

Data Path : Z:\HPCHEM1\BNA F\DATA\BF050815\
 Data File : BF079057.D
 Acq On : 9 May 2015 00:27
 Operator : TP/IZ
 Sample : PB83273BSD
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB83273BSD

Manual Integrations
 APPROVED

apatel
 5/12/2015 3:44:25 PM

Quant Time: May 09 01:17:17 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF043015.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri May 08 22:44:53 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.32	152	71198	20.00	ng	0.00
21) Naphthalene-d8	8.90	136	293812	20.00	ng	0.00
38) Acenaphthene-d10	11.07	164	144124	20.00	ng	0.00
63) Phenanthrene-d10	12.91	188	267058	20.00	ng	-0.01
75) Chrysene-d12	16.22	240	256947	20.00	ng	-0.03
86) Perylene-d12	17.93	264	258938	20.00	ng	-0.15

System Monitoring Compounds

5) 2-Fluorophenol	5.62	112	552887	133.67	ng	0.00
7) Phenol-d6	6.88	99	705076	128.96	ng	-0.01
23) Nitrobenzene-d5	8.01	82	399036	78.05	ng	0.00
41) 2,4,6-Tribromophenol	12.06	330	204666	131.27	ng	0.00
44) 2-Fluorobiphenyl	10.24	172	798362	77.18	ng	-0.01
78) Terphenyl-d14	14.91	244	903033	84.14	ng	-0.01

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.31	88	68968	38.35	ng	# 24
3) Pyridine	3.05	79	204818	38.92	ng	100
4) n-Nitrosodimethylamine	2.97	42	90255m	40.02	ng	
6) Aniline	6.91	93	184231	23.87	ng	98
8) 2-Chlorophenol	7.05	128	202108	43.08	ng	97
9) Benzaldehyde	6.76	77	107039	29.06	ng	91
10) Phenol	6.90	94	264157	41.65	ng	88
11) bis(2-Chloroethyl)ether	7.00	93	186132	40.03	ng	98
12) 1,3-Dichlorobenzene	7.24	146	218177	41.38	ng	95
13) 1,4-Dichlorobenzene	7.35	146	217387	40.06	ng	97
14) 1,2-Dichlorobenzene	7.53	146	207929	41.16	ng	97
15) Benzyl Alcohol	7.50	79	166125	40.93	ng	97
16) 2,2'-oxybis(1-Chloropropan	7.68	45	293768	41.30	ng	99
17) 2-Methylphenol	7.64	107	169265	44.84	ng	98
18) Hexachloroethane	7.94	117	76921	41.15	ng	# 87
19) n-Nitroso-di-n-propylamine	7.83	70	145101	39.29	ng	96
20) 3+4-Methylphenols	7.84	107	219469	43.63	ng	# 50
22) Acetophenone	7.83	105	288281	43.43	ng	96
24) Nitrobenzene	8.03	77	207542	40.96	ng	95
25) Isophorone	8.33	82	376049	39.68	ng	98
26) 2-Nitrophenol	8.43	139	99466	47.42	ng	95
27) 2,4-Dimethylphenol	8.49	122	194985	46.46	ng	97
28) bis(2-Chloroethoxy)methane	8.62	93	237265	40.61	ng	99
29) 2,4-Dichlorophenol	8.73	162	168673	45.20	ng	92
30) 1,2,4-Trichlorobenzene	8.83	180	169559	41.36	ng	97
31) Naphthalene	8.92	128	603406	43.10	ng	99
32) Benzoic acid	8.59	122	100999	39.68	ng	96
33) 4-Chloroaniline	8.99	127	129335	21.83	ng	99
34) Hexachlorobutadiene	9.07	225	91902	38.93	ng	96
35) Caprolactam	9.43	113	55244	45.78	ng	89
36) 4-Chloro-3-methylphenol	9.60	107	179644	45.83	ng	97
37) 2-Methylnaphthalene	9.78	142	402228	41.46	ng	99
39) 1,2,4,5-Tetrachlorobenzene	9.99	216	161521	39.79	ng	# 100
40) Hexachlorocyclopentadiene	9.98	237	165907	81.28	ng	97

