Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071222\
Data File : BG054257.D
Acq On : 12 Jul 2022 15:29
Operator : CG/JU
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
ICVBG071222

Quant Time: Jul 12 16:55:42 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG071222.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jul 12 15:49:20 2022
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	36.737	8.2	102	0.00
3		Pyridine	40.000	43.256	-8.1	102	0.00
4		n-Nitrosodimethylamine	40.000	43.485	-8.7	106	0.00
5	S	2-Fluorophenol	80.000	83.580	-4.5	102	0.00
6		Aniline	40.000	41.235	-3.1	102	0.00
7	S	Phenol-d6	80.000	81.928	-2.4	100	0.00
8		2-Chlorophenol	40.000	42.012	-5.0	103	0.00
9		Benzaldehyde	40.000	35.693	10.8	104	0.00
10	C	Phenol	40.000	40.673	-1.7	99	0.00
11		bis(2-Chloroethyl)ether	40.000	40.496	-1.2	97	0.00
12		1,3-Dichlorobenzene	40.000	40.767	-1.9	100	0.00
13	C	1,4-Dichlorobenzene	40.000	41.873	-4.7	104	0.00
14		1,2-Dichlorobenzene	40.000	40.330	-0.8	101	0.00
15		Benzyl Alcohol	40.000	40.081	-0.2	96	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	38.221	4.4	97	0.00
17		2-Methylphenol	40.000	40.330	-0.8	99	0.00
18		Hexachloroethane	40.000	41.867	-4.7	102	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	40.020	-0.1	100	0.00
20		3+4-Methylphenols	40.000	40.076	-0.2	99	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	97	0.00
22		Acetophenone	40.000	41.746	-4.4	98	0.00
23	S	Nitrobenzene-d5	80.000	84.841	-6.1	100	0.00
24		Nitrobenzene	40.000	42.155	-5.4	99	0.00
25		Isophorone	40.000	41.208	-3.0	97	0.00
26	C	2-Nitrophenol	40.000	42.797	-7.0	99	0.00
27		2,4-Dimethylphenol	40.000	42.625	-6.6	100	0.00
28		bis(2-Chloroethoxy)methane	40.000	41.105	-2.8	98	0.00
29	C	2,4-Dichlorophenol	40.000	43.252	-8.1	101	0.00
30		1,2,4-Trichlorobenzene	40.000	42.142	-5.4	99	0.00
31		Naphthalene	40.000	41.541	-3.9	98	0.00
32		Benzoic acid	40.000	37.496	6.3	96	0.00
33		4-Chloroaniline	40.000	42.361	-5.9	101	0.00
34	C	Hexachlorobutadiene	40.000	43.130	-7.8	100	0.00
35		Caprolactam	40.000	41.275	-3.2	98	0.00
36	C	4-Chloro-3-methylphenol	40.000	42.133	-5.3	100	0.00
37		2-Methylnaphthalene	40.000	41.539	-3.8	98	0.00
38		1-Methylnaphthalene	40.000	41.244	-3.1	97	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	100	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	41.527	-3.8	99	0.00
41	P	Hexachlorocyclopentadiene	40.000	40.628	-1.6	101	0.00
42	S	2,4,6-Tribromophenol	80.000	84.870	-6.1	103	0.00
43	C	2,4,6-Trichlorophenol	40.000	41.932	-4.8	100	0.00
44		2,4,5-Trichlorophenol	40.000	43.134	-7.8	103	0.00
45	S	2-Fluorobiphenyl	80.000	82.628	-3.3	99	0.00
46		1,1'-Biphenyl	40.000	40.827	-2.1	98	0.00
47		2-Chloronaphthalene	40.000	41.056	-2.6	99	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071222\
Data File : BG054257.D
Acq On : 12 Jul 2022 15:29
Operator : CG/JU
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
ICVBG071222

Quant Time: Jul 12 16:55:42 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG071222.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jul 12 15:49:20 2022
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	41.869	-4.7	101	0.00
49		Acenaphthylene	40.000	40.861	-2.2	99	0.00
50		Dimethylphthalate	40.000	40.558	-1.4	100	0.00
51		2,6-Dinitrotoluene	40.000	41.238	-3.1	98	0.00
52	C	Acenaphthene	40.000	40.603	-1.5	99	0.00
53		3-Nitroaniline	40.000	41.135	-2.8	99	0.00
54	P	2,4-Dinitrophenol	40.000	38.593	3.5	102	0.00
55		Dibenzofuran	40.000	41.110	-2.8	99	0.00
56	P	4-Nitrophenol	40.000	44.687	-11.7	107	0.00
57		2,4-Dinitrotoluene	40.000	41.293	-3.2	100	0.00
58		Fluorene	40.000	40.507	-1.3	99	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	41.562	-3.9	99	0.00
60		Diethylphthalate	40.000	40.806	-2.0	101	0.00
61		4-Chlorophenyl-phenylether	40.000	41.303	-3.3	100	0.00
62		4-Nitroaniline	40.000	42.282	-5.7	104	0.00
63		Azobenzene	40.000	40.209	-0.5	99	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	103	0.00
65		4,6-Dinitro-2-methylphenol	40.000	41.983	-5.0	101	0.00
66	c	n-Nitrosodiphenylamine	40.000	40.613	-1.5	101	0.00
67		4-Bromophenyl-phenylether	40.000	41.271	-3.2	102	0.00
68		Hexachlorobenzene	40.000	40.765	-1.9	101	0.00
69		Atrazine	40.000	42.372	-5.9	105	0.00
70	C	Pentachlorophenol	40.000	39.586	1.0	103	0.00
71		Phenanthrene	40.000	41.212	-3.0	102	0.00
72		Anthracene	40.000	41.041	-2.6	102	0.00
73		Carbazole	40.000	41.059	-2.6	102	0.00
74		Di-n-butylphthalate	40.000	40.959	-2.4	103	0.00
75	C	Fluoranthene	40.000	41.362	-3.4	103	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	106	0.00
77		Benzidine	40.000	38.329	4.2	95	0.00
78		Pyrene	40.000	41.235	-3.1	105	0.00
79	S	Terphenyl-d14	80.000	83.641	-4.6	107	0.00
80		Butylbenzylphthalate	40.000	41.846	-4.6	108	0.00
81		Benzo(a)anthracene	40.000	41.754	-4.4	108	0.00
82		3,3'-Dichlorobenzidine	40.000	41.799	-4.5	105	0.00
83		Chrysene	40.000	41.424	-3.6	107	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	41.212	-3.0	106	0.00
85	c	Di-n-octyl phthalate	40.000	41.352	-3.4	105	0.00
86	I	Perylene-d12	20.000	20.000	0.0	108	0.00
87		Indeno(1,2,3-cd)pyrene	40.000	42.058	-5.1	109	-0.01
88		Benzo(b)fluoranthene	40.000	42.195	-5.5	110	0.00
89		Benzo(k)fluoranthene	40.000	41.092	-2.7	107	0.00
90	C	Benzo(a)pyrene	40.000	41.638	-4.1	109	0.00
91		Dibenzo(a,h)anthracene	40.000	42.371	-5.9	110	0.00
92		Benzo(g,h,i)perylene	40.000	41.995	-5.0	110	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071222\
Data File : BG054257.D
Acq On : 12 Jul 2022 15:29
Operator : CG/JU
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
ICVBG071222

Quant Time: Jul 12 16:55:42 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG071222.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jul 12 15:49:20 2022
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