

Data Path : Z:\HPCHEM1\BNA F\DATA\BF010617\
 Data File : BF092126.D
 Acq On : 6 Jan 2017 19:31
 Operator : UM/SJ
 Sample : I1025-03MSD
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP7-8-COMP10MSD

Manual Integrations
 APPROVED

Sohil
 1/9/2017 10:48:08 AM

Quant Time: Jan 06 23:18:44 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF122116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 06 15:43:51 2017
 Response via : Initial Calibration

1/9/2017 10:48:08 AM

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.64	152	160289	20.00	ng	0.00
21) Naphthalene-d8	7.93	136	644857	20.00	ng	0.00
38) Acenaphthene-d10	9.68	164	279791	20.00	ng	0.00
63) Phenanthrene-d10	11.16	188	377039	20.00	ng	0.00
75) Chrysene-d12	13.79	240	307377	20.00	ng	0.00
86) Perylene-d12	15.16	264	327893	20.00	ng	0.01

System Monitoring Compounds

5) 2-Fluorophenol	5.27	112	1862778	194.05	ng	0.03
7) Phenol-d6	6.33	99	1978700	168.83	ng	0.02
23) Nitrobenzene-d5	7.22	82	1514153	130.57	ng	0.01
41) 2,4,6-Tribromophenol	10.47	330	285374	123.25	ng	0.00
44) 2-Fluorobiphenyl	9.01	172	1463389	111.13	ng	0.00
78) Terphenyl-d14	12.75	244	1092710	86.22	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.15	88	290058	61.37	ng	97
3) Pyridine	2.80	79	924368	56.04	ng	95
4) n-Nitrosodimethylamine	2.76	42	393100	63.18	ng	99
6) Aniline	6.31	93	323082	21.22	ng	# 44
8) 2-Chlorophenol	6.45	128	643460	58.93	ng	89
9) Benzaldehyde	6.18	77	65665	7.45	ng	99
10) Phenol	6.34	94	971863	72.77	ng	# 68
11) bis(2-Chloroethyl)ether	6.39	93	632120	59.48	ng	98
12) 1,3-Dichlorobenzene	6.58	146	646630	55.47	ng	# 93
13) 1,4-Dichlorobenzene	6.66	146	600904	50.64	ng	95
14) 1,2-Dichlorobenzene	6.81	146	626088	60.81	ng	97
15) Benzyl Alcohol	6.81	79	574114	63.04	ng	92
16) 2,2'-oxybis(1-Chloropropan	6.94	45	1132115	71.54	ng	74
17) 2-Methylphenol	6.95	107	573319	65.28	ng	97
18) Hexachloroethane	7.16	117	288650	65.62	ng	98
19) n-Nitroso-di-n-propylamine	7.08	70	567369	72.76	ng	98
20) 3+4-Methylphenols	7.11	107	789864	75.41	ng	# 70
22) Acetophenone	7.06	105	1049899	66.01	ng	# 92
24) Nitrobenzene	7.24	77	678683	54.95	ng	98
25) Isophorone	7.49	82	1304486	60.97	ng	97
26) 2-Nitrophenol	7.56	139	321703	52.81	ng	95
27) 2,4-Dimethylphenol	7.62	122	468467	46.37	ng	93
28) bis(2-Chloroethoxy)methane	7.69	93	672053	51.80	ng	98
29) 2,4-Dichlorophenol	7.82	162	426765	49.89	ng	94
30) 1,2,4-Trichlorobenzene	7.88	180	495482	52.85	ng	95
31) Naphthalene	7.96	128	1689184	57.23	ng	99
33) 4-Chloroaniline	8.02	127	161186	12.11	ng	# 94
34) Hexachlorobutadiene	8.07	225	269162	54.39	ng	99
35) Caprolactam	8.38	113	132899	47.27	ng	# 58
36) 4-Chloro-3-methylphenol	8.53	107	393843	40.01	ng	99
37) 2-Methylnaphthalene	8.64	142	936366	58.85	ng	99
39) 1,2,4,5-Tetrachlorobenzene	8.81	216	448042	63.38	ng	99
40) Hexachlorocyclopentadiene	8.80	237	206281	66.24	ng	97
42) 2,4,6-Trichlorophenol	8.94	196	263410	53.28	ng	97

Data Path : Z:\HPCHEM1\BNA F\DATA\BF010617\
 Data File : BF092126.D
 Acq On : 6 Jan 2017 19:31
 Operator : UM/SJ
 Sample : I1025-03MSD
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP7-8-COMP10MSD

