

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF070219\
 Data File : BF115376.D
 Acq On : 2 Jul 2019 22:40
 Operator : HP/JU
 Sample : K3629-01MS
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 NB-SOIL-PILEMS

Manual Integrations
 APPROVED

mohammad
 7/3/2019 10:51:56 PM

Quant Time: Jul 03 11:42:44 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF062119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 28 05:26:43 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.61	152	228434	20.00	ng	0.00
21) Naphthalene-d8	7.89	136	890192	20.00	ng	0.00
39) Acenaphthene-d10	9.64	164	432697	20.00	ng	0.01
64) Phenanthrene-d10	11.11	188	825743	20.00	ng	0.02
76) Chrysene-d12	13.74	240	461053	20.00	ng	0.02
87) Perylene-d12	15.11	264	505672	20.00	ng	0.03

System Monitoring Compounds

5) 2-Fluorophenol	5.21	112	1090350	73.66	ng	0.02
7) Phenol-d6	6.25	99	1390080	77.37	ng	0.01
23) Nitrobenzene-d5	7.17	82	842686	58.23	ng	0.00
42) 2,4,6-Tribromophenol	10.42	330	360162	78.93	ng	0.02
45) 2-Fluorobiphenyl	8.97	172	1618119	63.52	ng	0.01
79) Terphenyl-d14	12.70	244	1589005	73.28	ng	0.02

