

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF090619\
 Data File : BF116873.D
 Acq On : 7 Sep 2019 5:48
 Operator : HP/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 LabSampleId :
 SSTDCCC040

Quant Time: Sep 07 06:22:03 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF083019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 30 20:02:30 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	53	0.00
2	1,4-Dioxane	40.000	39.119	2.2	52	-0.02
3	Pyridine	40.000	37.680	5.8	49	-0.02
4	n-Nitrosodimethylamine	40.000	36.773	8.1	48	-0.02
5 S	2-Fluorophenol	80.000	87.824	-9.8	58	0.00
6	Aniline	40.000	39.638	0.9	51	0.01
7 S	Phenol-d6	80.000	85.843	-7.3	56	0.00
8	2-Chlorophenol	40.000	43.390	-8.5	57	0.00
9	Benzaldehyde	40.000	37.842	5.4	53	0.00
10 C	Phenol	40.000	42.298	-5.7	55	0.01
11	bis(2-Chloroethyl)ether	40.000	40.113	-0.3	53	0.00
12	1,3-Dichlorobenzene	40.000	42.448	-6.1	56	0.01
13 C	1,4-Dichlorobenzene	40.000	43.817	-9.5	57	0.00
14	1,2-Dichlorobenzene	40.000	44.080	-10.2	58	0.01
15	Benzyl Alcohol	40.000	38.479	3.8	50	0.00
16	2,2'-oxybis(1-Chloropropane)	40.000	34.652	13.4	45	0.00
17	2-Methylphenol	40.000	43.145	-7.9	57	0.01
18	Hexachloroethane	40.000	35.397	11.5	46	0.01
19 P	n-Nitroso-di-n-propylamine	40.000	40.040	-0.1	53	0.00
20	3+4-Methylphenols	40.000	44.856	-12.1	58	0.01
21 I	Naphthalene-d8	20.000	20.000	0.0	52	0.00
22	Acetophenone	40.000	45.124	-12.8	58	0.01
23 S	Nitrobenzene-d5	80.000	92.686	-15.9	60	0.00
24	Nitrobenzene	40.000	45.618	-14.0	58	0.00
25	Isophorone	40.000	39.767	0.6	52	0.01
26 C	2-Nitrophenol	40.000	47.204	-18.0	61	0.01
27	2,4-Dimethylphenol	40.000	40.851	-2.1	52	0.00
28	bis(2-Chloroethoxy)methane	40.000	41.426	-3.6	53	0.00
29 C	2,4-Dichlorophenol	40.000	43.864	-9.7	56	0.00
30	1,2,4-Trichlorobenzene	40.000	43.234	-8.1	55	0.01
31	Naphthalene	40.000	43.110	-7.8	55	0.00
32	Benzoic acid	40.000	34.014	15.0	51	-0.01
33	4-Chloroaniline	40.000	42.114	-5.3	54	0.01
34 C	Hexachlorobutadiene	40.000	45.366	-13.4	58	0.01
35	Caprolactam	40.000	38.570	3.6	50	0.00
36 C	4-Chloro-3-methylphenol	40.000	44.102	-10.3	56	0.01
37	2-Methylnaphthalene	40.000	43.831	-9.6	55	0.01
38	1-Methylnaphthalene	40.000	42.659	-6.6	55	0.00
39 I	Acenaphthene-d10	20.000	20.000	0.0	51	0.01
40	1,2,4,5-Tetrachlorobenzene	40.000	47.046	-17.6	59	0.01
41 P	Hexachlorocyclopentadiene	40.000	5.033	87.4#	6	0.01
42 S	2,4,6-Tribromophenol	80.000	93.330	-16.7	60	0.01
43 C	2,4,6-Trichlorophenol	40.000	44.968	-12.4	56	0.01
44	2,4,5-Trichlorophenol	40.000	45.738	-14.3	59	0.01

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF090619\
 Data File : BF116873.D
 Acq On : 7 Sep 2019 5:48
 Operator : HP/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 LabSampleId :
 SSTDCCC040

