

Data Path : Z:\HPCHEM1\BNA F\DATA\BF112715\
 Data File : BF083333.D
 Acq On : 27 Nov 2015 17:06
 Operator : UM/IZ
 Sample : SSTDICC050
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC050

Manual Integrations
 APPROVED

Mmdadoda
 12/1/2015 5:22:57 PM

Quant Time: Nov 27 18:11:28 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF112715.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 27 17:51:44 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.56	152	183613	20.00	ng	0.00
21) Naphthalene-d8	7.85	136	707278	20.00	ng	0.00
38) Acenaphthene-d10	9.60	164	343940	20.00	ng	0.00
63) Phenanthrene-d10	11.07	188	646067	20.00	ng	0.00
75) Chrysene-d12	13.69	240	443201	20.00	ng	0.00
86) Perylene-d12	15.04	264	356740	20.00	ng	0.01

System Monitoring Compounds

5) 2-Fluorophenol	5.14	112	1196347	47.27	ng	0.00
7) Phenol-d6	6.24	99	1534915	45.28	ng	0.00
23) Nitrobenzene-d5	7.14	82	1473678	45.46	ng	0.00
41) 2,4,6-Tribromophenol	10.39	330	313142	49.16	ng	0.00
44) 2-Fluorobiphenyl	8.94	172	2144444	43.16	ng	0.00
78) Terphenyl-d14	12.66	244	2203445	48.37	ng	0.01

Target Compounds

						Qvalue
3) Pyridine	2.59	79	826117	54.45	ng	92
4) n-Nitrosodimethylamine	2.56	42	407873	52.44	ng	89
6) Aniline	6.23	93	1110921	48.60	ng	99
8) 2-Chlorophenol	6.35	128	628496	45.56	ng	98
9) Benzaldehyde	6.10	77	376155	40.30	ng	97
10) Phenol	6.25	94	938625	46.39	ng	81
11) bis(2-Chloroethyl)ether	6.31	93	708622	47.39	ng	90
12) 1,3-Dichlorobenzene	6.50	146	721859	47.44	ng	95
13) 1,4-Dichlorobenzene	6.58	146	761773	47.91	ng	98
14) 1,2-Dichlorobenzene	6.74	146	679995	46.45	ng	96
15) Benzyl Alcohol	6.73	79	605196	46.62	ng	96
16) 2,2'-oxybis(1-Chloropropan	6.87	45	1247331	48.06	ng	# 66
17) 2-Methylphenol	6.86	107	537252	48.06	ng	98
18) Hexachloroethane	7.07	117	261271	46.60	ng	# 81
19) n-Nitroso-di-n-propylamine	7.00	70	564625	47.46	ng	99
20) 3+4-Methylphenols	7.02	107	701662	46.12	ng	96
22) Acetophenone	6.99	105	983455	46.77	ng	# 98
24) Nitrobenzene	7.16	77	888666	51.06	ng	95
25) Isophorone	7.40	82	1486705	47.99	ng	97
26) 2-Nitrophenol	7.48	139	342492	48.49	ng	# 85
27) 2,4-Dimethylphenol	7.54	122	595520	50.22	ng	95
28) bis(2-Chloroethoxy)methane	7.62	93	804224	46.24	ng	99
29) 2,4-Dichlorophenol	7.74	162	547920	48.27	ng	94
30) 1,2,4-Trichlorobenzene	7.80	180	600540	48.69	ng	98
31) Naphthalene	7.87	128	1739491	44.10	ng	99
32) Benzoic acid	7.70	122	335979	62.49	ng	96
33) 4-Chloroaniline	7.94	127	819963	49.80	ng	93
34) Hexachlorobutadiene	8.00	225	338452	47.56	ng	99
35) Caprolactam	8.33	113	175225	54.07	ng	95
36) 4-Chloro-3-methylphenol	8.44	107	623303	49.85	ng	93
37) 2-Methylnaphthalene	8.57	142	1245342	48.73	ng	96
39) 1,2,4,5-Tetrachlorobenzene	8.74	216	552101	49.53	ng	100
40) Hexachlorocyclopentadiene	8.73	237	287401	65.15	ng	96
42) 2,4,6-Trichlorophenol	8.86	196	387980	50.17	ng	99

