

Data Path : Z:\HPCHEM1\BNA F\DATA\BF081917\
 Data File : BF097878.D
 Acq On : 19 Aug 2017 14:57
 Operator : SJ/JU
 Sample : I4846-02MS
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PI-2(3)MS

Manual Integrations
 APPROVED

Sohil
 8/21/2017 5:19:54 PM

Quant Time: Aug 21 04:22:39 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF081517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 15 13:25:08 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.76	152	144746	20.00	ng	-0.02
21) Naphthalene-d8	8.05	136	594923	20.00	ng	-0.02
38) Acenaphthene-d10	9.80	164	265153	20.00	ng	-0.02
63) Phenanthrene-d10	11.28	188	459066	20.00	ng	-0.02
75) Chrysene-d12	13.91	240	261092	20.00	ng	-0.03
86) Perylene-d12	15.33	264	276142	20.00	ng	-0.03

System Monitoring Compounds

5) 2-Fluorophenol	5.38	112	1207246	126.86	ng	0.00
7) Phenol-d6	6.40	99	1685261	130.34	ng	0.00
23) Nitrobenzene-d5	7.33	82	1080478	98.66	ng	-0.02
41) 2,4,6-Tribromophenol	10.59	330	314839	136.85	ng	-0.02
44) 2-Fluorobiphenyl	9.13	172	1531336	84.85	ng	-0.02
78) Terphenyl-d14	12.86	244	1031053	82.35	ng	-0.03

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.48	88	116010	21.860	ng	88
3) Pyridine	3.20	79	356178	24.277	ng	88
4) n-Nitrosodimethylamine	3.14	42	154126	27.302	ng	86
6) Aniline	6.43	93	444559	24.475	ng	98
8) 2-Chlorophenol	6.55	128	309969	30.887	ng	96
9) Benzaldehyde	6.31	77	121235	14.803	ng	86
10) Phenol	6.42	94	452914	29.951	ng	93
11) bis(2-Chloroethyl)ether	6.50	93	364727	31.956	ng	97
12) 1,3-Dichlorobenzene	6.70	146	304155	28.176	ng	# 92
13) 1,4-Dichlorobenzene	6.78	146	304366	27.723	ng	# 93
14) 1,2-Dichlorobenzene	6.93	146	287578	28.408	ng	99
15) Benzyl Alcohol	6.91	79	327318	33.813	ng	96
16) 2,2'-oxybis(1-Chloropropan	7.05	45	478250	31.144	ng	91
17) 2-Methylphenol	7.02	107	294222	33.268	ng	96
18) Hexachloroethane	7.27	117	127556	29.634	ng	# 80
19) n-Nitroso-di-n-propylamine	7.19	70	271744	30.702	ng	89
20) 3+4-Methylphenols	7.17	107	345684	32.002	ng	93
22) Acetophenone	7.17	105	464630	29.563	ng	# 93
24) Nitrobenzene	7.35	77	410775	33.688	ng	95
25) Isophorone	7.59	82	769853	33.808	ng	96
26) 2-Nitrophenol	7.66	139	161307	37.772	ng	# 77
27) 2,4-Dimethylphenol	7.70	122	301461	31.743	ng	93
28) bis(2-Chloroethoxy)methane	7.80	93	478187	31.941	ng	99
29) 2,4-Dichlorophenol	7.90	162	260072	32.990	ng	93
30) 1,2,4-Trichlorobenzene	7.99	180	257877	29.508	ng	98
31) Naphthalene	8.07	128	920250	29.796	ng	100
32) Benzoic acid	7.82	122	192332	30.491	ng	# 88
33) 4-Chloroaniline	8.12	127	329628	25.306	ng	98
34) Hexachlorobutadiene	8.19	225	151248	29.641	ng	98
35) Caprolactam	8.49	113	89837	33.520	ng	98
36) 4-Chloro-3-methylphenol	8.60	107	321275	31.719	ng	# 77
37) 2-Methylnaphthalene	8.76	142	612401	32.341	ng	97
39) 1,2,4,5-Tetrachlorobenzene	8.92	216	247128	30.369	ng	# 98
40) Hexachlorocyclopentadiene	8.91	237	218552	55.905	ng	100

