

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN072518\
 Data File : BN002149.D
 Acq On : 25 Jul 2018 23:01
 Operator : JU/SJ
 Sample : J3762-03
 Misc : MDL 2 PPM
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 MDL-SOIL-03-QT3-2018

Manual Integrations
 APPROVED

Sohil
 7/27/2018 11:35:43 AM

Quant Time: Jul 26 06:50:15 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-BN072518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 25 20:08:02 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.69	152	211471	20.00	ng	0.00
21) Naphthalene-d8	10.46	136	877431	20.00	ng	0.00
38) Acenaphthene-d10	14.32	164	521219	20.00	ng	0.00
63) Phenanthrene-d10	17.07	188	1120640	20.00	ng	0.00
75) Chrysene-d12	21.27	240	972218	20.00	ng	0.00
86) Perylene-d12	23.49	264	875795	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.31	112	1589267	120.99	ng	0.00
7) Phenol-d6	6.88	99	2147504	116.50	ng	0.00
23) Nitrobenzene-d5	8.84	82	1498790	87.60	ng	0.00
41) 2,4,6-Tribromophenol	15.82	330	843006	124.55	ng	0.00
44) 2-Fluorobiphenyl	12.94	172	3440689	85.98	ng	0.00
78) Terphenyl-d14	19.71	244	5036879	108.05	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.29	88	8694	1.763	ng	# 65
3) Pyridine	3.68	79	21430	1.500	ng	# 68
4) n-Nitrosodimethylamine	3.59	42	12366	2.018	ng	# 44
6) Aniline	7.03	93	26148	1.161	ng	# 95
8) 2-Chlorophenol	7.26	128	27174	1.784	ng	96
9) Benzaldehyde	6.85	77	27718	2.442	ng	90
10) Phenol	6.90	94	35701	1.889	ng	80
11) bis(2-Chloroethyl)ether	7.13	93	26302	1.710	ng	87
12) 1,3-Dichlorobenzene	7.58	146	31543	1.817	ng	96
13) 1,4-Dichlorobenzene	7.73	146	32367	1.822	ng	97
14) 1,2-Dichlorobenzene	8.04	146	30460	1.778	ng	98
15) Benzyl Alcohol	7.93	79	17045	1.225	ng	95
16) 2,2'-oxybis(1-Chloropropan	8.22	45	46368	1.867	ng	92
17) 2-Methylphenol	8.13	107	22744	1.763	ng	97
18) Hexachloroethane	8.76	117	11479	1.766	ng	91
19) n-Nitroso-di-n-propylamine	8.49	70	21850	1.634	ng	95
20) 3+4-Methylphenols	8.46	107	28968	1.610	ng	88
22) Acetophenone	8.50	105	43024	1.891	ng	96
24) Nitrobenzene	8.88	77	33901	1.912	ng	# 95
25) Isophorone	9.40	82	58480	1.991	ng	# 95
26) 2-Nitrophenol	9.59	139	12129	1.416	ng	96
27) 2,4-Dimethylphenol	9.65	122	23659	1.986	ng	94
28) bis(2-Chloroethoxy)methane	9.89	93	36212	1.908	ng	99
29) 2,4-Dichlorophenol	10.13	162	22691	1.633	ng	95
30) 1,2,4-Trichlorobenzene	10.33	180	28716	1.825	ng	96
31) Naphthalene	10.51	128	91364	2.063	ng	97
32) Benzoic acid	9.72	122	6918m	0.684	ng	
33) 4-Chloroaniline	10.63	127	24692	1.253	ng	96
34) Hexachlorobutadiene	10.80	225	17418	1.808	ng	96
35) Caprolactam	11.38	113	6055	1.277	ng	# 55
36) 4-Chloro-3-methylphenol	11.76	107	26161	1.656	ng	# 93
37) 2-Methylnaphthalene	12.13	142	67368	2.157	ng	95
39) 1,2,4,5-Tetrachlorobenzene	12.50	216	32881	1.876	ng	99
40) Hexachlorocyclopentadiene	12.48	237	14795	1.495	ng	87

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	12.75	196	18757	1.595	ng	100
43) 2,4,5-Trichlorophenol	12.83	196	19349	1.553	ng	91
45) 1,1'-Biphenyl	13.15	154	91220	1.989	ng	99
46) 2-Chloronaphthalene	13.19	162	67030	1.893	ng	96
47) 2-Nitroaniline	13.40	65	16268	1.428	ng	91
48) Acenaphthylene	14.04	152	100710	1.875	ng	97
49) Dimethylphthalate	13.78	163	82507	1.906	ng	100
50) 2,6-Dinitrotoluene	13.90	165	15614	1.623	ng	# 90
51) Acenaphthene	14.39	154	63991	1.977	ng	95
52) 3-Nitroaniline	14.23	138	12977	1.245	ng	# 86
53) 2,4-Dinitrophenol	14.45	184	4179	0.821	ng	93
54) Dibenzofuran	14.72	168	102314	1.950	ng	96
55) 4-Nitrophenol	14.56	139	9534	1.192	ng	# 82
56) 2,4-Dinitrotoluene	14.70	165	18262	1.399	ng	# 95
57) Fluorene	15.37	166	77494	1.960	ng	99
58) 2,3,4,6-Tetrachlorophenol	14.95	232	17410	1.625	ng	# 95
59) Diethylphthalate	15.15	149	79240	1.798	ng	96
60) 4-Chlorophenyl-phenylether	15.37	204	40646	1.857	ng	96
61) 4-Nitroaniline	15.40	138	14409	1.356	ng	94
62) Azobenzene	15.66	77	74430	1.760	ng	98
64) 4,6-Dinitro-2-methylphenol	15.47	198	7277	0.962	ng	82
65) n-Nitrosodiphenylamine	15.59	169	66548	1.790	ng	93
66) 4-Bromophenyl-phenylether	16.27	248	23537	1.794	ng	95
67) Hexachlorobenzene	16.38	284	26996	1.860	ng	# 88
68) Atrazine	16.54	200	18903	1.510	ng	95
69) Pentachlorophenol	16.73	266	8503	0.910	ng	96
70) Phenanthrene	17.11	178	126148	2.062	ng	99
71) Anthracene	17.20	178	119390	1.972	ng	98
72) Carbazole	17.47	167	100097	1.648	ng	98
73) Di-n-butylphthalate	18.05	149	105503	1.489	ng	98
74) Fluoranthene	19.13	202	123048	1.853	ng	98
76) Benzidine	19.33	184	20905	0.659	ng	# 91
77) Pyrene	19.50	202	126845	1.892	ng	98
79) Butylbenzylphthalate	20.41	149	40058	1.373	ng	94
80) Benzo(a)anthracene	21.25	228	116661	1.969	ng	99
81) 3,3'-Dichlorobenzidine	21.19	252	25244	1.093	ng	# 94
82) Chrysene	21.30	228	111658	1.979	ng	98
83) Bis(2-ethylhexyl)phthalate	21.19	149	58465	1.437	ng	# 98
84) Di-n-octyl phthalate	22.07	149	90737	1.406	ng	98
85) Indeno(1,2,3-cd)pyrene	25.72	276	103566	1.883	ng	# 95
87) Benzo(b)fluoranthene	22.83	252	103997	2.012	ng	98
88) Benzo(k)fluoranthene	22.87	252	101341	1.991	ng	97
89) Benzo(a)pyrene	23.40	252	95510	1.979	ng	# 94
90) Dibenzo(a,h)anthracene	25.73	278	89909	1.984	ng	# 95
91) Benzo(g,h,i)perylene	26.40	276	87514	1.998	ng	98

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

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