

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF021219\
 Data File : BF112502.D
 Acq On : 12 Feb 2019 10:14
 Operator : JU/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Sohil
 2/13/2019 11:11:57 AM

Quant Time: Feb 13 09:00:57 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF012519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 28 10:38:18 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.83	152	175121	20.00	ng	-0.01
21) Naphthalene-d8	8.12	136	664518	20.00	ng	-0.01
39) Acenaphthene-d10	9.87	164	318649	20.00	ng	-0.01
64) Phenanthrene-d10	11.36	188	578483	20.00	ng	0.00
76) Chrysene-d12	14.01	240	415057	20.00	ng	0.00
87) Perylene-d12	15.47	264	340341	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.47	112	756435	91.75	ng	-0.01
7) Phenol-d6	6.49	99	937077	81.80	ng	0.00
23) Nitrobenzene-d5	7.40	82	903113	78.31	ng	0.00
42) 2,4,6-Tribromophenol	10.67	330	290028	78.20	ng	-0.01
45) 2-Fluorobiphenyl	9.19	172	1764730	78.33	ng	-0.01
79) Terphenyl-d14	12.94	244	1806193	81.96	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.59	88	147904	45.033	ng	95
3) Pyridine	3.35	79	383249	45.873	ng	96
4) n-Nitrosodimethylamine	3.31	42	165224	47.072	ng	88
6) Aniline	6.50	93	554232	40.668	ng	# 69
8) 2-Chlorophenol	6.63	128	450207	40.592	ng	96
9) Benzaldehyde	6.39	77	244415	40.818	ng	99
10) Phenol	6.50	94	510451	40.612	ng	92
11) bis(2-Chloroethyl)ether	6.57	93	399915	40.415	ng	99
12) 1,3-Dichlorobenzene	6.77	146	511355	39.589	ng	98
13) 1,4-Dichlorobenzene	6.85	146	513124	38.600	ng	98
14) 1,2-Dichlorobenzene	7.00	146	479753	38.546	ng	95
15) Benzyl Alcohol	6.99	79	370072	38.320	ng	96
16) 2,2'-oxybis(1-Chloropropan	7.11	45	495737	39.019	ng	# 52
17) 2-Methylphenol	7.10	107	339141	38.573	ng	# 87
18) Hexachloroethane	7.35	117	184822	39.593	ng	91
19) n-Nitroso-di-n-propylamine	7.25	70	308848	39.097	ng	99
20) 3+4-Methylphenols	7.26	107	445572	39.631	ng	# 76
22) Acetophenone	7.25	105	636966	39.365	ng	# 98
24) Nitrobenzene	7.42	77	451806	39.226	ng	98
25) Isophorone	7.66	82	712219	39.365	ng	99
26) 2-Nitrophenol	7.74	139	246835	40.101	ng	94
27) 2,4-Dimethylphenol	7.78	122	335714	39.586	ng	94
28) bis(2-Chloroethoxy)methane	7.86	93	488768	39.675	ng	98
29) 2,4-Dichlorophenol	7.99	162	383880	40.449	ng	98
30) 1,2,4-Trichlorobenzene	8.06	180	433787	39.764	ng	98
31) Naphthalene	8.14	128	1216033	39.131	ng	99
32) Benzoic acid	7.93	122	183888	38.013	ng	99
33) 4-Chloroaniline	8.19	127	521862	38.568	ng	98
34) Hexachlorobutadiene	8.25	225	256341	40.044	ng	99
35) Caprolactam	8.58	113	121427	36.269	ng	98
36) 4-Chloro-3-methylphenol	8.69	107	382910	38.113	ng	96
37) 2-Methylnaphthalene	8.83	142	823670	38.718	ng	97
38) 1-Methylnaphthalene	8.93	142	783294	38.414	ng	97
40) 1,2,4,5-Tetrachlorobenzene	9.00	216	417612	39.916	ng	98

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF021219\
 Data File : BF112502.D
 Acq On : 12 Feb 2019 10:14
 Operator : JU/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Sohil
 2/13/2019 11:11:57 AM

