Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100219\
Data File : BN007966.D
Acq On : 02 Oct 2019 15:30
Operator : JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_N
LabSampleId : SSTDCCC0.4

Quant Time: Oct 03 06:46:03 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN091819.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Sep 18 18:24:40 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	85	-0.02
2		1,4-Dioxane	0.400	0.468	-17.0	118	-0.02
3		n-Nitrosodimethylamine	0.400	0.640	-60.0#	90	-0.02
4	S	2-Fluorophenol	0.400	0.387	3.3	94	0.00
5	S	Phenol-d6	0.400	0.411	-2.7	96	0.00
6		bis(2-Chloroethyl)ether	0.400	0.516	-29.0#	125	-0.02
7	I	Naphthalene-d8	0.400	0.400	0.0	88	-0.01
8	S	Nitrobenzene-d5	0.400	0.454	-13.5	102	-0.01
9		Naphthalene	0.400	0.422	-5.5	93	-0.01
10		Hexachlorobutadiene	0.400	0.442	-10.5	96	-0.03
11	SURR	2-Methylnaphthalene-d10	0.400	0.437	-9.2	99	-0.01
12		2-Methylnaphthalene	0.400	0.440	-10.0	99	-0.02
13	I	Acenaphthene-d10	0.400	0.400	0.0	91	-0.01
14	S	2,4,6-Tribromophenol	0.400	0.451	-12.7	123	-0.01
15	S	2-Fluorobiphenyl	0.400	0.385	3.8	94	-0.01
16		Acenaphthylene	0.400	0.391	2.3	99	-0.02
17		Acenaphthene	0.400	0.374	6.5	93	-0.02
18		Fluorene	0.400	0.388	3.0	96	-0.01
19	I	Phenanthrene-d10	0.400	0.400	0.0	89	-0.01
20		4-Bromophenyl-phenylether	0.400	0.420	-5.0	95	-0.02
21		Hexachlorobenzene	0.400	0.430	-7.5	97	-0.01
22		Pentachlorophenol	0.400	0.591	-47.7#	151	-0.01
23		Phenanthrene	0.400	0.415	-3.7	93	-0.01
24		Anthracene	0.400	0.434	-8.5	100	-0.01
25	SURR	Fluoranthene-d10	0.400	0.438	-9.5	98	-0.01
26		Fluoranthene	0.400	0.434	-8.5	98	-0.01
27	I	Chrysene-d12	0.400	0.400	0.0	92	-0.02
28		Pyrene	0.400	0.420	-5.0	100	-0.01
29	S	Terphenyl-d14	0.400	0.427	-6.7	102	-0.01
30		Benzo(a)anthracene	0.400	0.420	-5.0	101	-0.01
31		Chrysene	0.400	0.424	-6.0	100	-0.01
32		Bis(2-ethylhexyl)phthalate	0.400	0.434	-8.5	112	-0.02
33		Indeno(1,2,3-cd)pyrene	0.400	0.395	1.3	93	-0.03
34	I	Perylene-d12	0.400	0.400	0.0	90	-0.02
35		Benzo(b)fluoranthene	0.400	0.406	-1.5	97	-0.02
36		Benzo(k)fluoranthene	0.400	0.410	-2.5	96	-0.02
37	C	Benzo(a)pyrene	0.400	0.412	-3.0	98	-0.02
38		Dibenzo(a,h)anthracene	0.400	0.392	2.0	93	-0.03
39		Benzo(g,h,i)perylene	0.400	0.397	0.8	93	-0.03

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

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