

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG070920\
 Data File : BG045992.D
 Acq On : 9 Jul 2020 13:01
 Operator : CG/JU
 Sample : L3216-02
 Misc : L00-SOIL-10PPM
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 LOQ-SOIL-02-QT3-2020

Manual Integrations
 APPROVED

mohammad
 7/10/2020 11:47:05 AM

Quant Time: Jul 09 13:39:19 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG070620.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 09 11:56:43 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.01	152	142048	20.00	ng	0.00
21) Naphthalene-d8	10.80	136	582463	20.00	ng	0.00
39) Acenaphthene-d10	14.63	164	349188	20.00	ng	0.00
64) Phenanthrene-d10	17.37	188	693489	20.00	ng	0.00
76) Chrysene-d12	21.62	240	623167	20.00	ng	0.00
86) Perylene-d12	24.78	264	617393	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.59	112	851217	97.18	ng	0.00
7) Phenol-d6	7.16	99	1247913	98.95	ng	0.00
23) Nitrobenzene-d5	9.16	82	841193	83.13	ng	0.00
42) 2,4,6-Tribromophenol	16.11	330	316730	97.73	ng	0.00
45) 2-Fluorobiphenyl	13.25	172	1648093	73.62	ng	0.00
79) Terphenyl-d14	19.99	244	1656323	59.15	ng	0.00

Target Compounds				Qvalue		
2) 1,4-Dioxane	3.51	88	47038	11.640	ng	99
3) Pyridine	3.90	79	98940	8.097	ng	99
4) n-Nitrosodimethylamine	3.82	42	40745	9.861	ng	91
6) Aniline	7.33	93	152336	9.443	ng	99
8) 2-Chlorophenol	7.57	128	86010	9.122	ng	96
9) Benzaldehyde	7.14	77	100135	12.533	ng	97
10) Phenol	7.19	94	115166	9.104	ng	99
11) bis(2-Chloroethyl)ether	7.43	93	95945	9.172	ng	97
12) 1,3-Dichlorobenzene	7.90	146	99156	9.123	ng	99
13) 1,4-Dichlorobenzene	8.04	146	99841	9.338	ng	98
14) 1,2-Dichlorobenzene	8.36	146	93156	9.053	ng	97
15) Benzyl Alcohol	8.24	79	86777	8.776	ng	94
16) 2,2'-oxybis(1-Chloropropan	8.54	45	122298	8.961	ng	98
17) 2-Methylphenol	8.44	107	80588	8.490	ng	99
18) Hexachloroethane	9.09	117	37466	9.107	ng	96
19) n-Nitroso-di-n-propylamine	8.81	70	83360	9.409	ng	97
20) 3+4-Methylphenols	8.77	107	108253	8.368	ng	97
22) Acetophenone	8.82	105	151061	9.708	ng	# 97
24) Nitrobenzene	9.20	77	115759	10.364	ng	99
25) Isophorone	9.72	82	218423	8.842	ng	99
26) 2-Nitrophenol	9.91	139	34389	9.446	ng	# 94
27) 2,4-Dimethylphenol	9.97	122	81755	9.297	ng	99
28) bis(2-Chloroethoxy)methane	10.20	93	128568	9.268	ng	95
29) 2,4-Dichlorophenol	10.45	162	78923	8.670	ng	95
30) 1,2,4-Trichlorobenzene	10.67	180	90576	8.980	ng	97
31) Naphthalene	10.86	128	290692	9.310	ng	99
32) Benzoic acid	10.03	122	32019m	6.725	ng	
33) 4-Chloroaniline	10.95	127	121544	8.711	ng	97
34) Hexachlorobutadiene	11.16	225	48294	8.201	ng	97
35) Caprolactam	11.68	113	27638	7.997	ng	91
36) 4-Chloro-3-methylphenol	12.07	107	86345	8.328	ng	99
37) 2-Methylnaphthalene	12.46	142	210649	9.490	ng	96
38) 1-Methylnaphthalene	12.68	142	197799m	9.442	ng	
40) 1,2,4,5-Tetrachlorobenzene	12.82	216	95303	9.682	ng	99

