Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120723\
Data File : BG059989.D
Acq On : 7 Dec 2023 16:42
Operator : MA/JU
Sample : SSTDICV040
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
ICVBG120723

Quant Time: Dec 08 02:37:23 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG120723.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Dec 08 02:35:41 2023
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	91	0.00
2		1,4-Dioxane	0.489	0.401	18.0	91	0.00
3		Pyridine	1.357	1.286	5.2	92	0.00
4		n-Nitrosodimethylamine	0.588	0.568	3.4	103	0.00
5	S	2-Fluorophenol	1.081	1.022	5.5	94	0.00
6		Aniline	1.621	1.721	-6.2	84	0.00
7	S	Phenol-d6	1.520	1.399	8.0	88	0.00
8		2-Chlorophenol	1.038	0.932	10.2	86	0.00
9		Benzaldehyde	0.927	0.882	4.9	90	0.00
10	C	Phenol	1.486	1.368	7.9	87	0.00
11		bis(2-Chloroethyl)ether	0.994	0.837	15.8	81	0.00
12		1,3-Dichlorobenzene	1.468	1.172	20.2	84	0.00
13	C	1,4-Dichlorobenzene	1.444	1.223	15.3	88	0.00
14		1,2-Dichlorobenzene	1.377	1.212	12.0	96	0.00
15		Benzyl Alcohol	1.357	1.305	3.8	87	0.00
16		2,2'-oxybis(1-Chloropropane	0.535	0.463	13.5	90	0.00
17		2-Methylphenol	1.004	0.993	1.1	94	0.00
18		Hexachloroethane	0.469	0.408	13.0	101	0.00
19	P	n-Nitroso-di-n-propylamine	1.025	0.943	8.0	93	0.00
20		3+4-Methylphenols	1.422	1.351	5.0	93	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	97	0.00
22		Acetophenone	0.585	0.562	3.9	96	0.00
23	S	Nitrobenzene-d5	0.453	0.457	-0.9	97	0.00
24		Nitrobenzene	0.468	0.453	3.2	92	0.00
25		Isophorone	0.815	0.781	4.2	92	0.00
26	C	2-Nitrophenol	0.182	0.169	7.1	92	0.00
27		2,4-Dimethylphenol	0.245	0.302	-23.3	94	0.00
28		bis(2-Chloroethoxy)methane	0.386	0.373	3.4	97	0.00
29	C	2,4-Dichlorophenol	0.462	0.449	2.8	94	0.00
30		1,2,4-Trichlorobenzene	0.642	0.602	6.2	97	0.00
31		Naphthalene	1.022	0.990	3.1	97	0.00
32		Benzoic acid	0.248	0.248	0.0	99	0.00
33		4-Chloroaniline	0.384	0.418	-8.9	96	0.00
34	C	Hexachlorobutadiene	0.699	0.645	7.7	98	0.00
35		Caprolactam	0.105	0.100	4.8	88	0.01
36	C	4-Chloro-3-methylphenol	0.406	0.393	3.2	89	0.00
37		2-Methylnaphthalene	0.885	0.863	2.5	88	0.00
38		1-Methylnaphthalene	0.841	0.827	1.7	102	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	97	0.00
40		1,2,4,5-Tetrachlorobenzene	1.279	1.291	-0.9	94	0.00
41	P	Hexachlorocyclopentadiene	0.598	0.706	-18.1	89	0.00
42	S	2,4,6-Tribromophenol	0.581	0.595	-2.4	102	0.00
43	C	2,4,6-Trichlorophenol	0.663	0.651	1.8	99	0.00
44		2,4,5-Trichlorophenol	0.748	0.776	-3.7	100	0.00
45	S	2-Fluorobiphenyl	1.664	1.691	-1.6	99	0.00
46		1,1'-Biphenyl	1.455	1.413	2.9	92	0.00
47		2-Chloronaphthalene	1.253	1.187	5.3	99	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120723\
Data File : BG059989.D
Acq On : 7 Dec 2023 16:42
Operator : MA/JU
Sample : SSTDICV040
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
ICVBG120723

