

Data Path : Z:\HPCHEM1\BNA F\DATA\BF092815\
 Data File : BF081850.D
 Acq On : 28 Sep 2015 18:42
 Operator : UM/IZ
 Sample : PB85617BS
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB85617BS

Manual Integrations
 APPROVED

MMDadoda
 9/29/2015 12:10:45 PM

Quant Time: Sep 29 03:39:30 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF092815.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Sep 28 18:10:42 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.94	152	95178	20.00	ng	0.00
21) Naphthalene-d8	8.22	136	388607	20.00	ng	0.00
38) Acenaphthene-d10	9.98	164	189440	20.00	ng	0.00
63) Phenanthrene-d10	11.46	188	366442	20.00	ng	0.00
75) Chrysene-d12	14.12	240	354787	20.00	ng	0.00
86) Perylene-d12	15.58	264	289293	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.53	112	450369	85.90	ng	0.01
7) Phenol-d6	6.56	99	582741	88.33	ng	0.00
23) Nitrobenzene-d5	7.50	82	484626	83.13	ng	0.00
41) 2,4,6-Tribromophenol	10.77	330	163918	85.44	ng	0.00
44) 2-Fluorobiphenyl	9.30	172	1008612	90.15	ng	0.00
78) Terphenyl-d14	13.05	244	950774	87.17	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.61	88	81442	35.72	ng	89
3) Pyridine	3.36	79	210897	35.53	ng	85
4) n-Nitrosodimethylamine	3.31	42	106271	39.41	ng	96
6) Aniline	6.59	93	231896	26.16	ng	# 84
8) 2-Chlorophenol	6.72	128	251530	43.44	ng	97
9) Benzaldehyde	6.48	77	84349	19.52	ng	# 81
10) Phenol	6.57	94	333367	45.05	ng	74
11) bis(2-Chloroethyl)ether	6.66	93	246059	42.87	ng	97
12) 1,3-Dichlorobenzene	6.87	146	291071	42.56	ng	# 94
13) 1,4-Dichlorobenzene	6.95	146	297960	41.72	ng	# 94
14) 1,2-Dichlorobenzene	7.10	146	264022m	41.49	ng	
15) Benzyl Alcohol	7.07	79	210951	44.29	ng	# 77
16) 2,2'-oxybis(1-Chloropropan	7.21	45	367813	45.33	ng	79
17) 2-Methylphenol	7.19	107	206257	43.94	ng	# 87
18) Hexachloroethane	7.44	117	93537	41.84	ng	# 88
19) n-Nitroso-di-n-propylamine	7.35	70	172587	43.04	ng	# 78
20) 3+4-Methylphenols	7.34	107	264891	44.68	ng	# 61
22) Acetophenone	7.35	105	345311	43.46	ng	# 85
24) Nitrobenzene	7.52	77	272913	43.11	ng	# 74
25) Isophorone	7.75	82	471335	40.79	ng	# 86
26) 2-Nitrophenol	7.84	139	133956	44.08	ng	# 81
27) 2,4-Dimethylphenol	7.87	122	236338	44.21	ng	# 85
28) bis(2-Chloroethoxy)methane	7.97	93	289307	42.16	ng	# 93
29) 2,4-Dichlorophenol	8.08	162	230560	44.48	ng	98
30) 1,2,4-Trichlorobenzene	8.16	180	242017	41.82	ng	99
31) Naphthalene	8.24	128	712166m	41.36	ng	
32) Benzoic acid	7.99	122	125682	40.64	ng	# 67
33) 4-Chloroaniline	8.30	127	124211	16.68	ng	# 93
34) Hexachlorobutadiene	8.35	225	144542	42.24	ng	98
35) Caprolactam	8.66	113	66893	46.62	ng	98
36) 4-Chloro-3-methylphenol	8.78	107	240898	47.53	ng	89
37) 2-Methylnaphthalene	8.94	142	515348	43.84	ng	# 91
39) 1,2,4,5-Tetrachlorobenzene	9.10	216	239546	41.19	ng	99
40) Hexachlorocyclopentadiene	9.09	237	289258	96.58	ng	97

