

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102924\
 Data File : BF140100.D
 Acq On : 29 Oct 2024 09:37
 Operator : RC/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 LabSampleId :
 SSTDCCC040

Quant Time: Oct 29 09:59:29 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF101824.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 18 15:07:50 2024
 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	76	0.00
2	1,4-Dioxane	40.000	36.065	9.8	69	-0.02
3	Pyridine	40.000	36.523	8.7	68	-0.01
4	n-Nitrosodimethylamine	40.000	35.174	12.1	66	-0.01
5 S	2-Fluorophenol	80.000	77.105	3.6	73	0.00
6	Aniline	40.000	35.714	10.7	66	0.00
7 S	Phenol-d6	80.000	75.076	6.2	72	0.00
8	2-Chlorophenol	40.000	38.460	3.8	73	0.00
9	Benzaldehyde	40.000	38.226	4.4	72	0.00
10 C	Phenol	40.000	37.804	5.5	72	0.00
11	bis(2-Chloroethyl)ether	40.000	38.288	4.3	74	0.00
12	1,3-Dichlorobenzene	40.000	39.591	1.0	75	0.00
13 C	1,4-Dichlorobenzene	40.000	39.522	1.2	76	0.00
14	1,2-Dichlorobenzene	40.000	39.670	0.8	76	0.00
15	Benzyl Alcohol	40.000	38.130	4.7	72	0.00
16	2,2'-oxybis(1-Chloropropane	40.000	36.318	9.2	69	0.00
17	2-Methylphenol	40.000	38.055	4.9	72	0.00
18	Hexachloroethane	40.000	40.062	-0.2	75	0.00
19 P	n-Nitroso-di-n-propylamine	40.000	36.422	8.9	71	0.00
20	3+4-Methylphenols	40.000	37.773	5.6	72	0.00
21 I	Naphthalene-d8	20.000	20.000	0.0	75	0.00
22	Acetophenone	40.000	38.706	3.2	74	0.00
23 S	Nitrobenzene-d5	80.000	84.711	-5.9	77	0.00
24	Nitrobenzene	40.000	40.486	-1.2	75	0.00
25	Isophorone	40.000	38.579	3.6	73	0.00
26 C	2-Nitrophenol	40.000	46.897	-17.2	84	0.00
27	2,4-Dimethylphenol	40.000	37.294	6.8	70	0.00
28	bis(2-Chloroethoxy)methane	40.000	38.395	4.0	73	0.00
29 C	2,4-Dichlorophenol	40.000	39.500	1.3	74	0.00
30	1,2,4-Trichlorobenzene	40.000	39.921	0.2	75	0.00
31	Naphthalene	40.000	39.161	2.1	74	0.00
32	Benzoic acid	40.000	41.140	-2.9	76	0.00
33	4-Chloroaniline	40.000	37.651	5.9	71	0.00
34 C	Hexachlorobutadiene	40.000	41.197	-3.0	78	0.00
35	Caprolactam	40.000	39.935	0.2	74	0.00
36 C	4-Chloro-3-methylphenol	40.000	39.486	1.3	74	0.00
37	2-Methylnaphthalene	40.000	39.026	2.4	75	0.00
38	1-Methylnaphthalene	40.000	38.730	3.2	74	0.00
39 I	Acenaphthene-d10	20.000	20.000	0.0	75	0.00
40	1,2,4,5-Tetrachlorobenzene	40.000	40.717	-1.8	77	0.00
41 P	Hexachlorocyclopentadiene	40.000	41.093	-2.7	73	0.00
42 S	2,4,6-Tribromophenol	80.000	86.905	-8.6	82	0.00
43 C	2,4,6-Trichlorophenol	40.000	39.415	1.5	73	0.00
44	2,4,5-Trichlorophenol	40.000	41.730	-4.3	77	0.00
45 S	2-Fluorobiphenyl	80.000	77.983	2.5	77	0.00
46	1,1'-Biphenyl	40.000	39.200	2.0	75	0.00
47	2-Chloronaphthalene	40.000	39.527	1.2	75	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102924\
 Data File : BF140100.D
 Acq On : 29 Oct 2024 09:37
 Operator : RC/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 LabSampleId :
 SSTDCCC040

