Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040121\
Data File : BM029308.D
Acq On : 01 Apr 2021 16:24
Operator : CG/JU
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : ICVBM040121

Quant Time: Apr 01 17:17:31 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM040121.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Apr 01 16:52:48 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	Amount	Calc.	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	20.000	20.000	0.0	104	0.00
2		1,4-Dioxane	40.000	38.146	4.6	99	0.00
3		Pyridine	40.000	40.416	-1.0	98	0.00
4		n-Nitrosodimethylamine	40.000	41.700	-4.3	101	0.00
5	S	2-Fluorophenol	80.000	78.725	1.6	102	0.00
6		Aniline	40.000	39.968	0.1	101	0.00
7	S	Phenol-d6	80.000	79.812	0.2	102	0.00
8		2-Chlorophenol	40.000	39.971	0.1	103	0.00
9		Benzaldehyde	40.000	38.814	3.0	104	0.00
10	C	Phenol	40.000	39.871	0.3	101	0.00
11		bis(2-Chloroethyl)ether	40.000	40.044	-0.1	101	0.00
12		1,3-Dichlorobenzene	40.000	39.191	2.0	100	0.00
13	C	1,4-Dichlorobenzene	40.000	38.984	2.5	101	0.00
14		1,2-Dichlorobenzene	40.000	39.434	1.4	101	0.00
15		Benzyl Alcohol	40.000	41.027	-2.6	102	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	40.382	-1.0	102	0.00
17		2-Methylphenol	40.000	40.360	-0.9	101	0.00
18		Hexachloroethane	40.000	38.964	2.6	101	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	41.080	-2.7	101	0.00
20		3+4-Methylphenols	40.000	40.454	-1.1	102	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	100	0.00
22		Acetophenone	40.000	41.058	-2.6	101	0.00
23	S	Nitrobenzene-d5	80.000	81.771	-2.2	101	0.00
24		Nitrobenzene	40.000	40.891	-2.2	102	0.00
25		Isophorone	40.000	41.595	-4.0	101	0.00
26	C	2-Nitrophenol	40.000	41.581	-4.0	102	0.00
27		2,4-Dimethylphenol	40.000	41.187	-3.0	101	0.00
28		bis(2-Chloroethoxy)methane	40.000	40.809	-2.0	100	0.00
29	C	2,4-Dichlorophenol	40.000	40.950	-2.4	101	0.00
30		1,2,4-Trichlorobenzene	40.000	40.243	-0.6	102	0.00
31		Naphthalene	40.000	40.336	-0.8	101	0.00
32		Benzoic acid	40.000	42.448	-6.1	104	0.00
33		4-Chloroaniline	40.000	41.416	-3.5	100	0.00
34	C	Hexachlorobutadiene	40.000	40.444	-1.1	103	0.00
35		Caprolactam	40.000	43.678	-9.2	102	0.00
36	C	4-Chloro-3-methylphenol	40.000	41.495	-3.7	100	0.00
37		2-Methylnaphthalene	40.000	41.212	-3.0	101	0.00
38		1-Methylnaphthalene	40.000	40.893	-2.2	101	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	100	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	40.284	-0.7	102	0.00
41	P	Hexachlorocyclopentadiene	40.000	40.399	-1.0	100	0.00
42	S	2,4,6-Tribromophenol	80.000	85.905	-7.4	104	0.00
43	C	2,4,6-Trichlorophenol	40.000	41.351	-3.4	101	0.00
44		2,4,5-Trichlorophenol	40.000	41.513	-3.8	102	0.00
45	S	2-Fluorobiphenyl	80.000	80.904	-1.1	102	0.00
46		1,1'-Biphenyl	40.000	40.418	-1.0	101	0.00
47		2-Chloronaphthalene	40.000	40.346	-0.9	101	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040121\
Data File : BM029308.D
Acq On : 01 Apr 2021 16:24
Operator : CG/JU
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : ICVBM040121

