Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN100419\
Data File : BN008021.D
Acq On : 05 Oct 2019 01:44
Operator : JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_N
LabSampleId : SSTDCCC0.4

Quant Time: Oct 05 04:31:03 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	88	-0.02
2		1,4-Dioxane	0.400	0.506	-26.5#	132	-0.03
3		n-Nitrosodimethylamine	0.400	0.645	-61.2#	93	-0.02
4	S	2-Fluorophenol	0.400	0.377	5.8	95	0.00
5	S	Phenol-d6	0.400	0.386	3.5	94	0.02
6		bis(2-Chloroethyl)ether	0.400	0.558	-39.5#	139	-0.03
7	I	Naphthalene-d8	0.400	0.400	0.0	88	-0.01
8	S	Nitrobenzene-d5	0.400	0.406	-1.5	92	-0.01
9		Naphthalene	0.400	0.411	-2.7	91	-0.01
10		Hexachlorobutadiene	0.400	0.458	-14.5	100	-0.03
11	SURR	2-Methylnaphthalene-d10	0.400	0.413	-3.2	94	-0.01
12		2-Methylnaphthalene	0.400	0.421	-5.2	95	-0.02
13	I	Acenaphthene-d10	0.400	0.400	0.0	81	-0.01
14	S	2,4,6-Tribromophenol	0.400	0.441	-10.2	107	-0.01
15	S	2-Fluorobiphenyl	0.400	0.392	2.0	86	-0.01
16		Acenaphthylene	0.400	0.584	-46.0#	132	-0.02
17		Acenaphthene	0.400	0.400	0.0	89	-0.02
18		Fluorene	0.400	0.446	-11.5	99	-0.01
19	I	Phenanthrene-d10	0.400	0.400	0.0	73	-0.01
20		4-Bromophenyl-phenylether	0.400	0.425	-6.2	79	-0.02
21		Hexachlorobenzene	0.400	0.429	-7.2	79	-0.01
22		Pentachlorophenol	0.400	0.512	-28.0#	108	0.00
23		Phenanthrene	0.400	0.422	-5.5	78	-0.01
24		Anthracene	0.400	0.445	-11.2	84	-0.01
25	SURR	Fluoranthene-d10	0.400	0.410	-2.5	76	0.00
26		Fluoranthene	0.400	0.428	-7.0	79	0.00
27	I	Chrysene-d12	0.400	0.400	0.0	67	-0.01
28		Pyrene	0.400	0.469	-17.2	81	-0.01
29	S	Terphenyl-d14	0.400	0.442	-10.5	76	-0.01
30		Benzo(a)anthracene	0.400	0.426	-6.5	74	-0.01
31		Chrysene	0.400	0.403	-0.8	69	-0.01
32		Bis(2-ethylhexyl)phthalate	0.400	0.554	-38.5#	103	-0.03
33		Indeno(1,2,3-cd)pyrene	0.400	0.393	1.8	67	-0.02
34	I	Perylene-d12	0.400	0.400	0.0	66	-0.02
35		Benzo(b)fluoranthene	0.400	0.399	0.3	70	-0.02
36		Benzo(k)fluoranthene	0.400	0.385	3.8	66	-0.02
37	C	Benzo(a)pyrene	0.400	0.413	-3.2	72	-0.02
38		Dibenzo(a,h)anthracene	0.400	0.390	2.5	68	-0.02
39		Benzo(g,h,i)perylene	0.400	0.398	0.5	69	-0.02

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

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