

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG051220\
 Data File : BG045353.D
 Acq On : 13 May 2020 8:23
 Operator : CG/JU
 Sample : L2575-02MSD
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 TP-28MSD

Manual Integrations
 APPROVED

mohammad
 5/14/2020 11:54:09 AM

Quant Time: May 13 09:03:57 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG041520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue May 12 15:48:20 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.97	152	63388	20.00	ng	0.00
21) Naphthalene-d8	10.78	136	259601	20.00	ng	0.00
39) Acenaphthene-d10	14.61	164	192052	20.00	ng	0.00
64) Phenanthrene-d10	17.36	188	456286	20.00	ng	0.00
76) Chrysene-d12	21.63	240	420013	20.00	ng	0.00
87) Perylene-d12	24.79	264	474151	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.54	112	449837	117.16	ng	0.00
7) Phenol-d6	7.14	99	652034	114.30	ng	0.00
23) Nitrobenzene-d5	9.13	82	405251	71.23	ng	0.00
42) 2,4,6-Tribromophenol	16.10	330	318349	107.30	ng	0.00
45) 2-Fluorobiphenyl	13.24	172	950217	72.55	ng	0.00
79) Terphenyl-d14	19.98	244	1681258	87.25	ng	0.00

Target Compounds					Qvalue	
2) 1,4-Dioxane	3.42	88	57664	33.794	ng	95
3) Pyridine	3.82	79	136936	26.064	ng	93
4) n-Nitrosodimethylamine	3.73	42	60791	37.772	ng	100
6) Aniline	7.29	93	142877	19.263	ng	98
8) 2-Chlorophenol	7.54	128	168769	40.900	ng	98
9) Benzaldehyde	7.10	77	128560	39.986	ng	99
10) Phenol	7.17	94	231634	40.578	ng	98
11) bis(2-Chloroethyl)ether	7.39	93	155929	34.308	ng	99
12) 1,3-Dichlorobenzene	7.86	146	176425	36.283	ng	98
13) 1,4-Dichlorobenzene	8.01	146	179700	37.089	ng	98
14) 1,2-Dichlorobenzene	8.33	146	167047	36.038	ng	94
15) Benzyl Alcohol	8.21	79	173981	36.377	ng	96
16) 2,2'-oxybis(1-Chloropropan	8.50	45	141447	31.705	ng	98
17) 2-Methylphenol	8.42	107	154460	35.070	ng	96
18) Hexachloroethane	9.06	117	70205	38.544	ng	97
19) n-Nitroso-di-n-propylamine	8.78	70	140881	34.921	ng	98
20) 3+4-Methylphenols	8.75	107	215872	35.017	ng	99
22) Acetophenone	8.79	105	262960	37.457	ng	98
24) Nitrobenzene	9.17	77	209934	35.998	ng	# 95
25) Isophorone	9.70	82	396011	33.989	ng	99
26) 2-Nitrophenol	9.89	139	97747	38.911	ng	96
27) 2,4-Dimethylphenol	9.95	122	156477	39.258	ng	96
28) bis(2-Chloroethoxy)methane	10.18	93	219696	35.888	ng	97
29) 2,4-Dichlorophenol	10.43	162	178792	38.847	ng	99
30) 1,2,4-Trichlorobenzene	10.65	180	186571	35.803	ng	97
31) Naphthalene	10.83	128	516160	36.346	ng	99
32) Benzoic acid	10.11	122	95046	31.972	ng	97
33) 4-Chloroaniline	10.94	127	66850	10.094	ng	98
34) Hexachlorobutadiene	11.13	225	130290	36.468	ng	97
35) Caprolactam	11.70	113	65882m	35.966	ng	
36) 4-Chloro-3-methylphenol	12.07	107	205423	37.994	ng	98
37) 2-Methylnaphthalene	12.44	142	392010	35.563	ng	98
38) 1-Methylnaphthalene	12.66	142	375381	36.073	ng	97
40) 1,2,4,5-Tetrachlorobenzene	12.81	216	216728	35.049	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.80	237	274167	70.939	nq	97
43) 2,4,6-Trichlorophenol	13.05	196	158270	36.266	nq	90
44) 2,4,5-Trichlorophenol	13.13	196	174189	35.065	nq	96
46) 1,1'-Biphenyl	13.45	154	512527	35.111	nq	100
47) 2-Chloronaphthalene	13.49	162	389571	34.927	nq	96
48) 2-Nitroaniline	13.68	65	130074	35.444	nq	98
49) Acenaphthylene	14.34	152	665167	36.612	nq	99
50) Dimethylphthalate	14.07	163	682164	43.623	nq	98
51) 2,6-Dinitrotoluene	14.18	165	125298	35.823	nq	96
52) Acenaphthene	14.68	154	499803m	40.660	nq	
53) 3-Nitroaniline	14.51	138	54174	14.122	nq	# 91
54) 2,4-Dinitrophenol	14.72	184	157140	75.554	nq	92
55) Dibenzofuran	15.02	168	639484	35.683	nq	99
56) 4-Nitrophenol	14.84	139	213036	71.623	nq	94
57) 2,4-Dinitrotoluene	14.98	165	180302	35.273	nq	96
58) Fluorene	15.66	166	526981	36.000	nq	100
59) 2,3,4,6-Tetrachlorophenol	15.25	232	162564	36.779	nq	98
60) Diethylphthalate	15.44	149	569189	34.911	nq	99
61) 4-Chlorophenyl-phenylether	15.66	204	291597	34.608	nq	98
62) 4-Nitroaniline	15.68	138	122470	30.780	nq	93
63) Azobenzene	15.95	77	537464	35.756	nq	98
65) 4,6-Dinitro-2-methylphenol	15.74	198	117030	39.021	nq	92
66) n-Nitrosodiphenylamine	15.87	169	474095	36.163	nq	99
67) 4-Bromophenyl-phenylether	16.55	248	203843	34.629	nq	99
68) Hexachlorobenzene	16.67	284	220646	36.142	nq	99
69) Atrazine	16.82	200	184287	37.773	nq	95
70) Pentachlorophenol	17.02	266	264662	70.668	nq	96
71) Phenanthrene	17.40	178	871246	37.158	nq	99
72) Anthracene	17.50	178	892914	37.964	nq	100
73) Carbazole	17.76	167	823825	34.515	nq	97
74) Di-n-butylphthalate	18.33	149	998253	37.395	nq	99
75) Fluoranthene	19.42	202	1099065	39.066	nq	98
77) Benzidine	19.59	184	484168	39.006	nq	94
78) Pyrene	19.78	202	1085821	37.499	nq	98
80) Butylbenzylphthalate	20.66	149	451624	37.627	nq	99
81) Benzo(a)anthracene	21.61	228	1041793	38.269	nq	98
82) 3,3'-Dichlorobenzidine	21.52	252	174048	16.063	nq	95
83) Chrysene	21.67	228	983670	37.608	nq	98
84) Bis(2-ethylhexyl)phthalate	21.52	149	624473	37.829	nq	97
85) Di-n-octyl phthalate	22.71	149	1064621	37.295	nq	100
86) Indeno(1,2,3-cd)pyrene	28.38	276	1318392	37.964	nq	# 97
88) Benzo(b)fluoranthene	23.79	252	1091682m	36.756	nq	
89) Benzo(k)fluoranthene	23.85	252	1079787	38.229	nq	97
90) Benzo(a)pyrene	24.64	252	1019329	36.350	nq	98
91) Dibenzo(a,h)anthracene	28.44	278	1075427	38.060	nq	97
92) Benzo(g,h,i)perylene	29.50	276	1085924	38.333	nq	99

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

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