Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP033121\
Data File : BP005117.D
Acq On : 31 Mar 2021 09:33
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_P
LabSampleId : SSTDCCC040

Quant Time: Mar 31 11:24:00 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP032321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Mar 26 10:07:03 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	35.926	10.2	91	0.00
3		Pyridine	40.000	38.477	3.8	85	0.00
4		n-Nitrosodimethylamine	40.000	37.990	5.0	88	0.00
5	S	2-Fluorophenol	80.000	76.038	5.0	90	0.00
6		Aniline	40.000	38.714	3.2	90	-0.01
7	S	Phenol-d6	80.000	78.852	1.4	92	0.00
8		2-Chlorophenol	40.000	38.791	3.0	91	-0.01
9		Benzaldehyde	40.000	35.018	12.5	87	0.00
10	C	Phenol	40.000	39.312	1.7	91	0.00
11		bis(2-Chloroethyl)ether	40.000	39.064	2.3	91	0.00
12		1,3-Dichlorobenzene	40.000	38.985	2.5	91	0.00
13	C	1,4-Dichlorobenzene	40.000	38.580	3.6	91	0.00
14		1,2-Dichlorobenzene	40.000	39.238	1.9	93	0.00
15		Benzyl Alcohol	40.000	39.048	2.4	88	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	40.228	-0.6	93	0.00
17		2-Methylphenol	40.000	40.819	-2.0	94	0.00
18		Hexachloroethane	40.000	38.886	2.8	92	-0.01
19	P	n-Nitroso-di-n-propylamine	40.000	38.831	2.9	88	0.00
20		3+4-Methylphenols	40.000	40.111	-0.3	91	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	93	-0.01
22		Acetophenone	40.000	38.417	4.0	90	0.00
23	S	Nitrobenzene-d5	80.000	76.560	4.3	89	-0.01
24		Nitrobenzene	40.000	37.867	5.3	89	-0.01
25		Isophorone	40.000	39.173	2.1	91	0.00
26	C	2-Nitrophenol	40.000	41.013	-2.5	94	0.00
27		2,4-Dimethylphenol	40.000	39.632	0.9	93	-0.01
28		bis(2-Chloroethoxy)methane	40.000	40.140	-0.4	91	0.00
29	C	2,4-Dichlorophenol	40.000	39.655	0.9	91	0.00
30		1,2,4-Trichlorobenzene	40.000	38.680	3.3	91	0.00
31		Naphthalene	40.000	39.112	2.2	92	0.00
32		Benzoic acid	40.000	39.289	1.8	87	0.00
33		4-Chloroaniline	40.000	40.305	-0.8	94	0.00
34	C	Hexachlorobutadiene	40.000	38.538	3.7	89	0.00
35		Caprolactam	40.000	40.595	-1.5	93	0.00
36	C	4-Chloro-3-methylphenol	40.000	39.891	0.3	92	0.00
37		2-Methylnaphthalene	40.000	39.570	1.1	92	-0.01
38		1-Methylnaphthalene	40.000	39.435	1.4	91	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	92	-0.01
40		1,2,4,5-Tetrachlorobenzene	40.000	38.235	4.4	88	-0.01
41	P	Hexachlorocyclopentadiene	40.000	38.425	3.9	86	0.00
42	S	2,4,6-Tribromophenol	80.000	77.298	3.4	89	-0.01
43	C	2,4,6-Trichlorophenol	40.000	39.999	0.0	90	0.00
44		2,4,5-Trichlorophenol	40.000	39.823	0.4	91	0.00
45	S	2-Fluorobiphenyl	80.000	77.100	3.6	90	0.00
46		1,1'-Biphenyl	40.000	38.492	3.8	89	-0.01
47		2-Chloronaphthalene	40.000	38.309	4.2	90	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP033121\
Data File : BP005117.D
Acq On : 31 Mar 2021 09:33
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_P
LabSampleId :
SSTDCCC040

