

Data Path : Z:\HPCHEM1\BNA F\DATA\BF040915\
 Data File : BF078359.D
 Acq On : 9 Apr 2015 14:02
 Operator : TP/IZ
 Sample : PB82625BS
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleID :
 PB82625BS

Manual Integrations
 APPROVED

apatel
 4/10/2015 5:12:51 PM

Quant Time: Apr 10 10:58:31 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF040115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 10 00:58:40 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.89	152	27783	20.00	ng	0.00
21) Naphthalene-d8	8.46	136	116709	20.00	ng	0.00
38) Acenaphthene-d10	10.62	164	60282	20.00	ng	0.00
63) Phenanthrene-d10	12.45	188	110663	20.00	ng	0.00
75) Chrysene-d12	15.77	240	110406m	20.00	ng	0.05
86) Perylene-d12	17.63	264	102013	20.00	ng	0.15

System Monitoring Compounds

5) 2-Fluorophenol	5.15	112	158558	94.10	ng	0.00
7) Phenol-d6	6.49	99	210176	99.53	ng	0.00
23) Nitrobenzene-d5	7.58	82	152359	79.13	ng	0.00
41) 2,4,6-Tribromophenol	11.61	330	51260	102.53	ng	0.00
44) 2-Fluorobiphenyl	9.81	172	335142	79.47	ng	0.00
78) Terphenyl-d14	14.44	244	372936	76.33	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	1.72	88	25336	33.25	ng	92
3) Pyridine	2.32	79	91255	45.72	ng	96
4) n-Nitrosodimethylamine	2.26	42	34945m	45.48	ng	
6) Aniline	6.48	93	71864	24.00	ng	93
8) 2-Chlorophenol	6.62	128	94287	47.32	ng	91
9) Benzaldehyde	6.33	77	30890	22.07	ng	98
10) Phenol	6.50	94	119956	47.41	ng	96
11) bis(2-Chloroethyl)ether	6.59	93	85384	46.92	ng	97
12) 1,3-Dichlorobenzene	6.81	146	97795	44.68	ng	# 91
13) 1,4-Dichlorobenzene	6.91	146	98285	44.61	ng	94
14) 1,2-Dichlorobenzene	7.09	146	90934	44.02	ng	92
15) Benzyl Alcohol	7.08	79	75941	51.17	ng	97
16) 2,2'-oxybis(1-Chloropropan	7.26	45	114580	47.21	ng	97
17) 2-Methylphenol	7.25	107	74931	48.44	ng	97
18) Hexachloroethane	7.50	117	35272	47.48	ng	# 87
19) n-Nitroso-di-n-propylamine	7.42	70	66323	47.60	ng	99
20) 3+4-Methylphenols	7.45	107	100624	48.76	ng	92
22) Acetophenone	7.40	105	131804	48.47	ng	# 90
24) Nitrobenzene	7.61	77	94186	46.79	ng	# 83
25) Isophorone	7.92	82	170984	46.79	ng	99
26) 2-Nitrophenol	8.01	139	46941	48.94	ng	97
27) 2,4-Dimethylphenol	8.09	122	84056	47.13	ng	95
28) bis(2-Chloroethoxy)methane	8.20	93	106237	45.34	ng	98
29) 2,4-Dichlorophenol	8.32	162	81163	50.02	ng	96
30) 1,2,4-Trichlorobenzene	8.40	180	80181	43.10	ng	# 95
31) Naphthalene	8.49	128	267834	44.09	ng	99
32) Benzoic acid	8.22	122	49174	57.34	ng	92
33) 4-Chloroaniline	8.58	127	70247	27.87	ng	99
34) Hexachlorobutadiene	8.65	225	46485	45.50	ng	98
35) Caprolactam	9.01	113	25785	53.23	ng	87