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.24	88	135640	19.415	ng	99
3) Pyridine	2.90	79	184602	9.375	ng	95
4) n-Nitrosodimethylamine	2.82	42	150602	19.510	ng	99
6) Aniline	6.26	93	296687	12.015	ng	# 1
8) 2-Chlorophenol	6.39	128	397490	23.880	ng	99
9) Benzaldehyde	6.14	77	239354	20.419	ng	98
10) Phenol	6.26	94	474862	22.563	ng	92
11) bis(2-Chloroethyl)ether	6.35	93	361166	23.280	ng	97
12) 1,3-Dichlorobenzene	6.54	146	433787	23.482	ng	98
13) 1,4-Dichlorobenzene	6.62	146	445993	24.076	ng	99
14) 1,2-Dichlorobenzene	6.78	146	416148	24.259	ng	98
15) Benzyl Alcohol	6.75	79	287352	21.964	ng	97
16) 2,2'-oxybis(1-Chloropropan	6.89	45	498064	22.135	ng	94
17) 2-Methylphenol	6.87	107	338596	24.644	ng	99
18) Hexachloroethane	7.12	117	151092	22.624	ng	99
19) n-Nitroso-di-n-propylamine	7.02	70	256029	22.258	ng	97
20) 3+4-Methylphenols	7.02	107	404609	25.504	ng	# 81
22) Acetophenone	7.02	105	546824	26.832	ng	# 95
24) Nitrobenzene	7.19	77	402103	25.222	ng	98
25) Isophorone	7.43	82	713318	24.062	ng	99
26) 2-Nitrophenol	7.51	139	210761	26.770	ng	98
27) 2,4-Dimethylphenol	7.55	122	334882	25.979	ng	98
28) bis(2-Chloroethoxy)methane	7.65	93	459512	24.525	ng	100
29) 2,4-Dichlorophenol	7.75	162	346021	26.011	ng	97
30) 1,2,4-Trichlorobenzene	7.84	180	377871	25.716	ng	98
31) Naphthalene	7.91	128	1162953	26.614	ng	98
32) Benzoic acid	7.67	122	8899m	0.967	ng	
33) 4-Chloroaniline	7.96	127	297684	14.906	ng	100
34) Hexachlorobutadiene	8.04	225	205111	25.472	ng	98
35) Caprolactam	8.31	113	84455m	18.883	ng	
36) 4-Chloro-3-methylphenol	8.45	107	345052	25.612	ng	99
37) 2-Methylnaphthalene	8.60	142	782285	26.551	ng	99
38) 1-Methylnaphthalene	8.70	142	731486	25.995	ng	99
40) 1,2,4,5-Tetrachlorobenzene	8.77	216	350322	28.651	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	8.76	237	210747	29.067	ng	98
43) 2,4,6-Trichlorophenol	8.88	196	232304	24.609	ng	99
44) 2,4,5-Trichlorophenol	8.92	196	258094	26.549	ng	96
46) 1,1'-Biphenyl	9.07	154	933992	26.977	ng	99
47) 2-Chloronaphthalene	9.09	162	731143	26.034	ng	100
48) 2-Nitroaniline	9.18	65	210418	25.700	ng	89
49) Acenaphthylene	9.50	152	1159239	26.494	ng	98
50) Dimethylphthalate	9.37	163	1071423	33.656	ng	100
51) 2,6-Dinitrotoluene	9.42	165	197111	26.819	ng	98
52) Acenaphthene	9.67	154	669160	24.440	ng	99
53) 3-Nitroaniline	9.59	138	198012	22.078	ng	98
54) 2,4-Dinitrophenol	9.70	184	13968	10.698	ng	# 86
55) Dibenzofuran	9.84	168	1013319	25.786	ng	98
56) 4-Nitrophenol	9.76	139	282918	39.347	ng	91
57) 2,4-Dinitrotoluene	9.82	165	253638	26.727	ng	92
58) Fluorene	10.18	166	734000	25.148	ng	98
59) 2,3,4,6-Tetrachlorophenol	9.97	232	187054	22.947	ng	96
60) Diethylphthalate	10.07	149	818578	25.580	ng	99
61) 4-Chlorophenyl-phenylether	10.18	204	357437	25.853	ng	98
62) 4-Nitroaniline	10.20	138	218458	24.280	ng	95
63) Azobenzene	10.34	77	737813	24.685	ng	99
65) 4,6-Dinitro-2-methylphenol	10.24	198	35700	12.246	ng	98
66) n-Nitrosodiphenylamine	10.30	169	680397	26.448	ng	99
67) 4-Bromophenyl-phenylether	10.67	248	238190	26.605	ng	98
68) Hexachlorobenzene	10.73	284	260966	26.855	ng	94
69) Atrazine	10.83	200	225004	27.189	ng	99
70) Pentachlorophenol	10.92	266	176801	28.558	ng	97
71) Phenanthrene	11.14	178	1130395	25.425	ng	98
72) Anthracene	11.19	178	1178732	26.265	ng	99
73) Carbazole	11.34	167	1033297	25.784	ng	99
74) Di-n-butylphthalate	11.70	149	1253371	27.066	ng	99
75) Fluoranthene	12.32	202	1075385	24.243	ng	98
77) Benzidine	12.45	184	90845	4.437	ng	99
78) Pyrene	12.55	202	1056734	29.824	ng	99
80) Butylbenzylphthalate	13.18	149	451461	28.872	ng	90
81) Benzo(a)anthracene	13.73	228	813452	24.589	ng	98
82) 3,3'-Dichlorobenzidine	13.70	252	264577	22.147	ng	99
83) Chrysene	13.77	228	786210	25.944	ng	98
84) Bis(2-ethylhexyl)phthalate	13.75	149	628434	30.672	ng	99
85) Di-n-octyl phthalate	14.36	149	928014	26.958	ng	99
86) Indeno(1,2,3-cd)pyrene	16.39	276	664278	18.525	ng	98
88) Benzo(b)fluoranthene	14.74	252	840649	26.643	ng	98
89) Benzo(k)fluoranthene	14.76	252	749988	26.505	ng	97
90) Benzo(a)pyrene	15.05	252	763108	26.833	ng	98
91) Dibenzo(a,h)anthracene	16.40	278	555053	20.906	ng	99
92) Benzo(g,h,i)perylene	16.77	276	517233	19.249	ng	100

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

