

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF030215\
 Data File : BF077531.D
 Acq On : 2 Mar 2015 19:24
 Operator : TP/IZ
 Sample : G1399-05MS
 Misc : S
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB13MS

Manual Integrations
 APPROVED

apatel
 3/3/2015 6:26:27 PM

Quant Time: Mar 03 04:10:07 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF021915.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 02 23:46:07 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.26	152	98544	20.00	ng	0.00
21) Naphthalene-d8	8.84	136	397874	20.00	ng	0.00
38) Acenaphthene-d10	11.01	164	193937	20.00	ng	0.00
63) Phenanthrene-d10	12.86	188	372820	20.00	ng	0.00
75) Chrysene-d12	16.15	240	414989	20.00	ng	0.02
86) Perylene-d12	17.92	264	410815	20.00	ng	0.09

System Monitoring Compounds

5) 2-Fluorophenol	5.56	112	482203	81.69	ng	0.01
7) Phenol-d6	6.83	99	627616	82.69	ng	0.01
23) Nitrobenzene-d5	7.95	82	335589	52.45	ng	0.00
41) 2,4,6-Tribromophenol	12.00	330	162830	87.20	ng	0.00
44) 2-Fluorobiphenyl	10.19	172	729833	58.68	ng	0.00
78) Terphenyl-d14	14.85	244	795598	44.82	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.21	88	52090	20.80	ng	94
3) Pyridine	2.96	79	151744	21.18	ng	98
4) n-Nitrosodimethylamine	2.86	42	76950	24.70	ng	99
6) Aniline	6.85	93	103338	9.85	ng	# 84
8) 2-Chlorophenol	6.99	128	163623	23.50	ng	97
9) Benzaldehyde	6.71	77	33895	7.36	ng	87
10) Phenol	6.84	94	222028	24.66	ng	89
11) bis(2-Chloroethyl)ether	6.95	93	153712	23.75	ng	100
12) 1,3-Dichlorobenzene	7.19	146	179455	23.67	ng	96
13) 1,4-Dichlorobenzene	7.28	146	183208	23.75	ng	98
14) 1,2-Dichlorobenzene	7.47	146	173098	23.59	ng	98
15) Benzyl Alcohol	7.44	79	135266	23.45	ng	93
16) 2,2'-oxybis(1-Chloropropan	7.62	45	212198	22.94	ng	96
17) 2-Methylphenol	7.59	107	127984	23.51	ng	98
18) Hexachloroethane	7.88	117	53396	20.72	ng	# 83
19) n-Nitroso-di-n-propylamine	7.78	70	118596	24.61	ng	94
20) 3+4-Methylphenols	7.78	107	176993	24.69	ng	# 79
22) Acetophenone	7.77	105	234080	25.01	ng	# 88
24) Nitrobenzene	7.98	77	159570	23.70	ng	# 82
25) Isophorone	8.28	82	305873	24.10	ng	96
26) 2-Nitrophenol	8.37	139	73555	26.66	ng	93
27) 2,4-Dimethylphenol	8.43	122	144685	23.44	ng	97
28) bis(2-Chloroethoxy)methane	8.56	93	198300	24.73	ng	98
29) 2,4-Dichlorophenol	8.67	162	143126	24.70	ng	95
30) 1,2,4-Trichlorobenzene	8.77	180	153103	24.03	ng	99
31) Naphthalene	8.87	128	486559	24.45	ng	99
32) Benzoic acid	8.52	122	28496	9.50	ng	# 81
33) 4-Chloroaniline	8.94	127	107874	12.19	ng	98
34) Hexachlorobutadiene	9.02	225	82804	22.77	ng	98
35) Caprolactam	9.36	113	41640	23.54	ng	90
36) 4-Chloro-3-methylphenol	9.55	107	139802	24.00	ng	90
37) 2-Methylnaphthalene	9.72	142	327179	23.16	ng	97
39) 1,2,4,5-Tetrachlorobenzene	9.93	216	147405	26.49	ng	99
40) Hexachlorocyclopentadiene	9.92	237	50451	16.91	ng	96

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	10.08	196	101700	27.60	ng	97
43) 2,4,5-Trichlorophenol	10.12	196	108139	27.44	ng	# 94
45) 1,1'-Biphenyl	10.31	154	385735	26.25	ng	97
46) 2-Chloronaphthalene	10.33	162	314545	27.30	ng	97
47) 2-Nitroaniline	10.45	65	90763	32.00	ng	89
48) Acenaphthylene	10.84	152	519137	28.46	ng	99
49) Dimethylphthalate	10.69	163	408189	29.94	ng	98
50) 2,6-Dinitrotoluene	10.77	165	69773	25.52	ng	# 49
51) Acenaphthene	11.05	154	287360	26.40	ng	99
52) 3-Nitroaniline	10.97	138	85023	26.97	ng	90
53) 2,4-Dinitrophenol	11.10	184	7623	9.16	ng	# 79
54) Dibenzofuran	11.27	168	440058	26.18	ng	98
55) 4-Nitrophenol	11.18	139	138239	54.53	ng	89
56) 2,4-Dinitrotoluene	11.26	165	96491	26.12	ng	# 84
57) Fluorene	11.69	166	369199	27.48	ng	99
58) 2,3,4,6-Tetrachlorophenol	11.42	232	87153	27.51	ng	99
59) Diethylphthalate	11.57	149	361057	26.86	ng	96
60) 4-Chlorophenyl-phenylether	11.70	204	162900	25.16	ng	# 90
61) 4-Nitroaniline	11.72	138	85563	28.32	ng	90
62) Azobenzene	11.90	77	349892	27.44	ng	94
64) 4,6-Dinitro-2-methylphenol	11.77	198	9515	9.61	ng	74
65) n-Nitrosodiphenylamine	11.85	169	310111	25.07	ng	99
66) 4-Bromophenyl-phenylether	12.31	248	100619	23.05	ng	# 90
67) Hexachlorobenzene	12.37	284	108195	23.09	ng	# 86
68) Atrazine	12.52	200	98824	24.57	ng	99
69) Pentachlorophenol	12.62	266	118780	48.23	ng	99
70) Phenanthrene	12.89	178	663112	30.69	ng	99
71) Anthracene	12.95	178	561866	26.20	ng	99
72) Carbazole	13.15	167	503229	24.68	ng	99
73) Di-n-butylphthalate	13.61	149	601200	25.11	ng	98
74) Fluoranthene	14.37	202	981406	41.58	ng	96
76) Benzidine	14.54	184	446850	31.87	ng	98
77) Pyrene	14.65	202	1034979	40.77	ng	99
79) Butylbenzylphthalate	15.47	149	268199	25.68	ng	92
80) Benzo(a)anthracene	16.14	228	888836	36.55	ng	99
81) 3,3'-Dichlorobenzidine	16.12	252	200926	22.55	ng	# 95
82) Chrysene	16.19	228	721783	31.60	ng	98
83) Bis(2-ethylhexyl)phthalate	16.20	149	398225	23.78	ng	# 95
84) Di-n-octyl phthalate	17.00	149	717042	25.64	ng	99
85) Indeno(1,2,3-cd)pyrene	19.38	276	723936m	27.80	ng	
87) Benzo(b)fluoranthene	17.46	252	991105m	41.49	ng	
88) Benzo(k)fluoranthene	17.49	252	497296m	21.24	ng	
89) Benzo(a)pyrene	17.85	252	735680	32.33	ng	# 98
90) Dibenzo(a,h)anthracene	19.40	278	490317	22.60	ng	98
91) Benzo(g,h,i)perylene	19.81	276	585387	26.73	ng	96

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

