

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM070618\
 Data File : BM015813.D
 Acq On : 06 Jul 2018 21:58
 Operator : SJ/JU
 Sample : J3845-03MS
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 GF-02MS

Manual Integrations
 APPROVED

Sohil
 7/9/2018 1:55:48 PM

Quant Time: Jul 07 02:37:36 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM070518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 05 17:39:50 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.30	152	126085	20.00	ng	-0.01
21) Naphthalene-d8	11.13	136	516671	20.00	ng	-0.01
38) Acenaphthene-d10	14.93	164	280133	20.00	ng	0.00
63) Phenanthrene-d10	17.65	188	530330	20.00	ng	0.00
75) Chrysene-d12	21.76	240	389094	20.00	ng	0.00
86) Perylene-d12	24.16	264	356066	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.81	112	903744	122.92	ng	0.00
7) Phenol-d6	7.46	99	1178764	117.11	ng	0.00
23) Nitrobenzene-d5	9.50	82	811011	76.59	ng	0.00
41) 2,4,6-Tribromophenol	16.41	330	409151	111.78	ng	0.00
44) 2-Fluorobiphenyl	13.55	172	1654018	73.30	ng	-0.01
78) Terphenyl-d14	20.21	244	1716992	81.85	ng	-0.01

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.56	88	84475	27.782	ng	# 100
3) Pyridine	4.01	79	182564	22.753	ng	# 77
4) n-Nitrosodimethylamine	3.92	42	207914	33.279	ng	# 43
6) Aniline	7.63	93	301760	24.729	ng	92
8) 2-Chlorophenol	7.87	128	327968	37.169	ng	97
9) Benzaldehyde	7.44	77	157460	24.821	ng	90
10) Phenol	7.49	94	407990	40.342	ng	93
11) bis(2-Chloroethyl)ether	7.72	93	273039	33.188	ng	88
12) 1,3-Dichlorobenzene	8.19	146	330102	33.742	ng	# 92
13) 1,4-Dichlorobenzene	8.34	146	339305	33.539	ng	# 94
14) 1,2-Dichlorobenzene	8.66	146	334083	34.146	ng	96
15) Benzyl Alcohol	8.56	79	297221	34.014	ng	89
16) 2,2'-oxybis(1-Chloropropan	8.83	45	432425	48.128	ng	97
17) 2-Methylphenol	8.75	107	260142	37.128	ng	# 85
18) Hexachloroethane	9.39	117	138430	34.605	ng	95
19) n-Nitroso-di-n-propylamine	9.12	70	250884	31.015	ng	# 75
20) 3+4-Methylphenols	9.09	107	349708	35.932	ng	89
22) Acetophenone	9.15	105	477638	35.487	ng	# 87
24) Nitrobenzene	9.54	77	389461	37.783	ng	98
25) Isophorone	10.06	82	646722	40.013	ng	# 94
26) 2-Nitrophenol	10.25	139	167672	37.775	ng	# 63
27) 2,4-Dimethylphenol	10.29	122	285939	42.405	ng	# 87
28) bis(2-Chloroethoxy)methane	10.53	93	383892	36.457	ng	98
29) 2,4-Dichlorophenol	10.78	162	300203	40.198	ng	96
30) 1,2,4-Trichlorobenzene	10.99	180	318339	36.700	ng	96
31) Naphthalene	11.19	128	973923	36.358	ng	99
32) Benzoic acid	10.44	122	91398m	15.309	ng	
33) 4-Chloroaniline	11.30	127	260311	23.835	ng	# 90
34) Hexachlorobutadiene	11.45	225	205205	35.448	ng	98
35) Caprolactam	12.10	113	75323	30.412	ng	90
36) 4-Chloro-3-methylphenol	12.40	107	321025	37.463	ng	88
37) 2-Methylnaphthalene	12.77	142	708667	37.190	ng	# 94
39) 1,2,4,5-Tetrachlorobenzene	13.13	216	348590	38.771	ng	# 100
40) Hexachlorocyclopentadiene	13.10	237	366177	86.368	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.37	196	226965	39.935	ng	99
43) 2,4,5-Trichlorophenol	13.45	196	235248	38.318	ng	# 85
45) 1,1'-Biphenyl	13.76	154	887200	36.631	ng	98
46) 2-Chloronaphthalene	13.82	162	683639	37.752	ng	# 90
47) 2-Nitroaniline	14.03	65	225684	35.738	ng	# 80
48) Acenaphthylene	14.66	152	1090118	37.005	ng	99
49) Dimethylphthalate	14.38	163	1137997	46.062	ng	99
50) 2,6-Dinitrotoluene	14.52	165	179299	37.222	ng	# 75
51) Acenaphthene	14.99	154	625372	34.116	ng	98
52) 3-Nitroaniline	14.84	138	135764	25.749	ng	# 78
53) 2,4-Dinitrophenol	15.06	184	142275	60.856	ng	93
54) Dibenzofuran	15.32	168	994685	35.215	ng	# 88
55) 4-Nitrophenol	15.14	139	257000	66.046	ng	# 69
56) 2,4-Dinitrotoluene	15.30	165	238053	34.230	ng	93
57) Fluorene	15.97	166	785233	33.711	ng	97
58) 2,3,4,6-Tetrachlorophenol	15.54	232	195421	37.721	ng	# 100
59) Diethylphthalate	15.72	149	867214	32.756	ng	96
60) 4-Chlorophenyl-phenylether	15.94	204	395017	32.957	ng	# 88
61) 4-Nitroaniline	16.00	138	152176	27.822	ng	# 36
62) Azobenzene	16.24	77	898367	36.199	ng	80
64) 4,6-Dinitro-2-methylphenol	16.06	198	111252	34.507	ng	89
65) n-Nitrosodiphenylamine	16.17	169	659872	36.313	ng	99
66) 4-Bromophenyl-phenylether	16.84	248	228727	38.628	ng	# 85
67) Hexachlorobenzene	16.96	284	247357	37.384	ng	# 78
68) Atrazine	17.10	200	227654	38.087	ng	98
69) Pentachlorophenol	17.30	266	249626	60.970	ng	98
70) Phenanthrene	17.70	178	1075062	35.433	ng	99
71) Anthracene	17.79	178	1086814	36.475	ng	99
72) Carbazole	18.05	167	933303	33.614	ng	98
73) Di-n-butylphthalate	18.57	149	1228019	32.814	ng	# 98
74) Fluoranthene	19.67	202	1038751	29.781	ng	99
76) Benzidine	19.85	184	204334	17.924	ng	98
77) Pyrene	20.03	202	1013488	41.815	ng	99
79) Butylbenzylphthalate	20.88	149	443566	37.860	ng	# 84
80) Benzo(a)anthracene	21.74	228	878591	35.614	ng	99
81) 3,3'-Dichlorobenzidine	21.67	252	232413	26.221	ng	100
82) Chrysene	21.80	228	816269	35.519	ng	100
83) Bis(2-ethylhexyl)phthalate	21.63	149	594156	34.322	ng	96
84) Di-n-octyl phthalate	22.54	149	924324	32.746	ng	# 90
85) Indeno(1,2,3-cd)pyrene	26.64	276	919855	43.082	ng	# 93
87) Benzo(b)fluoranthene	23.43	252	816866	34.362	ng	# 96
88) Benzo(k)fluoranthene	23.48	252	759529	32.890	ng	97
89) Benzo(a)pyrene	24.06	252	758941	36.267	ng	99
90) Dibenzo(a,h)anthracene	26.64	278	793994	42.171	ng	97
91) Benzo(g,h,i)perylene	27.40	276	742094	44.858	ng	# 94

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

