

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG073118\
 Data File : BG036017.D
 Acq On : 1 Aug 2018 1:17
 Operator : JU/SJ
 Sample : J3831-04MS
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SU-03-072718MS

Manual Integrations
 APPROVED

Sohil
 8/1/2018 5:08:57 PM

Quant Time: Aug 01 11:13:37 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	38879	20.00	ng	0.00
21) Naphthalene-d8	10.83	136	141057	20.00	ng	0.00
38) Acenaphthene-d10	14.65	164	105098	20.00	ng	0.00
63) Phenanthrene-d10	17.39	188	290865	20.00	ng	0.00
75) Chrysene-d12	21.65	240	313108	20.00	ng	0.00
86) Perylene-d12	24.85	264	304044	20.00	ng	-0.01

System Monitoring Compounds

5) 2-Fluorophenol	5.61	112	301399	128.71	ng	0.00
7) Phenol-d6	7.20	99	405566	116.08	ng	0.00
23) Nitrobenzene-d5	9.19	82	331268	98.11	ng	0.00
41) 2,4,6-Tribromophenol	16.14	330	235683	170.14	ng	0.00
44) 2-Fluorobiphenyl	13.27	172	733275	93.47	ng	0.00
78) Terphenyl-d14	20.00	244	1442290	101.54	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.53	88	36604	30.711	ng	# 81
3) Pyridine	3.93	79	82679	26.572	ng	# 79
4) n-Nitrosodimethylamine	3.83	42	47436	35.505	ng	# 86
6) Aniline	7.35	93	68305	15.083	ng	94
8) 2-Chlorophenol	7.60	128	108266	45.457	ng	91
9) Benzaldehyde	7.16	77	39825	18.577	ng	88
10) Phenol	7.22	94	136277	38.607	ng	93
11) bis(2-Chloroethyl)ether	7.44	93	112895	40.839	ng	89
12) 1,3-Dichlorobenzene	7.91	146	123801	43.044	ng	96
13) 1,4-Dichlorobenzene	8.06	146	125516	42.951	ng	96
14) 1,2-Dichlorobenzene	8.38	146	122185	43.523	ng	99
15) Benzyl Alcohol	8.27	79	127704	40.250	ng	92
16) 2,2'-oxybis(1-Chloropropan	8.55	45	131481	56.731	ng	78
17) 2-Methylphenol	8.47	107	105988	44.162	ng	# 92
18) Hexachloroethane	9.10	117	47754	42.739	ng	94
19) n-Nitroso-di-n-propylamine	8.82	70	107658	36.591	ng	# 83
20) 3+4-Methylphenols	8.79	107	146909	44.125	ng	95
22) Acetophenone	8.85	105	198258	45.198	ng	# 91
24) Nitrobenzene	9.23	77	157265	46.311	ng	96
25) Isophorone	9.75	82	316715	50.528	ng	98
26) 2-Nitrophenol	9.94	139	66607	55.228	ng	# 93
27) 2,4-Dimethylphenol	10.00	122	115647	56.648	ng	90
28) bis(2-Chloroethoxy)methane	10.23	93	165330	47.868	ng	98
29) 2,4-Dichlorophenol	10.48	162	133224	57.168	ng	92
30) 1,2,4-Trichlorobenzene	10.69	180	143110	50.081	ng	96
31) Naphthalene	10.87	128	357977	48.719	ng	98
32) Benzoic acid	10.15	122	18448m	13.424	ng	
33) 4-Chloroaniline	11.01	127	12949	4.043	ng	# 89
34) Hexachlorobutadiene	11.16	225	107888	48.418	ng	97
35) Caprolactam	11.77	113	33321m	33.319	ng	
36) 4-Chloro-3-methylphenol	12.11	107	167361	53.579	ng	93
37) 2-Methylnaphthalene	12.48	142	281430	51.410	ng	96
39) 1,2,4,5-Tetrachlorobenzene	12.84	216	189667	47.814	ng	98
40) Hexachlorocyclopentadiene	12.82	237	230106	110.610	ng	92

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.08	196	126581	54.566	nq	98
43) 2,4,5-Trichlorophenol	13.17	196	138504	53.371	nq	99
45) 1,1'-Biphenyl	13.48	154	389733	47.831	nq	97
46) 2-Chloronaphthalene	13.52	162	312174	49.254	nq	96
47) 2-Nitroaniline	13.73	65	111867	49.925	nq	97
48) Acenaphthylene	14.37	152	518869	48.872	nq	99
49) Dimethylphthalate	14.11	163	445952	48.430	nq	99
50) 2,6-Dinitrotoluene	14.22	165	103052	54.957	nq	92
51) Acenaphthene	14.71	154	305981	46.265	nq	98
52) 3-Nitroaniline	14.56	138	33192	17.319	nq #	94
53) 2,4-Dinitrophenol	14.76	184	53542	57.540	nq #	64
54) Dibenzofuran	15.05	168	519541	49.394	nq	93
55) 4-Nitrophenol	14.87	139	114819	84.125	nq #	84
56) 2,4-Dinitrotoluene	15.01	165	151648	56.372	nq	91
57) Fluorene	15.69	166	437826	49.664	nq	95
58) 2,3,4,6-Tetrachlorophenol	15.27	232	142866	57.418	nq	94
59) Diethylphthalate	15.46	149	453915	47.940	nq	98
60) 4-Chlorophenyl-phenylether	15.69	204	256993	49.599	nq	96
61) 4-Nitroaniline	15.72	138	94927	47.351	nq #	86
62) Azobenzene	15.97	77	456756	48.906	nq	94
64) 4,6-Dinitro-2-methylphenol	15.77	198	63368	39.137	nq	85
65) n-Nitrosodiphenylamine	15.90	169	393409	47.518	nq	97
66) 4-Bromophenyl-phenylether	16.58	248	171235	50.743	nq	97
67) Hexachlorobenzene	16.70	284	175543	50.514	nq	94
68) Atrazine	16.85	200	170299	49.959	nq	95
69) Pentachlorophenol	17.05	266	159618	75.410	nq	99
70) Phenanthrene	17.44	178	725779	50.007	nq	99
71) Anthracene	17.52	178	730761	50.566	nq	99
72) Carbazole	17.79	167	660994	48.162	nq	99
73) Di-n-butylphthalate	18.35	149	758297	46.755	nq #	98
74) Fluoranthene	19.44	202	939674	48.066	nq	96
76) Benzidine	19.63	184	66120m	8.032	nq	
77) Pyrene	19.80	202	946031	47.784	nq	97
79) Butylbenzylphthalate	20.68	149	357599	50.509	nq	96
80) Benzo(a)anthracene	21.64	228	974232	49.930	nq	99
81) 3,3'-Dichlorobenzidine	21.55	252	173426	25.491	nq	97
82) Chrysene	21.70	228	911508	50.412	nq	97
83) Bis(2-ethylhexyl)phthalate	21.54	149	501175	52.069	nq #	98
84) Di-n-octyl phthalate	22.73	149	858196	53.251	nq #	92
85) Indeno(1,2,3-cd)pyrene	28.50	276	1047989	52.351	nq #	93
87) Benzo(b)fluoranthene	23.84	252	925772	50.097	nq #	97
88) Benzo(k)fluoranthene	23.91	252	913504	50.563	nq #	97
89) Benzo(a)pyrene	24.70	252	906463	52.726	nq #	96
90) Dibenzo(a,h)anthracene	28.56	278	872604	51.700	nq #	95
91) Benzo(g,h,i)perylene	29.63	276	869516	52.646	nq #	94

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

