

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG013023\
 Data File : BG056471.D
 Acq On : 30 Jan 2023 15:26
 Operator : CG/JU
 Sample : N3980-02
 Misc : LOQ-SOIL-10ppm
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 LOQ-SOIL-02-QT3-2022

Quant Time: Jan 31 03:01:47 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG012723.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Jan 28 00:39:13 2023
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By :Christian
 Giraldo
 01/31/2023
 Supervised By :Jagrut
 Upadhyay
 01/31/2023

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.277	152	34209	20.000	ng	0.00
21) Naphthalene-d8	11.108	136	131501	20.000	ng	0.00
39) Acenaphthene-d10	14.898	164	109085	20.000	ng	0.00
64) Phenanthrene-d10	17.636	188	279855	20.000	ng	0.00
76) Chrysene-d12	21.925	240	273909	20.000	ng	0.00
86) Perylene-d12	25.345	264	297454	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.797	112	199802	107.443	ng	0.00
7) Phenol-d6	7.419	99	288964	104.875	ng	0.00
23) Nitrobenzene-d5	9.452	82	168344	72.695	ng	0.00
42) 2,4,6-Tribromophenol	16.379	330	153237	105.665	ng	0.00
45) 2-Fluorobiphenyl	13.523	172	570667	72.529	ng	0.00
79) Terphenyl-d14	20.210	244	1102513	70.696	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.600	88	7512	9.701	ng	# 97
3) Pyridine	4.023	79	19308	8.365	ng	97
4) n-Nitrosodimethylamine	3.929	42	4215	8.587	ng	# 89
6) Aniline	7.589	93	32306	8.955	ng	# 93
8) 2-Chlorophenol	7.836	128	17859	8.742	ng	87
9) Benzaldehyde	7.407	77	15788	9.885	ng	92
10) Phenol	7.448	94	26585	9.493	ng	96
11) bis(2-Chloroethyl)ether	7.677	93	18226	9.469	ng	93
12) 1,3-Dichlorobenzene	8.165	146	23852	10.141	ng	97
13) 1,4-Dichlorobenzene	8.312	146	23815	9.998	ng	95
14) 1,2-Dichlorobenzene	8.635	146	22734	9.837	ng	98
15) Benzyl Alcohol	8.512	79	16330	8.638	ng	95
16) 2,2'-oxybis(1-Chloropr...	8.788	45	3909	9.584	ng	96
17) 2-Methylphenol	8.717	107	16890	7.928	ng	91
18) Hexachloroethane	9.375	117	7090	9.841	ng	91
19) n-Nitroso-di-n-propyla...	9.076	70	14712	9.265	ng	# 97
20) 3+4-Methylphenols	9.046	107	25120	8.413	ng	97
22) Acetophenone	9.105	105	33708	9.647	ng	# 96
24) Nitrobenzene	9.499	77	21004	9.403	ng	93
25) Isophorone	10.016	82	42916	8.983	ng	99
26) 2-Nitrophenol	10.215	139	12364	9.123	ng	94
27) 2,4-Dimethylphenol	10.262	122	19443m	8.572	ng	
28) bis(2-Chloroethoxy)met...	10.492	93	24790	9.393	ng	# 97
29) 2,4-Dichlorophenol	10.756	162	22712	8.896	ng	96
30) 1,2,4-Trichlorobenzene	10.973	180	28469	10.034	ng	92
31) Naphthalene	11.161	128	65951	9.502	ng	99
32) Benzoic acid	10.374	122	6711m	5.289	ng	
33) 4-Chloroaniline	11.261	127	28928	8.929	ng	98
34) Hexachlorobutadiene	11.438	225	20108	10.097	ng	92
35) Caprolactam	12.013	113	7534	8.750	ng	86
36) 4-Chloro-3-methylphenol	12.366	107	22633	9.180	ng	98
37) 2-Methylnaphthalene	12.748	142	51669	9.030	ng	96
38) 1-Methylnaphthalene	12.965	142	49587	9.279	ng	98
40) 1,2,4,5-Tetrachloroben...	13.106	216	38860	9.751	ng	97
41) Hexachlorocyclopentadiene	13.083	237	11704	11.470	ng	95
43) 2,4,6-Trichlorophenol	13.341	196	22618	8.807	ng	92

