

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF090415\
 Data File : BF081425.D
 Acq On : 4 Sep 2015 18:18
 Operator : UM/IZ
 Sample : PB85409BSD
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB85409BSD

Manual Integrations
 APPROVED

MMDadoda
 9/8/2015 1:57:04 PM

Quant Time: Sep 07 01:32:34 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF090115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Sep 07 01:02:00 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.35	152	65497	20.00	ng	-0.01
21) Naphthalene-d8	7.64	136	259719	20.00	ng	-0.01
38) Acenaphthene-d10	9.39	164	113193	20.00	ng	0.00
63) Phenanthrene-d10	10.86	188	225955	20.00	ng	0.00
75) Chrysene-d12	13.46	240	173897	20.00	ng	0.00
86) Perylene-d12	14.77	264	111966	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	4.92	112	443206	104.99	ng	0.02
7) Phenol-d6	6.03	99	579090	103.63	ng	0.00
23) Nitrobenzene-d5	6.94	82	356929	66.30	ng	-0.01
41) 2,4,6-Tribromophenol	10.18	330	115938	115.03	ng	0.00
44) 2-Fluorobiphenyl	8.73	172	543484	61.53	ng	-0.01
78) Terphenyl-d14	12.43	244	496752	66.50	ng	-0.01

Target Compounds

						Qvalue
2) 1,4-Dioxane	1.77	88	55260	29.15	ng	# 89
3) Pyridine	2.31	79	173594	33.20	ng	# 89
4) n-Nitrosodimethylamine	2.27	42	106276	36.59	ng	# 78
6) Aniline	6.02	93	180423	22.87	ng	# 80
8) 2-Chlorophenol	6.15	128	170600	36.67	ng	# 87
9) Benzaldehyde	5.90	77	119126	34.02	ng	# 77
10) Phenol	6.04	94	199269	30.75	ng	# 44
11) bis(2-Chloroethyl)ether	6.11	93	165139	35.32	ng	# 86
12) 1,3-Dichlorobenzene	6.30	146	173692	34.95	ng	# 92
13) 1,4-Dichlorobenzene	6.38	146	168140	33.23	ng	# 89
14) 1,2-Dichlorobenzene	6.52	146	145938	32.03	ng	# 86
15) Benzyl Alcohol	6.52	79	153861	33.57	ng	# 71
16) 2,2'-oxybis(1-Chloropropan	6.66	45	334729	37.35	ng	# 1
17) 2-Methylphenol	6.66	107	136622	36.81	ng	# 74
18) Hexachloroethane	6.87	117	67149	34.11	ng	# 95
19) n-Nitroso-di-n-propylamine	6.80	70	127097	33.42	ng	# 87
20) 3+4-Methylphenols	6.82	107	179524	35.87	ng	# 89
22) Acetophenone	6.79	105	240095	33.85	ng	# 76
24) Nitrobenzene	6.96	77	215064	38.47	ng	# 67
25) Isophorone	7.20	82	346985	34.65	ng	# 82
26) 2-Nitrophenol	7.28	139	86329	35.43	ng	# 55
27) 2,4-Dimethylphenol	7.35	122	144669	34.38	ng	# 83
28) bis(2-Chloroethoxy)methane	7.43	93	191517	33.11	ng	# 92
29) 2,4-Dichlorophenol	7.53	162	136031	37.28	ng	# 97
30) 1,2,4-Trichlorobenzene	7.60	180	141540	35.70	ng	# 97
31) Naphthalene	7.67	128	452904	32.70	ng	# 99
32) Benzoic acid	7.50	122	73822	25.81	ng	# 75
33) 4-Chloroaniline	7.74	127	93880	16.62	ng	# 84
34) Hexachlorobutadiene	7.80	225	82347	34.13	ng	# 97
35) Caprolactam	8.11	113	41792m	37.34	ng	# 80
36) 4-Chloro-3-methylphenol	8.25	107	149006	36.48	ng	# 90
37) 2-Methylnaphthalene	8.36	142	303089	33.63	ng	# 99
39) 1,2,4,5-Tetrachlorobenzene	8.52	216	115090	32.15	ng	# 99
40) Hexachlorocyclopentadiene	8.51	237	119025	67.75	ng	# 95

