

Data Path : Z:\HPCHEM1\BNA F\DATA\BF021215\
 Data File : BF077300.D
 Acq On : 13 Feb 2015 00:36
 Operator : TP/IZ
 Sample : G1253-05MSD
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-SLDG-WT-70MSD

Manual Integrations
 APPROVED

mohammad
 2/13/2015 5:40:43 PM

Quant Time: Feb 13 12:05:21 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF021215.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Feb 13 03:50:26 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.50	152	63108	20.00	ng	0.01
21) Naphthalene-d8	10.41	136	264470	20.00	ng	0.00
38) Acenaphthene-d10	14.53	164	130591	20.00	ng	0.01
63) Phenanthrene-d10	17.44	188	254600	20.00	ng	0.00
75) Chrysene-d12	20.96	240	247018	20.00	ng	0.00
86) Perylene-d12	22.57	264	220273	20.00	ng	-0.07

System Monitoring Compounds

5) 2-Fluorophenol	4.51	112	205048	51.43	ng	0.02
7) Phenol-d6	6.99	99	247506	49.55	ng	0.06
23) Nitrobenzene-d5	8.82	82	185139	39.99	ng	0.00
41) 2,4,6-Tribromophenol	16.31	330	72843	68.40	ng	0.00
44) 2-Fluorobiphenyl	13.07	172	369770	42.62	ng	0.00
78) Terphenyl-d14	19.69	244	420278	38.22	ng	-0.01

