

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG072720\
 Data File : BG046091.D
 Acq On : 27 Jul 2020 16:59
 Operator : CG/JU
 Sample : L1955-19
 Misc : MDL-ME-05
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 MDL-ME-SOIL-05

Manual Integrations
 APPROVED

mohammad
 7/28/2020 11:42:53 AM

Quant Time: Jul 27 17:50:04 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG072320MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Jul 27 15:44:40 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.96	152	89931	20.00	ng/ul	0.00
18) Naphthalene-d8	10.76	136	384605	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.59	164	264627	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.33	188	603080	20.00	ng/ul	0.00
78) Chrysene-d12	21.58	240	635063	20.00	ng/ul	0.00
86) Perylene-d12	24.70	264	707051	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.43	96	3771	2.07	ng/uL	0.00
5) Phenol-d5	7.13	99	207786	29.44	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.29	67	130847	29.93	ng/ul	0.00
9) 2-Chlorophenol-d4	7.50	132	169373	29.98	ng/ul	0.00
13) 4-Methylphenol-d8	8.66	113	171969	28.88	ng/ul	0.00
19) Nitrobenzene-d5	9.11	128	82572	30.40	ng/ul	0.00
22) 2-Nitrophenol-d4	9.84	143	90092	32.46	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.38	165	178801	28.46	ng/ul	0.00
29) 4-Chloroaniline-d4	10.89	131	248034	35.06	ng/ul	0.00
44) Dimethylphthalate-d6	13.99	166	613148	29.71	ng/ul	0.00
47) Acenaphthylene-d8	14.28	160	743597	30.36	ng/ul	0.00
52) 4-Nitrophenol-d4	14.77	143	95280	28.27	ng/ul	0.00
58) Fluorene-d10	15.58	176	554706	29.53	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.69	200	84942	28.14	ng/ul	0.00
71) Anthracene-d10	17.43	188	844013	30.35	ng/ul	0.00
79) Pyrene-d10	19.71	212	1036035	30.84	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.48	264	1090052	27.82	ng/ul	0.00

Target Compounds

				Ovalue	
2) 1,4-Dioxane	3.46	88	3705	1.662	ng/uL# 92
4) Benzaldehyde	7.10	77	14343m	2.851	ng/ul
6) Phenol	7.15	94	30782	4.235	ng/ul 97
8) Bis(2-Chloroethyl)ether	7.38	93	24589	4.133	ng/ul 96
10) 2-Chlorophenol	7.53	128	22596	3.946	ng/ul 94
11) 2-Methylphenol	8.40	108	21139	3.609	ng/ul 98
12) 2,2'-oxybis(1-Chloropropan	8.49	45	42779	4.254	ng/ul 97
14) Acetophenone	8.78	105	37476	4.266	ng/ul 98
15) N-Nitroso-di-n-propylamine	8.76	70	20012	4.210	ng/ul 97
16) 4-Methylphenol	8.73	108	24748	3.944	ng/ul 99
17) Hexachloroethane	9.05	117	9894	4.072	ng/ul 96
20) Nitrobenzene	9.15	77	27921	4.002	ng/ul 95
21) Isophorone	9.68	82	52479	3.849	ng/ul 98
23) 2-Nitrophenol	9.87	139	13242	4.262	ng/ul 97
24) 2,4-Dimethylphenol	9.93	107	23211	3.371	ng/ul 97
25) Bis(2-Chloroethoxy)methane	10.17	93	33461	3.880	ng/ul 97
27) 2,4-Dichlorophenol	10.41	162	25500	4.119	ng/ul 98
28) Naphthalene	10.81	128	81760	3.972	ng/ul 99
30) 4-Chloroaniline	10.91	127	33012	4.852	ng/ul 90
31) Hexachlorobutadiene	11.11	225	18604	3.905	ng/ul 98
32) Caprolactam	11.68	113	7413m	3.552	ng/ul
33) 4-Chloro-3-methylphenol	12.03	107	23128	3.643	ng/ul 99
34) 2-Methylnaphthalene	12.42	142	65022	4.202	ng/ul 96

