

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF061915\
 Data File : BF080079.D
 Acq On : 20 Jun 2015 4:03
 Operator : TP/IZ
 Sample : G2683-02MS
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PE-2MS

Manual Integrations
 APPROVED

apatel
 6/22/2015 2:43:52 PM

Quant Time: Jun 20 04:47:42 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF061715.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 17 16:41:53 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.33	152	38037	20.00	ng	-0.01
21) Naphthalene-d8	8.62	136	159578	20.00	ng	-0.01
38) Acenaphthene-d10	10.39	164	86427	20.00	ng	-0.01
63) Phenanthrene-d10	11.90	188	184104	20.00	ng	-0.01
75) Chrysene-d12	14.81	240	209126	20.00	ng	-0.05
86) Perylene-d12	16.67	264	195609	20.00	ng	-0.08

System Monitoring Compounds

5) 2-Fluorophenol	5.94	112	323190	150.41	ng	0.00
7) Phenol-d6	6.93	99	411274	143.83	ng	-0.01
23) Nitrobenzene-d5	7.89	82	269740	98.40	ng	-0.01
41) 2,4,6-Tribromophenol	11.19	330	131098	155.12	ng	-0.01
44) 2-Fluorobiphenyl	9.69	172	496025	81.85	ng	-0.01
78) Terphenyl-d14	13.57	244	692616	77.35	ng	-0.02

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.22	88	32615	31.93	ng	89
3) Pyridine	4.04	79	103495	38.10	ng	# 81
4) n-Nitrosodimethylamine	3.94	42	48465	37.54	ng	95
6) Aniline	6.98	93	168195	42.75	ng	92
8) 2-Chlorophenol	7.11	128	118109	47.56	ng	85
9) Benzaldehyde	6.88	77	56406	34.17	ng	# 88
10) Phenol	6.95	94	144633	46.35	ng	86
11) bis(2-Chloroethyl)ether	7.05	93	108175	43.69	ng	87
12) 1,3-Dichlorobenzene	7.27	146	122796m	41.16	ng	
13) 1,4-Dichlorobenzene	7.35	146	139253m	49.08	ng	
14) 1,2-Dichlorobenzene	7.51	146	123517	44.74	ng	99
15) Benzyl Alcohol	7.45	79	105946	45.32	ng	# 75
16) 2,2'-oxybis(1-Chloropropan	7.60	45	188353	51.26	ng	84
17) 2-Methylphenol	7.56	107	91522	43.14	ng	# 82
18) Hexachloroethane	7.85	117	45078	47.72	ng	# 87
19) n-Nitroso-di-n-propylamine	7.73	70	87953	44.30	ng	# 77
20) 3+4-Methylphenols	7.72	107	131859	48.16	ng	92
22) Acetophenone	7.73	105	169233	45.51	ng	# 82
24) Nitrobenzene	7.91	77	123188	43.54	ng	# 77
25) Isophorone	8.14	82	231580	41.98	ng	# 83
26) 2-Nitrophenol	8.23	139	59868	50.55	ng	# 74
27) 2,4-Dimethylphenol	8.25	122	105972	42.91	ng	87
28) bis(2-Chloroethoxy)methane	8.34	93	150347	46.39	ng	# 92
29) 2,4-Dichlorophenol	8.47	162	106115	50.18	ng	97
30) 1,2,4-Trichlorobenzene	8.56	180	118053	49.16	ng	99
31) Naphthalene	8.64	128	341189m	40.89	ng	
32) Benzoic acid	8.25	122	105972	70.98	ng	# 25
33) 4-Chloroaniline	8.68	127	120157m	33.65	ng	
34) Hexachlorobutadiene	8.77	225	64171	43.39	ng	99
35) Caprolactam	9.01	113	29916	43.58	ng	# 69
36) 4-Chloro-3-methylphenol	9.14	107	105739	44.33	ng	# 71
37) 2-Methylnaphthalene	9.34	142	268479	49.75	ng	# 91
39) 1,2,4,5-Tetrachlorobenzene	9.51	216	121253	48.21	ng	99
40) Hexachlorocyclopentadiene	9.50	237	113000	86.74	ng	97

