

Data Path : Z:\HPCHEM1\BNA F\DATA\BF121416\
 Data File : BF091591.D
 Acq On : 14 Dec 2016 18:28
 Operator : UM/SJ
 Sample : PB95351BS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB95351BS

Manual Integrations
 APPROVED

sohil
 12/15/2016 5:39:54 PM

Quant Time: Dec 15 01:11:37 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF120916.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 09 19:00:21 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.80	152	277057	20.00	ng	-0.02
21) Naphthalene-d8	8.08	136	1122447	20.00	ng	-0.02
38) Acenaphthene-d10	9.84	164	617810	20.00	ng	-0.02
63) Phenanthrene-d10	11.31	188	970018	20.00	ng	-0.02
75) Chrysene-d12	13.95	240	699123	20.00	ng	-0.01
86) Perylene-d12	15.36	264	577154	20.00	ng	-0.02

System Monitoring Compounds

5) 2-Fluorophenol	5.40	112	2017599	122.69	ng	-0.01
7) Phenol-d6	6.44	99	2494761	121.69	ng	-0.01
23) Nitrobenzene-d5	7.36	82	1826462	90.82	ng	-0.02
41) 2,4,6-Tribromophenol	10.62	330	571171	111.31	ng	-0.01
44) 2-Fluorobiphenyl	9.16	172	2665647	74.55	ng	-0.02
78) Terphenyl-d14	12.90	244	2398253	77.84	ng	-0.02

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.42	88	300999	34.68	ng	# 86
3) Pyridine	3.12	79	844923	36.54	ng	# 80
4) n-Nitrosodimethylamine	3.06	42	413844	41.65	ng	# 63
6) Aniline	6.45	93	608677	22.64	ng	# 27
8) 2-Chlorophenol	6.58	128	762006	41.05	ng	86
9) Benzaldehyde	6.34	77	142836m	10.12	ng	
10) Phenol	6.45	94	878114	36.63	ng	91
11) bis(2-Chloroethyl)ether	6.53	93	786024	43.65	ng	89
12) 1,3-Dichlorobenzene	6.74	146	844600	41.96	ng	94
13) 1,4-Dichlorobenzene	6.81	146	810261	40.81	ng	94
14) 1,2-Dichlorobenzene	6.97	146	807052	45.11	ng	95
15) Benzyl Alcohol	6.94	79	727291	44.93	ng	98
16) 2,2'-oxybis(1-Chloropropan	7.08	45	1179539	42.27	ng	80
17) 2-Methylphenol	7.06	107	644090m	41.94	ng	
18) Hexachloroethane	7.31	117	324503	41.88	ng	93
19) n-Nitroso-di-n-propylamine	7.22	70	601728	46.94	ng	# 76
20) 3+4-Methylphenols	7.22	107	818118	49.65	ng	93
22) Acetophenone	7.21	105	1126735	41.90	ng	# 97
24) Nitrobenzene	7.38	77	870171	40.95	ng	# 89
25) Isophorone	7.62	82	1607921	44.98	ng	97
26) 2-Nitrophenol	7.70	139	457813	48.82	ng	# 80
27) 2,4-Dimethylphenol	7.74	122	767280	46.58	ng	95
28) bis(2-Chloroethoxy)methane	7.84	93	1006069	47.98	ng	100
29) 2,4-Dichlorophenol	7.94	162	643762	44.92	ng	95
30) 1,2,4-Trichlorobenzene	8.02	180	683318	41.95	ng	97
31) Naphthalene	8.10	128	2165601	41.47	ng	99
32) Benzoic acid	7.87	122	581527	49.08	ng	93
33) 4-Chloroaniline	8.16	127	214023	9.51	ng	# 93
34) Hexachlorobutadiene	8.22	225	380385	40.74	ng	98
35) Caprolactam	8.52	113	226231m	54.06	ng	
36) 4-Chloro-3-methylphenol	8.65	107	706510	43.96	ng	99
37) 2-Methylnaphthalene	8.80	142	1518027	44.91	ng	97
39) 1,2,4,5-Tetrachlorobenzene	8.97	216	631908	37.29	ng	99
40) Hexachlorocyclopentadiene	8.96	237	683350	90.45	ng	98

Data Path : Z:\HPCHEM1\BNA F\DATA\BF121416\
 Data File : BF091591.D
 Acq On : 14 Dec 2016 18:28
 Operator : UM/SJ
 Sample : PB95351BS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB95351BS

