

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF081915\
 Data File : BF081088.D
 Acq On : 19 Aug 2015 18:49
 Operator : UM/IZ
 Sample : PB85067BS
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB85067BS

Manual Integrations
 APPROVED

MMDadoda
 8/20/2015 3:41:22 PM

Quant Time: Aug 20 01:14:57 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF081715.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 17 17:57:25 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.58	152	60298	20.00	ng	-0.04
21) Naphthalene-d8	7.88	136	247716	20.00	ng	-0.03
38) Acenaphthene-d10	9.62	164	117937	20.00	ng	-0.03
63) Phenanthrene-d10	11.09	188	223923	20.00	ng	-0.04
75) Chrysene-d12	13.72	240	174261	20.00	ng	-0.03
86) Perylene-d12	15.04	264	164060	20.00	ng	-0.05

System Monitoring Compounds						
5) 2-Fluorophenol	5.17	112	427904	106.74	ng	-0.02
7) Phenol-d6	6.24	99	608260	116.29	ng	-0.02
23) Nitrobenzene-d5	7.16	82	401649	74.07	ng	-0.03
41) 2,4,6-Tribromophenol	10.41	330	127536	124.75	ng	-0.03
44) 2-Fluorobiphenyl	8.96	172	682872	75.87	ng	-0.03
78) Terphenyl-d14	12.68	244	606686	74.64	ng	-0.03

Target Compounds				Qvalue		
2) 1,4-Dioxane	2.08	88	54794	29.01	ng	95
3) Pyridine	2.70	79	182111	34.15	ng	# 82
4) n-Nitrosodimethylamine	2.65	42	102848	34.64	ng	# 79
6) Aniline	6.25	93	163717	21.53	ng	# 1
8) 2-Chlorophenol	6.37	128	168198	38.33	ng	91
9) Benzaldehyde	6.13	77	57061	16.92	ng	94
10) Phenol	6.25	94	254867	41.26	ng	88
11) bis(2-Chloroethyl)ether	6.33	93	169008	35.65	ng	91
12) 1,3-Dichlorobenzene	6.53	146	173629	36.83	ng	95
13) 1,4-Dichlorobenzene	6.61	146	182693	39.13	ng	97
14) 1,2-Dichlorobenzene	6.76	146	152905	34.99	ng	97
15) Benzyl Alcohol	6.74	79	162102	38.30	ng	95
16) 2,2'-oxybis(1-Chloropropan	6.88	45	344917	37.03	ng	100
17) 2-Methylphenol	6.87	107	148130	39.34	ng	93
18) Hexachloroethane	7.10	117	67434	35.75	ng	94
19) n-Nitroso-di-n-propylamine	7.02	70	132877	36.67	ng	92
20) 3+4-Methylphenols	7.02	107	173792	36.37	ng	# 59
22) Acetophenone	7.01	105	245074	37.68	ng	# 89
24) Nitrobenzene	7.18	77	203828	35.95	ng	98
25) Isophorone	7.42	82	365151	36.11	ng	99
26) 2-Nitrophenol	7.50	139	95541	40.52	ng	# 70
27) 2,4-Dimethylphenol	7.54	122	158062	37.89	ng	95
28) bis(2-Chloroethoxy)methane	7.65	93	232038	38.84	ng	98
29) 2,4-Dichlorophenol	7.74	162	129924	37.00	ng	89
30) 1,2,4-Trichlorobenzene	7.82	180	134373	34.43	ng	96
31) Naphthalene	7.90	128	518478	37.55	ng	99
32) Benzoic acid	7.68	122	127634	54.40	ng	92
33) 4-Chloroaniline	7.96	127	76347	13.10	ng	# 86
34) Hexachlorobutadiene	8.02	225	83717	37.93	ng	99
35) Caprolactam	8.33	113	55006m	44.81	ng	
36) 4-Chloro-3-methylphenol	8.45	107	163075	38.96	ng	99
37) 2-Methylnaphthalene	8.58	142	306431	35.13	ng	99
39) 1,2,4,5-Tetrachlorobenzene	8.76	216	132578	34.84	ng	# 97
40) Hexachlorocyclopentadiene	8.74	237	155402	78.90	ng	97

