

Data Path : Z:\HPCHEM1\BNA F\DATA\BF040115\
 Data File : BF078164.D
 Acq On : 1 Apr 2015 21:00
 Operator : TP/IZ
 Sample : G1614-06
 Misc : LOO-WATER-20PPM
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 LOQ-WATER2015-02

Manual Integrations
 APPROVED

apatel
 4/2/2015 5:51:03 PM

Quant Time: Apr 02 13:50:01 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF040115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 02 13:21:29 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.01	152	38348	20.00	ng	0.00
21) Naphthalene-d8	8.59	136	153937	20.00	ng	0.00
38) Acenaphthene-d10	10.76	164	79882	20.00	ng	0.00
63) Phenanthrene-d10	12.59	188	152595	20.00	ng	0.00
75) Chrysene-d12	15.87	240	156711	20.00	ng	0.00
86) Perylene-d12	17.60	264	147985	20.00	ng	0.01

System Monitoring Compounds

5) 2-Fluorophenol	5.30	112	350483	150.69	ng	0.00
7) Phenol-d6	6.60	99	433875	148.85	ng	0.00
23) Nitrobenzene-d5	7.71	82	253701	99.90	ng	0.00
41) 2,4,6-Tribromophenol	11.75	330	110518	166.82	ng	0.00
44) 2-Fluorobiphenyl	9.94	172	539785	96.59	ng	0.00
78) Terphenyl-d14	14.59	244	658144	94.90	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	1.88	88	19979	19.00	ng	97
3) Pyridine	2.54	79	48678	17.67	ng	98
4) n-Nitrosodimethylamine	2.46	42	19406m	18.30	ng	
6) Aniline	6.61	93	33994	8.22	ng	# 11
8) 2-Chlorophenol	6.75	128	60945	22.16	ng	99
9) Benzaldehyde	6.46	77	44254	22.91	ng	99
10) Phenol	6.62	94	69136	19.80	ng	98
11) bis(2-Chloroethyl)ether	6.72	93	56216	22.38	ng	98
12) 1,3-Dichlorobenzene	6.93	146	61843	20.47	ng	99
13) 1,4-Dichlorobenzene	7.04	146	64667	21.27	ng	99
14) 1,2-Dichlorobenzene	7.22	146	60979	21.39	ng	99
15) Benzyl Alcohol	7.21	79	47038	22.96	ng	98
16) 2,2'-oxybis(1-Chloropropan	7.39	45	73845	22.04	ng	100
17) 2-Methylphenol	7.37	107	47134	22.07	ng	99
18) Hexachloroethane	7.64	117	21832	21.29	ng	95
19) n-Nitroso-di-n-propylamine	7.54	70	38752	20.15	ng	# 86
20) 3+4-Methylphenols	7.56	107	62302	21.87	ng	99
22) Acetophenone	7.53	105	79322	22.12	ng	# 88
24) Nitrobenzene	7.73	77	57618	21.70	ng	# 78
25) Isophorone	8.04	82	105182	21.82	ng	97
26) 2-Nitrophenol	8.13	139	28478	22.51	ng	99
27) 2,4-Dimethylphenol	8.21	122	50555	21.49	ng	98
28) bis(2-Chloroethoxy)methane	8.33	93	69948	22.63	ng	99
29) 2,4-Dichlorophenol	8.44	162	47859	22.36	ng	96
30) 1,2,4-Trichlorobenzene	8.53	180	51775	21.10	ng	98
31) Naphthalene	8.62	128	170683	21.30	ng	99
32) Benzoic acid	8.32	122	18706	20.66	ng	90
33) 4-Chloroaniline	8.70	127	38274	11.51	ng	99
34) Hexachlorobutadiene	8.77	225	28501	21.15	ng	96
35) Caprolactam	9.12	113	13867	21.70	ng	94
