

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071119\
 Data File : BF115493.D
 Acq On : 11 Jul 2019 10:42
 Operator : HP/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 LabSampleId :
 SSTDCCC040

Quant Time: Jul 12 03:22:16 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF070519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 10 16:47:05 2019
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
 Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	20.000	20.000	0.0	109	0.00
2	1,4-Dioxane	40.000	35.877	10.3	102	0.00
3	Pyridine	40.000	35.598	11.0	102	0.00
4	n-Nitrosodimethylamine	40.000	32.221	19.4	91	0.00
5 S	2-Fluorophenol	80.000	72.863	8.9	102	0.00
6	Aniline	40.000	36.892	7.8	102	0.00
7 S	Phenol-d6	80.000	73.977	7.5	103	0.00
8	2-Chlorophenol	40.000	38.276	4.3	105	0.00
9	Benzaldehyde	40.000	38.488	3.8	107	0.00
10 C	Phenol	40.000	36.331	9.2	101	0.00
11	bis(2-Chloroethyl)ether	40.000	36.457	8.9	101	0.00
12	1,3-Dichlorobenzene	40.000	38.325	4.2	106	0.00
13 C	1,4-Dichlorobenzene	40.000	38.495	3.8	108	0.00
14	1,2-Dichlorobenzene	40.000	38.434	3.9	107	0.00
15	Benzyl Alcohol	40.000	38.054	4.9	105	0.00
16	2,2'-oxybis(1-Chloropropane)	40.000	35.285	11.8	99	0.00
17	2-Methylphenol	40.000	38.129	4.7	106	0.00
18	Hexachloroethane	40.000	38.387	4.0	107	0.00
19 P	n-Nitroso-di-n-propylamine	40.000	35.495	11.3	99	0.00
20	3+4-Methylphenols	40.000	37.753	5.6	102	0.00
21 I	Naphthalene-d8	20.000	20.000	0.0	108	0.00
22	Acetophenone	40.000	36.327	9.2	100	0.00
23 S	Nitrobenzene-d5	80.000	79.114	1.1	109	0.00
24	Nitrobenzene	40.000	38.213	4.5	105	0.00
25	Isophorone	40.000	36.249	9.4	102	0.00
26 C	2-Nitrophenol	40.000	48.169	-20.4#	131	0.00
27	2,4-Dimethylphenol	40.000	35.186	12.0	99	0.00
28	bis(2-Chloroethoxy)methane	40.000	36.583	8.5	103	0.00
29 C	2,4-Dichlorophenol	40.000	38.953	2.6	108	0.00
30	1,2,4-Trichlorobenzene	40.000	39.040	2.4	109	0.00
31	Naphthalene	40.000	37.959	5.1	105	0.00
32	Benzoic acid	40.000	32.603	18.5	85	0.00
33	4-Chloroaniline	40.000	38.345	4.1	106	0.00
34 C	Hexachlorobutadiene	40.000	39.353	1.6	109	0.00
35	Caprolactam	40.000	37.513	6.2	102	0.00
36 C	4-Chloro-3-methylphenol	40.000	38.285	4.3	105	0.00
37	2-Methylnaphthalene	40.000	37.988	5.0	104	0.00
38	1-Methylnaphthalene	40.000	37.804	5.5	104	0.00
39 I	Acenaphthene-d10	20.000	20.000	0.0	106	0.00
40	1,2,4,5-Tetrachlorobenzene	40.000	40.447	-1.1	109	0.00
41 P	Hexachlorocyclopentadiene	40.000	37.101	7.2	101	0.00
42 S	2,4,6-Tribromophenol	80.000	82.701	-3.4	109	0.00
43 C	2,4,6-Trichlorophenol	40.000	40.693	-1.7	109	0.00
44	2,4,5-Trichlorophenol	40.000	40.413	-1.0	108	0.00

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071119\
 Data File : BF115493.D
 Acq On : 11 Jul 2019 10:42
 Operator : HP/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 LabSampleId :
 SSTDCCC040