Quant Time: Feb 13 09:00:57 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF012519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 28 10:38:18 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	8.98	237	136239	39.781	nq	99
43) 2,4,6-Trichlorophenol	9.12	196	253886	39.368	nq	98
44) 2,4,5-Trichlorophenol	9.17	196	261199	38.753	nq	96
46) 1,1'-Biphenyl	9.29	154	1099650	39.474	nq	99
47) 2-Chloronaphthalene	9.32	162	853153	39.493	nq	95
48) 2-Nitroaniline	9.42	65	229432	38.966	nq	90
49) Acenaphthylene	9.74	152	1215084	38.376	nq	98
50) Dimethylphthalate	9.59	163	944232	38.139	nq	99
51) 2,6-Dinitrotoluene	9.66	165	212554	38.335	nq #	78
52) Acenaphthene	9.91	154	730876	38.641	nq	99
53) 3-Nitroaniline	9.84	138	226406	37.690	nq	91
54) 2,4-Dinitrophenol	9.96	184	68068m	38.252	nq	
55) Dibenzofuran	10.08	168	1088697	37.665	nq	93
56) 4-Nitrophenol	10.02	139	121707	38.323	nq	92
57) 2,4-Dinitrotoluene	10.07	165	263198	37.286	nq #	74
58) Fluorene	10.42	166	843511	38.997	nq	98
59) 2,3,4,6-Tetrachlorophenol	10.21	232	191052	37.508	nq #	89
60) Diethylphthalate	10.29	149	945003	38.463	nq	98
61) 4-Chlorophenyl-phenylether	10.41	204	456000	38.755	nq	91
62) 4-Nitroaniline	10.46	138	207407	35.201	nq	93
63) Azobenzene	10.57	77	847908	39.241	nq	96
65) 4,6-Dinitro-2-methylphenol	10.49	198	116475	35.912	nq	94
66) n-Nitrosodiphenylamine	10.53	169	794837	40.768	nq	100
67) 4-Bromophenyl-phenylether	10.90	248	276812	40.685	nq	96
68) Hexachlorobenzene	10.98	284	291858	39.982	nq	99
69) Atrazine	11.06	200	218347	34.708	nq	96
70) Pentachlorophenol	11.18	266	102912	36.232	nq	99
71) Phenanthrene	11.39	178	1170115	38.907	nq	99
72) Anthracene	11.44	178	1192447	39.804	nq	100
73) Carbazole	11.60	167	1161270	39.297	nq	99
74) Di-n-butylphthalate	11.91	149	1461097	39.834	nq	99
75) Fluoranthene	12.57	202	1184142	36.710	nq	99
77) Benzidine	12.69	184	378678m	33.148	nq	
78) Pyrene	12.81	202	1188351	41.354	nq	98
80) Butylbenzylphthalate	13.41	149	580780	41.773	nq	96
81) Benzo(a)anthracene	14.00	228	966064	39.594	nq	97
82) 3,3'-Dichlorobenzidine	13.95	252	360932	39.527	nq	98
83) Chrysene	14.03	228	930558	39.955	nq	98
84) Bis(2-ethylhexyl)phthalate	13.96	149	796409	40.355	nq #	96
85) Di-n-octyl phthalate	14.57	149	1187631	39.114	nq	98
86) Indeno(1,2,3-cd)pyrene	16.97	276	760472	41.207	nq	97
88) Benzo(b)fluoranthene	15.05	252	790427	38.834	nq	99
89) Benzo(k)fluoranthene	15.08	252	777780	39.894	nq	97
90) Benzo(a)pyrene	15.42	252	721727	39.408	nq	99
91) Dibenzo(a,h)anthracene	16.97	278	640497	43.393	nq	96
92) Benzo(g,h,i)perylene	17.42	276	614737	44.144	nq	96

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

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF021219\
 Data File : BF112502.D
 Acq On : 12 Feb 2019 10:14
 Operator : JU/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampled :
 SSTDCCC040

Manual Integrations
 APPROVED

Sohil
 2/13/2019 11:11:57 AM

Quant Time: Feb 13 09:00:57 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF012519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 28 10:38:18 2019
 Response via : Initial Calibration

