

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF081920\
 Data File : BF121332.D
 Acq On : 19 Aug 2020 22:02
 Operator : JU/CG
 Sample : SSTDICV040
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 ICVBF081920

Quant Time: Aug 20 05:30:05 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF081920.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 20 04:56:10 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	106	0.00
2	1,4-Dioxane	40.000	36.678	8.3	104	0.00
3	Pyridine	40.000	37.756	5.6	104	0.00
4	n-Nitrosodimethylamine	40.000	37.939	5.2	106	0.00
5 S	2-Fluorophenol	80.000	72.366	9.5	104	0.00
6	Aniline	40.000	36.615	8.5	104	0.00
7 S	Phenol-d6	80.000	72.140	9.8	103	0.00
8	2-Chlorophenol	40.000	37.013	7.5	105	0.00
9	Benzaldehyde	40.000	38.138	4.7	104	0.00
10 C	Phenol	40.000	34.423	13.9	104	0.00
11	bis(2-Chloroethyl)ether	40.000	36.455	8.9	104	0.00
12	1,3-Dichlorobenzene	40.000	36.417	9.0	104	0.00
13 C	1,4-Dichlorobenzene	40.000	36.486	8.8	105	0.00
14	1,2-Dichlorobenzene	40.000	34.770	13.1	103	0.00
15	Benzyl Alcohol	40.000	36.265	9.3	103	0.00
16	2,2'-oxybis(1-Chloropropane)	40.000	36.371	9.1	103	0.00
17	2-Methylphenol	40.000	36.212	9.5	104	0.00
18	Hexachloroethane	40.000	36.415	9.0	103	0.00
19 P	n-Nitroso-di-n-propylamine	40.000	34.741	13.1	103	0.00
20	3+4-Methylphenols	40.000	40.307	-0.8	102	0.00
21 I	Naphthalene-d8	20.000	20.000	0.0	103	0.00
22	Acetophenone	40.000	35.138	12.2	102	0.00
23 S	Nitrobenzene-d5	80.000	73.126	8.6	103	0.00
24	Nitrobenzene	40.000	36.961	7.6	103	0.00
25	Isophorone	40.000	36.139	9.7	102	0.00
26 C	2-Nitrophenol	40.000	37.801	5.5	103	0.00
27	2,4-Dimethylphenol	40.000	36.988	7.5	103	0.00
28	bis(2-Chloroethoxy)methane	40.000	36.431	8.9	102	0.00
29 C	2,4-Dichlorophenol	40.000	37.141	7.1	104	0.00
30	1,2,4-Trichlorobenzene	40.000	36.014	10.0	102	0.00
31	Naphthalene	40.000	35.843	10.4	102	0.00
32	Benzoic acid	40.000	39.405	1.5	107	0.00
33	4-Chloroaniline	40.000	36.439	8.9	103	0.00
34 C	Hexachlorobutadiene	40.000	36.243	9.4	101	0.00
35	Caprolactam	40.000	38.938	2.7	109	0.00
36 C	4-Chloro-3-methylphenol	40.000	37.069	7.3	104	0.00
37	2-Methylnaphthalene	40.000	36.260	9.4	102	0.00
38	1-Methylnaphthalene	40.000	36.321	9.2	103	0.00
39 I	Acenaphthene-d10	20.000	20.000	0.0	105	0.00
40	1,2,4,5-Tetrachlorobenzene	40.000	36.565	8.6	103	0.00
41 P	Hexachlorocyclopentadiene	40.000	38.875	2.8	105	0.00
42 S	2,4,6-Tribromophenol	80.000	74.625	6.7	105	0.00
43 C	2,4,6-Trichlorophenol	40.000	37.640	5.9	104	0.00
44	2,4,5-Trichlorophenol	40.000	37.405	6.5	104	0.00

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF081920\
 Data File : BF121332.D
 Acq On : 19 Aug 2020 22:02
 Operator : JU/CG
 Sample : SSTDICV040
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 ICVBF081920

