

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112524\
 Data File : BF140604.D
 Acq On : 25 Nov 2024 15:49
 Operator : RC/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 LabSampleId :
 SSTDCCC040

Quant Time: Nov 25 16:13:33 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF112124.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 21 15:23:48 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	99	0.00
2	1,4-Dioxane	40.000	41.024	-2.6	102	-0.01
3	Pyridine	40.000	46.035	-15.1	104	-0.01
4	n-Nitrosodimethylamine	40.000	37.334	6.7	91	-0.02
5 S	2-Fluorophenol	80.000	78.663	1.7	97	0.00
6	Aniline	40.000	45.917	-14.8	99	0.00
7 S	Phenol-d6	80.000	78.504	1.9	94	0.00
8	2-Chlorophenol	40.000	38.614	3.5	94	0.00
9	Benzaldehyde	40.000	41.353	-3.4	114	0.00
10 C	Phenol	40.000	39.924	0.2	94	0.00
11	bis(2-Chloroethyl)ether	40.000	39.593	1.0	95	0.00
12	1,3-Dichlorobenzene	40.000	39.031	2.4	98	0.00
13 C	1,4-Dichlorobenzene	40.000	39.049	2.4	97	0.00
14	1,2-Dichlorobenzene	40.000	39.050	2.4	96	0.00
15	Benzyl Alcohol	40.000	37.981	5.0	88	0.00
16	2,2'-oxybis(1-Chloropropane	40.000	39.966	0.1	97	0.00
17	2-Methylphenol	40.000	40.390	-1.0	97	0.00
18	Hexachloroethane	40.000	38.893	2.8	96	0.00
19 P	n-Nitroso-di-n-propylamine	40.000	38.269	4.3	92	0.00
20	3+4-Methylphenols	40.000	38.823	2.9	93	0.00
21 I	Naphthalene-d8	20.000	20.000	0.0	95	0.00
22	Acetophenone	40.000	38.065	4.8	92	0.00
23 S	Nitrobenzene-d5	80.000	77.261	3.4	92	0.00
24	Nitrobenzene	40.000	38.812	3.0	93	0.00
25	Isophorone	40.000	39.246	1.9	92	0.00
26 C	2-Nitrophenol	40.000	40.245	-0.6	95	0.00
27	2,4-Dimethylphenol	40.000	41.812	-4.5	95	0.00
28	bis(2-Chloroethoxy)methane	40.000	39.568	1.1	94	0.00
29 C	2,4-Dichlorophenol	40.000	39.979	0.1	96	0.00
30	1,2,4-Trichlorobenzene	40.000	39.526	1.2	96	0.00
31	Naphthalene	40.000	39.364	1.6	95	0.00
32	Benzoic acid	40.000	37.962	5.1	91	0.00
33	4-Chloroaniline	40.000	43.479	-8.7	98	0.00
34 C	Hexachlorobutadiene	40.000	39.213	2.0	95	0.00
35	Caprolactam	40.000	40.238	-0.6	93	-0.01
36 C	4-Chloro-3-methylphenol	40.000	39.361	1.6	91	0.00
37	2-Methylnaphthalene	40.000	39.171	2.1	94	0.00
38	1-Methylnaphthalene	40.000	39.666	0.8	95	0.00
39 I	Acenaphthene-d10	20.000	20.000	0.0	93	0.00
40	1,2,4,5-Tetrachlorobenzene	40.000	39.161	2.1	95	0.00
41 P	Hexachlorocyclopentadiene	40.000	36.552	8.6	87	0.00
42 S	2,4,6-Tribromophenol	80.000	76.932	3.8	89	0.00
43 C	2,4,6-Trichlorophenol	40.000	38.964	2.6	91	0.00
44	2,4,5-Trichlorophenol	40.000	39.616	1.0	91	0.00
45 S	2-Fluorobiphenyl	80.000	76.385	4.5	93	0.00
46	1,1'-Biphenyl	40.000	39.310	1.7	95	0.00
47	2-Chloronaphthalene	40.000	39.263	1.8	94	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112524\
 Data File : BF140604.D
 Acq On : 25 Nov 2024 15:49
 Operator : RC/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 LabSampleId :
 SSTDCCC040

