Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF123124\
Data File : BF141056.D
Acq On : 31 Dec 2024 14:18
Operator : RC/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Dec 31 14:40:26 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF122624.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Dec 27 00:31:16 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	94	0.00
2		1,4-Dioxane	40.000	37.601	6.0	89	0.00
3		Pyridine	40.000	39.135	2.2	94	0.00
4		n-Nitrosodimethylamine	40.000	37.516	6.2	88	0.00
5	S	2-Fluorophenol	80.000	76.722	4.1	93	0.00
6		Aniline	40.000	39.293	1.8	88	-0.02
7	S	Phenol-d6	80.000	74.706	6.6	90	0.00
8		2-Chlorophenol	40.000	38.916	2.7	93	0.00
9		Benzaldehyde	40.000	46.666	-16.7	104	0.00
10	C	Phenol	40.000	39.080	2.3	94	0.00
11		bis(2-Chloroethyl)ether	40.000	36.649	8.4	91	0.00
12		1,3-Dichlorobenzene	40.000	38.753	3.1	93	0.00
13	C	1,4-Dichlorobenzene	40.000	39.011	2.5	95	0.00
14		1,2-Dichlorobenzene	40.000	38.575	3.6	93	0.00
15		Benzyl Alcohol	40.000	38.046	4.9	90	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	36.192	9.5	87	0.00
17		2-Methylphenol	40.000	37.699	5.8	90	0.00
18		Hexachloroethane	40.000	39.952	0.1	96	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	35.625	10.9	87	0.00
20		3+4-Methylphenols	40.000	37.276	6.8	89	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	92	0.00
22		Acetophenone	40.000	36.896	7.8	89	0.00
23	S	Nitrobenzene-d5	80.000	94.938	-18.7	104	0.00
24		Nitrobenzene	40.000	44.276	-10.7	97	0.00
25		Isophorone	40.000	37.283	6.8	89	0.00
26	C	2-Nitrophenol	40.000	50.173	-25.4#	131	0.00
27		2,4-Dimethylphenol	40.000	38.304	4.2	91	0.00
28		bis(2-Chloroethoxy)methane	40.000	37.959	5.1	91	0.00
29	C	2,4-Dichlorophenol	40.000	40.267	-0.7	94	0.00
30		1,2,4-Trichlorobenzene	40.000	39.661	0.8	93	0.00
31		Naphthalene	40.000	38.303	4.2	91	0.00
32		Benzoic acid	40.000	48.603	-21.5	129	0.00
33		4-Chloroaniline	40.000	37.176	7.1	89	0.00
34	C	Hexachlorobutadiene	40.000	40.216	-0.5	96	0.00
35		Caprolactam	40.000	37.534	6.2	88	0.00
36	C	4-Chloro-3-methylphenol	40.000	38.425	3.9	90	0.00
37		2-Methylnaphthalene	40.000	38.195	4.5	92	0.00
38		1-Methylnaphthalene	40.000	37.935	5.2	91	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	90	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	41.182	-3.0	95	0.00
41	P	Hexachlorocyclopentadiene	40.000	47.466	-18.7	104	0.00
42	S	2,4,6-Tribromophenol	80.000	87.297	-9.1	97	0.00
43	C	2,4,6-Trichlorophenol	40.000	43.220	-8.0	98	0.00
44		2,4,5-Trichlorophenol	40.000	41.935	-4.8	93	0.00
45	S	2-Fluorobiphenyl	80.000	76.894	3.9	92	0.00
46		1,1'-Biphenyl	40.000	39.415	1.5	92	0.00
47		2-Chloronaphthalene	40.000	39.805	0.5	93	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF123124\
Data File : BF141056.D
Acq On : 31 Dec 2024 14:18
Operator : RC/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Dec 31 14:40:26 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF122624.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Dec 27 00:31:16 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	46.405	-16.0	112	0.00
49		Acenaphthylene	40.000	39.119	2.2	91	0.00
50		Dimethylphthalate	40.000	38.410	4.0	88	0.00
51		2,6-Dinitrotoluene	40.000	45.153	-12.9	106	0.00
52	C	Acenaphthene	40.000	40.559	-1.4	94	0.00
53		3-Nitroaniline	40.000	43.308	-8.3	100	0.00
54	P	2,4-Dinitrophenol	40.000	62.472	-56.2#	185	0.00
55		Dibenzofuran	40.000	38.614	3.5	90	0.00
56	P	4-Nitrophenol	40.000	44.114	-10.3	97	0.00
57		2,4-Dinitrotoluene	40.000	47.848	-19.6	114	0.00
58		Fluorene	40.000	38.136	4.7	91	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	41.813	-4.5	93	0.00
60		Diethylphthalate	40.000	37.678	5.8	87	0.00
61		4-Chlorophenyl-phenylether	40.000	38.896	2.8	92	0.00
62		4-Nitroaniline	40.000	41.190	-3.0	96	0.00
63		Azobenzene	40.000	36.490	8.8	85	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	83	0.00
65		4,6-Dinitro-2-methylphenol	40.000	58.512	-46.3#	152	0.00
66	c	n-Nitrosodiphenylamine	40.000	40.455	-1.1	87	0.00
67		4-Bromophenyl-phenylether	40.000	42.756	-6.9	90	0.00
68		Hexachlorobenzene	40.000	41.736	-4.3	89	0.00
69		Atrazine	40.000	30.966	22.6	72	0.00
70	C	Pentachlorophenol	40.000	45.157	-12.9	90	0.00
71		Phenanthrene	40.000	39.047	2.4	85	0.00
72		Anthracene	40.000	39.126	2.2	85	0.00
73		Carbazole	40.000	36.565	8.6	80	0.00
74		Di-n-butylphthalate	40.000	36.952	7.6	80	0.00
75	C	Fluoranthene	40.000	34.610	13.5	75	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	68	0.00
77		Benzidine	40.000	48.939	-22.3	85	0.00
78		Pyrene	40.000	41.154	-2.9	72	0.00
79	S	Terphenyl-d14	80.000	81.499	-1.9	73	0.00
80		Butylbenzylphthalate	40.000	38.729	3.2	64	0.00
81		Benzo(a)anthracene	40.000	36.987	7.5	64	0.00
82		3,3'-Dichlorobenzidine	40.000	44.728	-11.8	79	0.00
83		Chrysene	40.000	40.489	-1.2	72	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	36.806	8.0	63	0.00
85	c	Di-n-octyl phthalate	40.000	41.777	-4.4	71	0.00
86	I	Perylene-d12	20.000	20.000	0.0	99	0.00
87		Indeno(1,2,3-cd)pyrene	40.000	46.672	-16.7	111	0.01
88		Benzo(b)fluoranthene	40.000	32.178	19.6	78	0.00
89		Benzo(k)fluoranthene	40.000	39.512	1.2	100	0.00
90	C	Benzo(a)pyrene	40.000	38.024	4.9	95	0.00
91		Dibenzo(a,h)anthracene	40.000	46.181	-15.5	109	0.01
92		Benzo(g,h,i)perylene	40.000	47.475	-18.7	111	0.02

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF123124\
Data File : BF141056.D
Acq On : 31 Dec 2024 14:18
Operator : RC/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Dec 31 14:40:26 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF122624.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Dec 27 00:31:16 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 = 1