Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100519\
Data File : BN008042.D
Acq On : 06 Oct 2019 02:37
Operator : JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_N
LabSampleId : SSTDCCC0.4

Quant Time: Oct 07 07:33:48 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN091819.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Oct 03 18:39:49 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	64	-0.02
2		1,4-Dioxane	0.400	0.468	-17.0	89	-0.03
3		n-Nitrosodimethylamine	0.400	0.639	-59.7#	67	-0.03
4	S	2-Fluorophenol	0.400	0.368	8.0	68	-0.02
5	S	Phenol-d6	0.400	0.363	9.3	64	0.00
6		bis(2-Chloroethyl)ether	0.400	0.490	-22.5	89	-0.03
7	I	Naphthalene-d8	0.400	0.400	0.0	65	-0.03
8	S	Nitrobenzene-d5	0.400	0.401	-0.3	67	-0.01
9		Naphthalene	0.400	0.408	-2.0	66	-0.03
10		Hexachlorobutadiene	0.400	0.459	-14.8	74	-0.04
11	SURR	2-Methylnaphthalene-d10	0.400	0.418	-4.5	70	-0.02
12		2-Methylnaphthalene	0.400	0.417	-4.2	69	-0.03
13	I	Acenaphthene-d10	0.400	0.400	0.0	64	-0.02
14	S	2,4,6-Tribromophenol	0.400	0.390	2.5	74	-0.01
15	S	2-Fluorobiphenyl	0.400	0.368	8.0	63	-0.02
16		Acenaphthylene	0.400	0.397	0.8	70	-0.02
17		Acenaphthene	0.400	0.363	9.3	63	-0.02
18		Fluorene	0.400	0.367	8.3	64	-0.01
19	I	Phenanthrene-d10	0.400	0.400	0.0	57	-0.02
20		4-Bromophenyl-phenylether	0.400	0.404	-1.0	59	-0.02
21		Hexachlorobenzene	0.400	0.443	-10.7	64	-0.02
22		Pentachlorophenol	0.400	0.409	-2.2	67	0.00
23		Phenanthrene	0.400	0.391	2.3	57	-0.02
24		Anthracene	0.400	0.406	-1.5	61	-0.01
25	SURR	Fluoranthene-d10	0.400	0.400	0.0	58	-0.02
26		Fluoranthene	0.400	0.404	-1.0	59	-0.02
27	I	Chrysene-d12	0.400	0.400	0.0	53	-0.02
28		Pyrene	0.400	0.442	-10.5	61	-0.02
29	S	Terphenyl-d14	0.400	0.424	-6.0	58	-0.02
30		Benzo(a)anthracene	0.400	0.401	-0.3	55	-0.02
31		Chrysene	0.400	0.408	-2.0	55	-0.02
32		Bis(2-ethylhexyl)phthalate	0.400	0.489	-22.2	72	-0.03
33		Indeno(1,2,3-cd)pyrene	0.400	0.396	1.0	54	-0.03
34	I	Perylene-d12	0.400	0.400	0.0	52	-0.03
35		Benzo(b)fluoranthene	0.400	0.402	-0.5	55	-0.03
36		Benzo(k)fluoranthene	0.400	0.382	4.5	52	-0.03
37	C	Benzo(a)pyrene	0.400	0.399	0.3	55	-0.03
38		Dibenzo(a,h)anthracene	0.400	0.394	1.5	54	-0.04
39		Benzo(g,h,i)perylene	0.400	0.406	-1.5	55	-0.03

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

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