

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF050317\
 Data File : BF094863.D
 Acq On : 4 May 2017 7:48
 Operator : SJ/MA
 Sample : I2897-01MS
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OWS-1MS

Manual Integrations
 APPROVED

mohammad
 5/4/2017 3:35:30 PM

Quant Time: May 04 08:34:42 2017
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050217.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 03 17:24:51 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.73	152	97711	20.00	ng	0.00
21) Naphthalene-d8	9.75	136	357515	20.00	ng	0.00
38) Acenaphthene-d10	12.58	164	162610	20.00	ng	0.00
63) Phenanthrene-d10	14.98	188	254502	20.00	ng	0.00
75) Chrysene-d12	18.64	240	193920	20.00	ng	-0.01
86) Perylene-d12	20.31	264	157349	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.70	112	760171	129.71	ng	0.02
7) Phenol-d6	7.28	99	871156	123.88	ng	0.00
23) Nitrobenzene-d5	8.63	82	581690	102.60	ng	0.00
41) 2,4,6-Tribromophenol	13.88	330	189794	142.52	ng	0.00
44) 2-Fluorobiphenyl	11.52	172	1097917	97.18	ng	0.00
78) Terphenyl-d14	17.30	244	721768	81.98	ng	-0.01

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.24	88	94546	32.89	ng	# 80
3) Pyridine	2.96	79	176996m	26.57	ng	
4) n-Nitrosodimethylamine	2.79	42	113652	40.93	ng	86
6) Aniline	7.24	93	167930	18.92	ng	95
8) 2-Chlorophenol	7.42	128	310674	46.57	ng	94
9) Benzaldehyde	7.05	77	20817	4.85	ng	98
10) Phenol	7.30	94	324582	43.80	ng	90
11) bis(2-Chloroethyl)ether	7.37	93	281943	46.09	ng	94
12) 1,3-Dichlorobenzene	7.64	146	321022	42.42	ng	98
13) 1,4-Dichlorobenzene	7.76	146	335388	43.71	ng	97
14) 1,2-Dichlorobenzene	8.00	146	320878	44.80	ng	95
15) Benzyl Alcohol	8.02	79	244597	54.08	ng	98
16) 2,2'-oxybis(1-Chloropropan	8.22	45	368105	42.51	ng	75
17) 2-Methylphenol	8.23	107	246276	45.11	ng	97
18) Hexachloroethane	8.51	117	102318	39.36	ng	# 84
19) n-Nitroso-di-n-propylamine	8.44	70	215309	45.27	ng	95
20) 3+4-Methylphenols	8.48	107	400140	53.94	ng	# 58
22) Acetophenone	8.40	105	432826	50.46	ng	# 84
24) Nitrobenzene	8.66	77	297315	50.03	ng	94
25) Isophorone	9.06	82	548250	54.16	ng	# 94
26) 2-Nitrophenol	9.17	139	170273	53.10	ng	97
27) 2,4-Dimethylphenol	9.31	122	269464	51.97	ng	99
28) bis(2-Chloroethoxy)methane	9.43	93	354676	50.64	ng	98
29) 2,4-Dichlorophenol	9.58	162	264634	54.06	ng	99
30) 1,2,4-Trichlorobenzene	9.67	180	251470	47.95	ng	99
31) Naphthalene	9.78	128	870097	48.42	ng	100
32) Benzoic acid	9.77	122	599335m	162.51	ng	
33) 4-Chloroaniline	9.94	127	130794	19.53	ng	94
34) Hexachlorobutadiene	10.00	225	128306	46.37	ng	99
35) Caprolactam	10.67	113	60693	41.29	ng	91
36) 4-Chloro-3-methylphenol	10.78	107	258279	49.35	ng	# 88
37) 2-Methylnaphthalene	10.91	142	587950	49.86	ng	98
39) 1,2,4,5-Tetrachlorobenzene	11.18	216	236448	47.28	ng	99
40) Hexachlorocyclopentadiene	11.17	237	54174	29.87	ng	97

