

Data Path : Z:\HPCHEM1\BNA F\DATA\BF052016\
 Data File : BF087234.D
 Acq On : 20 May 2016 19:57
 Operator : UM/SJ
 Sample : PB90764BS
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB90764BS

Quant Time: May 23 16:40:42 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF051716.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri May 20 23:58:09 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.56	152	144161	20.00	ng	0.00
21) Naphthalene-d8	7.85	136	609198	20.00	ng	0.01
38) Acenaphthene-d10	9.59	164	299225	20.00	ng	0.00
63) Phenanthrene-d10	11.07	188	562724	20.00	ng	0.01
75) Chrysene-d12	13.70	240	375739	20.00	ng	0.01
86) Perylene-d12	15.04	264	306581	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.15	112	749253	87.36	ng	0.01
7) Phenol-d6	6.23	99	919896	79.97	ng	0.00
23) Nitrobenzene-d5	7.14	82	1018724	83.34	ng	0.01
41) 2,4,6-Tribromophenol	10.39	330	253384	76.63	ng	0.00
44) 2-Fluorobiphenyl	8.92	172	1434519	79.40	ng	0.00
78) Terphenyl-d14	12.65	244	1410783	84.93	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.06	88	146923	33.30	ng	# 59
3) Pyridine	2.67	79	368471	34.53	ng	# 69
4) n-Nitrosodimethylamine	2.65	42	316194	41.81	ng	# 50
6) Aniline	6.23	93	256502	17.55	ng	# 70
8) 2-Chlorophenol	6.34	128	447555	42.89	ng	# 78
9) Benzaldehyde	6.10	77	105065	14.07	ng	90
10) Phenol	6.24	94	542506	37.52	ng	# 64
11) bis(2-Chloroethyl)ether	6.31	93	457357	42.12	ng	93
12) 1,3-Dichlorobenzene	6.49	146	439884m	39.66	ng	
13) 1,4-Dichlorobenzene	6.57	146	452364m	39.86	ng	
14) 1,2-Dichlorobenzene	6.73	146	401946	40.45	ng	97
15) Benzyl Alcohol	6.73	79	422668	49.36	ng	91
16) 2,2'-oxybis(1-Chloropropan	6.86	45	866698	38.83	ng	53
17) 2-Methylphenol	6.86	107	379651	44.30	ng	# 80
18) Hexachloroethane	7.06	117	178270	41.09	ng	95
19) n-Nitroso-di-n-propylamine	7.14	70	147322	16.85	ng	# 78
20) 3+4-Methylphenols	7.02	107	518800	45.89	ng	# 59
22) Acetophenone	6.98	105	640908	42.99	ng	# 97
24) Nitrobenzene	7.15	77	497402	37.27	ng	94
25) Isophorone	7.14	82	1018724	45.53	ng	# 68
26) 2-Nitrophenol	7.47	139	245429	44.39	ng	# 76
27) 2,4-Dimethylphenol	7.53	122	391986	41.55	ng	86
28) bis(2-Chloroethoxy)methane	7.61	93	562442	45.79	ng	# 98
29) 2,4-Dichlorophenol	7.72	162	394798	47.48	ng	96
30) 1,2,4-Trichlorobenzene	7.79	180	402968	43.68	ng	98
31) Naphthalene	7.86	128	1264635	42.49	ng	99
32) Benzoic acid	7.71	122	220827m	39.27	ng	
33) 4-Chloroaniline	7.94	127	121273	9.80	ng	# 94
34) Hexachlorobutadiene	7.99	225	237463	45.35	ng	97
35) Caprolactam	8.32	113	118781m	42.90	ng	
36) 4-Chloro-3-methylphenol	8.44	107	418755	41.75	ng	88
37) 2-Methylnaphthalene	8.56	142	826218m	43.40	ng	
39) 1,2,4,5-Tetrachlorobenzene	8.73	216	386031	44.17	ng	98
40) Hexachlorocyclopentadiene	8.71	237	286127	84.31	ng	96

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42) 2,4,6-Trichlorophenol	8.86	196	248176m	43.79	nq	
43) 2,4,5-Trichlorophenol	8.90	196	254291	43.19	nq	# 88
45) 1,1'-Biphenyl	9.03	154	1145558	46.17	nq	99
46) 2-Chloronaphthalene	9.05	162	841449	44.18	nq	99
47) 2-Nitroaniline	9.16	65	323069m	44.26	nq	
48) Acenaphthylene	9.46	152	1244182	42.78	nq	99
49) Dimethylphthalate	9.34	163	998781	43.76	nq	100
50) 2,6-Dinitrotoluene	9.59	165	38866	7.84	nq	# 32
51) Acenaphthene	9.62	154	733474	40.37	nq	99
52) 3-Nitroaniline	9.58	138	79579m	14.83	nq	
53) 2,4-Dinitrophenol	9.69	184	198525	70.50	nq	# 70
54) Dibenzofuran	9.80	168	1151822	49.02	nq	98
55) 4-Nitrophenol	9.77	139	183434m	57.52	nq	
56) 2,4-Dinitrotoluene	9.59	165	38866	6.12	nq	# 26
57) Fluorene	10.14	166	807834	42.79	nq	99
58) 2,3,4,6-Tetrachlorophenol	9.93	232	213428	46.54	nq	97
59) Diethylphthalate	10.03	149	889815	40.83	nq	100
60) 4-Chlorophenyl-phenylether	10.14	204	402318	44.96	nq	94
61) 4-Nitroaniline	10.19	138	190494	35.96	nq	# 67
62) Azobenzene	10.30	77	1206942	48.07	nq	87
64) 4,6-Dinitro-2-methylphenol	10.23	198	150430	40.59	nq	90
65) n-Nitrosodiphenylamine	10.26	169	733550	41.42	nq	99
66) 4-Bromophenyl-phenylether	10.39	248	17297	2.99	nq	# 1
67) Hexachlorobenzene	10.68	284	259012	38.61	nq	# 73
68) Atrazine	10.80	200	247342	42.82	nq	93
69) Pentachlorophenol	10.89	266	202338	80.28	nq	97
70) Phenanthrene	11.10	178	1290952	42.30	nq	99
71) Anthracene	11.14	178	1261652	40.87	nq	100
72) Carbazole	11.31	167	1171942	41.56	nq	98
73) Di-n-butylphthalate	11.65	149	1482802	45.78	nq	# 98
74) Fluoranthene	12.27	202	1192082	38.98	nq	97
76) Benzidine	12.65	184	24727m	2.70	nq	
77) Pyrene	12.50	202	1331723	49.06	nq	99
79) Butylbenzylphthalate	13.13	149	633719	55.30	nq	# 86
80) Benzo(a)anthracene	13.68	228	968807m	44.59	nq	
81) 3,3'-Dichlorobenzidine	13.66	252	169012	12.23	nq	# 95
82) Chrysene	13.73	228	1053725	50.13	nq	99
83) Bis(2-ethylhexyl)phthalate	13.69	149	773029	55.43	nq	98
84) Di-n-octyl phthalate	14.30	149	1290131	53.70	nq	# 85
85) Indeno(1,2,3-cd)pyrene	16.27	276	741613	40.20	nq	96
87) Benzo(b)fluoranthene	14.69	252	827634m	40.69	nq	
88) Benzo(k)fluoranthene	14.71	252	731122m	48.28	nq	
89) Benzo(a)pyrene	14.99	252	740831	43.58	nq	98
90) Dibenzo(a,h)anthracene	16.27	278	624292	43.61	nq	96
91) Benzo(g,h,i)perylene	16.63	276	625954	43.30	nq	96

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

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