Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121621\
Data File : BM033488.D
Acq On : 16 Dec 2021 17:23
Operator : CG/JU
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
ICVBM121621

Quant Time: Dec 16 18:08:22 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM121621.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Dec 16 17:10:15 2021
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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	103	0.00
2		1,4-Dioxane	0.505	0.465	7.9	106	0.00
3		Pyridine	1.235	1.195	3.2	95	0.00
4		n-Nitrosodimethylamine	0.878	0.841	4.2	94	0.00
5	S	2-Fluorophenol	1.159	1.153	0.5	97	0.00
6		Aniline	1.776	1.824	-2.7	99	0.00
7	S	Phenol-d6	1.555	1.596	-2.6	99	0.00
8		2-Chlorophenol	1.256	1.248	0.6	97	0.00
9		Benzaldehyde	1.123	0.992	11.7	103	0.00
10	C	Phenol	1.571	1.580	-0.6	97	0.00
11		bis(2-Chloroethyl)ether	1.238	1.255	-1.4	101	0.00
12		1,3-Dichlorobenzene	1.526	1.507	1.2	97	0.00
13	C	1,4-Dichlorobenzene	1.567	1.546	1.3	98	0.00
14		1,2-Dichlorobenzene	1.537	1.515	1.4	96	0.00
15		Benzyl Alcohol	1.402	1.485	-5.9	98	0.00
16		2,2'-oxybis(1-Chloropropane	2.202	2.178	1.1	99	0.00
17		2-Methylphenol	1.117	1.133	-1.4	98	0.00
18		Hexachloroethane	0.653	0.641	1.8	97	0.00
19	P	n-Nitroso-di-n-propylamine	1.226	1.256	-2.4	101	0.00
20		3+4-Methylphenols	1.490	1.574	-5.6	99	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	107	0.00
22		Acetophenone	0.541	0.535	1.1	101	0.00
23	S	Nitrobenzene-d5	0.466	0.462	0.9	98	0.00
24		Nitrobenzene	0.458	0.449	2.0	98	0.00
25		Isophorone	0.754	0.731	3.1	100	0.00
26	C	2-Nitrophenol	0.160	0.168	-5.0	102	0.00
27		2,4-Dimethylphenol	0.263	0.260	1.1	98	0.00
28		bis(2-Chloroethoxy)methane	0.414	0.410	1.0	98	0.00
29	C	2,4-Dichlorophenol	0.300	0.304	-1.3	98	0.00
30		1,2,4-Trichlorobenzene	0.380	0.360	5.3	98	0.00
31		Naphthalene	1.083	1.040	4.0	97	0.00
32		Benzoic acid	0.164	0.182	-11.0	102	0.00
33		4-Chloroaniline	0.351	0.372	-6.0	98	0.00
34	C	Hexachlorobutadiene	0.265	0.252	4.9	97	0.00
35		Caprolactam	0.059	0.062	-5.1	101	0.00
36	C	4-Chloro-3-methylphenol	0.384	0.383	0.3	98	0.00
37		2-Methylnaphthalene	0.749	0.751	-0.3	101	0.00
38		1-Methylnaphthalene	0.748	0.724	3.2	98	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	107	0.00
40		1,2,4,5-Tetrachlorobenzene	0.576	0.553	4.0	100	0.00
41	P	Hexachlorocyclopentadiene	0.249	0.254	-2.0	98	0.00
42	S	2,4,6-Tribromophenol	0.225	0.218	3.1	97	0.00
43	C	2,4,6-Trichlorophenol	0.385	0.384	0.3	100	0.00
44		2,4,5-Trichlorophenol	0.411	0.403	1.9	98	0.00
45	S	2-Fluorobiphenyl	1.446	1.365	5.6	99	0.00
46		1,1'-Biphenyl	1.481	1.409	4.9	99	0.00
47		2-Chloronaphthalene	1.145	1.091	4.7	98	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121621\
Data File : BM033488.D
Acq On : 16 Dec 2021 17:23
Operator : CG/JU
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
ICVBM121621

