Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE011520\
Data File : BE101060.D
Acq On : 15 Jan 2020 16:23
Operator : JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Jan 16 04:19:09 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE010820.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jan 08 18:36:25 2020
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	95	-0.01
2		1,4-Dioxane	0.400	0.437	-9.2	103	0.00
3		n-Nitrosodimethylamine	0.400	0.458	-14.5	127	-0.01
4	S	2-Fluorophenol	0.400	0.499	-24.7	120	-0.01
5	S	Phenol-d6	0.400	0.465	-16.3	125	-0.01
6		bis(2-Chloroethyl)ether	0.400	0.377	5.8	98	-0.01
7	I	Naphthalene-d8	0.400	0.400	0.0	93	0.00
8	S	Nitrobenzene-d5	0.400	0.449	-12.2	117	0.00
9		Naphthalene	0.400	0.410	-2.5	94	0.00
10		Hexachlorobutadiene	0.400	0.420	-5.0	94	-0.01
11	SURR	2-Methylnaphthalene-d10	0.400	0.389	2.8	90	-0.02
12		2-Methylnaphthalene	0.400	0.381	4.8	90	-0.02
13	I	Acenaphthene-d10	0.400	0.400	0.0	85	0.00
14	S	2,4,6-Tribromophenol	0.400	0.421	-5.2	103	-0.01
15	S	2-Fluorobiphenyl	0.400	0.410	-2.5	86	-0.02
16		Acenaphthylene	0.400	0.425	-6.2	100	-0.02
17		Acenaphthene	0.400	0.409	-2.2	89	-0.02
18		Fluorene	0.400	0.387	3.3	84	-0.01
19	I	Phenanthrene-d10	0.400	0.400	0.0	79	0.00
20		4-Bromophenyl-phenylether	0.400	0.408	-2.0	85	-0.01
21		Hexachlorobenzene	0.400	0.424	-6.0	84	-0.01
22		Pentachlorophenol	0.400	0.543	-35.8#	108	-0.01
23		Phenanthrene	0.400	0.382	4.5	77	0.00
24		Anthracene	0.400	0.417	-4.2	87	0.00
25	SURR	Fluoranthene-d10	0.400	0.390	2.5	80	-0.01
26		Fluoranthene	0.400	0.380	5.0	81	-0.01
27	I	Chrysene-d12	0.400	0.400	0.0	86	-0.01
28		Pyrene	0.400	0.393	1.8	83	0.00
29	S	Terphenyl-d14	0.400	0.378	5.5	80	0.00
30		Benzo(a)anthracene	0.400	0.407	-1.7	92	0.00
31		Chrysene	0.400	0.430	-7.5	90	-0.01
32		Bis(2-ethylhexyl)phthalate	0.400	0.563	-40.7#	140	0.00
33		Indeno(1,2,3-cd)pyrene	0.400	0.546	-36.5#	121	-0.02
34	I	Perylene-d12	0.400	0.400	0.0	116	-0.01
35		Benzo(b)fluoranthene	0.400	0.341	14.8	106	0.00
36		Benzo(k)fluoranthene	0.400	0.336	16.0	97	0.00
37	C	Benzo(a)pyrene	0.400	0.406	-1.5	121	0.00
38		Dibenzo(a,h)anthracene	0.400	0.397	0.8	124	-0.02
39		Benzo(g,h,i)perylene	0.400	0.402	-0.5	113	-0.01

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

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