

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF111017\
 Data File : BF100408.D
 Acq On : 10 Nov 2017 23:25
 Operator : SJ/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

SOHIL
 11/13/2017 4:43:32 PM

Quant Time: Nov 11 00:29:06 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF110817.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 10 15:44:40 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.90	152	135700	20.00	ng	0.00
21) Naphthalene-d8	8.19	136	522074	20.00	ng	0.00
38) Acenaphthene-d10	9.94	164	213416	20.00	ng	0.00
63) Phenanthrene-d10	11.43	188	340319	20.00	ng	0.00
75) Chrysene-d12	14.07	240	269220	20.00	ng	0.00
86) Perylene-d12	15.54	264	216022	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.52	112	693238	81.89	ng	0.00
7) Phenol-d6	6.53	99	831990	79.70	ng	0.00
23) Nitrobenzene-d5	7.47	82	730014	92.45	ng	0.00
41) 2,4,6-Tribromophenol	10.73	330	156111	71.78	ng	0.00
44) 2-Fluorobiphenyl	9.26	172	1223420	87.29	ng	0.00
78) Terphenyl-d14	13.01	244	959867	81.92	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.69	88	165861	40.204	ng	94
3) Pyridine	3.47	79	446656	39.684	ng	94
4) n-Nitrosodimethylamine	3.42	42	218256	39.940	ng	98
6) Aniline	6.57	93	576946	39.121	ng	100
8) 2-Chlorophenol	6.69	128	363325	40.569	ng	98
9) Benzaldehyde	6.46	77	301602	40.457	ng	99
10) Phenol	6.54	94	437059	39.882	ng	99
11) bis(2-Chloroethyl)ether	6.64	93	368781	40.064	ng	93
12) 1,3-Dichlorobenzene	6.85	146	412468	39.948	ng	98
13) 1,4-Dichlorobenzene	6.92	146	420805	40.267	ng	98
14) 1,2-Dichlorobenzene	7.07	146	384660	39.811	ng	99
15) Benzyl Alcohol	7.04	79	315453	38.719	ng	98
16) 2,2'-oxybis(1-Chloropropan	7.17	45	484558	32.913	ng	98
17) 2-Methylphenol	7.15	107	299952	39.193	ng	99
18) Hexachloroethane	7.42	117	93154	27.036	ng	99
19) n-Nitroso-di-n-propylamine	7.32	70	268552	37.095	ng	93
20) 3+4-Methylphenols	7.30	107	372616	38.475	ng	# 86
22) Acetophenone	7.32	105	510448	40.096	ng	# 98
24) Nitrobenzene	7.49	77	381494	46.938	ng	99
25) Isophorone	7.72	82	644758	38.854	ng	99
26) 2-Nitrophenol	7.80	139	141202	38.281	ng	97
27) 2,4-Dimethylphenol	7.83	122	288845	38.829	ng	97
28) bis(2-Chloroethoxy)methane	7.93	93	447365	41.215	ng	99
29) 2,4-Dichlorophenol	8.04	162	283438	39.913	ng	97
30) 1,2,4-Trichlorobenzene	8.13	180	314339	40.921	ng	98
31) Naphthalene	8.21	128	1003389	39.903	ng	100
32) Benzoic acid	7.92	122	175055m	34.102	ng	
33) 4-Chloroaniline	8.25	127	417225	39.861	ng	98
34) Hexachlorobutadiene	8.32	225	186677	40.873	ng	99
35) Caprolactam	8.62	113	87771	36.015	ng	# 89
36) 4-Chloro-3-methylphenol	8.73	107	283266	37.140	ng	99
37) 2-Methylnaphthalene	8.90	142	640086	38.341	ng	99
39) 1,2,4,5-Tetrachlorobenzene	9.06	216	328426	46.401	ng	# 97
40) Hexachlorocyclopentadiene	9.05	237	9616	2.887	ng	97

