

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF032519\
 Data File : BF113281.D
 Acq On : 25 Mar 2019 16:07
 Operator : JU/SJ
 Sample : K1560-08
 Misc : L00 20 PPM
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 LOQ-WATER-02-QT1-2019

Manual Integrations
 APPROVED

Sohil
 3/26/2019 6:21:09 PM

Quant Time: Mar 26 04:51:36 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF032519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 26 03:32:08 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.70	152	159056	20.00	ng	0.00
21) Naphthalene-d8	7.99	136	673926	20.00	ng	0.00
39) Acenaphthene-d10	9.73	164	335342	20.00	ng	0.00
64) Phenanthrene-d10	11.22	188	671191	20.00	ng	0.00
76) Chrysene-d12	13.84	240	568376	20.00	ng	0.00
87) Perylene-d12	15.24	264	525784	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.32	112	982421	100.99	ng	0.00
7) Phenol-d6	6.35	99	1321400	107.38	ng	0.00
23) Nitrobenzene-d5	7.27	82	822276	72.42	ng	0.00
42) 2,4,6-Tribromophenol	10.53	330	435596	105.16	ng	0.00
45) 2-Fluorobiphenyl	9.06	172	1505685	66.55	ng	0.00
79) Terphenyl-d14	12.80	244	1716166	66.55	ng	0.00

Target Compounds				Qvalue		
2) 1,4-Dioxane	2.36	88	83210	16.860	ng	99
3) Pyridine	3.10	79	207145	17.025	ng	97
4) n-Nitrosodimethylamine	2.99	42	85163	15.745	ng	93
6) Aniline	6.37	93	273509	18.251	ng	# 79
8) 2-Chlorophenol	6.49	128	194583	17.226	ng	96
9) Benzaldehyde	6.25	77	157917	19.124	ng	97
10) Phenol	6.36	94	256698	18.272	ng	97
11) bis(2-Chloroethyl)ether	6.44	93	196066	18.207	ng	98
12) 1,3-Dichlorobenzene	6.65	146	212200	17.427	ng	98
13) 1,4-Dichlorobenzene	6.72	146	215865	18.361	ng	99
14) 1,2-Dichlorobenzene	6.87	146	206919	17.890	ng	98
15) Benzyl Alcohol	6.85	79	148375	18.277	ng	97
16) 2,2'-oxybis(1-Chloropropan	6.99	45	319751	19.765	ng	96
17) 2-Methylphenol	6.97	107	166862	18.849	ng	96
18) Hexachloroethane	7.22	117	84476	20.005	ng	98
19) n-Nitroso-di-n-propylamine	7.12	70	147087	19.029	ng	98
20) 3+4-Methylphenols	7.12	107	216081	16.511	ng	93
22) Acetophenone	7.11	105	283501	18.446	ng	98
24) Nitrobenzene	7.29	77	226291	19.099	ng	98
25) Isophorone	7.52	82	400790	18.214	ng	96
26) 2-Nitrophenol	7.60	139	115734	18.480	ng	94
27) 2,4-Dimethylphenol	7.65	122	168943	18.753	ng	98
28) bis(2-Chloroethoxy)methane	7.74	93	263914	19.046	ng	99
29) 2,4-Dichlorophenol	7.85	162	181849	18.627	ng	99
30) 1,2,4-Trichlorobenzene	7.93	180	197910	18.789	ng	100
31) Naphthalene	8.01	128	633789	19.245	ng	98
32) Benzoic acid	7.76	122	102748	14.169	ng	99
33) 4-Chloroaniline	8.06	127	249344	19.416	ng	99
34) Hexachlorobutadiene	8.13	225	119027	19.013	ng	97
35) Caprolactam	8.41	113	56989	18.188	ng	99
36) 4-Chloro-3-methylphenol	8.55	107	187570	19.102	ng	94
37) 2-Methylnaphthalene	8.70	142	414987	20.195	ng	98
38) 1-Methylnaphthalene	8.80	142	401922	20.335	ng	99
40) 1,2,4,5-Tetrachlorobenzene	8.87	216	198266	18.419	ng	99

