

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM050619\
 Data File : BM020145.D
 Acq On : 06 May 2019 20:50
 Operator : JU/SJ
 Sample : K2673-16MSD
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SP-02-COMPMSD

Manual Integrations
 APPROVED

mohammad
 5/7/2019 3:11:47 PM

Quant Time: May 07 07:02:26 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM043019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 01 03:58:37 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.79	152	175423	20.00	ng	0.00
21) Naphthalene-d8	10.59	136	587049	20.00	ng	0.00
39) Acenaphthene-d10	14.44	164	309548	20.00	ng	0.00
64) Phenanthrene-d10	17.19	188	682167	20.00	ng	0.00
76) Chrysene-d12	21.37	240	714542	20.00	ng	0.00
87) Perylene-d12	23.67	264	841964	20.00	ng	0.02

System Monitoring Compounds

5) 2-Fluorophenol	5.38	112	1797356	167.48	ng	0.00
7) Phenol-d6	6.97	99	2435555	166.12	ng	0.02
23) Nitrobenzene-d5	8.96	82	1633832	109.88	ng	0.00
42) 2,4,6-Tribromophenol	15.93	330	789994	164.42	ng	0.01
45) 2-Fluorobiphenyl	13.06	172	2509620	99.75	ng	0.00
79) Terphenyl-d14	19.81	244	3425840	83.63	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.30	88	203242	38.614	ng	93
3) Pyridine	3.70	79	625453	46.463	ng	96
4) n-Nitrosodimethylamine	3.62	42	222927	51.643	ng	# 77
6) Aniline	7.13	93	545805	29.574	ng	100
8) 2-Chlorophenol	7.36	128	632839	54.156	ng	98
9) Benzaldehyde	6.95	77	447493	40.327	ng	97
10) Phenol	7.00	94	892190	56.527	ng	93
11) bis(2-Chloroethyl)ether	7.23	93	647083	56.407	ng	97
12) 1,3-Dichlorobenzene	7.68	146	684714	50.362	ng	97
13) 1,4-Dichlorobenzene	7.83	146	705079	51.336	ng	98
14) 1,2-Dichlorobenzene	8.15	146	665688	50.874	ng	99
15) Benzyl Alcohol	8.04	79	695519	49.197	ng	97
16) 2,2'-oxybis(1-Chloropropan	8.32	45	469964	49.406	ng	98
17) 2-Methylphenol	8.24	107	545079	53.648	ng	95
18) Hexachloroethane	8.86	117	161511	28.203	ng	97
19) n-Nitroso-di-n-propylamine	8.60	70	610816	48.696	ng	96
20) 3+4-Methylphenols	8.57	107	717261	53.117	ng	97
22) Acetophenone	8.62	105	992756	61.877	ng	# 97
24) Nitrobenzene	9.00	77	857008	55.588	ng	96
25) Isophorone	9.52	82	1472055	57.408	ng	99
26) 2-Nitrophenol	9.71	139	182632	39.238	ng	92
27) 2,4-Dimethylphenol	9.76	122	534431	68.968	ng	99
28) bis(2-Chloroethoxy)methane	10.00	93	818084	58.578	ng	99
29) 2,4-Dichlorophenol	10.24	162	580144	59.373	ng	99
30) 1,2,4-Trichlorobenzene	10.45	180	690707	57.449	ng	98
31) Naphthalene	10.64	128	1731462	56.836	ng	99
32) Benzoic acid	9.86	122	58278	15.725	ng	# 89
33) 4-Chloroaniline	10.76	127	412837	32.925	ng	97
34) Hexachlorobutadiene	10.90	225	454891	53.484	ng	97
35) Caprolactam	11.56	113	156141m	53.442	ng	
36) 4-Chloro-3-methylphenol	11.88	107	595698	51.878	ng	95
