

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG071818\
 Data File : BG035732.D
 Acq On : 19 Jul 2018 00:45
 Operator : JU/SJ
 Sample : J3826-04MSD
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 OK-01-071618MSD

Manual Integrations
 APPROVED

Sohil
 7/19/2018 3:45:22 PM

Quant Time: Jul 19 03:26:44 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 18 17:45:13 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.04	152	51424	20.00	ng	0.00
21) Naphthalene-d8	10.84	136	190379	20.00	ng	0.00
38) Acenaphthene-d10	14.66	164	129836	20.00	ng	0.00
63) Phenanthrene-d10	17.40	188	351356	20.00	ng	0.00
75) Chrysene-d12	21.67	240	392647	20.00	ng	0.00
86) Perylene-d12	24.88	264	385207	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.63	112	403066	130.13	ng	0.00
7) Phenol-d6	7.21	99	528539	114.37	ng	0.00
23) Nitrobenzene-d5	9.20	82	409905	89.95	ng	0.00
41) 2,4,6-Tribromophenol	16.15	330	270245	157.92	ng	0.00
44) 2-Fluorobiphenyl	13.29	172	863638	89.12	ng	0.00
78) Terphenyl-d14	20.01	244	1706035	95.78	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.55	88	54478	34.557	ng	# 73
3) Pyridine	3.95	79	140186	34.063	ng	# 74
4) n-Nitrosodimethylamine	3.85	42	65563	37.102	ng	# 73
6) Aniline	7.37	93	111983	18.695	ng	96
8) 2-Chlorophenol	7.61	128	152691	48.470	ng	96
9) Benzaldehyde	7.18	77	80492	28.387	ng	92
10) Phenol	7.24	94	188348	40.341	ng	99
11) bis(2-Chloroethyl)ether	7.46	93	163038	44.590	ng	90
12) 1,3-Dichlorobenzene	7.93	146	168243m	44.226	ng	
13) 1,4-Dichlorobenzene	8.08	146	174087	45.039	ng	97
14) 1,2-Dichlorobenzene	8.39	146	163566	44.050	ng	97
15) Benzyl Alcohol	8.28	79	178982	42.650	ng	90
16) 2,2'-oxybis(1-Chloropropan	8.56	45	177389	57.868	ng	76
17) 2-Methylphenol	8.49	107	151693	47.787	ng	94
18) Hexachloroethane	9.12	117	63578	43.020	ng	91
19) n-Nitroso-di-n-propylamine	8.84	70	146694	37.696	ng	# 86
20) 3+4-Methylphenols	8.82	107	201340	45.721	ng	94
22) Acetophenone	8.86	105	272396	46.011	ng	# 90
24) Nitrobenzene	9.24	77	217753	47.511	ng	# 94
25) Isophorone	9.76	82	424221	50.145	ng	98
26) 2-Nitrophenol	9.96	139	89290	54.855	ng	# 93
27) 2,4-Dimethylphenol	10.01	122	156400	56.763	ng	92
28) bis(2-Chloroethoxy)methane	10.24	93	221980	47.619	ng	98
29) 2,4-Dichlorophenol	10.50	162	181083	57.574	ng	92
30) 1,2,4-Trichlorobenzene	10.70	180	188864	48.970	ng	97
31) Naphthalene	10.90	128	474423	47.839	ng	99
32) Benzoic acid	10.14	122	19782	10.665	ng	91
33) 4-Chloroaniline	11.02	127	17058	3.947	ng	# 91
34) Hexachlorobutadiene	11.18	225	140068	46.574	ng	95
35) Caprolactam	11.78	113	45878m	33.991	ng	
36) 4-Chloro-3-methylphenol	12.13	107	216080	51.254	ng	95
37) 2-Methylnaphthalene	12.49	142	369800m	50.052	ng	
39) 1,2,4,5-Tetrachlorobenzene	12.86	216	246682	50.339	ng	99
40) Hexachlorocyclopentadiene	12.84	237	294349	114.532	ng	91

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.10	196	163699	57.121	nq	95
43) 2,4,5-Trichlorophenol	13.18	196	179148	55.880	nq	99
45) 1,1'-Biphenyl	13.50	154	504582	50.127	nq	99
46) 2-Chloronaphthalene	13.54	162	396617	50.655	nq	97
47) 2-Nitroaniline	13.74	65	143479	51.833	nq	96
48) Acenaphthylene	14.38	152	654557	49.906	nq	99
49) Dimethylphthalate	14.12	163	559401	49.176	nq	99
50) 2,6-Dinitrotoluene	14.24	165	128824	55.612	nq	95
51) Acenaphthene	14.73	154	388155	47.508	nq	99
52) 3-Nitroaniline	14.57	138	33961	14.344	nq #	95
53) 2,4-Dinitrophenol	14.78	184	7896	12.665	nq #	79
54) Dibenzofuran	15.06	168	649036	49.949	nq	94
55) 4-Nitrophenol	14.89	139	144299	85.580	nq #	87
56) 2,4-Dinitrotoluene	15.03	165	189445	57.005	nq #	87
57) Fluorene	15.71	166	540230	49.604	nq	97
58) 2,3,4,6-Tetrachlorophenol	15.29	232	167485	54.487	nq	95
59) Diethylphthalate	15.48	149	565123	48.313	nq	97
60) 4-Chlorophenyl-phenylether	15.70	204	309664	48.377	nq	96
61) 4-Nitroaniline	15.74	138	126685	51.152	nq #	80
62) Azobenzene	15.99	77	566589	49.107	nq	95
64) 4,6-Dinitro-2-methylphenol	15.79	198	30580	15.635	nq	87
65) n-Nitrosodiphenylamine	15.92	169	485187	48.514	nq	97
66) 4-Bromophenyl-phenylether	16.59	248	213222	52.307	nq	98
67) Hexachlorobenzene	16.71	284	221987	52.881	nq	96
68) Atrazine	16.86	200	228519	55.497	nq	97
69) Pentachlorophenol	17.06	266	110056	43.043	nq	99
70) Phenanthrene	17.45	178	890070	50.768	nq	97
71) Anthracene	17.54	178	904535	51.815	nq	97
72) Carbazole	17.81	167	846855	51.081	nq	97
73) Di-n-butylphthalate	18.36	149	966756	49.346	nq #	98
74) Fluoranthene	19.45	202	1193266	50.529	nq	96
76) Benzidine	19.64	184	120863	11.708	nq	98
77) Pyrene	19.82	202	1196475	48.191	nq	96
79) Butylbenzylphthalate	20.70	149	463104	52.161	nq #	94
80) Benzo(a)anthracene	21.65	228	1223276	49.993	nq	100
81) 3,3'-Dichlorobenzidine	21.57	252	178037	20.868	nq #	97
82) Chrysene	21.72	228	1143052	50.411	nq	96
83) Bis(2-ethylhexyl)phthalate	21.55	149	636887	52.765	nq #	97
84) Di-n-octyl phthalate	22.76	149	1091963	54.031	nq #	93
85) Indeno(1,2,3-cd)pyrene	28.54	276	1338802	53.331	nq #	93
87) Benzo(b)fluoranthene	23.87	252	1179701	50.387	nq #	97
88) Benzo(k)fluoranthene	23.94	252	1154531	50.440	nq #	96
89) Benzo(a)pyrene	24.74	252	1152116	52.894	nq #	97
90) Dibenzo(a,h)anthracene	28.60	278	1118816	52.321	nq #	96
91) Benzo(g,h,i)perylene	29.70	276	1110315	53.062	nq #	94

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

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