

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF013123\
 Data File : BF132242.D
 Acq On : 31 Jan 2023 21:49
 Operator : CG\JU
 Sample : SSTDICV040
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 ICVBF013123

Quant Time: Feb 01 05:32:16 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF013123.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 01 05:24:37 2023
 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	100	0.00
2	1,4-Dioxane	40.000	39.750	0.6	97	0.00
3	Pyridine	40.000	40.190	-0.5	98	0.00
4	n-Nitrosodimethylamine	40.000	39.320	1.7	97	0.00
5 S	2-Fluorophenol	80.000	79.590	0.5	100	0.00
6	Aniline	40.000	39.351	1.6	98	0.00
7 S	Phenol-d6	80.000	79.479	0.7	99	0.00
8	2-Chlorophenol	40.000	40.310	-0.8	99	0.00
9	Benzaldehyde	40.000	40.441	-1.1	99	0.00
10 C	Phenol	40.000	39.592	1.0	98	0.00
11	bis(2-Chloroethyl)ether	40.000	40.015	-0.0	99	0.00
12	1,3-Dichlorobenzene	40.000	40.115	-0.3	99	0.00
13 C	1,4-Dichlorobenzene	40.000	40.655	-1.6	100	0.00
14	1,2-Dichlorobenzene	40.000	40.138	-0.3	99	0.00
15	Benzyl Alcohol	40.000	39.944	0.1	97	0.00
16	2,2'-oxybis(1-Chloropropane	40.000	39.108	2.2	96	0.00
17	2-Methylphenol	40.000	40.359	-0.9	100	0.00
18	Hexachloroethane	40.000	41.103	-2.8	101	0.00
19 P	n-Nitroso-di-n-propylamine	40.000	37.866	5.3	95	0.00
20	3+4-Methylphenols	40.000	40.090	-0.2	97	0.00
21 I	Naphthalene-d8	20.000	20.000	0.0	99	0.00
22	Acetophenone	40.000	38.970	2.6	97	0.00
23 S	Nitrobenzene-d5	80.000	79.213	1.0	98	0.00
24	Nitrobenzene	40.000	39.689	0.8	98	0.00
25	Isophorone	40.000	39.191	2.0	99	0.00
26 C	2-Nitrophenol	40.000	42.532	-6.3	101	0.00
27	2,4-Dimethylphenol	40.000	39.886	0.3	99	0.00
28	bis(2-Chloroethoxy)methane	40.000	39.646	0.9	98	0.00
29 C	2,4-Dichlorophenol	40.000	40.257	-0.6	99	0.00
30	1,2,4-Trichlorobenzene	40.000	40.220	-0.5	99	0.00
31	Naphthalene	40.000	39.757	0.6	99	0.00
32	Benzoic acid	40.000	41.648	-4.1	100	0.00
33	4-Chloroaniline	40.000	39.711	0.7	98	0.00
34 C	Hexachlorobutadiene	40.000	40.742	-1.9	101	0.00
35	Caprolactam	40.000	39.862	0.3	97	0.00
36 C	4-Chloro-3-methylphenol	40.000	40.380	-1.0	99	0.00
37	2-Methylnaphthalene	40.000	39.849	0.4	98	0.00
38	1-Methylnaphthalene	40.000	39.986	0.0	99	0.00
39 I	Acenaphthene-d10	20.000	20.000	0.0	99	0.00
40	1,2,4,5-Tetrachlorobenzene	40.000	39.994	0.0	99	0.00
41 P	Hexachlorocyclopentadiene	40.000	42.218	-5.5	101	0.00
42 S	2,4,6-Tribromophenol	80.000	79.927	0.1	100	0.00
43 C	2,4,6-Trichlorophenol	40.000	39.928	0.2	97	0.00
44	2,4,5-Trichlorophenol	40.000	40.408	-1.0	99	0.00
45 S	2-Fluorobiphenyl	80.000	73.224	8.5	98	0.00
46	1,1'-Biphenyl	40.000	39.484	1.3	97	0.00
47	2-Chloronaphthalene	40.000	39.228	1.9	97	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF013123\
 Data File : BF132242.D
 Acq On : 31 Jan 2023 21:49
 Operator : CG\JU
 Sample : SSTDICV040
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 ICVBF013123