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	8.87	196	100955	37.55	ng	98
43) 2,4,5-Trichlorophenol	8.92	196	100824	37.53	ng	96
45) 1,1'-Biphenyl	9.05	154	379842	36.56	ng	97
46) 2-Chloronaphthalene	9.08	162	297583	35.66	ng	96
47) 2-Nitroaniline	9.18	65	128746	37.70	ng	# 64
48) Acenaphthylene	9.49	152	528885	35.43	ng	99
49) Dimethylphthalate	9.36	163	336207	34.61	ng	99
50) 2,6-Dinitrotoluene	9.42	165	70645	33.88	ng	# 67
51) Acenaphthene	9.66	154	316123	35.37	ng	99
52) 3-Nitroaniline	9.58	138	44970	17.23	ng	96
53) 2,4-Dinitrophenol	9.69	184	90781	78.32	ng	85
54) Dibenzofuran	9.83	168	459775	38.39	ng	97
55) 4-Nitrophenol	9.76	139	135169	73.75	ng	# 80
56) 2,4-Dinitrotoluene	9.82	165	92496	33.46	ng	# 70
57) Fluorene	10.16	166	322227	34.97	ng	99
58) 2,3,4,6-Tetrachlorophenol	9.94	232	77245m	37.15	ng	
59) Diethylphthalate	10.06	149	362024	35.21	ng	98
60) 4-Chlorophenyl-phenylether	10.16	204	142819	36.63	ng	94
61) 4-Nitroaniline	10.20	138	84565	31.71	ng	92
62) Azobenzene	10.32	77	444777	37.55	ng	97
64) 4,6-Dinitro-2-methylphenol	10.23	198	64218	48.02	ng	# 44
65) n-Nitrosodiphenylamine	10.29	169	309624	38.74	ng	99
66) 4-Bromophenyl-phenylether	10.65	248	89488	40.68	ng	# 87
67) Hexachlorobenzene	10.71	284	92728	41.62	ng	94
68) Atrazine	10.82	200	96916	43.52	ng	94
69) Pentachlorophenol	10.90	266	106311	79.53	ng	100
70) Phenanthrene	11.12	178	512931	38.65	ng	100
71) Anthracene	11.17	178	470697	36.45	ng	99
72) Carbazole	11.33	167	506433	40.73	ng	99
73) Di-n-butylphthalate	11.67	149	562944	37.42	ng	98
74) Fluoranthene	12.29	202	554378	38.71	ng	92
76) Benzidine	12.43	184	181637	27.59	ng	98
77) Pyrene	12.52	202	527443	37.23	ng	99
79) Butylbenzylphthalate	13.16	149	269468	44.25	ng	# 78
80) Benzo(a)anthracene	13.71	228	497772	42.59	ng	97
81) 3,3'-Dichlorobenzidine	13.67	252	113017	29.86	ng	98
82) Chrysene	13.74	228	479776	45.55	ng	100
83) Bis(2-ethylhexyl)phthalate	13.72	149	349556	44.82	ng	# 97
84) Di-n-octyl phthalate	14.32	149	605655	51.09	ng	99
85) Indeno(1,2,3-cd)pyrene	16.27	276	361169	48.84	ng	# 94
87) Benzo(b)fluoranthene	14.70	252	536289m	46.21	ng	
88) Benzo(k)fluoranthene	14.72	252	309759m	28.65	ng	
89) Benzo(a)pyrene	15.00	252	391078	38.68	ng	# 94
90) Dibenzo(a,h)anthracene	16.28	278	305177	38.73	ng	# 97
91) Benzo(g,h,i)perylene	16.62	276	290776	36.38	ng	# 92

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

