Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN100319\
Data File : BN008002.D
Acq On : 04 Oct 2019 04:10
Operator : JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_N
LabSampleId : SSTDCCC0.4

Quant Time: Oct 04 05:43:10 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN091819.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Oct 03 18:39:49 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	81	-0.02
2		1,4-Dioxane	0.400	0.367	8.3	88	-0.02
3		n-Nitrosodimethylamine	0.400	0.604	-51.0#	81	-0.02
4	S	2-Fluorophenol	0.400	0.362	9.5	84	-0.02
5	S	Phenol-d6	0.400	0.400	0.0	89	0.00
6		bis(2-Chloroethyl)ether	0.400	0.487	-21.7	112	-0.02
7	I	Naphthalene-d8	0.400	0.400	0.0	89	-0.01
8	S	Nitrobenzene-d5	0.400	0.402	-0.5	92	-0.01
9		Naphthalene	0.400	0.411	-2.7	92	-0.01
10		Hexachlorobutadiene	0.400	0.431	-7.7	96	-0.03
11	SURR	2-Methylnaphthalene-d10	0.400	0.424	-6.0	98	-0.02
12		2-Methylnaphthalene	0.400	0.425	-6.2	98	-0.02
13	I	Acenaphthene-d10	0.400	0.400	0.0	89	-0.01
14	S	2,4,6-Tribromophenol	0.400	0.382	4.5	102	-0.01
15	S	2-Fluorobiphenyl	0.400	0.380	5.0	91	-0.01
16		Acenaphthylene	0.400	0.395	1.3	98	-0.02
17		Acenaphthene	0.400	0.369	7.8	90	-0.02
18		Fluorene	0.400	0.376	6.0	91	-0.01
19	I	Phenanthrene-d10	0.400	0.400	0.0	81	-0.01
20		4-Bromophenyl-phenylether	0.400	0.409	-2.2	84	-0.02
21		Hexachlorobenzene	0.400	0.436	-9.0	90	-0.02
22		Pentachlorophenol	0.400	0.448	-12.0	104	-0.01
23		Phenanthrene	0.400	0.403	-0.8	83	-0.02
24		Anthracene	0.400	0.423	-5.7	89	-0.01
25	SURR	Fluoranthene-d10	0.400	0.404	-1.0	83	-0.02
26		Fluoranthene	0.400	0.404	-1.0	83	-0.02
27	I	Chrysene-d12	0.400	0.400	0.0	73	-0.02
28		Pyrene	0.400	0.445	-11.2	85	-0.02
29	S	Terphenyl-d14	0.400	0.438	-9.5	83	-0.02
30		Benzo(a)anthracene	0.400	0.408	-2.0	78	-0.01
31		Chrysene	0.400	0.410	-2.5	77	-0.01
32		Bis(2-ethylhexyl)phthalate	0.400	0.479	-19.7	98	-0.03
33		Indeno(1,2,3-cd)pyrene	0.400	0.375	6.3	71	-0.03
34	I	Perylene-d12	0.400	0.400	0.0	70	-0.02
35		Benzo(b)fluoranthene	0.400	0.404	-1.0	75	-0.02
36		Benzo(k)fluoranthene	0.400	0.385	3.8	70	-0.02
37	C	Benzo(a)pyrene	0.400	0.400	0.0	74	-0.02
38		Dibenzo(a,h)anthracene	0.400	0.384	4.0	70	-0.03
39		Benzo(g,h,i)perylene	0.400	0.391	2.3	71	-0.03

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

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