

Data Path : Z:\HPCHEM1\BNA N\DATA\BN021518\
 Data File : BN000094.D
 Acq On : 15 Feb 2018 13:32
 Operator : JU/SJ
 Sample : J1161-03
 Misc : 2 PPM MDL
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 MDL-SOIL-03-QT1-2018

Quant Time: Feb 16 09:57:18 2018
 Quant Method : Z:\HPCHEM1\BNA N\METHODS\8270-BN020718.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 07 18:24:45 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.46	152	384449	20.00	ng	0.00
21) Naphthalene-d8	11.30	136	1913340	20.00	ng	-0.01
38) Acenaphthene-d10	15.07	164	1146902	20.00	ng	-0.01
63) Phenanthrene-d10	17.79	188	2556080	20.00	ng	-0.01
75) Chrysene-d12	21.90	240	2722328	20.00	ng	-0.02
86) Perylene-d12	24.39	264	2821325	20.00	ng	-0.03

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.93	112	3093863	131.16	ng	-0.01
7) Phenol-d6	7.58	99	4379836	122.57	ng	-0.01
23) Nitrobenzene-d5	9.64	82	3066116	95.28	ng	-0.01
41) 2,4,6-Tribromophenol	16.54	330	1467288	129.25	ng	-0.01
44) 2-Fluorobiphenyl	13.70	172	7780983	95.62	ng	-0.01
78) Terphenyl-d14	20.35	244	9677826	95.99	ng	-0.01

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	3.66	88	19108	2.013	ng	# 93
3) Pyridine	4.14	79	40783	1.380	ng	# 95
4) n-Nitrosodimethylamine	4.02	42	13932	1.667	ng	# 91
6) Aniline	7.77	93	77492	1.574	ng	# 85
8) 2-Chlorophenol	8.01	128	52932	1.919	ng	92
9) Benzaldehyde	7.58	77	34399	1.547	ng	92
10) Phenol	7.61	94	72250	1.915	ng	93
11) bis(2-Chloroethyl)ether	7.86	93	60046	1.936	ng	# 82
12) 1,3-Dichlorobenzene	8.35	146	61916	1.975	ng	99
13) 1,4-Dichlorobenzene	8.50	146	63140	1.973	ng	94
14) 1,2-Dichlorobenzene	8.82	146	62877	1.996	ng	88
15) Benzyl Alcohol	8.69	79	39013	1.519	ng	93
16) 2,2'-oxybis(1-Chloropropan	8.99	45	61039	1.859	ng	81
17) 2-Methylphenol	8.89	107	41168	1.536	ng	96
18) Hexachloroethane	9.57	117	19752	1.773	ng	# 78
19) n-Nitroso-di-n-propylamine	9.26	70	34878	1.457	ng	# 82
20) 3+4-Methylphenols	9.22	107	54779	1.461	ng	92
22) Acetophenone	9.29	105	87289	1.628	ng	# 87
24) Nitrobenzene	9.68	77	68460	1.946	ng	# 89
25) Isophorone	10.20	82	94023	1.413	ng	96
26) 2-Nitrophenol	10.40	139	14903	8.071	ng	90
27) 2,4-Dimethylphenol	10.44	122	45350	1.572	ng	94
28) bis(2-Chloroethoxy)methane	10.69	93	72764	1.706	ng	98
29) 2,4-Dichlorophenol	10.93	162	32141	1.204	ng	96
30) 1,2,4-Trichlorobenzene	11.16	180	59536	1.937	ng	90
31) Naphthalene	11.35	128	203382	1.939	ng	97
32) Benzoic acid	10.46	122	8642m	10.858	ng	
33) 4-Chloroaniline	11.45	127	66224	1.437	ng	94
34) Hexachlorobutadiene	11.63	225	31811	1.996	ng	95
35) Caprolactam	12.16	113	9495	4.764	ng	# 90
36) 4-Chloro-3-methylphenol	12.53	107	36316	1.072	ng	87
37) 2-Methylnaphthalene	12.93	142	129514	1.775	ng	95
39) 1,2,4,5-Tetrachlorobenzene	13.29	216	64950	1.919	ng	# 96
40) Hexachlorocyclopentadiene	13.26	237	20253	1.435	ng	96

