

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF081115\
 Data File : BF080964.D
 Acq On : 11 Aug 2015 16:28
 Operator : UM/IZ
 Sample : PB84907BSD
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB84907BSD

Manual Integrations
 APPROVED

apatel
 8/12/2015 8:33:55 AM

Quant Time: Aug 12 15:04:56 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF080715.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 12 14:59:04 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.78	152	61189	20.00	ng	-0.02
21) Naphthalene-d8	8.08	136	242429	20.00	ng	-0.02
38) Acenaphthene-d10	9.82	164	107536	20.00	ng	-0.02
63) Phenanthrene-d10	11.30	188	199078	20.00	ng	-0.02
75) Chrysene-d12	13.93	240	180927	20.00	ng	-0.03
86) Perylene-d12	15.32	264	169465	20.00	ng	-0.03

System Monitoring Compounds

5) 2-Fluorophenol	5.39	112	431359	115.18	ng	0.01
7) Phenol-d6	6.43	99	666241	129.82	ng	-0.01
23) Nitrobenzene-d5	7.36	82	435659	85.00	ng	-0.02
41) 2,4,6-Tribromophenol	10.61	330	131972	111.71	ng	-0.02
44) 2-Fluorobiphenyl	9.15	172	620501	81.23	ng	-0.02
78) Terphenyl-d14	12.89	244	657913	75.11	ng	-0.02

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.44	88	56854	32.01	ng	99
3) Pyridine	3.13	79	181012	35.93	ng	89
4) n-Nitrosodimethylamine	3.08	42	95513	37.14	ng	99
6) Aniline	6.45	93	205607	27.27	ng	93
8) 2-Chlorophenol	6.57	128	155449	35.87	ng	# 81
9) Benzaldehyde	6.33	77	125766	36.36	ng	97
10) Phenol	6.44	94	249878	41.04	ng	80
11) bis(2-Chloroethyl)ether	6.52	93	162359	35.71	ng	95
12) 1,3-Dichlorobenzene	6.73	146	172832	37.58	ng	98
13) 1,4-Dichlorobenzene	6.81	146	171639	35.92	ng	98
14) 1,2-Dichlorobenzene	6.96	146	158907	36.78	ng	99
15) Benzyl Alcohol	6.93	79	158057	37.63	ng	94
16) 2,2'-oxybis(1-Chloropropan	7.07	45	326779	35.73	ng	98
17) 2-Methylphenol	7.05	107	141044	38.13	ng	# 92
18) Hexachloroethane	7.30	117	65669	35.87	ng	96
19) n-Nitroso-di-n-propylamine	7.21	70	129244	34.46	ng	100
20) 3+4-Methylphenols	7.21	107	181998	39.17	ng	# 79
22) Acetophenone	7.20	105	232765	39.18	ng	# 95
24) Nitrobenzene	7.37	77	175325	33.22	ng	97
25) Isophorone	7.61	82	354149	35.93	ng	# 93
26) 2-Nitrophenol	7.69	139	84982	38.27	ng	99
27) 2,4-Dimethylphenol	7.73	122	169015	40.66	ng	96
28) bis(2-Chloroethoxy)methane	7.82	93	200778	33.93	ng	97
29) 2,4-Dichlorophenol	7.93	162	128747	37.69	ng	98
30) 1,2,4-Trichlorobenzene	8.02	180	138304	37.04	ng	98
31) Naphthalene	8.09	128	470853	35.24	ng	99
32) Benzoic acid	7.86	122	100213	40.27	ng	93
33) 4-Chloroaniline	8.14	127	120739	22.23	ng	94
34) Hexachlorobutadiene	8.21	225	80583	36.27	ng	97
35) Caprolactam	8.52	113	51417m	41.84	ng	
36) 4-Chloro-3-methylphenol	8.64	107	164285	39.41	ng	90
37) 2-Methylnaphthalene	8.78	142	320017	37.62	ng	100
39) 1,2,4,5-Tetrachlorobenzene	8.96	216	124224	37.47	ng	99
40) Hexachlorocyclopentadiene	8.94	237	169199	93.53	ng	98

