

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF063015\
 Data File : BF080241.D
 Acq On : 30 Jun 2015 17:34
 Operator : TP/IZ
 Sample : PB84281BS
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB84281BS

Manual Integrations
 APPROVED

apatel
 7/1/2015 4:16:36 PM

Quant Time: Jul 01 11:09:51 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF061715.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 23 00:53:56 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.28	152	36096	20.00	ng	-0.05
21) Naphthalene-d8	8.57	136	147040	20.00	ng	-0.05
38) Acenaphthene-d10	10.34	164	75327	20.00	ng	-0.05
63) Phenanthrene-d10	11.85	188	154981	20.00	ng	-0.06
75) Chrysene-d12	14.83	240	160775	20.00	ng	-0.17
86) Perylene-d12	16.70	264	153572	20.00	ng	-0.23

System Monitoring Compounds

5) 2-Fluorophenol	5.90	112	240284	117.84	ng	-0.03
7) Phenol-d6	6.89	99	337585	124.41	ng	-0.03
23) Nitrobenzene-d5	7.84	82	195145	77.26	ng	-0.03
41) 2,4,6-Tribromophenol	11.14	330	86660	117.65	ng	-0.03
44) 2-Fluorobiphenyl	9.65	172	375373	71.07	ng	-0.05
78) Terphenyl-d14	13.54	244	475007	69.00	ng	-0.13

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.13	88	28450m	29.35	ng	
3) Pyridine	3.97	79	78244	30.36	ng	# 81
4) n-Nitrosodimethylamine	3.86	42	43019	35.12	ng	96
6) Aniline	6.94	93	101863	27.29	ng	95
8) 2-Chlorophenol	7.06	128	89086	37.80	ng	87
9) Benzaldehyde	6.82	77	45500	29.05	ng	# 66
10) Phenol	6.90	94	123264	41.63	ng	81
11) bis(2-Chloroethyl)ether	7.01	93	80224	34.15	ng	87
12) 1,3-Dichlorobenzene	7.22	146	95487m	33.73	ng	
13) 1,4-Dichlorobenzene	7.30	146	103123	38.30	ng	96
14) 1,2-Dichlorobenzene	7.46	146	90457	34.52	ng	98
15) Benzyl Alcohol	7.41	79	74346m	33.51	ng	
16) 2,2'-oxybis(1-Chloropropan	7.56	45	138811	39.81	ng	85
17) 2-Methylphenol	7.52	107	70189	34.87	ng	# 89
18) Hexachloroethane	7.81	117	33914	37.83	ng	# 79
19) n-Nitroso-di-n-propylamine	7.68	70	62973	33.43	ng	# 77
20) 3+4-Methylphenols	7.67	107	97104	37.38	ng	91
22) Acetophenone	7.68	105	125341	36.58	ng	# 82
24) Nitrobenzene	7.86	77	91027	34.92	ng	# 78
25) Isophorone	8.09	82	164060	32.28	ng	# 83
26) 2-Nitrophenol	8.18	139	44768	41.03	ng	# 79
27) 2,4-Dimethylphenol	8.21	122	88271	38.79	ng	# 84
28) bis(2-Chloroethoxy)methane	8.30	93	106730	35.74	ng	# 90
29) 2,4-Dichlorophenol	8.42	162	80635	41.38	ng	98
30) 1,2,4-Trichlorobenzene	8.52	180	85032	38.43	ng	100
31) Naphthalene	8.60	128	262214m	34.11	ng	
32) Benzoic acid	8.28	122	33193	24.13	ng	# 70
33) 4-Chloroaniline	8.63	127	57383	17.44	ng	# 85
34) Hexachlorobutadiene	8.72	225	43655	32.04	ng	99
35) Caprolactam	8.97	113	22391	35.40	ng	86
36) 4-Chloro-3-methylphenol	9.11	107	76517	34.81	ng	93
37) 2-Methylnaphthalene	9.29	142	190828	38.37	ng	# 92
39) 1,2,4,5-Tetrachlorobenzene	9.46	216	88811	40.52	ng	97
40) Hexachlorocyclopentadiene	9.45	237	80401	71.96	ng	97