Target Compounds

						Qvalue
2) 1,4-Dioxane	1.05	88	18145	11.92	ng	92
3) Pyridine	1.47	79	44798	11.10	ng	97
4) n-Nitrosodimethylamine	1.41	42	37270	10.72	ng	# 97
6) Aniline	6.80	93	43185	6.54	ng	95
8) 2-Chlorophenol	7.05	128	79628	17.39	ng	91
10) Phenol	7.04	94	84458m	16.45	ng	
11) bis(2-Chloroethyl)ether	7.05	93	78173	19.11	ng	90
12) 1,3-Dichlorobenzene	7.34	146	86276	17.04	ng	97
13) 1,4-Dichlorobenzene	7.54	146	87040	16.67	ng	95
14) 1,2-Dichlorobenzene	7.85	146	84841	17.15	ng	96
15) Benzyl Alcohol	7.97	79	115658	29.23	ng	92
16) 2,2'-oxybis(1-Chloropropan	8.30	45	99383	17.33	ng	92
17) 2-Methylphenol	8.33	107	62928	17.77	ng	94
18) Hexachloroethane	8.59	117	31134	16.42	ng	97
19) n-Nitroso-di-n-propylamine	8.60	70	65916	18.77	ng	99
20) 3+4-Methylphenols	8.73	107	81027	17.67	ng	96
22) Acetophenone	8.51	105	126202	19.87	ng	96
24) Nitrobenzene	8.85	77	83773	18.25	ng	100
25) Isophorone	9.47	82	164554	19.72	ng	# 94
26) 2-Nitrophenol	9.60	139	41074	18.93	ng	97
27) 2,4-Dimethylphenol	9.90	122	74710	19.92	ng	97
28) bis(2-Chloroethoxy)methane	10.10	93	99718	19.63	ng	100
29) 2,4-Dichlorophenol	10.22	162	65883m	19.22	ng	
30) 1,2,4-Trichlorobenzene	10.33	180	68922	17.79	ng	99
31) Naphthalene	10.45	128	294079	22.53	ng	99
32) Benzoic acid	10.58	122	43995m	63.51	ng	
33) 4-Chloroaniline	10.74	127	23456m	4.45	ng	
34) Hexachlorobutadiene	10.83	225	40632	18.46	ng	97
35) Caprolactam	11.54	113	2902	10.32	ng	# 64
36) 4-Chloro-3-methylphenol	12.04	107	72665	18.95	ng	93
37) 2-Methylnaphthalene	12.11	142	176545	18.04	ng	98
39) 1,2,4,5-Tetrachlorobenzene	12.51	216	67187	20.44	ng	# 100
40) Hexachlorocyclopentadiene	12.50	237	62296	42.74	ng	93
42) 2,4,6-Trichlorophenol	12.88	196	46058	21.55	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
43) 2,4,5-Trichlorophenol	12.99	196	49429	23.51	nq	# 92
45) 1,1'-Biphenyl	13.27	154	197745	19.65	nq	98
46) 2-Chloronaphthalene	13.24	162	160998	20.03	nq	98
47) 2-Nitroaniline	13.60	65	54087	23.03	nq	97
48) Acenaphthylene	14.18	152	263089	20.75	nq	99
49) Dimethylphthalate	14.16	163	200705	21.25	nq	98
50) 2,6-Dinitrotoluene	14.25	165	40208	22.02	nq	98
51) Acenaphthene	14.60	154	144105	19.62	nq	99
52) 3-Nitroaniline	14.60	138	11180	5.01	nq	92
53) 2,4-Dinitrophenol	14.89	184	32600	56.85	nq	# 71
54) Dibenzofuran	15.03	168	231630	20.40	nq	97
55) 4-Nitrophenol	15.24	139	48906	44.99	nq	99
56) 2,4-Dinitrotoluene	15.19	165	58056	22.41	nq	# 93
57) Fluorene	15.80	166	193945	20.86	nq	97
58) 2,3,4,6-Tetrachlorophenol	15.40	232	35168	24.60	nq	# 100
59) Diethylphthalate	15.83	149	210986	21.45	nq	99
60) 4-Chlorophenyl-phenylether	15.92	204	81293	21.03	nq	96
61) 4-Nitroaniline	15.99	138	32133	15.25	nq	95
62) Azobenzene	16.21	77	199672m	21.28	nq	
64) 4,6-Dinitro-2-methylphenol	16.07	198	26281	22.28	nq	100
65) n-Nitrosodiphenylamine	16.18	169	163440	18.00	nq	98
66) 4-Bromophenyl-phenylether	16.80	248	47464	17.84	nq	96
67) Hexachlorobenzene	16.81	284	48594	17.94	nq	92
68) Atrazine	17.20	200	53187	18.49	nq	88
69) Pentachlorophenol	17.20	266	41756	53.34	nq	95
70) Phenanthrene	17.47	178	281931	18.36	nq	99
71) Anthracene	17.55	178	284518	18.47	nq	99
72) Carbazole	17.85	167	278048	18.89	nq	99
73) Di-n-butylphthalate	18.45	149	427013	22.80	nq	100
74) Fluoranthene	19.13	202	295718	18.70	nq	99
76) Benzidine	19.38	184	77296	7.72	nq	99
77) Pyrene	19.41	202	316092	19.19	nq	99
79) Butylbenzylphthalate	20.36	149	219490	26.31	nq	97
80) Benzo(a)anthracene	20.95	228	305896	20.12	nq	99
81) 3,3'-Dichlorobenzidine	20.96	252	51445	9.65	nq	97
82) Chrysene	20.98	228	250095	18.15	nq	99
83) Bis(2-ethylhexyl)phthalate	21.11	149	259250	20.89	nq	97
84) Di-n-octyl phthalate	21.86	149	443744	20.79	nq	100
85) Indeno(1,2,3-cd)pyrene	23.61	276	275790m	20.59	nq	
87) Benzo(b)fluoranthene	22.18	252	275892	17.59	nq	99
88) Benzo(k)fluoranthene	22.20	252	255611m	21.57	nq	
89) Benzo(a)pyrene	22.51	252	252857	19.36	nq	# 97
90) Dibenzo(a,h)anthracene	23.63	278	231862m	19.99	nq	
91) Benzo(g,h,i)perylene	23.87	276	229940m	20.03	nq	

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

