

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF070920\
 Data File : BF120811.D
 Acq On : 9 Jul 2020 11:32
 Operator : JU/CG
 Sample : L3216-01
 Misc : LOD-MDL-SOIL-8PPM
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 LOD-MDL-SOIL-01-QT3-2020

Manual Integrations
 APPROVED

mohammad
 7/10/2020 11:49:31 AM

Quant Time: Jul 09 12:17:40 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF070220.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 09 11:36:34 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	136434	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	507072	20.00	ng	0.00
39) Acenaphthene-d10	9.88	164	247149	20.00	ng	0.00
64) Phenanthrene-d10	11.37	188	435077	20.00	ng	0.00
76) Chrysene-d12	14.00	240	312942	20.00	ng	0.00
86) Perylene-d12	15.46	264	259786	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.46	112	820735	92.86	ng	0.00
7) Phenol-d6	6.47	99	1046128	92.68	ng	0.00
23) Nitrobenzene-d5	7.41	82	697105	79.65	ng	0.00
42) 2,4,6-Tribromophenol	10.67	330	245573	98.90	ng	0.00
45) 2-Fluorobiphenyl	9.21	172	1273040	77.93	ng	0.00
79) Terphenyl-d14	12.96	244	1007034	63.00	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.58	88	32317	8.005	ng	98
3) Pyridine	3.36	79	57713	5.219	ng	98
4) n-Nitrosodimethylamine	3.26	42	35462	6.410	ng	95
6) Aniline	6.50	93	101872	7.028	ng	99
8) 2-Chlorophenol	6.63	128	65941	7.008	ng	92
9) Benzaldehyde	6.39	77	57571	8.576	ng	95
10) Phenol	6.48	94	82284	7.087	ng	95
11) bis(2-Chloroethyl)ether	6.58	93	64920	6.871	ng	99
12) 1,3-Dichlorobenzene	6.79	146	71910	6.943	ng	96
13) 1,4-Dichlorobenzene	6.86	146	73070	7.044	ng	98
14) 1,2-Dichlorobenzene	7.02	146	69841	7.076	ng	99
15) Benzyl Alcohol	6.98	79	64687	7.628	ng	98
16) 2,2'-oxybis(1-Chloropropan	7.12	45	97296	6.602	ng	99
17) 2-Methylphenol	7.09	107	50377	5.984	ng	95
18) Hexachloroethane	7.36	117	26537	6.806	ng	94
19) n-Nitroso-di-n-propylamine	7.25	70	45396	6.680	ng	99
20) 3+4-Methylphenols	7.24	107	66495	6.413	ng	93
22) Acetophenone	7.25	105	97049	7.700	ng	95
24) Nitrobenzene	7.42	77	74370	7.568	ng	94
25) Isophorone	7.66	82	122902	6.512	ng	98
26) 2-Nitrophenol	7.74	139	21830	6.077	ng	93
27) 2,4-Dimethylphenol	7.77	122	53830	7.144	ng	98
28) bis(2-Chloroethoxy)methane	7.87	93	78415	7.020	ng	# 97
29) 2,4-Dichlorophenol	7.98	162	48932	6.505	ng	98
30) 1,2,4-Trichlorobenzene	8.07	180	56407	6.677	ng	96
31) Naphthalene	8.15	128	196196	7.240	ng	98
32) Benzoic acid	7.83	122	17881	4.385	ng	93
33) 4-Chloroaniline	8.19	127	75157	6.536	ng	95
34) Hexachlorobutadiene	8.27	225	33011	6.497	ng	97
35) Caprolactam	8.52	113	12661	5.777	ng	96
36) 4-Chloro-3-methylphenol	8.67	107	49640	6.344	ng	95
37) 2-Methylnaphthalene	8.84	142	123011	6.983	ng	98
38) 1-Methylnaphthalene	8.94	142	116866	7.082	ng	100
40) 1,2,4,5-Tetrachlorobenzene	9.00	216	54382	7.281	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	9.00	237	26175	5.931	nq	97
43) 2,4,6-Trichlorophenol	9.12	196	30275	5.948	nq	98
44) 2,4,5-Trichlorophenol	9.15	196	32953	5.834	nq	97
46) 1,1'-Biphenyl	9.30	154	155066	7.778	nq	99
47) 2-Chloronaphthalene	9.33	162	109796	6.875	nq	99
48) 2-Nitroaniline	9.42	65	28311	6.625	nq	86
49) Acenaphthylene	9.75	152	174607	6.973	nq	98
50) Dimethylphthalate	9.60	163	119820	6.646	nq	99
51) 2,6-Dinitrotoluene	9.66	165	21292	6.537	nq	89
52) Acenaphthene	9.92	154	105738	6.807	nq	99
53) 3-Nitroaniline	9.83	138	25115	6.241	nq	97
54) 2,4-Dinitrophenol	9.93	184	3076	2.860	nq	# 1
55) Dibenzofuran	10.09	168	155687	6.953	nq	99
56) 4-Nitrophenol	9.97	139	15827	5.355	nq	93
57) 2,4-Dinitrotoluene	10.06	165	25319	6.475	nq	98
58) Fluorene	10.43	166	117158	7.030	nq	98
59) 2,3,4,6-Tetrachlorophenol	10.20	232	25095m	5.970	nq	
60) Diethylphthalate	10.30	149	119208	6.347	nq	98
61) 4-Chlorophenyl-phenylether	10.42	204	57125	6.818	nq	95
62) 4-Nitroaniline	10.43	138	23271	6.075	nq	90
63) Azobenzene	10.58	77	130430	6.789	nq	92
65) 4,6-Dinitro-2-methylphenol	10.46	198	5757	4.118	nq	95
66) n-Nitrosodiphenylamine	10.54	169	99377	6.815	nq	99
67) 4-Bromophenyl-phenylether	10.92	248	30105	5.965	nq	95
68) Hexachlorobenzene	10.97	284	33935	6.408	nq	97
69) Atrazine	11.06	200	27010	7.190	nq	98
70) Pentachlorophenol	11.17	266	13487	4.501	nq	97
71) Phenanthrene	11.39	178	163685	6.934	nq	98
72) Anthracene	11.44	178	162568	6.815	nq	98
73) Carbazole	11.59	167	145824	6.361	nq	98
74) Di-n-butylphthalate	11.93	149	161691	5.697	nq	99
75) Fluoranthene	12.57	202	159238	6.332	nq	98
77) Benzidine	12.70	184	56305	6.374	nq	96
78) Pyrene	12.80	202	166179	6.649	nq	98
80) Butylbenzylphthalate	13.43	149	51470	4.770	nq	97
81) Benzo(a)anthracene	13.99	228	134635	6.712	nq	99
82) 3,3'-Dichlorobenzidine	13.96	252	37984	5.611	nq	# 99
83) Chrysene	14.03	228	129874	6.347	nq	97
84) Bis(2-ethylhexyl)phthalate	13.99	149	82552	5.834	nq	98
85) Di-n-octyl phthalate	14.60	149	100274	4.006	nq	97
87) Indeno(1,2,3-cd)pyrene	16.90	276	97341	5.842	nq	97
88) Benzo(b)fluoranthene	15.04	252	101477	5.942	nq	97
89) Benzo(k)fluoranthene	15.06	252	108672	6.539	nq	98
90) Benzo(a)pyrene	15.40	252	88742	5.825	nq	98
91) Dibenzo(a,h)anthracene	16.93	278	83567	6.017	nq	95
92) Benzo(g,h,i)perylene	17.34	276	81349	6.409	nq	98

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

