Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG102121\
Data File : BG050660.D
Acq On : 21 Oct 2021 9:39
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Oct 21 11:11:18 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG100621.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Oct 14 10:56:22 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	94	0.00
2		1,4-Dioxane	40.000	31.745	20.6	75	0.00
3		Pyridine	40.000	34.119	14.7	81	0.00
4		n-Nitrosodimethylamine	40.000	35.323	11.7	86	0.00
5	S	2-Fluorophenol	80.000	74.383	7.0	90	0.00
6		Aniline	40.000	38.024	4.9	94	0.00
7	S	Phenol-d6	80.000	77.815	2.7	95	0.00
8		2-Chlorophenol	40.000	39.950	0.1	98	0.00
9		Benzaldehyde	40.000	39.886	0.3	89	0.00
10	C	Phenol	40.000	38.855	2.9	96	0.00
11		bis(2-Chloroethyl)ether	40.000	37.727	5.7	93	0.00
12		1,3-Dichlorobenzene	40.000	38.241	4.4	95	0.00
13	C	1,4-Dichlorobenzene	40.000	37.849	5.4	94	0.00
14		1,2-Dichlorobenzene	40.000	38.097	4.8	95	0.00
15		Benzyl Alcohol	40.000	40.508	-1.3	100	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	36.996	7.5	92	0.00
17		2-Methylphenol	40.000	40.151	-0.4	100	0.00
18		Hexachloroethane	40.000	39.158	2.1	95	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	38.133	4.7	94	0.00
20		3+4-Methylphenols	40.000	39.940	0.2	97	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	104	0.00
22		Acetophenone	40.000	35.907	10.2	96	0.00
23	S	Nitrobenzene-d5	80.000	72.067	9.9	94	0.00
24		Nitrobenzene	40.000	35.657	10.9	94	0.00
25		Isophorone	40.000	36.400	9.0	97	0.00
26	C	2-Nitrophenol	40.000	40.008	-0.0	103	0.00
27		2,4-Dimethylphenol	40.000	36.734	8.2	96	0.00
28		bis(2-Chloroethoxy)methane	40.000	37.127	7.2	99	0.00
29	C	2,4-Dichlorophenol	40.000	39.081	2.3	102	0.00
30		1,2,4-Trichlorobenzene	40.000	37.655	5.9	101	0.00
31		Naphthalene	40.000	36.727	8.2	98	0.00
32		Benzoic acid	40.000	37.114	7.2	93	0.02
33		4-Chloroaniline	40.000	37.241	6.9	99	0.00
34	C	Hexachlorobutadiene	40.000	37.151	7.1	99	0.00
35		Caprolactam	40.000	37.876	5.3	99	0.00
36	C	4-Chloro-3-methylphenol	40.000	37.716	5.7	100	0.00
37		2-Methylnaphthalene	40.000	36.938	7.7	99	0.00
38		1-Methylnaphthalene	40.000	36.739	8.2	99	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	105	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	37.218	7.0	100	0.00
41	P	Hexachlorocyclopentadiene	40.000	37.373	6.6	96	0.00
42	S	2,4,6-Tribromophenol	80.000	79.024	1.2	105	0.00
43	C	2,4,6-Trichlorophenol	40.000	38.078	4.8	100	0.00
44		2,4,5-Trichlorophenol	40.000	38.683	3.3	102	0.00
45	S	2-Fluorobiphenyl	80.000	73.870	7.7	99	0.00
46		1,1'-Biphenyl	40.000	37.003	7.5	100	0.00
47		2-Chloronaphthalene	40.000	37.720	5.7	101	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG102121\
Data File : BG050660.D
Acq On : 21 Oct 2021 9:39
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Oct 21 11:11:18 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG100621.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Oct 14 10:56:22 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	39.317	1.7	102	0.00
49		Acenaphthylene	40.000	36.977	7.6	99	0.00
50		Dimethylphthalate	40.000	37.281	6.8	100	0.00
51		2,6-Dinitrotoluene	40.000	38.080	4.8	100	0.00
52	C	Acenaphthene	40.000	37.089	7.3	98	0.00
53		3-Nitroaniline	40.000	38.585	3.5	102	0.00
54	P	2,4-Dinitrophenol	40.000	35.335	11.7	91	0.00
55		Dibenzofuran	40.000	36.623	8.4	100	0.00
56	P	4-Nitrophenol	40.000	35.930	10.2	95	0.00
57		2,4-Dinitrotoluene	40.000	39.369	1.6	104	0.00
58		Fluorene	40.000	36.790	8.0	99	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	35.665	10.8	94	0.00
60		Diethylphthalate	40.000	36.871	7.8	100	0.00
61		4-Chlorophenyl-phenylether	40.000	37.673	5.8	101	0.00
62		4-Nitroaniline	40.000	37.986	5.0	101	0.00
63		Azobenzene	40.000	35.917	10.2	97	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	101	0.00
65		4,6-Dinitro-2-methylphenol	40.000	40.786	-2.0	104	0.00
66	c	n-Nitrosodiphenylamine	40.000	38.036	4.9	98	0.00
67		4-Bromophenyl-phenylether	40.000	39.589	1.0	102	0.00
68		Hexachlorobenzene	40.000	39.339	1.7	102	0.00
69		Atrazine	40.000	37.127	7.2	95	0.00
70	C	Pentachlorophenol	40.000	35.941	10.1	90	0.00
71		Phenanthrene	40.000	37.049	7.4	96	0.00
72		Anthracene	40.000	37.022	7.4	96	0.00
73		Carbazole	40.000	36.413	9.0	95	0.00
74		Di-n-butylphthalate	40.000	36.890	7.8	96	0.00
75	C	Fluoranthene	40.000	33.303	16.7	87	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	85	0.00
77		Benzidine	40.000	37.007	7.5	82	0.00
78		Pyrene	40.000	39.947	0.1	87	0.00
79	S	Terphenyl-d14	80.000	82.288	-2.9	90	0.00
80		Butylbenzylphthalate	40.000	39.261	1.8	85	0.00
81		Benzo(a)anthracene	40.000	38.514	3.7	85	0.00
82		3,3'-Dichlorobenzidine	40.000	37.527	6.2	82	0.00
83		Chrysene	40.000	38.456	3.9	85	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	38.276	4.3	83	0.00
85	c	Di-n-octyl phthalate	40.000	38.656	3.4	83	0.00
86	I	Perylene-d12	20.000	20.000	0.0	84	0.01
87		Indeno(1,2,3-cd)pyrene	40.000	38.758	3.1	84	0.01
88		Benzo(b)fluoranthene	40.000	38.475	3.8	83	0.00
89		Benzo(k)fluoranthene	40.000	38.151	4.6	82	0.00
90	C	Benzo(a)pyrene	40.000	38.145	4.6	82	0.00
91		Dibenzo(a,h)anthracene	40.000	38.973	2.6	84	0.01
92		Benzo(g,h,i)perylene	40.000	38.321	4.2	82	0.02

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG102121\
Data File : BG050660.D
Acq On : 21 Oct 2021 9:39
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_G
LabSampleId :
SSTDCCC040

Quant Time: Oct 21 11:11:18 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG100621.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Oct 14 10:56:22 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