

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF092915\
 Data File : BF081881.D
 Acq On : 29 Sep 2015 15:55
 Operator : UM/IZ
 Sample : PB85717BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB85717BS

Manual Integrations
 APPROVED

apatel
 9/30/2015 9:09:16 AM

Quant Time: Sep 30 01:50:14 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092815.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Sep 28 18:10:42 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.93	152	84766	20.00	ng	-0.01
21) Naphthalene-d8	8.21	136	321323	20.00	ng	-0.01
38) Acenaphthene-d10	9.97	164	168726	20.00	ng	-0.01
63) Phenanthrene-d10	11.45	188	420682	20.00	ng	-0.01
75) Chrysene-d12	14.10	240	412183	20.00	ng	-0.01
86) Perylene-d12	15.57	264	294439	20.00	ng	-0.01

System Monitoring Compounds						
5) 2-Fluorophenol	5.52	112	504565	108.06	ng	0.00
7) Phenol-d6	6.55	99	614117	104.52	ng	-0.01
23) Nitrobenzene-d5	7.49	82	399443	82.87	ng	-0.01
41) 2,4,6-Tribromophenol	10.77	330	234282	137.10	ng	0.00
44) 2-Fluorobiphenyl	9.29	172	862640	85.69	ng	-0.01
78) Terphenyl-d14	13.04	244	1019663	78.14	ng	-0.01

Target Compounds						Qvalue
2) 1,4-Dioxane	2.59	88	50279	24.76	ng	92
3) Pyridine	3.35	79	122677	23.20	ng	90
4) n-Nitrosodimethylamine	3.30	42	65814	27.41	ng	95
6) Aniline	6.59	93	191458	24.25	ng	# 89
8) 2-Chlorophenol	6.71	128	188541	36.56	ng	99
9) Benzaldehyde	6.47	77	55573	14.44	ng	# 76
10) Phenol	6.56	94	207847	31.53	ng	# 52
11) bis(2-Chloroethyl)ether	6.66	93	193946	37.94	ng	94
12) 1,3-Dichlorobenzene	6.86	146	209406	34.38	ng	# 91
13) 1,4-Dichlorobenzene	6.94	146	226761m	35.65	ng	
14) 1,2-Dichlorobenzene	7.10	146	204868	36.15	ng	99
15) Benzyl Alcohol	7.07	79	151699	35.77	ng	# 74
16) 2,2'-oxybis(1-Chloropropan	7.20	45	259601	35.93	ng	76
17) 2-Methylphenol	7.18	107	152502	36.48	ng	# 83
18) Hexachloroethane	7.43	117	73189	36.76	ng	90
19) n-Nitroso-di-n-propylamine	7.34	70	133079	37.26	ng	# 78
20) 3+4-Methylphenols	7.33	107	187904	35.58	ng	# 67
22) Acetophenone	7.34	105	266022	40.49	ng	# 81
24) Nitrobenzene	7.51	77	202400	38.67	ng	# 68
25) Isophorone	7.75	82	363623	38.06	ng	# 93
26) 2-Nitrophenol	7.83	139	100435	39.97	ng	# 75
27) 2,4-Dimethylphenol	7.86	122	172581	39.05	ng	# 79
28) bis(2-Chloroethoxy)methane	7.97	93	216211	38.11	ng	# 95
29) 2,4-Dichlorophenol	8.07	162	176092	41.09	ng	98
30) 1,2,4-Trichlorobenzene	8.15	180	185807	38.83	ng	99
31) Naphthalene	8.23	128	551082	38.71	ng	98
32) Benzoic acid	7.98	122	81451	33.44	ng	# 66
33) 4-Chloroaniline	8.29	127	128349	20.85	ng	# 92
34) Hexachlorobutadiene	8.34	225	109328	38.64	ng	99
35) Caprolactam	8.65	113	48806m	41.13	ng	
36) 4-Chloro-3-methylphenol	8.77	107	179961	42.94	ng	86
37) 2-Methylnaphthalene	8.93	142	389585	40.09	ng	# 90
39) 1,2,4,5-Tetrachlorobenzene	9.09	216	183957	35.51	ng	98
40) Hexachlorocyclopentadiene	9.07	237	239065	89.62	ng	97

