Evaluate Continuing Calibration Report

Data Path : \\74.0.250.170\terastorage\svoasrv\HPCHEM1\BNA_N\Data\BN062819\
Data File : BN006637.D
Acq On : 29 Jun 2019 05:05
Operator : HP/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_N
LabSampleId : SSTDCCC0.4

Quant Time: Jun 29 05:42:23 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN062819.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat Jun 29 01:19:02 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	95	0.00
2		1,4-Dioxane	0.400	0.401	-0.3	102	0.00
3		n-Nitrosodimethylamine	0.400	0.378	5.5	102	0.00
4	S	2-Fluorophenol	0.400	0.412	-3.0	103	0.00
5	S	Phenol-d6	0.400	0.424	-6.0	107	0.00
6		bis(2-Chloroethyl)ether	0.400	0.420	-5.0	96	0.00
7	I	Naphthalene-d8	0.400	0.400	0.0	97	0.00
8	S	Nitrobenzene-d5	0.400	0.402	-0.5	105	0.00
9		Naphthalene	0.400	0.426	-6.5	101	0.00
10		Hexachlorobutadiene	0.400	0.429	-7.2	101	0.00
11	SURR	2-Methylnaphthalene-d10	0.400	0.429	-7.2	102	0.00
12		2-Methylnaphthalene	0.400	0.436	-9.0	103	0.00
13	I	Acenaphthene-d10	0.400	0.400	0.0	100	-0.01
14	S	2,4,6-Tribromophenol	0.400	0.419	-4.7	117	0.00
15	S	2-Fluorobiphenyl	0.400	0.410	-2.5	103	-0.01
16		Acenaphthylene	0.400	0.414	-3.5	106	0.00
17		Acenaphthene	0.400	0.430	-7.5	105	0.00
18		Fluorene	0.400	0.425	-6.2	105	0.00
19	I	Phenanthrene-d10	0.400	0.400	0.0	99	-0.01
20		4-Bromophenyl-phenylether	0.400	0.425	-6.2	106	-0.01
21		Hexachlorobenzene	0.400	0.430	-7.5	103	0.00
22		Pentachlorophenol	0.400	0.544	-36.0#	157	-0.02
23		Phenanthrene	0.400	0.424	-6.0	104	0.00
24		Anthracene	0.400	0.413	-3.2	105	0.00
25	SURR	Fluoranthene-d10	0.400	0.423	-5.7	104	0.00
26		Fluoranthene	0.400	0.432	-8.0	104	0.00
27	I	Chrysene-d12	0.400	0.400	0.0	94	-0.02
28		Pyrene	0.400	0.450	-12.5	103	0.00
29	S	Terphenyl-d14	0.400	0.439	-9.7	103	0.00
30		Benzo(a)anthracene	0.400	0.421	-5.2	100	-0.02
31		Chrysene	0.400	0.434	-8.5	96	-0.02
32		Bis(2-ethylhexyl)phthalate	0.400	0.536	-34.0#	140	-0.02
33		Indeno(1,2,3-cd)pyrene	0.400	0.376	6.0	76	-0.05
34	I	Perylene-d12	0.400	0.400	0.0	80	-0.05
35		Benzo(b)fluoranthene	0.400	0.441	-10.2	87	-0.04
36		Benzo(k)fluoranthene	0.400	0.442	-10.5	86	-0.04
37	C	Benzo(a)pyrene	0.400	0.427	-6.7	82	-0.04
38		Dibenzo(a,h)anthracene	0.400	0.405	-1.3	76	-0.04
39		Benzo(g,h,i)perylene	0.400	0.405	-1.3	76	-0.05

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

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