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.03	196	176570	32.319	nq	97
43) 2,4,5-Trichlorophenol	9.07	196	187404	34.577	nq	95
45) 1,1'-Biphenyl	9.22	154	735050	30.400	nq	96
46) 2-Chloronaphthalene	9.25	162	533173	29.843	nq #	91
47) 2-Nitroaniline	9.34	65	212721	35.517	nq	93
48) Acenaphthylene	9.66	152	859686	30.642	nq	100
49) Dimethylphthalate	9.53	163	814635	42.031	nq	100
50) 2,6-Dinitrotoluene	9.59	165	137506	35.542	nq #	61
51) Acenaphthene	9.83	154	533955	30.177	nq	97
52) 3-Nitroaniline	9.75	138	133104	26.666	nq	84
53) 2,4-Dinitrophenol	9.86	184	95000	55.219	nq #	79
54) Dibenzofuran	10.00	168	774252	32.649	nq	100
55) 4-Nitrophenol	9.92	139	214641	53.906	nq #	36
56) 2,4-Dinitrotoluene	9.99	165	175847	36.218	nq #	53
57) Fluorene	10.35	166	558570	30.465	nq	96
58) 2,3,4,6-Tetrachlorophenol	10.12	232	147405	35.127	nq	95
59) Diethylphthalate	10.23	149	696033	34.341	nq	99
60) 4-Chlorophenyl-phenylether	10.34	204	269148	31.599	nq	89
61) 4-Nitroaniline	10.36	138	137392	28.961	nq #	68
62) Azobenzene	10.50	77	809169	33.609	nq	93
64) 4,6-Dinitro-2-methylphenol	10.39	198	65729	33.438	nq	85
65) n-Nitrosodiphenylamine	10.46	169	521326	33.349	nq	100
66) 4-Bromophenyl-phenylether	10.83	248	167113	35.055	nq	97
67) Hexachlorobenzene	10.89	284	152137	31.311	nq #	88
68) Atrazine	10.99	200	153358	36.992	nq	98
69) Pentachlorophenol	11.09	266	157732	55.676	nq	99
70) Phenanthrene	11.30	178	901424	36.849	nq	99
71) Anthracene	11.36	178	797211	32.131	nq	99
72) Carbazole	11.51	167	688453	29.232	nq	99
73) Di-n-butylphthalate	11.84	149	1020669	37.625	nq	99
74) Fluoranthene	12.49	202	908101	35.566	nq	98
76) Benzidine	12.61	184	142103m	16.600	nq	
77) Pyrene	12.72	202	858820	40.218	nq	99
79) Butylbenzylphthalate	13.34	149	337099	41.173	nq #	69
80) Benzo(a)anthracene	13.90	228	539632	34.485	nq	99
81) 3,3'-Dichlorobenzidine	13.86	252	161261	30.540	nq #	97
82) Chrysene	13.94	228	557228	35.797	nq	99
83) Bis(2-ethylhexyl)phthalate	13.90	149	431001	47.507	nq #	100
84) Di-n-octyl phthalate	14.51	149	753913	47.759	nq	99
85) Indeno(1,2,3-cd)pyrene	16.72	276	592188	51.714	nq	93
87) Benzo(b)fluoranthene	14.93	252	591579	33.987	nq	99
88) Benzo(k)fluoranthene	14.96	252	525515m	32.136	nq	
89) Benzo(a)pyrene	15.27	252	565508	36.699	nq	99
90) Dibenzo(a,h)anthracene	16.74	278	451013	35.952	nq	97
91) Benzo(g,h,i)perylene	17.14	276	492794	37.648	nq	94

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