Quant Time: Mar 31 11:24:00 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP032321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Mar 26 10:07:03 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	39.457	1.4	88	0.00
49		Acenaphthylene	40.000	39.415	1.5	91	0.00
50		Dimethylphthalate	40.000	38.601	3.5	91	-0.01
51		2,6-Dinitrotoluene	40.000	38.981	2.5	89	0.00
52	C	Acenaphthene	40.000	38.833	2.9	91	-0.01
53		3-Nitroaniline	40.000	40.061	-0.2	88	0.00
54	P	2,4-Dinitrophenol	40.000	37.412	6.5	84	0.00
55		Dibenzofuran	40.000	38.595	3.5	90	0.00
56	P	4-Nitrophenol	40.000	40.182	-0.5	88	-0.01
57		2,4-Dinitrotoluene	40.000	40.183	-0.5	90	0.00
58		Fluorene	40.000	38.848	2.9	90	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	38.202	4.5	87	0.00
60		Diethylphthalate	40.000	39.322	1.7	90	-0.01
61		4-Chlorophenyl-phenylether	40.000	38.119	4.7	91	-0.01
62		4-Nitroaniline	40.000	39.874	0.3	88	0.00
63		Azobenzene	40.000	38.026	4.9	88	-0.01
64	I	Phenanthrene-d10	20.000	20.000	0.0	91	0.00
65		4,6-Dinitro-2-methylphenol	40.000	41.333	-3.3	87	0.00
66	c	n-Nitrosodiphenylamine	40.000	40.432	-1.1	91	0.00
67		4-Bromophenyl-phenylether	40.000	40.048	-0.1	88	0.00
68		Hexachlorobenzene	40.000	39.944	0.1	89	0.00
69		Atrazine	40.000	41.205	-3.0	88	0.00
70	C	Pentachlorophenol	40.000	40.859	-2.1	91	0.00
71		Phenanthrene	40.000	39.445	1.4	88	0.00
72		Anthracene	40.000	39.951	0.1	89	-0.01
73		Carbazole	40.000	40.715	-1.8	91	-0.01
74		Di-n-butylphthalate	40.000	41.763	-4.4	91	-0.01
75	C	Fluoranthene	40.000	39.581	1.0	89	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	91	-0.01
77		Benzidine	40.000	40.213	-0.5	79	0.00
78		Pyrene	40.000	40.129	-0.3	89	0.00
79	S	Terphenyl-d14	80.000	79.589	0.5	89	-0.01
80		Butylbenzylphthalate	40.000	42.713	-6.8	91	-0.01
81		Benzo(a)anthracene	40.000	39.575	1.1	90	-0.01
82		3,3'-Dichlorobenzidine	40.000	39.942	0.1	87	0.00
83		Chrysene	40.000	39.358	1.6	89	-0.01
84		Bis(2-ethylhexyl)phthalate	40.000	42.414	-6.0	90	-0.01
85	c	Di-n-octyl phthalate	40.000	42.975	-7.4	93	-0.01
86	I	Perylene-d12	20.000	20.000	0.0	92	-0.02
87		Indeno(1,2,3-cd)pyrene	40.000	39.086	2.3	89	-0.03
88		Benzo(b)fluoranthene	40.000	40.064	-0.2	92	-0.01
89		Benzo(k)fluoranthene	40.000	38.128	4.7	86	-0.01
90	C	Benzo(a)pyrene	40.000	39.913	0.2	91	-0.02
91		Dibenzo(a,h)anthracene	40.000	39.341	1.6	90	-0.02
92		Benzo(g,h,i)perylene	40.000	39.529	1.2	90	-0.03

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP033121\
Data File : BP005117.D
Acq On : 31 Mar 2021 09:33
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_P
LabSampleId :
SSTDCCC040

Quant Time: Mar 31 11:24:00 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP032321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Mar 26 10:07:03 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