Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN100319\
Data File : BN007985.D
Acq On : 03 Oct 2019 18:00
Operator : JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_N
LabSampleId : SSTDCCC0.4

Quant Time: Oct 04 08:29:32 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	66	-0.02
2		1,4-Dioxane	0.400	0.420	-5.0	82	-0.02
3		n-Nitrosodimethylamine	0.400	0.641	-60.2#	69	-0.02
4	S	2-Fluorophenol	0.400	0.369	7.8	70	-0.02
5	S	Phenol-d6	0.400	0.379	5.3	69	0.00
6		bis(2-Chloroethyl)ether	0.400	0.432	-8.0	81	-0.02
7	I	Naphthalene-d8	0.400	0.400	0.0	68	-0.01
8	S	Nitrobenzene-d5	0.400	0.394	1.5	69	-0.01
9		Naphthalene	0.400	0.416	-4.0	71	-0.03
10		Hexachlorobutadiene	0.400	0.439	-9.7	75	-0.03
11	SURR	2-Methylnaphthalene-d10	0.400	0.426	-6.5	75	-0.02
12		2-Methylnaphthalene	0.400	0.429	-7.2	75	-0.02
13	I	Acenaphthene-d10	0.400	0.400	0.0	71	-0.02
14	S	2,4,6-Tribromophenol	0.400	0.391	2.3	83	-0.01
15	S	2-Fluorobiphenyl	0.400	0.380	5.0	72	-0.01
16		Acenaphthylene	0.400	0.375	6.3	74	-0.02
17		Acenaphthene	0.400	0.363	9.3	70	-0.02
18		Fluorene	0.400	0.381	4.8	74	-0.01
19	I	Phenanthrene-d10	0.400	0.400	0.0	70	-0.02
20		4-Bromophenyl-phenylether	0.400	0.411	-2.7	73	-0.02
21		Hexachlorobenzene	0.400	0.427	-6.7	75	-0.02
22		Pentachlorophenol	0.400	0.425	-6.2	85	-0.01
23		Phenanthrene	0.400	0.403	-0.8	71	-0.02
24		Anthracene	0.400	0.413	-3.2	75	-0.01
25	SURR	Fluoranthene-d10	0.400	0.426	-6.5	75	-0.02
26		Fluoranthene	0.400	0.421	-5.2	74	-0.02
27	I	Chrysene-d12	0.400	0.400	0.0	71	-0.02
28		Pyrene	0.400	0.414	-3.5	76	-0.02
29	S	Terphenyl-d14	0.400	0.421	-5.2	77	-0.02
30		Benzo(a)anthracene	0.400	0.403	-0.8	75	-0.02
31		Chrysene	0.400	0.412	-3.0	75	-0.02
32		Bis(2-ethylhexyl)phthalate	0.400	0.462	-15.5	92	-0.03
33		Indeno(1,2,3-cd)pyrene	0.400	0.388	3.0	71	-0.04
34	I	Perylene-d12	0.400	0.400	0.0	69	-0.03
35		Benzo(b)fluoranthene	0.400	0.409	-2.2	75	-0.03
36		Benzo(k)fluoranthene	0.400	0.400	0.0	72	-0.03
37	C	Benzo(a)pyrene	0.400	0.402	-0.5	73	-0.03
38		Dibenzo(a,h)anthracene	0.400	0.391	2.3	71	-0.04
39		Benzo(g,h,i)perylene	0.400	0.394	1.5	71	-0.04

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

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