

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF013023\
 Data File : BF132212.D
 Acq On : 30 Jan 2023 15:41
 Operator : CG\JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 LabSampleId :
 SSTDCCC040

Quant Time: Jan 31 00:30:07 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF012723.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Jan 28 02:25:45 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	98	0.00
2	1,4-Dioxane	40.000	40.148	-0.4	93	0.00
3	Pyridine	40.000	39.256	1.9	92	0.00
4	n-Nitrosodimethylamine	40.000	37.643	5.9	87	0.00
5 S	2-Fluorophenol	80.000	79.294	0.9	94	0.00
6	Aniline	40.000	39.729	0.7	93	0.00
7 S	Phenol-d6	80.000	78.043	2.4	92	0.00
8	2-Chlorophenol	40.000	40.660	-1.6	94	0.00
9	Benzaldehyde	40.000	37.900	5.3	94	0.00
10 C	Phenol	40.000	39.291	1.8	92	0.00
11	bis(2-Chloroethyl)ether	40.000	39.142	2.1	94	0.00
12	1,3-Dichlorobenzene	40.000	39.841	0.4	95	0.00
13 C	1,4-Dichlorobenzene	40.000	39.939	0.2	94	0.00
14	1,2-Dichlorobenzene	40.000	39.983	0.0	95	0.00
15	Benzyl Alcohol	40.000	39.572	1.1	90	0.00
16	2,2'-oxybis(1-Chloropropane	40.000	37.755	5.6	89	0.00
17	2-Methylphenol	40.000	39.670	0.8	93	0.00
18	Hexachloroethane	40.000	40.658	-1.6	94	0.00
19 P	n-Nitroso-di-n-propylamine	40.000	37.561	6.1	90	0.00
20	3+4-Methylphenols	40.000	40.224	-0.6	94	0.00
21 I	Naphthalene-d8	20.000	20.000	0.0	97	0.00
22	Acetophenone	40.000	38.947	2.6	95	0.00
23 S	Nitrobenzene-d5	80.000	78.087	2.4	91	0.00
24	Nitrobenzene	40.000	39.130	2.2	91	0.00
25	Isophorone	40.000	38.444	3.9	91	0.00
26 C	2-Nitrophenol	40.000	37.650	5.9	92	0.00
27	2,4-Dimethylphenol	40.000	40.425	-1.1	95	0.00
28	bis(2-Chloroethoxy)methane	40.000	39.247	1.9	94	0.00
29 C	2,4-Dichlorophenol	40.000	41.098	-2.7	95	0.00
30	1,2,4-Trichlorobenzene	40.000	40.567	-1.4	96	0.00
31	Naphthalene	40.000	39.403	1.5	95	0.00
32	Benzoic acid	40.000	35.415	11.5	92	-0.01
33	4-Chloroaniline	40.000	40.486	-1.2	96	0.00
34 C	Hexachlorobutadiene	40.000	40.573	-1.4	96	0.00
35	Caprolactam	40.000	42.015	-5.0	94	0.00
36 C	4-Chloro-3-methylphenol	40.000	40.529	-1.3	94	0.00
37	2-Methylnaphthalene	40.000	39.647	0.9	95	0.00
38	1-Methylnaphthalene	40.000	39.666	0.8	96	0.00
39 I	Acenaphthene-d10	20.000	20.000	0.0	98	0.00
40	1,2,4,5-Tetrachlorobenzene	40.000	40.224	-0.6	97	0.00
41 P	Hexachlorocyclopentadiene	40.000	41.342	-3.4	94	0.00
42 S	2,4,6-Tribromophenol	80.000	80.407	-0.5	94	0.00
43 C	2,4,6-Trichlorophenol	40.000	41.495	-3.7	95	0.00
44	2,4,5-Trichlorophenol	40.000	41.177	-2.9	95	0.00
45 S	2-Fluorobiphenyl	80.000	72.362	9.5	95	0.00
46	1,1'-Biphenyl	40.000	39.434	1.4	96	0.00
47	2-Chloronaphthalene	40.000	39.304	1.7	96	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF013023\
 Data File : BF132212.D
 Acq On : 30 Jan 2023 15:41
 Operator : CG\JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 LabSampleId :
 SSTDCCC040