Quant Time: Nov 25 16:13:33 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF112124.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 21 15:23:48 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	38.667	3.3	89	0.00
49	Acenaphthylene	40.000	39.111	2.2	94	0.00
50	Dimethylphthalate	40.000	38.266	4.3	91	0.00
51	2,6-Dinitrotoluene	40.000	39.685	0.8	92	0.00
52 C	Acenaphthene	40.000	38.661	3.3	92	0.00
53	3-Nitroaniline	40.000	38.687	3.3	88	0.00
54 P	2,4-Dinitrophenol	40.000	36.476	8.8	82	0.00
55	Dibenzofuran	40.000	38.157	4.6	91	0.00
56 P	4-Nitrophenol	40.000	34.868	12.8	78	0.00
57	2,4-Dinitrotoluene	40.000	39.205	2.0	90	0.00
58	Fluorene	40.000	38.375	4.1	92	0.00
59	2,3,4,6-Tetrachlorophenol	40.000	40.946	-2.4	94	0.00
60	Diethylphthalate	40.000	38.281	4.3	91	0.00
61	4-Chlorophenyl-phenylether	40.000	38.445	3.9	92	0.00
62	4-Nitroaniline	40.000	38.184	4.5	87	0.00
63	Azobenzene	40.000	38.931	2.7	93	0.00
64 I	Phenanthrene-d10	20.000	20.000	0.0	91	0.00
65	4,6-Dinitro-2-methylphenol	40.000	42.537	-6.3	90	0.00
66 c	n-Nitrosodiphenylamine	40.000	39.439	1.4	91	0.00
67	4-Bromophenyl-phenylether	40.000	39.900	0.3	93	0.00
68	Hexachlorobenzene	40.000	39.651	0.9	92	0.00
69	Atrazine	40.000	39.313	1.7	106	0.00
70 C	Pentachlorophenol	40.000	40.265	-0.7	82	0.00
71	Phenanthrene	40.000	39.222	1.9	90	0.00
72	Anthracene	40.000	39.611	1.0	91	0.00
73	Carbazole	40.000	39.233	1.9	90	0.00
74	Di-n-butylphthalate	40.000	39.448	1.4	91	0.00
75 C	Fluoranthene	40.000	38.587	3.5	90	0.00
76 I	Chrysene-d12	20.000	20.000	0.0	91	0.00
77	Benzydine	40.000	54.347	-35.9#	125	0.00
78	Pyrene	40.000	40.033	-0.1	91	0.00
79 S	Terphenyl-d14	80.000	78.343	2.1	89	-0.01
80	Butylbenzylphthalate	40.000	40.617	-1.5	91	0.00
81	Benzo(a)anthracene	40.000	38.832	2.9	91	0.00
82	3,3'-Dichlorobenzidine	40.000	40.346	-0.9	88	0.00
83	Chrysene	40.000	39.966	0.1	92	0.00
84	Bis(2-ethylhexyl)phthalate	40.000	40.221	-0.6	92	0.00
85 c	Di-n-octyl phthalate	40.000	40.002	-0.0	92	0.02
86 I	Perylene-d12	20.000	20.000	0.0	91	0.01
87	Indeno(1,2,3-cd)pyrene	40.000	38.092	4.8	87	0.01
88	Benzo(b)fluoranthene	40.000	36.487	8.8	88	0.01
89	Benzo(k)fluoranthene	40.000	41.760	-4.4	95	0.01
90 C	Benzo(a)pyrene	40.000	40.250	-0.6	93	0.02
91	Dibenzo(a,h)anthracene	40.000	38.118	4.7	87	0.01
92	Benzo(g,h,i)perylene	40.000	37.523	6.2	86	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112524\
Data File : BF140604.D
Acq On : 25 Nov 2024 15:49
Operator : RC/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Nov 25 16:13:33 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF112124.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Nov 21 15:23:48 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