

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071719\
 Data File : BF115625.D
 Acq On : 17 Jul 2019 16:39
 Operator : HP/JU
 Sample : K3838-01MS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WARINANCO-CITY-SEWER-INSTAL

Manual Integrations
 APPROVED

mohammad
 7/19/2019 2:25:52 PM

Quant Time: Jul 18 05:12:08 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071219.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 12 14:03:52 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.88	152	190811	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	659260	20.00	ng	0.00
39) Acenaphthene-d10	9.92	164	321655	20.00	ng	-0.01
64) Phenanthrene-d10	11.40	188	589759	20.00	ng	-0.01
76) Chrysene-d12	14.03	240	307427	20.00	ng	-0.02
87) Perylene-d12	15.50	264	323097	20.00	ng	-0.02

System Monitoring Compounds

5) 2-Fluorophenol	5.52	112	1133356	97.16	ng	0.01
7) Phenol-d6	6.52	99	1540039	104.04	ng	0.00
23) Nitrobenzene-d5	7.45	82	887985	78.32	ng	0.00
42) 2,4,6-Tribromophenol	10.70	330	363550	96.83	ng	-0.01
45) 2-Fluorobiphenyl	9.24	172	1600725	87.10	ng	-0.01
79) Terphenyl-d14	12.98	244	1455446	87.36	ng	-0.02

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.76	88	193819	33.309	ng	98
3) Pyridine	3.52	79	213663	13.534	ng	99
4) n-Nitrosodimethylamine	3.45	42	220414	38.485	ng	98
6) Aniline	6.55	93	516345	25.066	ng	97
8) 2-Chlorophenol	6.67	128	507812	37.863	ng	96
9) Benzaldehyde	6.43	77	266945	28.859	ng	99
10) Phenol	6.53	94	619235	36.038	ng	86
11) bis(2-Chloroethyl)ether	6.62	93	474789	36.292	ng	100
12) 1,3-Dichlorobenzene	6.83	146	526536	34.918	ng	98
13) 1,4-Dichlorobenzene	6.90	146	551979	36.688	ng	97
14) 1,2-Dichlorobenzene	7.06	146	512952	37.296	ng	97
15) Benzyl Alcohol	7.02	79	446948	38.064	ng	99
16) 2,2'-oxybis(1-Chloropropan	7.16	45	660050	38.323	ng	97
17) 2-Methylphenol	7.13	107	441744	38.219	ng	99
18) Hexachloroethane	7.40	117	160680	28.773	ng	98
19) n-Nitroso-di-n-propylamine	7.30	70	308563	31.963	ng	99
20) 3+4-Methylphenols	7.29	107	441202	32.137	ng	94
22) Acetophenone	7.29	105	610704	41.004	ng	# 93
24) Nitrobenzene	7.46	77	473043	38.682	ng	97
25) Isophorone	7.70	82	865061	38.130	ng	100
26) 2-Nitrophenol	7.78	139	212730	33.797	ng	94
27) 2,4-Dimethylphenol	7.82	122	366996	38.968	ng	99
28) bis(2-Chloroethoxy)methane	7.91	93	535671	38.487	ng	98
29) 2,4-Dichlorophenol	8.02	162	368548	37.627	ng	98
30) 1,2,4-Trichlorobenzene	8.10	180	424293	38.323	ng	99
31) Naphthalene	8.19	128	1270408	39.790	ng	97
33) 4-Chloroaniline	8.23	127	392721	27.769	ng	99
34) Hexachlorobutadiene	8.31	225	240525	37.883	ng	98
35) Caprolactam	8.59	113	65351m	21.592	ng	
36) 4-Chloro-3-methylphenol	8.71	107	393272	38.708	ng	99
37) 2-Methylnaphthalene	8.87	142	848491	40.091	ng	99
38) 1-Methylnaphthalene	8.97	142	795454	39.570	ng	100
40) 1,2,4,5-Tetrachlorobenzene	9.05	216	387845	40.041	ng	98
41) Hexachlorocyclopentadiene	9.03	237	328801	59.508	ng	99

