Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE060419\
Data File : BE099792.D
Acq On : 4 Jun 2019 9:04
Operator : JU/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Jun 04 11:53:45 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	68	0.00
2		1,4-Dioxane	0.400	0.359	10.3	63	-0.02
3		n-Nitrosodimethylamine	0.400	0.319	20.3	61	-0.02
4	S	2-Fluorophenol	0.400	0.359	10.3	69	0.00
5	S	Phenol-d6	0.400	0.345	13.8	70	0.00
6		bis(2-Chloroethyl)ether	0.400	0.363	9.3	70	0.00
7	I	Naphthalene-d8	0.400	0.400	0.0	70	0.00
8	S	Nitrobenzene-d5	0.400	0.338	15.5	69	-0.01
9		Naphthalene	0.400	0.359	10.3	71	-0.01
10		Hexachlorobutadiene	0.400	0.365	8.8	72	0.00
11	SURR	2-Methylnaphthalene-d10	0.400	0.357	10.8	71	-0.01
12		2-Methylnaphthalene	0.400	0.340	15.0	69	0.00
13	I	Acenaphthene-d10	0.400	0.400	0.0	71	-0.01
14	S	2,4,6-Tribromophenol	0.400	0.384	4.0	75	-0.01
15	S	2-Fluorobiphenyl	0.400	0.355	11.3	71	0.00
16		Acenaphthylene	0.400	0.374	6.5	78	-0.01
17		Acenaphthene	0.400	0.350	12.5	71	0.00
18		Fluorene	0.400	0.356	11.0	73	-0.01
19	I	Phenanthrene-d10	0.400	0.400	0.0	79	0.00
20		4-Bromophenyl-phenylether	0.400	0.319	20.3	72	-0.01
21		Hexachlorobenzene	0.400	0.361	9.8	80	-0.01
22		Pentachlorophenol	0.400	0.366	8.5	67	0.01
23		Phenanthrene	0.400	0.349	12.8	78	0.00
24		Anthracene	0.400	0.331	17.3	78	0.00
25	SURR	Fluoranthene-d10	0.400	0.383	4.3	87	0.00
26		Fluoranthene	0.400	0.389	2.8	88	0.00
27	I	Chrysene-d12	0.400	0.400	0.0	95	-0.01
28		Pyrene	0.400	0.371	7.3	90	-0.01
29	S	Terphenyl-d14	0.400	0.351	12.3	87	-0.01
30		Benzo(a)anthracene	0.400	0.372	7.0	93	-0.01
31		Chrysene	0.400	0.394	1.5	95	0.00
32		Bis(2-ethylhexyl)phthalate	0.400	0.362	9.5	96	-0.01
33		Indeno(1,2,3-cd)pyrene	0.400	0.390	2.5	99	-0.02
34	I	Perylene-d12	0.400	0.400	0.0	96	-0.02
35		Benzo(b)fluoranthene	0.400	0.368	8.0	98	-0.01
36		Benzo(k)fluoranthene	0.400	0.360	10.0	96	-0.02
37	C	Benzo(a)pyrene	0.400	0.362	9.5	98	-0.01
38		Dibenzo(a,h)anthracene	0.400	0.352	12.0	97	-0.02
39		Benzo(g,h,i)perylene	0.400	0.358	10.5	97	-0.03

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

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