

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM032122\
 Data File : BM034292.D
 Acq On : 21 Mar 2022 17:40
 Operator : CG/JU
 Sample : M4068-02
 Misc : LOQ-SOIL-5ppm
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 LOQ-SOIL-02-QT4-2021

Quant Time: Mar 22 04:09:59 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM031522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 16 02:38:26 2022
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By :Christian
 Giraldo

Supervised By :Jagrut
 Upadhyay
 03/22/2022

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.942	152	193576	20.000	ng	0.00
21) Naphthalene-d8	10.754	136	792805	20.000	ng	-0.01
39) Acenaphthene-d10	14.583	164	555707	20.000	ng	0.00
64) Phenanthrene-d10	17.324	188	1180612	20.000	ng	0.00
76) Chrysene-d12	21.495	240	1185763	20.000	ng	0.00
86) Perylene-d12	23.906	264	1136280	20.000	ng	0.00

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System Monitoring Compounds						
5) 2-Fluorophenol	5.496	112	1122721	104.048	ng	0.00
7) Phenol-d6	7.101	99	1573351	101.839	ng	0.00
23) Nitrobenzene-d5	9.107	82	1193387	73.287	ng	0.00
42) 2,4,6-Tribromophenol	16.065	330	806725	107.704	ng	0.00
45) 2-Fluorobiphenyl	13.213	172	2793960	68.704	ng	0.00
79) Terphenyl-d14	19.942	244	4793580	70.245	ng	0.00

Target Compounds						Qvalue
2) 1,4-Dioxane	3.355	88	22958	4.703	ng	94
3) Pyridine	3.772	79	52717	3.990	ng	# 92
4) n-Nitrosodimethylamine	3.672	42	29823	4.754	ng	# 93
6) Aniline	7.266	93	90552	5.129	ng	98
8) 2-Chlorophenol	7.501	128	59392	4.706	ng	94
9) Benzaldehyde	7.072	77	57146m	6.866	ng	
10) Phenol	7.125	94	71591	4.659	ng	89
11) bis(2-Chloroethyl)ether	7.360	93	61569	4.882	ng	98
12) 1,3-Dichlorobenzene	7.831	146	69298	4.841	ng	97
13) 1,4-Dichlorobenzene	7.978	146	70003	4.776	ng	99
14) 1,2-Dichlorobenzene	8.301	146	65938	4.615	ng	96
15) Benzyl Alcohol	8.178	79	51991	4.229	ng	96
16) 2,2'-oxybis(1-Chloropr...	8.484	45	86406	4.992	ng	99
17) 2-Methylphenol	8.384	107	47063	4.373	ng	96
18) Hexachloroethane	9.037	117	24445	4.522	ng	95
19) n-Nitroso-di-n-propyla...	8.754	70	52325	4.682	ng	100
20) 3+4-Methylphenols	8.713	107	63300	4.253	ng	99
22) Acetophenone	8.766	105	97036	4.769	ng	97
24) Nitrobenzene	9.148	77	79934	5.258	ng	98
25) Isophorone	9.678	82	137179	5.034	ng	99
26) 2-Nitrophenol	9.866	139	27692	4.107	ng	96
27) 2,4-Dimethylphenol	9.925	122	51638	4.891	ng	98
28) bis(2-Chloroethoxy)met...	10.160	93	81239	4.657	ng	99
29) 2,4-Dichlorophenol	10.401	162	50732	4.136	ng	97
30) 1,2,4-Trichlorobenzene	10.619	180	65349	4.687	ng	98
31) Naphthalene	10.807	128	200206	4.847	ng	100
32) Benzoic acid	9.984	122	16529	1.893	ng	93
33) 4-Chloroaniline	10.913	127	69821	4.285	ng	96
34) Hexachlorobutadiene	11.101	225	41046	4.648	ng	94
35) Caprolactam	11.666	113	17385	4.077	ng	91
36) 4-Chloro-3-methylphenol	12.036	107	59655	4.244	ng	98
37) 2-Methylnaphthalene	12.419	142	141335	4.796	ng	96
38) 1-Methylnaphthalene	12.636	142	137126	4.756	ng	97
40) 1,2,4,5-Tetrachloroben...	12.783	216	76254	4.547	ng	100
41) Hexachlorocyclopentadiene	12.772	237	40991	4.221	ng	100
43) 2,4,6-Trichlorophenol	13.019	196	45635	3.942	ng	99

