Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE070219\
Data File : BE100304.D
Acq On : 2 Jul 2019 16:14
Operator : JU/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Jul 03 02:13:02 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\8270-SIM-BE062819.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 28 14:06:53 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	96	0.00
2		1,4-Dioxane	0.400	0.393	1.8	91	0.00
3		n-Nitrosodimethylamine	0.400	0.362	9.5	105	0.00
4	S	2-Fluorophenol	0.400	0.430	-7.5	102	0.00
5	S	Phenol-d6	0.400	0.375	6.3	90	0.00
6		bis(2-Chloroethyl)ether	0.400	0.397	0.8	86	-0.01
7	I	Naphthalene-d8	0.400	0.400	0.0	88	-0.01
8	S	Nitrobenzene-d5	0.400	0.402	-0.5	90	0.00
9		Naphthalene	0.400	0.412	-3.0	89	0.00
10		Hexachlorobutadiene	0.400	0.431	-7.7	92	0.00
11	SURR	2-Methylnaphthalene-d10	0.400	0.408	-2.0	88	0.00
12		2-Methylnaphthalene	0.400	0.381	4.8	88	0.00
13	I	Acenaphthene-d10	0.400	0.400	0.0	92	0.00
14	S	2,4,6-Tribromophenol	0.400	0.426	-6.5	100	0.01
15	S	2-Fluorobiphenyl	0.400	0.402	-0.5	90	0.00
16		Acenaphthylene	0.400	0.418	-4.5	93	0.00
17		Acenaphthene	0.400	0.411	-2.7	90	0.00
18		Fluorene	0.400	0.418	-4.5	93	0.01
19	I	Phenanthrene-d10	0.400	0.400	0.0	89	0.01
20		4-Bromophenyl-phenylether	0.400	0.432	-8.0	92	0.00
21		Hexachlorobenzene	0.400	0.436	-9.0	93	0.01
22		Pentachlorophenol	0.400	0.027	93.3#	7	0.04
23		Phenanthrene	0.400	0.422	-5.5	93	0.00
24		Anthracene	0.400	0.433	-8.2	97	0.00
25	SURR	Fluoranthene-d10	0.400	0.423	-5.7	93	0.00
26		Fluoranthene	0.400	0.424	-6.0	92	0.00
27	I	Chrysene-d12	0.400	0.400	0.0	81	0.00
28		Pyrene	0.400	0.468	-17.0	93	0.00
29	S	Terphenyl-d14	0.400	0.459	-14.8	93	0.00
30		Benzo(a)anthracene	0.400	0.420	-5.0	84	0.01
31		Chrysene	0.400	0.444	-11.0	88	0.00
32		Bis(2-ethylhexyl)phthalate	0.400	0.405	-1.3	80	0.00
33		Indeno(1,2,3-cd)pyrene	0.400	0.434	-8.5	82	0.02
34	I	Perylene-d12	0.400	0.400	0.0	81	0.01
35		Benzo(b)fluoranthene	0.400	0.411	-2.7	82	0.00
36		Benzo(k)fluoranthene	0.400	0.412	-3.0	80	0.01
37	C	Benzo(a)pyrene	0.400	0.437	-9.2	85	0.01
38		Dibenzo(a,h)anthracene	0.400	0.423	-5.7	85	0.01
39		Benzo(g,h,i)perylene	0.400	0.423	-5.7	82	0.02

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

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