

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF122419\
 Data File : BF118307.D
 Acq On : 24 Dec 2019 15:55
 Operator : JU
 Sample : SSTDICV040
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 ICVBF122419

Quant Time: Dec 24 16:34:29 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF122419.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 24 15:58:52 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	107	0.00
2	1,4-Dioxane	40.000	37.585	6.0	105	0.00
3	Pyridine	40.000	38.509	3.7	107	0.00
4	n-Nitrosodimethylamine	40.000	38.528	3.7	104	0.00
5 S	2-Fluorophenol	80.000	75.899	5.1	104	0.00
6	Aniline	40.000	38.076	4.8	105	0.00
7 S	Phenol-d6	80.000	76.366	4.5	107	0.00
8	2-Chlorophenol	40.000	38.067	4.8	105	0.00
9	Benzaldehyde	40.000	35.640	10.9	103	0.00
10 C	Phenol	40.000	38.444	3.9	107	0.00
11	bis(2-Chloroethyl)ether	40.000	37.914	5.2	105	0.00
12	1,3-Dichlorobenzene	40.000	37.912	5.2	106	0.00
13 C	1,4-Dichlorobenzene	40.000	37.937	5.2	105	0.00
14	1,2-Dichlorobenzene	40.000	38.125	4.7	105	0.00
15	Benzyl Alcohol	40.000	38.560	3.6	106	0.00
16	2,2'-oxybis(1-Chloropropane)	40.000	37.978	5.1	106	0.00
17	2-Methylphenol	40.000	37.812	5.5	105	0.00
18	Hexachloroethane	40.000	37.848	5.4	105	0.00
19 P	n-Nitroso-di-n-propylamine	40.000	38.069	4.8	107	0.00
20	3+4-Methylphenols	40.000	38.221	4.4	104	0.00
21 I	Naphthalene-d8	20.000	20.000	0.0	104	0.00
22	Acetophenone	40.000	38.016	5.0	105	0.00
23 S	Nitrobenzene-d5	80.000	76.329	4.6	106	0.00
24	Nitrobenzene	40.000	38.255	4.4	106	0.00
25	Isophorone	40.000	37.584	6.0	105	0.00
26 C	2-Nitrophenol	40.000	38.581	3.5	107	0.00
27	2,4-Dimethylphenol	40.000	37.483	6.3	105	0.00
28	bis(2-Chloroethoxy)methane	40.000	38.002	5.0	106	0.00
29 C	2,4-Dichlorophenol	40.000	37.189	7.0	104	0.00
30	1,2,4-Trichlorobenzene	40.000	37.528	6.2	105	0.00
31	Naphthalene	40.000	38.306	4.2	106	0.00
32	Benzoic acid	40.000	38.225	4.4	110	0.01
33	4-Chloroaniline	40.000	38.636	3.4	107	0.00
34 C	Hexachlorobutadiene	40.000	37.994	5.0	105	0.00
35	Caprolactam	40.000	38.904	2.7	106	0.00
36 C	4-Chloro-3-methylphenol	40.000	37.877	5.3	106	0.00
37	2-Methylnaphthalene	40.000	37.860	5.4	104	0.00
38	1-Methylnaphthalene	40.000	37.674	5.8	105	0.00
39 I	Acenaphthene-d10	20.000	20.000	0.0	104	0.00
40	1,2,4,5-Tetrachlorobenzene	40.000	38.430	3.9	105	0.00
41 P	Hexachlorocyclopentadiene	40.000	35.961	10.1	105	0.00
42 S	2,4,6-Tribromophenol	80.000	76.525	4.3	103	0.00
43 C	2,4,6-Trichlorophenol	40.000	38.605	3.5	106	0.00
44	2,4,5-Trichlorophenol	40.000	38.520	3.7	105	0.00

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF122419\
 Data File : BF118307.D
 Acq On : 24 Dec 2019 15:55
 Operator : JU
 Sample : SSTDICV040
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 ICVBF122419

