

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090324\
 Data File : BG062622.D
 Acq On : 3 Sep 2024 12:29
 Operator : MA/JU
 Sample : P3390-02
 Misc : LOQ--SOIL-10ng
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 LOQ-SOIL-02-QT3-2024

Quant Time: Sep 04 04:53:33 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG083024.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Aug 31 02:36:01 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.924	152	61755	20.000	ng	0.00
21) Naphthalene-d8	10.715	136	275764	20.000	ng	0.00
39) Acenaphthene-d10	14.546	164	227463	20.000	ng	0.00
64) Phenanthrene-d10	17.290	188	597466	20.000	ng	0.00
76) Chrysene-d12	21.538	240	609044	20.000	ng	0.00
86) Perylene-d12	24.605	264	663688	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.504	112	490533	137.301	ng	0.00
7) Phenol-d6	7.090	99	690042	133.697	ng	0.00
23) Nitrobenzene-d5	9.076	82	499928	97.635	ng	0.00
42) 2,4,6-Tribromophenol	16.038	330	467654	146.091	ng	0.00
45) 2-Fluorobiphenyl	13.171	172	1320443	94.030	ng	0.00
79) Terphenyl-d14	19.910	244	2635508	85.803	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.412	88	19165	12.877	ng	97
3) Pyridine	3.806	79	48643	11.343	ng	96
4) n-Nitrosodimethylamine	3.717	42	24949	11.967	ng	# 99
6) Aniline	7.249	93	64205	13.315	ng	# 89
8) 2-Chlorophenol	7.490	128	43678	11.387	ng	90
9) Benzaldehyde	7.061	77	44426	14.798	ng	97
10) Phenol	7.114	94	59537	11.395	ng	96
11) bis(2-Chloroethyl)ether	7.343	93	46335	11.583	ng	98
12) 1,3-Dichlorobenzene	7.813	146	48332	11.482	ng	97
13) 1,4-Dichlorobenzene	7.960	146	50216	11.592	ng	98
14) 1,2-Dichlorobenzene	8.277	146	47643	11.189	ng	97
15) Benzyl Alcohol	8.159	79	43504	11.392	ng	95
16) 2,2'-oxybis(1-Chloropr...	8.441	45	85033	11.153	ng	99
17) 2-Methylphenol	8.365	107	37562	10.313	ng	94
18) Hexachloroethane	9.000	117	17741	10.923	ng	96
19) n-Nitroso-di-n-propyla...	8.717	70	45326	11.145	ng	92
20) 3+4-Methylphenols	8.688	107	52819	9.830	ng	98
22) Acetophenone	8.735	105	84948	11.419	ng	# 97
24) Nitrobenzene	9.117	77	62304	11.444	ng	98
25) Isophorone	9.640	82	130853	11.642	ng	99
26) 2-Nitrophenol	9.828	139	22716	11.043	ng	90
27) 2,4-Dimethylphenol	9.887	122	20707	6.870	ng	93
28) bis(2-Chloroethoxy)met...	10.122	93	69515	11.553	ng	99
29) 2,4-Dichlorophenol	10.363	162	46319	10.991	ng	97
30) 1,2,4-Trichlorobenzene	10.580	180	57164	11.442	ng	97
31) Naphthalene	10.768	128	155574	11.382	ng	99
32) Benzoic acid	9.992	122	50430	15.556	ng	97
33) 4-Chloroaniline	10.874	127	65509	12.179	ng	97
34) Hexachlorobutadiene	11.056	225	37433	10.728	ng	94
35) Caprolactam	11.626	113	17991	11.074	ng	92
36) 4-Chloro-3-methylphenol	11.996	107	54534	10.728	ng	99
37) 2-Methylnaphthalene	12.372	142	119941	11.314	ng	95
38) 1-Methylnaphthalene	12.589	142	116742	10.967	ng	96
40) 1,2,4,5-Tetrachloroben...	12.742	216	80605	11.783	ng	98
41) Hexachlorocyclopentadiene	12.725	237	21121	10.822	ng	92
43) 2,4,6-Trichlorophenol	12.977	196	49675	11.455	ng	97

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.048	196	53302	10.854	ng	96
46) 1,1'-Biphenyl	13.383	154	178840	11.785	ng	99
47) 2-Chloronaphthalene	13.424	162	139980	11.382	ng	99
48) 2-Nitroaniline	13.624	65	34645	10.710	ng	97
49) Acenaphthylene	14.270	152	224114	12.305	ng	99
50) Dimethylphthalate	14.000	163	192738	11.588	ng	98
51) 2,6-Dinitrotoluene	14.117	165	33452	10.633	ng	97
52) Acenaphthene	14.611	154	130784	11.487	ng	97
53) 3-Nitroaniline	14.446	138	33913	11.140	ng	97
54) 2,4-Dinitrophenol	14.652	184	22677	12.844	ng	91
55) Dibenzofuran	14.945	168	231497	11.856	ng	96
56) 4-Nitrophenol	14.757	139	63503	24.615	ng	94
57) 2,4-Dinitrotoluene	14.904	165	45543	10.610	ng	97
58) Fluorene	15.598	166	178623	11.657	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.175	232	54920	11.805	ng	# 98
60) Diethylphthalate	15.363	149	187657	11.383	ng	98
61) 4-Chlorophenyl-phenyle...	15.592	204	102887	11.696	ng	95
62) 4-Nitroaniline	15.604	138	37827	11.270	ng	90
63) Azobenzene	15.880	77	179315	11.584	ng	99
65) 4,6-Dinitro-2-methylph...	15.668	198	26085	9.783	ng	89
66) n-Nitrosodiphenylamine	15.803	169	163070	11.352	ng	99
67) 4-Bromophenyl-phenylether	16.485	248	69998	11.352	ng	95
68) Hexachlorobenzene	16.596	284	84858	11.620	ng	96
69) Atrazine	16.749	200	72112	14.780	ng	97
70) Pentachlorophenol	16.943	266	100040	21.389	ng	97
71) Phenanthrene	17.331	178	339325	12.168	ng	99
72) Anthracene	17.425	178	323625	11.803	ng	97
73) Carbazole	17.689	167	295543	11.725	ng	99
74) Di-n-butylphthalate	18.259	149	322060	11.329	ng	100
75) Fluoranthene	19.346	202	417690	11.843	ng	99
77) Benzidine	19.522	184	114917	10.469	ng	98
78) Pyrene	19.705	202	439108	11.342	ng	99
80) Butylbenzylphthalate	20.598	149	116834	10.274	ng	97
81) Benzo(a)anthracene	21.514	228	450402	11.725	ng	98
82) 3,3'-Dichlorobenzidine	21.438	252	153915	12.299	ng	98
83) Chrysene	21.579	228	425569	11.566	ng	100
84) Bis(2-ethylhexyl)phtha...	21.438	149	185597	10.743	ng	99
85) Di-n-octyl phthalate	22.601	149	314453	10.396	ng	99
87) Indeno(1,2,3-cd)pyrene	28.071	276	507740	11.931	ng	99
88) Benzo(b)fluoranthene	23.635	252	420532	11.472	ng	98
89) Benzo(k)fluoranthene	23.700	252	430477	12.005	ng	99
90) Benzo(a)pyrene	24.464	252	386712	11.941	ng	99
91) Dibenzo(a,h)anthracene	28.124	278	416025	11.867	ng	97
92) Benzo(g,h,i)perylene	29.158	276	396358	11.212	ng	95

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

