

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF062515\
 Data File : BF080171.D
 Acq On : 25 Jun 2015 17:42
 Operator : TP/IZ
 Sample : G2748-01MSD
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC062315MSD

Manual Integrations
 APPROVED

apatel
 6/26/2015 2:05:17 PM

Quant Time: Jun 26 02:00:23 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF061715.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 26 01:01:50 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.29	152	28450	20.00	ng	0.00
21) Naphthalene-d8	8.58	136	113763	20.00	ng	0.00
38) Acenaphthene-d10	10.36	164	61128	20.00	ng	0.00
63) Phenanthrene-d10	11.86	188	120947	20.00	ng	0.00
75) Chrysene-d12	14.88	240	125778	20.00	ng	0.06
86) Perylene-d12	16.78	264	119318	20.00	ng	0.07

System Monitoring Compounds

5) 2-Fluorophenol	5.91	112	256868	159.82	ng	0.01
7) Phenol-d6	6.90	99	324407	151.68	ng	0.00
23) Nitrobenzene-d5	7.85	82	233884	119.69	ng	0.00
41) 2,4,6-Tribromophenol	11.16	330	102124	170.85	ng	0.00
44) 2-Fluorobiphenyl	9.66	172	427865	99.82	ng	-0.01
78) Terphenyl-d14	13.58	244	536382	99.60	ng	0.03

