Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN090119\
Data File : BN007527.D
Acq On : 01 Sep 2019 16:14
Operator : HP/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_N
LabSampleId : SSTDCCC0.4

Quant Time: Sep 02 08:19:14 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	165	-0.02
2		1,4-Dioxane	0.400	0.399	0.3	158	0.00
3		n-Nitrosodimethylamine	0.400	0.315	21.3	125	0.00
4	S	2-Fluorophenol	0.400	0.313	21.8	129	-0.02
5	S	Phenol-d6	0.400	0.313	21.8	136	0.00
6		bis(2-Chloroethyl)ether	0.400	0.417	-4.2	173	0.00
7	I	Naphthalene-d8	0.400	0.400	0.0	182	-0.01
8	S	Nitrobenzene-d5	0.400	0.335	16.3	159	0.00
9		Naphthalene	0.400	0.394	1.5	174	-0.01
10		Hexachlorobutadiene	0.400	0.357	10.8	156	-0.01
11	SURR	2-Methylnaphthalene-d10	0.400	0.402	-0.5	180	-0.01
12		2-Methylnaphthalene	0.400	0.405	-1.3	179	-0.01
13	I	Acenaphthene-d10	0.400	0.400	0.0	206	-0.02
14	S	2,4,6-Tribromophenol	0.400	0.297	25.8#	156	-0.02
15	S	2-Fluorobiphenyl	0.400	0.385	3.8	194	0.00
16		Acenaphthylene	0.400	0.335	16.3	170	-0.02
17		Acenaphthene	0.400	0.385	3.8	194	0.00
18		Fluorene	0.400	0.398	0.5	200	-0.02
19	I	Phenanthrene-d10	0.400	0.400	0.0	213	0.00
20		4-Bromophenyl-phenylether	0.400	0.155	61.3#	96	0.00
21		Hexachlorobenzene	0.400	0.400	0.0	211	-0.02
22		Pentachlorophenol	0.400	0.228	43.0#	130	0.00
23		Phenanthrene	0.400	0.410	-2.5	218	-0.02
24		Anthracene	0.400	0.356	11.0	188	0.00
25	SURR	Fluoranthene-d10	0.400	0.380	5.0	196	-0.01
26		Fluoranthene	0.400	0.388	3.0	202	-0.01
27	I	Chrysene-d12	0.400	0.400	0.0	208	-0.01
28		Pyrene	0.400	0.395	1.3	201	-0.01
29	S	Terphenyl-d14	0.400	0.389	2.8	196	-0.01
30		Benzo(a)anthracene	0.400	0.360	10.0	184	-0.01
31		Chrysene	0.400	0.405	-1.3	202	-0.01
32		Bis(2-ethylhexyl)phthalate	0.400	0.267	33.3#	136	-0.01
33		Indeno(1,2,3-cd)pyrene	0.400	0.333	16.8	167	-0.03
34	I	Perylene-d12	0.400	0.400	0.0	189	-0.02
35		Benzo(b)fluoranthene	0.400	0.400	0.0	185	-0.02
36		Benzo(k)fluoranthene	0.400	0.436	-9.0	198	-0.02
37	C	Benzo(a)pyrene	0.400	0.375	6.3	172	-0.02
38		Dibenzo(a,h)anthracene	0.400	0.357	10.8	163	-0.03
39		Benzo(g,h,i)perylene	0.400	0.359	10.3	165	-0.03

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

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