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF032519\
 Data File : BF113281.D
 Acq On : 25 Mar 2019 16:07
 Operator : JU/SJ
 Sample : K1560-08
 Misc : L00 20 PPM
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 LOQ-WATER-02-QT1-2019

Manual Integrations
 APPROVED

Sohil
 3/26/2019 6:21:09 PM

Quant Time: Mar 26 04:51:36 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF032519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 26 03:32:08 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	8.86	237	57788	12.683	nq	97
43) 2,4,6-Trichlorophenol	8.98	196	120280	17.159	nq	98
44) 2,4,5-Trichlorophenol	9.03	196	130982	16.884	nq	98
46) 1,1'-Biphenyl	9.16	154	523442	18.689	nq	99
47) 2-Chloronaphthalene	9.19	162	390866	18.126	nq	97
48) 2-Nitroaniline	9.28	65	117247	17.325	nq	98
49) Acenaphthylene	9.60	152	664217	19.050	nq	99
50) Dimethylphthalate	9.46	163	447081	17.990	nq	100
51) 2,6-Dinitrotoluene	9.52	165	100439	17.915	nq	91
52) Acenaphthene	9.77	154	364398	18.558	nq	99
53) 3-Nitroaniline	9.69	138	112125	17.734	nq	92
54) 2,4-Dinitrophenol	9.80	184	40868	14.656	nq	95
55) Dibenzofuran	9.94	168	568301	19.252	nq	98
56) 4-Nitrophenol	9.87	139	70675	15.679	nq	92
57) 2,4-Dinitrotoluene	9.93	165	130555	18.311	nq	93
58) Fluorene	10.28	166	438369	19.116	nq	100
59) 2,3,4,6-Tetrachlorophenol	10.06	232	118703	18.094	nq	97
60) Diethylphthalate	10.16	149	444617	18.238	nq	99
61) 4-Chlorophenyl-phenylether	10.27	204	214793	19.218	nq	93
62) 4-Nitroaniline	10.30	138	116694	17.927	nq	97
63) Azobenzene	10.43	77	429705	18.724	nq	97
65) 4,6-Dinitro-2-methylphenol	10.34	198	58411	15.908	nq	98
66) n-Nitrosodiphenylamine	10.39	169	381562	18.526	nq	99
67) 4-Bromophenyl-phenylether	10.76	248	137777	18.834	nq	93
68) Hexachlorobenzene	10.83	284	158808	18.702	nq	97
69) Atrazine	10.92	200	120578	18.548	nq	95
70) Pentachlorophenol	11.03	266	67733	15.310	nq	97
71) Phenanthrene	11.24	178	642953	19.290	nq	99
72) Anthracene	11.29	178	656928	19.404	nq	99
73) Carbazole	11.44	167	624319	19.325	nq	98
74) Di-n-butylphthalate	11.78	149	719594	19.208	nq	98
75) Fluoranthene	12.42	202	716587	19.017	nq	98
77) Benzidine	12.54	184	334848m	15.384	nq	
78) Pyrene	12.64	202	734908	18.659	nq	99
80) Butylbenzylphthalate	13.26	149	299510	17.255	nq	97
81) Benzo(a)anthracene	13.83	228	606466	18.641	nq	98
82) 3,3'-Dichlorobenzidine	13.79	252	237042	18.161	nq	99
83) Chrysene	13.87	228	626794	18.967	nq	98
84) Bis(2-ethylhexyl)phthalate	13.83	149	387295	18.201	nq	99
85) Di-n-octyl phthalate	14.43	149	678081	18.773	nq	100
86) Indeno(1,2,3-cd)pyrene	16.59	276	518822	18.293	nq	98
88) Benzo(b)fluoranthene	14.84	252	575044	17.606	nq	99
89) Benzo(k)fluoranthene	14.87	252	550987	19.210	nq	100
90) Benzo(a)pyrene	15.19	252	518464	17.905	nq	98
91) Dibenzo(a,h)anthracene	16.60	278	435584	17.398	nq	100
92) Benzo(g,h,i)perylene	16.99	276	416741	16.508	nq	98

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

