Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN091319\
Data File : BN007627.D
Acq On : 13 Sep 2019 15:15
Operator : HP/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_N
LabSampleId : SSTDCCC0.4

Quant Time: Sep 14 02:59:54 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN091219.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Sep 12 14:27:51 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	0.385	3.8	68	0.00
3		n-Nitrosodimethylamine	0.400	0.581	-45.2#	69	0.00
4	S	2-Fluorophenol	0.400	0.335	16.3	62	0.00
5	S	Phenol-d6	0.400	0.356	11.0	62	0.00
6		bis(2-Chloroethyl)ether	0.400	0.363	9.3	61	0.00
7	I	Naphthalene-d8	0.400	0.400	0.0	67	-0.01
8	S	Nitrobenzene-d5	0.400	0.383	4.3	62	0.00
9		Naphthalene	0.400	0.426	-6.5	67	-0.01
10		Hexachlorobutadiene	0.400	0.418	-4.5	66	-0.01
11	SURR	2-Methylnaphthalene-d10	0.400	0.415	-3.7	66	0.00
12		2-Methylnaphthalene	0.400	0.419	-4.7	66	0.00
13	I	Acenaphthene-d10	0.400	0.400	0.0	66	0.00
14	S	2,4,6-Tribromophenol	0.400	0.296	26.0#	57	0.00
15	S	2-Fluorobiphenyl	0.400	0.404	-1.0	67	0.00
16		Acenaphthylene	0.400	0.363	9.3	61	0.00
17		Acenaphthene	0.400	0.388	3.0	65	0.00
18		Fluorene	0.400	0.388	3.0	65	-0.01
19	I	Phenanthrene-d10	0.400	0.400	0.0	64	0.00
20		4-Bromophenyl-phenylether	0.400	0.429	-7.2	65	0.00
21		Hexachlorobenzene	0.400	0.433	-8.2	65	-0.01
22		Pentachlorophenol	0.400	0.345	13.8	60	-0.01
23		Phenanthrene	0.400	0.428	-7.0	64	0.00
24		Anthracene	0.400	0.399	0.3	61	0.00
25	SURR	Fluoranthene-d10	0.400	0.408	-2.0	62	0.00
26		Fluoranthene	0.400	0.412	-3.0	62	0.00
27	I	Chrysene-d12	0.400	0.400	0.0	62	-0.01
28		Pyrene	0.400	0.427	-6.7	63	0.00
29	S	Terphenyl-d14	0.400	0.423	-5.7	63	0.00
30		Benzo(a)anthracene	0.400	0.380	5.0	57	0.00
31		Chrysene	0.400	0.432	-8.0	64	0.00
32		Bis(2-ethylhexyl)phthalate	0.400	0.313	21.8	51	-0.01
33		Indeno(1,2,3-cd)pyrene	0.400	0.440	-10.0	65	-0.01
34	I	Perylene-d12	0.400	0.400	0.0	62	0.00
35		Benzo(b)fluoranthene	0.400	0.399	0.3	61	0.00
36		Benzo(k)fluoranthene	0.400	0.415	-3.7	65	0.00
37	C	Benzo(a)pyrene	0.400	0.380	5.0	60	-0.01
38		Dibenzo(a,h)anthracene	0.400	0.416	-4.0	64	-0.01
39		Benzo(g,h,i)perylene	0.400	0.402	-0.5	62	-0.01

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

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