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.61	196	81410	47.57	ng	98
43) 2,4,5-Trichlorophenol	9.65	196	82073	41.63	ng #	90
45) 1,1'-Biphenyl	9.81	154	292098	41.54	ng	93
46) 2-Chloronaphthalene	9.83	162	236228	41.40	ng	93
47) 2-Nitroaniline	9.91	65	66907	42.72	ng #	57
48) Acenaphthylene	10.25	152	405546	42.59	ng	98
49) Dimethylphthalate	10.09	163	310657	47.81	ng #	97
50) 2,6-Dinitrotoluene	10.15	165	57727	44.32	ng #	47
51) Acenaphthene	10.42	154	254252	43.64	ng	94
52) 3-Nitroaniline	10.32	138	58488	38.92	ng #	52
53) 2,4-Dinitrophenol	10.44	184	45336	85.11	ng #	1
54) Dibenzofuran	10.60	168	348466	42.15	ng #	87
55) 4-Nitrophenol	10.47	139	116124	96.66	ng #	67
56) 2,4-Dinitrotoluene	10.56	165	87052	52.66	ng #	91
57) Fluorene	10.95	166	295109	44.99	ng	93
58) 2,3,4,6-Tetrachlorophenol	10.71	232	71055	42.69	ng	99
59) Diethylphthalate	10.79	149	284060	43.42	ng	98
60) 4-Chlorophenyl-phenylether	10.93	204	137190	43.52	ng #	84
61) 4-Nitroaniline	10.94	138	70485	45.41	ng #	48
62) Azobenzene	11.09	77	302371	45.40	ng	86
64) 4,6-Dinitro-2-methylphenol	10.97	198	40352	44.69	ng	94
65) n-Nitrosodiphenylamine	11.04	169	255941	42.69	ng	91
66) 4-Bromophenyl-phenylether	11.43	248	83266	42.54	ng #	85
67) Hexachlorobenzene	11.51	284	92865	42.69	ng #	78
68) Atrazine	11.57	200	76093	40.18	ng	86
69) Pentachlorophenol	11.69	266	102018	82.75	ng	97
70) Phenanthrene	11.92	178	467056	41.90	ng	96
71) Anthracene	11.98	178	452418	42.16	ng	99
72) Carbazole	12.13	167	408515	39.87	ng	97
73) Di-n-butylphthalate	12.45	149	524674	44.32	ng #	99
74) Fluoranthene	13.17	202	503312	42.40	ng	90
76) Benzidine	13.28	184	317504	54.32	ng #	98
77) Pyrene	13.42	202	513588	39.89	ng	96
79) Butylbenzylphthalate	14.09	149	215133	41.61	ng #	78
80) Benzo(a)anthracene	14.80	228	519484	40.78	ng	96
81) 3,3'-Dichlorobenzidine	14.75	252	167323	38.08	ng #	93
82) Chrysene	14.85	228	478638	40.74	ng	94
83) Bis(2-ethylhexyl)phthalate	14.77	149	343074	41.98	ng #	92
84) Di-n-octyl phthalate	15.56	149	553397	39.52	ng	95
85) Indeno(1,2,3-cd)pyrene	18.52	276	552775	37.31	ng #	88
87) Benzo(b)fluoranthene	16.13	252	470604	39.74	ng #	86
88) Benzo(k)fluoranthene	16.16	252	506441m	41.99	ng	
89) Benzo(a)pyrene	16.59	252	480654	42.73	ng #	87
90) Dibenzo(a,h)anthracene	18.54	278	469707	40.80	ng #	80
91) Benzo(g,h,i)perylene	19.08	276	467928	40.26	ng #	76

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