Quant Time: Jul 12 03:22:16 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF070519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 10 16:47:05 2019
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
 Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
45 S	2-Fluorobiphenyl	80.000	80.898	-1.1	104	0.00
46	1,1'-Biphenyl	40.000	38.591	3.5	105	0.00
47	2-Chloronaphthalene	40.000	38.932	2.7	105	0.00
48	2-Nitroaniline	40.000	42.074	-5.2	109	0.00
49	Acenaphthylene	40.000	38.101	4.7	103	0.00
50	Dimethylphthalate	40.000	39.279	1.8	105	0.00
51	2,6-Dinitrotoluene	40.000	41.654	-4.1	110	0.00
52 C	Acenaphthene	40.000	38.839	2.9	104	0.00
53	3-Nitroaniline	40.000	42.011	-5.0	109	0.00
54 P	2,4-Dinitrophenol	40.000	49.905	-24.8	152	0.00
55	Dibenzofuran	40.000	38.325	4.2	103	0.00
56 P	4-Nitrophenol	40.000	41.211	-3.0	106	0.00
57	2,4-Dinitrotoluene	40.000	43.778	-9.4	113	0.00
58	Fluorene	40.000	35.858	10.4	102	0.00
59	2,3,4,6-Tetrachlorophenol	40.000	40.640	-1.6	107	0.00
60	Diethylphthalate	40.000	38.203	4.5	101	0.00
61	4-Chlorophenyl-phenylether	40.000	38.235	4.4	105	0.00
62	4-Nitroaniline	40.000	41.680	-4.2	107	0.00
63	Azobenzene	40.000	35.998	10.0	95	0.00
64 I	Phenanthrene-d10	20.000	20.000	0.0	103	0.00
65	4,6-Dinitro-2-methylphenol	40.000	48.292	-20.7	137	0.00
66 c	n-Nitrosodiphenylamine	40.000	38.743	3.1	103	0.00
67	4-Bromophenyl-phenylether	40.000	39.596	1.0	106	0.00
68	Hexachlorobenzene	40.000	40.236	-0.6	107	0.00
69	Atrazine	40.000	40.067	-0.2	102	0.00
70 C	Pentachlorophenol	40.000	40.646	-1.6	104	0.00
71	Phenanthrene	40.000	37.682	5.8	100	0.00
72	Anthracene	40.000	38.521	3.7	102	0.00
73	Carbazole	40.000	37.465	6.3	98	0.00
74	Di-n-butylphthalate	40.000	37.375	6.6	98	0.00
75 C	Fluoranthene	40.000	36.991	7.5	97	0.00
76 I	Chrysene-d12	20.000	20.000	0.0	90	0.00
77	Benzidine	40.000	36.001	10.0	83	0.00
78	Pyrene	40.000	41.134	-2.8	96	0.00
79 S	Terphenyl-d14	80.000	82.195	-2.7	98	0.00
80	Butylbenzylphthalate	40.000	40.705	-1.8	93	0.00
81	Benzo(a)anthracene	40.000	39.909	0.2	92	0.00
82	3,3'-Dichlorobenzidine	40.000	39.855	0.4	87	0.00
83	Chrysene	40.000	39.301	1.7	90	0.00
84	Bis(2-ethylhexyl)phthalate	40.000	40.307	-0.8	95	-0.01
85 c	Di-n-octyl phthalate	40.000	37.844	5.4	86	-0.03
86	Indeno(1,2,3-cd)pyrene	40.000	44.805	-12.0	99	-0.03
87 I	Perylene-d12	20.000	20.000	0.0	84	-0.02

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071119\
Data File : BF115493.D
Acq On : 11 Jul 2019 10:42
Operator : HP/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Jul 12 03:22:16 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF070519.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jul 10 16:47:05 2019
Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
88	Benzo(b)fluoranthene	40.000	39.062	2.3	83	-0.03
89	Benzo(k)fluoranthene	40.000	38.235	4.4	86	-0.02
90 C	Benzo(a)pyrene	40.000	39.049	2.4	85	-0.02
91	Dibenzo(a,h)anthracene	40.000	42.407	-6.0	98	-0.04
92	Benzo(q,h,i)perylene	40.000	43.955	-9.9	105	-0.03

(#) = Out of Range

SPCC's out = 0 CCC's out = 1