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.57	196	59436	39.85	ng	99
43) 2,4,5-Trichlorophenol	9.60	196	56578	32.92	ng #	89
45) 1,1'-Biphenyl	9.76	154	206641	33.72	ng	95
46) 2-Chloronaphthalene	9.78	162	175503	35.29	ng	94
47) 2-Nitroaniline	9.86	65	46761	34.26	ng #	57
48) Acenaphthylene	10.21	152	303116	36.52	ng	98
49) Dimethylphthalate	10.05	163	202692	35.79	ng #	97
50) 2,6-Dinitrotoluene	10.10	165	39928	35.17	ng #	43
51) Acenaphthene	10.38	154	193149	38.04	ng	90
52) 3-Nitroaniline	10.28	138	30588	23.35	ng #	49
53) 2,4-Dinitrophenol	10.39	184	42027	90.15	ng #	32
54) Dibenzofuran	10.55	168	255873	35.51	ng #	89
55) 4-Nitrophenol	10.42	139	68978	65.88	ng #	32
56) 2,4-Dinitrotoluene	10.52	165	58759	40.79	ng #	79
57) Fluorene	10.90	166	200063	34.99	ng	92
58) 2,3,4,6-Tetrachlorophenol	10.66	232	48365	33.34	ng #	98
59) Diethylphthalate	10.74	149	189350	33.21	ng	96
60) 4-Chlorophenyl-phenylether	10.88	204	98616	35.89	ng #	88
61) 4-Nitroaniline	10.89	138	45840	33.88	ng #	50
62) Azobenzene	11.04	77	200045	34.46	ng	86
64) 4,6-Dinitro-2-methylphenol	10.93	198	24937	35.61	ng	77
65) n-Nitrosodiphenylamine	11.00	169	176927	35.06	ng	93
66) 4-Bromophenyl-phenylether	11.38	248	58430	35.46	ng #	85
67) Hexachlorobenzene	11.46	284	62273	34.00	ng #	78
68) Atrazine	11.52	200	57620	36.14	ng	83
69) Pentachlorophenol	11.65	266	62683	60.40	ng	98
70) Phenanthrene	11.88	178	305114	32.52	ng	97
71) Anthracene	11.93	178	329975	36.52	ng	98
72) Carbazole	12.08	167	286993	33.27	ng	96
73) Di-n-butylphthalate	12.41	149	333186	33.43	ng #	99
74) Fluoranthene	13.14	202	341365	34.16	ng	89
76) Benzidine	13.26	184	209315	46.58	ng	98
77) Pyrene	13.40	202	347129	35.07	ng	97
79) Butylbenzylphthalate	14.09	149	135416	34.07	ng #	82
80) Benzo(a)anthracene	14.81	228	340928	34.81	ng	97
81) 3,3'-Dichlorobenzidine	14.76	252	82606	24.45	ng #	95
82) Chrysene	14.86	228	316631	35.06	ng	95
83) Bis(2-ethylhexyl)phthalate	14.79	149	208957	33.26	ng #	93
84) Di-n-octyl phthalate	15.60	149	335896	31.20	ng	97
85) Indeno(1,2,3-cd)pyrene	18.52	276	367281	32.25	ng #	88
87) Benzo(h)fluoranthene	16.16	252	343589	36.95	ng #	88
88) Benzo(k)fluoranthene	16.20	252	308210	32.55	ng #	89
89) Benzo(a)pyrene	16.62	252	323690	36.66	ng #	87
90) Dibenzo(a,h)anthracene	18.54	278	310461	34.35	ng #	82
91) Benzo(g,h,i)perylene	19.07	276	310301	34.00	ng #	78

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

