

Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG051119\
 Data File : BG040905.D
 Acq On : 12 May 2019 12:16
 Operator : HP/JU
 Sample : K2768-14MS
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 P026-SS015-1824-01MS

Manual Integrations
 APPROVED

Sohil
 5/14/2019 8:01:51 PM

Quant Time: May 13 05:07:19 2019

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG042319.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Wed Apr 24 05:35:33 2019

Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.14	152	52562	20.00	ng	-0.02
21) Naphthalene-d8	10.96	136	220155	20.00	ng	-0.03
39) Acenaphthene-d10	14.76	164	160913	20.00	ng	-0.03
64) Phenanthrene-d10	17.51	188	400285	20.00	ng	-0.03
76) Chrysene-d12	21.80	240	401358	20.00	ng	-0.03
87) Perylene-d12	25.14	264	445586	20.00	ng	-0.06

System Monitoring Compounds

5) 2-Fluorophenol	5.67	112	414845	139.13	ng	-0.02
7) Phenol-d6	7.28	99	620822	135.22	ng	-0.02
23) Nitrobenzene-d5	9.31	82	419749	88.10	ng	-0.03
42) 2,4,6-Tribromophenol	16.25	330	314544	142.21	ng	-0.02
45) 2-Fluorobiphenyl	13.39	172	888812	79.77	ng	-0.03
79) Terphenyl-d14	20.10	244	1517722	83.78	ng	-0.03

