Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN042820\
Data File : BN010640.D
Acq On : 29 Apr 2020 01:28
Operator : CG/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_N
LabSampleId : SSTDCCC0.4

Quant Time: Apr 29 02:04:41 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN041320.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Apr 28 11:13:37 2020
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	102	0.00
2		1,4-Dioxane	0.400	0.403	-0.8	103	0.00
3		n-Nitrosodimethylamine	0.400	0.432	-8.0	130	0.00
4	S	2-Fluorophenol	0.400	0.422	-5.5	107	0.00
5	S	Phenol-d6	0.400	0.435	-8.7	110	0.00
6		bis(2-Chloroethyl)ether	0.400	0.413	-3.2	103	0.00
7	I	Naphthalene-d8	0.400	0.400	0.0	108	0.00
8	S	Nitrobenzene-d5	0.400	0.407	-1.7	115	0.00
9		Naphthalene	0.400	0.410	-2.5	112	0.00
10		Hexachlorobutadiene	0.400	0.398	0.5	107	0.00
11	SURR	2-Methylnaphthalene-d10	0.400	0.462	-15.5	126	0.00
12		2-Methylnaphthalene	0.400	0.414	-3.5	113	0.00
13	I	Acenaphthene-d10	0.400	0.400	0.0	111	0.00
14	S	2,4,6-Tribromophenol	0.400	0.346	13.5	100	0.00
15	S	2-Fluorobiphenyl	0.400	0.394	1.5	108	0.00
16		Acenaphthylene	0.400	0.381	4.8	108	0.00
17		Acenaphthene	0.400	0.411	-2.7	115	0.00
18		Fluorene	0.400	0.402	-0.5	112	0.00
19	I	Phenanthrene-d10	0.400	0.400	0.0	111	0.00
20		4-Bromophenyl-phenylether	0.400	0.386	3.5	110	0.00
21		Hexachlorobenzene	0.400	0.415	-3.7	114	0.00
22		Pentachlorophenol	0.400	0.329	17.8	108	0.00
23		Phenanthrene	0.400	0.406	-1.5	113	0.00
24		Anthracene	0.400	0.380	5.0	108	0.00
25	SURR	Fluoranthene-d10	0.400	0.428	-7.0	119	0.00
26		Fluoranthene	0.400	0.382	4.5	106	0.00
27	I	Chrysene-d12	0.400	0.400	0.0	90	0.00
28		Pyrene	0.400	0.458	-14.5	105	0.00
29	S	Terphenyl-d14	0.400	0.460	-15.0	107	0.00
30		Benzo(a)anthracene	0.400	0.396	1.0	91	-0.01
31		Chrysene	0.400	0.397	0.8	91	0.00
32		Bis(2-ethylhexyl)phthalate	0.400	0.394	1.5	94	0.00
33		Indeno(1,2,3-cd)pyrene	0.400	0.463	-15.8	105	-0.02
34	I	Perylene-d12	0.400	0.400	0.0	96	-0.02
35		Benzo(b)fluoranthene	0.400	0.396	1.0	96	0.00
36		Benzo(k)fluoranthene	0.400	0.403	-0.8	97	0.00
37	C	Benzo(a)pyrene	0.400	0.395	1.3	96	-0.02
38		Dibenzo(a,h)anthracene	0.400	0.432	-8.0	104	-0.01
39		Benzo(g,h,i)perylene	0.400	0.432	-8.0	103	-0.02

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Compound	Amount	Calc.	%Dev	Area%	Dev(min)

(#) = Out of Range SPCC's out = 0 CCC's out = 0