

Data Path : Z:\HPCHEM1\BNA F\DATA\BF020717\
 Data File : BF092831.D
 Acq On : 7 Feb 2017 16:39
 Operator : SJ/MA
 Sample : PB96775BS
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB96775BS

Manual Integrations
 APPROVED

mohammad
 2/8/2017 10:42:58 AM

Quant Time: Feb 08 00:03:08 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF013017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 07 23:52:21 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.91	152	175849	20.00	ng	0.00
21) Naphthalene-d8	8.19	136	685465	20.00	ng	0.00
38) Acenaphthene-d10	9.95	164	358622	20.00	ng	0.00
63) Phenanthrene-d10	11.44	188	514490	20.00	ng	0.00
75) Chrysene-d12	14.08	240	392661	20.00	ng	0.00
86) Perylene-d12	15.53	264	330235	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.52	112	1119230	109.19	ng	0.01
7) Phenol-d6	6.56	99	1310266	99.35	ng	0.01
23) Nitrobenzene-d5	7.48	82	896087	72.80	ng	0.01
41) 2,4,6-Tribromophenol	10.74	330	276639	106.65	ng	0.00
44) 2-Fluorobiphenyl	9.28	172	1612578	81.46	ng	0.01
78) Terphenyl-d14	13.03	244	1172264	70.15	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.56	88	137249	24.78	ng	# 84
3) Pyridine	3.29	79	423776	30.09	ng	# 79
4) n-Nitrosodimethylamine	3.24	42	214247	32.40	ng	# 84
6) Aniline	6.58	93	403619	22.44	ng	# 47
8) 2-Chlorophenol	6.69	128	378946	33.51	ng	# 86
9) Benzaldehyde	6.45	77	164372	17.40	ng	# 96
10) Phenol	6.57	94	501131	32.48	ng	# 89
11) bis(2-Chloroethyl)ether	6.65	93	399169	32.41	ng	# 95
12) 1,3-Dichlorobenzene	6.85	146	451894	34.12	ng	# 97
13) 1,4-Dichlorobenzene	6.92	146	426551	31.82	ng	# 99
14) 1,2-Dichlorobenzene	7.08	146	443970	34.64	ng	# 96
15) Benzyl Alcohol	7.06	79	354445	36.30	ng	# 94
16) 2,2'-oxybis(1-Chloropropan	7.18	45	681793	31.86	ng	# 96
17) 2-Methylphenol	7.17	107	341662	35.22	ng	# 97
18) Hexachloroethane	7.42	117	131888	28.82	ng	# 84
19) n-Nitroso-di-n-propylamine	7.32	70	264834	31.55	ng	# 93
20) 3+4-Methylphenols	7.32	107	413022	35.61	ng	# 67
22) Acetophenone	7.32	105	519153	33.17	ng	# 94
24) Nitrobenzene	7.49	77	405218	32.06	ng	# 97
25) Isophorone	7.73	82	772161	33.66	ng	# 98
26) 2-Nitrophenol	7.81	139	229135	38.14	ng	# 87
27) 2,4-Dimethylphenol	7.85	122	374530	35.49	ng	# 99
28) bis(2-Chloroethoxy)methane	7.94	93	494377	32.54	ng	# 98
29) 2,4-Dichlorophenol	8.05	162	331862	33.72	ng	# 89
30) 1,2,4-Trichlorobenzene	8.13	180	351458	33.14	ng	# 95
31) Naphthalene	8.21	128	1067465	30.72	ng	# 99
32) Benzoic acid	7.97	122	207262	36.66	ng	# 95
33) 4-Chloroaniline	8.27	127	141173	10.36	ng	# 95
34) Hexachlorobutadiene	8.33	225	187264	32.09	ng	# 99
35) Caprolactam	8.64	113	113309	36.60	ng	# 39
36) 4-Chloro-3-methylphenol	8.76	107	384562	37.48	ng	# 89
37) 2-Methylnaphthalene	8.91	142	791065	35.46	ng	# 96
39) 1,2,4,5-Tetrachlorobenzene	9.07	216	323270	32.11	ng	# 99
40) Hexachlorocyclopentadiene	9.06	237	76341	16.81	ng	# 97

