Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE032319\
Data File : BE099122.D
Acq On : 23 Mar 2019 22:01
Operator : JU/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Mar 23 22:43:59 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE032219.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat Mar 23 04:40:11 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	100	0.00
2		1,4-Dioxane	0.400	0.422	-5.5	121	-0.01
3		n-Nitrosodimethylamine	0.400	0.424	-6.0	115	-0.01
4	S	2-Fluorophenol	0.400	0.464	-16.0	130	0.00
5	S	Phenol-d6	0.400	0.483	-20.7	138	0.00
6		bis(2-Chloroethyl)ether	0.400	0.480	-20.0	134	-0.01
7	I	Naphthalene-d8	0.400	0.400	0.0	112	-0.01
8	S	Nitrobenzene-d5	0.400	0.776	-94.0#	272	0.00
9		Naphthalene	0.400	0.441	-10.2	140	0.00
10		Hexachlorobutadiene	0.400	0.468	-17.0	146	0.00
11	SURR	2-Methylnaphthalene-d10	0.400	0.459	-14.8	147	-0.02
12		2-Methylnaphthalene	0.400	0.464	-16.0	150	-0.02
13	I	Acenaphthene-d10	0.400	0.400	0.0	131	0.00
14	S	2,4,6-Tribromophenol	0.400	0.773	-93.3#	326	0.00
15	S	2-Fluorobiphenyl	0.400	0.445	-11.2	159	-0.02
16		Acenaphthylene	0.400	0.423	-5.7	164	-0.02
17		Acenaphthene	0.400	0.433	-8.2	166	0.00
18		Fluorene	0.400	0.426	-6.5	162	0.00
19	I	Phenanthrene-d10	0.400	0.400	0.0	126	0.00
20		4-Bromophenyl-phenylether	0.400	0.462	-15.5	172	0.00
21		Hexachlorobenzene	0.400	0.488	-22.0	183	0.00
22		Pentachlorophenol	0.400	0.654	-63.5#	257	-0.01
23		Phenanthrene	0.400	0.424	-6.0	159	0.00
24		Anthracene	0.400	0.411	-2.7	156	0.00
25	SURR	Fluoranthene-d10	0.400	0.400	0.0	146	0.00
26		Fluoranthene	0.400	0.389	2.8	147	0.00
27	I	Chrysene-d12	0.400	0.400	0.0	117	-0.01
28		Pyrene	0.400	0.423	-5.7	147	0.00
29	S	Terphenyl-d14	0.400	0.473	-18.2	158	0.00
30		Benzo(a)anthracene	0.400	0.408	-2.0	141	0.00
31		Chrysene	0.400	0.428	-7.0	149	0.00
32		Bis(2-ethylhexyl)phthalate	0.400	0.430	-7.5	178	-0.01
33		Indeno(1,2,3-cd)pyrene	0.400	0.435	-8.7	152	-0.01
34	I	Perylene-d12	0.400	0.400	0.0	125	0.00
35		Benzo(b)fluoranthene	0.400	0.401	-0.3	151	0.00
36		Benzo(k)fluoranthene	0.400	0.405	-1.3	151	0.00
37	C	Benzo(a)pyrene	0.400	0.401	-0.3	152	0.00
38		Dibenzo(a,h)anthracene	0.400	0.410	-2.5	155	-0.02
39		Benzo(g,h,i)perylene	0.400	0.406	-1.5	152	-0.02

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

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