Quant Time: Aug 20 05:30:05 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF081920.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 20 04:56:10 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	79.946	0.1	101	0.00
46	1,1'-Biphenyl	40.000	36.175	9.6	102	0.00
47	2-Chloronaphthalene	40.000	36.263	9.3	102	0.00
48	2-Nitroaniline	40.000	37.502	6.2	105	0.00
49	Acenaphthylene	40.000	36.439	8.9	103	0.00
50	Dimethylphthalate	40.000	36.384	9.0	103	0.00
51	2,6-Dinitrotoluene	40.000	37.212	7.0	104	0.00
52 C	Acenaphthene	40.000	36.176	9.6	104	0.00
53	3-Nitroaniline	40.000	37.068	7.3	104	0.00
54 P	2,4-Dinitrophenol	40.000	42.842	-7.1	109	0.00
55	Dibenzofuran	40.000	40.510	-1.3	103	0.00
56 P	4-Nitrophenol	40.000	39.821	0.4	108	0.00
57	2,4-Dinitrotoluene	40.000	37.834	5.4	104	0.00
58	Fluorene	40.000	40.783	-2.0	102	0.00
59	2,3,4,6-Tetrachlorophenol	40.000	38.277	4.3	106	0.00
60	Diethylphthalate	40.000	36.252	9.4	103	0.00
61	4-Chlorophenyl-phenylether	40.000	39.674	0.8	101	0.00
62	4-Nitroaniline	40.000	38.263	4.3	108	0.00
63	Azobenzene	40.000	36.567	8.6	103	0.00
64 I	Phenanthrene-d10	20.000	20.000	0.0	106	0.00
65	4,6-Dinitro-2-methylphenol	40.000	40.999	-2.5	106	0.00
66 c	n-Nitrosodiphenylamine	40.000	36.602	8.5	105	0.00
67	4-Bromophenyl-phenylether	40.000	36.798	8.0	103	0.00
68	Hexachlorobenzene	40.000	36.243	9.4	104	0.00
69	Atrazine	40.000	37.028	7.4	106	0.00
70 C	Pentachlorophenol	40.000	41.354	-3.4	106	0.00
71	Phenanthrene	40.000	35.032	12.4	106	0.00
72	Anthracene	40.000	35.961	10.1	105	0.00
73	Carbazole	40.000	36.267	9.3	106	0.00
74	Di-n-butylphthalate	40.000	36.996	7.5	107	0.00
75 C	Fluoranthene	40.000	35.358	11.6	107	0.00
76 I	Chrysene-d12	20.000	20.000	0.0	114	0.00
77	Benzidine	40.000	36.553	8.6	106	0.00
78	Pyrene	40.000	35.821	10.4	108	0.00
79 S	Terphenyl-d14	80.000	63.637	20.5	108	0.00
80	Butylbenzylphthalate	40.000	37.277	6.8	112	0.00
81	Benzo(a)anthracene	40.000	33.650	15.9	111	0.00
82	3,3'-Dichlorobenzidine	40.000	37.417	6.5	114	0.00
83	Chrysene	40.000	35.919	10.2	112	0.00
84	Bis(2-ethylhexyl)phthalate	40.000	35.001	12.5	109	0.00
85 c	Di-n-octyl phthalate	40.000	38.077	4.8	117	0.00
86 I	Perylene-d12	20.000	20.000	0.0	127	0.00
87	Indeno(1,2,3-cd)pyrene	40.000	36.589	8.5	127	0.00

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF081920\
Data File : BF121332.D
Acq On : 19 Aug 2020 22:02
Operator : JU/CG
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
ICVBF081920

Quant Time: Aug 20 05:30:05 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF081920.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Aug 20 04:56:10 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	35.660	10.9	120	0.00
89	Benzo(k)fluoranthene	40.000	36.229	9.4	123	0.00
90 C	Benzo(a)pyrene	40.000	36.684	8.3	125	0.00
91	Dibenzo(a,h)anthracene	40.000	36.038	9.9	125	0.00
92	Benzo(q,h,i)perylene	40.000	36.683	8.3	128	0.00

(#) = Out of Range

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