

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF040320\
 Data File : BF119759.D
 Acq On : 4 Apr 2020 11:38
 Operator : MA/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 100 Sample Multiplier: 1

Instrument :
 BNA_F
 LabSampleId :
 SSTDCCC040

Quant Time: Apr 06 02:54:55 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF031820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 01 11:59:45 2020
 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	71	0.02
2	1,4-Dioxane	40.000	36.023	9.9	62	0.04
3	Pyridine	40.000	35.198	12.0	60	0.05
4	n-Nitrosodimethylamine	40.000	33.265	16.8	57	0.05
5 S	2-Fluorophenol	80.000	78.669	1.7	67	0.03
6	Aniline	40.000	36.116	9.7	60	0.02
7 S	Phenol-d6	80.000	76.720	4.1	64	0.02
8	2-Chlorophenol	40.000	39.806	0.5	67	0.02
9	Benzaldehyde	40.000	35.353	11.6	61	0.02
10 C	Phenol	40.000	40.411	-1.0	68	0.02
11	bis(2-Chloroethyl)ether	40.000	38.885	2.8	66	0.02
12	1,3-Dichlorobenzene	40.000	40.391	-1.0	69	0.02
13 C	1,4-Dichlorobenzene	40.000	40.186	-0.5	69	0.02
14	1,2-Dichlorobenzene	40.000	40.474	-1.2	68	0.02
15	Benzyl Alcohol	40.000	37.227	6.9	62	0.02
16	2,2'-oxybis(1-Chloropropane)	40.000	34.616	13.5	59	0.02
17	2-Methylphenol	40.000	38.596	3.5	66	0.02
18	Hexachloroethane	40.000	32.690	18.3	56	0.02
19 P	n-Nitroso-di-n-propylamine	40.000	37.071	7.3	63	0.02
20	3+4-Methylphenols	40.000	37.665	5.8	63	0.02
21 I	Naphthalene-d8	20.000	20.000	0.0	67	0.02
22	Acetophenone	40.000	39.326	1.7	64	0.02
23 S	Nitrobenzene-d5	80.000	83.289	-4.1	67	0.02
24	Nitrobenzene	40.000	40.123	-0.3	65	0.02
25	Isophorone	40.000	38.289	4.3	63	0.02
26 C	2-Nitrophenol	40.000	35.407	11.5	57	0.02
27	2,4-Dimethylphenol	40.000	36.467	8.8	59	0.02
28	bis(2-Chloroethoxy)methane	40.000	39.937	0.2	65	0.02
29 C	2,4-Dichlorophenol	40.000	38.936	2.7	64	0.02
30	1,2,4-Trichlorobenzene	40.000	40.794	-2.0	67	0.02
31	Naphthalene	40.000	39.868	0.3	64	0.02
32	Benzoic acid	40.000	30.834	22.9	50	0.00
33	4-Chloroaniline	40.000	37.183	7.0	60	0.02
34 C	Hexachlorobutadiene	40.000	43.519	-8.8	71	0.02
35	Caprolactam	40.000	33.850	15.4	54	0.02
36 C	4-Chloro-3-methylphenol	40.000	36.149	9.6	59	0.02
37	2-Methylnaphthalene	40.000	39.093	2.3	64	0.02
38	1-Methylnaphthalene	40.000	38.348	4.1	62	0.02
39 I	Acenaphthene-d10	20.000	20.000	0.0	59	0.02
40	1,2,4,5-Tetrachlorobenzene	40.000	46.415	-16.0	66	0.02
41 P	Hexachlorocyclopentadiene	40.000	7.863	80.3#	11	0.02
42 S	2,4,6-Tribromophenol	80.000	79.848	0.2	56	0.02
43 C	2,4,6-Trichlorophenol	40.000	41.569	-3.9	59	0.02
44	2,4,5-Trichlorophenol	40.000	39.915	0.2	56	0.02

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF040320\
 Data File : BF119759.D
 Acq On : 4 Apr 2020 11:38
 Operator : MA/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 100 Sample Multiplier: 1

