Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE052419\
Data File : BE099736.D
Acq On : 24 May 2019 11:23
Operator : JU/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: May 24 21:37:55 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE052019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon May 20 19:29:40 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	60	-0.01
2		1,4-Dioxane	0.400	0.432	-8.0	68	-0.02
3		n-Nitrosodimethylamine	0.400	0.466	-16.5	80	-0.03
4	S	2-Fluorophenol	0.400	0.453	-13.2	78	-0.01
5	S	Phenol-d6	0.400	0.452	-13.0	82	-0.01
6		bis(2-Chloroethyl)ether	0.400	0.433	-8.2	75	-0.01
7	I	Naphthalene-d8	0.400	0.400	0.0	63	0.00
8	S	Nitrobenzene-d5	0.400	0.462	-15.5	85	-0.01
9		Naphthalene	0.400	0.411	-2.7	73	-0.01
10		Hexachlorobutadiene	0.400	0.410	-2.5	73	0.00
11	SURR	2-Methylnaphthalene-d10	0.400	0.416	-4.0	74	-0.02
12		2-Methylnaphthalene	0.400	0.426	-6.5	77	-0.02
13	I	Acenaphthene-d10	0.400	0.400	0.0	65	0.00
14	S	2,4,6-Tribromophenol	0.400	0.541	-35.3#	112	-0.01
15	S	2-Fluorobiphenyl	0.400	0.412	-3.0	75	0.00
16		Acenaphthylene	0.400	0.423	-5.7	80	-0.02
17		Acenaphthene	0.400	0.416	-4.0	76	0.00
18		Fluorene	0.400	0.426	-6.5	79	-0.01
19	I	Phenanthrene-d10	0.400	0.400	0.0	68	0.00
20		4-Bromophenyl-phenylether	0.400	0.400	0.0	78	-0.01
21		Hexachlorobenzene	0.400	0.400	0.0	77	-0.01
22		Pentachlorophenol	0.400	0.720	-80.0#	176	-0.03
23		Phenanthrene	0.400	0.423	-5.7	81	0.00
24		Anthracene	0.400	0.420	-5.0	85	0.00
25	SURR	Fluoranthene-d10	0.400	0.425	-6.2	83	0.00
26		Fluoranthene	0.400	0.425	-6.2	82	0.00
27	I	Chrysene-d12	0.400	0.400	0.0	83	0.00
28		Pyrene	0.400	0.407	-1.7	86	-0.01
29	S	Terphenyl-d14	0.400	0.401	-0.3	86	-0.01
30		Benzo(a)anthracene	0.400	0.425	-6.2	92	-0.01
31		Chrysene	0.400	0.411	-2.7	86	0.00
32		Bis(2-ethylhexyl)phthalate	0.400	0.539	-34.8#	124	-0.01
33		Indeno(1,2,3-cd)pyrene	0.400	0.406	-1.5	90	-0.02
34	I	Perylene-d12	0.400	0.400	0.0	76	-0.02
35		Benzo(b)fluoranthene	0.400	0.409	-2.2	86	-0.01
36		Benzo(k)fluoranthene	0.400	0.398	0.5	84	-0.02
37	C	Benzo(a)pyrene	0.400	0.412	-3.0	89	-0.01
38		Dibenzo(a,h)anthracene	0.400	0.401	-0.3	88	-0.02
39		Benzo(g,h,i)perylene	0.400	0.402	-0.5	86	-0.02

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

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