Quant Time: Dec 08 02:37:23 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG120723.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Dec 08 02:35:41 2023
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	0.245	0.261	-6.5	105	0.00
49		Acenaphthylene	1.689	1.710	-1.2	97	0.00
50		Dimethylphthalate	1.594	1.635	-2.6	99	0.00
51		2,6-Dinitrotoluene	0.351	0.342	2.6	99	0.00
52	C	Acenaphthene	1.224	1.210	1.1	100	0.00
53		3-Nitroaniline	0.240	0.237	1.3	89	0.00
54	P	2,4-Dinitrophenol	0.292	0.300	-2.7	99	0.00
55		Dibenzofuran	1.989	2.008	-1.0	99	0.00
56	P	4-Nitrophenol	0.203	0.197	3.0	93	0.00
57		2,4-Dinitrotoluene	0.503	0.501	0.4	104	0.00
58		Fluorene	1.733	1.710	1.3	98	0.00
59		2,3,4,6-Tetrachlorophenol	0.839	0.846	-0.8	101	0.00
60		Diethylphthalate	1.433	1.447	-1.0	100	0.00
61		4-Chlorophenyl-phenylether	1.455	1.461	-0.4	99	0.00
62		4-Nitroaniline	0.266	0.270	-1.5	97	0.00
63		Azobenzene	1.252	1.236	1.3	95	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	98	0.00
65		4,6-Dinitro-2-methylphenol	0.143	0.141	1.4	97	0.00
66	c	n-Nitrosodiphenylamine	0.475	0.485	-2.1	101	0.00
67		4-Bromophenyl-phenylether	0.314	0.322	-2.5	104	0.00
68		Hexachlorobenzene	0.365	0.339	7.1	99	0.00
69		Atrazine	0.207	0.163	21.3	98	0.00
70	C	Pentachlorophenol	0.255	0.251	1.6	103	0.00
71		Phenanthrene	0.916	0.942	-2.8	99	0.00
72		Anthracene	0.925	0.971	-5.0	100	0.00
73		Carbazole	0.755	0.730	3.3	95	0.00
74		Di-n-butylphthalate	0.674	0.674	0.0	99	0.00
75	C	Fluoranthene	1.434	1.372	4.3	99	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	95	0.00
77		Benzidine	0.235	0.169	28.1#	80	0.00
78		Pyrene	1.071	1.023	4.5	98	0.00
79	S	Terphenyl-d14	0.949	0.807	15.0	99	0.00
80		Butylbenzylphthalate	0.189	0.187	1.1	94	0.00
81		Benzo(a)anthracene	1.276	1.238	3.0	97	0.00
82		3,3'-Dichlorobenzidine	0.455	0.489	-7.5	94	0.00
83		Chrysene	1.205	1.171	2.8	95	0.00
84		Bis(2-ethylhexyl)phthalate	0.278	0.287	-3.2	97	0.00
85	c	Di-n-octyl phthalate	0.480	0.485	-1.0	96	0.00
86	I	Perylene-d12	1.000	1.000	0.0	94	0.00
87		Indeno(1,2,3-cd)pyrene	1.457	1.419	2.6	96	0.00
88		Benzo(b)fluoranthene	1.164	1.102	5.3	96	0.00
89		Benzo(k)fluoranthene	1.127	1.135	-0.7	95	0.00
90	C	Benzo(a)pyrene	1.061	1.097	-3.4	96	0.00
91		Dibenzo(a,h)anthracene	1.227	1.219	0.7	95	0.00
92		Benzo(g,h,i)perylene	1.196	1.153	3.6	96	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120723\
Data File : BG059989.D
Acq On : 7 Dec 2023 16:42
Operator : MA/JU
Sample : SSTDICV040
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
ICVBG120723

Quant Time: Dec 08 02:37:23 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG120723.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Dec 08 02:35:41 2023
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