

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF070615\
 Data File : BF080352.D
 Acq On : 6 Jul 2015 18:52
 Operator : TP/IZ
 Sample : PB84363BS
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB84363BS

Manual Integrations
 APPROVED

apatel
 7/7/2015 4:25:53 PM

Quant Time: Jul 07 04:53:24 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF070615.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 07 03:05:40 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.23	152	37027	20.00	ng	0.00
21) Naphthalene-d8	8.52	136	139085	20.00	ng	0.00
38) Acenaphthene-d10	10.29	164	66676	20.00	ng	0.00
63) Phenanthrene-d10	11.79	188	139960	20.00	ng	0.00
75) Chrysene-d12	14.84	240	155033	20.00	ng	0.10
86) Perylene-d12	16.75	264	142936	20.00	ng	0.19

System Monitoring Compounds

5) 2-Fluorophenol	5.84	112	273335	135.20	ng	0.00
7) Phenol-d6	6.83	99	341613	127.47	ng	0.00
23) Nitrobenzene-d5	7.79	82	207809	87.15	ng	0.00
41) 2,4,6-Tribromophenol	11.08	330	96336	138.08	ng	0.00
44) 2-Fluorobiphenyl	9.60	172	441550	89.37	ng	0.00
78) Terphenyl-d14	13.53	244	566966	85.30	ng	0.06

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.04	88	32512	34.06	ng	94
3) Pyridine	3.88	79	88616	36.48	ng	99
4) n-Nitrosodimethylamine	3.78	42	48854	42.80	ng	98
6) Aniline	6.89	93	98836	26.64	ng	99
8) 2-Chlorophenol	7.01	128	99782	42.58	ng	99
9) Benzaldehyde	6.77	77	58811	39.46	ng	99
10) Phenol	6.85	94	121503	41.81	ng	97
11) bis(2-Chloroethyl)ether	6.94	93	88998	40.60	ng	97
12) 1,3-Dichlorobenzene	7.17	146	114549	39.81	ng	99
13) 1,4-Dichlorobenzene	7.25	146	104118	38.75	ng	97
14) 1,2-Dichlorobenzene	7.40	146	111803	41.11	ng	98
15) Benzyl Alcohol	7.36	79	87547	42.20	ng	98
16) 2,2'-oxybis(1-Chloropropan	7.49	45	146873	41.43	ng	99
17) 2-Methylphenol	7.46	107	71879m	39.57	ng	
18) Hexachloroethane	7.74	117	33621	38.95	ng	# 61
19) n-Nitroso-di-n-propylamine	7.62	70	66234	38.09	ng	# 77
20) 3+4-Methylphenols	7.62	107	104085	41.11	ng	96
22) Acetophenone	7.63	105	143356	44.68	ng	# 98
24) Nitrobenzene	7.80	77	102537	42.81	ng	98
25) Isophorone	8.04	82	193692	42.20	ng	98
26) 2-Nitrophenol	8.13	139	43231	41.70	ng	99
27) 2,4-Dimethylphenol	8.16	122	89697	43.98	ng	98
28) bis(2-Chloroethoxy)methane	8.25	93	124449	43.66	ng	99
29) 2,4-Dichlorophenol	8.36	162	80358	44.38	ng	# 83
30) 1,2,4-Trichlorobenzene	8.46	180	85123	39.55	ng	99
31) Naphthalene	8.54	128	289212m	40.54	ng	
32) Benzoic acid	8.22	122	43135	41.46	ng	95
33) 4-Chloroaniline	8.58	127	73699	25.92	ng	97
34) Hexachlorobutadiene	8.66	225	54189	41.39	ng	100
35) Caprolactam	8.92	113	24983	44.67	ng	88
36) 4-Chloro-3-methylphenol	9.05	107	77807	41.64	ng	97
37) 2-Methylnaphthalene	9.24	142	192043	42.22	ng	98
39) 1,2,4,5-Tetrachlorobenzene	9.40	216	86707	41.03	ng	98
40) Hexachlorocyclopentadiene	9.39	237	75570	85.50	ng	100

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42) 2,4,6-Trichlorophenol	9.52	196	58021	41.68	ng	97
43) 2,4,5-Trichlorophenol	9.55	196	71259	43.92	ng	97
45) 1,1'-Biphenyl	9.70	154	255732	43.21	ng	99
46) 2-Chloronaphthalene	9.73	162	188953	41.87	ng	99
47) 2-Nitroaniline	9.81	65	57394	45.45	ng	99
48) Acenaphthylene	10.14	152	301290	40.11	ng	98
49) Dimethylphthalate	9.98	163	217632	40.65	ng	100
50) 2,6-Dinitrotoluene	10.05	165	50676	45.28	ng	93
51) Acenaphthene	10.33	154	196134	43.62	ng	94
52) 3-Nitroaniline	10.22	138	38591	30.60	ng	95
53) 2,4-Dinitrophenol	10.34	184	41189	84.14	ng	97
54) Dibenzofuran	10.50	168	276033	43.18	ng	99
55) 4-Nitrophenol	10.37	139	87643	94.82	ng	90
56) 2,4-Dinitrotoluene	10.46	165	64176	44.59	ng	100
57) Fluorene	10.84	166	235597	42.80	ng	99
58) 2,3,4,6-Tetrachlorophenol	10.61	232	57147	43.97	ng	99
59) Diethylphthalate	10.69	149	242343	45.32	ng	100
60) 4-Chlorophenyl-phenylether	10.83	204	104752	43.00	ng	99
61) 4-Nitroaniline	10.84	138	52677	43.03	ng	98
62) Azobenzene	10.99	77	222886	43.88	ng	99
64) 4,6-Dinitro-2-methylphenol	10.88	198	33077	41.42	ng	94
65) n-Nitrosodiphenylamine	10.94	169	192810	43.25	ng	99
66) 4-Bromophenyl-phenylether	11.32	248	71301	44.47	ng	97
67) Hexachlorobenzene	11.40	284	73286	43.57	ng	97
68) Atrazine	11.46	200	54522	41.08	ng	97
69) Pentachlorophenol	11.60	266	74700	89.82	ng	95
70) Phenanthrene	11.82	178	352071	41.94	ng	99
71) Anthracene	11.87	178	351966	42.76	ng	100
72) Carbazole	12.03	167	328931	42.85	ng	98
73) Di-n-butylphthalate	12.36	149	364421	41.50	ng	99
74) Fluoranthene	13.10	202	380445	42.18	ng	97
76) Benzidine	13.23	184	212922	54.15	ng	97
77) Pyrene	13.37	202	399275	42.58	ng	99
79) Butylbenzylphthalate	14.09	149	162146	43.33	ng	# 78
80) Benzo(a)anthracene	14.83	228	385982	41.31	ng	99
81) 3,3'-Dichlorobenzidine	14.77	252	105461	34.04	ng	# 99
82) Chrysene	14.88	228	359658	41.14	ng	99
83) Bis(2-ethylhexyl)phthalate	14.82	149	249035	43.66	ng	99
84) Di-n-octyl phthalate	15.67	149	410685	43.59	ng	100
85) Indeno(1,2,3-cd)pyrene	18.55	276	389840	37.92	ng	99
87) Benzo(b)fluoranthene	16.23	252	376612	41.93	ng	99
88) Benzo(k)fluoranthene	16.26	252	347053	40.81	ng	100
89) Benzo(a)pyrene	16.67	252	349780	42.58	ng	# 97
90) Dibenzo(a,h)anthracene	18.57	278	329087	40.68	ng	98
91) Benzo(g,h,i)perylene	19.07	276	330131	40.24	ng	98

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