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG070920\
 Data File : BG045992.D
 Acq On : 9 Jul 2020 13:01
 Operator : CG/JU
 Sample : L3216-02
 Misc : L00-SOIL-10PPM
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 LOQ-SOIL-02-QT3-2020

Manual Integrations
 APPROVED

mohammad
 7/10/2020 11:47:05 AM

Quant Time: Jul 09 13:39:19 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG070620.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 09 11:56:43 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.81	237	50923	8.719	ng	95
43) 2,4,6-Trichlorophenol	13.06	196	53966	8.107	ng	90
44) 2,4,5-Trichlorophenol	13.12	196	62515	8.306	ng	95
46) 1,1'-Biphenyl	13.46	154	263818	10.043	ng	97
47) 2-Chloronaphthalene	13.51	162	187703	9.057	ng	98
48) 2-Nitroaniline	13.69	65	53751	11.081	ng	94
49) Acenaphthylene	14.35	152	321319	9.609	ng	96
50) Dimethylphthalate	14.08	163	230370	8.791	ng	98
51) 2,6-Dinitrotoluene	14.19	165	43790	10.769	ng	99
52) Acenaphthene	14.69	154	189690	8.993	ng	98
53) 3-Nitroaniline	14.51	138	50909	10.103	ng	# 97
54) 2,4-Dinitrophenol	14.72	184	9012	6.283	ng	# 35
55) Dibenzofuran	15.02	168	293201	9.483	ng	96
56) 4-Nitrophenol	14.82	139	36550	8.561	ng	94
57) 2,4-Dinitrotoluene	14.97	165	54886	10.924	ng	94
58) Fluorene	15.67	166	233512	9.207	ng	98
59) 2,3,4,6-Tetrachlorophenol	15.24	232	51320m	8.831	ng	
60) Diethylphthalate	15.45	149	242608	8.828	ng	96
61) 4-Chlorophenyl-phenylether	15.67	204	110441	8.752	ng	95
62) 4-Nitroaniline	15.67	138	51773	9.559	ng	94
63) Azobenzene	15.96	77	279251	9.467	ng	95
65) 4,6-Dinitro-2-methylphenol	15.74	198	16722	8.233	ng	94
66) n-Nitrosodiphenylamine	15.88	169	203601	9.474	ng	98
67) 4-Bromophenyl-phenylether	16.56	248	66305	8.561	ng	97
68) Hexachlorobenzene	16.68	284	69949	9.000	ng	99
69) Atrazine	16.82	200	66752	10.506	ng	98
70) Pentachlorophenol	17.02	266	31165	7.212	ng	99
71) Phenanthrene	17.41	178	357098	9.740	ng	97
72) Anthracene	17.50	178	363453	9.768	ng	97
73) Carbazole	17.76	167	339514	9.104	ng	96
74) Di-n-butylphthalate	18.35	149	407000	9.031	ng	# 95
75) Fluoranthene	19.42	202	405277	9.562	ng	97
77) Benzidine	19.59	184	203149	11.401	ng	98
78) Pyrene	19.78	202	415868	9.777	ng	94
80) Butylbenzylphthalate	20.68	149	160993	8.009	ng	95
81) Benzo(a)anthracene	21.61	228	381007	9.298	ng	93
82) 3,3'-Dichlorobenzidine	21.52	252	138899	9.174	ng	100
83) Chrysene	21.67	228	352336	9.026	ng	96
84) Bis(2-ethylhexyl)phthalate	21.55	149	244059	8.364	ng	# 93
85) Di-n-octyl phthalate	22.75	149	400000	7.899	ng	99
87) Indeno(1,2,3-cd)pyrene	28.36	276	400368	8.864	ng	# 69
88) Benzo(b)fluoranthene	23.78	252	363883	9.284	ng	98
89) Benzo(k)fluoranthene	23.85	252	359968	9.713	ng	98
90) Benzo(a)pyrene	24.63	252	332390	9.019	ng	97
91) Dibenzo(a,h)anthracene	28.43	278	336693	9.242	ng	98
92) Benzo(g,h,i)perylene	29.45	276	333969	9.110	ng	100

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

