

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF121817\
 Data File : BF101473.D
 Acq On : 18 Dec 2017 11:48
 Operator : SJ/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Quant Time: Dec 18 15:39:20 2017
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF121417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 18 15:38:59 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.81	152	133520	20.00	ng	0.00
21) Naphthalene-d8	8.09	136	541109	20.00	ng	0.00
38) Acenaphthene-d10	9.85	164	253507	20.00	ng	0.00
63) Phenanthrene-d10	11.34	188	497279	20.00	ng	0.00
75) Chrysene-d12	13.99	240	404367	20.00	ng	0.00
86) Perylene-d12	15.45	264	367317	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.42	112	786249	87.72	ng	0.00
7) Phenol-d6	6.45	99	914948	82.02	ng	0.00
23) Nitrobenzene-d5	7.38	82	835866	85.99	ng	0.00
41) 2,4,6-Tribromophenol	10.65	330	236137	96.67	ng	0.00
44) 2-Fluorobiphenyl	9.17	172	1417230	86.72	ng	0.00
78) Terphenyl-d14	12.92	244	1399937	83.46	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.53	88	189221	41.083	ng	96
3) Pyridine	3.30	79	522527	40.833	ng	95
4) n-Nitrosodimethylamine	3.26	42	236356	40.115	ng	100
6) Aniline	6.47	93	642332	41.434	ng	92
8) 2-Chlorophenol	6.60	128	403653	42.809	ng	99
9) Benzaldehyde	6.37	77	324453	39.382	ng	97
10) Phenol	6.46	94	532183	41.856	ng	92
11) bis(2-Chloroethyl)ether	6.55	93	421033	42.216	ng	93
12) 1,3-Dichlorobenzene	6.75	146	466770	43.861	ng	98
13) 1,4-Dichlorobenzene	6.83	146	476576	43.854	ng	98
14) 1,2-Dichlorobenzene	6.98	146	431919	44.155	ng	97
15) Benzyl Alcohol	6.96	79	314547	40.965	ng	99
16) 2,2'-oxybis(1-Chloropropan	7.08	45	639574	41.901	ng	82
17) 2-Methylphenol	7.07	107	357270	41.191	ng	96
18) Hexachloroethane	7.31	117	161386	42.714	ng	98
19) n-Nitroso-di-n-propylamine	7.23	70	285480	38.968	ng	95
20) 3+4-Methylphenols	7.22	107	399027	43.415	ng	# 65
22) Acetophenone	7.23	105	527588	42.731	ng	# 90
24) Nitrobenzene	7.40	77	447682	41.613	ng	98
25) Isophorone	7.63	82	795159	40.777	ng	98
26) 2-Nitrophenol	7.72	139	225844	48.863	ng	95
27) 2,4-Dimethylphenol	7.75	122	339302	43.212	ng	96
28) bis(2-Chloroethoxy)methane	7.84	93	518750	41.879	ng	99
29) 2,4-Dichlorophenol	7.96	162	350021	45.120	ng	97
30) 1,2,4-Trichlorobenzene	8.03	180	362257	44.250	ng	96
31) Naphthalene	8.12	128	1155979	42.435	ng	99
32) Benzoic acid	7.91	122	252772	47.813	ng	97
33) 4-Chloroaniline	8.17	127	508893	42.981	ng	98
34) Hexachlorobutadiene	8.22	225	211718	44.715	ng	98
35) Caprolactam	8.56	113	118894	44.138	ng	98
36) 4-Chloro-3-methylphenol	8.66	107	366481	42.982	ng	97
37) 2-Methylnaphthalene	8.80	142	755676	42.577	ng	97
39) 1,2,4,5-Tetrachlorobenzene	8.97	216	372303	43.498	ng	97
40) Hexachlorocyclopentadiene	8.95	237	66843	49.331	ng	96

