

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM072318\
 Data File : BM016035.D
 Acq On : 23 Jul 2018 20:51
 Operator : SJ/JU
 Sample : J3762-02
 Misc : L00 20 PPM
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 LOQ-SOIL-02-QT3-2018

Quant Time: Jul 24 06:05:33 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM072318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 23 16:08:38 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.23	152	41298	20.00	ng	0.00
21) Naphthalene-d8	11.05	136	157206	20.00	ng	0.00
38) Acenaphthene-d10	14.86	164	82748	20.00	ng	0.00
63) Phenanthrene-d10	17.59	188	179548	20.00	ng	0.00
75) Chrysene-d12	21.70	240	194373	20.00	ng	0.00
86) Perylene-d12	24.07	264	199592	20.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.74	112	321452	129.30	ng	0.00
7) Phenol-d6	7.39	99	398586	122.76	ng	0.00
23) Nitrobenzene-d5	9.42	82	302101	89.38	ng	0.00
41) 2,4,6-Tribromophenol	16.34	330	142450	135.73	ng	0.00
44) 2-Fluorobiphenyl	13.48	172	611439	89.91	ng	0.00
78) Terphenyl-d14	20.16	244	1109535	119.49	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	3.52	88	19068	16.370	ng	# 100
3) Pyridine	3.97	79	44901	16.001	ng	# 79
4) n-Nitrosodimethylamine	3.88	42	40226	18.379	ng	# 38
6) Aniline	7.56	93	48117	12.871	ng	92
8) 2-Chlorophenol	7.79	128	52995	17.935	ng	99
9) Benzaldehyde	7.37	77	46781	20.508	ng	89
10) Phenol	7.42	94	58826	18.120	ng	96
11) bis(2-Chloroethyl)ether	7.65	93	42532	16.583	ng	90
12) 1,3-Dichlorobenzene	8.12	146	58748	16.986	ng	# 85
13) 1,4-Dichlorobenzene	8.27	146	60178	17.156	ng	# 93
14) 1,2-Dichlorobenzene	8.59	146	58641	17.085	ng	# 93
15) Benzyl Alcohol	8.49	79	40493	15.663	ng	89
16) 2,2'-oxybis(1-Chloropropan	8.75	45	65596	16.711	ng	97
17) 2-Methylphenol	8.68	107	40673	18.329	ng	# 86
18) Hexachloroethane	9.31	117	26039	17.173	ng	97
19) n-Nitroso-di-n-propylamine	9.05	70	38039	16.029	ng	# 75
20) 3+4-Methylphenols	9.01	107	52552	16.460	ng	89
22) Acetophenone	9.07	105	75379	19.043	ng	# 80
24) Nitrobenzene	9.46	77	64942	19.088	ng	95
25) Isophorone	9.98	82	94880	20.522	ng	# 94
26) 2-Nitrophenol	10.17	139	23774	16.740	ng	# 56
27) 2,4-Dimethylphenol	10.22	122	41842	21.147	ng	# 78
28) bis(2-Chloroethoxy)methane	10.46	93	58982	19.629	ng	99
29) 2,4-Dichlorophenol	10.71	162	44988	19.134	ng	92
30) 1,2,4-Trichlorobenzene	10.92	180	52384	18.337	ng	# 95
31) Naphthalene	11.10	128	159587	20.058	ng	98
32) Benzoic acid	10.35	122	25026	16.592	ng	# 85
33) 4-Chloroaniline	11.23	127	41597	12.812	ng	# 89
34) Hexachlorobutadiene	11.36	225	36779	18.560	ng	96
35) Caprolactam	12.02	113	10007	15.367	ng	93
36) 4-Chloro-3-methylphenol	12.33	107	47834	18.497	ng	86
37) 2-Methylnaphthalene	12.70	142	113003	21.202	ng	# 90
39) 1,2,4,5-Tetrachlorobenzene	13.06	216	55052	19.542	ng	# 100
40) Hexachlorocyclopentadiene	13.02	237	23125	22.792	ng	95

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42) 2,4,6-Trichlorophenol	13.30	196	33733	19.139	nq	99
43) 2,4,5-Trichlorophenol	13.38	196	35827	19.706	nq #	83
45) 1,1'-Biphenyl	13.69	154	139360	18.661	nq	96
46) 2-Chloronaphthalene	13.74	162	109197	18.800	nq #	89
47) 2-Nitroaniline	13.96	65	32181	16.759	nq #	75
48) Acenaphthylene	14.59	152	173215	20.371	nq	99
49) Dimethylphthalate	14.31	163	140648	19.422	nq	99
50) 2,6-Dinitrotoluene	14.44	165	27851	19.114	nq #	51
51) Acenaphthene	14.92	154	101400	19.390	nq	98
52) 3-Nitroaniline	14.78	138	21002	13.345	nq #	68
53) 2,4-Dinitrophenol	15.00	184	9193	19.688	nq #	83
54) Dibenzofuran	15.25	168	164946	19.141	nq #	79
55) 4-Nitrophenol	15.10	139	18363	16.276	nq #	82
56) 2,4-Dinitrotoluene	15.24	165	37133	17.823	nq #	85
57) Fluorene	15.90	166	130145	20.323	nq	97
58) 2,3,4,6-Tetrachlorophenol	15.48	232	30955	19.809	nq #	100
59) Diethylphthalate	15.65	149	147397	18.884	nq	95
60) 4-Chlorophenyl-phenylether	15.88	204	64887	18.202	nq #	80
61) 4-Nitroaniline	15.94	138	23437	15.521	nq #	10
62) Azobenzene	16.17	77	159447	19.397	nq #	76
64) 4,6-Dinitro-2-methylphenol	16.00	198	16140	14.810	nq	85
65) n-Nitrosodiphenylamine	16.10	169	110167	18.634	nq	95
66) 4-Bromophenyl-phenylether	16.77	248	37318	18.355	nq #	75
67) Hexachlorobenzene	16.89	284	41078	18.347	nq #	67
68) Atrazine	17.03	200	39676	18.724	nq	97
69) Pentachlorophenol	17.24	266	17560	12.666	nq	96
70) Phenanthrene	17.63	178	200626	19.958	nq	99
71) Anthracene	17.72	178	200001	20.472	nq	98
72) Carbazole	17.99	167	179246	17.710	nq	98
73) Di-n-butylphthalate	18.51	149	225573	17.886	nq #	96
74) Fluoranthene	19.61	202	219751	19.598	nq	99
76) Benzidine	19.79	184	72240	13.300	nq	100
77) Pyrene	19.97	202	228913	19.652	nq	99
79) Butylbenzylphthalate	20.83	149	107928	18.239	nq #	84
80) Benzo(a)anthracene	21.69	228	238198	19.897	nq	99
81) 3,3'-Dichlorobenzidine	21.62	252	48562	10.071	nq	95
82) Chrysene	21.74	228	229142	19.953	nq	99
83) Bis(2-ethylhexyl)phthalate	21.57	149	150825	17.587	nq	91
84) Di-n-octyl phthalate	22.48	149	264474	18.349	nq #	89
85) Indeno(1,2,3-cd)pyrene	26.50	276	279498	20.510	nq #	94
87) Benzo(b)fluoranthene	23.35	252	241442	20.546	nq #	95
88) Benzo(k)fluoranthene	23.40	252	236134	19.827	nq	97
89) Benzo(a)pyrene	23.97	252	233110	20.634	nq	100
90) Dibenzo(a,h)anthracene	26.50	278	235904	20.155	nq #	96
91) Benzo(g,h,i)perylene	27.25	276	226988	20.504	nq #	94

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

