

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF053019\
 Data File : BF114656.D
 Acq On : 30 May 2019 13:08
 Operator : HP/JU
 Sample : K2241-01
 Misc : LOD-SOIL-5PPM
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 LOD-MDL-SOIL-01-QT2-2019

Manual Integrations
 APPROVED

mohammad
 6/3/2019 8:47:03 AM

Quant Time: May 30 14:46:06 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF052919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu May 30 14:17:39 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.69	152	412684	20.00	ng	0.00
21) Naphthalene-d8	7.97	136	1555639	20.00	ng	0.00
39) Acenaphthene-d10	9.72	164	786814	20.00	ng	0.00
64) Phenanthrene-d10	11.19	188	1488563	20.00	ng	0.00
76) Chrysene-d12	13.81	240	1126995	20.00	ng	0.00
87) Perylene-d12	15.20	264	1192429	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.30	112	3009285	128.03	ng	0.00
7) Phenol-d6	6.33	99	3575339	126.23	ng	0.00
23) Nitrobenzene-d5	7.26	82	2189545	87.81	ng	0.00
42) 2,4,6-Tribromophenol	10.50	330	1064043	141.93	ng	0.00
45) 2-Fluorobiphenyl	9.05	172	3751008	79.20	ng	0.00
79) Terphenyl-d14	12.78	244	3993089	66.62	ng	0.00

