

Data Path : Z:\HPCHEM1\BNA F\DATA\BF040216\
 Data File : BF085979.D
 Acq On : 1 Apr 2016 23:41
 Operator : UM/SJ
 Sample : H1824-03
 Misc : LOD 1PPM
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 LOD-SOIL2016-01

Manual Integrations
 APPROVED

Sohil
 4/4/2016 7:38:47 PM

Quant Time: Apr 02 01:26:56 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF040216.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 01 23:17:06 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.77	152	106199	20.00	ng	0.00
21) Naphthalene-d8	8.05	136	416393	20.00	ng	-0.01
38) Acenaphthene-d10	9.80	164	187970	20.00	ng	-0.01
63) Phenanthrene-d10	11.29	188	389206	20.00	ng	-0.01
75) Chrysene-d12	13.93	240	343299	20.00	ng	0.00
86) Perylene-d12	15.33	264	295413	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.37	112	745296	119.96	ng	0.00
7) Phenol-d6	6.42	99	980583	124.26	ng	0.00
23) Nitrobenzene-d5	7.34	82	598002	84.69	ng	0.00
41) 2,4,6-Tribromophenol	10.60	330	279945	158.67	ng	0.00
44) 2-Fluorobiphenyl	9.14	172	1057347	93.53	ng	0.00
78) Terphenyl-d14	12.88	244	1088733	74.55	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.42	88	1318	0.45	ng	# 61
6) Aniline	6.45	93	5438	0.53	ng	# 88
8) 2-Chlorophenol	6.56	128	6080	0.82	ng	88
9) Benzaldehyde	6.37	77	3640	0.86	ng	95
10) Phenol	6.43	94	6730	0.75	ng	# 72
11) bis(2-Chloroethyl)ether	6.52	93	5947	0.83	ng	98
12) 1,3-Dichlorobenzene	6.73	146	6062	0.77	ng	# 92
13) 1,4-Dichlorobenzene	6.80	146	6567	0.81	ng	95
14) 1,2-Dichlorobenzene	6.94	146	5961	0.80	ng	98
15) Benzyl Alcohol	6.92	79	8917	1.75	ng	# 48
16) 2,2'-oxybis(1-Chloropropan	7.06	45	6647	0.76	ng	# 42
17) 2-Methylphenol	7.05	107	5016	0.86	ng	# 82
18) Hexachloroethane	7.28	117	2022	0.68	ng	# 75
19) n-Nitroso-di-n-propylamine	7.18	70	3599	0.74	ng	# 76
20) 3+4-Methylphenols	7.24	107	5173	0.73	ng	# 67
22) Acetophenone	7.21	105	7856	0.80	ng	# 89
24) Nitrobenzene	7.36	77	8015	1.04	ng	98
25) Isophorone	7.61	82	10489m	0.75	ng	
26) 2-Nitrophenol	7.71	139	1682	0.46	ng	# 77
27) 2,4-Dimethylphenol	7.74	122	4216	0.64	ng	98
28) bis(2-Chloroethoxy)methane	7.82	93	6084	0.74	ng	93
29) 2,4-Dichlorophenol	7.96	162	3338	0.59	ng	93
30) 1,2,4-Trichlorobenzene	8.01	180	4815	0.76	ng	98
31) Naphthalene	8.08	128	17499	0.84	ng	99
33) 4-Chloroaniline	8.18	127	4989m	0.59	ng	
34) Hexachlorobutadiene	8.20	225	2267	0.67	ng	96
35) Caprolactam	8.50	113	546	0.31	ng	# 83
36) 4-Chloro-3-methylphenol	8.65	107	3945	0.63	ng	# 89
37) 2-Methylnaphthalene	8.77	142	11221	0.82	ng	98
39) 1,2,4,5-Tetrachlorobenzene	8.94	216	4521	0.80	ng	# 93
42) 2,4,6-Trichlorophenol	9.07	196	1745	0.48	ng	# 82
43) 2,4,5-Trichlorophenol	9.15	196	2041m	0.55	ng	
45) 1,1'-Biphenyl	9.24	154	14032	0.87	ng	94
46) 2-Chloronaphthalene	9.26	162	10286	0.88	ng	96

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47) 2-Nitroaniline	9.39	65	2020	0.59	nq	95
48) Acenaphthylene	9.68	152	15755	0.84	nq	98
49) Dimethylphthalate	9.54	163	10929	0.78	nq	100
50) 2,6-Dinitrotoluene	9.61	165	1418	0.48	nq	89
51) Acenaphthene	9.84	154	10869	0.97	nq	97
52) 3-Nitroaniline	9.80	138	1603	0.45	nq	# 1
54) Dibenzofuran	10.02	168	15132	1.02	nq	96
56) 2,4-Dinitrotoluene	10.03	165	1660	0.45	nq	# 33
57) Fluorene	10.36	166	11333	0.90	nq	98
58) 2,3,4,6-Tetrachlorophenol	10.16	232	1559	0.57	nq	# 90
59) Diethylphthalate	10.24	149	11200	0.80	nq	97
60) 4-Chlorophenyl-phenylether	10.35	204	5605	0.95	nq	98
61) 4-Nitroaniline	10.41	138	1046	0.30	nq	# 68
62) Azobenzene	10.51	77	11394	0.85	nq	97
65) n-Nitrosodiphenylamine	10.48	169	9361	0.77	nq	98
66) 4-Bromophenyl-phenylether	10.84	248	2804	0.71	nq	90
67) Hexachlorobenzene	10.91	284	3091	0.75	nq	94
68) Atrazine	11.00	200	2526	0.64	nq	90
70) Phenanthrene	11.32	178	18324	0.86	nq	100
71) Anthracene	11.38	178	17841	0.80	nq	100
72) Carbazole	11.54	167	14698	0.77	nq	99
73) Di-n-butylphthalate	11.86	149	15669	0.66	nq	96
74) Fluoranthene	12.51	202	16309	0.74	nq	91
77) Pyrene	12.74	202	17557	0.73	nq	99
79) Butylbenzylphthalate	13.35	149	5230	0.48	nq	# 74
80) Benzo(a)anthracene	13.92	228	16934	0.84	nq	99
81) 3,3'-Dichlorobenzidine	13.89	252	2288	0.15	nq	# 90
82) Chrysene	13.95	228	14680	0.75	nq	98
83) Bis(2-ethylhexyl)phthalate	13.91	149	7295	0.50	nq	# 99
84) Di-n-octyl phthalate	14.51	149	10272	0.44	nq	# 90
85) Indeno(1,2,3-cd)pyrene	16.71	276	10324	0.65	nq	96
87) Benzo(b)fluoranthene	14.95	252	13142	0.71	nq	# 95
88) Benzo(k)fluoranthene	14.97	252	12776m	0.78	nq	
89) Benzo(a)pyrene	15.28	252	11239	0.68	nq	# 97
90) Dibenzo(a,h)anthracene	16.71	278	8473	0.64	nq	97
91) Benzo(a,h,i)perylene	17.12	276	8597	0.67	nq	# 92

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

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