Quant Time: Dec 24 16:34:29 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF122419.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 24 15:58:52 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	76.885	3.9	103	0.00
46	1,1'-Biphenyl	40.000	38.622	3.4	105	0.00
47	2-Chloronaphthalene	40.000	38.597	3.5	104	0.00
48	2-Nitroaniline	40.000	38.120	4.7	104	0.00
49	Acenaphthylene	40.000	38.691	3.3	104	0.00
50	Dimethylphthalate	40.000	37.952	5.1	104	0.00
51	2,6-Dinitrotoluene	40.000	38.745	3.1	106	0.00
52 C	Acenaphthene	40.000	38.938	2.7	106	0.00
53	3-Nitroaniline	40.000	38.341	4.1	103	0.00
54 P	2,4-Dinitrophenol	40.000	37.757	5.6	106	0.00
55	Dibenzofuran	40.000	38.676	3.3	104	0.00
56 P	4-Nitrophenol	40.000	39.547	1.1	106	0.00
57	2,4-Dinitrotoluene	40.000	39.059	2.4	106	0.00
58	Fluorene	40.000	38.244	4.4	103	0.00
59	2,3,4,6-Tetrachlorophenol	40.000	38.168	4.6	102	0.00
60	Diethylphthalate	40.000	38.715	3.2	105	0.00
61	4-Chlorophenyl-phenylether	40.000	38.132	4.7	102	0.00
62	4-Nitroaniline	40.000	38.647	3.4	103	0.00
63	Azobenzene	40.000	38.624	3.4	105	0.00
64 I	Phenanthrene-d10	20.000	20.000	0.0	104	0.00
65	4,6-Dinitro-2-methylphenol	40.000	39.153	2.1	105	0.00
66 c	n-Nitrosodiphenylamine	40.000	38.248	4.4	104	0.00
67	4-Bromophenyl-phenylether	40.000	37.617	6.0	104	0.00
68	Hexachlorobenzene	40.000	37.333	6.7	103	0.00
69	Atrazine	40.000	37.940	5.2	100	0.00
70 C	Pentachlorophenol	40.000	39.102	2.2	105	0.00
71	Phenanthrene	40.000	38.022	4.9	104	0.00
72	Anthracene	40.000	37.890	5.3	103	0.00
73	Carbazole	40.000	38.531	3.7	106	0.00
74	Di-n-butylphthalate	40.000	38.560	3.6	104	0.00
75 C	Fluoranthene	40.000	38.659	3.4	104	0.00
76 I	Chrysene-d12	20.000	20.000	0.0	112	0.00
77	Benzidine	40.000	33.938	15.2	92	0.00
78	Pyrene	40.000	34.580	13.6	106	0.00
79 S	Terphenyl-d14	80.000	69.826	12.7	105	0.00
80	Butylbenzylphthalate	40.000	36.694	8.3	110	0.00
81	Benzo(a)anthracene	40.000	37.340	6.6	111	0.00
82	3,3'-Dichlorobenzidine	40.000	37.914	5.2	108	0.00
83	Chrysene	40.000	37.377	6.6	111	0.00
84	Bis(2-ethylhexyl)phthalate	40.000	38.079	4.8	113	0.00
85 c	Di-n-octyl phthalate	40.000	41.205	-3.0	123	0.00
86	Indeno(1,2,3-cd)pyrene	40.000	37.107	7.2	118	0.02
87 I	Perylene-d12	20.000	20.000	0.0	123	0.01

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF122419\
Data File : BF118307.D
Acq On : 24 Dec 2019 15:55
Operator : JU
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
ICVBF122419

Quant Time: Dec 24 16:34:29 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF122419.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Dec 24 15:58:52 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	36.861	7.8	122	0.01
89	Benzo(k)fluoranthene	40.000	37.569	6.1	118	0.01
90 C	Benzo(a)pyrene	40.000	37.459	6.4	121	0.02
91	Dibenzo(a,h)anthracene	40.000	36.591	8.5	121	0.02
92	Benzo(q,h,i)perylene	40.000	35.868	10.3	118	0.02

(#) = Out of Range

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