Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE062119\
Data File : BE100139.D
Acq On : 21 Jun 2019 16:51
Operator : JU/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Jun 22 02:03:37 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE061919.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jun 19 15:07:24 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	91	-0.01
2		1,4-Dioxane	0.400	0.425	-6.2	110	-0.01
3		n-Nitrosodimethylamine	0.400	0.594	-48.5#	349	-0.04
4	S	2-Fluorophenol	0.400	0.568	-42.0#	138	-0.01
5	S	Phenol-d6	0.400	0.576	-44.0#	131	-0.01
6		bis(2-Chloroethyl)ether	0.400	0.421	-5.2	104	-0.03
7	I	Naphthalene-d8	0.400	0.400	0.0	106	-0.01
8	S	Nitrobenzene-d5	0.400	0.404	-1.0	107	-0.01
9		Naphthalene	0.400	0.401	-0.3	107	-0.01
10		Hexachlorobutadiene	0.400	0.349	12.8	95	-0.01
11	SURR	2-Methylnaphthalene-d10	0.400	0.441	-10.2	122	-0.01
12		2-Methylnaphthalene	0.400	0.449	-12.2	125	-0.01
13	I	Acenaphthene-d10	0.400	0.400	0.0	136	-0.01
14	S	2,4,6-Tribromophenol	0.400	0.456	-14.0	174	-0.01
15	S	2-Fluorobiphenyl	0.400	0.390	2.5	122	-0.01
16		Acenaphthylene	0.400	0.419	-4.7	141	-0.01
17		Acenaphthene	0.400	0.418	-4.5	139	0.00
18		Fluorene	0.400	0.441	-10.2	147	-0.01
19	I	Phenanthrene-d10	0.400	0.400	0.0	148	-0.01
20		4-Bromophenyl-phenylether	0.400	0.402	-0.5	155	-0.01
21		Hexachlorobenzene	0.400	0.378	5.5	141	-0.01
22		Pentachlorophenol	0.400	0.507	-26.7#	218	-0.04
23		Phenanthrene	0.400	0.390	2.5	147	0.00
24		Anthracene	0.400	0.403	-0.8	161	0.00
25	SURR	Fluoranthene-d10	0.400	0.386	3.5	145	0.00
26		Fluoranthene	0.400	0.380	5.0	143	-0.01
27	I	Chrysene-d12	0.400	0.400	0.0	120	0.00
28		Pyrene	0.400	0.494	-23.5	149	0.00
29	S	Terphenyl-d14	0.400	0.471	-17.7	140	-0.01
30		Benzo(a)anthracene	0.400	0.421	-5.2	137	0.00
31		Chrysene	0.400	0.401	-0.3	116	-0.01
32		Bis(2-ethylhexyl)phthalate	0.400	0.720	-80.0#	249	0.00
33		Indeno(1,2,3-cd)pyrene	0.400	0.336	16.0	97	0.00
34	I	Perylene-d12	0.400	0.400	0.0	120	-0.01
35		Benzo(b)fluoranthene	0.400	0.373	6.8	113	0.00
36		Benzo(k)fluoranthene	0.400	0.357	10.8	111	-0.01
37	C	Benzo(a)pyrene	0.400	0.372	7.0	112	0.00
38		Dibenzo(a,h)anthracene	0.400	0.293	26.8#	91	0.00
39		Benzo(g,h,i)perylene	0.400	0.362	9.5	110	-0.02

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

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