

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF081715\
 Data File : BF081007.D
 Acq On : 17 Aug 2015 17:04
 Operator : UM/IZ
 Sample : SSTDICC025
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC025

Manual Integrations
 APPROVED

MMDadoda
 8/18/2015 5:11:56 PM

Quant Time: Aug 17 17:45:49 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 17:27:51 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.63	152	55211	20.00	ng	0.00
21) Naphthalene-d8	7.92	136	217193	20.00	ng	0.01
38) Acenaphthene-d10	9.66	164	97126	20.00	ng	0.00
63) Phenanthrene-d10	11.13	188	198652	20.00	ng	0.00
75) Chrysene-d12	13.76	240	164334	20.00	ng	0.01
86) Perylene-d12	15.10	264	124774	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.19	112	186322	25.38	ng	0.00
7) Phenol-d6	6.26	99	233439	24.37	ng	0.00
23) Nitrobenzene-d5	7.20	82	258877	27.22	ng	0.01
41) 2,4,6-Tribromophenol	10.44	330	45314	26.91	ng	0.00
44) 2-Fluorobiphenyl	8.99	172	400243	27.00	ng	0.00
78) Terphenyl-d14	12.72	244	426871	27.84	ng	0.01

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.04	88	42549	24.60	ng	92
3) Pyridine	2.68	79	130377	26.70	ng	82
4) n-Nitrosodimethylamine	2.63	42	71955	26.47	ng	# 80
6) Aniline	6.28	93	182328	26.18	ng	# 91
8) 2-Chlorophenol	6.41	128	102392	25.48	ng	# 79
9) Benzaldehyde	6.17	77	86390	26.85	ng	84
10) Phenol	6.27	94	127405	22.52	ng	97
11) bis(2-Chloroethyl)ether	6.36	93	107572	24.78	ng	90
12) 1,3-Dichlorobenzene	6.56	146	107241m	24.84	ng	
13) 1,4-Dichlorobenzene	6.64	146	108483	25.38	ng	93
14) 1,2-Dichlorobenzene	6.80	146	108396	27.09	ng	92
15) Benzyl Alcohol	6.78	79	104277	26.91	ng	94
16) 2,2'-oxybis(1-Chloropropan	6.92	45	226949	26.61	ng	99
17) 2-Methylphenol	6.89	107	85311	24.75	ng	96
18) Hexachloroethane	7.14	117	44393	25.70	ng	92
19) n-Nitroso-di-n-propylamine	7.05	70	85283	25.71	ng	93
20) 3+4-Methylphenols	7.05	107	111094	25.39	ng	# 65
22) Acetophenone	7.04	105	140517	24.64	ng	# 94
24) Nitrobenzene	7.21	77	116307	23.40	ng	95
25) Isophorone	7.46	82	235621	26.57	ng	# 94
26) 2-Nitrophenol	7.53	139	52620	25.45	ng	# 85
27) 2,4-Dimethylphenol	7.58	122	87849	24.02	ng	97
28) bis(2-Chloroethoxy)methane	7.68	93	146583	27.98	ng	99
29) 2,4-Dichlorophenol	7.77	162	76114	24.72	ng	88
30) 1,2,4-Trichlorobenzene	7.86	180	89985	26.29	ng	98
31) Naphthalene	7.93	128	291792	24.10	ng	98
32) Benzoic acid	7.69	122	57136	27.77	ng	# 78
33) 4-Chloroaniline	7.99	127	124566	24.38	ng	96
34) Hexachlorobutadiene	8.06	225	47813	24.71	ng	99
35) Caprolactam	8.35	113	27775	25.81	ng	# 86
36) 4-Chloro-3-methylphenol	8.48	107	98866	26.94	ng	# 89
37) 2-Methylnaphthalene	8.63	142	215827	28.22	ng	96
39) 1,2,4,5-Tetrachlorobenzene	8.79	216	78116	24.93	ng	97
40) Hexachlorocyclopentadiene	8.79	237	41329	25.48	ng	96

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	8.90	196	56085	25.33	ng	96
43) 2,4,5-Trichlorophenol	8.95	196	57995	26.21	ng #	92
45) 1,1'-Biphenyl	9.08	154	210507	24.61	ng	94
46) 2-Chloronaphthalene	9.11	162	169513	24.66	ng #	92
47) 2-Nitroaniline	9.21	65	84674	30.11	ng #	84
48) Acenaphthylene	9.52	152	304244	24.75	ng	98
49) Dimethylphthalate	9.39	163	190973	23.87	ng	100
50) 2,6-Dinitrotoluene	9.45	165	41992	24.45	ng #	66
51) Acenaphthene	9.69	154	183842	24.98	ng	99
52) 3-Nitroaniline	9.62	138	63248	29.43	ng #	72
53) 2,4-Dinitrophenol	9.72	184	22120	27.30	ng #	57
54) Dibenzofuran	9.86	168	244242	24.76	ng	92
55) 4-Nitrophenol	9.78	139	43683	28.94	ng #	76
56) 2,4-Dinitrotoluene	9.85	165	58023	25.49	ng #	83
57) Fluorene	10.20	166	205239	27.05	ng	99
58) 2,3,4,6-Tetrachlorophenol	9.99	232	47258	27.60	ng	97
59) Diethylphthalate	10.09	149	212369	25.08	ng	99
60) 4-Chlorophenyl-phenylether	10.20	204	93284	27.84	ng	93
61) 4-Nitroaniline	10.23	138	65621	29.87	ng	85
62) Azobenzene	10.36	77	264926	27.16	ng	90
64) 4,6-Dinitro-2-methylphenol	10.26	198	28837	24.31	ng #	40
65) n-Nitrosodiphenylamine	10.32	169	169910	23.96	ng	98
66) 4-Bromophenyl-phenylether	10.68	248	46205	23.68	ng #	62
67) Hexachlorobenzene	10.75	284	49190	24.89	ng #	86
68) Atrazine	10.86	200	58203	29.46	ng	98
69) Pentachlorophenol	10.94	266	28541	24.07	ng	97
70) Phenanthrene	11.15	178	281780	23.93	ng	100
71) Anthracene	11.21	178	320693	27.99	ng	98
72) Carbazole	11.37	167	297503	26.97	ng	97
73) Di-n-butylphthalate	11.71	149	366078	27.43	ng	99
74) Fluoranthene	12.33	202	315506	24.83	ng	92
76) Benzidine	12.47	184	189883	30.59	ng	98
77) Pyrene	12.56	202	330838	24.76	ng	99
79) Butylbenzylphthalate	13.20	149	152556	26.57	ng #	84
80) Benzo(a)anthracene	13.75	228	283769	25.75	ng	98
81) 3,3'-Dichlorobenzidine	13.71	252	90615	25.38	ng	96
82) Chrysene	13.78	228	283518	28.55	ng	99
83) Bis(2-ethylhexyl)phthalate	13.76	149	188754	25.66	ng #	96
84) Di-n-octyl phthalate	14.37	149	279504	25.00	ng	100
85) Indeno(1,2,3-cd)pyrene	16.33	276	182689	26.20	ng #	94
87) Benzo(b)fluoranthene	14.74	252	257120m	27.74	ng	
88) Benzo(k)fluoranthene	14.77	252	187226m	23.87	ng	
89) Benzo(a)pyrene	15.05	252	197397	25.67	ng	98
90) Dibenzo(a,h)anthracene	16.34	278	147188	24.56	ng #	92
91) Benzo(g,h,i)perylene	16.70	276	150142	24.70	ng #	88

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