37) 2-Methylnaphthalene	12.25	142	1203060	54.169	ng	96
38) 1-Methylnaphthalene	12.47	142	1140142	52.498	ng	96
40) 1,2,4,5-Tetrachlorobenzene	12.62	216	738845	61.008	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.59	237	29731	4.169	ng	95
43) 2,4,6-Trichlorophenol	12.87	196	470367	68.339	ng	98
44) 2,4,5-Trichlorophenol	12.94	196	491309	63.897	ng	98
46) 1,1'-Biphenyl	13.27	154	1504591	59.053	ng	98
47) 2-Chloronaphthalene	13.32	162	1188081	58.403	ng	99
48) 2-Nitroaniline	13.53	65	479526	66.212	ng	93
49) Acenaphthylene	14.16	152	1853866	56.943	ng	99
50) Dimethylphthalate	13.90	163	1814792	68.979	ng	99
51) 2,6-Dinitrotoluene	14.03	165	238395	43.022	ng	92
52) Acenaphthene	14.50	154	1044893	55.967	ng	99
53) 3-Nitroaniline	14.36	138	306755	59.646	ng	96
54) 2,4-Dinitrophenol	14.56	184	2316	9.218	ng	91
55) Dibenzofuran	14.84	168	1710199	54.307	ng	99
56) 4-Nitrophenol	14.67	139	455001	103.692	ng	93
57) 2,4-Dinitrotoluene	14.82	165	281702	37.413	ng	94
58) Fluorene	15.49	166	1359226	52.555	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.06	232	442047	60.497	ng	97
60) Diethylphthalate	15.26	149	1418418	53.513	ng	100
61) 4-Chlorophenyl-phenylether	15.48	204	776006	52.955	ng	97
62) 4-Nitroaniline	15.53	138	337013	59.378	ng	94
63) Azobenzene	15.77	77	1498427	47.919	ng	99
65) 4,6-Dinitro-2-methylphenol	15.57	198	2776	7.848	ng	# 50
66) n-Nitrosodiphenylamine	15.70	169	1187144	59.141	ng	100
67) 4-Bromophenyl-phenylether	16.38	248	468271	55.979	ng	95
68) Hexachlorobenzene	16.49	284	533298	55.401	ng	96
69) Atrazine	16.65	200	345649	44.063	ng	97
70) Pentachlorophenol	16.83	266	688879	125.076	ng	98
71) Phenanthrene	17.23	178	2290525	60.827	ng	99
72) Anthracene	17.32	178	2254549	59.883	ng	100
73) Carbazole	17.60	167	1864605	54.887	ng	99
74) Di-n-butylphthalate	18.14	149	2161635	52.874	ng	98
75) Fluoranthene	19.24	202	3224321	67.674	ng	100
77) Benzidine	19.44	184	81862	3.583	ng	# 93
78) Pyrene	19.61	202	3245219	67.646	ng	99
80) Butylbenzylphthalate	20.50	149	1022501	58.611	ng	97
81) Benzo(a)anthracene	21.36	228	3153526	62.931	ng	98
82) 3,3'-Dichlorobenzidine	21.29	252	753687	39.407	ng	99
83) Chrysene	21.41	228	2932654	60.344	ng	99
84) Bis(2-ethylhexyl)phthalate	21.27	149	2526783	93.339	ng	98
85) Di-n-octyl phthalate	22.17	149	2807858	62.405	ng	# 99
86) Indeno(1,2,3-cd)pyrene	26.03	276	3725212	64.442	ng	# 100
88) Benzo(b)fluoranthene	22.98	252	3553047	63.905	ng	99
89) Benzo(k)fluoranthene	23.03	252	2977074	58.700	ng	99
90) Benzo(a)pyrene	23.57	252	3232316	64.004	ng	98
91) Dibenzo(a,h)anthracene	26.03	278	2900874	55.234	ng	99
92) Benzo(g,h,i)perylene	26.75	276	3145142	63.076	ng	99

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

