

Data Path : Z:\HPCHEM1\BNA F\DATA\BF040215\
 Data File : BF078235.D
 Acq On : 3 Apr 2015 9:50
 Operator : TP/IZ
 Sample : G1731-04MSD
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 COMP-4MSD

Manual Integrations
 APPROVED

apatel
 4/6/2015 12:06:20 PM

Quant Time: Apr 03 12:34:32 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	7.02	152	48843	20.00	ng	0.01
21) Naphthalene-d8	8.60	136	201123	20.00	ng	0.01
38) Acenaphthene-d10	10.76	164	103264	20.00	ng	0.00
63) Phenanthrene-d10	12.60	188	199663	20.00	ng	0.01
75) Chrysene-d12	15.88	240	214501	20.00	ng	0.01
86) Perylene-d12	17.68	264	210732	20.00	ng	0.09

System Monitoring Compounds

5) 2-Fluorophenol	5.32	112	233403	78.79	ng	0.02
7) Phenol-d6	6.64	99	300872	81.04	ng	0.03
23) Nitrobenzene-d5	7.72	82	175105	52.77	ng	0.01
41) 2,4,6-Tribromophenol	11.74	330	71526	83.52	ng	0.00
44) 2-Fluorobiphenyl	9.95	172	366907	50.79	ng	0.01
78) Terphenyl-d14	14.59	244	411371	43.34	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	1.87	88	22298	16.65	ng	87
3) Pyridine	2.52	79	63937	18.22	ng	93
4) n-Nitrosodimethylamine	2.44	42	37657m	27.88	ng	
6) Aniline	6.61	93	79735	15.14	ng	# 86
8) 2-Chlorophenol	6.76	128	92744	26.48	ng	98
10) Phenol	6.65	94	117227	26.35	ng	96
11) bis(2-Chloroethyl)ether	6.72	93	79793	24.94	ng	97
12) 1,3-Dichlorobenzene	6.94	146	89250	23.20	ng	# 90
13) 1,4-Dichlorobenzene	7.04	146	91127	23.53	ng	98
14) 1,2-Dichlorobenzene	7.22	146	88104	24.26	ng	98
15) Benzyl Alcohol	7.22	79	72002	27.60	ng	94
16) 2,2'-oxybis(1-Chloropropan	7.39	45	110230	25.83	ng	# 5
17) 2-Methylphenol	7.39	107	73776	27.13	ng	94
18) Hexachloroethane	7.64	117	30209	23.13	ng	93
19) n-Nitroso-di-n-propylamine	7.55	70	64156	26.19	ng	92
20) 3+4-Methylphenols	7.58	107	95467	26.31	ng	97
22) Acetophenone	7.54	105	122500	26.14	ng	# 97
24) Nitrobenzene	7.74	77	91669	26.43	ng	95
25) Isophorone	8.04	82	166202	26.39	ng	93
26) 2-Nitrophenol	8.13	139	39538	23.92	ng	92
27) 2,4-Dimethylphenol	8.22	122	78739	25.62	ng	94
28) bis(2-Chloroethoxy)methane	8.33	93	102947	25.49	ng	98
29) 2,4-Dichlorophenol	8.45	162	76091	27.21	ng	97
30) 1,2,4-Trichlorobenzene	8.53	180	77885	24.29	ng	99
31) Naphthalene	8.62	128	266236	25.43	ng	99
32) Benzoic acid	8.34	122	31045	24.68	ng	91
33) 4-Chloroaniline	8.72	127	74296	17.10	ng	96
34) Hexachlorobutadiene	8.78	225	42284	24.02	ng	98
35) Caprolactam	9.14	113	22012	26.37	ng	94
36) 4-Chloro-3-methylphenol	9.34	107	81862	29.01	ng	91
37) 2-Methylnaphthalene	9.48	142	182528	25.68	ng	98
39) 1,2,4,5-Tetrachlorobenzene	9.69	216	78489	24.53	ng	99
40) Hexachlorocyclopentadiene	9.68	237	39850	33.28	ng	96
42) 2,4,6-Trichlorophenol	9.85	196	54348	26.93	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
43) 2,4,5-Trichlorophenol	9.90	196	55765	26.45	nq	# 83
45) 1,1'-Biphenyl	10.06	154	214783	25.43	nq	99
46) 2-Chloronaphthalene	10.08	162	167809	25.25	nq	97
47) 2-Nitroaniline	10.22	65	49840	30.31	nq	# 84
48) Acenaphthylene	10.59	152	271744	25.32	nq	100
49) Dimethylphthalate	10.46	163	210954	27.19	nq	# 97
50) 2,6-Dinitrotoluene	10.53	165	42820	26.83	nq	# 76
51) Acenaphthene	10.81	154	153149	25.35	nq	99
52) 3-Nitroaniline	10.74	138	45111	24.86	nq	93
53) 2,4-Dinitrophenol	10.88	184	8644	24.93	nq	# 90
54) Dibenzofuran	11.01	168	233495	25.30	nq	95
55) 4-Nitrophenol	10.99	139	64503m	49.36	nq	
56) 2,4-Dinitrotoluene	11.02	165	55637	26.91	nq	# 80
57) Fluorene	11.44	166	195717	26.16	nq	99
58) 2,3,4,6-Tetrachlorophenol	11.18	232	43305	27.71	nq	99
59) Diethylphthalate	11.33	149	194007	25.94	nq	98
60) 4-Chlorophenyl-phenylether	11.46	204	86871	25.24	nq	94
61) 4-Nitroaniline	11.49	138	46260	25.21	nq	97
62) Azobenzene	11.65	77	201876	29.57	nq	85
64) 4,6-Dinitro-2-methylphenol	11.53	198	9557	16.49	nq	# 58
65) n-Nitrosodiphenylamine	11.61	169	171933	24.89	nq	99
66) 4-Bromophenyl-phenylether	12.05	248	52619	24.88	nq	93
67) Hexachlorobenzene	12.11	284	55132	23.79	nq	96
68) Atrazine	12.28	200	49624	24.81	nq	93
69) Pentachlorophenol	12.37	266	48751	50.60	nq	99
70) Phenanthrene	12.62	178	284133	24.60	nq	100
71) Anthracene	12.69	178	282523	24.82	nq	99
72) Carbazole	12.91	167	289317	27.13	nq	97
73) Di-n-butylphthalate	13.36	149	319863	26.47	nq	99
74) Fluoranthene	14.10	202	314260	25.80	nq	92
76) Benzidine	14.29	184	148921	18.03	nq	97
77) Pyrene	14.37	202	333336	24.75	nq	99
79) Butylbenzylphthalate	15.22	149	148577	28.95	nq	# 81
80) Benzo(a)anthracene	15.87	228	307858	24.12	nq	99
81) 3,3'-Dichlorobenzidine	15.86	252	103574	24.23	nq	# 99
82) Chrysene	15.92	228	294476	24.56	nq	99
83) Bis(2-ethylhexyl)phthalate	15.95	149	224159	27.85	nq	97
84) Di-n-octyl phthalate	16.76	149	392957	29.73	nq	# 100
85) Indeno(1,2,3-cd)pyrene	19.01	276	346688	25.48	nq	# 86
87) Benzo(b)fluoranthene	17.21	252	298459m	22.92	nq	
88) Benzo(k)fluoranthene	17.24	252	308594m	26.66	nq	
89) Benzo(a)pyrene	17.61	252	305959	26.92	nq	99
90) Dibenzo(a,h)anthracene	19.05	278	277702	24.40	nq	100
91) Benzo(g,h,i)perylene	19.40	276	286897	25.15	nq	99

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

