Evaluate Continuing Calibration Report

Data Path : \\74.0.250.170\terastorage\svoasrv\HPCHEM1\BNA_N\Data\BN062819\
Data File : BN006639.D
Acq On : 29 Jun 2019 06:48
Operator : HP/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_N
LabSampleId : SSTDCCC0.4

Quant Time: Jul 01 08:14:29 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN062819.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 28 13:47:56 2019
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	122	0.00
2		1,4-Dioxane	0.400	0.382	4.5	125	0.00
3		n-Nitrosodimethylamine	0.400	0.344	14.0	119	0.00
4	S	2-Fluorophenol	0.400	0.378	5.5	121	0.00
5	S	Phenol-d6	0.400	0.396	1.0	128	0.00
6		bis(2-Chloroethyl)ether	0.400	0.410	-2.5	120	0.00
7	I	Naphthalene-d8	0.400	0.400	0.0	125	0.00
8	S	Nitrobenzene-d5	0.400	0.368	8.0	122	0.00
9		Naphthalene	0.400	0.411	-2.7	125	0.00
10		Hexachlorobutadiene	0.400	0.413	-3.2	125	0.00
11	SURR	2-Methylnaphthalene-d10	0.400	0.413	-3.2	125	0.00
12		2-Methylnaphthalene	0.400	0.416	-4.0	125	0.00
13	I	Acenaphthene-d10	0.400	0.400	0.0	126	-0.01
14	S	2,4,6-Tribromophenol	0.400	0.374	6.5	131	0.00
15	S	2-Fluorobiphenyl	0.400	0.399	0.3	126	-0.01
16		Acenaphthylene	0.400	0.394	1.5	126	0.00
17		Acenaphthene	0.400	0.412	-3.0	127	0.00
18		Fluorene	0.400	0.405	-1.3	126	-0.01
19	I	Phenanthrene-d10	0.400	0.400	0.0	125	-0.01
20		4-Bromophenyl-phenylether	0.400	0.395	1.3	125	-0.01
21		Hexachlorobenzene	0.400	0.412	-3.0	125	0.00
22		Pentachlorophenol	0.400	0.436	-9.0	138	-0.01
23		Phenanthrene	0.400	0.403	-0.8	125	0.00
24		Anthracene	0.400	0.384	4.0	124	0.00
25	SURR	Fluoranthene-d10	0.400	0.394	1.5	123	0.00
26		Fluoranthene	0.400	0.403	-0.8	123	0.00
27	I	Chrysene-d12	0.400	0.400	0.0	115	-0.01
28		Pyrene	0.400	0.434	-8.5	122	0.00
29	S	Terphenyl-d14	0.400	0.429	-7.2	123	0.00
30		Benzo(a)anthracene	0.400	0.410	-2.5	120	-0.02
31		Chrysene	0.400	0.411	-2.7	112	-0.02
32		Bis(2-ethylhexyl)phthalate	0.400	0.403	-0.8	129	-0.01
33		Indeno(1,2,3-cd)pyrene	0.400	0.373	6.8	92	-0.04
34	I	Perylene-d12	0.400	0.400	0.0	99	-0.04
35		Benzo(b)fluoranthene	0.400	0.427	-6.7	105	-0.03
36		Benzo(k)fluoranthene	0.400	0.434	-8.5	105	-0.04
37	C	Benzo(a)pyrene	0.400	0.424	-6.0	101	-0.04
38		Dibenzo(a,h)anthracene	0.400	0.395	1.3	92	-0.04
39		Benzo(g,h,i)perylene	0.400	0.391	2.3	91	-0.04

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

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