Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE102919\
Data File : BE100694.D
Acq On : 29 Oct 2019 14:26
Operator : JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Oct 29 16:45:00 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	67	0.00
2		1,4-Dioxane	0.400	0.438	-9.5	74	0.00
3		n-Nitrosodimethylamine	0.400	0.411	-2.7	75	0.00
4	S	2-Fluorophenol	0.400	0.417	-4.2	74	0.00
5	S	Phenol-d6	0.400	0.440	-10.0	78	0.00
6		bis(2-Chloroethyl)ether	0.400	0.455	-13.7	81	0.00
7	I	Naphthalene-d8	0.400	0.400	0.0	72	0.00
8	S	Nitrobenzene-d5	0.400	0.517	-29.3#	106	0.00
9		Naphthalene	0.400	0.444	-11.0	82	0.00
10		Hexachlorobutadiene	0.400	0.456	-14.0	85	0.00
11	SURR	2-Methylnaphthalene-d10	0.400	0.447	-11.7	84	0.00
12		2-Methylnaphthalene	0.400	0.451	-12.7	84	0.00
13	I	Acenaphthene-d10	0.400	0.400	0.0	73	0.00
14	S	2,4,6-Tribromophenol	0.400	0.504	-26.0#	105	0.00
15	S	2-Fluorobiphenyl	0.400	0.435	-8.7	82	0.00
16		Acenaphthylene	0.400	0.403	-0.8	77	-0.01
17		Acenaphthene	0.400	0.431	-7.7	82	0.00
18		Fluorene	0.400	0.454	-13.5	86	0.00
19	I	Phenanthrene-d10	0.400	0.400	0.0	78	-0.01
20		4-Bromophenyl-phenylether	0.400	0.434	-8.5	88	0.00
21		Hexachlorobenzene	0.400	0.443	-10.7	91	-0.01
22		Pentachlorophenol	0.400	0.330	17.5	73	0.00
23		Phenanthrene	0.400	0.450	-12.5	88	-0.01
24		Anthracene	0.400	0.415	-3.7	84	0.00
25	SURR	Fluoranthene-d10	0.400	0.435	-8.7	89	0.00
26		Fluoranthene	0.400	0.445	-11.2	87	0.00
27	I	Chrysene-d12	0.400	0.400	0.0	69	0.00
28		Pyrene	0.400	0.485	-21.2	85	0.00
29	S	Terphenyl-d14	0.400	0.469	-17.2	86	0.00
30		Benzo(a)anthracene	0.400	0.435	-8.7	76	0.00
31		Chrysene	0.400	0.444	-11.0	78	0.00
32		Bis(2-ethylhexyl)phthalate	0.400	0.347	13.3	60	-0.01
33		Indeno(1,2,3-cd)pyrene	0.400	0.425	-6.2	76	-0.01
34	I	Perylene-d12	0.400	0.400	0.0	66	-0.01
35		Benzo(b)fluoranthene	0.400	0.414	-3.5	73	0.00
36		Benzo(k)fluoranthene	0.400	0.446	-11.5	78	-0.01
37	C	Benzo(a)pyrene	0.400	0.406	-1.5	72	-0.01
38		Dibenzo(a,h)anthracene	0.400	0.415	-3.7	75	-0.01
39		Benzo(g,h,i)perylene	0.400	0.430	-7.5	76	-0.02

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

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