

Data Path : Z:\HPCHEM1\BNA F\DATA\BF040115\
 Data File : BF078172.D
 Acq On : 2 Apr 2015 00:49
 Operator : TP/IZ
 Sample : PB82482BS
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB82482BS

Manual Integrations
 APPROVED

apatel
 4/2/2015 5:51:23 PM

Quant Time: Apr 02 13:55:54 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF040115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 02 13:21:29 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.01	152	35628	20.00	ng	0.00
21) Naphthalene-d8	8.59	136	150058	20.00	ng	0.00
38) Acenaphthene-d10	10.76	164	75286	20.00	ng	0.00
63) Phenanthrene-d10	12.59	188	143283	20.00	ng	0.00
75) Chrysene-d12	15.87	240	144698	20.00	ng	0.00
86) Perylene-d12	17.56	264	138781	20.00	ng	-0.02

System Monitoring Compounds

5) 2-Fluorophenol	5.30	112	269816	124.86	ng	0.00
7) Phenol-d6	6.60	99	331208	122.30	ng	0.00
23) Nitrobenzene-d5	7.71	82	190949	77.13	ng	0.00
41) 2,4,6-Tribromophenol	11.74	330	84949	136.05	ng	0.00
44) 2-Fluorobiphenyl	9.94	172	434343	82.47	ng	0.00
78) Terphenyl-d14	14.59	244	504131	78.73	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	1.88	88	30409	31.12	ng	97
3) Pyridine	2.52	79	92878	36.29	ng	97
4) n-Nitrosodimethylamine	2.45	42	42488m	43.12	ng	
6) Aniline	6.61	93	61140	15.92	ng	# 39
8) 2-Chlorophenol	6.75	128	109843	42.99	ng	99
10) Phenol	6.62	94	132169	40.73	ng	99
11) bis(2-Chloroethyl)ether	6.72	93	102870	44.08	ng	99
12) 1,3-Dichlorobenzene	6.93	146	114170	40.68	ng	99
13) 1,4-Dichlorobenzene	7.04	146	115009	40.71	ng	100
14) 1,2-Dichlorobenzene	7.22	146	109596	41.37	ng	99
15) Benzyl Alcohol	7.21	79	81477	42.81	ng	99
16) 2,2'-oxybis(1-Chloropropan	7.39	45	133942	43.03	ng	100
17) 2-Methylphenol	7.37	107	87183	43.95	ng	99
18) Hexachloroethane	7.64	117	39608	41.57	ng	95
19) n-Nitroso-di-n-propylamine	7.55	70	74161	41.51	ng	98
20) 3+4-Methylphenols	7.56	107	114109	43.12	ng	97
22) Acetophenone	7.54	105	145527	41.62	ng	98
24) Nitrobenzene	7.73	77	103310	39.92	ng	# 79
25) Isophorone	8.04	82	197255	41.98	ng	99
26) 2-Nitrophenol	8.13	139	53637	43.49	ng	99
27) 2,4-Dimethylphenol	8.21	122	99022	43.18	ng	98
28) bis(2-Chloroethoxy)methane	8.33	93	131246	43.56	ng	99
29) 2,4-Dichlorophenol	8.44	162	89396	42.85	ng	99
30) 1,2,4-Trichlorobenzene	8.53	180	95116	39.76	ng	100
31) Naphthalene	8.62	128	312963	40.07	ng	99
32) Benzoic acid	8.34	122	39320	37.85	ng	97
33) 4-Chloroaniline	8.70	127	57162	17.64	ng	100
34) Hexachlorobutadiene	8.78	225	52580	40.03	ng	98
35) Caprolactam	9.13	113	26772	42.99	ng	99
36) 4-Chloro-3-methylphenol	9.32	107	92799	44.08	ng	97
37) 2-Methylnaphthalene	9.48	142	222132	41.89	ng	99
39) 1,2,4,5-Tetrachlorobenzene	9.69	216	95731	41.04	ng	98
40) Hexachlorocyclopentadiene	9.68	237	90803	89.53	ng	97
42) 2,4,6-Trichlorophenol	9.84	196	62383	42.40	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
43) 2,4,5-Trichlorophenol	9.89	196	65541	42.64	nq	# 85
45) 1,1'-Biphenyl	10.06	154	256582	41.67	nq	99
46) 2-Chloronaphthalene	10.08	162	202914	41.88	nq	99
47) 2-Nitroaniline	10.21	65	51835	43.24	nq	99
48) Acenaphthylene	10.59	152	329149	42.07	nq	100
49) Dimethylphthalate	10.45	163	233054	41.20	nq	99
50) 2,6-Dinitrotoluene	10.53	165	54256	46.62	nq	98
51) Acenaphthene	10.80	154	183218	41.59	nq	99
52) 3-Nitroaniline	10.73	138	32986	24.93	nq	99
53) 2,4-Dinitrophenol	10.86	184	37980	82.59	nq	97
54) Dibenzofuran	11.01	168	293311	43.59	nq	100
55) 4-Nitrophenol	10.97	139	80025	82.06	nq	92
56) 2,4-Dinitrotoluene	11.02	165	70192	46.57	nq	95
57) Fluorene	11.44	166	234805	43.05	nq	99
58) 2,3,4,6-Tetrachlorophenol	11.17	232	54479m	47.81	nq	
59) Diethylphthalate	11.33	149	234109	42.93	nq	99
60) 4-Chlorophenyl-phenylether	11.46	204	106130	42.30	nq	96
61) 4-Nitroaniline	11.48	138	54320	40.61	nq	99
62) Azobenzene	11.64	77	221603	44.52	nq	99
64) 4,6-Dinitro-2-methylphenol	11.52	198	26884	39.62	nq	95
65) n-Nitrosodiphenylamine	11.61	169	209908	42.35	nq	98
66) 4-Bromophenyl-phenylether	12.05	248	65454	43.13	nq	97
67) Hexachlorobenzene	12.11	284	71154	42.79	nq	99
68) Atrazine	12.28	200	65449	45.59	nq	98
69) Pentachlorophenol	12.36	266	60994	82.45	nq	98
70) Phenanthrene	12.63	178	361970	43.67	nq	99
71) Anthracene	12.68	178	354883	43.45	nq	98
72) Carbazole	12.90	167	338874	44.27	nq	99
73) Di-n-butylphthalate	13.36	149	381540	44.00	nq	99
74) Fluoranthene	14.09	202	373176	42.69	nq	98
76) Benzidine	14.28	184	137452	24.67	nq	99
77) Pyrene	14.37	202	404785	44.56	nq	99
79) Butylbenzylphthalate	15.21	149	161694	46.70	nq	98
80) Benzo(a)anthracene	15.86	228	362349	42.09	nq	99
81) 3,3'-Dichlorobenzidine	15.84	252	47974	16.64	nq	# 96
82) Chrysene	15.89	228	335598	41.49	nq	99
83) Bis(2-ethylhexyl)phthalate	15.93	149	246128	45.33	nq	# 96
84) Di-n-octyl phthalate	16.71	149	423273	47.47	nq	# 100
85) Indeno(1,2,3-cd)pyrene	18.87	276	391896	42.70	nq	# 87
87) Benzo(b)fluoranthene	17.13	252	348227	40.60	nq	99
88) Benzo(k)fluoranthene	17.16	252	357127m	46.85	nq	
89) Benzo(a)pyrene	17.49	252	338651	45.24	nq	# 95
90) Dibenzo(a,h)anthracene	18.90	278	323842	43.21	nq	97
91) Benzo(g,h,i)perylene	19.25	276	325316	43.29	nq	96

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

