

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM082118\
 Data File : BM016502.D
 Acq On : 22 Aug 2018 02:19
 Operator : SJ/JU
 Sample : J4539-05MS
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 MS-RO-27186MS

Manual Integrations
 APPROVED

Sohil
 8/22/2018 6:07:03 PM

Quant Time: Aug 22 04:08:19 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM082118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 21 16:31:22 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.02	152	195217	20.00	ng	0.00
21) Naphthalene-d8	10.83	136	678834	20.00	ng	0.00
38) Acenaphthene-d10	14.66	164	571212	20.00	ng	0.00
63) Phenanthrene-d10	17.40	188	1707349	20.00	ng	0.00
75) Chrysene-d12	21.54	240	2003760	20.00	ng	0.00
86) Perylene-d12	23.83	264	1851598	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.57	112	1389308	139.46	ng	0.01
7) Phenol-d6	7.20	99	1591469	120.72	ng	0.00
23) Nitrobenzene-d5	9.21	82	1729102	116.30	ng	0.00
41) 2,4,6-Tribromophenol	16.16	330	3295288	206.37	ng	0.00
44) 2-Fluorobiphenyl	13.28	172	7520617	126.56	ng	0.00
78) Terphenyl-d14	19.99	244	11908924	125.79	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.40	88	181123	32.910	ng	# 100
3) Pyridine	3.84	79	273064	23.847	ng	# 83
4) n-Nitrosodimethylamine	3.75	42	230711	45.226	ng	# 68
6) Aniline	7.36	93	209049	14.171	ng	92
8) 2-Chlorophenol	7.59	128	527133	44.468	ng	98
9) Benzaldehyde	7.19	77	106947	11.690	ng	84
10) Phenol	7.23	94	456979	35.052	ng	90
11) bis(2-Chloroethyl)ether	7.46	93	407673	41.701	ng	98
12) 1,3-Dichlorobenzene	7.90	146	638892	40.651	ng	# 86
13) 1,4-Dichlorobenzene	8.06	146	647930	40.172	ng	98
14) 1,2-Dichlorobenzene	8.37	146	662360	41.800	ng	98
15) Benzyl Alcohol	8.29	79	487603	39.830	ng	# 80
16) 2,2'-oxybis(1-Chloropropan	8.55	45	250522	37.216	ng	# 41
17) 2-Methylphenol	8.48	107	401378	42.516	ng	# 75
18) Hexachloroethane	9.08	117	312809	42.098	ng	# 36
19) n-Nitroso-di-n-propylamine	8.85	70	429797	41.584	ng	# 86
20) 3+4-Methylphenols	8.82	107	532670	40.808	ng	# 82
22) Acetophenone	8.87	105	853622	52.221	ng	# 95
24) Nitrobenzene	9.25	77	724919	48.800	ng	94
25) Isophorone	9.77	82	1179086	59.338	ng	# 94
26) 2-Nitrophenol	9.96	139	328195	51.680	ng	# 49
27) 2,4-Dimethylphenol	10.01	122	484786	58.577	ng	# 76
28) bis(2-Chloroethoxy)methane	10.25	93	626593	52.826	ng	# 96
29) 2,4-Dichlorophenol	10.49	162	767380	61.954	ng	93
30) 1,2,4-Trichlorobenzene	10.69	180	1178108	62.683	ng	# 94
31) Naphthalene	10.88	128	1737943	53.736	ng	98
32) Benzoic acid	10.25	122	311601	45.000	ng	# 76
33) 4-Chloroaniline	11.03	127	124012m	9.111	ng	
34) Hexachlorobutadiene	11.13	225	1082388	60.338	ng	97
35) Caprolactam	11.85	113	101269m	34.073	ng	
36) 4-Chloro-3-methylphenol	12.14	107	620715	52.127	ng	92
