Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN090319\
Data File : BN007534.D
Acq On : 03 Sep 2019 14:37
Operator : HP/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_N
LabSampleId : SSTDCCC0.4

Quant Time: Sep 03 16:50:07 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN082319.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Aug 29 12:18:45 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	197	-0.02
2		1,4-Dioxane	0.400	0.381	4.8	181	-0.02
3		n-Nitrosodimethylamine	0.400	0.309	22.8	147	-0.02
4	S	2-Fluorophenol	0.400	0.345	13.8	169	-0.02
5	S	Phenol-d6	0.400	0.359	10.3	185	-0.02
6		bis(2-Chloroethyl)ether	0.400	0.414	-3.5	206	-0.02
7	I	Naphthalene-d8	0.400	0.400	0.0	227	-0.03
8	S	Nitrobenzene-d5	0.400	0.350	12.5	207	-0.01
9		Naphthalene	0.400	0.394	1.5	218	-0.01
10		Hexachlorobutadiene	0.400	0.344	14.0	187	-0.01
11	SURR	2-Methylnaphthalene-d10	0.400	0.405	-1.3	226	-0.02
12		2-Methylnaphthalene	0.400	0.410	-2.5	225	-0.02
13	I	Acenaphthene-d10	0.400	0.400	0.0	249	-0.02
14	S	2,4,6-Tribromophenol	0.400	0.338	15.5	214	-0.02
15	S	2-Fluorobiphenyl	0.400	0.392	2.0	239	0.00
16		Acenaphthylene	0.400	0.367	8.3	226	-0.02
17		Acenaphthene	0.400	0.385	3.8	235	-0.02
18		Fluorene	0.400	0.407	-1.7	248	-0.02
19	I	Phenanthrene-d10	0.400	0.400	0.0	250	-0.02
20		4-Bromophenyl-phenylether	0.400	0.093	76.8#	67	0.00
21		Hexachlorobenzene	0.400	0.390	2.5	241	-0.02
22		Pentachlorophenol	0.400	0.357	10.8	238	-0.02
23		Phenanthrene	0.400	0.440	-10.0	275	-0.02
24		Anthracene	0.400	0.386	3.5	239	0.00
25	SURR	Fluoranthene-d10	0.400	0.372	7.0	225	-0.01
26		Fluoranthene	0.400	0.410	-2.5	250	-0.01
27	I	Chrysene-d12	0.400	0.400	0.0	223	0.00
28		Pyrene	0.400	0.469	-17.2	255	0.00
29	S	Terphenyl-d14	0.400	0.417	-4.2	225	-0.01
30		Benzo(a)anthracene	0.400	0.421	-5.2	230	-0.01
31		Chrysene	0.400	0.402	-0.5	215	-0.01
32		Bis(2-ethylhexyl)phthalate	0.400	0.341	14.8	183	-0.01
33		Indeno(1,2,3-cd)pyrene	0.400	0.291	27.3#	157	-0.02
34	I	Perylene-d12	0.400	0.400	0.0	181	-0.01
35		Benzo(b)fluoranthene	0.400	0.437	-9.2	193	-0.01
36		Benzo(k)fluoranthene	0.400	0.424	-6.0	184	0.00
37	C	Benzo(a)pyrene	0.400	0.400	0.0	175	-0.02
38		Dibenzo(a,h)anthracene	0.400	0.348	13.0	152	-0.02
39		Benzo(g,h,i)perylene	0.400	0.378	5.5	166	-0.02

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

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