

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG060917\
 Data File : BG027356.D
 Acq On : 9 Jun 2017 20:46
 Operator : SJ/MA
 Sample : I3541-03MSD
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 RT-4215MSD

Manual Integrations
 APPROVED

Quant Time: Jun 10 06:41:10 2017
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 05 16:00:35 2017
 Response via : Initial Calibration

monammad
 6/12/2017 3:58:23 PM

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.56	152	45165	20.00	ng	-0.02
21) Naphthalene-d8	11.38	136	213895	20.00	ng	-0.02
38) Acenaphthene-d10	15.14	164	132427	20.00	ng	-0.02
63) Phenanthrene-d10	17.89	188	315237	20.00	ng	-0.03
75) Chrysene-d12	22.28	240	384211	20.00	ng	-0.03
86) Perylene-d12	25.96	264	317100	20.00	ng	-0.07

System Monitoring Compounds

5) 2-Fluorophenol	6.14	112	452884	153.01	ng	-0.02
7) Phenol-d6	7.69	99	621846	143.13	ng	0.00
23) Nitrobenzene-d5	9.72	82	405980	119.37	ng	-0.02
41) 2,4,6-Tribromophenol	16.62	330	302244	173.32	ng	-0.02
44) 2-Fluorobiphenyl	13.76	172	884515	93.44	ng	-0.03
78) Terphenyl-d14	20.46	244	1508760	100.18	ng	-0.03

Target Compounds

						Qvalue
2) 1,4-Dioxane	4.17	88	47657	38.832	ng	90
3) Pyridine	4.55	79	194035	51.288	ng	# 78
4) n-Nitrosodimethylamine	4.45	42	68103	48.618	ng	# 46
6) Aniline	7.89	93	75213	13.747	ng	96
8) 2-Chlorophenol	8.12	128	160423	51.330	ng	95
9) Benzaldehyde	7.70	77	59482	19.801	ng	92
10) Phenol	7.72	94	219713	49.451	ng	# 76
11) bis(2-Chloroethyl)ether	7.96	93	168200	46.146	ng	90
12) 1,3-Dichlorobenzene	8.45	146	141939	40.266	ng	94
13) 1,4-Dichlorobenzene	8.59	146	144331	39.896	ng	99
14) 1,2-Dichlorobenzene	8.92	146	143603	40.698	ng	97
15) Benzyl Alcohol	8.79	79	156918	51.261	ng	# 86
16) 2,2'-oxybis(1-Chloropropan	9.07	45	280583	53.799	ng	95
17) 2-Methylphenol	8.97	107	154946	49.336	ng	# 88
18) Hexachloroethane	9.64	117	53014	43.502	ng	# 83
19) n-Nitroso-di-n-propylamine	9.35	70	151032	49.692	ng	# 86
20) 3+4-Methylphenols	9.30	107	212214	48.779	ng	91
22) Acetophenone	9.37	105	259406	47.423	ng	# 87
24) Nitrobenzene	9.76	77	209995	54.195	ng	# 90
25) Isophorone	10.28	82	408012	51.017	ng	# 96
26) 2-Nitrophenol	10.47	139	85570	52.733	ng	# 83
27) 2,4-Dimethylphenol	10.51	122	177781	51.702	ng	95
28) bis(2-Chloroethoxy)methane	10.75	93	246025	47.510	ng	100
29) 2,4-Dichlorophenol	11.01	162	159316	50.752	ng	95
30) 1,2,4-Trichlorobenzene	11.23	180	149883	43.120	ng	97
31) Naphthalene	11.42	128	518731	44.374	ng	100
32) Benzoic acid	10.63	122	109002	46.192	ng	# 85
34) Hexachlorobutadiene	11.70	225	90428	42.330	ng	99
35) Caprolactam	12.29	113	54793m	50.894	ng	
36) 4-Chloro-3-methylphenol	12.60	107	209825	53.330	ng	96
37) 2-Methylnaphthalene	12.99	142	374517	43.920	ng	94
39) 1,2,4,5-Tetrachlorobenzene	13.35	216	180734	43.809	ng	97
40) Hexachlorocyclopentadiene	13.32	237	242516	110.448	ng	96
42) 2,4,6-Trichlorophenol	13.58	196	132136	52.807	ng	94