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.21	196	169564	42.53	nq	99
43) 2,4,5-Trichlorophenol	9.26	196	182406	44.49	nq #	94
45) 1,1'-Biphenyl	9.39	154	604239	41.34	nq	94
46) 2-Chloronaphthalene	9.43	162	499122	43.49	nq	98
47) 2-Nitroaniline	9.52	65	143350	42.55	nq #	65
48) Acenaphthylene	9.84	152	773080	42.09	nq	96
49) Dimethylphthalate	9.70	163	598496	45.77	nq #	98
50) 2,6-Dinitrotoluene	9.76	165	124458	43.84	nq #	55
51) Acenaphthene	10.01	154	516369	47.34	nq	88
52) 3-Nitroaniline	9.93	138	84506	25.73	nq #	62
53) 2,4-Dinitrophenol	10.05	184	109574	79.54	nq #	79
54) Dibenzofuran	10.18	168	673459	43.69	nq #	87
55) 4-Nitrophenol	10.09	139	207813	87.57	nq #	41
56) 2,4-Dinitrotoluene	10.17	165	184870	51.08	nq #	89
57) Fluorene	10.53	166	515617	46.86	nq	93
58) 2,3,4,6-Tetrachlorophenol	10.30	232	147896	45.47	nq	97
59) Diethylphthalate	10.40	149	571711	46.81	nq	99
60) 4-Chlorophenyl-phenylether	10.53	204	239981	42.53	nq	88
61) 4-Nitroaniline	10.55	138	133818	40.64	nq #	45
62) Azobenzene	10.67	77	548013	45.97	nq	87
64) 4,6-Dinitro-2-methylphenol	10.57	198	80649	48.40	nq	90
65) n-Nitrosodiphenylamine	10.64	169	481730	43.45	nq	91
66) 4-Bromophenyl-phenylether	11.01	248	161582	43.73	nq #	85
67) Hexachlorobenzene	11.07	284	189790	45.51	nq #	80
68) Atrazine	11.17	200	156585	45.52	nq	83
69) Pentachlorophenol	11.27	266	214594	90.32	nq	99
70) Phenanthrene	11.50	178	819283	46.52	nq	97
71) Anthracene	11.54	178	821090	44.09	nq	96
72) Carbazole	11.70	167	793500	47.07	nq	96
73) Di-n-butylphthalate	12.02	149	838612	48.96	nq #	98
74) Fluoranthene	12.69	202	872324	44.44	nq	86
76) Benzidine	12.81	184	252897	23.66	nq	98
77) Pyrene	12.92	202	932397	43.99	nq	96
79) Butylbenzylphthalate	13.53	149	383486	48.08	nq	96
80) Benzo(a)anthracene	14.10	228	803156	42.04	nq	95
81) 3,3'-Dichlorobenzidine	14.07	252	164123	25.44	nq #	92
82) Chrysene	14.14	228	769078	47.46	nq	94
83) Bis(2-ethylhexyl)phthalate	14.08	149	464015	46.38	nq #	90
84) Di-n-octyl phthalate	14.70	149	780768	47.96	nq #	92
85) Indeno(1,2,3-cd)pyrene	17.05	276	654364	43.78	nq #	86
87) Benzo(b)fluoranthene	15.16	252	780969	47.28	nq #	85
88) Benzo(k)fluoranthene	15.18	252	660348m	43.62	nq	
89) Benzo(a)pyrene	15.52	252	671204	46.85	nq #	86
90) Dibenzo(a,h)anthracene	17.06	278	535721	44.57	nq #	81
91) Benzo(g,h,i)perylene	17.50	276	547469	44.44	nq #	74

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