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42) 2,4,6-Trichlorophenol	13.51	196	21271	1.048	nq	88
43) 2,4,5-Trichlorophenol	13.57	196	26022	1.067	nq #	93
45) 1,1'-Biphenyl	13.91	154	207601	2.055	nq	98
46) 2-Chloronaphthalene	13.96	162	142457	1.880	nq	99
47) 2-Nitroaniline	14.15	65	19617	5.671	nq	91
48) Acenaphthylene	14.80	152	183946	1.506	nq	98
49) Dimethylphthalate	14.50	163	154233	1.609	nq	99
50) 2,6-Dinitrotoluene	14.63	165	18145	0.917	nq	93
51) Acenaphthene	15.13	154	146134	1.839	nq	96
52) 3-Nitroaniline	14.95	138	22440	0.933	nq #	94
53) 2,4-Dinitrophenol	15.15	184	5037	9.863	nq #	67
54) Dibenzofuran	15.46	168	214387	1.904	nq	92
55) 4-Nitrophenol	15.23	139	13181	5.604	nq	89
56) 2,4-Dinitrotoluene	15.40	165	20341	4.838	nq #	85
57) Fluorene	16.10	166	167431	1.822	nq	97
58) 2,3,4,6-Tetrachlorophenol	15.67	232	19873m	1.032	nq	
59) Diethylphthalate	15.85	149	144654	1.485	nq	97
60) 4-Chlorophenyl-phenylether	16.09	204	79523	1.898	nq	94
61) 4-Nitroaniline	16.10	138	23976	0.970	nq	83
62) Azobenzene	16.38	77	134228	1.384	nq	92
64) 4,6-Dinitro-2-methylphenol	16.16	198	7494	9.214	nq	89
65) n-Nitrosodiphenylamine	16.30	169	128479	1.545	nq	98
66) 4-Bromophenyl-phenylether	16.97	248	41828	1.721	nq #	85
67) Hexachlorobenzene	17.10	284	51868	1.827	nq #	90
68) Atrazine	17.22	200	29252	1.081	nq	91
69) Pentachlorophenol	17.43	266	16648	6.620	nq	94
70) Phenanthrene	17.83	178	301030	1.994	nq	97
71) Anthracene	17.92	178	235597	1.612	nq	98
72) Carbazole	18.18	167	228914	1.554	nq	97
73) Di-n-butylphthalate	18.71	149	173370	1.066	nq #	96
74) Fluoranthene	19.80	202	259547	1.639	nq	94
76) Benzidine	19.97	184	46787	0.539	nq	98
77) Pyrene	20.16	202	297808	1.721	nq	99
79) Butylbenzylphthalate	21.02	149	59381	6.018	nq	94
80) Benzo(a)anthracene	21.89	228	289438	1.730	nq	96
81) 3,3'-Dichlorobenzidine	21.81	252	59710	0.972	nq #	92
82) Chrysene	21.95	228	335041	2.009	nq	100
83) Bis(2-ethylhexyl)phthalate	21.79	149	89420	3.593	nq #	90
84) Di-n-octyl phthalate	22.74	149	119210	7.169	nq	97
85) Indeno(1,2,3-cd)pyrene	26.94	276	309642	1.512	nq #	86
87) Benzo(b)fluoranthene	23.63	252	278059	1.684	nq #	95
88) Benzo(k)fluoranthene	23.68	252	291311	1.751	nq #	97
89) Benzo(a)pyrene	24.27	252	253549	1.601	nq #	94
90) Dibenzo(a,h)anthracene	26.95	278	260098	1.571	nq #	89
91) Benzo(g,h,i)perylene	27.72	276	260814	1.557	nq #	87

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

