Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE070819\
Data File : BE100414.D
Acq On : 8 Jul 2019 16:29
Operator : JU/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Jul 08 19:05:04 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\8270-SIM-BE062819.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jul 08 15:19:44 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	72	-0.02
2		1,4-Dioxane	0.400	0.429	-7.2	74	-0.05
3		n-Nitrosodimethylamine	0.400	0.569	-42.2#	124	-0.02
4	S	2-Fluorophenol	0.400	0.389	2.8	69	-0.02
5	S	Phenol-d6	0.400	0.366	8.5	66	-0.02
6		bis(2-Chloroethyl)ether	0.400	0.307	23.3	50	-0.04
7	I	Naphthalene-d8	0.400	0.400	0.0	70	-0.03
8	S	Nitrobenzene-d5	0.400	0.350	12.5	63	-0.03
9		Naphthalene	0.400	0.436	-9.0	75	-0.01
10		Hexachlorobutadiene	0.400	0.538	-34.5#	92	-0.03
11	SURR	2-Methylnaphthalene-d10	0.400	0.433	-8.2	75	-0.01
12		2-Methylnaphthalene	0.400	0.399	0.3	73	-0.01
13	I	Acenaphthene-d10	0.400	0.400	0.0	70	-0.01
14	S	2,4,6-Tribromophenol	0.400	0.469	-17.2	83	-0.03
15	S	2-Fluorobiphenyl	0.400	0.459	-14.8	78	-0.01
16		Acenaphthylene	0.400	0.448	-12.0	75	-0.01
17		Acenaphthene	0.400	0.441	-10.2	73	-0.03
18		Fluorene	0.400	0.466	-16.5	79	-0.01
19	I	Phenanthrene-d10	0.400	0.400	0.0	75	-0.01
20		4-Bromophenyl-phenylether	0.400	0.440	-10.0	79	-0.03
21		Hexachlorobenzene	0.400	0.496	-24.0	89	-0.01
22		Pentachlorophenol	0.400	0.397	0.8	84	-0.04
23		Phenanthrene	0.400	0.421	-5.2	79	-0.03
24		Anthracene	0.400	0.397	0.8	76	-0.03
25	SURR	Fluoranthene-d10	0.400	0.418	-4.5	78	-0.02
26		Fluoranthene	0.400	0.416	-4.0	76	-0.02
27	I	Chrysene-d12	0.400	0.400	0.0	66	-0.01
28		Pyrene	0.400	0.478	-19.5	78	-0.02
29	S	Terphenyl-d14	0.400	0.451	-12.7	74	-0.02
30		Benzo(a)anthracene	0.400	0.430	-7.5	70	-0.01
31		Chrysene	0.400	0.422	-5.5	68	-0.01
32		Bis(2-ethylhexyl)phthalate	0.400	0.331	17.3	55	-0.01
33		Indeno(1,2,3-cd)pyrene	0.400	0.445	-11.2	68	-0.03
34	I	Perylene-d12	0.400	0.400	0.0	64	-0.02
35		Benzo(b)fluoranthene	0.400	0.403	-0.8	64	-0.02
36		Benzo(k)fluoranthene	0.400	0.494	-23.5	76	-0.02
37	C	Benzo(a)pyrene	0.400	0.421	-5.2	65	-0.02
38		Dibenzo(a,h)anthracene	0.400	0.405	-1.3	65	-0.04
39		Benzo(g,h,i)perylene	0.400	0.465	-16.3	72	-0.04

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

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