Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE110819\
Data File : BE100711.D
Acq On : 8 Nov 2019 10:49
Operator : JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Nov 08 12:40:17 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	41	0.01
2		1,4-Dioxane	0.400	0.322	19.5	33	0.00
3		n-Nitrosodimethylamine	0.400	0.356	11.0	39	0.00
4	S	2-Fluorophenol	0.400	0.442	-10.5	47	0.01
5	S	Phenol-d6	0.400	0.486	-21.5	52	0.02
6		bis(2-Chloroethyl)ether	0.400	0.384	4.0	42	0.00
7	I	Naphthalene-d8	0.400	0.400	0.0	46	0.01
8	S	Nitrobenzene-d5	0.400	0.583	-45.7#	77	0.01
9		Naphthalene	0.400	0.385	3.8	46	0.01
10		Hexachlorobutadiene	0.400	0.401	-0.3	48	0.00
11	SURR	2-Methylnaphthalene-d10	0.400	0.403	-0.8	49	0.01
12		2-Methylnaphthalene	0.400	0.397	0.8	48	0.01
13	I	Acenaphthene-d10	0.400	0.400	0.0	50	0.01
14	S	2,4,6-Tribromophenol	0.400	0.174	56.5#	24	0.03
15	S	2-Fluorobiphenyl	0.400	0.408	-2.0	52	0.01
16		Acenaphthylene	0.400	0.380	5.0	49	0.00
17		Acenaphthene	0.400	0.367	8.3	47	0.01
18		Fluorene	0.400	0.377	5.8	48	0.01
19	I	Phenanthrene-d10	0.400	0.400	0.0	48	0.00
20		4-Bromophenyl-phenylether	0.400	0.403	-0.8	50	0.00
21		Hexachlorobenzene	0.400	0.403	-0.8	51	0.00
22		Pentachlorophenol	0.400	0.004	99.0#	1	0.00
23		Phenanthrene	0.400	0.397	0.8	47	0.00
24		Anthracene	0.400	0.395	1.3	49	0.01
25	SURR	Fluoranthene-d10	0.400	0.377	5.8	47	0.01
26		Fluoranthene	0.400	0.385	3.8	46	0.01
27	I	Chrysene-d12	0.400	0.400	0.0	41	0.01
28		Pyrene	0.400	0.445	-11.2	46	0.00
29	S	Terphenyl-d14	0.400	0.474	-18.5	52	0.00
30		Benzo(a)anthracene	0.400	0.399	0.3	42	0.01
31		Chrysene	0.400	0.385	3.8	40	0.01
32		Bis(2-ethylhexyl)phthalate	0.400	0.484	-21.0	50	0.00
33		Indeno(1,2,3-cd)pyrene	0.400	0.379	5.3	41	0.04
34	I	Perylene-d12	0.400	0.400	0.0	38	0.02
35		Benzo(b)fluoranthene	0.400	0.385	3.8	39	0.02
36		Benzo(k)fluoranthene	0.400	0.379	5.3	38	0.01
37	C	Benzo(a)pyrene	0.400	0.402	-0.5	41	0.02
38		Dibenzo(a,h)anthracene	0.400	0.394	1.5	41	0.04
39		Benzo(g,h,i)perylene	0.400	0.384	4.0	39	0.04

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

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