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF111017\
 Data File : BF100408.D
 Acq On : 10 Nov 2017 23:25
 Operator : SJ/JU
 Sample : SSTDC0040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDC0040

Manual Integrations
 APPROVED

SOHIL
 11/13/2017 4:43:32 PM

Quant Time: Nov 11 00:29:06 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF110817.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 10 15:44:40 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.17	196	188949	42.121	nq	98
43) 2,4,5-Trichlorophenol	9.21	196	192382	41.393	nq	96
45) 1,1'-Biphenyl	9.36	154	802049	43.452	nq	99
46) 2-Chloronaphthalene	9.39	162	572891	42.782	nq	98
47) 2-Nitroaniline	9.48	65	194301	49.042	nq	92
48) Acenaphthylene	9.80	152	891689	40.181	nq	99
49) Dimethylphthalate	9.66	163	704967	43.264	nq	99
50) 2,6-Dinitrotoluene	9.72	165	131695	40.830	nq	96
51) Acenaphthene	9.97	154	524549	38.865	nq	100
52) 3-Nitroaniline	9.89	138	173443	45.538	nq	95
53) 2,4-Dinitrophenol	9.99	184	1769	9.539	nq #	42
54) Dibenzofuran	10.15	168	751064	39.705	nq	99
55) 4-Nitrophenol	10.04	139	98240	31.766	nq	93
56) 2,4-Dinitrotoluene	10.12	165	155367	38.183	nq #	96
57) Fluorene	10.49	166	540452	39.001	nq	99
58) 2,3,4,6-Tetrachlorophenol	10.26	232	132476	34.778	nq #	88
59) Diethylphthalate	10.36	149	682440	41.851	nq	96
60) 4-Chlorophenyl-phenylether	10.48	204	283862	41.757	nq	93
61) 4-Nitroaniline	10.50	138	154379	42.746	nq	91
62) Azobenzene	10.64	77	587231	38.236	nq	96
64) 4,6-Dinitro-2-methylphenol	10.53	198	5804	11.345	nq	87
65) n-Nitrosodiphenylamine	10.60	169	553702	46.792	nq	96
66) 4-Bromophenyl-phenylether	10.97	248	168986	46.321	nq	91
67) Hexachlorobenzene	11.03	284	168526	44.168	nq	92
68) Atrazine	11.12	200	155872	43.516	nq	96
69) Pentachlorophenol	11.23	266	97896	35.781	nq	98
70) Phenanthrene	11.45	178	728615	40.663	nq	98
71) Anthracene	11.50	178	749376	41.222	nq	99
72) Carbazole	11.66	167	670213	39.956	nq	99
73) Di-n-butylphthalate	11.99	149	937695	47.770	nq	99
74) Fluoranthene	12.64	202	747688	39.373	nq	97
76) Benzidine	12.76	184	400989	37.192	nq	99
77) Pyrene	12.87	202	773146	40.287	nq	99
79) Butylbenzylphthalate	13.48	149	354080	45.242	nq	99
80) Benzo(a)anthracene	14.06	228	666631	41.119	nq	99
81) 3,3'-Dichlorobenzidine	14.02	252	263453	45.355	nq	99
82) Chrysene	14.09	228	635967	40.509	nq	100
83) Bis(2-ethylhexyl)phthalate	14.04	149	501523	48.722	nq	99
84) Di-n-octyl phthalate	14.65	149	849825	48.057	nq	99
85) Indeno(1,2,3-cd)pyrene	17.04	276	466851	36.764	nq	96
87) Benzo(b)fluoranthene	15.12	252	555814	43.429	nq	98
88) Benzo(k)fluoranthene	15.14	252	521490	43.125	nq #	97
89) Benzo(a)pyrene	15.49	252	472637	41.657	nq	99
90) Dibenzo(a,h)anthracene	17.06	278	395375	42.610	nq	98
91) Benzo(g,h,i)perylene	17.50	276	376693	41.472	nq	94

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

