

Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG101918\
 Data File : BG037448.D
 Acq On : 19 Oct 2018 11:24
 Operator : JU/SJ
 Sample : J5164-01
 Misc : LOD 1PPM
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 LOD-MDL-SOIL-01-QT4-2018

Manual Integrations
 APPROVED

Sohil
 10/22/2018 3:38:03 PM

Quant Time: Oct 20 00:43:29 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG101818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 18 14:06:42 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.11	152	50287	20.00	ng	0.00
21) Naphthalene-d8	10.93	136	233142	20.00	ng	0.00
39) Acenaphthene-d10	14.74	164	155854	20.00	ng	0.00
64) Phenanthrene-d10	17.48	188	385113	20.00	ng	0.00
76) Chrysene-d12	21.77	240	359834	20.00	ng	0.00
87) Perylene-d12	25.11	264	350533	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.64	112	399936	132.96	ng	0.00
7) Phenol-d6	7.24	99	598020	131.48	ng	0.00
23) Nitrobenzene-d5	9.28	82	348958	83.85	ng	0.00
42) 2,4,6-Tribromophenol	16.23	330	229562	135.05	ng	0.00
45) 2-Fluorobiphenyl	13.36	172	864307	83.62	ng	0.00
79) Terphenyl-d14	20.09	244	1323264	78.58	ng	0.00

Target Compounds

						Qvalue
6) Aniline	7.43	93	3031	0.546	ng	94
8) 2-Chlorophenol	7.67	128	3469	0.986	ng	84
9) Benzaldehyde	7.24	77	3105	1.091	ng	# 69
10) Phenol	7.27	94	4699	0.979	ng	81
11) bis(2-Chloroethyl)ether	7.53	93	3513	0.964	ng	85
12) 1,3-Dichlorobenzene	8.00	146	3904m	0.997	ng	
13) 1,4-Dichlorobenzene	8.14	146	3856	0.962	ng	# 86
14) 1,2-Dichlorobenzene	8.46	146	3497m	0.907	ng	
15) Benzyl Alcohol	8.35	79	2755	0.820	ng	# 91
16) 2,2'-oxybis(1-Chloropropan	8.62	45	6492	0.995	ng	91
18) Hexachloroethane	9.20	117	1300	0.952	ng	# 33
24) Nitrobenzene	9.33	77	4566	1.056	ng	97
31) Naphthalene	10.98	128	12381	1.081	ng	# 93
34) Hexachlorobutadiene	11.25	225	2286	0.916	ng	# 75
37) 2-Methylnaphthalene	12.58	142	9364	1.122	ng	97
40) 1,2,4,5-Tetrachlorobenzene	12.93	216	5006	0.996	ng	96
46) 1,1'-Biphenyl	13.58	154	13654	1.058	ng	98
47) 2-Chloronaphthalene	13.62	162	9768	1.000	ng	94
48) 2-Nitroaniline	13.82	65	2138	0.722	ng	99
49) Acenaphthylene	14.47	152	14071	0.958	ng	97
50) Dimethylphthalate	14.18	163	12519	1.038	ng	99
51) 2,6-Dinitrotoluene	14.31	165	1996	0.748	ng	# 87
52) Acenaphthene	14.80	154	9857	1.090	ng	86
55) Dibenzofuran	15.13	168	16084	1.073	ng	91
58) Fluorene	15.79	166	12977	1.156	ng	# 91
59) 2,3,4,6-Tetrachlorophenol	15.36	232	2215	0.707	ng	# 89
60) Diethylphthalate	15.54	149	12452	1.006	ng	99
63) Azobenzene	16.06	77	11613	0.983	ng	95
66) n-Nitrosodiphenylamine	15.99	169	10301	0.889	ng	97
67) 4-Bromophenyl-phenylether	16.67	248	3872	0.916	ng	# 76
68) Hexachlorobenzene	16.78	284	4017	0.916	ng	# 89
69) Atrazine	16.92	200	3740	0.893	ng	76
71) Phenanthrene	17.53	178	21656	1.106	ng	94
72) Anthracene	17.62	178	20460	1.065	ng	97

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)

73) Carbazole	17.89	167	17214	0.896	ng	98
74) Di-n-butylphthalate	18.43	149	18981	0.888	ng #	99
75) Fluoranthene	19.53	202	21404	0.964	ng	98
78) Pyrene	19.89	202	23435	0.996	ng	96
80) Butylbenzylphthalate	20.77	149	7263	0.754	ng	92
81) Benzo(a)anthracene	21.75	228	22850	1.074	ng	92
83) Chrysene	21.82	228	21959	1.073	ng	94
84) Bis(2-ethylhexyl)phthalate	21.63	149	10416	0.772	ng #	91
85) Di-n-octyl phthalate	22.88	149	15105	0.686	ng #	74
86) Indeno(1,2,3-cd)pyrene	28.92	276	11409	0.520	ng #	78
88) Benzo(b)fluoranthene	24.03	252	18758m	0.976	ng	
89) Benzo(k)fluoranthene	24.10	252	20329	1.080	ng	99
90) Benzo(a)pyrene	24.95	252	18690	1.023	ng	97
91) Dibenzo(a,h)anthracene	29.00	278	6288	0.373	ng	95
92) Benzo(g,h,i)perylene	30.13	276	12931m	0.759	ng	

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

