

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032916\
 Data File : BF085857.D
 Acq On : 29 Mar 2016 19:08
 Operator : UM/SJ
 Sample : H1970-07MS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-7MS

Manual Integrations
 APPROVED

Sohil
 3/30/2016 7:05:03 PM

Quant Time: Mar 30 01:14:41 2016
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032416.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 28 15:27:05 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.80	152	111767	20.00	ng	-0.02
21) Naphthalene-d8	8.09	136	468561	20.00	ng	-0.02
38) Acenaphthene-d10	9.85	164	215179	20.00	ng	-0.01
63) Phenanthrene-d10	11.33	188	400211	20.00	ng	-0.02
75) Chrysene-d12	13.97	240	302642	20.00	ng	-0.01
86) Perylene-d12	15.39	264	211711	20.00	ng	-0.01

System Monitoring Compounds

5) 2-Fluorophenol	5.41	112	862993	128.59	ng	0.00
7) Phenol-d6	6.45	99	1241369	145.48	ng	-0.01
23) Nitrobenzene-d5	7.37	82	770437	92.59	ng	-0.02
41) 2,4,6-Tribromophenol	10.64	330	288394	141.06	ng	-0.01
44) 2-Fluorobiphenyl	9.17	172	1169997	90.84	ng	-0.02
78) Terphenyl-d14	12.92	244	1058327	71.92	ng	-0.01