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.17	88	32070	41.98	ng	83
3) Pyridine	4.00	79	74065	36.46	ng	# 82
4) n-Nitrosodimethylamine	3.89	42	52187	54.05	ng	98
6) Aniline	6.95	93	86338	29.34	ng	92
8) 2-Chlorophenol	7.08	128	105729	56.92	ng	84
9) Benzaldehyde	6.85	77	40842	33.08	ng	92
10) Phenol	6.92	94	126942	54.39	ng	84
11) bis(2-Chloroethyl)ether	7.02	93	108259	58.46	ng	86
12) 1,3-Dichlorobenzene	7.24	146	112024m	50.20	ng	
13) 1,4-Dichlorobenzene	7.32	146	126814m	59.76	ng	
14) 1,2-Dichlorobenzene	7.48	146	108662	52.62	ng	98
15) Benzyl Alcohol	7.42	79	97678	55.86	ng	# 74
16) 2,2'-oxybis(1-Chloropropan	7.57	45	176828	64.33	ng	85
17) 2-Methylphenol	7.53	107	86914	54.78	ng	# 89
18) Hexachloroethane	7.82	117	41208	58.32	ng	# 85
19) n-Nitroso-di-n-propylamine	7.69	70	85193	57.37	ng	# 77
20) 3+4-Methylphenols	7.68	107	118921	58.08	ng	91
22) Acetophenone	7.69	105	154403	58.24	ng	# 81
24) Nitrobenzene	7.88	77	116619	57.82	ng	# 77
25) Isophorone	8.10	82	212534	54.04	ng	# 83
26) 2-Nitrophenol	8.20	139	58995	69.88	ng	# 77
27) 2,4-Dimethylphenol	8.22	122	108795	61.80	ng	# 85
28) bis(2-Chloroethoxy)methane	8.31	93	128820	55.75	ng	# 90
29) 2,4-Dichlorophenol	8.44	162	103277	68.50	ng	97
30) 1,2,4-Trichlorobenzene	8.53	180	106551	62.23	ng	100
31) Naphthalene	8.61	128	318756m	53.59	ng	
32) Benzoic acid	8.29	122	58952	55.39	ng	# 71
33) 4-Chloroaniline	8.64	127	47660	18.72	ng	# 82
34) Hexachlorobutadiene	8.73	225	57471	54.51	ng	98
35) Caprolactam	8.98	113	24906	50.90	ng	94
36) 4-Chloro-3-methylphenol	9.12	107	97531	57.35	ng	95
37) 2-Methylnaphthalene	9.30	142	243773	63.36	ng	# 92
39) 1,2,4,5-Tetrachlorobenzene	9.48	216	107000	60.15	ng	99
40) Hexachlorocyclopentadiene	9.46	237	105071	112.07	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.58	196	76978	63.60	ng	100
43) 2,4,5-Trichlorophenol	9.61	196	73598	52.78	ng	# 89
45) 1,1'-Biphenyl	9.77	154	262790	52.84	ng	94
46) 2-Chloronaphthalene	9.80	162	223903	55.49	ng	94
47) 2-Nitroaniline	9.88	65	65086	58.75	ng	# 61
48) Acenaphthylene	10.22	152	386494	57.38	ng	97
49) Dimethylphthalate	10.06	163	259845	56.54	ng	# 98
50) 2,6-Dinitrotoluene	10.12	165	52049	56.49	ng	# 44
51) Acenaphthene	10.39	154	241952	58.72	ng	90
52) 3-Nitroaniline	10.29	138	30739	28.92	ng	# 45
53) 2,4-Dinitrophenol	10.40	184	60814	156.18	ng	# 37
54) Dibenzofuran	10.56	168	314258	53.75	ng	# 88
55) 4-Nitrophenol	10.44	139	86766	102.12	ng	# 44
56) 2,4-Dinitrotoluene	10.53	165	79380	67.90	ng	# 82
57) Fluorene	10.92	166	259834	56.01	ng	92
58) 2,3,4,6-Tetrachlorophenol	10.68	232	63493	53.93	ng	98
59) Diethylphthalate	10.76	149	248298	53.66	ng	95
60) 4-Chlorophenyl-phenylether	10.89	204	124140	55.68	ng	# 86
61) 4-Nitroaniline	10.90	138	58100	52.92	ng	# 47
62) Azobenzene	11.05	77	244822	51.97	ng	86
64) 4,6-Dinitro-2-methylphenol	10.94	198	35859	55.56	ng	85
65) n-Nitrosodiphenylamine	11.01	169	222346	56.46	ng	92
66) 4-Bromophenyl-phenylether	11.40	248	76946	59.83	ng	# 88
67) Hexachlorobenzene	11.48	284	83220	58.23	ng	# 80
68) Atrazine	11.53	200	70456	56.62	ng	81
69) Pentachlorophenol	11.66	266	88357	109.09	ng	97
70) Phenanthrene	11.90	178	392595	53.62	ng	98
71) Anthracene	11.94	178	404970	57.44	ng	97
72) Carbazole	12.09	167	353969	52.58	ng	95
73) Di-n-butylphthalate	12.44	149	422532	54.33	ng	# 99
74) Fluoranthene	13.17	202	436726	56.00	ng	91
76) Benzidine	13.29	184	120850	34.37	ng	97
77) Pyrene	13.43	202	455297	58.79	ng	96
79) Butylbenzylphthalate	14.14	149	184017	59.17	ng	92
80) Benzo(a)anthracene	14.87	228	426441	55.65	ng	97
81) 3,3'-Dichlorobenzidine	14.81	252	101280	38.32	ng	# 94
82) Chrysene	14.92	228	387262	54.81	ng	95
83) Bis(2-ethylhexyl)phthalate	14.85	149	268583	54.65	ng	# 94
84) Di-n-octyl phthalate	15.67	149	452503	53.73	ng	94
85) Indeno(1,2,3-cd)pyrene	18.62	276	466750	52.38	ng	# 89
87) Benzo(b)fluoranthene	16.24	252	405483	56.13	ng	# 87
88) Benzo(k)fluoranthene	16.28	252	417609	56.77	ng	# 87
89) Benzo(a)pyrene	16.70	252	403628	58.83	ng	# 86
90) Dibenzo(a,h)anthracene	18.64	278	389666	55.49	ng	# 82
91) Benzo(g,h,i)perylene	19.17	276	397748	56.10	ng	# 79

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

