

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF070519\
 Data File : BF115391.D
 Acq On : 5 Jul 2019 16:15
 Operator : HP/JU
 Sample : SSTDICV040
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 ICVBF070519

Quant Time: Jul 05 16:55:01 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF070519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 05 16:50:10 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	111	0.00
2	1,4-Dioxane	40.000	37.752	5.6	109	0.00
3	Pyridine	40.000	38.245	4.4	112	0.00
4	n-Nitrosodimethylamine	40.000	37.902	5.2	108	0.00
5 S	2-Fluorophenol	80.000	75.994	5.0	108	0.00
6	Aniline	40.000	38.682	3.3	108	0.00
7 S	Phenol-d6	80.000	77.628	3.0	110	0.00
8	2-Chlorophenol	40.000	38.885	2.8	108	0.00
9	Benzaldehyde	40.000	40.944	-2.4	114	0.00
10 C	Phenol	40.000	36.295	9.3	103	0.00
11	bis(2-Chloroethyl)ether	40.000	38.215	4.5	108	0.00
12	1,3-Dichlorobenzene	40.000	38.422	3.9	108	0.00
13 C	1,4-Dichlorobenzene	40.000	38.340	4.1	109	0.00
14	1,2-Dichlorobenzene	40.000	38.951	2.6	110	0.00
15	Benzyl Alcohol	40.000	39.388	1.5	110	0.00
16	2,2'-oxybis(1-Chloropropane)	40.000	38.178	4.6	108	0.00
17	2-Methylphenol	40.000	38.449	3.9	108	0.00
18	Hexachloroethane	40.000	39.724	0.7	113	0.00
19 P	n-Nitroso-di-n-propylamine	40.000	37.760	5.6	107	0.00
20	3+4-Methylphenols	40.000	39.305	1.7	108	0.00
21 I	Naphthalene-d8	20.000	20.000	0.0	108	0.00
22	Acetophenone	40.000	38.825	2.9	107	0.00
23 S	Nitrobenzene-d5	80.000	82.448	-3.1	114	0.00
24	Nitrobenzene	40.000	40.588	-1.5	112	0.00
25	Isophorone	40.000	38.170	4.6	107	0.00
26 C	2-Nitrophenol	40.000	46.560	-16.4	127	0.00
27	2,4-Dimethylphenol	40.000	39.194	2.0	110	0.00
28	bis(2-Chloroethoxy)methane	40.000	38.469	3.8	108	0.00
29 C	2,4-Dichlorophenol	40.000	39.402	1.5	109	0.00
30	1,2,4-Trichlorobenzene	40.000	39.160	2.1	109	0.00
31	Naphthalene	40.000	38.521	3.7	106	0.00
32	Benzoic acid	40.000	36.559	8.6	97	0.00
33	4-Chloroaniline	40.000	38.503	3.7	107	0.00
34 C	Hexachlorobutadiene	40.000	39.465	1.3	109	0.00
35	Caprolactam	40.000	38.092	4.8	104	0.00
36 C	4-Chloro-3-methylphenol	40.000	38.909	2.7	106	0.00
37	2-Methylnaphthalene	40.000	38.282	4.3	105	0.00
38	1-Methylnaphthalene	40.000	38.572	3.6	106	0.00
39 I	Acenaphthene-d10	20.000	20.000	0.0	105	0.00
40	1,2,4,5-Tetrachlorobenzene	40.000	39.851	0.4	107	0.00
41 P	Hexachlorocyclopentadiene	40.000	41.492	-3.7	112	0.00
42 S	2,4,6-Tribromophenol	80.000	79.340	0.8	104	0.00
43 C	2,4,6-Trichlorophenol	40.000	40.583	-1.5	108	0.00
44	2,4,5-Trichlorophenol	40.000	39.686	0.8	105	0.00

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF070519\
 Data File : BF115391.D
 Acq On : 5 Jul 2019 16:15
 Operator : HP/JU
 Sample : SSTDICV040
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 ICVBF070519

