

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF050517\
 Data File : BF094946.D
 Acq On : 6 May 2017 6:57
 Operator : SJ/MA
 Sample : PB98737BS
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB98737BS

Manual Integrations
 APPROVED

mohammad
 5/8/2017 2:48:45 PM

Quant Time: May 08 01:49:38 2017
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050217.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 03 17:24:51 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.72	152	99662	20.00	ng	0.00
21) Naphthalene-d8	9.74	136	410050	20.00	ng	-0.01
38) Acenaphthene-d10	12.57	164	196389	20.00	ng	-0.01
63) Phenanthrene-d10	14.97	188	333261	20.00	ng	-0.01
75) Chrysene-d12	18.64	240	247448	20.00	ng	-0.01
86) Perylene-d12	20.31	264	200631	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.68	112	813637	136.11	ng	0.00
7) Phenol-d6	7.27	99	955433	133.21	ng	0.00
23) Nitrobenzene-d5	8.63	82	556103	85.52	ng	0.00
41) 2,4,6-Tribromophenol	13.88	330	207539	129.04	ng	0.00
44) 2-Fluorobiphenyl	11.52	172	1123546	82.35	ng	-0.01
78) Terphenyl-d14	17.30	244	968543	86.21	ng	-0.01

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.16	88	99939	34.09	ng	93
3) Pyridine	2.84	79	235148	34.61	ng	97
4) n-Nitrosodimethylamine	2.76	42	116084	40.99	ng	87
6) Aniline	7.23	93	225898	24.95	ng	96
8) 2-Chlorophenol	7.41	128	317093	46.60	ng	96
9) Benzaldehyde	7.05	77	40611	9.27	ng	91
10) Phenol	7.29	94	348078	46.05	ng	90
11) bis(2-Chloroethyl)ether	7.37	93	271753	43.55	ng	95
12) 1,3-Dichlorobenzene	7.63	146	326462	42.30	ng	99
13) 1,4-Dichlorobenzene	7.75	146	322200	41.16	ng	97
14) 1,2-Dichlorobenzene	7.99	146	313902	42.97	ng	96
15) Benzyl Alcohol	8.01	79	241746	52.40	ng	97
16) 2,2'-oxybis(1-Chloropropan	8.21	45	363978	41.21	ng	52
17) 2-Methylphenol	8.22	107	253158	45.47	ng	# 88
18) Hexachloroethane	8.51	117	106912	40.32	ng	# 86
19) n-Nitroso-di-n-propylamine	8.44	70	208300	42.94	ng	92
20) 3+4-Methylphenols	8.48	107	330030	43.62	ng	# 63
22) Acetophenone	8.40	105	412938	41.97	ng	# 84
24) Nitrobenzene	8.65	77	289816	42.52	ng	94
25) Isophorone	9.05	82	553983	47.71	ng	# 94
26) 2-Nitrophenol	9.16	139	174132	47.35	ng	98
27) 2,4-Dimethylphenol	9.30	122	269339	45.29	ng	97
28) bis(2-Chloroethoxy)methane	9.42	93	351570	43.77	ng	98
29) 2,4-Dichlorophenol	9.58	162	244487	43.55	ng	98
30) 1,2,4-Trichlorobenzene	9.67	180	247132	41.08	ng	99
31) Naphthalene	9.78	128	830969	40.32	ng	100
32) Benzoic acid	9.62	122	164308	42.71	ng	97
33) 4-Chloroaniline	9.94	127	47925	6.24	ng	96
34) Hexachlorobutadiene	10.00	225	126291	39.79	ng	98
35) Caprolactam	10.55	113	83052m	49.26	ng	
36) 4-Chloro-3-methylphenol	10.77	107	279726	46.60	ng	89
37) 2-Methylnaphthalene	10.90	142	609248	45.04	ng	98
39) 1,2,4,5-Tetrachlorobenzene	11.18	216	233130	38.60	ng	99
40) Hexachlorocyclopentadiene	11.16	237	206968	81.40	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	11.40	196	173115	42.93	ng	97
43) 2,4,5-Trichlorophenol	11.48	196	176898	40.82	ng #	94
45) 1,1'-Biphenyl	11.67	154	683018	41.31	ng	99
46) 2-Chloronaphthalene	11.68	162	541124	40.53	ng	95
47) 2-Nitroaniline	11.90	65	153880	42.11	ng #	83
48) Acenaphthylene	12.34	152	870575	41.63	ng	99
49) Dimethylphthalate	12.22	163	648643	40.23	ng	98
50) 2,6-Dinitrotoluene	12.31	165	144992	44.08	ng	96
51) Acenaphthene	12.62	154	509887	40.82	ng	99
52) 3-Nitroaniline	12.58	138	58720	16.63	ng #	94
53) 2,4-Dinitrophenol	12.77	184	143527	79.29	ng #	66
54) Dibenzofuran	12.91	168	786928	45.53	ng	98
55) 4-Nitrophenol	12.95	139	188476	88.89	ng #	77
56) 2,4-Dinitrotoluene	12.97	165	194665	46.06	ng #	92
57) Fluorene	13.47	166	624845	41.57	ng	99
58) 2,3,4,6-Tetrachlorophenol	13.15	232	137569	50.44	ng #	93
59) Diethylphthalate	13.37	149	636692	41.20	ng	99
60) 4-Chlorophenyl-phenylether	13.49	204	276668	41.89	ng	90
61) 4-Nitroaniline	13.58	138	132252	40.81	ng	96
62) Azobenzene	13.75	77	586836	47.26	ng	93
64) 4,6-Dinitro-2-methylphenol	13.64	198	101856	45.59	ng #	66
65) n-Nitrosodiphenylamine	13.70	169	532943	44.06	ng	99
66) 4-Bromophenyl-phenylether	14.28	248	158123	44.91	ng	96
67) Hexachlorobenzene	14.36	284	158531	43.93	ng #	83
68) Atrazine	14.62	200	163532	47.89	ng	96
69) Pentachlorophenol	14.71	266	142090	86.57	ng	99
70) Phenanthrene	15.01	178	828891	43.29	ng	98
71) Anthracene	15.09	178	869921	44.48	ng	99
72) Carbazole	15.38	167	805949	45.29	ng	97
73) Di-n-butylphthalate	15.94	149	994492	44.81	ng	99
74) Fluoranthene	16.75	202	889621	45.29	ng	99
76) Benzidine	16.97	184	276849	42.28	ng #	93
77) Pyrene	17.05	202	891079	48.15	ng	99
79) Butylbenzylphthalate	17.96	149	373918	43.44	ng	90
80) Benzo(a)anthracene	18.63	228	625813	43.46	ng	99
81) 3,3'-Dichlorobenzidine	18.61	252	123639	24.86	ng #	94
82) Chrysene	18.68	228	600682	43.14	ng	98
83) Bis(2-ethylhexyl)phthalate	18.71	149	550337	44.87	ng #	99
84) Di-n-octyl phthalate	19.47	149	953888	47.44	ng	96
85) Indeno(1,2,3-cd)pyrene	21.59	276	529216	46.80	ng	100
87) Benzo(b)fluoranthene	19.90	252	544469	43.92	ng	99
88) Benzo(k)fluoranthene	19.93	252	533442	44.61	ng #	94
89) Benzo(a)pyrene	20.25	252	510540	44.63	ng	99
90) Dibenzo(a,h)anthracene	21.60	278	433416	45.88	ng	97
91) Benzo(g,h,i)perylene	21.97	276	465290	50.19	ng	96

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