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42) 2,4,6-Trichlorophenol	9.07	196	88741	36.52	ng	94
43) 2,4,5-Trichlorophenol	9.10	196	103103	42.11	ng #	86
45) 1,1'-Biphenyl	9.25	154	391981	42.47	ng	99
46) 2-Chloronaphthalene	9.28	162	281592	39.13	ng	99
47) 2-Nitroaniline	9.37	65	111130	38.52	ng	97
48) Acenaphthylene	9.69	152	480695	38.43	ng	100
49) Dimethylphthalate	9.56	163	329272	36.86	ng #	98
50) 2,6-Dinitrotoluene	9.62	165	71906	38.96	ng #	46
51) Acenaphthene	9.86	154	326773	42.65	ng	99
52) 3-Nitroaniline	9.78	138	55271	24.29	ng	99
53) 2,4-Dinitrophenol	9.88	184	78565	70.84	ng #	70
54) Dibenzofuran	10.03	168	404098	39.01	ng	99
55) 4-Nitrophenol	9.95	139	139160	81.71	ng #	73
56) 2,4-Dinitrotoluene	10.02	165	107628	43.59	ng #	72
57) Fluorene	10.37	166	325330	40.60	ng	100
58) 2,3,4,6-Tetrachlorophenol	10.14	232	71412m	36.85	ng	
59) Diethylphthalate	10.26	149	349786	37.62	ng	96
60) 4-Chlorophenyl-phenylether	10.36	204	127838	32.74	ng	96
61) 4-Nitroaniline	10.40	138	80481	34.38	ng	96
62) Azobenzene	10.52	77	364879	35.09	ng	98
64) 4,6-Dinitro-2-methylphenol	10.42	198	47909	38.83	ng #	70
65) n-Nitrosodiphenylamine	10.49	169	281761	40.74	ng	99
66) 4-Bromophenyl-phenylether	10.85	248	91164	43.24	ng	97
67) Hexachlorobenzene	10.91	284	95542	40.76	ng #	56
68) Atrazine	11.01	200	83011	40.94	ng	98
69) Pentachlorophenol	11.10	266	106053	77.80	ng	99
70) Phenanthrene	11.32	178	405291	35.95	ng	99
71) Anthracene	11.38	178	452264	40.52	ng	99
72) Carbazole	11.53	167	412233	37.94	ng #	97
73) Di-n-butylphthalate	11.87	149	569725	41.93	ng	99
74) Fluoranthene	12.51	202	496464	40.74	ng	96
76) Benzidine	12.64	184	371074	52.04	ng	97
77) Pyrene	12.74	202	519382	38.86	ng	99
79) Butylbenzylphthalate	13.36	149	212695	33.01	ng #	67
80) Benzo(a)anthracene	13.92	228	436639	38.21	ng	99
81) 3,3'-Dichlorobenzidine	13.88	252	79092	19.03	ng #	98
82) Chrysene	13.95	228	376942	34.45	ng	99
83) Bis(2-ethylhexyl)phthalate	13.92	149	303872	33.89	ng #	96
84) Di-n-octyl phthalate	14.53	149	521627	34.32	ng	98
85) Indeno(1,2,3-cd)pyrene	16.67	276	388141	37.27	ng	99
87) Benzo(b)fluoranthene	14.93	252	443002m	37.59	ng	
88) Benzo(k)fluoranthene	14.96	252	322354m	31.27	ng	
89) Benzo(a)pyrene	15.27	252	402603	39.75	ng #	95
90) Dibenzo(a,h)anthracene	16.68	278	327428	36.55	ng	99
91) Benzo(g,h,i)perylene	17.07	276	290456	33.34	ng	96

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