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43) 2,4,5-Trichlorophenol	13.65	196	144770	51.703	ng	96
45) 1,1'-Biphenyl	13.97	154	494596	45.757	ng	96
46) 2-Chloronaphthalene	14.02	162	376584	46.469	ng	98
47) 2-Nitroaniline	14.22	65	146567	61.377	ng	88
48) Acenaphthylene	14.86	152	614710	44.819	ng	97
49) Dimethylphthalate	14.57	163	518965	49.401	ng	97
50) 2,6-Dinitrotoluene	14.70	165	113098	52.411	ng #	82
51) Acenaphthene	15.20	154	458723	51.419	ng	98
52) 3-Nitroaniline	15.03	138	16944	12.383	ng	97
53) 2,4-Dinitrophenol	15.23	184	136187	112.207	ng	95
54) Dibenzofuran	15.53	168	612395	49.122	ng	94
55) 4-Nitrophenol	15.32	139	196869	97.741	ng #	59
56) 2,4-Dinitrotoluene	15.48	165	159684	56.960	ng #	89
57) Fluorene	16.18	166	504389	45.551	ng	98
58) 2,3,4,6-Tetrachlorophenol	15.75	232	140617	56.258	ng	96
59) Diethylphthalate	15.93	149	540082	51.543	ng	97
60) 4-Chlorophenyl-phenylether	16.16	204	258243	46.298	ng	96
61) 4-Nitroaniline	16.20	138	88870	40.310	ng #	67
62) Azobenzene	16.45	77	568167	58.504	ng	89
64) 4,6-Dinitro-2-methylphenol	16.24	198	93233	60.798	ng	84
65) n-Nitrosodiphenylamine	16.37	169	377016	38.148	ng	99
66) 4-Bromophenyl-phenylether	17.06	248	171696	46.018	ng	97
67) Hexachlorobenzene	17.19	284	185285	46.170	ng	90
68) Atrazine	17.31	200	41220	12.472	ng	96
69) Pentachlorophenol	17.52	266	241999	102.876	ng	96
70) Phenanthrene	17.93	178	830673	46.835	ng	96
71) Anthracene	18.02	178	845221	47.569	ng	98
72) Carbazole	18.28	167	548704	32.741	ng	96
73) Di-n-butylphthalate	18.81	149	953915	58.700	ng #	94
74) Fluoranthene	19.92	202	978118	51.623	ng	93
76) Benzdine	20.13	184	26842m	2.319	ng	
77) Pyrene	20.29	202	1019580	43.920	ng	97
79) Butylbenzylphthalate	21.17	149	446847	54.476	ng	94
80) Benzo(a)anthracene	22.25	228	1047549	47.810	ng	95
81) 3,3'-Dichlorobenzidine	22.15	252	32555	3.819	ng #	98
82) Chrysene	22.33	228	977698	46.477	ng	95
83) Bis(2-ethylhexyl)phthalate	22.11	149	637714	51.445	ng	99
84) Di-n-octyl phthalate	23.49	149	1083055	52.753	ng #	91
85) Indeno(1,2,3-cd)pyrene	30.21	276	1173085	46.046	ng #	94
87) Benzo(b)fluoranthene	24.77	252	1051870	53.486	ng #	96
88) Benzo(k)fluoranthene	24.86	252	1029181	53.343	ng #	94
89) Benzo(a)pyrene	25.78	252	960166	51.864	ng #	93
90) Dibenzo(a,h)anthracene	30.29	278	1025702	52.926	ng #	94
91) Benzo(g,h,i)perylene	31.54	276	925903	49.333	ng #	89

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

