

Data Path : Z:\HPCHEM1\BNA F\DATA\BF050815\
 Data File : BF079045.D
 Acq On : 8 May 2015 18:45
 Operator : TP/IZ
 Sample : PB83253BS
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB83253BS

Manual Integrations
 APPROVED

apatel
 5/12/2015 3:43:56 PM

Quant Time: May 08 23:26:51 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF043015.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri May 08 22:44:53 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.32	152	78035	20.00	ng	0.00
21) Naphthalene-d8	8.90	136	323094	20.00	ng	0.00
38) Acenaphthene-d10	11.07	164	158472	20.00	ng	0.00
63) Phenanthrene-d10	12.93	188	298627	20.00	ng	0.00
75) Chrysene-d12	16.22	240	292372	20.00	ng	-0.03
86) Perylene-d12	17.91	264	294113	20.00	ng	-0.17

System Monitoring Compounds

5) 2-Fluorophenol	5.62	112	368809	81.36	ng	0.00
7) Phenol-d6	6.88	99	468143	78.12	ng	-0.01
23) Nitrobenzene-d5	8.01	82	449459	79.95	ng	0.00
41) 2,4,6-Tribromophenol	12.06	330	130609	76.18	ng	0.00
44) 2-Fluorobiphenyl	10.24	172	912874	80.26	ng	-0.01
78) Terphenyl-d14	14.91	244	1046997	85.74	ng	-0.01

