Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100219\
Data File : BN007983.D
Acq On : 03 Oct 2019 01:40
Operator : JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_N
LabSampleId : SSTDCCC0.4

Quant Time: Oct 03 07:28:43 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	99	-0.02
2		1,4-Dioxane	0.400	0.427	-6.7	125	-0.02
3		n-Nitrosodimethylamine	0.400	0.541	-35.3#	88	0.00
4	S	2-Fluorophenol	0.400	0.376	6.0	107	0.00
5	S	Phenol-d6	0.400	0.417	-4.2	114	0.00
6		bis(2-Chloroethyl)ether	0.400	0.513	-28.2#	144	-0.02
7	I	Naphthalene-d8	0.400	0.400	0.0	105	-0.01
8	S	Nitrobenzene-d5	0.400	0.424	-6.0	115	-0.01
9		Naphthalene	0.400	0.425	-6.2	112	-0.01
10		Hexachlorobutadiene	0.400	0.438	-9.5	115	-0.03
11	SURR	2-Methylnaphthalene-d10	0.400	0.436	-9.0	119	-0.02
12		2-Methylnaphthalene	0.400	0.435	-8.7	118	-0.02
13	I	Acenaphthene-d10	0.400	0.400	0.0	104	-0.01
14	S	2,4,6-Tribromophenol	0.400	0.394	1.5	122	-0.01
15	S	2-Fluorobiphenyl	0.400	0.391	2.3	109	-0.01
16		Acenaphthylene	0.400	0.397	0.8	115	-0.02
17		Acenaphthene	0.400	0.378	5.5	107	-0.02
18		Fluorene	0.400	0.379	5.3	107	-0.01
19	I	Phenanthrene-d10	0.400	0.400	0.0	94	-0.01
20		4-Bromophenyl-phenylether	0.400	0.424	-6.0	102	-0.02
21		Hexachlorobenzene	0.400	0.447	-11.7	107	-0.02
22		Pentachlorophenol	0.400	0.472	-18.0	128	-0.01
23		Phenanthrene	0.400	0.418	-4.5	100	-0.01
24		Anthracene	0.400	0.431	-7.7	105	-0.01
25	SURR	Fluoranthene-d10	0.400	0.419	-4.7	100	-0.01
26		Fluoranthene	0.400	0.414	-3.5	99	-0.02
27	I	Chrysene-d12	0.400	0.400	0.0	92	-0.02
28		Pyrene	0.400	0.422	-5.5	101	-0.02
29	S	Terphenyl-d14	0.400	0.426	-6.5	102	-0.02
30		Benzo(a)anthracene	0.400	0.409	-2.2	98	-0.02
31		Chrysene	0.400	0.419	-4.7	99	-0.02
32		Bis(2-ethylhexyl)phthalate	0.400	0.410	-2.5	105	-0.03
33		Indeno(1,2,3-cd)pyrene	0.400	0.382	4.5	90	-0.04
34	I	Perylene-d12	0.400	0.400	0.0	87	-0.03
35		Benzo(b)fluoranthene	0.400	0.419	-4.7	97	-0.03
36		Benzo(k)fluoranthene	0.400	0.403	-0.8	92	-0.03
37	C	Benzo(a)pyrene	0.400	0.411	-2.7	95	-0.03
38		Dibenzo(a,h)anthracene	0.400	0.393	1.8	90	-0.04
39		Benzo(g,h,i)perylene	0.400	0.397	0.8	90	-0.04

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

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