Quant Time: Apr 01 17:17:31 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM040121.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Apr 01 16:52:48 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	Amount	Calc.	%Dev	Area%	Dev(min)
48		2-Nitroaniline	40.000	42.486	-6.2	102	0.00
49		Acenaphthylene	40.000	41.305	-3.3	102	0.00
50		Dimethylphthalate	40.000	40.887	-2.2	101	0.00
51		2,6-Dinitrotoluene	40.000	41.822	-4.6	101	0.00
52	C	Acenaphthene	40.000	40.742	-1.9	101	0.00
53		3-Nitroaniline	40.000	42.961	-7.4	102	0.00
54	P	2,4-Dinitrophenol	40.000	44.599	-11.5	102	0.00
55		Dibenzofuran	40.000	40.947	-2.4	102	0.00
56	P	4-Nitrophenol	40.000	44.398	-11.0	101	0.00
57		2,4-Dinitrotoluene	40.000	43.204	-8.0	102	0.00
58		Fluorene	40.000	41.007	-2.5	101	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	42.214	-5.5	102	0.00
60		Diethylphthalate	40.000	41.285	-3.2	101	0.00
61		4-Chlorophenyl-phenylether	40.000	41.295	-3.2	102	0.00
62		4-Nitroaniline	40.000	43.456	-8.6	102	0.00
63		Azobenzene	40.000	41.667	-4.2	101	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	101	0.00
65		4,6-Dinitro-2-methylphenol	40.000	41.356	-3.4	101	0.00
66	c	n-Nitrosodiphenylamine	40.000	39.965	0.1	100	0.00
67		4-Bromophenyl-phenylether	40.000	40.463	-1.2	101	0.00
68		Hexachlorobenzene	40.000	39.994	0.0	103	0.00
69		Atrazine	40.000	41.243	-3.1	102	0.00
70	C	Pentachlorophenol	40.000	42.765	-6.9	112	0.00
71		Phenanthrene	40.000	40.256	-0.6	102	0.00
72		Anthracene	40.000	40.952	-2.4	103	0.00
73		Carbazole	40.000	41.048	-2.6	103	0.00
74		Di-n-butylphthalate	40.000	41.132	-2.8	100	0.00
75	C	Fluoranthene	40.000	40.797	-2.0	102	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	102	0.00
77		Benzidine	40.000	43.629	-9.1	104	0.00
78		Pyrene	40.000	39.733	0.7	102	0.00
79	S	Terphenyl-d14	80.000	80.553	-0.7	103	0.00
80		Butylbenzylphthalate	40.000	40.202	-0.5	101	0.00
81		Benzo(a)anthracene	40.000	40.136	-0.3	102	0.00
82		3,3'-Dichlorobenzidine	40.000	40.213	-0.5	101	0.00
83		Chrysene	40.000	40.035	-0.1	102	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	40.587	-1.5	102	0.00
85	c	Di-n-octyl phthalate	40.000	40.135	-0.3	100	0.00
86	I	Perylene-d12	20.000	20.000	0.0	101	0.00
87		Indeno(1,2,3-cd)pyrene	40.000	41.146	-2.9	103	0.00
88		Benzo(b)fluoranthene	40.000	41.443	-3.6	106	0.00
89		Benzo(k)fluoranthene	40.000	40.216	-0.5	98	0.00
90	C	Benzo(a)pyrene	40.000	40.740	-1.9	102	0.00
91		Dibenzo(a,h)anthracene	40.000	41.054	-2.6	102	0.00
92		Benzo(g,h,i)perylene	40.000	40.338	-0.8	102	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040121\
Data File : BM029308.D
Acq On : 01 Apr 2021 16:24
Operator : CG/JU
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
ICVBM040121

Quant Time: Apr 01 17:17:31 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM040121.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Apr 01 16:52:48 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	Amount	Calc.	%Dev Area%	Dev(min)

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