

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF090815\
 Data File : BF081523.D
 Acq On : 8 Sep 2015 15:31
 Operator : UM/IZ
 Sample : PB85437BSD
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB85437BSD

Manual Integrations
 APPROVED

MMDadoda
 9/9/2015 2:57:41 PM

Quant Time: Sep 09 00:27:08 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF090115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 08 23:43:58 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.34	152	47990	20.00	ng	0.00
21) Naphthalene-d8	7.63	136	184424	20.00	ng	-0.01
38) Acenaphthene-d10	9.38	164	93015	20.00	ng	0.00
63) Phenanthrene-d10	10.84	188	194870	20.00	ng	0.00
75) Chrysene-d12	13.45	240	161303	20.00	ng	0.00
86) Perylene-d12	14.75	264	99166	20.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	4.90	112	224321	72.52	ng	0.02
7) Phenol-d6	6.02	99	322342	78.73	ng	0.00
23) Nitrobenzene-d5	6.92	82	266755	69.79	ng	0.00
41) 2,4,6-Tribromophenol	10.16	330	57001	68.82	ng	0.00
44) 2-Fluorobiphenyl	8.72	172	469502	64.68	ng	0.00
78) Terphenyl-d14	12.42	244	441832	63.76	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	1.75	88	40841	29.40	ng	# 79
3) Pyridine	2.27	79	130065	33.95	ng	# 75
4) n-Nitrosodimethylamine	2.24	42	80539	37.85	ng	# 67
6) Aniline	6.01	93	139648	24.15	ng	# 81
8) 2-Chlorophenol	6.14	128	123770	36.31	ng	# 96
9) Benzaldehyde	5.88	77	71472	27.86	ng	# 83
10) Phenol	6.03	94	191331	40.30	ng	# 60
11) bis(2-Chloroethyl)ether	6.10	93	133692	39.03	ng	# 94
12) 1,3-Dichlorobenzene	6.28	146	131451	36.10	ng	# 93
13) 1,4-Dichlorobenzene	6.36	146	137194	37.00	ng	# 89
14) 1,2-Dichlorobenzene	6.51	146	115236m	34.52	ng	# 89
15) Benzyl Alcohol	6.51	79	119243	35.50	ng	# 71
16) 2,2'-oxybis(1-Chloropropan	6.65	45	260406	39.66	ng	# 12
17) 2-Methylphenol	6.65	107	103731	38.14	ng	# 76
18) Hexachloroethane	6.86	117	52370	36.30	ng	# 90
19) n-Nitroso-di-n-propylamine	6.79	70	95811	34.38	ng	# 89
20) 3+4-Methylphenols	6.81	107	138815	37.86	ng	# 88
22) Acetophenone	6.78	105	185706	36.87	ng	# 76
24) Nitrobenzene	6.95	77	160524	40.44	ng	# 68
25) Isophorone	7.19	82	260838	36.69	ng	# 82
26) 2-Nitrophenol	7.27	139	66307	38.32	ng	# 55
27) 2,4-Dimethylphenol	7.34	122	107540	35.99	ng	# 85
28) bis(2-Chloroethoxy)methane	7.42	93	150931	36.75	ng	# 90
29) 2,4-Dichlorophenol	7.52	162	99680	38.47	ng	# 98
30) 1,2,4-Trichlorobenzene	7.59	180	108272	38.46	ng	# 98
31) Naphthalene	7.66	128	346523	35.23	ng	# 98
32) Benzoic acid	7.47	122	47354	23.56	ng	# 61
33) 4-Chloroaniline	7.72	127	76603	19.09	ng	# 83
34) Hexachlorobutadiene	7.78	225	63762	37.22	ng	# 97
35) Caprolactam	8.10	113	31024m	39.04	ng	# 84
36) 4-Chloro-3-methylphenol	8.24	107	121036	41.73	ng	# 90
37) 2-Methylnaphthalene	8.35	142	256520	40.08	ng	# 98
39) 1,2,4,5-Tetrachlorobenzene	8.51	216	95779	32.56	ng	# 98
40) Hexachlorocyclopentadiene	8.50	237	99678	69.05	ng	# 95

