Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN103118\
Data File : BN003368.D
Acq On : 31 Oct 2018 15:29
Operator : MJ/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_N
LabSampleId : SSTDCCC0.4

Quant Time: Oct 31 16:34:47 2018
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN102618.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Oct 29 16:15:49 2018
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	98	-0.02
2		1,4-Dioxane	0.400	0.401	-0.3	105	0.00
3		n-Nitrosodimethylamine	0.400	0.457	-14.2	117	-0.02
4	S	2-Fluorophenol	0.400	0.381	4.8	97	-0.02
5	S	Phenol-d6	0.400	0.373	6.8	95	-0.02
6		bis(2-Chloroethyl)ether	0.400	0.403	-0.8	99	-0.02
7	I	Naphthalene-d8	0.400	0.400	0.0	99	-0.03
8	S	Nitrobenzene-d5	0.400	0.498	-24.5	135	-0.01
9		Naphthalene	0.400	0.390	2.5	99	-0.03
10		Hexachlorobutadiene	0.400	0.387	3.3	98	-0.01
11	SURR	2-Methylnaphthalene-d10	0.400	0.389	2.8	100	-0.02
12		2-Methylnaphthalene	0.400	0.388	3.0	99	-0.01
13	I	Acenaphthene-d10	0.400	0.400	0.0	101	-0.01
14	S	2,4,6-Tribromophenol	0.400	0.397	0.8	118	-0.01
15	S	2-Fluorobiphenyl	0.400	0.374	6.5	100	-0.02
16		Acenaphthylene	0.400	0.358	10.5	96	-0.01
17		Acenaphthene	0.400	0.385	3.8	100	-0.01
18		Fluorene	0.400	0.374	6.5	97	-0.01
19	I	Phenanthrene-d10	0.400	0.400	0.0	100	-0.01
20		4-Bromophenyl-phenylether	0.400	0.385	3.8	101	-0.02
21		Hexachlorobenzene	0.400	0.400	0.0	103	-0.01
22		Pentachlorophenol	0.400	0.505	-26.2#	137	-0.01
23		Phenanthrene	0.400	0.384	4.0	101	-0.01
24		Anthracene	0.400	0.363	9.3	97	-0.01
25	SURR	Fluoranthene-d10	0.400	0.373	6.8	98	0.00
26		Fluoranthene	0.400	0.375	6.3	99	0.00
27	I	Chrysene-d12	0.400	0.400	0.0	100	0.00
28		Pyrene	0.400	0.374	6.5	100	0.00
29	S	Terphenyl-d14	0.400	0.362	9.5	99	-0.01
30		Benzo(a)anthracene	0.400	0.361	9.8	101	-0.01
31		Chrysene	0.400	0.366	8.5	97	-0.01
32		Bis(2-ethylhexyl)phthalate	0.400	0.385	3.8	110	-0.02
33		Indeno(1,2,3-cd)pyrene	0.400	0.347	13.3	92	-0.02
34	I	Perylene-d12	0.400	0.400	0.0	95	0.00
35		Benzo(b)fluoranthene	0.400	0.400	0.0	103	-0.01
36		Benzo(k)fluoranthene	0.400	0.377	5.8	96	-0.01
37	C	Benzo(a)pyrene	0.400	0.385	3.8	97	-0.01
38		Dibenzo(a,h)anthracene	0.400	0.359	10.3	94	-0.02
39		Benzo(g,h,i)perylene	0.400	0.350	12.5	92	-0.02

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

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