Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE102819\
Data File : BE100692.D
Acq On : 29 Oct 2019 2:26
Operator : JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Oct 29 11:55:12 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE101019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Oct 23 16:49:33 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	57	0.00
2		1,4-Dioxane	0.400	0.499	-24.7	72	0.00
3		n-Nitrosodimethylamine	0.400	0.434	-8.5	68	0.00
4	S	2-Fluorophenol	0.400	0.451	-12.7	68	0.00
5	S	Phenol-d6	0.400	0.469	-17.2	71	0.00
6		bis(2-Chloroethyl)ether	0.400	0.475	-18.7	73	0.00
7	I	Naphthalene-d8	0.400	0.400	0.0	64	0.00
8	S	Nitrobenzene-d5	0.400	0.555	-38.8#	100	0.00
9		Naphthalene	0.400	0.445	-11.2	73	0.00
10		Hexachlorobutadiene	0.400	0.452	-13.0	75	0.00
11	SURR	2-Methylnaphthalene-d10	0.400	0.439	-9.7	73	0.00
12		2-Methylnaphthalene	0.400	0.436	-9.0	72	0.00
13	I	Acenaphthene-d10	0.400	0.400	0.0	66	0.00
14	S	2,4,6-Tribromophenol	0.400	0.545	-36.3#	101	0.00
15	S	2-Fluorobiphenyl	0.400	0.423	-5.7	71	0.00
16		Acenaphthylene	0.400	0.413	-3.2	70	-0.02
17		Acenaphthene	0.400	0.437	-9.2	74	0.00
18		Fluorene	0.400	0.442	-10.5	75	0.00
19	I	Phenanthrene-d10	0.400	0.400	0.0	69	-0.01
20		4-Bromophenyl-phenylether	0.400	0.436	-9.0	78	0.00
21		Hexachlorobenzene	0.400	0.427	-6.7	78	-0.01
22		Pentachlorophenol	0.400	0.356	11.0	69	0.00
23		Phenanthrene	0.400	0.445	-11.2	77	-0.01
24		Anthracene	0.400	0.417	-4.2	75	0.00
25	SURR	Fluoranthene-d10	0.400	0.437	-9.2	79	0.00
26		Fluoranthene	0.400	0.443	-10.7	77	0.00
27	I	Chrysene-d12	0.400	0.400	0.0	61	0.00
28		Pyrene	0.400	0.493	-23.2	75	0.00
29	S	Terphenyl-d14	0.400	0.468	-17.0	75	0.00
30		Benzo(a)anthracene	0.400	0.446	-11.5	68	0.00
31		Chrysene	0.400	0.442	-10.5	68	0.00
32		Bis(2-ethylhexyl)phthalate	0.400	0.454	-13.5	69	-0.01
33		Indeno(1,2,3-cd)pyrene	0.400	0.423	-5.7	66	-0.01
34	I	Perylene-d12	0.400	0.400	0.0	56	0.00
35		Benzo(b)fluoranthene	0.400	0.438	-9.5	66	0.00
36		Benzo(k)fluoranthene	0.400	0.444	-11.0	66	0.00
37	C	Benzo(a)pyrene	0.400	0.429	-7.2	65	-0.01
38		Dibenzo(a,h)anthracene	0.400	0.435	-8.7	67	-0.02
39		Benzo(g,h,i)perylene	0.400	0.437	-9.2	66	-0.01

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

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