Quant Time: Dec 16 18:08:22 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM121621.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Dec 16 17:10:15 2021
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	AvgRF	CCRF	%Dev	Area%	Dev(min)
48		2-Nitroaniline	0.390	0.392	-0.5	99	0.00
49		Acenaphthylene	1.814	1.754	3.3	99	0.00
50		Dimethylphthalate	1.515	1.431	5.5	98	0.00
51		2,6-Dinitrotoluene	0.286	0.276	3.5	94	0.00
52	C	Acenaphthene	1.102	1.055	4.3	99	0.00
53		3-Nitroaniline	0.269	0.268	0.4	94	0.00
54	P	2,4-Dinitrophenol	0.125	0.135	-8.0	112	0.00
55		Dibenzofuran	1.867	1.761	5.7	98	0.00
56	P	4-Nitrophenol	0.189	0.186	1.6	91	0.00
57		2,4-Dinitrotoluene	0.392	0.385	1.8	97	0.00
58		Fluorene	1.498	1.408	6.0	97	0.00
59		2,3,4,6-Tetrachlorophenol	0.374	0.365	2.4	98	0.00
60		Diethylphthalate	1.629	1.552	4.7	97	0.00
61		4-Chlorophenyl-phenylether	0.778	0.738	5.1	96	0.00
62		4-Nitroaniline	0.244	0.254	-4.1	95	0.00
63		Azobenzene	1.828	1.752	4.2	97	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	106	0.00
65		4,6-Dinitro-2-methylphenol	0.102	0.109	-6.9	105	0.00
66	c	n-Nitrosodiphenylamine	0.594	0.583	1.9	98	0.00
67		4-Bromophenyl-phenylether	0.218	0.211	3.2	98	0.00
68		Hexachlorobenzene	0.234	0.224	4.3	97	0.00
69		Atrazine	0.126	0.134	-6.3	99	0.00
70	C	Pentachlorophenol	0.132	0.134	-1.5	97	0.00
71		Phenanthrene	1.113	1.061	4.7	96	0.00
72		Anthracene	1.088	1.042	4.2	95	0.00
73		Carbazole	0.961	0.935	2.7	95	0.00
74		Di-n-butylphthalate	1.336	1.281	4.1	95	0.00
75	C	Fluoranthene	1.280	1.205	5.9	94	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	100	0.00
77		Benzidine	0.094	0.097	-3.2	101	0.00
78		Pyrene	1.578	1.532	2.9	93	0.00
79	S	Terphenyl-d14	1.270	1.240	2.4	94	0.00
80		Butylbenzylphthalate	0.706	0.673	4.7	91	0.00
81		Benzo(a)anthracene	1.340	1.289	3.8	94	0.00
82		3,3'-Dichlorobenzidine	0.526	0.530	-0.8	96	0.00
83		Chrysene	1.291	1.247	3.4	93	0.00
84		Bis(2-ethylhexyl)phthalate	1.035	0.977	5.6	92	0.00
85	c	Di-n-octyl phthalate	1.683	1.594	5.3	94	0.00
86	I	Perylene-d12	1.000	1.000	0.0	99	0.00
87		Indeno(1,2,3-cd)pyrene	0.922	1.001	-8.6	103	0.00
88		Benzo(b)fluoranthene	1.282	1.261	1.6	96	0.00
89		Benzo(k)fluoranthene	1.297	1.311	-1.1	96	0.00
90	C	Benzo(a)pyrene	0.960	0.991	-3.2	98	0.00
91		Dibenzo(a,h)anthracene	0.692	0.808	-16.8	108	0.00
92		Benzo(g,h,i)perylene	0.889	0.920	-3.5	101	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121621\
Data File : BM033488.D
Acq On : 16 Dec 2021 17:23
Operator : CG/JU
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
ICVBM121621

Quant Time: Dec 16 18:08:22 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM121621.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Dec 16 17:10:15 2021
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	AvgRF	CCRF	%Dev Area%	Dev(min)

(#) = Out of Range SPCC's out = 0 CCC's out = 0