Quant Time: Oct 29 09:59:29 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF101824.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 18 15:07:50 2024
 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	44.929	-12.3	80	0.00
49	Acenaphthylene	40.000	39.600	1.0	75	0.00
50	Dimethylphthalate	40.000	39.396	1.5	75	0.00
51	2,6-Dinitrotoluene	40.000	43.705	-9.3	79	0.00
52 C	Acenaphthene	40.000	40.465	-1.2	77	0.00
53	3-Nitroaniline	40.000	42.683	-6.7	76	0.00
54 P	2,4-Dinitrophenol	40.000	57.194	-43.0#	115	0.00
55	Dibenzofuran	40.000	39.241	1.9	75	0.00
56 P	4-Nitrophenol	40.000	44.022	-10.1	77	0.00
57	2,4-Dinitrotoluene	40.000	46.765	-16.9	82	0.00
58	Fluorene	40.000	39.507	1.2	77	0.00
59	2,3,4,6-Tetrachlorophenol	40.000	40.906	-2.3	77	0.00
60	Diethylphthalate	40.000	39.485	1.3	75	0.00
61	4-Chlorophenyl-phenylether	40.000	40.286	-0.7	77	0.00
62	4-Nitroaniline	40.000	44.623	-11.6	80	0.00
63	Azobenzene	40.000	38.698	3.3	72	0.00
64 I	Phenanthrene-d10	20.000	20.000	0.0	77	0.00
65	4,6-Dinitro-2-methylphenol	40.000	58.416	-46.0#	106	0.00
66 c	n-Nitrosodiphenylamine	40.000	37.816	5.5	74	0.00
67	4-Bromophenyl-phenylether	40.000	40.226	-0.6	79	0.00
68	Hexachlorobenzene	40.000	40.035	-0.1	78	0.00
69	Atrazine	40.000	39.389	1.5	71	0.00
70 C	Pentachlorophenol	40.000	43.806	-9.5	79	0.00
71	Phenanthrene	40.000	38.654	3.4	76	0.00
72	Anthracene	40.000	39.128	2.2	76	0.00
73	Carbazole	40.000	39.205	2.0	77	0.00
74	Di-n-butylphthalate	40.000	40.805	-2.0	77	0.00
75 C	Fluoranthene	40.000	39.514	1.2	77	0.00
76 I	Chrysene-d12	20.000	20.000	0.0	76	0.00
77	Benzydine	40.000	47.078	-17.7	81	0.00
78	Pyrene	40.000	41.982	-5.0	78	0.00
79 S	Terphenyl-d14	80.000	83.670	-4.6	79	0.00
80	Butylbenzylphthalate	40.000	44.907	-12.3	81	0.00
81	Benzo(a)anthracene	40.000	41.194	-3.0	77	0.00
82	3,3'-Dichlorobenzidine	40.000	42.900	-7.2	80	0.00
83	Chrysene	40.000	37.838	5.4	74	0.00
84	Bis(2-ethylhexyl)phthalate	40.000	45.315	-13.3	82	0.00
85 c	Di-n-octyl phthalate	40.000	39.177	2.1	71	-0.01
86 I	Perylene-d12	20.000	20.000	0.0	72	0.00
87	Indeno(1,2,3-cd)pyrene	40.000	41.128	-2.8	71	0.00
88	Benzo(b)fluoranthene	40.000	38.915	2.7	72	0.00
89	Benzo(k)fluoranthene	40.000	39.346	1.6	70	0.00
90 C	Benzo(a)pyrene	40.000	39.861	0.3	71	0.00
91	Dibenzo(a,h)anthracene	40.000	40.482	-1.2	70	0.00
92	Benzo(g,h,i)perylene	40.000	42.618	-6.5	73	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102924\
Data File : BF140100.D
Acq On : 29 Oct 2024 09:37
Operator : RC/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Oct 29 09:59:29 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF101824.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Oct 18 15:07:50 2024
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