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.31	88	62655	31.79	ng	# 33
3) Pyridine	3.05	79	201732	34.97	ng	95
4) n-Nitrosodimethylamine	2.97	42	85145m	34.45	ng	
6) Aniline	6.91	93	168662	19.94	ng	99
8) 2-Chlorophenol	7.05	128	188064	36.58	ng	97
9) Benzaldehyde	6.78	77	111938	27.73	ng	95
10) Phenol	6.90	94	256624	36.92	ng	92
11) bis(2-Chloroethyl)ether	7.00	93	175857	34.51	ng	96
12) 1,3-Dichlorobenzene	7.24	146	200269	34.65	ng	93
13) 1,4-Dichlorobenzene	7.35	146	206412	34.71	ng	96
14) 1,2-Dichlorobenzene	7.53	146	193610	34.97	ng	97
15) Benzyl Alcohol	7.50	79	157008	35.30	ng	96
16) 2,2'-oxybis(1-Chloropropan	7.68	45	277121	35.54	ng	99
17) 2-Methylphenol	7.64	107	155139	37.49	ng	97
18) Hexachloroethane	7.94	117	72109	35.20	ng	# 88
19) n-Nitroso-di-n-propylamine	7.84	70	134156	33.15	ng	# 83
20) 3+4-Methylphenols	7.84	107	203675	36.94	ng	# 49
22) Acetophenone	7.83	105	268439	36.77	ng	96
24) Nitrobenzene	8.03	77	189740	34.05	ng	98
25) Isophorone	8.33	82	344399	33.05	ng	99
26) 2-Nitrophenol	8.43	139	88989	38.58	ng	96
27) 2,4-Dimethylphenol	8.49	122	181527	39.33	ng	97
28) bis(2-Chloroethoxy)methane	8.62	93	223530	34.79	ng	99
29) 2,4-Dichlorophenol	8.73	162	159329	38.82	ng	95
30) 1,2,4-Trichlorobenzene	8.83	180	158909	35.25	ng	98
31) Naphthalene	8.92	128	545537	35.44	ng	99
32) Benzoic acid	8.59	122	93242	34.32	ng	96
33) 4-Chloroaniline	8.99	127	118819	18.24	ng	97
34) Hexachlorobutadiene	9.07	225	85778	33.04	ng	97
35) Caprolactam	9.43	113	48896	36.84	ng	92
36) 4-Chloro-3-methylphenol	9.60	107	162717	37.75	ng	96
37) 2-Methylnaphthalene	9.78	142	370246	34.71	ng	98
39) 1,2,4,5-Tetrachlorobenzene	9.99	216	145419	32.58	ng	# 100
40) Hexachlorocyclopentadiene	9.98	237	150456	67.04	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	10.14	196	108971	35.51	nq	97
43) 2,4,5-Trichlorophenol	10.18	196	117007	37.72	nq	# 92
45) 1,1'-Biphenyl	10.37	154	442176	35.10	nq	98
46) 2-Chloronaphthalene	10.39	162	352599	35.88	nq	96
47) 2-Nitroaniline	10.51	65	100220	35.20	nq	99
48) Acenaphthylene	10.90	152	575101	35.64	nq	99
49) Dimethylphthalate	10.75	163	410344	35.28	nq	100
50) 2,6-Dinitrotoluene	10.82	165	78450	34.18	nq	98
51) Acenaphthene	11.12	154	321909	33.46	nq	99
52) 3-Nitroaniline	11.03	138	70938	24.74	nq	# 96
53) 2,4-Dinitrophenol	11.17	184	65445	72.11	nq	# 84
54) Dibenzofuran	11.34	168	493072	35.48	nq	94
55) 4-Nitrophenol	11.25	139	170791	75.26	nq	92
56) 2,4-Dinitrotoluene	11.33	165	115008	37.90	nq	# 76
57) Fluorene	11.76	166	404769	35.48	nq	99
58) 2,3,4,6-Tetrachlorophenol	11.49	232	91864	36.24	nq	# 100
59) Diethylphthalate	11.63	149	396921	34.45	nq	99
60) 4-Chlorophenyl-phenylether	11.76	204	171381	33.87	nq	# 87
61) 4-Nitroaniline	11.78	138	95569	33.32	nq	93
62) Azobenzene	11.95	77	399133	35.15	nq	93
64) 4,6-Dinitro-2-methylphenol	11.83	198	46147	39.84	nq	# 64
65) n-Nitrosodiphenylamine	11.91	169	338855	34.07	nq	98
66) 4-Bromophenyl-phenylether	12.38	248	110647	35.55	nq	# 84
67) Hexachlorobenzene	12.43	284	125594	35.30	nq	# 87
68) Atrazine	12.58	200	107766	37.26	nq	100
69) Pentachlorophenol	12.69	266	116351	57.96	nq	98
70) Phenanthrene	12.95	178	570662	34.16	nq	99
71) Anthracene	13.02	178	599225	36.56	nq	99
72) Carbazole	13.22	167	580288	37.15	nq	98
73) Di-n-butylphthalate	13.67	149	677352	36.16	nq	# 98
74) Fluoranthene	14.43	202	630350	36.21	nq	94
76) Benzidine	14.61	184	250058	31.73	nq	99
77) Pyrene	14.71	202	637184	37.82	nq	99
79) Butylbenzylphthalate	15.53	149	289831	39.35	nq	95
80) Benzo(a)anthracene	16.21	228	573942	35.85	nq	99
81) 3,3'-Dichlorobenzidine	16.17	252	150073	26.31	nq	# 97
82) Chrysene	16.24	228	532852	34.60	nq	99
83) Bis(2-ethylhexyl)phthalate	16.25	149	437166	39.19	nq	99
84) Di-n-octyl phthalate	17.03	149	733975	37.36	nq	# 100
85) Indeno(1,2,3-cd)pyrene	19.39	276	677605	34.54	nq	# 100
87) Benzo(b)fluoranthene	17.47	252	593343	36.86	nq	# 96
88) Benzo(k)fluoranthene	17.50	252	539888m	34.64	nq	
89) Benzo(a)pyrene	17.85	252	552203	37.25	nq	99
90) Dibenzo(a,h)anthracene	19.42	278	543385	34.49	nq	99
91) Benzo(g,h,i)perylene	19.83	276	545142	34.53	nq	98

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

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