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF050317\
 Data File : BF094863.D
 Acq On : 4 May 2017 7:48
 Operator : SJ/MA
 Sample : I2897-01MS
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OWS-1MS

Manual Integrations
 APPROVED

mohammad
 5/4/2017 3:35:30 PM

Quant Time: May 04 08:34:42 2017
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050217.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 03 17:24:51 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	11.40	196	166338	49.82	ng	98
43) 2,4,5-Trichlorophenol	11.49	196	177254	49.39	ng #	91
45) 1,1'-Biphenyl	11.68	154	669956	48.93	ng	97
46) 2-Chloronaphthalene	11.69	162	520490	47.08	ng	95
47) 2-Nitroaniline	11.90	65	149670	49.47	ng #	79
48) Acenaphthylene	12.35	152	812781	46.94	ng	99
49) Dimethylphthalate	12.22	163	657456	49.25	ng	99
50) 2,6-Dinitrotoluene	12.31	165	133707	49.10	ng	93
51) Acenaphthene	12.63	154	495185	47.88	ng	99
52) 3-Nitroaniline	12.58	138	47836	16.36	ng #	86
53) 2,4-Dinitrophenol	12.77	184	72063	50.54	ng #	66
54) Dibenzofuran	12.92	168	745038	52.06	ng	98
55) 4-Nitrophenol	12.96	139	168658	96.07	ng #	72
56) 2,4-Dinitrotoluene	12.97	165	171598	49.03	ng #	84
57) Fluorene	13.47	166	588970	47.33	ng	99
58) 2,3,4,6-Tetrachlorophenol	13.16	232	132806	58.80	ng #	93
59) Diethylphthalate	13.37	149	595469	46.54	ng	99
60) 4-Chlorophenyl-phenylether	13.50	204	265981	48.63	ng	90
61) 4-Nitroaniline	13.57	138	117093	43.64	ng	98
62) Azobenzene	13.75	77	518280	50.41	ng	95
64) 4,6-Dinitro-2-methylphenol	13.63	198	50761	29.75	ng #	73
65) n-Nitrosodiphenylamine	13.71	169	475177	51.44	ng	99
66) 4-Bromophenyl-phenylether	14.28	248	144217	53.64	ng	99
67) Hexachlorobenzene	14.37	284	139477	50.62	ng #	88
68) Atrazine	14.62	200	118387	45.39	ng	96
69) Pentachlorophenol	14.72	266	136714	107.55	ng	99
70) Phenanthrene	15.01	178	723451	49.47	ng	99
71) Anthracene	15.09	178	717226	48.02	ng	98
72) Carbazole	15.38	167	669943	49.29	ng	97
73) Di-n-butylphthalate	15.95	149	845912	49.91	ng	99
74) Fluoranthene	16.75	202	691942	46.13	ng	99
76) Benzidine	16.98	184	32359	11.91	ng #	94
77) Pyrene	17.05	202	687292	47.39	ng	99
79) Butylbenzylphthalate	17.97	149	326827	48.45	ng	92
80) Benzo(a)anthracene	18.63	228	555349	49.21	ng	98
81) 3,3'-Dichlorobenzidine	18.61	252	146802	37.66	ng #	96
82) Chrysene	18.68	228	517054	47.38	ng	99
83) Bis(2-ethylhexyl)phthalate	18.71	149	444311	46.23	ng #	99
84) Di-n-octyl phthalate	19.48	149	735182	46.65	ng	97
85) Indeno(1,2,3-cd)pyrene	21.59	276	409667	46.23	ng	100
87) Benzo(b)fluoranthene	19.90	252	483871	49.77	ng	99
88) Benzo(k)fluoranthene	19.93	252	495305	52.81	ng #	95
89) Benzo(a)pyrene	20.25	252	461639	51.45	ng	99
90) Dibenzo(a,h)anthracene	21.61	278	349772	47.21	ng	99
91) Benzo(g,h,i)perylene	21.96	276	336662	46.30	ng	99

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