36) 4-Chloro-3-methylphenol	9.32	107	50894	23.57	ng	98
37) 2-Methylnaphthalene	9.48	142	120842	22.21	ng	99
39) 1,2,4,5-Tetrachlorobenzene	9.69	216	51802	20.93	ng	100
40) Hexachlorocyclopentadiene	9.68	237	18517	22.71	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.84	196	33167	21.25	nq	98
43) 2,4,5-Trichlorophenol	9.88	196	36159	22.17	nq	98
45) 1,1'-Biphenyl	10.07	154	146551	22.43	nq	98
46) 2-Chloronaphthalene	10.08	162	111312	21.65	nq	99
47) 2-Nitroaniline	10.21	65	28449	22.37	nq	97
48) Acenaphthylene	10.59	152	178161	21.46	nq	99
49) Dimethylphthalate	10.45	163	129608	21.60	nq	99
50) 2,6-Dinitrotoluene	10.52	165	27130	21.97	nq	# 46
51) Acenaphthene	10.80	154	100944	21.60	nq	99
52) 3-Nitroaniline	10.73	138	25567	18.21	nq	94
53) 2,4-Dinitrophenol	10.87	184	4804	20.46	nq	98
54) Dibenzofuran	11.01	168	161853	22.67	nq	99
55) 4-Nitrophenol	10.97	139	19182	20.68	nq	98
56) 2,4-Dinitrotoluene	11.03	165	35685	22.31	nq	94
57) Fluorene	11.44	166	130697	22.59	nq	100
58) 2,3,4,6-Tetrachlorophenol	11.17	232	27605	22.83	nq	99
59) Diethylphthalate	11.33	149	128063	22.13	nq	98
60) 4-Chlorophenyl-phenylether	11.45	204	60091	22.57	nq	# 75
61) 4-Nitroaniline	11.48	138	28417	20.02	nq	93
62) Azobenzene	11.64	77	117244	22.20	nq	98
64) 4,6-Dinitro-2-methylphenol	11.52	198	12241	21.84	nq	83
65) n-Nitrosodiphenylamine	11.60	169	111095	21.05	nq	98
66) 4-Bromophenyl-phenylether	12.05	248	35183	21.77	nq	97
67) Hexachlorobenzene	12.11	284	37998	21.46	nq	94
68) Atrazine	12.28	200	35057	22.93	nq	98
69) Pentachlorophenol	12.36	266	11372	20.84	nq	98
70) Phenanthrene	12.63	178	187211	21.21	nq	99
71) Anthracene	12.68	178	188292	21.65	nq	100
72) Carbazole	12.90	167	179844	22.06	nq	99
73) Di-n-butylphthalate	13.36	149	201738	21.85	nq	100
74) Fluoranthene	14.09	202	202237	21.73	nq	97
76) Benzidine	14.28	184	81935	13.58	nq	99
77) Pyrene	14.36	202	211473	21.50	nq	98
79) Butylbenzylphthalate	15.21	149	83955	22.39	nq	90
80) Benzo(a)anthracene	15.85	228	194665	20.88	nq	98
81) 3,3'-Dichlorobenzidine	15.84	252	32104	10.28	nq	# 94
82) Chrysene	15.89	228	190538	21.75	nq	99
83) Bis(2-ethylhexyl)phthalate	15.93	149	129268	21.98	nq	# 95
84) Di-n-octyl phthalate	16.73	149	214570	22.22	nq	# 100
85) Indeno(1,2,3-cd)pyrene	18.91	276	208057	20.93	nq	# 87
87) Benzo(b)fluoranthene	17.15	252	189004	20.67	nq	99
88) Benzo(k)fluoranthene	17.19	252	190668m	23.46	nq	
89) Benzo(a)pyrene	17.53	252	178601	22.38	nq	# 95
90) Dibenzo(a,h)anthracene	18.93	278	169180	21.17	nq	97
91) Benzo(g,h,i)perylene	19.29	276	171060	21.35	nq	98

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

