Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN092719\
Data File : BN007882.D
Acq On : 27 Sep 2019 11:22
Operator : JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_N
LabSampleId : SSTDCCC0.4

Quant Time: Sep 27 15:51:33 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN091819.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Sep 18 18:24:40 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	136	0.00
2		1,4-Dioxane	0.400	0.412	-3.0	166	0.00
3		n-Nitrosodimethylamine	0.400	0.628	-57.0#	141	0.00
4	S	2-Fluorophenol	0.400	0.367	8.3	144	0.00
5	S	Phenol-d6	0.400	0.374	6.5	141	0.00
6		bis(2-Chloroethyl)ether	0.400	0.421	-5.2	163	-0.02
7	I	Naphthalene-d8	0.400	0.400	0.0	135	-0.01
8	S	Nitrobenzene-d5	0.400	0.421	-5.2	147	0.00
9		Naphthalene	0.400	0.426	-6.5	145	-0.01
10		Hexachlorobutadiene	0.400	0.443	-10.7	149	-0.01
11	SURR	2-Methylnaphthalene-d10	0.400	0.420	-5.0	147	-0.01
12		2-Methylnaphthalene	0.400	0.427	-6.7	148	-0.01
13	I	Acenaphthene-d10	0.400	0.400	0.0	131	-0.01
14	S	2,4,6-Tribromophenol	0.400	0.348	13.0	136	-0.01
15	S	2-Fluorobiphenyl	0.400	0.404	-1.0	142	-0.01
16		Acenaphthylene	0.400	0.388	3.0	141	-0.01
17		Acenaphthene	0.400	0.386	3.5	138	-0.01
18		Fluorene	0.400	0.387	3.3	138	-0.01
19	I	Phenanthrene-d10	0.400	0.400	0.0	118	-0.01
20		4-Bromophenyl-phenylether	0.400	0.447	-11.7	134	-0.01
21		Hexachlorobenzene	0.400	0.450	-12.5	134	-0.01
22		Pentachlorophenol	0.400	0.401	-0.3	136	-0.01
23		Phenanthrene	0.400	0.431	-7.7	129	-0.01
24		Anthracene	0.400	0.427	-6.7	130	-0.01
25	SURR	Fluoranthene-d10	0.400	0.401	-0.3	119	-0.01
26		Fluoranthene	0.400	0.407	-1.7	122	-0.01
27	I	Chrysene-d12	0.400	0.400	0.0	109	-0.01
28		Pyrene	0.400	0.433	-8.2	122	-0.01
29	S	Terphenyl-d14	0.400	0.439	-9.7	124	-0.01
30		Benzo(a)anthracene	0.400	0.412	-3.0	117	-0.01
31		Chrysene	0.400	0.417	-4.2	117	-0.01
32		Bis(2-ethylhexyl)phthalate	0.400	0.415	-3.7	126	-0.02
33		Indeno(1,2,3-cd)pyrene	0.400	0.399	0.3	111	-0.03
34	I	Perylene-d12	0.400	0.400	0.0	107	-0.02
35		Benzo(b)fluoranthene	0.400	0.409	-2.2	116	-0.01
36		Benzo(k)fluoranthene	0.400	0.417	-4.2	116	-0.02
37	C	Benzo(a)pyrene	0.400	0.410	-2.5	115	-0.02
38		Dibenzo(a,h)anthracene	0.400	0.400	0.0	112	-0.02
39		Benzo(g,h,i)perylene	0.400	0.397	0.8	111	-0.03

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

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