Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_N\DATA\BN022818\
Data File : BN000198.D
Acq On : 28 Feb 2018 09:03
Operator : JU/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_N
LabSampleId : SSTDCCC0.4

Quant Time: Feb 28 16:59:48 2018
Quant Method : Z:\HPCHEM1\BNA_N\METHODS\8270-SIM-BN022718.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Feb 27 15:07:57 2018
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	0.400	0.400	0.0	92	0.00
2		1,4-Dioxane	0.400	0.428	-7.0	104	0.00
3		n-Nitrosodimethylamine	0.400	0.374	6.5	86	0.00
4	S	2-Fluorophenol	0.400	0.435	-8.7	97	0.00
5	S	Phenol-d6	0.400	0.417	-4.2	96	0.00
6		bis(2-Chloroethyl)ether	0.400	0.415	-3.7	91	0.00
7	I	Naphthalene-d8	0.400	0.400	0.0	92	0.00
8	S	Nitrobenzene-d5	0.400	0.394	1.5	90	0.00
9		Naphthalene	0.400	0.405	-1.3	91	0.00
10		Hexachlorobutadiene	0.400	0.411	-2.7	91	0.00
11	SURR	2-Methylnaphthalene-d10	0.400	0.404	-1.0	91	0.00
12		2-Methylnaphthalene	0.400	0.402	-0.5	91	0.00
13	I	Acenaphthene-d10	0.400	0.400	0.0	96	0.00
14	S	2,4,6-Tribromophenol	0.400	0.385	3.8	95	0.00
15	S	2-Fluorobiphenyl	0.400	0.394	1.5	93	0.00
16		Acenaphthylene	0.400	0.358	10.5	88	0.00
17		Acenaphthene	0.400	0.384	4.0	92	-0.01
18		Fluorene	0.400	0.380	5.0	92	0.00
19	I	Phenanthrene-d10	0.400	0.400	0.0	97	0.00
20		4-Bromophenyl-phenylether	0.400	0.377	5.8	92	0.00
21		Hexachlorobenzene	0.400	0.385	3.8	93	0.00
22		Pentachlorophenol	0.400	0.381	4.8	93	0.00
23		Phenanthrene	0.400	0.375	6.3	90	0.00
24		Anthracene	0.400	0.356	11.0	88	0.00
25	SURR	Fluoranthene-d10	0.400	0.359	10.3	88	0.00
26		Fluoranthene	0.400	0.357	10.8	87	0.00
27	I	Chrysene-d12	0.400	0.400	0.0	88	0.00
28		Pyrene	0.400	0.401	-0.3	88	0.00
29	S	Terphenyl-d14	0.400	0.396	1.0	86	0.00
30		Benzo(a)anthracene	0.400	0.377	5.8	83	0.00
31		Chrysene	0.400	0.386	3.5	83	0.00
32		Bis(2-ethylhexyl)phthalate	0.400	0.373	6.8	89	0.00
33		Indeno(1,2,3-cd)pyrene	0.400	0.341	14.8	72	0.00
34	I	Perylene-d12	0.400	0.400	0.0	81	0.00
35		Benzo(b)fluoranthene	0.400	0.381	4.8	78	0.00
36		Benzo(k)fluoranthene	0.400	0.387	3.3	77	0.00
37	C	Benzo(a)pyrene	0.400	0.370	7.5	76	0.00
38		Dibenzo(a,h)anthracene	0.400	0.350	12.5	72	0.00
39		Benzo(g,h,i)perylene	0.400	0.363	9.3	73	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_N\DATA\BN022818\
Data File : BN000198.D
Acq On : 28 Feb 2018 09:03
Operator : JU/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_N
LabSampleId : SSTDCCC0.4

Quant Time: Feb 28 16:59:48 2018
Quant Method : Z:\HPCHEM1\BNA_N\METHODS\8270-SIM-BN022718.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Feb 27 15:07:57 2018
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