Instrument :
 BNA_F
 LabSampleId :
 SSTDCCC040

Quant Time: Apr 06 02:54:55 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF031820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 01 11:59:45 2020
 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	89.640	-12.1	66	0.02
46	1,1'-Biphenyl	40.000	44.988	-12.5	63	0.02
47	2-Chloronaphthalene	40.000	42.204	-5.5	60	0.02
48	2-Nitroaniline	40.000	45.213	-13.0	62	0.02
49	Acenaphthylene	40.000	40.444	-1.1	57	0.02
50	Dimethylphthalate	40.000	44.668	-11.7	63	0.01
51	2,6-Dinitrotoluene	40.000	38.068	4.8	53	0.02
52 C	Acenaphthene	40.000	38.555	3.6	55	0.02
53	3-Nitroaniline	40.000	42.544	-6.4	59	0.02
54 P	2,4-Dinitrophenol	40.000	9.967	75.1#	5	0.02
55	Dibenzofuran	40.000	39.544	1.1	56	0.02
56 P	4-Nitrophenol	40.000	29.891	25.3#	41	0.02
57	2,4-Dinitrotoluene	40.000	35.775	10.6	50	0.02
58	Fluorene	40.000	39.344	1.6	55	0.02
59	2,3,4,6-Tetrachlorophenol	40.000	36.504	8.7	50	0.02
60	Diethylphthalate	40.000	42.972	-7.4	61	0.01
61	4-Chlorophenyl-phenylether	40.000	43.040	-7.6	61	0.02
62	4-Nitroaniline	40.000	42.204	-5.5	58	0.02
63	Azobenzene	40.000	36.244	9.4	51	0.02
64 I	Phenanthrene-d10	20.000	20.000	0.0	49	0.02
65	4,6-Dinitro-2-methylphenol	40.000	5.987	85.0#	7	0.01
66 c	n-Nitrosodiphenylamine	40.000	44.975	-12.4	54	0.02
67	4-Bromophenyl-phenylether	40.000	45.820	-14.6	54	0.02
68	Hexachlorobenzene	40.000	43.772	-9.4	53	0.02
69	Atrazine	40.000	36.838	7.9	44	0.01
70 C	Pentachlorophenol	40.000	38.730	3.2	46	0.02
71	Phenanthrene	40.000	41.652	-4.1	50	0.02
72	Anthracene	40.000	41.577	-3.9	49	0.02
73	Carbazole	40.000	40.033	-0.1	47	0.02
74	Di-n-butylphthalate	40.000	47.199	-18.0	56	0.02
75 C	Fluoranthene	40.000	41.052	-2.6	49	0.02
76 I	Chrysene-d12	20.000	20.000	0.0	54	0.02
77	Benzidine	40.000	33.029	17.4	44	0.02
78	Pyrene	40.000	36.221	9.4	49	0.02
79 S	Terphenyl-d14	80.000	80.378	-0.5	55	0.02
80	Butylbenzylphthalate	40.000	41.362	-3.4	56	0.02
81	Benzo(a)anthracene	40.000	40.218	-0.5	52	0.02
82	3,3'-Dichlorobenzidine	40.000	42.264	-5.7	54	0.02
83	Chrysene	40.000	39.351	1.6	51	0.02
84	Bis(2-ethylhexyl)phthalate	40.000	46.077	-15.2	62	0.02
85 c	Di-n-octyl phthalate	40.000	47.847	-19.6	62	0.02
86	Indeno(1,2,3-cd)pyrene	40.000	32.347	19.1	45	0.05
87 I	Perylene-d12	20.000	20.000	0.0	47	0.03

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF040320\
Data File : BF119759.D
Acq On : 4 Apr 2020 11:38
Operator : MA/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 100 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Apr 06 02:54:55 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF031820.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Apr 01 11:59:45 2020
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	41.255	-3.1	48	0.02
89	Benzo(k)fluoranthene	40.000	42.848	-7.1	48	0.02
90 C	Benzo(a)pyrene	40.000	40.040	-0.1	46	0.02
91	Dibenzo(a,h)anthracene	40.000	38.031	4.9	46	0.04
92	Benzo(q,h,i)perylene	40.000	35.091	12.3	44	0.04

(#) = Out of Range

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