Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE032919\
Data File : BE099265.D
Acq On : 29 Mar 2019 22:46
Operator : JU/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Mar 29 23:38:33 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\8270-SIM-BE032519.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Mar 26 07:02:06 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	89	-0.01
2		1,4-Dioxane	0.400	0.421	-5.2	106	0.00
3		n-Nitrosodimethylamine	0.400	0.431	-7.7	125	-0.01
4	S	2-Fluorophenol	0.400	0.468	-17.0	114	-0.01
5	S	Phenol-d6	0.400	0.471	-17.7	117	0.00
6		bis(2-Chloroethyl)ether	0.400	0.440	-10.0	111	-0.01
7	I	Naphthalene-d8	0.400	0.400	0.0	89	0.00
8	S	Nitrobenzene-d5	0.400	0.480	-20.0	131	-0.01
9		Naphthalene	0.400	0.451	-12.7	114	-0.01
10		Hexachlorobutadiene	0.400	0.484	-21.0	125	0.00
11	SURR	2-Methylnaphthalene-d10	0.400	0.457	-14.2	114	-0.01
12		2-Methylnaphthalene	0.400	0.451	-12.7	114	0.00
13	I	Acenaphthene-d10	0.400	0.400	0.0	92	-0.01
14	S	2,4,6-Tribromophenol	0.400	0.513	-28.2#	161	0.00
15	S	2-Fluorobiphenyl	0.400	0.471	-17.7	120	0.00
16		Acenaphthylene	0.400	0.441	-10.2	120	0.00
17		Acenaphthene	0.400	0.445	-11.2	117	-0.01
18		Fluorene	0.400	0.437	-9.2	116	-0.01
19	I	Phenanthrene-d10	0.400	0.400	0.0	92	-0.01
20		4-Bromophenyl-phenylether	0.400	0.432	-8.0	119	-0.01
21		Hexachlorobenzene	0.400	0.432	-8.0	119	-0.01
22		Pentachlorophenol	0.400	0.571	-42.7#	206	-0.01
23		Phenanthrene	0.400	0.421	-5.2	113	-0.01
24		Anthracene	0.400	0.423	-5.7	120	0.00
25	SURR	Fluoranthene-d10	0.400	0.466	-16.5	125	-0.01
26		Fluoranthene	0.400	0.436	-9.0	121	0.00
27	I	Chrysene-d12	0.400	0.400	0.0	98	-0.01
28		Pyrene	0.400	0.438	-9.5	124	0.00
29	S	Terphenyl-d14	0.400	0.448	-12.0	122	0.00
30		Benzo(a)anthracene	0.400	0.442	-10.5	130	0.00
31		Chrysene	0.400	0.429	-7.2	122	0.00
32		Bis(2-ethylhexyl)phthalate	0.400	0.520	-30.0#	181	-0.01
33		Indeno(1,2,3-cd)pyrene	0.400	0.369	7.8	114	-0.01
34	I	Perylene-d12	0.400	0.400	0.0	92	0.00
35		Benzo(b)fluoranthene	0.400	0.407	-1.7	114	0.00
36		Benzo(k)fluoranthene	0.400	0.455	-13.7	123	0.00
37	C	Benzo(a)pyrene	0.400	0.432	-8.0	120	0.00
38		Dibenzo(a,h)anthracene	0.400	0.384	4.0	114	-0.02
39		Benzo(g,h,i)perylene	0.400	0.342	14.5	99	-0.02

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

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