Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN070119\
Data File : BN006652.D
Acq On : 01 Jul 2019 08:39
Operator : HP/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_N
LabSampleId : SSTDCCC0.4

Quant Time: Jul 02 05:03:21 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	108	0.00
2		1,4-Dioxane	0.400	0.423	-5.7	122	0.00
3		n-Nitrosodimethylamine	0.400	0.378	5.5	115	0.00
4	S	2-Fluorophenol	0.400	0.410	-2.5	115	0.00
5	S	Phenol-d6	0.400	0.395	1.3	112	0.00
6		bis(2-Chloroethyl)ether	0.400	0.403	-0.8	104	0.00
7	I	Naphthalene-d8	0.400	0.400	0.0	97	0.00
8	S	Nitrobenzene-d5	0.400	0.406	-1.5	106	0.00
9		Naphthalene	0.400	0.418	-4.5	99	-0.01
10		Hexachlorobutadiene	0.400	0.439	-9.7	104	-0.01
11	SURR	2-Methylnaphthalene-d10	0.400	0.392	2.0	93	0.00
12		2-Methylnaphthalene	0.400	0.399	0.3	94	0.00
13	I	Acenaphthene-d10	0.400	0.400	0.0	85	-0.01
14	S	2,4,6-Tribromophenol	0.400	0.379	5.3	90	0.00
15	S	2-Fluorobiphenyl	0.400	0.422	-5.5	90	-0.01
16		Acenaphthylene	0.400	0.408	-2.0	88	0.00
17		Acenaphthene	0.400	0.423	-5.7	88	0.00
18		Fluorene	0.400	0.403	-0.8	84	-0.01
19	I	Phenanthrene-d10	0.400	0.400	0.0	80	0.00
20		4-Bromophenyl-phenylether	0.400	0.424	-6.0	85	-0.01
21		Hexachlorobenzene	0.400	0.436	-9.0	84	0.00
22		Pentachlorophenol	0.400	0.430	-7.5	86	0.00
23		Phenanthrene	0.400	0.413	-3.2	82	0.00
24		Anthracene	0.400	0.404	-1.0	83	0.00
25	SURR	Fluoranthene-d10	0.400	0.410	-2.5	81	0.00
26		Fluoranthene	0.400	0.417	-4.2	81	0.00
27	I	Chrysene-d12	0.400	0.400	0.0	74	0.00
28		Pyrene	0.400	0.449	-12.2	81	0.00
29	S	Terphenyl-d14	0.400	0.435	-8.7	80	0.00
30		Benzo(a)anthracene	0.400	0.424	-6.0	80	-0.01
31		Chrysene	0.400	0.429	-7.2	75	-0.01
32		Bis(2-ethylhexyl)phthalate	0.400	0.469	-17.2	97	0.00
33		Indeno(1,2,3-cd)pyrene	0.400	0.533	-33.3#	85	-0.01
34	I	Perylene-d12	0.400	0.400	0.0	76	-0.01
35		Benzo(b)fluoranthene	0.400	0.404	-1.0	76	-0.01
36		Benzo(k)fluoranthene	0.400	0.402	-0.5	75	-0.01
37	C	Benzo(a)pyrene	0.400	0.428	-7.0	78	-0.01
38		Dibenzo(a,h)anthracene	0.400	0.474	-18.5	84	-0.01
39		Benzo(g,h,i)perylene	0.400	0.488	-22.0	87	-0.01

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

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