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071719\
 Data File : BF115625.D
 Acq On : 17 Jul 2019 16:39
 Operator : HP/JU
 Sample : K3838-01MS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WARINANCO-CITY-SEWER-INSTAL

Manual Integrations
 APPROVED

mohammad
 7/19/2019 2:25:52 PM

Quant Time: Jul 18 05:12:08 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071219.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 12 14:03:52 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
43) 2,4,6-Trichlorophenol	9.15	196	254265	35.896	nq	98
44) 2,4,5-Trichlorophenol	9.19	196	260531	35.915	nq	99
46) 1,1'-Biphenyl	9.34	154	1036786	41.230	nq	99
47) 2-Chloronaphthalene	9.36	162	814936	38.921	nq	99
48) 2-Nitroaniline	9.46	65	247694	38.220	nq	93
49) Acenaphthylene	9.78	152	1230331	39.655	nq	98
50) Dimethylphthalate	9.64	163	1202457	49.688	nq	99
51) 2,6-Dinitrotoluene	9.70	165	216172	39.306	nq	93
52) Acenaphthene	9.95	154	734985	36.693	nq	98
53) 3-Nitroaniline	9.86	138	203114	32.318	nq	99
54) 2,4-Dinitrophenol	9.97	184	14935	5.289	nq	93
55) Dibenzofuran	10.12	168	1092657	38.412	nq	99
56) 4-Nitrophenol	10.02	139	301448	62.900	nq	98
57) 2,4-Dinitrotoluene	10.10	165	276607	38.821	nq	97
58) Fluorene	10.46	166	786256	38.664	nq	99
59) 2,3,4,6-Tetrachlorophenol	10.24	232	190346	31.389	nq	99
60) Diethylphthalate	10.33	149	893566	39.408	nq	98
61) 4-Chlorophenyl-phenylether	10.46	204	415439	36.839	nq	99
62) 4-Nitroaniline	10.47	138	207179	34.367	nq	97
63) Azobenzene	10.62	77	861450	39.168	nq	98
65) 4,6-Dinitro-2-methylphenol	10.51	198	57817	16.046	nq	89
66) n-Nitrosodiphenylamine	10.57	169	734003	40.010	nq	98
67) 4-Bromophenyl-phenylether	10.95	248	263962	39.166	nq	96
68) Hexachlorobenzene	11.02	284	286044	37.165	nq	# 94
69) Atrazine	11.10	200	240572	39.974	nq	99
70) Pentachlorophenol	11.20	266	180563	38.357	nq	100
71) Phenanthrene	11.43	178	1153106	38.144	nq	97
72) Anthracene	11.47	178	1193240	38.193	nq	100
73) Carbazole	11.63	167	1023490	37.358	nq	98
74) Di-n-butylphthalate	11.96	149	1298094	38.466	nq	99
75) Fluoranthene	12.61	202	1107398	34.665	nq	99
77) Benzidine	12.72	184	103120	9.469	nq	100
78) Pyrene	12.84	202	1071963	41.156	nq	98
80) Butylbenzylphthalate	13.45	149	464869	41.867	nq	98
81) Benzo(a)anthracene	14.03	228	796394	39.322	nq	96
82) 3,3'-Dichlorobenzidine	13.98	252	268381	37.275	nq	97
83) Chrysene	14.06	228	783064	38.515	nq	97
84) Bis(2-ethylhexyl)phthalate	14.02	149	605227	44.005	nq	99
85) Di-n-octyl phthalate	14.62	149	935676	42.758	nq	98
86) Indeno(1,2,3-cd)pyrene	16.97	276	836029	48.400	nq	100
88) Benzo(b)fluoranthene	15.07	252	756340	36.354	nq	99
89) Benzo(k)fluoranthene	15.10	252	717256	38.669	nq	98
90) Benzo(a)pyrene	15.44	252	714331	39.948	nq	99
91) Dibenzo(a,h)anthracene	16.99	278	702653	44.760	nq	99
92) Benzo(g,h,i)perylene	17.42	276	687153	43.890	nq	99

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

