Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG061721\
Data File : BG049008.D
Acq On : 17 Jun 2021 9:49
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jun 17 12:35:20 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG061121.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jun 16 12:04:50 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	74	0.00
2		1,4-Dioxane	40.000	34.841	12.9	74	0.00
3		Pyridine	40.000	36.242	9.4	73	0.00
4		n-Nitrosodimethylamine	40.000	42.499	-6.2	83	0.00
5	S	2-Fluorophenol	80.000	72.574	9.3	75	0.00
6		Aniline	40.000	34.538	13.7	71	0.00
7	S	Phenol-d6	80.000	70.755	11.6	72	0.00
8		2-Chlorophenol	40.000	35.098	12.3	74	0.00
9		Benzaldehyde	40.000	33.757	15.6	73	0.00
10	C	Phenol	40.000	35.723	10.7	75	0.00
11		bis(2-Chloroethyl)ether	40.000	32.973	17.6	68	0.00
12		1,3-Dichlorobenzene	40.000	35.404	11.5	73	0.00
13	C	1,4-Dichlorobenzene	40.000	35.660	10.9	75	0.00
14		1,2-Dichlorobenzene	40.000	35.424	11.4	75	0.00
15		Benzyl Alcohol	40.000	34.108	14.7	72	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	32.911	17.7	69	0.00
17		2-Methylphenol	40.000	34.534	13.7	71	0.00
18		Hexachloroethane	40.000	36.328	9.2	75	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	33.085	17.3	70	0.00
20		3+4-Methylphenols	40.000	34.923	12.7	72	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	72	0.00
22		Acetophenone	40.000	34.644	13.4	72	0.00
23	S	Nitrobenzene-d5	80.000	72.934	8.8	75	0.00
24		Nitrobenzene	40.000	36.144	9.6	75	0.00
25		Isophorone	40.000	32.431	18.9	68	0.00
26	C	2-Nitrophenol	40.000	39.303	1.7	80	0.00
27		2,4-Dimethylphenol	40.000	33.319	16.7	70	0.00
28		bis(2-Chloroethoxy)methane	40.000	32.503	18.7	67	0.00
29	C	2,4-Dichlorophenol	40.000	34.945	12.6	72	0.00
30		1,2,4-Trichlorobenzene	40.000	35.381	11.5	74	0.00
31		Naphthalene	40.000	33.526	16.2	70	0.00
32		Benzoic acid	40.000	40.865	-2.2	82	-0.01
33		4-Chloroaniline	40.000	34.875	12.8	73	0.00
34	C	Hexachlorobutadiene	40.000	34.258	14.4	72	0.00
35		Caprolactam	40.000	40.813	-2.0	85	0.00
36	C	4-Chloro-3-methylphenol	40.000	34.872	12.8	72	0.00
37		2-Methylnaphthalene	40.000	32.888	17.8	69	0.00
38		1-Methylnaphthalene	40.000	33.144	17.1	69	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	71	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	33.776	15.6	72	0.00
41	P	Hexachlorocyclopentadiene	40.000	34.618	13.5	72	0.00
42	S	2,4,6-Tribromophenol	80.000	72.024	10.0	74	0.00
43	C	2,4,6-Trichlorophenol	40.000	37.127	7.2	76	0.00
44		2,4,5-Trichlorophenol	40.000	39.847	0.4	82	0.00
45	S	2-Fluorobiphenyl	80.000	67.734	15.3	70	0.00
46		1,1'-Biphenyl	40.000	33.306	16.7	69	0.00
47		2-Chloronaphthalene	40.000	34.629	13.4	73	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG061721\
Data File : BG049008.D
Acq On : 17 Jun 2021 9:49
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jun 17 12:35:20 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG061121.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jun 16 12:04:50 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	40.670	-1.7	83	0.00
49		Acenaphthylene	40.000	33.111	17.2	69	0.00
50		Dimethylphthalate	40.000	34.993	12.5	73	0.00
51		2,6-Dinitrotoluene	40.000	39.366	1.6	80	0.00
52	C	Acenaphthene	40.000	34.442	13.9	71	0.00
53		3-Nitroaniline	40.000	38.211	4.5	78	0.00
54	P	2,4-Dinitrophenol	40.000	44.204	-10.5	89	0.00
55		Dibenzofuran	40.000	33.438	16.4	70	0.00
56	P	4-Nitrophenol	40.000	39.564	1.1	78	0.00
57		2,4-Dinitrotoluene	40.000	45.655	-14.1	83	0.00
58		Fluorene	40.000	33.563	16.1	71	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	36.713	8.2	75	0.00
60		Diethylphthalate	40.000	34.171	14.6	72	0.00
61		4-Chlorophenyl-phenylether	40.000	34.140	14.6	73	0.00
62		4-Nitroaniline	40.000	39.603	1.0	80	0.00
63		Azobenzene	40.000	32.148	19.6	68	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	72	0.00
65		4,6-Dinitro-2-methylphenol	40.000	47.766	-19.4	92	0.00
66	c	n-Nitrosodiphenylamine	40.000	32.714	18.2	69	0.00
67		4-Bromophenyl-phenylether	40.000	32.947	17.6	69	0.00
68		Hexachlorobenzene	40.000	34.383	14.0	74	0.00
69		Atrazine	40.000	37.459	6.4	73	0.00
70	C	Pentachlorophenol	40.000	35.376	11.6	73	0.00
71		Phenanthrene	40.000	34.087	14.8	73	0.00
72		Anthracene	40.000	33.638	15.9	73	0.00
73		Carbazole	40.000	34.629	13.4	74	0.00
74		Di-n-butylphthalate	40.000	36.471	8.8	79	0.00
75	C	Fluoranthene	40.000	36.247	9.4	79	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	78	0.00
77		Benzidine	40.000	43.994	-10.0	87	0.00
78		Pyrene	40.000	34.721	13.2	79	0.00
79	S	Terphenyl-d14	80.000	73.139	8.6	84	0.00
80		Butylbenzylphthalate	40.000	37.210	7.0	85	0.00
81		Benzo(a)anthracene	40.000	35.362	11.6	80	0.00
82		3,3'-Dichlorobenzidine	40.000	39.670	0.8	89	0.00
83		Chrysene	40.000	35.443	11.4	81	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	37.645	5.9	87	0.00
85	c	Di-n-octyl phthalate	40.000	37.412	6.5	85	0.00
86	I	Perylene-d12	20.000	20.000	0.0	81	0.00
87		Indeno(1,2,3-cd)pyrene	40.000	38.895	2.8	88	0.00
88		Benzo(b)fluoranthene	40.000	37.155	7.1	85	0.00
89		Benzo(k)fluoranthene	40.000	36.034	9.9	81	0.00
90	C	Benzo(a)pyrene	40.000	36.896	7.8	84	0.00
91		Dibenzo(a,h)anthracene	40.000	39.325	1.7	88	0.00
92		Benzo(g,h,i)perylene	40.000	37.790	5.5	85	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG061721\
Data File : BG049008.D
Acq On : 17 Jun 2021 9:49
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_G
LabSampleId :
SSTDCCC040

Quant Time: Jun 17 12:35:20 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG061121.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jun 16 12:04:50 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