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.53	88	42705	31.492	ng	92
3) Pyridine	3.94	79	86450	21.994	ng	98
4) n-Nitrosodimethylamine	3.86	42	73633	33.774	ng	93
6) Aniline	7.45	93	86868	14.275	ng	99
8) 2-Chlorophenol	7.69	128	153996	42.296	ng	98
9) Benzaldehyde	7.27	77	82637	27.126	ng	96
10) Phenol	7.30	94	214573	43.478	ng	96
11) bis(2-Chloroethyl)ether	7.55	93	142356	38.466	ng	97
12) 1,3-Dichlorobenzene	8.02	146	155975	39.249	ng	96
13) 1,4-Dichlorobenzene	8.17	146	159899	39.691	ng	99
14) 1,2-Dichlorobenzene	8.49	146	153012	39.384	ng	99
15) Benzyl Alcohol	8.37	79	178576	37.382	ng	95
16) 2,2'-oxybis(1-Chloropropan	8.65	45	234110	31.773	ng	91
17) 2-Methylphenol	8.56	107	150250	43.020	ng	94
18) Hexachloroethane	9.22	117	60714	36.740	ng	98
19) n-Nitroso-di-n-propylamine	8.94	70	142721	34.534	ng	91
20) 3+4-Methylphenols	8.89	107	208390	42.831	ng	98
22) Acetophenone	8.96	105	254055	41.499	ng	# 96
24) Nitrobenzene	9.36	77	214584	42.211	ng	97
25) Isophorone	9.87	82	407125	38.360	ng	99
26) 2-Nitrophenol	10.06	139	95055	52.164	ng	94
27) 2,4-Dimethylphenol	10.10	122	161965	47.372	ng	96
28) bis(2-Chloroethoxy)methane	10.35	93	214150	40.217	ng	97
29) 2,4-Dichlorophenol	10.59	162	187840	48.325	ng	95
30) 1,2,4-Trichlorobenzene	10.81	180	194005	45.008	ng	97
31) Naphthalene	11.01	128	484051	41.385	ng	98
32) Benzoic acid	10.23	122	89951	38.431	ng	95
33) 4-Chloroaniline	11.11	127	54819	10.571	ng	95
34) Hexachlorobutadiene	11.27	225	137362	43.165	ng	98
35) Caprolactam	11.90	113	56632m	36.903	ng	
36) 4-Chloro-3-methylphenol	12.22	107	218585	44.577	ng	95
37) 2-Methylnaphthalene	12.60	142	367214	41.992	ng	98
38) 1-Methylnaphthalene	12.82	142	353577	41.365	ng	98
40) 1,2,4,5-Tetrachlorobenzene	12.96	216	264412	44.110	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.93	237	324570	91.759	ng	99
43) 2,4,6-Trichlorophenol	13.19	196	181030	49.974	ng	99
44) 2,4,5-Trichlorophenol	13.27	196	193372	49.003	ng	96
46) 1,1'-Biphenyl	13.60	154	494446	39.577	ng	96
47) 2-Chloronaphthalene	13.65	162	409986	41.933	ng	98
48) 2-Nitroaniline	13.85	65	154142	44.290	ng	92
49) Acenaphthylene	14.49	152	645574	38.918	ng	99
50) Dimethylphthalate	14.21	163	716202	53.197	ng	99
51) 2,6-Dinitrotoluene	14.34	165	125326	47.696	ng	89
52) Acenaphthene	14.83	154	380376	39.375	ng	96
53) 3-Nitroaniline	14.67	138	66681	24.767	ng	# 91
54) 2,4-Dinitrophenol	14.87	184	159608	105.272	ng	92
55) Dibenzofuran	15.16	168	641252	40.527	ng	99
56) 4-Nitrophenol	14.97	139	207019	88.677	ng	# 84
57) 2,4-Dinitrotoluene	15.13	165	175388	49.121	ng	90
58) Fluorene	15.81	166	500413	39.128	ng	100
59) 2,3,4,6-Tetrachlorophenol	15.38	232	199559	50.243	ng	94
60) Diethylphthalate	15.56	149	557759	39.265	ng	100
61) 4-Chlorophenyl-phenylether	15.80	204	320107	43.035	ng	95
62) 4-Nitroaniline	15.84	138	120749	40.094	ng	92
63) Azobenzene	16.08	77	518080	35.415	ng	97
65) 4,6-Dinitro-2-methylphenol	15.88	198	106994	53.324	ng	91
66) n-Nitrosodiphenylamine	16.01	169	460182	39.075	ng	98
67) 4-Bromophenyl-phenylether	16.69	248	208093	41.727	ng	92
68) Hexachlorobenzene	16.80	284	213369	41.378	ng	96
69) Atrazine	16.95	200	207851	44.547	ng	95
70) Pentachlorophenol	17.15	266	285005	94.452	ng	97
71) Phenanthrene	17.55	178	859823	39.880	ng	99
72) Anthracene	17.64	178	875311	40.389	ng	100
73) Carbazole	17.91	167	789117	39.966	ng	99
74) Di-n-butylphthalate	18.45	149	899917	37.762	ng	99
75) Fluoranthene	19.55	202	1100041	40.943	ng	99
77) Benzidine	19.73	184	135575	12.337	ng	98
78) Pyrene	19.92	202	1127640	44.827	ng	98
80) Butylbenzylphthalate	20.78	149	416783	42.510	ng	94
81) Benzo(a)anthracene	21.77	228	1065651	42.011	ng	99
82) 3,3'-Dichlorobenzidine	21.68	252	227879	25.205	ng	98
83) Chrysene	21.84	228	1045420	43.335	ng	100
84) Bis(2-ethylhexyl)phthalate	21.64	149	590265	41.707	ng	96
85) Di-n-octyl phthalate	22.89	149	1010903	42.412	ng	96
86) Indeno(1,2,3-cd)pyrene	28.99	276	1303227	44.682	ng	# 92
88) Benzo(b)fluoranthene	24.07	252	1113744	40.903	ng	98
89) Benzo(k)fluoranthene	24.14	252	1049936	41.288	ng	# 98
90) Benzo(a)pyrene	24.98	252	1073096	41.634	ng	# 99
91) Dibenzo(a,h)anthracene	29.06	278	1062158	40.722	ng	97
92) Benzo(g,h,i)perylene	30.22	276	1076168	41.457	ng	99

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