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.18	196	239255	36.17	nq	91
43) 2,4,5-Trichlorophenol	9.23	196	245197	33.83	nq	93
45) 1,1'-Biphenyl	9.37	154	864922	33.31	nq	96
46) 2-Chloronaphthalene	9.40	162	738683	36.36	nq	95
47) 2-Nitroaniline	9.49	65	228152	33.40	nq	96
48) Acenaphthylene	9.81	152	1130209	32.21	nq	99
49) Dimethylphthalate	9.68	163	934254	38.68	nq	100
50) 2,6-Dinitrotoluene	9.73	165	174783	33.50	nq	# 72
51) Acenaphthene	9.98	154	697457	33.24	nq	96
52) 3-Nitroaniline	9.90	138	124467	20.85	nq	96
53) 2,4-Dinitrophenol	10.02	184	53643	23.75	nq	95
54) Dibenzofuran	10.16	168	986357	35.15	nq	94
55) 4-Nitrophenol	10.09	139	323690	75.28	nq	# 68
56) 2,4-Dinitrotoluene	10.14	165	262516	37.80	nq	# 85
57) Fluorene	10.50	166	725707	33.61	nq	98
58) 2,3,4,6-Tetrachlorophenol	10.28	232	188048	38.89	nq	92
59) Diethylphthalate	10.37	149	851579	37.41	nq	97
60) 4-Chlorophenyl-phenylether	10.49	204	340398	32.82	nq	92
61) 4-Nitroaniline	10.52	138	187395	30.65	nq	87
62) Azobenzene	10.65	77	836571	35.57	nq	97
64) 4,6-Dinitro-2-methylphenol	10.56	198	47364	16.95	nq	94
65) n-Nitrosodiphenylamine	10.61	169	708722	43.12	nq	99
66) 4-Bromophenyl-phenylether	10.98	248	200869	37.37	nq	98
67) Hexachlorobenzene	11.05	284	184664	34.76	nq	# 88
68) Atrazine	11.14	200	209026	39.63	nq	97
69) Pentachlorophenol	11.24	266	178712	66.51	nq	96
70) Phenanthrene	11.46	178	935871	31.75	nq	98
71) Anthracene	11.52	178	1101051	37.20	nq	98
72) Carbazole	11.66	167	882523	31.19	nq	98
73) Di-n-butylphthalate	12.00	149	1263825	41.30	nq	# 97
74) Fluoranthene	12.65	202	890491	29.29	nq	97
76) Benzidine	12.77	184	540459	32.30	nq	96
77) Pyrene	12.88	202	895033	30.46	nq	99
79) Butylbenzylphthalate	13.49	149	459656	38.52	nq	86
80) Benzo(a)anthracene	14.07	228	796934	33.18	nq	99
81) 3,3'-Dichlorobenzidine	14.03	252	308928	36.09	nq	# 95
82) Chrysene	14.10	228	680572	30.27	nq	99
83) Bis(2-ethylhexyl)phthalate	14.05	149	635137	38.92	nq	# 95
84) Di-n-octyl phthalate	14.66	149	1026902	38.64	nq	100
85) Indeno(1,2,3-cd)pyrene	16.98	276	601653	34.54	nq	95
87) Benzo(b)fluoranthene	15.11	252	897465m	41.29	nq	
88) Benzo(k)fluoranthene	15.14	252	532842m	28.39	nq	
89) Benzo(a)pyrene	15.47	252	655807	34.88	nq	99
90) Dibenzo(a,h)anthracene	16.99	278	516833	34.01	nq	99
91) Benzo(g,h,i)perylene	17.41	276	485961	31.66	nq	# 93

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

