Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM060525\
Data File : BM050202.D
Acq On : 05 Jun 2025 14:36
Operator : RC/JU
Sample : SSTDICV040
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
ICVBM060525

Quant Time: Jun 05 15:10:28 2025
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM060525.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jun 05 14:55:18 2025
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	103	0.00
2		1,4-Dioxane	40.000	34.075	14.8	96	0.00
3		Pyridine	40.000	37.216	7.0	100	0.00
4		n-Nitrosodimethylamine	40.000	38.276	4.3	102	0.00
5	S	2-Fluorophenol	80.000	71.321	10.8	96	0.00
6		Aniline	40.000	38.184	4.5	100	0.00
7	S	Phenol-d6	80.000	72.656	9.2	95	0.00
8		2-Chlorophenol	40.000	38.620	3.5	102	0.00
9		Benzaldehyde	40.000	41.383	-3.5	122	0.00
10	C	Phenol	40.000	38.313	4.2	100	0.00
11		bis(2-Chloroethyl)ether	40.000	38.237	4.4	101	0.00
12		1,3-Dichlorobenzene	40.000	37.856	5.4	102	0.00
13	C	1,4-Dichlorobenzene	40.000	37.381	6.5	102	0.00
14		1,2-Dichlorobenzene	40.000	37.789	5.5	102	0.00
15		Benzyl Alcohol	40.000	39.129	2.2	100	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	37.404	6.5	99	0.00
17		2-Methylphenol	40.000	39.079	2.3	100	0.00
18		Hexachloroethane	40.000	38.024	4.9	102	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	39.802	0.5	99	0.00
20		3+4-Methylphenols	40.000	39.372	1.6	99	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	102	0.00
22		Acetophenone	40.000	37.622	5.9	101	0.00
23	S	Nitrobenzene-d5	80.000	65.977	17.5	86	0.00
24		Nitrobenzene	40.000	38.131	4.7	100	0.00
25		Isophorone	40.000	38.498	3.8	99	0.00
26	C	2-Nitrophenol	40.000	40.867	-2.2	102	0.00
27		2,4-Dimethylphenol	40.000	37.886	5.3	99	0.00
28		bis(2-Chloroethoxy)methane	40.000	38.061	4.8	99	0.00
29	C	2,4-Dichlorophenol	40.000	38.988	2.5	100	0.00
30		1,2,4-Trichlorobenzene	40.000	37.573	6.1	100	0.00
31		Naphthalene	40.000	36.633	8.4	99	0.00
32		Benzoic acid	40.000	39.598	1.0	97	-0.04
33		4-Chloroaniline	40.000	38.018	5.0	99	0.00
34	C	Hexachlorobutadiene	40.000	37.511	6.2	100	0.00
35		Caprolactam	40.000	38.807	3.0	93	-0.02
36	C	4-Chloro-3-methylphenol	40.000	38.568	3.6	96	0.00
37		2-Methylnaphthalene	40.000	37.986	5.0	98	0.00
38		1-Methylnaphthalene	40.000	37.722	5.7	98	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	101	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	37.845	5.4	102	0.00
41	P	Hexachlorocyclopentadiene	40.000	39.386	1.5	102	0.00
42	S	2,4,6-Tribromophenol	80.000	71.554	10.6	89	0.00
43	C	2,4,6-Trichlorophenol	40.000	38.255	4.4	98	0.00
44		2,4,5-Trichlorophenol	40.000	38.297	4.3	97	0.00
45	S	2-Fluorobiphenyl	80.000	63.026	21.2	85	0.00
46		1,1'-Biphenyl	40.000	36.765	8.1	99	0.00
47		2-Chloronaphthalene	40.000	36.269	9.3	98	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM060525\
Data File : BM050202.D
Acq On : 05 Jun 2025 14:36
Operator : RC/JU
Sample : SSTDICV040
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
ICVBM060525

Quant Time: Jun 05 15:10:28 2025
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM060525.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jun 05 14:55:18 2025
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	40.253	-0.6	97	0.00
49		Acenaphthylene	40.000	36.877	7.8	96	0.00
50		Dimethylphthalate	40.000	37.379	6.6	95	0.00
51		2,6-Dinitrotoluene	40.000	39.718	0.7	94	0.00
52	C	Acenaphthene	40.000	36.444	8.9	95	0.00
53		3-Nitroaniline	40.000	39.347	1.6	93	0.00
54	P	2,4-Dinitrophenol	40.000	34.596	13.5	92	0.00
55		Dibenzofuran	40.000	36.397	9.0	95	0.00
56	P	4-Nitrophenol	40.000	38.588	3.5	92	0.00
57		2,4-Dinitrotoluene	40.000	39.946	0.1	91	0.00
58		Fluorene	40.000	36.130	9.7	95	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	37.508	6.2	93	0.00
60		Diethylphthalate	40.000	36.717	8.2	92	0.00
61		4-Chlorophenyl-phenylether	40.000	36.132	9.7	95	0.00
62		4-Nitroaniline	40.000	38.283	4.3	89	-0.01
63		Azobenzene	40.000	36.142	9.6	93	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	93	0.00
65		4,6-Dinitro-2-methylphenol	40.000	38.808	3.0	88	0.00
66	c	n-Nitrosodiphenylamine	40.000	37.727	5.7	93	0.00
67		4-Bromophenyl-phenylether	40.000	38.453	3.9	93	0.00
68		Hexachlorobenzene	40.000	37.911	5.2	92	0.00
69		Atrazine	40.000	39.186	2.0	88	0.00
70	C	Pentachlorophenol	40.000	38.925	2.7	89	0.00
71		Phenanthrene	40.000	36.170	9.6	89	0.00
72		Anthracene	40.000	36.806	8.0	90	0.00
73		Carbazole	40.000	36.195	9.5	87	0.00
74		Di-n-butylphthalate	40.000	37.377	6.6	85	0.00
75	C	Fluoranthene	40.000	35.092	12.3	83	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	81	0.00
77		Benzidine	40.000	48.798	-22.0	90	0.00
78		Pyrene	40.000	40.348	-0.9	83	0.00
79	S	Terphenyl-d14	80.000	68.934	13.8	73	0.00
80		Butylbenzylphthalate	40.000	40.599	-1.5	75	0.00
81		Benzo(a)anthracene	40.000	37.193	7.0	78	0.00
82		3,3'-Dichlorobenzidine	40.000	40.289	-0.7	76	0.00
83		Chrysene	40.000	36.579	8.6	77	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	38.675	3.3	73	0.00
85	c	Di-n-octyl phthalate	40.000	38.642	3.4	72	0.00
86	I	Perylene-d12	20.000	20.000	0.0	84	0.00
87		Indeno(1,2,3-cd)pyrene	40.000	38.807	3.0	88	-0.01
88		Benzo(b)fluoranthene	40.000	37.732	5.7	79	0.00
89		Benzo(k)fluoranthene	40.000	36.362	9.1	78	0.00
90	C	Benzo(a)pyrene	40.000	37.840	5.4	81	0.00
91		Dibenzo(a,h)anthracene	40.000	38.821	2.9	88	0.00
92		Benzo(g,h,i)perylene	40.000	38.771	3.1	90	-0.02

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM060525\
Data File : BM050202.D
Acq On : 05 Jun 2025 14:36
Operator : RC/JU
Sample : SSTDICV040
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
ICVBM060525

Quant Time: Jun 05 15:10:28 2025
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM060525.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jun 05 14:55:18 2025
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