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF090815\
 Data File : BF081523.D
 Acq On : 8 Sep 2015 15:31
 Operator : UM/IZ
 Sample : PB85437BSD
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB85437BSD

Manual Integrations
 APPROVED

MMDadoda
 9/9/2015 2:57:41 PM

Quant Time: Sep 09 00:27:08 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF090115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 08 23:43:58 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	8.64	196	68553	33.77	ng	99
43) 2,4,5-Trichlorophenol	8.68	196	78765	38.03	ng #	91
45) 1,1'-Biphenyl	8.82	154	277132	32.38	ng	95
46) 2-Chloronaphthalene	8.83	162	225538	36.50	ng #	91
47) 2-Nitroaniline	8.95	65	93294	37.04	ng #	78
48) Acenaphthylene	9.23	152	358521	32.70	ng	97
49) Dimethylphthalate	9.13	163	251166	34.75	ng #	97
50) 2,6-Dinitrotoluene	9.19	165	50005	30.49	ng #	5
51) Acenaphthene	9.42	154	213329	34.20	ng	92
52) 3-Nitroaniline	9.35	138	37941	20.32	ng #	29
53) 2,4-Dinitrophenol	9.47	184	50935	73.73	ng #	76
54) Dibenzofuran	9.59	168	322447	35.41	ng	92
55) 4-Nitrophenol	9.55	139	92595	78.93	ng #	33
56) 2,4-Dinitrotoluene	9.59	165	69087	33.08	ng #	18
57) Fluorene	9.92	166	247933	35.87	ng	97
58) 2,3,4,6-Tetrachlorophenol	9.71	232	66142	41.83	ng	96
59) Diethylphthalate	9.83	149	281436	37.02	ng	96
60) 4-Chlorophenyl-phenylether	9.93	204	118387	36.75	ng	92
61) 4-Nitroaniline	9.96	138	65833	34.90	ng #	29
62) Azobenzene	10.08	77	312195	36.94	ng	81
64) 4,6-Dinitro-2-methylphenol	10.00	198	36962	33.91	ng	84
65) n-Nitrosodiphenylamine	10.04	169	215227	30.35	ng	91
66) 4-Bromophenyl-phenylether	10.41	248	71857	36.47	ng #	85
67) Hexachlorobenzene	10.47	284	69619	33.79	ng #	75
68) Atrazine	10.58	200	67718	33.00	ng	81
69) Pentachlorophenol	10.66	266	66274	67.60	ng	99
70) Phenanthrene	10.87	178	401398	37.05	ng	96
71) Anthracene	10.91	178	356799	32.06	ng	98
72) Carbazole	11.08	167	390125	37.35	ng	97
73) Di-n-butylphthalate	11.44	149	489127	39.66	ng #	99
74) Fluoranthene	12.03	202	387303	33.02	ng	92
76) Benzidine	12.18	184	199645	28.83	ng	97
77) Pyrene	12.26	202	456804	38.23	ng	98
79) Butylbenzylphthalate	12.91	149	194243	37.38	ng	93
80) Benzo(a)anthracene	13.44	228	354146	34.48	ng	98
81) 3,3'-Dichlorobenzidine	13.42	252	56378	16.27	ng #	96
82) Chrysene	13.47	228	262537	28.51	ng	94
83) Bis(2-ethylhexyl)phthalate	13.47	149	228030	33.25	ng #	89
84) Di-n-octyl phthalate	14.08	149	339048	30.03	ng #	94
85) Indeno(1,2,3-cd)pyrene	15.83	276	207204	30.67	ng #	85
87) Benzo(b)fluoranthene	14.43	252	307041m	43.37	ng	
88) Benzo(k)fluoranthene	14.46	252	214865m	36.58	ng	
89) Benzo(a)pyrene	14.71	252	219091	36.52	ng #	86
90) Dibenzo(a,h)anthracene	15.84	278	174903	37.73	ng #	78
91) Benzo(g,h,i)perylene	16.14	276	173246	39.16	ng #	76

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