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
35) 1-Methylnaphthalene	12.63	142	63074	4.127	ng/uL	99
37) 1,2,4,5-Tetrachlorobenzene	12.79	216	38325	4.189	ng/uL	97
38) Hexachlorocyclopentadiene	12.77	237	12402	2.368	ng/uL	98
39) 2,4,6-Trichlorophenol	13.02	196	18596	3.477	ng/uL	94
40) 2,4,5-Trichlorophenol	13.09	196	20265	3.486	ng/uL	95
41) 1,1'-Biphenyl	13.42	154	87841	4.290	ng/uL	98
42) 2-Chloronaphthalene	13.46	162	69068	4.228	ng/uL	97
43) 2-Nitroaniline	13.66	65	14661	3.807	ng/uL	94
45) Dimethylphthalate	14.04	163	85988	4.154	ng/uL	97
46) 2,6-Dinitrotoluene	14.15	165	14491	3.671	ng/uL	99
48) Acenaphthylene	14.31	152	104479	4.169	ng/uL	98
49) 3-Nitroaniline	14.48	138	12333	3.317	ng/uL#	95
50) Acenaphthene	14.65	153	73882	4.094	ng/uL	99
51) 2,4-Dinitrophenol	14.68	184	3681	2.107	ng/uL#	69
53) 4-Nitrophenol	14.79	109	9311	3.850	ng/uL	93
54) Dibenzofuran	14.98	168	108044	4.391	ng/uL	100
55) 2,4-Dinitrotoluene	14.94	165	22213	3.952	ng/uL	95
56) 2,3,4,6-Tetrachlorophenol	15.21	232	20555	3.792	ng/uL	97
57) Diethylphthalate	15.41	149	85094	4.162	ng/uL	99
59) Fluorene	15.64	166	86571	4.149	ng/uL	93
60) 4-Chlorophenyl-phenylether	15.63	204	45020	4.107	ng/uL	94
61) 4-Nitroaniline	15.64	138	13264m	3.200	ng/uL	
64) 4,6-Dinitro-2-methylphenol	15.70	198	12836	3.992	ng/uL#	87
65) N-Nitrosodiphenylamine	15.83	169	74805	4.266	ng/uL	99
66) 4-Bromophenyl-phenylether	16.52	248	31704	4.230	ng/uL	98
67) Hexachlorobenzene	16.64	284	35852	4.238	ng/uL	96
68) Atrazine	16.79	200	28335	4.042	ng/uL	99
69) Pentachlorophenol	16.98	266	13425m	2.977	ng/uL	
70) Phenanthrene	17.37	178	147725	4.389	ng/uL	99
72) Anthracene	17.46	178	145031	4.325	ng/uL	99
73) 1,2,3,4-Tetrachlorobenzene	13.38	216	40040	4.421	ng/uL	91
74) Pentachlorobenzene	14.91	250	38684	4.230	ng/uL	97
75) Carbazole	17.72	167	120858	4.491	ng/uL	99
76) Di-n-butylphthalate	18.30	149	134728	4.069	ng/uL	98
77) Fluoranthene	19.38	202	184485	4.997	ng/uL	98
80) Pvrene	19.74	202	188215	4.383	ng/uL	99
81) Butylbenzylphthalate	20.64	149	56675	3.845	ng/uL	93
82) 3,3'-Dichlorobenzidine	21.48	252	53210	3.904	ng/uL	99
83) Benzo(a)anthracene	21.56	228	190135	4.546	ng/uL	98
84) Bis(2-ethylhexyl)phthalate	21.49	149	82113	3.830	ng/uL	95
85) Chrysene	21.62	228	183079	4.503	ng/uL	99
87) Di-n-octyl phthalate	22.67	149	137193	3.713	ng/uL	100
88) Benzo(b)fluoranthene	23.70	252	186714	3.909	ng/uL	99
89) Benzo(k)fluoranthene	23.77	252	176383	3.761	ng/uL	100
91) Benzo(a)pyrene	24.54	252	185613m	4.237	ng/uL	
92) Indeno(1,2,3-cd)pyrene	28.20	276	200084	3.658	ng/uL	99
93) Dibenzo(a,h)anthracene	28.26	278	165582	3.700	ng/uL	98
94) Benzo(a,h,i)perylene	29.30	276	175369	3.793	ng/uL	96

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

