzaldehyde	7.19	77	119243	37.486	ng	96
10) Phenol	7.24	94	216632	41.360	ng	98
11) bis(2-Chloroethyl)ether	7.46	93	161463	39.363	ng	89
12) 1,3-Dichlorobenzene	7.93	146	173435	40.639	ng	95
13) 1,4-Dichlorobenzene	8.08	146	174090	40.148	ng	99
14) 1,2-Dichlorobenzene	8.40	146	165303	39.682	ng	94
15) Benzyl Alcohol	8.28	79	178420	37.898	ng	93
16) 2,2'-oxybis(1-Chloropropan	8.56	45	153017m	44.495	ng	
17) 2-Methylphenol	8.48	107	139256	39.104	ng	94
18) Hexachloroethane	9.12	117	66192	39.924	ng	91
19) n-Nitroso-di-n-propylamine	8.84	70	157295	36.030	ng	# 87
20) 3+4-Methylphenols	8.81	107	199071	40.295	ng	91
22) Acetophenone	8.87	105	281547	38.211	ng	# 90
24) Nitrobenzene	9.25	77	212017	37.168	ng	92
25) Isophorone	9.76	82	395909	37.601	ng	98
26) 2-Nitrophenol	9.96	139	90234	44.541	ng	# 92
27) 2,4-Dimethylphenol	10.02	122	137195	40.007	ng	90
28) bis(2-Chloroethoxy)methane	10.25	93	223983	38.606	ng	98
29) 2,4-Dichlorophenol	10.49	162	161106	41.156	ng	90
30) 1,2,4-Trichlorobenzene	10.71	180	198086	41.267	ng	96
31) Naphthalene	10.90	128	478017	38.729	ng	97
32) Benzoic acid	10.18	122	88870	38.496	ng	98
33) 4-Chloroaniline	11.00	127	207631	38.597	ng	96
34) Hexachlorobutadiene	11.18	225	154644	41.315	ng	98
35) Caprolactam	11.78	113	60302	35.897	ng	# 71
36) 4-Chloro-3-methylphenol	12.13	107	204969	39.064	ng	96
37) 2-Methylnaphthalene	12.50	142	359639	39.110	ng	95
39) 1,2,4,5-Tetrachlorobenzene	12.86	216	248948	40.493	ng	98
40) Hexachlorocyclopentadiene	12.84	237	137623	42.684	ng	91

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.10	196	153479	42.688	nq	98
43) 2,4,5-Trichlorophenol	13.18	196	169057	42.032	nq	97
45) 1,1'-Biphenyl	13.50	154	508664	40.279	nq	98
46) 2-Chloronaphthalene	13.55	162	394650	40.176	nq	99
47) 2-Nitroaniline	13.75	65	135808	39.106	nq	91
48) Acenaphthylene	14.39	152	647997	39.380	nq	98
49) Dimethylphthalate	14.12	163	548764	38.452	nq	# 98
50) 2,6-Dinitrotoluene	14.24	165	122678	42.213	nq	92
51) Acenaphthene	14.73	154	398893	38.915	nq	96
52) 3-Nitroaniline	14.57	138	117612	39.596	nq	# 95
53) 2,4-Dinitrophenol	14.82	184	3169	8.528	nq	# 43
54) Dibenzofuran	15.06	168	635259	38.969	nq	93
55) 4-Nitrophenol	14.89	139	85173	40.264	nq	# 86
56) 2,4-Dinitrotoluene	15.03	165	171378	41.105	nq	# 84
57) Fluorene	15.71	166	526659	38.546	nq	97
58) 2,3,4,6-Tetrachlorophenol	15.29	232	150460	39.016	nq	94
59) Diethylphthalate	15.48	149	551280	37.567	nq	97
60) 4-Chlorophenyl-phenylether	15.70	204	316418	39.402	nq	97
61) 4-Nitroaniline	15.73	138	120760	38.866	nq	87
62) Azobenzene	16.00	77	520748	35.976	nq	95
64) 4,6-Dinitro-2-methylphenol	15.79	198	59869	25.557	nq	87
65) n-Nitrosodiphenylamine	15.91	169	485642	40.544	nq	100
66) 4-Bromophenyl-phenylether	16.60	248	205457	42.083	nq	98
67) Hexachlorobenzene	16.72	284	212167	42.199	nq	94
68) Atrazine	16.87	200	195467	39.635	nq	98
69) Pentachlorophenol	17.07	266	155620	50.817	nq	100
70) Phenanthrene	17.45	178	814560	38.792	nq	99
71) Anthracene	17.54	178	830150	39.704	nq	99
72) Carbazole	17.81	167	765997	38.577	nq	98
73) Di-n-butylphthalate	18.36	149	887577	37.826	nq	# 99
74) Fluoranthene	19.46	202	1061597	37.533	nq	97
76) Benzo(a)pyrene	19.64	184	424713	36.271	nq	99
77) Pyrene	19.82	202	1082860	38.451	nq	96
79) Butylbenzylphthalate	20.70	149	410340	40.745	nq	96
80) Benzo(a)anthracene	21.65	228	1069203	38.522	nq	99
81) 3,3'-Dichlorobenzidine	21.57	252	409279	42.291	nq	# 99
82) Chrysene	21.72	228	1000710	38.908	nq	98
83) Bis(2-ethylhexyl)phthalate	21.55	149	573076	41.856	nq	99
84) Di-n-octyl phthalate	22.76	149	974051	42.489	nq	# 92
85) Indeno(1,2,3-cd)pyrene	28.55	276	1169453	41.068	nq	# 87
87) Benzo(b)fluoranthene	23.87	252	1024582	38.131	nq	# 97
88) Benzo(k)fluoranthene	23.94	252	1033858	39.357	nq	# 95
89) Benzo(a)pyrene	24.74	252	977812	39.116	nq	# 97
90) Dibenzo(a,h)anthracene	28.61	278	972313	39.620	nq	# 96
91) Benzo(g,h,i)perylene	29.70	276	948160	39.483	nq	# 95

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

