Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE060319\
Data File : BE099784.D
Acq On : 3 Jun 2019 13:50
Operator : JU/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Jun 04 03:46:50 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE052019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon May 20 19:29:40 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	75	0.00
2		1,4-Dioxane	0.400	0.391	2.3	76	-0.02
3		n-Nitrosodimethylamine	0.400	0.355	11.3	75	-0.02
4	S	2-Fluorophenol	0.400	0.372	7.0	79	0.00
5	S	Phenol-d6	0.400	0.383	4.3	86	0.00
6		bis(2-Chloroethyl)ether	0.400	0.370	7.5	79	0.00
7	I	Naphthalene-d8	0.400	0.400	0.0	74	0.00
8	S	Nitrobenzene-d5	0.400	0.388	3.0	84	-0.01
9		Naphthalene	0.400	0.405	-1.3	84	-0.01
10		Hexachlorobutadiene	0.400	0.403	-0.8	84	0.00
11	SURR	2-Methylnaphthalene-d10	0.400	0.407	-1.7	85	-0.01
12		2-Methylnaphthalene	0.400	0.407	-1.7	87	-0.01
13	I	Acenaphthene-d10	0.400	0.400	0.0	76	0.00
14	S	2,4,6-Tribromophenol	0.400	0.424	-6.0	92	0.00
15	S	2-Fluorobiphenyl	0.400	0.398	0.5	85	0.00
16		Acenaphthylene	0.400	0.420	-5.0	93	-0.01
17		Acenaphthene	0.400	0.406	-1.5	87	0.00
18		Fluorene	0.400	0.416	-4.0	90	-0.01
19	I	Phenanthrene-d10	0.400	0.400	0.0	80	0.00
20		4-Bromophenyl-phenylether	0.400	0.389	2.8	90	-0.01
21		Hexachlorobenzene	0.400	0.416	-4.0	95	-0.01
22		Pentachlorophenol	0.400	0.355	11.3	64	0.00
23		Phenanthrene	0.400	0.400	0.0	91	0.00
24		Anthracene	0.400	0.379	5.3	91	0.00
25	SURR	Fluoranthene-d10	0.400	0.407	-1.7	94	0.00
26		Fluoranthene	0.400	0.407	-1.7	94	0.00
27	I	Chrysene-d12	0.400	0.400	0.0	89	0.00
28		Pyrene	0.400	0.420	-5.0	95	-0.01
29	S	Terphenyl-d14	0.400	0.404	-1.0	93	-0.01
30		Benzo(a)anthracene	0.400	0.418	-4.5	97	-0.01
31		Chrysene	0.400	0.391	2.3	88	0.00
32		Bis(2-ethylhexyl)phthalate	0.400	0.356	11.0	88	-0.01
33		Indeno(1,2,3-cd)pyrene	0.400	0.404	-1.0	95	-0.02
34	I	Perylene-d12	0.400	0.400	0.0	83	-0.02
35		Benzo(b)fluoranthene	0.400	0.406	-1.5	93	0.00
36		Benzo(k)fluoranthene	0.400	0.402	-0.5	93	-0.01
37	C	Benzo(a)pyrene	0.400	0.408	-2.0	95	-0.01
38		Dibenzo(a,h)anthracene	0.400	0.397	0.8	95	-0.02
39		Benzo(g,h,i)perylene	0.400	0.405	-1.3	95	-0.01

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

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