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	10.14	196	117683	42.17	nq	97
43) 2,4,5-Trichlorophenol	10.18	196	125617	44.52	nq #	89
45) 1,1'-Biphenyl	10.36	154	486887	42.49	nq	98
46) 2-Chloronaphthalene	10.39	162	381233	42.66	nq	96
47) 2-Nitroaniline	10.51	65	111875	43.20	nq	96
48) Acenaphthylene	10.90	152	622113	42.40	nq	99
49) Dimethylphthalate	10.75	163	445766	42.14	nq	100
50) 2,6-Dinitrotoluene	10.82	165	89825	43.03	nq	94
51) Acenaphthene	11.12	154	352642	40.30	nq	98
52) 3-Nitroaniline	11.03	138	78118	29.96	nq #	97
53) 2,4-Dinitrophenol	11.16	184	66944	77.31	nq #	75
54) Dibenzofuran	11.32	168	532354	42.12	nq	97
55) 4-Nitrophenol	11.24	139	188490	91.33	nq	95
56) 2,4-Dinitrotoluene	11.32	165	129395	46.89	nq #	72
57) Fluorene	11.76	166	437929	42.21	nq	99
58) 2,3,4,6-Tetrachlorophenol	11.48	232	96913	42.04	nq #	100
59) Diethylphthalate	11.63	149	445598	42.52	nq	99
60) 4-Chlorophenyl-phenylether	11.76	204	187762	40.80	nq	90
61) 4-Nitroaniline	11.78	138	105728	40.53	nq	92
62) Azobenzene	11.95	77	437966	42.42	nq	94
64) 4,6-Dinitro-2-methylphenol	11.83	198	47730	44.27	nq #	63
65) n-Nitrosodiphenylamine	11.91	169	373452	41.99	nq	99
66) 4-Bromophenyl-phenylether	12.36	248	117147	42.08	nq #	83
67) Hexachlorobenzene	12.43	284	137783	43.31	nq #	82
68) Atrazine	12.58	200	120227	46.49	nq	98
69) Pentachlorophenol	12.69	266	125338	68.42	nq	98
70) Phenanthrene	12.95	178	641060	42.91	nq	100
71) Anthracene	13.02	178	651705	44.46	nq	100
72) Carbazole	13.22	167	620773	44.44	nq	98
73) Di-n-butylphthalate	13.67	149	743070	44.35	nq #	97
74) Fluoranthene	14.43	202	665726	42.76	nq	91
76) Benzidine	14.61	184	332523	48.02	nq	98
77) Pyrene	14.71	202	701172	47.36	nq	99
79) Butylbenzylphthalate	15.53	149	315824	48.80	nq	91
80) Benzo(a)anthracene	16.19	228	624861	44.41	nq	99
81) 3,3'-Dichlorobenzidine	16.17	252	146270	29.18	nq #	98
82) Chrysene	16.24	228	573330	42.36	nq	100
83) Bis(2-ethylhexyl)phthalate	16.25	149	461973	47.12	nq	100
84) Di-n-octyl phthalate	17.03	149	805666	46.66	nq #	100
85) Indeno(1,2,3-cd)pyrene	19.42	276	734888	42.62	nq #	100
87) Benzo(b)fluoranthene	17.49	252	711363	50.19	nq	99
88) Benzo(k)fluoranthene	17.52	252	517146m	37.69	nq	
89) Benzo(a)pyrene	17.87	252	601460	46.09	nq	98
90) Dibenzo(a,h)anthracene	19.44	278	593139	42.77	nq	97
91) Benzo(g,h,i)perylene	19.86	276	604838	43.51	nq	98

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

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