

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG072919\
 Data File : BG041787.D
 Acq On : 29 Jul 2019 13:36
 Operator : HP/JU
 Sample : K3591-01
 Misc : LOD-SOIL-5PPM
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 LOD-MDL-SOIL-01-QT3-2019

Manual Integrations
 APPROVED

mohammad
 7/30/2019 4:52:24 PM

Quant Time: Jul 30 07:45:12 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG072719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 29 12:47:21 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.07	152	89531	20.00	ng	0.02
21) Naphthalene-d8	10.89	136	385114	20.00	ng	0.02
39) Acenaphthene-d10	14.72	164	285710	20.00	ng	0.01
64) Phenanthrene-d10	17.46	188	645774	20.00	ng	0.00
76) Chrysene-d12	21.74	240	674819	20.00	ng	0.00
87) Perylene-d12	25.02	264	754207	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.61	112	591573	128.56	ng	0.02
7) Phenol-d6	7.22	99	780301	125.78	ng	0.02
23) Nitrobenzene-d5	9.23	82	480030	85.78	ng	0.02
42) 2,4,6-Tribromophenol	16.20	330	423237	130.42	ng	0.01
45) 2-Fluorobiphenyl	13.34	172	1347997	83.21	ng	0.01
79) Terphenyl-d14	20.07	244	2150364	72.92	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.48	88	12619	5.826	ng	89
3) Pyridine	3.90	79	25019	4.212	ng	88
4) n-Nitrosodimethylamine	3.80	42	10236	4.347	ng	# 97
6) Aniline	7.38	93	42749	4.894	ng	97
8) 2-Chlorophenol	7.62	128	26778	5.082	ng	98
9) Benzaldehyde	7.19	77	26027	5.962	ng	99
10) Phenol	7.24	94	32735	5.072	ng	99
11) bis(2-Chloroethyl)ether	7.48	93	26630	5.019	ng	90
12) 1,3-Dichlorobenzene	7.96	146	32866	4.779	ng	96
13) 1,4-Dichlorobenzene	8.11	146	34582	5.025	ng	93
14) 1,2-Dichlorobenzene	8.42	146	30814	4.693	ng	96
15) Benzyl Alcohol	8.30	79	22702	4.651	ng	97
16) 2,2'-oxybis(1-Chloropropan	8.59	45	44271	4.690	ng	94
17) 2-Methylphenol	8.51	107	22255	4.741	ng	94
18) Hexachloroethane	9.16	117	11601	4.923	ng	96
19) n-Nitroso-di-n-propylamine	8.88	70	21503	4.774	ng	# 89
20) 3+4-Methylphenols	8.83	107	30694	4.779	ng	97
22) Acetophenone	8.89	105	42776	4.966	ng	# 91
24) Nitrobenzene	9.27	77	30975	5.254	ng	92
25) Isophorone	9.80	82	60120	4.938	ng	98
26) 2-Nitrophenol	9.99	139	13158	4.809	ng	92
27) 2,4-Dimethylphenol	10.04	122	26475	5.482	ng	92
28) bis(2-Chloroethoxy)methane	10.29	93	37335	4.908	ng	98
29) 2,4-Dichlorophenol	10.53	162	27378	4.655	ng	95
30) 1,2,4-Trichlorobenzene	10.75	180	35126	4.710	ng	# 93
31) Naphthalene	10.94	128	90236	5.120	ng	95
32) Benzoic acid	10.14	122	2939m	9.425	ng	
33) 4-Chloroaniline	11.04	127	39493	4.725	ng	94
34) Hexachlorobutadiene	11.24	225	23298	4.629	ng	99
35) Caprolactam	11.78	113	9517	4.656	ng	# 81
36) 4-Chloro-3-methylphenol	12.15	107	28264	4.848	ng	97
37) 2-Methylnaphthalene	12.55	142	73061	5.237	ng	99
38) 1-Methylnaphthalene	12.76	142	69859	5.284	ng	99
40) 1,2,4,5-Tetrachlorobenzene	12.91	216	45601	5.042	ng	97

