

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM122118\
 Data File : BM018374.D
 Acq On : 22 Dec 2018 01:30
 Operator : JU/SJ
 Sample : J6500-02MS
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SOIL-CUTTINGSMS

Manual Integrations
 APPROVED

Sohil
 12/26/2018 3:20:23 PM

Quant Time: Dec 22 05:56:27 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM121518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Dec 15 21:39:37 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.48	152	103314	20.00	ng	-0.02
21) Naphthalene-d8	10.25	136	418890	20.00	ng	-0.02
39) Acenaphthene-d10	14.15	164	224893	20.00	ng	-0.02
64) Phenanthrene-d10	16.91	188	402205	20.00	ng	-0.01
76) Chrysene-d12	21.13	240	357996	20.00	ng	-0.01
87) Perylene-d12	23.30	264	415924	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.13	112	526862	85.19	ng	0.00
7) Phenol-d6	6.70	99	702813	81.76	ng	0.00
23) Nitrobenzene-d5	8.64	82	475582	58.24	ng	-0.02
42) 2,4,6-Tribromophenol	15.66	330	200037	60.60	ng	-0.01
45) 2-Fluorobiphenyl	12.75	172	913147	52.59	ng	-0.02
79) Terphenyl-d14	19.56	244	921501	58.22	ng	-0.01

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.12	88	51786	21.138	ng	97
3) Pyridine	3.50	79	177497	24.627	ng	95
4) n-Nitrosodimethylamine	3.42	42	77888	31.396	ng	92
6) Aniline	6.83	93	172571	16.003	ng	97
8) 2-Chlorophenol	7.06	128	181350	24.818	ng	92
9) Benzaldehyde	6.65	77	150917	28.803	ng	88
10) Phenol	6.73	94	295139	32.419	ng	94
11) bis(2-Chloroethyl)ether	6.93	93	168645	23.309	ng	97
12) 1,3-Dichlorobenzene	7.37	146	191707	23.494	ng	95
13) 1,4-Dichlorobenzene	7.52	146	192444	23.225	ng	97
14) 1,2-Dichlorobenzene	7.82	146	186486	23.346	ng	97
15) Benzyl Alcohol	7.74	79	196772	28.092	ng	93
16) 2,2'-oxybis(1-Chloropropan	8.00	45	128750	19.774	ng	87
17) 2-Methylphenol	7.95	107	157044	25.835	ng	90
18) Hexachloroethane	8.53	117	56487	18.599	ng	92
19) n-Nitroso-di-n-propylamine	8.29	70	155588	24.241	ng	97
20) 3+4-Methylphenols	8.28	107	207194	24.325	ng	94
22) Acetophenone	8.30	105	281073	25.356	ng	94
24) Nitrobenzene	8.68	77	232723	28.520	ng	93
25) Isophorone	9.20	82	391488	28.794	ng	96
26) 2-Nitrophenol	9.38	139	57452	14.355	ng	# 88
27) 2,4-Dimethylphenol	9.46	122	161913	28.071	ng	95
28) bis(2-Chloroethoxy)methane	9.69	93	225350	25.427	ng	97
29) 2,4-Dichlorophenol	9.92	162	164238	25.681	ng	98
30) 1,2,4-Trichlorobenzene	10.12	180	187942	25.836	ng	96
31) Naphthalene	10.30	128	530627	25.906	ng	99
32) Benzoic acid	9.59	122	38528m	7.054	ng	
33) 4-Chloroaniline	10.43	127	140695	15.681	ng	93
34) Hexachlorobutadiene	10.56	225	112973	24.555	ng	98
35) Caprolactam	11.24	113	48310	19.824	ng	89
36) 4-Chloro-3-methylphenol	11.59	107	185006	24.075	ng	91
37) 2-Methylnaphthalene	11.92	142	381606	26.418	ng	96
38) 1-Methylnaphthalene	12.15	142	360219	25.556	ng	# 98
40) 1,2,4,5-Tetrachlorobenzene	12.30	216	203066	26.258	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.26	237	6108	9.711	nq	92
43) 2,4,6-Trichlorophenol	12.57	196	125746	25.281	nq	93
44) 2,4,5-Trichlorophenol	12.65	196	126184	23.896	nq	95
46) 1,1'-Biphenyl	12.96	154	484714	23.958	nq	98
47) 2-Chloronaphthalene	13.00	162	372836	25.037	nq	97
48) 2-Nitroaniline	13.24	65	135470	29.532	nq	90
49) Acenaphthylene	13.86	152	577169	25.720	nq	99
50) Dimethylphthalate	13.62	163	554091	29.635	nq	99
51) 2,6-Dinitrotoluene	13.75	165	83439	20.301	nq #	81
52) Acenaphthene	14.21	154	340141	24.802	nq	98
53) 3-Nitroaniline	14.09	138	103432	23.913	nq	89
54) 2,4-Dinitrophenol	14.33	184	128	0.056	nq #	33
55) Dibenzofuran	14.55	168	539104	24.464	nq	96
56) 4-Nitrophenol	14.45	139	93709	28.575	nq #	69
57) 2,4-Dinitrotoluene	14.56	165	110401	19.432	nq #	86
58) Fluorene	15.20	166	424050	24.915	nq	99
59) 2,3,4,6-Tetrachlorophenol	14.80	232	101487	21.340	nq	95
60) Diethylphthalate	14.99	149	464917	23.042	nq	98
61) 4-Chlorophenyl-phenylether	15.20	204	223279	23.196	nq	98
62) 4-Nitroaniline	15.26	138	95134	23.180	nq #	73
63) Azobenzene	15.50	77	480911	26.139	nq	94
65) 4,6-Dinitro-2-methylphenol	15.34	198	1085	0.366	nq #	22
66) n-Nitrosodiphenylamine	15.43	169	365724	27.498	nq	98
67) 4-Bromophenyl-phenylether	16.10	248	131032	26.900	nq	96
68) Hexachlorobenzene	16.21	284	133830	25.079	nq #	93
69) Atrazine	16.40	200	125695	27.994	nq	98
70) Pentachlorophenol	16.57	266	115387	32.740	nq	97
71) Phenanthrene	16.96	178	583370	26.703	nq	100
72) Anthracene	17.04	178	580611	26.789	nq	98
73) Carbazole	17.33	167	489180	21.990	nq	99
74) Di-n-butylphthalate	17.89	149	696983	25.393	nq	99
75) Fluoranthene	18.99	202	584709	23.020	nq	98
77) Benzidine	19.20	184	425721	46.895	nq	100
78) Pyrene	19.35	202	575586	26.983	nq	99
80) Butylbenzylphthalate	20.27	149	264480	26.063	nq	91
81) Benzo(a)anthracene	21.12	228	564844	27.010	nq	99
82) 3,3'-Dichlorobenzidine	21.06	252	215186	25.756	nq	99
83) Chrysene	21.17	228	528381	26.269	nq	100
84) Bis(2-ethylhexyl)phthalate	21.06	149	388985	25.720	nq	99
85) Di-n-octyl phthalate	21.92	149	679402	27.429	nq	98
86) Indeno(1,2,3-cd)pyrene	25.44	276	712724	35.571	nq	97
88) Benzo(b)fluoranthene	22.66	252	622633	26.650	nq	98
89) Benzo(k)fluoranthene	22.70	252	573130	24.632	nq	97
90) Benzo(a)pyrene	23.20	252	591134	26.603	nq	97
91) Dibenzo(a,h)anthracene	25.44	278	600805	29.188	nq	99
92) Benzo(g,h,i)perylene	26.09	276	591440	30.397	nq	97

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

