

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF033015\
 Data File : BF078084.D
 Acq On : 30 Mar 2015 16:28
 Operator : TP/IZ
 Sample : PB82403BS
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB82403BS

Manual Integrations
 APPROVED

apatel
 3/31/2015 2:16:58 PM

Quant Time: Mar 31 01:48:39 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 31 01:02:18 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.06	152	33238	20.00	ng	0.00
21) Naphthalene-d8	8.64	136	137809	20.00	ng	0.00
38) Acenaphthene-d10	10.81	164	65363	20.00	ng	0.00
63) Phenanthrene-d10	12.65	188	133700	20.00	ng	0.00
75) Chrysene-d12	15.94	240	141352	20.00	ng	0.00
86) Perylene-d12	17.71	264	138614	20.00	ng	0.01

System Monitoring Compounds

5) 2-Fluorophenol	5.36	112	262233	137.71	ng	0.01
7) Phenol-d6	6.65	99	318117	135.22	ng	0.00
23) Nitrobenzene-d5	7.76	82	178752	85.03	ng	0.00
41) 2,4,6-Tribromophenol	11.79	330	80721	129.08	ng	0.00
44) 2-Fluorobiphenyl	9.98	172	400880	90.13	ng	0.00
78) Terphenyl-d14	14.65	244	478772	82.74	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	1.93	88	33107	38.29	ng	# 100
3) Pyridine	2.58	79	100903	43.44	ng	99
4) n-Nitrosodimethylamine	2.52	42	38663m	38.23	ng	
6) Aniline	6.65	93	66428	19.74	ng	# 72
8) 2-Chlorophenol	6.80	128	105097	46.37	ng	99
9) Benzaldehyde	6.51	77	3017	3.13	ng	97
10) Phenol	6.67	94	127837	45.97	ng	# 69
11) bis(2-Chloroethyl)ether	6.76	93	97068	46.60	ng	98
12) 1,3-Dichlorobenzene	6.98	146	107977	42.83	ng	# 96
13) 1,4-Dichlorobenzene	7.08	146	111806	42.68	ng	96
14) 1,2-Dichlorobenzene	7.26	146	106146	44.20	ng	96
15) Benzyl Alcohol	7.25	79	79408	43.24	ng	97
16) 2,2'-oxybis(1-Chloropropan	7.44	45	122601	43.67	ng	99
17) 2-Methylphenol	7.41	107	84176	45.62	ng	97
18) Hexachloroethane	7.68	117	36499	42.49	ng	# 90
19) n-Nitroso-di-n-propylamine	7.58	70	67533	41.49	ng	97
20) 3+4-Methylphenols	7.61	107	107096	44.23	ng	95
22) Acetophenone	7.57	105	140562	46.07	ng	# 98
24) Nitrobenzene	7.78	77	97478	43.04	ng	97
25) Isophorone	8.09	82	176178	41.50	ng	97
26) 2-Nitrophenol	8.18	139	47527	42.86	ng	97
27) 2,4-Dimethylphenol	8.26	122	87920	43.52	ng	97
28) bis(2-Chloroethoxy)methane	8.37	93	110692	42.95	ng	# 96
29) 2,4-Dichlorophenol	8.48	162	84292	43.78	ng	92
30) 1,2,4-Trichlorobenzene	8.57	180	86152	39.99	ng	# 95
31) Naphthalene	8.66	128	289814	40.95	ng	99
32) Benzoic acid	8.37	122	43266	39.02	ng	97
33) 4-Chloroaniline	8.75	127	55780	19.42	ng	96
34) Hexachlorobutadiene	8.82	225	49148	40.25	ng	98
35) Caprolactam	9.17	113	25697	45.66	ng	# 79
36) 4-Chloro-3-methylphenol	9.37	107	90321	46.62	ng	88
37) 2-Methylnaphthalene	9.53	142	203772	42.86	ng	98
39) 1,2,4,5-Tetrachlorobenzene	9.73	216	84497	43.34	ng	# 100
40) Hexachlorocyclopentadiene	9.71	237	82033	83.89	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.88	196	60113	44.51	ng	98
43) 2,4,5-Trichlorophenol	9.94	196	62748	43.01	ng	96
45) 1,1'-Biphenyl	10.11	154	238831	45.71	ng	98
46) 2-Chloronaphthalene	10.12	162	191321	45.06	ng	93
47) 2-Nitroaniline	10.26	65	47112	44.67	ng	85
48) Acenaphthylene	10.64	152	303386	42.97	ng	99
49) Dimethylphthalate	10.50	163	219256	45.23	ng	99
50) 2,6-Dinitrotoluene	10.57	165	47188	46.96	ng	# 86
51) Acenaphthene	10.84	154	170513	42.12	ng	97
52) 3-Nitroaniline	10.77	138	32217	27.61	ng	86
53) 2,4-Dinitrophenol	10.91	184	30072	81.31	ng	# 79
54) Dibenzofuran	11.06	168	265947	44.29	ng	92
55) 4-Nitrophenol	11.01	139	74918	86.66	ng	# 79
56) 2,4-Dinitrotoluene	11.07	165	65726	50.16	ng	# 64
57) Fluorene	11.48	166	220932	44.31	ng	98
58) 2,3,4,6-Tetrachlorophenol	11.22	232	48517	43.19	ng	# 100
59) Diethylphthalate	11.38	149	215393	45.60	ng	97
60) 4-Chlorophenyl-phenylether	11.51	204	100823	44.31	ng	93
61) 4-Nitroaniline	11.53	138	48251	41.48	ng	80
62) Azobenzene	11.70	77	223761	49.56	ng	95
64) 4,6-Dinitro-2-methylphenol	11.57	198	22078	36.14	ng	# 68
65) n-Nitrosodiphenylamine	11.65	169	196267	44.09	ng	99
66) 4-Bromophenyl-phenylether	12.11	248	59684	41.83	ng	94
67) Hexachlorobenzene	12.16	284	64229	40.97	ng	# 85
68) Atrazine	12.33	200	55506	44.49	ng	96
69) Pentachlorophenol	12.42	266	55395	67.94	ng	98
70) Phenanthrene	12.67	178	323313	42.12	ng	99
71) Anthracene	12.74	178	330906	43.64	ng	99
72) Carbazole	12.96	167	315365	43.72	ng	98
73) Di-n-butylphthalate	13.41	149	344887	42.64	ng	98
74) Fluoranthene	14.15	202	358172	42.31	ng	95
76) Benzidine	14.34	184	150218	43.00	ng	100
77) Pyrene	14.43	202	385184	43.33	ng	99
79) Butylbenzylphthalate	15.27	149	151943	43.36	ng	# 78
80) Benzo(a)anthracene	15.93	228	353266	42.16	ng	100
81) 3,3'-Dichlorobenzidine	15.92	252	69636	25.19	ng	97
82) Chrysene	15.97	228	345720	45.89	ng	100
83) Bis(2-ethylhexyl)phthalate	16.01	149	226787	42.72	ng	98
84) Di-n-octyl phthalate	16.82	149	380670	41.94	ng	100
85) Indeno(1,2,3-cd)pyrene	19.07	276	392615	41.75	ng	# 100
87) Benzo(b)fluoranthene	17.25	252	404210	44.67	ng	98
88) Benzo(k)fluoranthene	17.29	252	276675	39.14	ng	98
89) Benzo(a)pyrene	17.64	252	333104	44.12	ng	# 96
90) Dibenzo(a,h)anthracene	19.09	278	321196	42.89	ng	# 96
91) Benzo(g,h,i)perylene	19.47	276	323456	42.43	ng	96

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