Manual Integrations
 APPROVED

Sohil
 1/9/2017 10:48:08 AM

Quant Time: Jan 06 23:18:44 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF122116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 06 15:43:51 2017
 Response via : Initial Calibration

1/9/2017 10:48:08 AM

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
43) 2,4,5-Trichlorophenol	9.00	196	238219	46.82	nq	# 85
45) 1,1'-Biphenyl	9.11	154	1173307	61.25	nq	99
46) 2-Chloronaphthalene	9.13	162	897443	59.15	nq	98
47) 2-Nitroaniline	9.24	65	244952	45.35	nq	99
48) Acenaphthylene	9.54	152	1304275	55.90	nq	99
49) Dimethylphthalate	9.42	163	1123163	62.76	nq	99
50) 2,6-Dinitrotoluene	9.48	165	192158	49.06	nq	# 66
51) Acenaphthene	9.72	154	837256	52.93	nq	99
52) 3-Nitroaniline	9.65	138	152337	31.43	nq	92
53) 2,4-Dinitrophenol	9.76	184	15306	7.02	nq	# 74
54) Dibenzofuran	9.89	168	1130846	52.94	nq	97
55) 4-Nitrophenol	9.85	139	267141m	81.17	nq	
56) 2,4-Dinitrotoluene	9.88	165	238676	48.55	nq	# 69
57) Fluorene	10.23	166	831793	53.52	nq	99
58) 2,3,4,6-Tetrachlorophenol	10.01	232	184503	45.85	nq	# 99
59) Diethylphthalate	10.12	149	897167	51.62	nq	98
60) 4-Chlorophenyl-phenylether	10.22	204	342957	50.40	nq	98
61) 4-Nitroaniline	10.26	138	184340	36.70	nq	76
62) Azobenzene	10.38	77	955397	49.93	nq	99
64) 4,6-Dinitro-2-methylphenol	10.29	198	57638	25.28	nq	# 65
65) n-Nitrosodiphenylamine	10.34	169	687660	61.46	nq	99
66) 4-Bromophenyl-phenylether	10.71	248	241121	61.48	nq	# 89
67) Hexachlorobenzene	10.78	284	238359	56.86	nq	# 83
68) Atrazine	10.88	200	220749	61.17	nq	96
69) Pentachlorophenol	10.97	266	192198	93.43	nq	98
70) Phenanthrene	11.18	178	1047165	51.22	nq	99
71) Anthracene	11.24	178	1166945	Below	Cal	98
72) Carbazole	11.40	167	1016260	Below	Cal	98
73) Di-n-butylphthalate	11.73	149	1210266	64.23	nq	99
74) Fluoranthene	12.37	202	1125806	64.42	nq	91
76) Benzidine	12.49	184	423680	57.70	nq	96
77) Pyrene	12.59	202	1104841	48.99	nq	99
79) Butylbenzylphthalate	13.23	149	557076	54.31	nq	# 81
80) Benzo(a)anthracene	13.77	228	1024271	59.03	nq	100
81) 3,3'-Dichlorobenzidine	13.75	252	256827	39.03	nq	# 95
82) Chrysene	13.81	228	895261	51.44	nq	99
83) Bis(2-ethylhexyl)phthalate	13.79	149	714067	59.11	nq	98
84) Di-n-octyl phthalate	14.39	149	1315604	58.79	nq	98
85) Indeno(1,2,3-cd)pyrene	16.43	276	907397	60.18	nq	99
87) Benzo(b)fluoranthene	14.78	252	1013905m	49.67	nq	
88) Benzo(k)fluoranthene	14.80	252	964020m	55.38	nq	
89) Benzo(a)pyrene	15.10	252	941243	51.95	nq	99
90) Dibenzo(a,h)anthracene	16.44	278	734495	49.32	nq	99
91) Benzo(g,h,i)perylene	16.80	276	754328	48.71	nq	# 95

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

Data Path : Z:\HPCHEM1\BNA F\DATA\BF010617\
 Data File : BF092126.D
 Acq On : 6 Jan 2017 19:31
 Operator : UM/SJ
 Sample : I1025-03MSD
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP7-8-COMP10MSD

Manual Integrations
 APPROVED

Sohil
 1/9/2017 10:48:08 AM

Quant Time: Jan 06 23:18:44 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF122116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 06 15:43:51 2017
 Response via : Initial Calibration

1/9/2017 10:48:08 AM

