Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE012020\
Data File : BE101096.D
Acq On : 20 Jan 2020 10:05
Operator : JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Jan 21 02:45:44 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE010820.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jan 08 18:36:25 2020
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	58	-0.01
2		1,4-Dioxane	0.400	0.403	-0.8	59	0.00
3		n-Nitrosodimethylamine	0.400	0.438	-9.5	73	-0.01
4	S	2-Fluorophenol	0.400	0.500	-25.0	73	-0.01
5	S	Phenol-d6	0.400	0.392	2.0	64	-0.01
6		bis(2-Chloroethyl)ether	0.400	0.417	-4.2	66	-0.01
7	I	Naphthalene-d8	0.400	0.400	0.0	61	0.00
8	S	Nitrobenzene-d5	0.400	0.363	9.3	61	0.00
9		Naphthalene	0.400	0.387	3.3	57	0.00
10		Hexachlorobutadiene	0.400	0.403	-0.8	59	-0.01
11	SURR	2-Methylnaphthalene-d10	0.400	0.410	-2.5	62	-0.02
12		2-Methylnaphthalene	0.400	0.390	2.5	60	0.00
13	I	Acenaphthene-d10	0.400	0.400	0.0	72	0.00
14	S	2,4,6-Tribromophenol	0.400	0.352	12.0	73	0.00
15	S	2-Fluorobiphenyl	0.400	0.360	10.0	64	-0.02
16		Acenaphthylene	0.400	0.365	8.8	72	0.00
17		Acenaphthene	0.400	0.372	7.0	69	-0.02
18		Fluorene	0.400	0.395	1.3	72	-0.01
19	I	Phenanthrene-d10	0.400	0.400	0.0	82	0.00
20		4-Bromophenyl-phenylether	0.400	0.384	4.0	83	-0.01
21		Hexachlorobenzene	0.400	0.360	10.0	73	-0.01
22		Pentachlorophenol	0.400	0.402	-0.5	58	0.00
23		Phenanthrene	0.400	0.371	7.3	77	0.00
24		Anthracene	0.400	0.363	9.3	78	0.00
25	SURR	Fluoranthene-d10	0.400	0.420	-5.0	89	-0.01
26		Fluoranthene	0.400	0.415	-3.7	90	-0.01
27	I	Chrysene-d12	0.400	0.400	0.0	124	-0.01
28		Pyrene	0.400	0.302	24.5	91	0.00
29	S	Terphenyl-d14	0.400	0.321	19.8	98	0.00
30		Benzo(a)anthracene	0.400	0.374	6.5	120	0.00
31		Chrysene	0.400	0.394	1.5	118	-0.01
32		Bis(2-ethylhexyl)phthalate	0.400	0.441	-10.2	157	0.00
33		Indeno(1,2,3-cd)pyrene	0.400	0.319	20.3	101	-0.01
34	I	Perylene-d12	0.400	0.400	0.0	132	0.00
35		Benzo(b)fluoranthene	0.400	0.378	5.5	134	-0.01
36		Benzo(k)fluoranthene	0.400	0.356	11.0	118	0.00
37	C	Benzo(a)pyrene	0.400	0.382	4.5	131	0.00
38		Dibenzo(a,h)anthracene	0.400	0.305	23.8	109	0.00
39		Benzo(g,h,i)perylene	0.400	0.287	28.3#	92	-0.01

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

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