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.20	196	123957	34.91	ng	99
43) 2,4,5-Trichlorophenol	9.25	196	154151	42.21	ng	97
45) 1,1'-Biphenyl	9.39	154	476266	36.59	ng	95
46) 2-Chloronaphthalene	9.42	162	399738	39.11	ng	96
47) 2-Nitroaniline	9.52	65	120365	40.12	ng	89
48) Acenaphthylene	9.83	152	603624	36.90	ng	96
49) Dimethylphthalate	9.69	163	489499	42.03	ng	# 98
50) 2,6-Dinitrotoluene	9.76	165	104881	41.48	ng	96
51) Acenaphthene	10.00	154	417234	42.95	ng	89
52) 3-Nitroaniline	9.93	138	72613	24.82	ng	# 77
53) 2,4-Dinitrophenol	10.03	184	98515	79.98	ng	# 72
54) Dibenzofuran	10.17	168	523339	38.12	ng	# 84
55) 4-Nitrophenol	10.09	139	181406	85.82	ng	# 61
56) 2,4-Dinitrotoluene	10.16	165	148983	46.22	ng	# 89
57) Fluorene	10.51	166	428111	43.53	ng	91
58) 2,3,4,6-Tetrachlorophenol	10.29	232	127470	44.00	ng	94
59) Diethylphthalate	10.39	149	468264	43.04	ng	97
60) 4-Chlorophenyl-phenylether	10.51	204	206660	41.12	ng	95
61) 4-Nitroaniline	10.55	138	112068	38.21	ng	# 44
62) Azobenzene	10.67	77	474912	44.73	ng	96
64) 4,6-Dinitro-2-methylphenol	10.56	198	62793	35.19	ng	# 77
65) n-Nitrosodiphenylamine	10.63	169	385121	30.26	ng	90
66) 4-Bromophenyl-phenylether	11.01	248	159369	37.57	ng	# 88
67) Hexachlorobenzene	11.06	284	181515	37.92	ng	# 81
68) Atrazine	11.15	200	149798	37.93	ng	84
69) Pentachlorophenol	11.26	266	215799	79.12	ng	99
70) Phenanthrene	11.49	178	836051	41.15	ng	96
71) Anthracene	11.53	178	856138	40.04	ng	96
72) Carbazole	11.69	167	792903	40.97	ng	95
73) Di-n-butylphthalate	12.01	149	893762	45.36	ng	# 97
74) Fluoranthene	12.67	202	897139	39.81	ng	86
76) Benzidine	12.80	184	325985	26.25	ng	98
77) Pyrene	12.90	202	899414	36.53	ng	96
79) Butylbenzylphthalate	13.52	149	430169	46.43	ng	98
80) Benzo(a)anthracene	14.09	228	859325	38.72	ng	96
81) 3,3'-Dichlorobenzidine	14.06	252	204561	27.29	ng	# 94
82) Chrysene	14.13	228	686438m	30.57	ng	
83) Bis(2-ethylhexyl)phthalate	14.07	149	570011	49.04	ng	# 89
84) Di-n-octyl phthalate	14.69	149	888110	46.95	ng	# 93
85) Indeno(1,2,3-cd)pyrene	17.03	276	712104	41.01	ng	# 86
87) Benzo(b)fluoranthene	15.14	252	748209	44.51	ng	# 85
88) Benzo(k)fluoranthene	15.17	252	674574m	43.78	ng	
89) Benzo(a)pyrene	15.51	252	602748	41.34	ng	# 85
90) Dibenzo(a,h)anthracene	17.05	278	591319	48.33	ng	# 79
91) Benzo(g,h,i)perylene	17.48	276	566463	45.18	ng	# 72

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