Data Path : Z:\HPCHEM1\BNA F\DATA\BF112715\
 Data File : BF083333.D
 Acq On : 27 Nov 2015 17:06
 Operator : UM/IZ
 Sample : SSTDICC050
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC050

Manual Integrations
 APPROVED

Mmdadoda
 12/1/2015 5:22:57 PM

Quant Time: Nov 27 18:11:28 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF112715.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 27 17:51:44 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
43) 2,4,5-Trichlorophenol	8.90	196	403987	51.79	nq	# 89
45) 1,1'-Biphenyl	9.04	154	1545841	49.49	nq	99
46) 2-Chloronaphthalene	9.05	162	1100880	46.94	nq	# 93
47) 2-Nitroaniline	9.16	65	492579	55.06	nq	# 81
48) Acenaphthylene	9.46	152	1677324	45.24	nq	98
49) Dimethylphthalate	9.35	163	1283770	48.06	nq	99
50) 2,6-Dinitrotoluene	9.40	165	278181	48.37	nq	96
51) Acenaphthene	9.63	154	973983	45.69	nq	99
52) 3-Nitroaniline	9.58	138	339497	51.68	nq	# 70
53) 2,4-Dinitrophenol	9.68	184	136647	86.94	nq	95
54) Dibenzofuran	9.80	168	1373031	44.25	nq	96
55) 4-Nitrophenol	9.76	139	200881	63.00	nq	89
56) 2,4-Dinitrotoluene	9.80	165	348387	47.93	nq	# 88
57) Fluorene	10.15	166	1176379	46.37	nq	99
58) 2,3,4,6-Tetrachlorophenol	9.93	232	305466	49.59	nq	98
59) Diethylphthalate	10.04	149	1348539	52.98	nq	99
60) 4-Chlorophenyl-phenylether	10.15	204	567422	48.82	nq	# 86
61) 4-Nitroaniline	10.19	138	348670	55.71	nq	95
62) Azobenzene	10.31	77	1634307	51.15	nq	91
64) 4,6-Dinitro-2-methylphenol	10.22	198	222687	67.19	nq	91
65) n-Nitrosodiphenylamine	10.27	169	1177594	49.02	nq	97
66) 4-Bromophenyl-phenylether	10.63	248	331699	45.06	nq	# 84
67) Hexachlorobenzene	10.70	284	403403	48.57	nq	# 85
68) Atrazine	10.80	200	299413	45.11	nq	96
69) Pentachlorophenol	10.89	266	214209	63.48	nq	96
70) Phenanthrene	11.10	178	1661783	42.41	nq	98
71) Anthracene	11.15	178	1907216	47.23	nq	99
72) Carbazole	11.31	167	1673481	49.62	nq	99
73) Di-n-butylphthalate	11.66	149	2041362	50.97	nq	99
74) Fluoranthene	12.27	202	1671113	45.76	nq	96
76) Benzidine	12.41	184	825447	50.66	nq	96
77) Pyrene	12.50	202	1860623	49.16	nq	100
79) Butylbenzylphthalate	13.14	149	768290	56.92	nq	# 82
80) Benzo(a)anthracene	13.68	228	1435682	50.89	nq	100
81) 3,3'-Dichlorobenzidine	13.66	252	446444	50.94	nq	# 97
82) Chrysene	13.73	228	1397782	52.08	nq	99
83) Bis(2-ethylhexyl)phthalate	13.70	149	929144	54.88	nq	# 98
84) Di-n-octyl phthalate	14.31	149	1313802	48.35	nq	# 93
85) Indeno(1,2,3-cd)pyrene	16.25	276	1279374	42.84	nq	96
87) Benzo(b)fluoranthene	14.69	252	1274612m	51.55	nq	
88) Benzo(k)fluoranthene	14.71	252	803902m	40.76	nq	
89) Benzo(a)pyrene	14.98	252	948888	47.91	nq	# 95
90) Dibenzo(a,h)anthracene	16.26	278	1036212	46.70	nq	# 95
91) Benzo(q,h,i)perylene	16.62	276	1071349	47.97	nq	97

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