Quant Time: Sep 07 06:22:03 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF083019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 30 20:02:30 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	93.655	-17.1	60	0.01
46	1,1'-Biphenyl	40.000	45.174	-12.9	57	0.01
47	2-Chloronaphthalene	40.000	43.981	-10.0	55	0.01
48	2-Nitroaniline	40.000	51.668	-29.2#	66	0.00
49	Acenaphthylene	40.000	43.401	-8.5	54	0.01
50	Dimethylphthalate	40.000	44.930	-12.3	57	0.01
51	2,6-Dinitrotoluene	40.000	47.569	-18.9	59	0.01
52 C	Acenaphthene	40.000	41.799	-4.5	53	0.01
53	3-Nitroaniline	40.000	48.675	-21.7	60	0.01
54 P	2,4-Dinitrophenol	40.000	12.399	69.0#	6	0.01
55	Dibenzofuran	40.000	43.565	-8.9	54	0.01
56 P	4-Nitrophenol	40.000	39.799	0.5	50	0.01
57	2,4-Dinitrotoluene	40.000	50.374	-25.9#	63	0.01
58	Fluorene	40.000	44.412	-11.0	57	0.01
59	2,3,4,6-Tetrachlorophenol	40.000	43.898	-9.7	56	0.01
60	Diethylphthalate	40.000	44.073	-10.2	55	0.01
61	4-Chlorophenyl-phenylether	40.000	46.405	-16.0	59	0.01
62	4-Nitroaniline	40.000	50.957	-27.4#	65	0.01
63	Azobenzene	40.000	41.561	-3.9	53	0.01
64 I	Phenanthrene-d10	20.000	20.000	0.0	47	0.02
65	4,6-Dinitro-2-methylphenol	40.000	14.711	63.2#	11	0.01
66 c	n-Nitrosodiphenylamine	40.000	47.173	-17.9	55	0.01
67	4-Bromophenyl-phenylether	40.000	47.422	-18.6	55	0.01
68	Hexachlorobenzene	40.000	46.600	-16.5	54	0.01
69	Atrazine	40.000	45.069	-12.7	53	0.01
70 C	Pentachlorophenol	40.000	37.505	6.2	44	0.02
71	Phenanthrene	40.000	43.439	-8.6	50	0.01
72	Anthracene	40.000	43.614	-9.0	51	0.01
73	Carbazole	40.000	42.179	-5.4	50	0.01
74	Di-n-butylphthalate	40.000	46.360	-15.9	54	0.01
75 C	Fluoranthene	40.000	40.833	-2.1	48	0.02
76 I	Chrysene-d12	20.000	20.000	0.0	56	0.02
77	Benzidine	40.000	38.437	3.9	55	0.02
78	Pyrene	40.000	36.594	8.5	48	0.02
79 S	Terphenyl-d14	80.000	80.688	-0.9	54	0.02
80	Butylbenzylphthalate	40.000	39.357	1.6	54	0.01
81	Benzo(a)anthracene	40.000	44.632	-11.6	62	0.02
82	3,3'-Dichlorobenzidine	40.000	46.059	-15.1	62	0.01
83	Chrysene	40.000	42.257	-5.6	59	0.02
84	Bis(2-ethylhexyl)phthalate	40.000	44.295	-10.7	62	0.02
85 c	Di-n-octyl phthalate	40.000	45.204	-13.0	65	0.02
86	Indeno(1,2,3-cd)pyrene	40.000	38.256	4.4	53	0.03
87 I	Perylene-d12	20.000	20.000	0.0	63	0.02

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF090619\
Data File : BF116873.D
Acq On : 7 Sep 2019 5:48
Operator : HP/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Sep 07 06:22:03 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF083019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Aug 30 20:02:30 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	42.192	-5.5	68	0.02
89	Benzo(k)fluoranthene	40.000	44.602	-11.5	65	0.02
90 C	Benzo(a)pyrene	40.000	42.509	-6.3	66	0.02
91	Dibenzo(a,h)anthracene	40.000	37.092	7.3	56	0.03
92	Benzo(q,h,i)perylene	40.000	32.023	19.9	48	0.04

(#) = Out of Range

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