

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF073119\
 Data File : BF115867.D
 Acq On : 31 Jul 2019 15:49
 Operator : HP/JU
 Sample : K3591-08
 Misc : LOO-WATER-10PPM
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 LOQ-WATER-02-QT3-2019

Manual Integrations
 APPROVED

mohammad
 8/1/2019 5:06:37 PM

Quant Time: Aug 01 11:50:32 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 01 09:04:21 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	133931	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	529340	20.00	ng	0.00
39) Acenaphthene-d10	9.89	164	275476	20.00	ng	0.00
64) Phenanthrene-d10	11.38	188	480570	20.00	ng	0.00
76) Chrysene-d12	14.02	240	369135	20.00	ng	0.00
87) Perylene-d12	15.49	264	341556	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.47	112	954257	119.28	ng	0.00
7) Phenol-d6	6.49	99	1189596	117.85	ng	0.00
23) Nitrobenzene-d5	7.42	82	748920	81.67	ng	0.00
42) 2,4,6-Tribromophenol	10.68	330	312089	116.85	ng	0.00
45) 2-Fluorobiphenyl	9.22	172	1259036	85.61	ng	0.00
79) Terphenyl-d14	12.96	244	1345757	76.82	ng	0.00

Target Compounds					Qvalue
2) 1,4-Dioxane	2.61	88	42121	10.422 ng	99
3) Pyridine	3.42	79	89863	7.959 ng	99
4) n-Nitrosodimethylamine	3.30	42	41191	9.245 ng	97
6) Aniline	6.52	93	145308	10.052 ng	99
8) 2-Chlorophenol	6.64	128	92888	10.500 ng	97
9) Benzaldehyde	6.40	77	79486	11.413 ng	99
10) Phenol	6.49	94	125180	10.531 ng	92
11) bis(2-Chloroethyl)ether	6.59	93	87730	10.039 ng	96
12) 1,3-Dichlorobenzene	6.79	146	104606	10.231 ng	95
13) 1,4-Dichlorobenzene	6.87	146	104726	10.230 ng	98
14) 1,2-Dichlorobenzene	7.02	146	97976	10.394 ng	98
15) Benzyl Alcohol	6.99	79	81790	10.677 ng	99
16) 2,2'-oxybis(1-Chloropropan	7.13	45	116302	9.757 ng	99
17) 2-Methylphenol	7.10	107	75053	10.019 ng	98
18) Hexachloroethane	7.36	117	35856	9.513 ng	93
19) n-Nitroso-di-n-propylamine	7.26	70	65207	10.425 ng	98
20) 3+4-Methylphenols	7.26	107	95706	10.796 ng	# 56
22) Acetophenone	7.26	105	132324	11.398 ng	98
24) Nitrobenzene	7.43	77	107664	10.873 ng	99
25) Isophorone	7.67	82	176678	10.076 ng	99
26) 2-Nitrophenol	7.75	139	43576	9.336 ng	98
27) 2,4-Dimethylphenol	7.78	122	77602	11.051 ng	98
28) bis(2-Chloroethoxy)methane	7.87	93	107281	9.871 ng	98
29) 2,4-Dichlorophenol	7.99	162	75110	10.130 ng	97
30) 1,2,4-Trichlorobenzene	8.07	180	84092	9.950 ng	98
31) Naphthalene	8.15	128	275117	11.045 ng	98
32) Benzoic acid	7.87	122	26003m	5.252 ng	
33) 4-Chloroaniline	8.20	127	112239	10.085 ng	98
34) Hexachlorobutadiene	8.27	225	47117	9.678 ng	98
35) Caprolactam	8.55	113	22621	9.433 ng	97
36) 4-Chloro-3-methylphenol	8.69	107	78020	10.115 ng	98
37) 2-Methylnaphthalene	8.85	142	181003	11.033 ng	97
38) 1-Methylnaphthalene	8.95	142	169033	10.891 ng	100
40) 1,2,4,5-Tetrachlorobenzene	9.02	216	78656	10.680 ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	9.00	237	15843	5.504	ng	97
43) 2,4,6-Trichlorophenol	9.13	196	46482	9.170	ng	98
44) 2,4,5-Trichlorophenol	9.17	196	51252	9.781	ng	96
46) 1,1'-Biphenyl	9.31	154	221788	11.404	ng	98
47) 2-Chloronaphthalene	9.34	162	168526	10.411	ng	100
48) 2-Nitroaniline	9.43	65	49222	9.200	ng	# 81
49) Acenaphthylene	9.75	152	266306	11.277	ng	98
50) Dimethylphthalate	9.60	163	183935	9.806	ng	99
51) 2,6-Dinitrotoluene	9.67	165	39828	9.771	ng	88
52) Acenaphthene	9.92	154	156678	10.815	ng	98
53) 3-Nitroaniline	9.84	138	45010	8.937	ng	93
54) 2,4-Dinitrophenol	9.96	184	7654	4.422	ng	90
55) Dibenzofuran	10.10	168	236169	11.209	ng	100
56) 4-Nitrophenol	10.01	139	22216	6.886	ng	91
57) 2,4-Dinitrotoluene	10.07	165	50593	9.322	ng	# 86
58) Fluorene	10.44	166	183017	11.291	ng	100
59) 2,3,4,6-Tetrachlorophenol	10.22	232	40582	9.205	ng	97
60) Diethylphthalate	10.31	149	174307	9.954	ng	99
61) 4-Chlorophenyl-phenylether	10.43	204	86085	10.486	ng	95
62) 4-Nitroaniline	10.45	138	47129	9.054	ng	97
63) Azobenzene	10.59	77	191155	10.613	ng	97
65) 4,6-Dinitro-2-methylphenol	10.49	198	16899	6.636	ng	96
66) n-Nitrosodiphenylamine	10.55	169	153861	10.637	ng	99
67) 4-Bromophenyl-phenylether	10.92	248	51518	10.014	ng	93
68) Hexachlorobenzene	10.99	284	56704	9.532	ng	98
69) Atrazine	11.07	200	44928	9.441	ng	98
70) Pentachlorophenol	11.19	266	17938	6.211	ng	# 87
71) Phenanthrene	11.40	178	269201	10.957	ng	98
72) Anthracene	11.45	178	274859	11.055	ng	99
73) Carbazole	11.61	167	243743	10.805	ng	96
74) Di-n-butylphthalate	11.94	149	247670	9.412	ng	99
75) Fluoranthene	12.59	202	272693	10.602	ng	98
77) Benzidine	12.71	184	125874	10.093	ng	100
78) Pyrene	12.82	202	283459	10.187	ng	98
80) Butylbenzylphthalate	13.43	149	94789	8.053	ng	96
81) Benzo(a)anthracene	14.01	228	232146	9.839	ng	99
82) 3,3'-Dichlorobenzidine	13.97	252	79966	9.756	ng	98
83) Chrysene	14.04	228	233961	10.214	ng	98
84) Bis(2-ethylhexyl)phthalate	13.99	149	126406	8.264	ng	98
85) Di-n-octyl phthalate	14.60	149	188907	7.776	ng	99
86) Indeno(1,2,3-cd)pyrene	16.96	276	167484	8.706	ng	99
88) Benzo(b)fluoranthene	15.06	252	185467	9.391	ng	98
89) Benzo(k)fluoranthene	15.09	252	205976	10.708	ng	97
90) Benzo(a)pyrene	15.42	252	179799	10.184	ng	98
91) Dibenzo(a,h)anthracene	16.97	278	138741	9.079	ng	98
92) Benzo(g,h,i)perylene	17.40	276	134628	8.970	ng	97

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

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