Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA E\Data\BE122719\
Data File : BE100987.D
Acq On : 27 Dec 2019 9:58
Operator : JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Dec 27 23:56:57 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE122419.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Dec 27 17:39:05 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	103	0.00
2		1,4-Dioxane	0.400	0.370	7.5	102	0.00
3		n-Nitrosodimethylamine	0.400	0.622	-55.5#	201	-0.02
4	S	2-Fluorophenol	0.400	0.421	-5.2	120	0.00
5	S	Phenol-d6	0.400	0.420	-5.0	102	-0.04
6		bis(2-Chloroethyl)ether	0.400	0.346	13.5	106	0.00
7	I	Naphthalene-d8	0.400	0.400	0.0	108	0.00
8	S	Nitrobenzene-d5	0.400	0.451	-12.7	115	-0.01
9		Naphthalene	0.400	0.412	-3.0	110	0.00
10		Hexachlorobutadiene	0.400	0.415	-3.7	109	0.00
11	SURR	2-Methylnaphthalene-d10	0.400	0.434	-8.5	113	0.00
12		2-Methylnaphthalene	0.400	0.372	7.0	98	0.00
13	I	Acenaphthene-d10	0.400	0.400	0.0	105	-0.01
14	S	2,4,6-Tribromophenol	0.400	0.452	-13.0	118	-0.01
15	S	2-Fluorobiphenyl	0.400	0.435	-8.7	112	0.00
16		Acenaphthylene	0.400	0.363	9.3	102	0.00
17		Acenaphthene	0.400	0.417	-4.2	108	0.00
18		Fluorene	0.400	0.416	-4.0	112	0.00
19	I	Phenanthrene-d10	0.400	0.400	0.0	104	-0.01
20		4-Bromophenyl-phenylether	0.400	0.382	4.5	113	0.00
21		Hexachlorobenzene	0.400	0.390	2.5	105	0.00
22		Pentachlorophenol	0.400	0.383	4.3	92	-0.01
23		Phenanthrene	0.400	0.398	0.5	116	0.00
24		Anthracene	0.400	0.377	5.8	108	0.00
25	SURR	Fluoranthene-d10	0.400	0.399	0.3	112	0.00
26		Fluoranthene	0.400	0.388	3.0	112	0.00
27	I	Chrysene-d12	0.400	0.400	0.0	101	0.00
28		Pyrene	0.400	0.450	-12.5	111	0.00
29	S	Terphenyl-d14	0.400	0.452	-13.0	116	0.00
30		Benzo(a)anthracene	0.400	0.422	-5.5	122	0.00
31		Chrysene	0.400	0.435	-8.7	99	0.00
32		Bis(2-ethylhexyl)phthalate	0.400	0.393	1.8	101	-0.01
33		Indeno(1,2,3-cd)pyrene	0.400	0.414	-3.5	109	-0.01
34	I	Perylene-d12	0.400	0.400	0.0	92	0.00
35		Benzo(b)fluoranthene	0.400	0.481	-20.2	113	0.00
36		Benzo(k)fluoranthene	0.400	0.422	-5.5	96	0.00
37	C	Benzo(a)pyrene	0.400	0.429	-7.2	100	0.00
38		Dibenzo(a,h)anthracene	0.400	0.426	-6.5	118	-0.01
39		Benzo(g,h,i)perylene	0.400	0.445	-11.2	102	-0.01

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

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