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF090415\
 Data File : BF081425.D
 Acq On : 4 Sep 2015 18:18
 Operator : UM/IZ
 Sample : PB85409BSD
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB85409BSD

Manual Integrations
 APPROVED

MMDadoda
 9/8/2015 1:57:04 PM

Quant Time: Sep 07 01:32:34 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF090115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Sep 07 01:02:00 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	8.65	196	83081	33.63	ng	98
43) 2,4,5-Trichlorophenol	8.70	196	92406	36.66	ng #	90
45) 1,1'-Biphenyl	8.83	154	344812	33.11	ng	94
46) 2-Chloronaphthalene	8.84	162	260266	34.61	ng #	90
47) 2-Nitroaniline	8.96	65	110661	36.10	ng #	79
48) Acenaphthylene	9.26	152	446006	33.43	ng	97
49) Dimethylphthalate	9.15	163	305889	34.78	ng #	96
50) 2,6-Dinitrotoluene	9.20	165	58126	29.12	ng #	11
51) Acenaphthene	9.43	154	280401	36.94	ng	90
52) 3-Nitroaniline	9.37	138	42085	18.52	ng #	85
53) 2,4-Dinitrophenol	9.47	184	55555	66.08	ng #	1
54) Dibenzofuran	9.60	168	427726	38.59	ng #	90
55) 4-Nitrophenol	9.56	139	114518	80.22	ng #	23
56) 2,4-Dinitrotoluene	9.60	165	83788	32.97	ng #	14
57) Fluorene	9.93	166	300028	35.67	ng	96
58) 2,3,4,6-Tetrachlorophenol	9.72	232	76957	39.99	ng	92
59) Diethylphthalate	9.84	149	325116	35.14	ng	95
60) 4-Chlorophenyl-phenylether	9.94	204	151919	38.76	ng	99
61) 4-Nitroaniline	9.98	138	78313	34.11	ng #	33
62) Azobenzene	10.09	77	372234	36.19	ng #	80
64) 4,6-Dinitro-2-methylphenol	10.01	198	43369	34.32	ng #	75
65) n-Nitrosodiphenylamine	10.06	169	265610	32.30	ng	91
66) 4-Bromophenyl-phenylether	10.42	248	88856	38.89	ng	93
67) Hexachlorobenzene	10.48	284	82954	34.73	ng #	77
68) Atrazine	10.60	200	83078	34.92	ng	84
69) Pentachlorophenol	10.67	266	78112	68.57	ng	99
70) Phenanthrene	10.88	178	434855	34.62	ng	97
71) Anthracene	10.92	178	412945	32.00	ng	98
72) Carbazole	11.10	167	434956	35.91	ng	96
73) Di-n-butylphthalate	11.45	149	517228	36.17	ng #	98
74) Fluoranthene	12.05	202	444256	32.67	ng	92
76) Benzidine	12.19	184	182592	24.46	ng	98
77) Pyrene	12.27	202	489771	38.02	ng	99
79) Butylbenzylphthalate	12.93	149	204764	36.55	ng	93
80) Benzo(a)anthracene	13.45	228	376491	34.00	ng	97
81) 3,3'-Dichlorobenzidine	13.43	252	68866	18.44	ng #	95
82) Chrysene	13.49	228	280918	28.30	ng	95
83) Bis(2-ethylhexyl)phthalate	13.49	149	236524	31.99	ng #	88
84) Di-n-octyl phthalate	14.09	149	314441	25.83	ng	98
85) Indeno(1,2,3-cd)pyrene	15.84	276	248882	34.17	ng #	87
87) Benzo(b)fluoranthene	14.45	252	316859m	39.64	ng	
88) Benzo(k)fluoranthene	14.47	252	178874m	26.97	ng	
89) Benzo(a)pyrene	14.72	252	238744	35.25	ng #	85
90) Dibenzo(a,h)anthracene	15.85	278	211393	40.39	ng #	81
91) Benzo(g,h,i)perylene	16.16	276	211716	42.38	ng #	73

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