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.90	237	22606	4.446	nq	96
43) 2,4,6-Trichlorophenol	13.15	196	24958	4.366	nq	98
44) 2,4,5-Trichlorophenol	13.22	196	27493	4.628	nq	# 94
46) 1,1'-Biphenyl	13.55	154	95323	5.190	nq	96
47) 2-Chloronaphthalene	13.59	162	78380	5.013	nq	94
48) 2-Nitroaniline	13.78	65	15983	4.030	nq	95
49) Acenaphthylene	14.43	152	125908	5.383	nq	96
50) Dimethylphthalate	14.16	163	102695	5.263	nq	97
51) 2,6-Dinitrotoluene	14.27	165	20473	5.018	nq	# 87
52) Acenaphthene	14.78	154	75537	4.965	nq	97
53) 3-Nitroaniline	14.60	138	19610	4.493	nq	# 89
54) 2,4-Dinitrophenol	14.81	184	4227	5.627	nq	# 89
55) Dibenzofuran	15.11	168	130452	5.606	nq	# 88
56) 4-Nitrophenol	14.90	139	12845	3.877	nq	# 82
57) 2,4-Dinitrotoluene	15.06	165	24388	4.346	nq	96
58) Fluorene	15.76	166	102590	5.494	nq	98
59) 2,3,4,6-Tetrachlorophenol	15.33	232	26781	4.552	nq	# 96
60) Diethylphthalate	15.53	149	102320	5.341	nq	98
61) 4-Chlorophenyl-phenylether	15.76	204	59063	5.181	nq	96
62) 4-Nitroaniline	15.76	138	22156	4.686	nq	89
63) Azobenzene	16.04	77	82515	5.273	nq	95
65) 4,6-Dinitro-2-methylphenol	15.83	198	11411	9.947	nq	92
66) n-Nitrosodiphenylamine	15.96	169	93538	5.318	nq	97
67) 4-Bromophenyl-phenylether	16.65	248	38074	4.786	nq	97
68) Hexachlorobenzene	16.77	284	39448	4.740	nq	92
69) Atrazine	16.91	200	33119	4.792	nq	97
70) Pentachlorophenol	17.11	266	17074	3.612	nq	91
71) Phenanthrene	17.51	178	174940	5.652	nq	# 91
72) Anthracene	17.59	178	176057	5.703	nq	# 89
73) Carbazole	17.86	167	152069	5.488	nq	# 93
74) Di-n-butylphthalate	18.43	149	166960	5.315	nq	# 91
75) Fluoranthene	19.52	202	217133	5.771	nq	87
77) Benzidine	19.69	184	96878	4.398	nq	97
78) Pyrene	19.87	202	228472	5.714	nq	# 89
80) Butylbenzylphthalate	20.76	149	69345	4.524	nq	89
81) Benzo(a)anthracene	21.72	228	216295m	5.417	nq	
82) 3,3'-Dichlorobenzidine	21.64	252	78271	4.883	nq	# 97
83) Chrysene	21.79	228	210021	5.488	nq	90
84) Bis(2-ethylhexyl)phthalate	21.64	149	104155	4.926	nq	95
85) Di-n-octyl phthalate	22.88	149	167668	4.712	nq	# 89
86) Indeno(1,2,3-cd)pyrene	28.76	276	234882	4.636	nq	# 89
88) Benzo(b)fluoranthene	23.98	252	208301	5.049	nq	# 94
89) Benzo(k)fluoranthene	24.05	252	220556	5.488	nq	# 92
90) Benzo(a)pyrene	24.87	252	206912	5.229	nq	# 94
91) Dibenzo(a,h)anthracene	28.84	278	193413	4.978	nq	96
92) Benzo(g,h,i)perylene	29.93	276	192226	4.952	nq	# 90

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