Quant Time: Jul 05 16:55:01 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF070519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 05 16:50:10 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	80.236	-0.3	102	0.00
46	1,1'-Biphenyl	40.000	39.535	1.2	106	0.00
47	2-Chloronaphthalene	40.000	39.499	1.3	106	0.00
48	2-Nitroaniline	40.000	42.960	-7.4	111	0.00
49	Acenaphthylene	40.000	38.792	3.0	104	0.00
50	Dimethylphthalate	40.000	38.582	3.5	103	0.00
51	2,6-Dinitrotoluene	40.000	40.702	-1.8	106	0.00
52 C	Acenaphthene	40.000	39.767	0.6	106	0.00
53	3-Nitroaniline	40.000	40.985	-2.5	106	0.00
54 P	2,4-Dinitrophenol	40.000	45.967	-14.9	136	0.00
55	Dibenzofuran	40.000	38.765	3.1	103	0.00
56 P	4-Nitrophenol	40.000	42.392	-6.0	108	0.00
57	2,4-Dinitrotoluene	40.000	42.159	-5.4	108	0.00
58	Fluorene	40.000	36.200	9.5	102	0.00
59	2,3,4,6-Tetrachlorophenol	40.000	39.512	1.2	103	0.00
60	Diethylphthalate	40.000	38.814	3.0	102	0.00
61	4-Chlorophenyl-phenylether	40.000	37.498	6.3	103	0.00
62	4-Nitroaniline	40.000	40.934	-2.3	104	0.00
63	Azobenzene	40.000	38.435	3.9	101	0.00
64 I	Phenanthrene-d10	20.000	20.000	0.0	100	0.00
65	4,6-Dinitro-2-methylphenol	40.000	45.237	-13.1	123	0.00
66 c	n-Nitrosodiphenylamine	40.000	39.247	1.9	101	0.00
67	4-Bromophenyl-phenylether	40.000	39.279	1.8	102	0.00
68	Hexachlorobenzene	40.000	39.472	1.3	102	0.00
69	Atrazine	40.000	40.852	-2.1	101	0.00
70 C	Pentachlorophenol	40.000	41.368	-3.4	103	0.00
71	Phenanthrene	40.000	38.949	2.6	100	0.00
72	Anthracene	40.000	39.570	1.1	102	0.00
73	Carbazole	40.000	39.183	2.0	100	0.00
74	Di-n-butylphthalate	40.000	39.630	0.9	101	0.00
75 C	Fluoranthene	40.000	38.604	3.5	98	0.00
76 I	Chrysene-d12	20.000	20.000	0.0	95	-0.02
77	Benzidine	40.000	36.826	7.9	89	0.00
78	Pyrene	40.000	40.018	-0.0	99	0.00
79 S	Terphenyl-d14	80.000	78.933	1.3	99	0.00
80	Butylbenzylphthalate	40.000	41.368	-3.4	100	0.00
81	Benzo(a)anthracene	40.000	38.400	4.0	94	-0.02
82	3,3'-Dichlorobenzidine	40.000	41.120	-2.8	94	-0.02
83	Chrysene	40.000	38.696	3.3	94	-0.02
84	Bis(2-ethylhexyl)phthalate	40.000	39.817	0.5	99	-0.02
85 c	Di-n-octyl phthalate	40.000	39.375	1.6	94	-0.05
86	Indeno(1,2,3-cd)pyrene	40.000	38.839	2.9	92	-0.05
87 I	Perylene-d12	20.000	20.000	0.0	85	-0.04

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF070519\
Data File : BF115391.D
Acq On : 5 Jul 2019 16:15
Operator : HP/JU
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
ICVBF070519

Quant Time: Jul 05 16:55:01 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF070519.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jul 05 16:50:10 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	38.226	4.4	82	-0.04
89	Benzo(k)fluoranthene	40.000	40.559	-1.4	92	-0.05
90 C	Benzo(a)pyrene	40.000	39.221	1.9	86	-0.05
91	Dibenzo(a,h)anthracene	40.000	39.926	0.2	92	-0.05
92	Benzo(q,h,i)perylene	40.000	40.695	-1.7	98	-0.05

(#) = Out of Range

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