Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN061620\
Data File : BN010919.D
Acq On : 16 Jun 2020 22:27
Operator : CG/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument : BNA_N
LabSampleId : SSTDCCC0.4

Quant Time: Jun 17 03:37:45 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN061020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jun 16 08:49:35 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	118	0.00
2		1,4-Dioxane	0.400	0.511	-27.7#	144	0.00
3		n-Nitrosodimethylamine	0.400	0.395	1.3	132	0.02
4	S	2-Fluorophenol	0.400	0.531	-32.8#	153	0.00
5	S	Phenol-d6	0.400	0.490	-22.5	145	0.00
6		bis(2-Chloroethyl)ether	0.400	0.480	-20.0	142	0.00
7	I	Naphthalene-d8	0.400	0.400	0.0	126	-0.01
8	S	Nitrobenzene-d5	0.400	0.425	-6.2	129	0.00
9		Naphthalene	0.400	0.426	-6.5	128	0.00
10		Hexachlorobutadiene	0.400	0.404	-1.0	118	-0.01
11	SURR	2-Methylnaphthalene-d10	0.400	0.456	-14.0	137	0.00
12		2-Methylnaphthalene	0.400	0.481	-20.2	148	0.00
13	I	Acenaphthene-d10	0.400	0.400	0.0	128	0.00
14	S	2,4,6-Tribromophenol	0.400	0.395	1.3	130	0.00
15	S	2-Fluorobiphenyl	0.400	0.430	-7.5	129	0.00
16		Acenaphthylene	0.400	0.417	-4.2	132	0.00
17		Acenaphthene	0.400	0.421	-5.2	129	-0.01
18		Fluorene	0.400	0.419	-4.7	130	-0.01
19	I	Phenanthrene-d10	0.400	0.400	0.0	128	-0.01
20		4-Bromophenyl-phenylether	0.400	0.394	1.5	123	0.00
21		Hexachlorobenzene	0.400	0.412	-3.0	126	0.00
22		Pentachlorophenol	0.400	0.387	3.3	128	0.00
23		Phenanthrene	0.400	0.418	-4.5	130	-0.01
24		Anthracene	0.400	0.408	-2.0	132	0.00
25	SURR	Fluoranthene-d10	0.400	0.461	-15.3	146	0.00
26		Fluoranthene	0.400	0.466	-16.5	149	0.00
27	I	Chrysene-d12	0.400	0.400	0.0	132	-0.01
28		Pyrene	0.400	0.459	-14.8	146	0.00
29	S	Terphenyl-d14	0.400	0.444	-11.0	140	0.00
30		Benzo(a)anthracene	0.400	0.416	-4.0	139	-0.01
31		Chrysene	0.400	0.432	-8.0	132	-0.01
32		Bis(2-ethylhexyl)phthalate	0.400	0.444	-11.0	150	-0.01
33		Indeno(1,2,3-cd)pyrene	0.400	0.483	-20.7	152	-0.02
34	I	Perylene-d12	0.400	0.400	0.0	141	-0.01
35		Benzo(b)fluoranthene	0.400	0.403	-0.8	137	-0.01
36		Benzo(k)fluoranthene	0.400	0.410	-2.5	137	0.00
37	C	Benzo(a)pyrene	0.400	0.417	-4.2	146	-0.01
38		Dibenzo(a,h)anthracene	0.400	0.439	-9.7	152	-0.02
39		Benzo(g,h,i)perylene	0.400	0.398	0.5	132	-0.02

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

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