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44) 2,4,5-Trichlorophenol	13.089	196	48271	3.884	ng	96
46) 1,1'-Biphenyl	13.419	154	199511	4.709	ng	98
47) 2-Chloronaphthalene	13.460	162	147017	4.520	ng	99
48) 2-Nitroaniline	13.660	65	37439	3.765	ng	94
49) Acenaphthylene	14.307	152	239298	4.721	ng	99
50) Dimethylphthalate	14.036	163	192365	4.548	ng	100
51) 2,6-Dinitrotoluene	14.154	165	38391	4.425	ng	99
52) Acenaphthene	14.648	154	150446	4.406	ng	98
53) 3-Nitroaniline	14.483	138	31306	3.597	ng	# 99
54) 2,4-Dinitrophenol	14.683	184	12202	2.510	ng	94
55) Dibenzofuran	14.977	168	232825	4.576	ng	98
56) 4-Nitrophenol	14.783	139	18791	2.663	ng	93
57) 2,4-Dinitrotoluene	14.936	165	49120	4.189	ng	96
58) Fluorene	15.624	166	190098	4.602	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.201	232	45285	4.041	ng	96
60) Diethylphthalate	15.395	149	200160	4.613	ng	99
61) 4-Chlorophenyl-phenyle...	15.618	204	94886	4.499	ng	96
62) 4-Nitroaniline	15.636	138	30378m	3.628	ng	
63) Azobenzene	15.913	77	189491	4.529	ng	99
65) 4,6-Dinitro-2-methylph...	15.695	198	24436	3.183	ng	91
66) n-Nitrosodiphenylamine	15.830	169	165025	4.400	ng	98
67) 4-Bromophenyl-phenylether	16.513	248	58212	4.289	ng	100
68) Hexachlorobenzene	16.636	284	69322	4.579	ng	99
69) Atrazine	16.777	200	59304	4.797	ng	99
70) Pentachlorophenol	16.971	266	35070	3.623	ng	98
71) Phenanthrene	17.365	178	304412	4.604	ng	100
72) Anthracene	17.454	178	285966	4.444	ng	99
73) Carbazole	17.724	167	239016	3.982	ng	99
74) Di-n-butylphthalate	18.289	149	321599	4.310	ng	99
75) Fluoranthene	19.377	202	343329	4.632	ng	99
77) Benzidine	19.559	184	81976m	4.718	ng	
78) Pyrene	19.736	202	365652	4.377	ng	99
80) Butylbenzylphthalate	20.630	149	135116	3.824	ng	99
81) Benzo(a)anthracene	21.477	228	343517	4.590	ng	98
82) 3,3'-Dichlorobenzidine	21.412	252	97367	3.895	ng	98
83) Chrysene	21.530	228	336273	4.400	ng	99
84) Bis(2-ethylhexyl)phtha...	21.406	149	208123	3.989	ng	99
85) Di-n-octyl phthalate	22.336	149	337278	4.072	ng	99
87) Indeno(1,2,3-cd)pyrene	26.418	276	263501	3.605	ng	98
88) Benzo(b)fluoranthene	23.171	252	313697	4.617	ng	99
89) Benzo(k)fluoranthene	23.218	252	337785	4.693	ng	99
90) Benzo(a)pyrene	23.800	252	279620	4.854	ng	97
91) Dibenzo(a,h)anthracene	26.435	278	213567	3.603	ng	98
92) Benzo(g,h,i)perylene	27.194	276	243027	3.962	ng	97

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

