Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF122624\
Data File : BF140978.D
Acq On : 26 Dec 2024 19:52
Operator : RC/JU
Sample : SSTDICV040
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
ICVBF122624

Quant Time: Dec 27 00:40:05 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	106	0.00
2		1,4-Dioxane	40.000	39.078	2.3	104	0.00
3		Pyridine	40.000	39.221	1.9	106	0.00
4		n-Nitrosodimethylamine	40.000	39.126	2.2	103	0.01
5	S	2-Fluorophenol	80.000	77.230	3.5	105	0.00
6		Aniline	40.000	40.214	-0.5	100	-0.02
7	S	Phenol-d6	80.000	77.619	3.0	105	0.00
8		2-Chlorophenol	40.000	45.402	-13.5	121	0.00
9		Benzaldehyde	40.000	42.275	-5.7	107	0.00
10	C	Phenol	40.000	46.698	-16.7	125	0.00
11		bis(2-Chloroethyl)ether	40.000	39.366	1.6	109	0.00
12		1,3-Dichlorobenzene	40.000	38.861	2.8	105	0.00
13	C	1,4-Dichlorobenzene	40.000	40.612	-1.5	110	0.00
14		1,2-Dichlorobenzene	40.000	38.772	3.1	105	0.00
15		Benzyl Alcohol	40.000	39.751	0.6	105	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	38.326	4.2	103	0.00
17		2-Methylphenol	40.000	39.617	1.0	106	0.00
18		Hexachloroethane	40.000	40.027	-0.1	107	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	40.828	-2.1	112	0.00
20		3+4-Methylphenols	40.000	39.323	1.7	105	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	105	0.00
22		Acetophenone	40.000	37.920	5.2	104	0.00
23	S	Nitrobenzene-d5	80.000	91.412	-14.3	115	0.00
24		Nitrobenzene	40.000	44.060	-10.2	110	0.00
25		Isophorone	40.000	38.534	3.7	105	0.00
26	C	2-Nitrophenol	40.000	43.185	-8.0	126	0.00
27		2,4-Dimethylphenol	40.000	39.926	0.2	108	0.00
28		bis(2-Chloroethoxy)methane	40.000	38.323	4.2	104	0.00
29	C	2,4-Dichlorophenol	40.000	40.280	-0.7	107	0.00
30		1,2,4-Trichlorobenzene	40.000	41.204	-3.0	110	0.00
31		Naphthalene	40.000	38.493	3.8	104	0.00
32		Benzoic acid	40.000	43.158	-7.9	127	0.00
33		4-Chloroaniline	40.000	39.203	2.0	107	0.00
34	C	Hexachlorobutadiene	40.000	39.095	2.3	106	0.00
35		Caprolactam	40.000	40.139	-0.3	107	0.00
36	C	4-Chloro-3-methylphenol	40.000	42.714	-6.8	113	0.00
37		2-Methylnaphthalene	40.000	38.340	4.1	105	0.00
38		1-Methylnaphthalene	40.000	37.974	5.1	104	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	106	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	38.527	3.7	105	0.00
41	P	Hexachlorocyclopentadiene	40.000	40.560	-1.4	105	0.00
42	S	2,4,6-Tribromophenol	80.000	82.713	-3.4	108	0.00
43	C	2,4,6-Trichlorophenol	40.000	40.180	-0.4	107	0.00
44		2,4,5-Trichlorophenol	40.000	40.663	-1.7	106	0.00
45	S	2-Fluorobiphenyl	80.000	74.821	6.5	106	0.00
46		1,1'-Biphenyl	40.000	37.916	5.2	104	0.00
47		2-Chloronaphthalene	40.000	38.442	3.9	106	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF122624\
Data File : BF140978.D
Acq On : 26 Dec 2024 19:52
Operator : RC/JU
Sample : SSTDICV040
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
ICVBF122624

Quant Time: Dec 27 00:40:05 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	43.198	-8.0	122	0.00
49		Acenaphthylene	40.000	38.272	4.3	104	0.00
50		Dimethylphthalate	40.000	38.422	3.9	104	0.00
51		2,6-Dinitrotoluene	40.000	42.386	-6.0	116	0.00
52	C	Acenaphthene	40.000	40.168	-0.4	110	0.00
53		3-Nitroaniline	40.000	42.522	-6.3	115	0.00
54	P	2,4-Dinitrophenol	40.000	45.538	-13.8	138	0.00
55		Dibenzofuran	40.000	38.318	4.2	105	0.00
56	P	4-Nitrophenol	40.000	48.593	-21.5	126	0.00
57		2,4-Dinitrotoluene	40.000	44.506	-11.3	124	0.00
58		Fluorene	40.000	37.668	5.8	105	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	41.681	-4.2	109	0.00
60		Diethylphthalate	40.000	38.629	3.4	105	0.00
61		4-Chlorophenyl-phenylether	40.000	37.928	5.2	105	0.00
62		4-Nitroaniline	40.000	41.571	-3.9	114	0.00
63		Azobenzene	40.000	38.471	3.8	105	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	105	0.00
65		4,6-Dinitro-2-methylphenol	40.000	43.296	-8.2	132	0.00
66	c	n-Nitrosodiphenylamine	40.000	38.777	3.1	106	0.00
67		4-Bromophenyl-phenylether	40.000	39.247	1.9	104	0.00
68		Hexachlorobenzene	40.000	38.890	2.8	105	0.00
69		Atrazine	40.000	35.858	10.4	106	0.00
70	C	Pentachlorophenol	40.000	45.869	-14.7	115	0.00
71		Phenanthrene	40.000	38.373	4.1	105	0.00
72		Anthracene	40.000	38.325	4.2	104	0.00
73		Carbazole	40.000	38.096	4.8	104	0.00
74		Di-n-butylphthalate	40.000	38.673	3.3	105	0.00
75	C	Fluoranthene	40.000	37.724	5.7	103	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	94	0.00
77		Benzidine	40.000	40.101	-0.3	96	0.00
78		Pyrene	40.000	43.575	-8.9	106	0.00
79	S	Terphenyl-d14	80.000	83.347	-4.2	104	0.00
80		Butylbenzylphthalate	40.000	43.715	-9.3	100	0.00
81		Benzo(a)anthracene	40.000	37.941	5.1	91	0.00
82		3,3'-Dichlorobenzidine	40.000	35.805	10.5	88	0.00
83		Chrysene	40.000	39.695	0.8	97	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	40.713	-1.8	97	0.00
85	c	Di-n-octyl phthalate	40.000	39.285	1.8	93	0.00
86	I	Perylene-d12	20.000	20.000	0.0	102	0.00
87		Indeno(1,2,3-cd)pyrene	40.000	44.255	-10.6	109	0.00
88		Benzo(b)fluoranthene	40.000	36.115	9.7	91	0.00
89		Benzo(k)fluoranthene	40.000	40.555	-1.4	106	0.00
90	C	Benzo(a)pyrene	40.000	38.510	3.7	100	0.00
91		Dibenzo(a,h)anthracene	40.000	44.756	-11.9	109	0.00
92		Benzo(g,h,i)perylene	40.000	45.122	-12.8	108	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF122624\
Data File : BF140978.D
Acq On : 26 Dec 2024 19:52
Operator : RC/JU
Sample : SSTDICV040
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
ICVBF122624

Quant Time: Dec 27 00:40:05 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 = 0