

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM041519\
 Data File : BM019919.D
 Acq On : 15 Apr 2019 18:35
 Operator : JU/SJ
 Sample : K1560-01
 Misc : LOD 16PPM
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 LOD-MDL-SOIL-01-QT1-2019

Manual Integrations
 APPROVED

Sohil
 4/16/2019 5:25:46 PM

Quant Time: Apr 16 03:31:07 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\8270-BM041519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Apr 15 16:16:29 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.53	152	111173	20.00	ng	0.00
21) Naphthalene-d8	10.31	136	497739	20.00	ng	0.00
39) Acenaphthene-d10	14.20	164	302366	20.00	ng	0.00
64) Phenanthrene-d10	16.96	188	666579	20.00	ng	0.00
76) Chrysene-d12	21.17	240	681723	20.00	ng	0.00
87) Perylene-d12	23.37	264	636892	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.14	112	909342	132.56	ng	0.00
7) Phenol-d6	6.72	99	1326454	128.48	ng	0.00
23) Nitrobenzene-d5	8.70	82	929260	83.49	ng	0.00
42) 2,4,6-Tribromophenol	15.71	330	682339	140.19	ng	0.00
45) 2-Fluorobiphenyl	12.80	172	2036799	83.95	ng	0.00
79) Terphenyl-d14	19.61	244	3388231	87.66	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.12	88	47658	16.099	ng	98
3) Pyridine	3.53	79	136251	16.435	ng	97
4) n-Nitrosodimethylamine	3.46	42	38548	14.434	ng	# 99
6) Aniline	6.88	93	201825	15.136	ng	99
8) 2-Chlorophenol	7.10	128	131128	15.369	ng	98
9) Benzaldehyde	6.70	77	122143	17.351	ng	99
10) Phenol	6.75	94	174254	15.449	ng	94
11) bis(2-Chloroethyl)ether	6.98	93	128107	15.523	ng	100
12) 1,3-Dichlorobenzene	7.42	146	140863	15.727	ng	97
13) 1,4-Dichlorobenzene	7.56	146	141637	15.585	ng	96
14) 1,2-Dichlorobenzene	7.88	146	139225	15.451	ng	100
15) Benzyl Alcohol	7.80	79	144458	15.392	ng	98
16) 2,2'-oxybis(1-Chloropropan	8.05	45	123604	15.711	ng	99
17) 2-Methylphenol	7.99	107	110575	14.812	ng	98
18) Hexachloroethane	8.58	117	54151	15.329	ng	98
19) n-Nitroso-di-n-propylamine	8.34	70	134756	14.849	ng	98
20) 3+4-Methylphenols	8.33	107	148041	14.436	ng	98
22) Acetophenone	8.37	105	216825	15.994	ng	97
24) Nitrobenzene	8.75	77	194397	16.861	ng	100
25) Isophorone	9.26	82	347394	16.034	ng	99
26) 2-Nitrophenol	9.45	139	74792	15.222	ng	97
27) 2,4-Dimethylphenol	9.51	122	102797	14.797	ng	99
28) bis(2-Chloroethoxy)methane	9.75	93	190948	15.739	ng	98
29) 2,4-Dichlorophenol	9.99	162	120730	15.302	ng	99
30) 1,2,4-Trichlorobenzene	10.18	180	138377	16.295	ng	99
31) Naphthalene	10.36	128	435516	16.313	ng	99
32) Benzoic acid	9.67	122	78336m	15.779	ng	
33) 4-Chloroaniline	10.51	127	167825	15.259	ng	99
34) Hexachlorobutadiene	10.61	225	81480	16.269	ng	99
35) Caprolactam	11.33	113	42970	14.388	ng	87
36) 4-Chloro-3-methylphenol	11.64	107	155140	15.825	ng	97
37) 2-Methylnaphthalene	11.99	142	314668	16.189	ng	98
38) 1-Methylnaphthalene	12.20	142	311436	16.228	ng	98
40) 1,2,4,5-Tetrachlorobenzene	12.36	216	146883	15.747	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.31	237	18387	16.420	ng	95
43) 2,4,6-Trichlorophenol	12.63	196	98798	15.623	ng	99
44) 2,4,5-Trichlorophenol	12.71	196	96155	14.609	ng	99
46) 1,1'-Biphenyl	13.02	154	426143	16.272	ng	99
47) 2-Chloronaphthalene	13.06	162	325140	16.276	ng	99
48) 2-Nitroaniline	13.31	65	110084	15.289	ng	97
49) Acenaphthylene	13.92	152	557452	16.039	ng	99
50) Dimethylphthalate	13.67	163	437048	16.387	ng	100
51) 2,6-Dinitrotoluene	13.80	165	94721	16.061	ng	88
52) Acenaphthene	14.26	154	312372	16.209	ng	99
53) 3-Nitroaniline	14.15	138	89879	15.262	ng	98
54) 2,4-Dinitrophenol	14.39	184	33487	16.098	ng	91
55) Dibenzofuran	14.60	168	507879	16.116	ng	99
56) 4-Nitrophenol	14.49	139	51248	16.081	ng	98
57) 2,4-Dinitrotoluene	14.62	165	127679	15.061	ng	# 98
58) Fluorene	15.26	166	416229	15.587	ng	100
59) 2,3,4,6-Tetrachlorophenol	14.84	232	94972	15.814	ng	98
60) Diethylphthalate	15.04	149	448060	16.153	ng	99
61) 4-Chlorophenyl-phenylether	15.25	204	199317	15.503	ng	94
62) 4-Nitroaniline	15.33	138	81602	14.064	ng	96
63) Azobenzene	15.55	77	482712	16.282	ng	98
65) 4,6-Dinitro-2-methylphenol	15.39	198	57930	13.621	ng	98
66) n-Nitrosodiphenylamine	15.48	169	349482	16.037	ng	99
67) 4-Bromophenyl-phenylether	16.15	248	122441	15.977	ng	98
68) Hexachlorobenzene	16.26	284	147030	15.989	ng	99
69) Atrazine	16.44	200	125273	16.934	ng	99
70) Pentachlorophenol	16.63	266	66343	14.120	ng	99
71) Phenanthrene	17.00	178	638953	16.153	ng	100
72) Anthracene	17.10	178	622816	15.817	ng	99
73) Carbazole	17.39	167	578471	15.735	ng	99
74) Di-n-butylphthalate	17.93	149	740966	15.347	ng	100
75) Fluoranthene	19.03	202	718609	15.786	ng	99
77) Benzidine	19.25	184	249533	13.266	ng	98
78) Pyrene	19.40	202	739411	16.378	ng	99
80) Butylbenzylphthalate	20.31	149	341438	15.600	ng	98
81) Benzo(a)anthracene	21.16	228	701288	16.239	ng	99
82) 3,3'-Dichlorobenzidine	21.10	252	265423	16.519	ng	98
83) Chrysene	21.22	228	680534	16.432	ng	100
84) Bis(2-ethylhexyl)phthalate	21.07	149	491765	15.092	ng	99
85) Di-n-octyl phthalate	21.94	149	806912	15.261	ng	100
86) Indeno(1,2,3-cd)pyrene	25.58	276	705856	17.325	ng	100
88) Benzo(b)fluoranthene	22.72	252	657863	15.556	ng	100
89) Benzo(k)fluoranthene	22.76	252	647087	16.253	ng	99
90) Benzo(a)pyrene	23.28	252	604599	16.135	ng	99
91) Dibenzo(a,h)anthracene	25.59	278	583708	16.157	ng	100
92) Benzo(g,h,i)perylene	26.26	276	550441	16.822	ng	100

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

