

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF053019\
 Data File : BF114660.D
 Acq On : 30 May 2019 14:53
 Operator : HP/JU
 Sample : K2241-07
 Misc : LOD-WATER-5PPM
 ALS Vial : 12 Sample Multiplier: 1

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

Quant Time: May 31 11:28:47 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF052919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 29 19:08:51 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.69	152	413346	20.00	ng	0.00
21) Naphthalene-d8	7.97	136	1553928	20.00	ng	0.00
39) Acenaphthene-d10	9.72	164	776410	20.00	ng	0.00
64) Phenanthrene-d10	11.19	188	1475869	20.00	ng	0.00
76) Chrysene-d12	13.81	240	1108099	20.00	ng	0.00
87) Perylene-d12	15.20	264	1119457	20.00	ng	-0.01

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.30	112	3014289	128.03	ng	0.01
7) Phenol-d6	6.33	99	3575569	126.03	ng	0.00
23) Nitrobenzene-d5	7.26	82	2201636	88.39	ng	0.00
42) 2,4,6-Tribromophenol	10.50	330	1045052	141.26	ng	0.00
45) 2-Fluorobiphenyl	9.05	172	3696294	79.09	ng	0.00
79) Terphenyl-d14	12.78	244	3896988	66.12	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	2.38	88	59372	5.255	ng	98
3) Pyridine	3.14	79	136567	4.153	ng	95
4) n-Nitrosodimethylamine	2.98	42	55369	4.161	ng	# 100
6) Aniline	6.36	93	219294	5.117	ng	97
8) 2-Chlorophenol	6.48	128	148635	5.467	ng	95
9) Benzaldehyde	6.23	77	120624	7.452	ng	98
10) Phenol	6.34	94	181148	5.286	ng	91
11) bis(2-Chloroethyl)ether	6.43	93	140074	5.343	ng	95
12) 1,3-Dichlorobenzene	6.63	146	165876	5.355	ng	99
13) 1,4-Dichlorobenzene	6.71	146	169631	5.459	ng	98
14) 1,2-Dichlorobenzene	6.86	146	156819	5.543	ng	99
15) Benzyl Alcohol	6.83	79	147033	6.589	ng	98
16) 2,2'-oxybis(1-Chloropropan	6.97	45	195274	4.857	ng	99
17) 2-Methylphenol	6.94	107	116350	5.060	ng	98
18) Hexachloroethane	7.21	117	56521	4.804	ng	90
19) n-Nitroso-di-n-propylamine	7.10	70	100486	5.131	ng	97
20) 3+4-Methylphenols	7.10	107	151117	5.587	ng	92
22) Acetophenone	7.10	105	206649	6.241	ng	97
24) Nitrobenzene	7.27	77	153716	5.810	ng	98
25) Isophorone	7.51	82	270139	5.404	ng	100
26) 2-Nitrophenol	7.59	139	61139	4.747	ng	98
27) 2,4-Dimethylphenol	7.63	122	125802	5.873	ng	98
28) bis(2-Chloroethoxy)methane	7.73	93	169586	5.330	ng	98
29) 2,4-Dichlorophenol	7.82	162	121886	5.591	ng	96
30) 1,2,4-Trichlorobenzene	7.92	180	139937	5.634	ng	98
31) Naphthalene	7.99	128	448854	6.271	ng	97
32) Benzoic acid	7.68	122	44905	3.228	ng	98
33) 4-Chloroaniline	8.04	127	183296	5.375	ng	96
34) Hexachlorobutadiene	8.12	225	76195	5.522	ng	99
35) Caprolactam	8.36	113	36155	4.914	ng	96
36) 4-Chloro-3-methylphenol	8.52	107	120289	5.502	ng	98
37) 2-Methylnaphthalene	8.68	142	303592	6.384	ng	99
38) 1-Methylnaphthalene	8.78	142	286944	6.368	ng	98
40) 1,2,4,5-Tetrachlorobenzene	8.85	216	132923	6.429	ng	98

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41) Hexachlorocyclopentadiene	8.85	237	74551	5.456	ng	98
43) 2,4,6-Trichlorophenol	8.96	196	84534	5.196	ng	100
44) 2,4,5-Trichlorophenol	8.99	196	88069	5.425	ng	98
46) 1,1'-Biphenyl	9.15	154	369691	6.415	ng	95
47) 2-Chloronaphthalene	9.17	162	286310	6.056	ng	96
48) 2-Nitroaniline	9.25	65	66954	4.589	ng	91
49) Acenaphthylene	9.57	152	444927	6.092	ng	96
50) Dimethylphthalate	9.44	163	312976	5.785	ng	98
51) 2,6-Dinitrotoluene	9.50	165	63831	5.078	ng	97
52) Acenaphthene	9.75	154	265098	5.918	ng	98
53) 3-Nitroaniline	9.66	138	71102	4.621	ng	97
54) 2,4-Dinitrophenol	9.76	184	12242	2.469	ng	# 53
55) Dibenzofuran	9.92	168	393590	6.173	ng	92
56) 4-Nitrophenol	9.81	139	48533	4.257	ng	93
57) 2,4-Dinitrotoluene	9.89	165	78399	4.859	ng	93
58) Fluorene	10.26	166	302929	7.037	ng	98
59) 2,3,4,6-Tetrachlorophenol	10.03	232	71885	5.410	ng	98
60) Diethylphthalate	10.15	149	311845	5.644	ng	98
61) 4-Chlorophenyl-phenylether	10.26	204	146821	6.816	ng	99
62) 4-Nitroaniline	10.26	138	67503	4.502	ng	98
63) Azobenzene	10.42	77	288801	5.693	ng	96
65) 4,6-Dinitro-2-methylphenol	10.30	198	21650	3.308	ng	89
66) n-Nitrosodiphenylamine	10.37	169	262532	6.021	ng	98
67) 4-Bromophenyl-phenylether	10.75	248	87835	5.550	ng	99
68) Hexachlorobenzene	10.81	284	96276	5.616	ng	94
69) Atrazine	10.90	200	75949	5.528	ng	98
70) Pentachlorophenol	11.00	266	42988	4.350	ng	99
71) Phenanthrene	11.22	178	450260	5.981	ng	96
72) Anthracene	11.26	178	460042	6.038	ng	97
73) Carbazole	11.42	167	397025	5.929	ng	97
74) Di-n-butylphthalate	11.76	149	471051	5.789	ng	96
75) Fluoranthene	12.39	202	447723	5.743	ng	95
77) Benzidine	12.52	184	189820	4.160	ng	98
78) Pyrene	12.62	202	459078	4.898	ng	96
80) Butylbenzylphthalate	13.25	149	179456	3.809	ng	99
81) Benzo(a)anthracene	13.80	228	412966	4.850	ng	97
82) 3,3'-Dichlorobenzidine	13.76	252	141224	4.294	ng	99
83) Chrysene	13.83	228	404429	4.971	ng	96
84) Bis(2-ethylhexyl)phthalate	13.82	149	267198	4.588	ng	98
85) Di-n-octyl phthalate	14.43	149	420827	4.159	ng	97
86) Indeno(1,2,3-cd)pyrene	16.51	276	341410	4.190	ng	97
88) Benzo(b)fluoranthene	14.81	252	387279	5.493	ng	96
89) Benzo(k)fluoranthene	14.83	252	358043	5.842	ng	97
90) Benzo(a)pyrene	15.14	252	343692	5.549	ng	98
91) Dibenzo(a,h)anthracene	16.53	278	284681	5.201	ng	100
92) Benzo(g,h,i)perylene	16.90	276	268381	5.052	ng	98

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