Target Compounds				Qvalue		
2) 1,4-Dioxane	2.37	88	58981	5.229	ng	98
3) Pyridine	3.12	79	145928	4.445	ng	97
4) n-Nitrosodimethylamine	2.98	42	56345	4.241	ng	# 97
6) Aniline	6.36	93	214926	5.023	ng	99
8) 2-Chlorophenol	6.47	128	142828	5.262	ng	91
9) Benzaldehyde	6.23	77	117912	7.296	ng	99
10) Phenol	6.34	94	180252	5.268	ng	92
11) bis(2-Chloroethyl)ether	6.43	93	136599	5.219	ng	95
12) 1,3-Dichlorobenzene	6.63	146	164938	5.333	ng	97
13) 1,4-Dichlorobenzene	6.71	146	167729	5.406	ng	99
14) 1,2-Dichlorobenzene	6.86	146	156248	5.532	ng	97
15) Benzyl Alcohol	6.83	79	114281m	5.129	ng	
16) 2,2'-oxybis(1-Chloropropan	6.97	45	196813	4.903	ng	98
17) 2-Methylphenol	6.94	107	115582	5.034	ng	99
18) Hexachloroethane	7.21	117	57076	4.859	ng	96
19) n-Nitroso-di-n-propylamine	7.10	70	99886	5.109	ng	95
20) 3+4-Methylphenols	7.10	107	150230	5.563	ng	92
22) Acetophenone	7.10	105	204484	6.169	ng	99
24) Nitrobenzene	7.27	77	151734	5.729	ng	98
25) Isophorone	7.51	82	270441	5.404	ng	98
26) 2-Nitrophenol	7.59	139	61751	4.790	ng	99
27) 2,4-Dimethylphenol	7.63	122	128854	6.009	ng	99
28) bis(2-Chloroethoxy)methane	7.73	93	170353	5.348	ng	98
29) 2,4-Dichlorophenol	7.82	162	119094	5.457	ng	98
30) 1,2,4-Trichlorobenzene	7.92	180	140078	5.633	ng	99
31) Naphthalene	7.99	128	444260	6.200	ng	96
32) Benzoic acid	7.68	122	48707	3.497	ng	95
33) 4-Chloroaniline	8.04	127	181535	5.318	ng	97
34) Hexachlorobutadiene	8.12	225	74749	5.411	ng	98
35) Caprolactam	8.36	113	35684	4.845	ng	96
36) 4-Chloro-3-methylphenol	8.52	107	119697	5.469	ng	99
37) 2-Methylnaphthalene	8.68	142	304622	6.398	ng	98
38) 1-Methylnaphthalene	8.78	142	287167	6.366	ng	99
40) 1,2,4,5-Tetrachlorobenzene	8.85	216	131779	6.289	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	8.85	237	79235	5.722	nq	98
43) 2,4,6-Trichlorophenol	8.96	196	84960	5.153	nq	99
44) 2,4,5-Trichlorophenol	8.99	196	89637	5.448	nq	99
46) 1,1'-Biphenyl	9.15	154	372588	6.379	nq	96
47) 2-Chloronaphthalene	9.17	162	282259	5.891	nq	94
48) 2-Nitroaniline	9.25	65	67631	4.574	nq	91
49) Acenaphthylene	9.57	152	451089	6.095	nq	97
50) Dimethylphthalate	9.44	163	318399	5.807	nq	98
51) 2,6-Dinitrotoluene	9.50	165	64025	5.026	nq	99
52) Acenaphthene	9.75	154	268684	5.919	nq	99
53) 3-Nitroaniline	9.66	138	74128	4.754	nq	93
54) 2,4-Dinitrophenol	9.76	184	13849	2.756	nq	# 42
55) Dibenzofuran	9.92	168	399945	6.190	nq	95
56) 4-Nitrophenol	9.81	139	51851	4.487	nq	92
57) 2,4-Dinitrotoluene	9.89	165	79460	4.860	nq	97
58) Fluorene	10.26	166	306417	7.023	nq	100
59) 2,3,4,6-Tetrachlorophenol	10.03	232	72926	5.415	nq	99
60) Diethylphthalate	10.15	149	316500	5.652	nq	98
61) 4-Chlorophenyl-phenylether	10.26	204	147113	6.740	nq	99
62) 4-Nitroaniline	10.26	138	71893	4.732	nq	98
63) Azobenzene	10.42	77	296363	5.765	nq	96
65) 4,6-Dinitro-2-methylphenol	10.30	198	23949	3.628	nq	95
66) n-Nitrosodiphenylamine	10.37	169	265562	6.038	nq	99
67) 4-Bromophenyl-phenylether	10.74	248	88277	5.530	nq	98
68) Hexachlorobenzene	10.81	284	97396	5.633	nq	95
69) Atrazine	10.90	200	76406	5.514	nq	98
70) Pentachlorophenol	11.00	266	45402	4.555	nq	97
71) Phenanthrene	11.22	178	461369	6.076	nq	96
72) Anthracene	11.26	178	466423	6.069	nq	97
73) Carbazole	11.42	167	405267	6.000	nq	97
74) Di-n-butylphthalate	11.76	149	483766	5.895	nq	97
75) Fluoranthene	12.39	202	462194	5.878	nq	95
77) Benzidine	12.52	184	237783	5.124	nq	97
78) Pyrene	12.62	202	468548	4.915	nq	97
80) Butylbenzylphthalate	13.25	149	187848	3.921	nq	98
81) Benzo(a)anthracene	13.80	228	426114	4.920	nq	97
82) 3,3'-Dichlorobenzidine	13.77	252	151398	4.527	nq	98
83) Chrysene	13.83	228	405268	4.898	nq	96
84) Bis(2-ethylhexyl)phthalate	13.82	149	268164	4.527	nq	98
85) Di-n-octyl phthalate	14.43	149	446956	4.343	nq	98
86) Indeno(1,2,3-cd)pyrene	16.51	276	374317	4.517	nq	99
88) Benzo(b)fluoranthene	14.81	252	396279	5.276	nq	97
89) Benzo(k)fluoranthene	14.83	252	390905	5.988	nq	97
90) Benzo(a)pyrene	15.14	252	367593	5.571	nq	99
91) Dibenzo(a,h)anthracene	16.53	278	317594	5.447	nq	99
92) Benzo(g,h,i)perylene	16.90	276	299748	5.297	nq	98

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