Quant Time: Jan 31 00:30:07 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF012723.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Jan 28 02:25:45 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	40.274	-0.7	89	0.00
49	Acenaphthylene	40.000	39.809	0.5	96	0.00
50	Dimethylphthalate	40.000	40.015	-0.0	97	0.00
51	2,6-Dinitrotoluene	40.000	41.945	-4.9	95	0.00
52 C	Acenaphthene	40.000	39.194	2.0	95	0.00
53	3-Nitroaniline	40.000	41.590	-4.0	96	0.00
54 P	2,4-Dinitrophenol	40.000	35.930	10.2	90	0.00
55	Dibenzofuran	40.000	39.255	1.9	96	0.00
56 P	4-Nitrophenol	40.000	40.311	-0.8	93	0.00
57	2,4-Dinitrotoluene	40.000	42.431	-6.1	95	0.00
58	Fluorene	40.000	39.096	2.3	96	0.00
59	2,3,4,6-Tetrachlorophenol	40.000	41.904	-4.8	95	0.00
60	Diethylphthalate	40.000	39.926	0.2	94	0.00
61	4-Chlorophenyl-phenylether	40.000	39.786	0.5	97	0.00
62	4-Nitroaniline	40.000	40.990	-2.5	95	0.00
63	Azobenzene	40.000	38.421	3.9	93	0.00
64 I	Phenanthrene-d10	20.000	20.000	0.0	99	0.00
65	4,6-Dinitro-2-methylphenol	40.000	37.158	7.1	94	0.00
66 c	n-Nitrosodiphenylamine	40.000	38.999	2.5	97	0.00
67	4-Bromophenyl-phenylether	40.000	39.956	0.1	98	0.00
68	Hexachlorobenzene	40.000	40.016	-0.0	100	0.00
69	Atrazine	40.000	42.023	-5.1	97	0.00
70 C	Pentachlorophenol	40.000	40.307	-0.8	95	0.00
71	Phenanthrene	40.000	39.448	1.4	99	0.00
72	Anthracene	40.000	39.073	2.3	97	0.00
73	Carbazole	40.000	39.883	0.3	97	0.00
74	Di-n-butylphthalate	40.000	40.234	-0.6	94	0.00
75 C	Fluoranthene	40.000	40.487	-1.2	98	0.00
76 I	Chrysene-d12	20.000	20.000	0.0	102	0.00
77	Benzidine	40.000	45.887	-14.7	105	0.00
78	Pyrene	40.000	40.427	-1.1	98	0.00
79 S	Terphenyl-d14	80.000	79.207	1.0	98	0.00
80	Butylbenzylphthalate	40.000	38.770	3.1	94	0.00
81	Benzo(a)anthracene	40.000	40.120	-0.3	100	0.00
82	3,3'-Dichlorobenzidine	40.000	43.721	-9.3	100	0.00
83	Chrysene	40.000	40.414	-1.0	100	0.00
84	Bis(2-ethylhexyl)phthalate	40.000	38.711	3.2	93	0.00
85 c	Di-n-octyl phthalate	40.000	34.810	13.0	87	0.00
86 I	Perylene-d12	20.000	20.000	0.0	100	0.00
87	Indeno(1,2,3-cd)pyrene	40.000	40.900	-2.2	94	-0.01
88	Benzo(b)fluoranthene	40.000	40.843	-2.1	96	0.00
89	Benzo(k)fluoranthene	40.000	40.866	-2.2	103	0.00
90 C	Benzo(a)pyrene	40.000	41.621	-4.1	97	0.00
91	Dibenzo(a,h)anthracene	40.000	41.213	-3.0	94	0.00
92	Benzo(g,h,i)perylene	40.000	40.608	-1.5	95	-0.01

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF013023\
Data File : BF132212.D
Acq On : 30 Jan 2023 15:41
Operator : CG\JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Jan 31 00:30:07 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF012723.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat Jan 28 02:25:45 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