

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG013023\
 Data File : BG056469.D
 Acq On : 30 Jan 2023 11:31
 Operator : CG/JU
 Sample : N3980-02
 Misc : LOQ-SOIL-5ppm
 ALS Vial : 4 Sample Multiplier: 1

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

Manual Integrations
 APPROVED

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

Quant Time: Jan 31 03:00:27 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

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.276	152	29710	20.000	ng	0.00
21) Naphthalene-d8	11.113	136	121869	20.000	ng	0.00
39) Acenaphthene-d10	14.897	164	101378	20.000	ng	0.00
64) Phenanthrene-d10	17.635	188	265860	20.000	ng	0.00
76) Chrysene-d12	21.924	240	260219	20.000	ng	0.00
86) Perylene-d12	25.344	264	285878	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.796	112	201475	124.749	ng	0.00
7) Phenol-d6	7.418	99	283647	118.534	ng	0.00
23) Nitrobenzene-d5	9.451	82	159476	74.308	ng	0.00
42) 2,4,6-Tribromophenol	16.378	330	154915	114.943	ng	0.00
45) 2-Fluorobiphenyl	13.522	172	548690	75.038	ng	0.00
79) Terphenyl-d14	20.215	244	1062962	71.745	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.599	88	4048	6.019	ng	94
3) Pyridine	4.028	79	8813	4.396	ng	# 93
4) n-Nitrosodimethylamine	3.934	42	2270	5.325	ng	# 78
6) Aniline	7.594	93	16496	5.265	ng	96
8) 2-Chlorophenol	7.835	128	9659	5.444	ng	91
9) Benzaldehyde	7.406	77	8026	5.441	ng	95
10) Phenol	7.447	94	12583	5.174	ng	90
11) bis(2-Chloroethyl)ether	7.682	93	9317	5.573	ng	88
12) 1,3-Dichlorobenzene	8.170	146	10874	5.323	ng	97
13) 1,4-Dichlorobenzene	8.317	146	11824	5.716	ng	96
14) 1,2-Dichlorobenzene	8.640	146	11024	5.492	ng	91
15) Benzyl Alcohol	8.516	79	7881	4.800	ng	94
16) 2,2'-oxybis(1-Chloropr...	8.799	45	1783	5.034	ng	87
17) 2-Methylphenol	8.716	107	8608	4.652	ng	96
18) Hexachloroethane	9.374	117	3490	5.578	ng	# 85
19) n-Nitroso-di-n-propyla...	9.075	70	6889	4.996	ng	# 91
20) 3+4-Methylphenols	9.051	107	11730	4.524	ng	93
22) Acetophenone	9.104	105	16522	5.102	ng	# 97
24) Nitrobenzene	9.498	77	10254	4.953	ng	# 91
25) Isophorone	10.015	82	21450	4.845	ng	# 91
26) 2-Nitrophenol	10.214	139	6358	5.062	ng	97
27) 2,4-Dimethylphenol	10.262	122	9214	4.383	ng	92
28) bis(2-Chloroethoxy)met...	10.491	93	12935	5.288	ng	# 96
29) 2,4-Dichlorophenol	10.755	162	11439	4.835	ng	93
30) 1,2,4-Trichlorobenzene	10.967	180	13731	5.222	ng	# 93
31) Naphthalene	11.160	128	33190	5.160	ng	97
33) 4-Chloroaniline	11.260	127	15122	5.036	ng	94
34) Hexachlorobutadiene	11.437	225	9652	5.230	ng	95
35) Caprolactam	12.007	113	3474	4.354	ng	# 78
36) 4-Chloro-3-methylphenol	12.365	107	11065	4.842	ng	91
37) 2-Methylnaphthalene	12.747	142	25179	4.748	ng	94
38) 1-Methylnaphthalene	12.964	142	24986m	5.045	ng	
40) 1,2,4,5-Tetrachloroben...	13.105	216	18626	5.029	ng	98
41) Hexachlorocyclopentadiene	13.082	237	3983	8.239	ng	85
43) 2,4,6-Trichlorophenol	13.346	196	10960	4.592	ng	92
44) 2,4,5-Trichlorophenol	13.422	196	12240	4.380	ng	98

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
46) 1,1'-Biphenyl	13.734	154	37801	5.141	ng	99
47) 2-Chloronaphthalene	13.781	162	30274	5.267	ng	97
48) 2-Nitroaniline	13.981	65	5957	5.014	ng	95
49) Acenaphthylene	14.621	152	46649	4.989	ng	98
50) Dimethylphthalate	14.333	163	40392	4.925	ng	98
51) 2,6-Dinitrotoluene	14.462	165	8466	4.733	ng	98
52) Acenaphthene	14.956	154	28010	5.051	ng	94
53) 3-Nitroaniline	14.797	138	7963	4.592	ng	94
55) Dibenzofuran	15.291	168	50547	5.083	ng	97
56) 4-Nitrophenol	15.115	139	4489m	3.789	ng	
57) 2,4-Dinitrotoluene	15.250	165	11967	4.750	ng	# 99
58) Fluorene	15.937	166	40464	5.113	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.520	232	11097	4.466	ng	# 99
60) Diethylphthalate	15.685	149	38606	5.013	ng	95
61) 4-Chlorophenyl-phenyle...	15.920	204	23870	4.943	ng	97
62) 4-Nitroaniline	15.949	138	8126	4.683	ng	94
63) Azobenzene	16.213	77	27407	5.172	ng	98
65) 4,6-Dinitro-2-methylph...	16.014	198	5740	3.071	ng	92
66) n-Nitrosodiphenylamine	16.131	169	35957	4.777	ng	97
67) 4-Bromophenyl-phenylether	16.813	248	16439	4.830	ng	94
68) Hexachlorobenzene	16.942	284	17421	5.226	ng	97
69) Atrazine	17.065	200	15150	5.217	ng	99
71) Phenanthrene	17.676	178	70910	5.096	ng	96
72) Anthracene	17.770	178	71466	5.003	ng	99
73) Carbazole	18.035	167	60091	4.940	ng	98
74) Di-n-butylphthalate	18.564	149	61287	4.798	ng	# 96
75) Fluoranthene	19.668	202	90043	5.127	ng	99
77) Benzidine	19.839	184	44289	6.289	ng	99
78) Pyrene	20.032	202	91509	4.944	ng	97
80) Butylbenzylphthalate	20.884	149	25975	4.596	ng	99
81) Benzo(a)anthracene	21.901	228	90492	4.974	ng	98
82) 3,3'-Dichlorobenzidine	21.807	252	35746	5.456	ng	98
83) Chrysene	21.971	228	86017	5.033	ng	98
84) Bis(2-ethylhexyl)phtha...	21.766	149	39696	4.899	ng	97
85) Di-n-octyl phthalate	23.035	149	64958	4.813	ng	99
87) Indeno(1,2,3-cd)pyrene	29.280	276	100133	4.785	ng	# 91
88) Benzo(b)fluoranthene	24.245	252	91271	5.144	ng	96
89) Benzo(k)fluoranthene	24.316	252	92223	5.148	ng	# 98
90) Benzo(a)pyrene	25.173	252	82143	4.700	ng	100
91) Dibenzo(a,h)anthracene	29.333	278	89058	5.070	ng	96
92) Benzo(g,h,i)perylene	30.508	276	85790	5.062	ng	# 96

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