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.41	88	130452	40.71	ng	99
3) Pyridine	3.12	79	373134	48.57	ng	99
4) n-Nitrosodimethylamine	3.08	42	154943	50.38	ng	98
6) Aniline	6.47	93	419918	36.56	ng	# 64
8) 2-Chlorophenol	6.60	128	392418	50.33	ng	98
9) Benzaldehyde	6.36	77	43399	8.44	ng	98
10) Phenol	6.47	94	421011	43.55	ng	# 71
11) bis(2-Chloroethyl)ether	6.55	93	373996m	50.27	ng	
12) 1,3-Dichlorobenzene	6.74	146	374923	44.65	ng	# 93
13) 1,4-Dichlorobenzene	6.82	146	399140	46.87	ng	96
14) 1,2-Dichlorobenzene	6.97	146	343345	43.26	ng	94
15) Benzyl Alcohol	6.96	79	299317	52.58	ng	95
16) 2,2'-oxybis(1-Chloropropan	7.09	45	431894	43.76	ng	95
17) 2-Methylphenol	7.08	107	330395	52.93	ng	99
18) Hexachloroethane	7.32	117	139564	44.76	ng	# 88
19) n-Nitroso-di-n-propylamine	7.22	70	241440	47.26	ng	93
20) 3+4-Methylphenols	7.22	107	350920	46.76	ng	98
22) Acetophenone	7.22	105	474146	43.44	ng	# 98
24) Nitrobenzene	7.40	77	412900	48.14	ng	96
25) Isophorone	7.64	82	727070	45.36	ng	98
26) 2-Nitrophenol	7.72	139	220470	51.38	ng	95
27) 2,4-Dimethylphenol	7.75	122	338426	45.12	ng	99
28) bis(2-Chloroethoxy)methane	7.85	93	458561	50.21	ng	99
29) 2,4-Dichlorophenol	7.96	162	307707	48.80	ng	99
30) 1,2,4-Trichlorobenzene	8.04	180	318318	45.47	ng	98
31) Naphthalene	8.12	128	1069229	45.29	ng	100
32) Benzoic acid	7.75	122	338426	67.94	ng	# 15
33) 4-Chloroaniline	8.17	127	276259	28.66	ng	97
34) Hexachlorobutadiene	8.23	225	156117	41.03	ng	99
35) Caprolactam	8.55	113	103348m	46.27	ng	
36) 4-Chloro-3-methylphenol	8.66	107	362278	49.19	ng	92
37) 2-Methylnaphthalene	8.80	142	704169	53.58	ng	97
39) 1,2,4,5-Tetrachlorobenzene	8.97	216	270732	42.43	ng	98
40) Hexachlorocyclopentadiene	8.96	237	220205	82.48	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.09	196	199276	46.24	ng	100
43) 2,4,5-Trichlorophenol	9.14	196	217679	48.16	ng	97
45) 1,1'-Biphenyl	9.27	154	748764	41.24	ng	98
46) 2-Chloronaphthalene	9.29	162	596716	44.69	ng	# 90
47) 2-Nitroaniline	9.40	65	197018	48.26	ng	93
48) Acenaphthylene	9.72	152	1029752	47.33	ng	99
49) Dimethylphthalate	9.58	163	887973	54.39	ng	100
50) 2,6-Dinitrotoluene	9.65	165	176958	50.36	ng	# 60
51) Acenaphthene	9.89	154	609506	45.87	ng	100
52) 3-Nitroaniline	9.81	138	144327	35.88	ng	94
53) 2,4-Dinitrophenol	9.92	184	76693	45.98	ng	# 81
54) Dibenzofuran	10.06	168	894605	52.01	ng	97
55) 4-Nitrophenol	9.99	139	262041	100.00	ng	# 74
56) 2,4-Dinitrotoluene	10.05	165	209148	46.83	ng	# 62
57) Fluorene	10.40	166	682087	47.71	ng	99
58) 2,3,4,6-Tetrachlorophenol	10.17	232	154225	49.10	ng	98
59) Diethylphthalate	10.28	149	736586	45.67	ng	99
60) 4-Chlorophenyl-phenylether	10.39	204	329023	47.07	ng	89
61) 4-Nitroaniline	10.42	138	179614	46.69	ng	89
62) Azobenzene	10.55	77	809607	52.23	ng	97
64) 4,6-Dinitro-2-methylphenol	10.46	198	108839	46.44	ng	# 76
65) n-Nitrosodiphenylamine	10.52	169	630315	49.73	ng	99
66) 4-Bromophenyl-phenylether	10.88	248	200122	48.88	ng	# 87
67) Hexachlorobenzene	10.95	284	217844	50.87	ng	# 91
68) Atrazine	11.04	200	199018	46.50	ng	99
69) Pentachlorophenol	11.14	266	211805	100.76	ng	100
70) Phenanthrene	11.36	178	1057147	51.44	ng	99
71) Anthracene	11.41	178	948102	43.95	ng	100
72) Carbazole	11.57	167	960822	53.05	ng	100
73) Di-n-butylphthalate	11.90	149	1179716	47.47	ng	99
74) Fluoranthene	12.54	202	923300	42.27	ng	97
76) Benzidine	12.67	184	391626	36.74	ng	98
77) Pyrene	12.77	202	943514	37.93	ng	99
79) Butylbenzylphthalate	13.39	149	420435	40.36	ng	# 73
80) Benzo(a)anthracene	13.96	228	759046	43.99	ng	99
81) 3,3'-Dichlorobenzidine	13.92	252	160543	13.40	ng	# 97
82) Chrysene	13.99	228	652934	37.56	ng	98
83) Bis(2-ethylhexyl)phthalate	13.95	149	514729	41.26	ng	# 98
84) Di-n-octyl phthalate	14.55	149	844807	45.55	ng	99
85) Indeno(1,2,3-cd)pyrene	16.77	276	616435	45.93	ng	99
87) Benzo(b)fluoranthene	14.99	252	710157m	53.25	ng	
88) Benzo(k)fluoranthene	15.01	252	484732m	40.32	ng	
89) Benzo(a)pyrene	15.33	252	566628	47.85	ng	# 97
90) Dibenzo(a,h)anthracene	16.78	278	520178	47.28	ng	96
91) Benzo(g,h,i)perylene	17.18	276	518644	46.55	ng	98

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

