

Data File : BG041026.D
Acq On : 3 Jun 2019 9:55
Operator : HP/JU
Sample : K3084-01MSD
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
TEST-PIT-6MSD

Manual Integrations
APPROVED

mohammad
6/5/2019 8:35:17 AM

Quant Time: Jun 03 15:45:23 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG052919.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed May 29 18:39:57 2019
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.12	152	60661	20.00	ng	-0.01
21) Naphthalene-d8	10.94	136	245042	20.00	ng	0.00
39) Acenaphthene-d10	14.75	164	177664	20.00	ng	0.00
64) Phenanthrene-d10	17.50	188	449288	20.00	ng	0.00
76) Chrysene-d12	21.78	240	443484	20.00	ng	0.00
87) Perylene-d12	25.11	264	470863	20.00	ng	-0.02

System Monitoring Compounds

5) 2-Fluorophenol	5.66	112	488096	145.09	ng	0.00
7) Phenol-d6	7.27	99	710603	136.77	ng	0.00
23) Nitrobenzene-d5	9.30	82	485288	87.92	ng	0.00
42) 2,4,6-Tribromophenol	16.24	330	364877	138.34	ng	0.00
45) 2-Fluorobiphenyl	13.38	172	995818	80.72	ng	0.00
79) Terphenyl-d14	20.09	244	1633105	78.15	ng	-0.01

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.52	88	55969	36.664	ng	95
3) Pyridine	3.92	79	152575	33.778	ng	95
4) n-Nitrosodimethylamine	3.85	42	95078	45.353	ng	94
6) Aniline	7.45	93	118973	17.156	ng	98
8) 2-Chlorophenol	7.68	128	186230	46.435	ng	94
9) Benzaldehyde	7.26	77	87564	28.253	ng	92
10) Phenol	7.30	94	266044	47.759	ng	99
11) bis(2-Chloroethyl)ether	7.54	93	174230	43.114	ng	98
12) 1,3-Dichlorobenzene	8.01	146	195219	40.755	ng	96
13) 1,4-Dichlorobenzene	8.16	146	204773	42.233	ng	94
14) 1,2-Dichlorobenzene	8.48	146	196321	41.701	ng	99
15) Benzyl Alcohol	8.36	79	211721	44.054	ng	95
16) 2,2'-oxybis(1-Chloropropan	8.64	45	276322	41.845	ng	97
17) 2-Methylphenol	8.55	107	179218	44.434	ng	100
18) Hexachloroethane	9.21	117	76665	41.025	ng	98
19) n-Nitroso-di-n-propylamine	8.93	70	167934	42.003	ng	98
20) 3+4-Methylphenols	8.88	107	240125	42.980	ng	98
22) Acetophenone	8.95	105	310951	45.237	ng	# 97
24) Nitrobenzene	9.34	77	260704	46.347	ng	99
25) Isophorone	9.86	82	471834	44.917	ng	99
26) 2-Nitrophenol	10.05	139	114002	49.161	ng	90
27) 2,4-Dimethylphenol	10.09	122	192367	50.228	ng	93
28) bis(2-Chloroethoxy)methane	10.34	93	249651	44.034	ng	97
29) 2,4-Dichlorophenol	10.58	162	223986	46.055	ng	97
30) 1,2,4-Trichlorobenzene	10.80	180	240040	43.386	ng	97
31) Naphthalene	11.00	128	588416	44.724	ng	97
32) Benzoic acid	10.16	122	12047	11.490	ng	90
33) 4-Chloroaniline	11.10	127	82228	13.470	ng	96
34) Hexachlorobutadiene	11.26	225	176370	41.662	ng	98
35) Caprolactam	11.89	113	67779m	42.202	ng	
36) 4-Chloro-3-methylphenol	12.21	107	245882	44.268	ng	97
37) 2-Methylnaphthalene	12.59	142	435591	42.988	ng	98
38) 1-Methylnaphthalene	12.81	142	413483	43.303	ng	96
40) 1,2,4,5-Tetrachlorobenzene	12.95	216	318295	44.449	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.92	237	312552	81.074	ng	97
43) 2,4,6-Trichlorophenol	13.19	196	211825	45.727	ng	95
44) 2,4,5-Trichlorophenol	13.26	196	222905	45.626	ng	98
46) 1,1'-Biphenyl	13.59	154	582585	44.300	ng	99
47) 2-Chloronaphthalene	13.64	162	486227	43.067	ng	99
48) 2-Nitroaniline	13.84	65	180570	44.582	ng	98
49) Acenaphthylene	14.48	152	758531	44.640	ng	98
50) Dimethylphthalate	14.21	163	801323	53.282	ng	98
51) 2,6-Dinitrotoluene	14.33	165	147925	45.619	ng	95
52) Acenaphthene	14.82	154	443140	43.050	ng	97
53) 3-Nitroaniline	14.66	138	73869	22.782	ng	97
54) 2,4-Dinitrophenol	14.86	184	106714	67.861	ng	97
55) Dibenzofuran	15.15	168	763020	44.053	ng	99
56) 4-Nitrophenol	14.96	139	226224	101.717	ng	94
57) 2,4-Dinitrotoluene	15.12	165	210476	45.951	ng	# 98
58) Fluorene	15.80	166	595532	43.174	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.37	232	228033	47.495	ng	99
60) Diethylphthalate	15.55	149	657161	42.817	ng	98
61) 4-Chlorophenyl-phenylether	15.78	204	379574	42.336	ng	97
62) 4-Nitroaniline	15.83	138	141021	41.081	ng	96
63) Azobenzene	16.07	77	607031	43.516	ng	99
65) 4,6-Dinitro-2-methylphenol	15.87	198	115508	47.981	ng	96
66) n-Nitrosodiphenylamine	16.00	169	554343	43.891	ng	98
67) 4-Bromophenyl-phenylether	16.68	248	247693	40.851	ng	95
68) Hexachlorobenzene	16.79	284	260865	40.404	ng	98
69) Atrazine	16.94	200	245045	50.283	ng	99
70) Pentachlorophenol	17.14	266	318276	91.381	ng	97
71) Phenanthrene	17.54	178	1001928	42.812	ng	98
72) Anthracene	17.63	178	1032498	44.733	ng	100
73) Carbazole	17.90	167	923101	46.610	ng	99
74) Di-n-butylphthalate	18.43	149	1050257	42.667	ng	97
75) Fluoranthene	19.54	202	1289537	44.611	ng	98
77) Benzidine	19.72	184	297085	28.081	ng	98
78) Pyrene	19.90	202	1278802	44.358	ng	99
80) Butylbenzylphthalate	20.77	149	500949	44.651	ng	99
81) Benzo(a)anthracene	21.75	228	1260476	42.698	ng	99
82) 3,3'-Dichlorobenzidine	21.67	252	227265	21.888	ng	97
83) Chrysene	21.83	228	1223159	43.297	ng	97
84) Bis(2-ethylhexyl)phthalate	21.63	149	681766	43.168	ng	98
85) Di-n-octyl phthalate	22.87	149	1173825	44.627	ng	99
86) Indeno(1,2,3-cd)pyrene	28.95	276	1526547	44.112	ng	# 98
88) Benzo(b)fluoranthene	24.05	252	1293166	45.280	ng	98
89) Benzo(k)fluoranthene	24.12	252	1251332	44.277	ng	99
90) Benzo(a)pyrene	24.96	252	1256103	46.099	ng	98
91) Dibenzo(a,h)anthracene	29.02	278	1233547	45.307	ng	99
92) Benzo(g,h,i)perylene	30.17	276	1242470	46.972	ng	99

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

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