

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF062615\
 Data File : BF080194.D
 Acq On : 26 Jun 2015 17:31
 Operator : TP/IZ
 Sample : PB84223BSD
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB84223BSD

Manual Integrations
 APPROVED

apatel
 6/29/2015 11:38:22 AM

Quant Time: Jun 27 00:04:45 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF061715.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 26 23:02:05 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.29	152	28852	20.00	ng	0.00
21) Naphthalene-d8	8.58	136	112075	20.00	ng	0.00
38) Acenaphthene-d10	10.36	164	58690	20.00	ng	0.00
63) Phenanthrene-d10	11.86	188	117285	20.00	ng	-0.02
75) Chrysene-d12	14.85	240	122410	20.00	ng	-0.18
86) Perylene-d12	16.73	264	116945	20.00	ng	-0.29

System Monitoring Compounds

5) 2-Fluorophenol	5.90	112	205713	126.21	ng	0.00
7) Phenol-d6	6.89	99	268204	123.66	ng	0.00
23) Nitrobenzene-d5	7.85	82	170337	88.48	ng	0.00
41) 2,4,6-Tribromophenol	11.16	330	69145	120.48	ng	0.00
44) 2-Fluorobiphenyl	9.66	172	320835	77.96	ng	0.00
78) Terphenyl-d14	13.57	244	400243	76.36	ng	-0.09