37) 2-Methylnaphthalene	12.49	142	1604667	63.377	ng	# 90
39) 1,2,4,5-Tetrachlorobenzene	12.85	216	1769003	54.835	ng	# 100
40) Hexachlorocyclopentadiene	12.80	237	2051429	118.177	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.10	196	1086530	56.322	nq	97
43) 2,4,5-Trichlorophenol	13.19	196	1124389m	59.667	nq	
45) 1,1'-Biphenyl	13.49	154	2406083	48.038	nq	98
46) 2-Chloronaphthalene	13.55	162	2007516	49.334	nq	98
47) 2-Nitroaniline	13.77	65	444335	40.996	nq	82
48) Acenaphthylene	14.39	152	3010876	52.418	nq	99
49) Dimethylphthalate	14.13	163	2816583	51.324	nq	# 98
50) 2,6-Dinitrotoluene	14.27	165	554426	51.796	nq	# 89
51) Acenaphthene	14.73	154	1744128	49.425	nq	95
52) 3-Nitroaniline	14.60	138	186215	19.120	nq	# 69
53) 2,4-Dinitrophenol	14.83	184	802017	149.724	nq	# 71
54) Dibenzofuran	15.06	168	3780529	55.964	nq	98
55) 4-Nitrophenol	14.92	139	485923	78.223	nq	# 82
56) 2,4-Dinitrotoluene	15.07	165	947526	53.086	nq	# 83
57) Fluorene	15.71	166	3019656	59.172	nq	99
58) 2,3,4,6-Tetrachlorophenol	15.30	232	1042692	57.570	nq	# 100
59) Diethylphthalate	15.48	149	2671334	46.685	nq	94
60) 4-Chlorophenyl-phenylether	15.70	204	2339460	54.666	nq	# 86
61) 4-Nitroaniline	15.77	138	339430	35.660	nq	# 1
62) Azobenzene	15.99	77	2677495	43.932	nq	87
64) 4,6-Dinitro-2-methylphenol	15.83	198	639346	49.283	nq	# 55
65) n-Nitrosodiphenylamine	15.92	169	2486909	51.777	nq	99
66) 4-Bromophenyl-phenylether	16.59	248	1436442	54.755	nq	99
67) Hexachlorobenzene	16.71	284	1642526	53.688	nq	96
68) Atrazine	16.87	200	1138854	51.955	nq	90
69) Pentachlorophenol	17.06	266	1587882	98.094	nq	100
70) Phenanthrene	17.44	178	5031417	57.016	nq	99
71) Anthracene	17.54	178	5145244	58.717	nq	98
72) Carbazole	17.82	167	3646689	46.308	nq	98
73) Di-n-butylphthalate	18.34	149	4519276	47.073	nq	# 97
74) Fluoranthene	19.44	202	6974916	53.907	nq	94
76) Benzidine	19.64	184	164737	4.275	nq	97
77) Pyrene	19.80	202	6968287	63.990	nq	99
79) Butylbenzylphthalate	20.67	149	1966805	52.325	nq	# 83
80) Benzo(a)anthracene	21.53	228	6624648	56.235	nq	99
81) 3,3'-Dichlorobenzidine	21.46	252	1111745	21.201	nq	# 95
82) Chrysene	21.59	228	6099323	54.264	nq	98
83) Bis(2-ethylhexyl)phthalate	21.42	149	2806348	50.221	nq	# 97
84) Di-n-octyl phthalate	22.31	149	4436827	47.779	nq	# 92
85) Indeno(1,2,3-cd)pyrene	26.16	276	7473695	46.235	nq	# 92
87) Benzo(b)fluoranthene	23.14	252	6726134	63.017	nq	# 90
88) Benzo(k)fluoranthene	23.19	252	6124462	56.394	nq	# 93
89) Benzo(a)pyrene	23.74	252	6032184	58.083	nq	# 94
90) Dibenzo(a,h)anthracene	26.16	278	6365714	54.692	nq	# 88
91) Benzo(g,h,i)perylene	26.87	276	5741386	54.953	nq	# 81

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

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