Manual Integrations
 APPROVED

sohil
 12/15/2016 5:39:54 PM

Quant Time: Dec 15 01:11:37 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF120916.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 09 19:00:21 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.08	196	474735	43.60	nq	97
43) 2,4,5-Trichlorophenol	9.13	196	480436	42.94	nq	# 91
45) 1,1'-Biphenyl	9.26	154	1884084	42.10	nq	95
46) 2-Chloronaphthalene	9.29	162	1449550	41.83	nq	94
47) 2-Nitroaniline	9.38	65	441823	37.92	nq	89
48) Acenaphthylene	9.70	152	2199101	41.63	nq	98
49) Dimethylphthalate	9.56	163	1541914	39.42	nq	98
50) 2,6-Dinitrotoluene	9.62	165	362599	42.23	nq	89
51) Acenaphthene	9.87	154	1596047	43.50	nq	97
52) 3-Nitroaniline	9.79	138	204579	20.28	nq	93
53) 2,4-Dinitrophenol	9.89	184	407155	80.35	nq	# 81
54) Dibenzofuran	10.04	168	1851696	40.79	nq	99
55) 4-Nitrophenol	9.96	139	657404	87.61	nq	93
56) 2,4-Dinitrotoluene	10.03	165	546377	50.77	nq	89
57) Fluorene	10.38	166	1510177	46.75	nq	98
58) 2,3,4,6-Tetrachlorophenol	10.17	232	384034	43.84	nq	96
59) Diethylphthalate	10.27	149	1666984	44.25	nq	# 93
60) 4-Chlorophenyl-phenylether	10.37	204	596361	36.54	nq	# 76
61) 4-Nitroaniline	10.41	138	395671	39.48	nq	95
62) Azobenzene	10.53	77	1794683	42.04	nq	88
64) 4,6-Dinitro-2-methylphenol	10.43	198	248892	42.60	nq	# 79
65) n-Nitrosodiphenylamine	10.50	169	1337850	46.26	nq	100
66) 4-Bromophenyl-phenylether	10.86	248	417624	41.99	nq	100
67) Hexachlorobenzene	10.93	284	470084	45.46	nq	98
68) Atrazine	11.02	200	387314	41.65	nq	99
69) Pentachlorophenol	11.13	266	527107	90.57	nq	94
70) Phenanthrene	11.35	178	2321849	48.34	nq	99
71) Anthracene	11.39	178	2072673	40.00	nq	99
72) Carbazole	11.55	167	2069050	45.56	nq	99
73) Di-n-butylphthalate	11.88	149	2320940	43.86	nq	98
74) Fluoranthene	12.52	202	2099822	41.75	nq	94
76) Benzidine	12.65	184	675534	31.25	nq	98
77) Pyrene	12.75	202	2110459	38.05	nq	98
79) Butylbenzylphthalate	13.38	149	1049750	48.55	nq	98
80) Benzo(a)anthracene	13.94	228	1644088	40.49	nq	100
81) 3,3'-Dichlorobenzidine	13.91	252	316586	21.71	nq	100
82) Chrysene	13.97	228	1611632	37.94	nq	99
83) Bis(2-ethylhexyl)phthalate	13.94	149	1229269	44.31	nq	98
84) Di-n-octyl phthalate	14.56	149	2211295	46.96	nq	# 90
85) Indeno(1,2,3-cd)pyrene	16.72	276	1415011	37.02	nq	98
87) Benzo(b)fluoranthene	14.96	252	1447718	42.08	nq	99
88) Benzo(k)fluoranthene	14.99	252	1419442m	46.92	nq	
89) Benzo(a)pyrene	15.30	252	1340739	43.26	nq	98
90) Dibenzo(a,h)anthracene	16.74	278	1162902	41.88	nq	97
91) Benzo(g,h,i)perylene	17.13	276	1233751	43.66	nq	# 94

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

Data Path : Z:\HPCHEM1\BNA F\DATA\BF121416\
 Data File : BF091591.D
 Acq On : 14 Dec 2016 18:28
 Operator : UM/SJ
 Sample : PB95351BS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 Client Sample ID :
 PB95351BS

Manual Integrations
 APPROVED

sohil
 12/15/2016 5:39:54 PM

Quant Time: Dec 15 01:11:37 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF120916.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 09 19:00:21 2016
 Response via : Initial Calibration