Quant Time: Feb 01 05:32:16 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF013123.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 01 05:24:37 2023
 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)
48	2-Nitroaniline	40.000	39.902	0.2	95	0.00
49	Acenaphthylene	40.000	39.493	1.3	97	0.00
50	Dimethylphthalate	40.000	39.399	1.5	97	0.00
51	2,6-Dinitrotoluene	40.000	41.158	-2.9	99	0.00
52 C	Acenaphthene	40.000	39.930	0.2	99	0.00
53	3-Nitroaniline	40.000	40.904	-2.3	99	0.00
54 P	2,4-Dinitrophenol	40.000	39.829	0.4	100	0.00
55	Dibenzofuran	40.000	39.579	1.1	98	0.00
56 P	4-Nitrophenol	40.000	42.042	-5.1	99	0.00
57	2,4-Dinitrotoluene	40.000	41.568	-3.9	99	0.00
58	Fluorene	40.000	39.156	2.1	99	0.00
59	2,3,4,6-Tetrachlorophenol	40.000	40.809	-2.0	98	0.00
60	Diethylphthalate	40.000	40.256	-0.6	98	0.00
61	4-Chlorophenyl-phenylether	40.000	39.350	1.6	98	0.00
62	4-Nitroaniline	40.000	40.094	-0.2	98	0.00
63	Azobenzene	40.000	38.966	2.6	95	0.00
64 I	Phenanthrene-d10	20.000	20.000	0.0	99	0.00
65	4,6-Dinitro-2-methylphenol	40.000	43.395	-8.5	100	0.00
66 c	n-Nitrosodiphenylamine	40.000	38.561	3.6	97	0.00
67	4-Bromophenyl-phenylether	40.000	39.197	2.0	98	0.00
68	Hexachlorobenzene	40.000	38.879	2.8	98	0.00
69	Atrazine	40.000	41.088	-2.7	98	0.00
70 C	Pentachlorophenol	40.000	41.408	-3.5	100	0.00
71	Phenanthrene	40.000	38.693	3.3	98	0.00
72	Anthracene	40.000	38.808	3.0	97	0.00
73	Carbazole	40.000	38.859	2.9	98	0.00
74	Di-n-butylphthalate	40.000	39.844	0.4	99	0.00
75 C	Fluoranthene	40.000	39.691	0.8	100	0.00
76 I	Chrysene-d12	20.000	20.000	0.0	100	0.00
77	Benzydine	40.000	33.982	15.0	89	0.00
78	Pyrene	40.000	39.181	2.0	99	0.00
79 S	Terphenyl-d14	80.000	76.022	5.0	99	0.00
80	Butylbenzylphthalate	40.000	40.242	-0.6	99	0.00
81	Benzo(a)anthracene	40.000	39.822	0.4	99	0.00
82	3,3'-Dichlorobenzidine	40.000	40.347	-0.9	99	0.00
83	Chrysene	40.000	39.156	2.1	99	0.00
84	Bis(2-ethylhexyl)phthalate	40.000	40.413	-1.0	101	0.00
85 c	Di-n-octyl phthalate	40.000	40.944	-2.4	101	0.00
86 I	Perylene-d12	20.000	20.000	0.0	99	0.00
87	Indeno(1,2,3-cd)pyrene	40.000	40.426	-1.1	104	0.00
88	Benzo(b)fluoranthene	40.000	39.441	1.4	94	0.00
89	Benzo(k)fluoranthene	40.000	42.124	-5.3	104	0.00
90 C	Benzo(a)pyrene	40.000	40.252	-0.6	98	0.00
91	Dibenzo(a,h)anthracene	40.000	40.858	-2.1	103	0.00
92	Benzo(g,h,i)perylene	40.000	39.954	0.1	104	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF013123\
Data File : BF132242.D
Acq On : 31 Jan 2023 21:49
Operator : CG\JU
Sample : SSTDICV040
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
ICVBF013123

Quant Time: Feb 01 05:32:16 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF013123.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Feb 01 05:24:37 2023
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)
----------	--------	-------	------	-------	----------

(#) = Out of Range SPCC's out = 0 CCC's out = 0