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF121817\
 Data File : BF101473.D
 Acq On : 18 Dec 2017 11:48
 Operator : SJ/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Quant Time: Dec 18 15:39:20 2017
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF121417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 18 15:38:59 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.09	196	234613	45.531	ng	98
43) 2,4,5-Trichlorophenol	9.14	196	253360	44.906	ng	93
45) 1,1'-Biphenyl	9.27	154	938418	41.740	ng	99
46) 2-Chloronaphthalene	9.30	162	729132	42.385	ng	96
47) 2-Nitroaniline	9.40	65	258349	43.474	ng	98
48) Acenaphthylene	9.71	152	1124419	41.383	ng	100
49) Dimethylphthalate	9.57	163	845114	41.445	ng	99
50) 2,6-Dinitrotoluene	9.65	165	187065	45.137	ng	93
51) Acenaphthene	9.88	154	685981	42.022	ng	98
52) 3-Nitroaniline	9.82	138	235114	45.503	ng	98
53) 2,4-Dinitrophenol	9.94	184	60852	47.039	ng	# 20
54) Dibenzofuran	10.05	168	910340	43.882	ng	97
55) 4-Nitrophenol	9.99	139	100999	44.828	ng	94
56) 2,4-Dinitrotoluene	10.06	165	218016	46.691	ng	# 92
57) Fluorene	10.40	166	773767	44.091	ng	98
58) 2,3,4,6-Tetrachlorophenol	10.18	232	190189	46.919	ng	# 87
59) Diethylphthalate	10.27	149	866283	42.900	ng	96
60) 4-Chlorophenyl-phenylether	10.39	204	382504	45.478	ng	92
61) 4-Nitroaniline	10.43	138	227357	43.776	ng	96
62) Azobenzene	10.55	77	844356	40.365	ng	96
64) 4,6-Dinitro-2-methylphenol	10.47	198	135191	53.274	ng	79
65) n-Nitrosodiphenylamine	10.51	169	764490	41.635	ng	96
66) 4-Bromophenyl-phenylether	10.87	248	244343	43.730	ng	# 90
67) Hexachlorobenzene	10.95	284	260340	44.023	ng	# 84
68) Atrazine	11.03	200	244778	44.708	ng	97
69) Pentachlorophenol	11.15	266	97250	45.379	ng	98
70) Phenanthrene	11.36	178	1154225	41.737	ng	99
71) Anthracene	11.42	178	1175089	41.599	ng	99
72) Carbazole	11.57	167	1110199	41.977	ng	99
73) Di-n-butylphthalate	11.89	149	1381192	43.027	ng	100
74) Fluoranthene	12.56	202	1232697	42.467	ng	98
76) Benzidine	12.67	184	698621	40.906	ng	98
77) Pyrene	12.79	202	1281870	42.240	ng	99
79) Butylbenzylphthalate	13.39	149	602005	43.841	ng	97
80) Benzo(a)anthracene	13.98	228	1122860	42.053	ng	99
81) 3,3'-Dichlorobenzidine	13.93	252	376604	43.257	ng	98
82) Chrysene	14.02	228	1035118	41.059	ng	100
83) Bis(2-ethylhexyl)phthalate	13.94	149	752421	43.261	ng	98
84) Di-n-octyl phthalate	14.55	149	1338430	43.150	ng	99
85) Indeno(1,2,3-cd)pyrene	16.94	276	819564	36.506	ng	96
87) Benzo(b)fluoranthene	15.03	252	955514	42.678	ng	98
88) Benzo(k)fluoranthene	15.06	252	940839	43.221	ng	98
89) Benzo(a)pyrene	15.39	252	881273	42.699	ng	99
90) Dibenzo(a,h)anthracene	16.94	278	691364	39.015	ng	97
91) Benzo(g,h,i)perylene	17.38	276	687505	39.181	ng	95

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

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF121817\
Data File : BF101473.D
Acq On : 18 Dec 2017 11:48
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SSTDCCC040

Quant Time: Dec 18 15:39:20 2017
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF121417.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Dec 18 15:38:59 2017
Response via : Initial Calibration