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.14	88	24040	31.03	ng	88
3) Pyridine	3.98	79	72038	34.96	ng	# 81
4) n-Nitrosodimethylamine	3.89	42	35978	36.74	ng	95
6) Aniline	6.95	93	84008	28.15	ng	97
8) 2-Chlorophenol	7.08	128	78842	41.85	ng	90
9) Benzaldehyde	6.84	77	33038	26.39	ng	# 69
10) Phenol	6.92	94	90169	38.10	ng	88
11) bis(2-Chloroethyl)ether	7.01	93	68679	36.57	ng	94
12) 1,3-Dichlorobenzene	7.24	146	86172	38.08	ng	# 93
13) 1,4-Dichlorobenzene	7.32	146	87788	40.79	ng	95
14) 1,2-Dichlorobenzene	7.46	146	79143m	37.79	ng	
15) Benzyl Alcohol	7.42	79	68230	38.47	ng	# 76
16) 2,2'-oxybis(1-Chloropropan	7.56	45	116452	41.78	ng	89
17) 2-Methylphenol	7.52	107	58237	36.19	ng	# 83
18) Hexachloroethane	7.82	117	29445	41.09	ng	# 79
19) n-Nitroso-di-n-propylamine	7.69	70	51623	34.28	ng	# 79
20) 3+4-Methylphenols	7.68	107	81599	39.29	ng	92
22) Acetophenone	7.69	105	111268	42.60	ng	# 84
24) Nitrobenzene	7.88	77	75885	38.19	ng	# 81
25) Isophorone	8.10	82	146807	37.89	ng	# 83
26) 2-Nitrophenol	8.20	139	36958	44.44	ng	# 81
27) 2,4-Dimethylphenol	8.22	122	70357	40.57	ng	87
28) bis(2-Chloroethoxy)methane	8.31	93	93657	41.14	ng	# 91
29) 2,4-Dichlorophenol	8.44	162	65061	43.81	ng	96
30) 1,2,4-Trichlorobenzene	8.53	180	72994	43.28	ng	99
31) Naphthalene	8.61	128	226761m	38.70	ng	
32) Benzoic acid	8.28	122	15525	14.81	ng	# 59
33) 4-Chloroaniline	8.64	127	57951m	23.11	ng	
34) Hexachlorobutadiene	8.73	225	39173	37.72	ng	99
35) Caprolactam	8.98	113	19329	40.10	ng	# 83
36) 4-Chloro-3-methylphenol	9.11	107	61903	36.95	ng	# 73
37) 2-Methylnaphthalene	9.30	142	164009	43.27	ng	# 92
39) 1,2,4,5-Tetrachlorobenzene	9.48	216	71002	41.57	ng	98
40) Hexachlorocyclopentadiene	9.46	237	61200	70.45	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.58	196	47500	40.88	ng	98
43) 2,4,5-Trichlorophenol	9.61	196	50030	37.37	ng #	93
45) 1,1'-Biphenyl	9.76	154	175865	36.83	ng	92
46) 2-Chloronaphthalene	9.80	162	149623	38.62	ng	95
47) 2-Nitroaniline	9.88	65	41570	39.09	ng #	61
48) Acenaphthylene	10.22	152	255307	39.48	ng	97
49) Dimethylphthalate	10.05	163	158836	36.00	ng #	97
50) 2,6-Dinitrotoluene	10.12	165	35607	40.25	ng #	60
51) Acenaphthene	10.39	154	160760	40.64	ng	91
52) 3-Nitroaniline	10.29	138	32503	31.85	ng #	56
53) 2,4-Dinitrophenol	10.40	184	29867	82.74	ng #	45
54) Dibenzofuran	10.56	168	220100	39.21	ng #	90
55) 4-Nitrophenol	10.44	139	66096	81.02	ng #	56
56) 2,4-Dinitrotoluene	10.53	165	53360	47.54	ng #	92
57) Fluorene	10.92	166	170704	38.32	ng	92
58) 2,3,4,6-Tetrachlorophenol	10.68	232	43433	38.43	ng	97
59) Diethylphthalate	10.76	149	165713	37.30	ng	97
60) 4-Chlorophenyl-phenylether	10.89	204	81909	38.27	ng #	88
61) 4-Nitroaniline	10.90	138	42356	40.18	ng #	57
62) Azobenzene	11.05	77	172592	38.16	ng	85
64) 4,6-Dinitro-2-methylphenol	10.94	198	22387	40.39	ng	93
65) n-Nitrosodiphenylamine	11.01	169	155273	40.66	ng	92
66) 4-Bromophenyl-phenylether	11.40	248	47679	38.23	ng #	82
67) Hexachlorobenzene	11.48	284	52049	37.55	ng #	77
68) Atrazine	11.53	200	44659	37.01	ng	84
69) Pentachlorophenol	11.66	266	53831	68.54	ng	100
70) Phenanthrene	11.89	178	262499	36.97	ng	98
71) Anthracene	11.94	178	275867	40.35	ng	99
72) Carbazole	12.09	167	252660	38.71	ng	97
73) Di-n-butylphthalate	12.42	149	272865	36.18	ng #	99
74) Fluoranthene	13.16	202	284656	37.64	ng	90
76) Benzidine	13.27	184	137012	40.04	ng	98
77) Pyrene	13.41	202	293520	38.95	ng	97
79) Butylbenzylphthalate	14.12	149	114337	37.78	ng	95
80) Benzo(a)anthracene	14.84	228	286405	38.41	ng	97
81) 3,3'-Dichlorobenzidine	14.78	252	59753	23.23	ng #	93
82) Chrysene	14.88	228	262305	38.14	ng	95
83) Bis(2-ethylhexyl)phthalate	14.81	149	164010	34.29	ng #	93
84) Di-n-octyl phthalate	15.64	149	270299	32.98	ng	96
85) Indeno(1,2,3-cd)pyrene	18.56	276	315850	36.42	ng #	89
87) Benzo(b)fluoranthene	16.20	252	268277	37.89	ng #	86
88) Benzo(k)fluoranthene	16.23	252	285442	39.59	ng #	88
89) Benzo(a)pyrene	16.65	252	271940	40.44	ng #	86
90) Dibenzo(a,h)anthracene	18.59	278	263815	38.33	ng #	83
91) Benzo(g,h,i)perylene	19.11	276	270205	38.88	ng #	78

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