36) 4-Chloro-3-methylphenol	9.21	107	85807	52.41	ng	# 87
37) 2-Methylnaphthalene	9.34	142	184315	44.69	ng	98
39) 1,2,4,5-Tetrachlorobenzene	9.55	216	78016	41.77	ng	# 97
40) Hexachlorocyclopentadiene	9.54	237	76165	93.47	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.71	196	58471	49.64	nq	99
43) 2,4,5-Trichlorophenol	9.76	196	59775	48.57	nq	95
45) 1,1'-Biphenyl	9.93	154	220948	44.81	nq	97
46) 2-Chloronaphthalene	9.95	162	171763	44.27	nq	94
47) 2-Nitroaniline	10.09	65	49066	51.12	nq	# 82
48) Acenaphthylene	10.45	152	283575	45.26	nq	99
49) Dimethylphthalate	10.33	163	206532	45.60	nq	98
50) 2,6-Dinitrotoluene	10.40	165	43050	46.20	nq	# 46
51) Acenaphthene	10.66	154	161813	45.88	nq	98
52) 3-Nitroaniline	10.60	138	35249	33.27	nq	84
53) 2,4-Dinitrophenol	10.74	184	34135	89.17	nq	# 88
54) Dibenzofuran	10.88	168	254156	47.18	nq	97
55) 4-Nitrophenol	10.84	139	72274	92.21	nq	86
56) 2,4-Dinitrotoluene	10.89	165	57602	47.73	nq	# 55
57) Fluorene	11.30	166	206891	47.38	nq	100
58) 2,3,4,6-Tetrachlorophenol	11.05	232	45054	49.38	nq	98
59) Diethylphthalate	11.21	149	199213	45.62	nq	98
60) 4-Chlorophenyl-phenylether	11.32	204	95521	47.55	nq	92
61) 4-Nitroaniline	11.34	138	42892	40.05	nq	92
62) Azobenzene	11.52	77	212113	53.22	nq	85
64) 4,6-Dinitro-2-methylphenol	11.39	198	23968	44.41	nq	# 74
65) n-Nitrosodiphenylamine	11.47	169	177762	46.44	nq	99
66) 4-Bromophenyl-phenylether	11.92	248	54459	46.47	nq	97
67) Hexachlorobenzene	11.97	284	61196	47.65	nq	# 92
68) Atrazine	12.15	200	54607	49.25	nq	97
69) Pentachlorophenol	12.23	266	55513	95.78	nq	100
70) Phenanthrene	12.48	178	295669	46.19	nq	99
71) Anthracene	12.55	178	307748	48.79	nq	100
72) Carbazole	12.76	167	288269	48.76	nq	99
73) Di-n-butylphthalate	13.22	149	326174	48.71	nq	99
74) Fluoranthene	13.95	202	317837	47.08	nq	90
76) Benzidine	14.13	184	141184	33.21	nq	98
77) Pyrene	14.23	202	341458	49.27	nq	100
79) Butylbenzylphthalate	15.09	149	138669	52.49	nq	# 77
80) Benzo(a)anthracene	15.76	228	303660	46.23	nq	99
81) 3,3'-Dichlorobenzidine	15.75	252	66935	30.42	nq	# 97
82) Chrysene	15.80	228	289684	46.94	nq	99
83) Bis(2-ethylhexyl)phthalate	15.86	149	201177	48.56	nq	98
84) Di-n-octyl phthalate	16.72	149	336914m	49.53	nq	
85) Indeno(1,2,3-cd)pyrene	18.99	276	313074	44.70	nq	# 85
87) Benzo(b)fluoranthene	17.15	252	286508m	45.44	nq	
88) Benzo(k)fluoranthene	17.19	252	287345m	51.29	nq	
89) Benzo(a)pyrene	17.56	252	279956	50.88	nq	# 95
90) Dibenzo(a,h)anthracene	19.01	278	268504m	48.74	nq	
91) Benzo(g,h,i)perylene	19.36	276	267882m	48.50	nq	

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