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44) 2,4,5-Trichlorophenol	13.418	196	25320	8.421	ng	97
46) 1,1'-Biphenyl	13.735	154	75842	9.585	ng	97
47) 2-Chloronaphthalene	13.782	162	59690	9.652	ng	98
48) 2-Nitroaniline	13.982	65	11215	8.772	ng	87
49) Acenaphthylene	14.622	152	92060	9.149	ng	99
50) Dimethylphthalate	14.334	163	79122	8.965	ng	100
51) 2,6-Dinitrotoluene	14.463	165	16975	8.820	ng	98
52) Acenaphthene	14.963	154	53992	9.049	ng	93
53) 3-Nitroaniline	14.798	138	15579	8.350	ng	99
54) 2,4-Dinitrophenol	15.010	184	10051	8.330	ng	95
55) Dibenzofuran	15.292	168	98786	9.232	ng	98
56) 4-Nitrophenol	15.104	139	10064m	7.894	ng	
57) 2,4-Dinitrotoluene	15.251	165	24131	8.901	ng	# 97
58) Fluorene	15.938	166	78965	9.274	ng	100
59) 2,3,4,6-Tetrachlorophenol	15.521	232	22360	8.364	ng	# 98
60) Diethylphthalate	15.685	149	76320	9.210	ng	99
61) 4-Chlorophenyl-phenyle...	15.921	204	48199	9.277	ng	99
62) 4-Nitroaniline	15.950	138	16624	8.904	ng	98
63) Azobenzene	16.214	77	53279	9.344	ng	97
65) 4,6-Dinitro-2-methylph...	16.015	198	19764	10.045	ng	94
66) n-Nitrosodiphenylamine	16.132	169	72747	9.181	ng	97
67) 4-Bromophenyl-phenylether	16.814	248	32389	9.041	ng	96
68) Hexachlorobenzene	16.943	284	33614	9.580	ng	99
69) Atrazine	17.066	200	31071	10.165	ng	99
70) Pentachlorophenol	17.290	266	21359m	10.060	ng	
71) Phenanthrene	17.677	178	136830	9.342	ng	99
72) Anthracene	17.765	178	141491	9.410	ng	98
73) Carbazole	18.036	167	121499	9.488	ng	98
74) Di-n-butylphthalate	18.559	149	120554	8.966	ng	99
75) Fluoranthene	19.669	202	176338	9.539	ng	99
77) Benzidine	19.839	184	95271	12.853	ng	97
78) Pyrene	20.027	202	179960	9.238	ng	100
80) Butylbenzylphthalate	20.885	149	53864	9.053	ng	98
81) Benzo(a)anthracene	21.902	228	178777	9.336	ng	99
82) 3,3'-Dichlorobenzidine	21.802	252	73550	10.666	ng	99
83) Chrysene	21.972	228	167808	9.328	ng	99
84) Bis(2-ethylhexyl)phtha...	21.761	149	75758	8.881	ng	# 98
85) Di-n-octyl phthalate	23.030	149	128691	9.059	ng	100
87) Indeno(1,2,3-cd)pyrene	29.281	276	200032	9.187	ng	100
88) Benzo(b)fluoranthene	24.246	252	182275	9.873	ng	99
89) Benzo(k)fluoranthene	24.317	252	185917	9.974	ng	98
90) Benzo(a)pyrene	25.186	252	163429	8.986	ng	95
91) Dibenzo(a,h)anthracene	29.334	278	175339	9.593	ng	99
92) Benzo(g,h,i)perylene	30.509	276	167980	9.525	ng	96

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

