Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE070119\
Data File : BE100291.D
Acq On : 2 Jul 2019 7:45
Operator : JU/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Jul 02 08:28:53 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE062819.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 28 14:06:53 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	340	0.00
2		1,4-Dioxane	0.400	0.249	37.8#	205	0.01
3		n-Nitrosodimethylamine	0.400	0.506	-26.5#	522	-0.01
4	S	2-Fluorophenol	0.400	0.397	0.8	336	0.01
5	S	Phenol-d6	0.400	0.524	-31.0#	447	-0.01
6		bis(2-Chloroethyl)ether	0.400	0.587	-46.7#	451	-0.01
7	I	Naphthalene-d8	0.400	0.400	0.0	456	-0.01
8	S	Nitrobenzene-d5	0.400	0.449	-12.2	521	-0.01
9		Naphthalene	0.400	0.406	-1.5	455	0.00
10		Hexachlorobutadiene	0.400	0.401	-0.3	443	0.00
11	SURR	2-Methylnaphthalene-d10	0.400	0.426	-6.5	477	-0.01
12		2-Methylnaphthalene	0.400	0.404	-1.0	482	0.00
13	I	Acenaphthene-d10	0.400	0.400	0.0	459	0.00
14	S	2,4,6-Tribromophenol	0.400	0.450	-12.5	524	-0.01
15	S	2-Fluorobiphenyl	0.400	0.411	-2.7	457	-0.01
16		Acenaphthylene	0.400	0.403	-0.8	445	0.00
17		Acenaphthene	0.400	0.415	-3.7	450	0.00
18		Fluorene	0.400	0.395	1.3	437	0.00
19	I	Phenanthrene-d10	0.400	0.400	0.0	399	0.01
20		4-Bromophenyl-phenylether	0.400	0.437	-9.2	418	-0.01
21		Hexachlorobenzene	0.400	0.426	-6.5	407	0.01
22		Pentachlorophenol	0.400	0.499	-24.7	557	-0.03
23		Phenanthrene	0.400	0.422	-5.5	418	0.00
24		Anthracene	0.400	0.461	-15.3	465	0.00
25	SURR	Fluoranthene-d10	0.400	0.395	1.3	389	0.00
26		Fluoranthene	0.400	0.389	2.8	377	0.00
27	I	Chrysene-d12	0.400	0.400	0.0	325	0.00
28		Pyrene	0.400	0.480	-20.0	382	0.00
29	S	Terphenyl-d14	0.400	0.477	-19.2	383	0.00
30		Benzo(a)anthracene	0.400	0.445	-11.2	354	0.01
31		Chrysene	0.400	0.435	-8.7	344	0.00
32		Bis(2-ethylhexyl)phthalate	0.400	0.434	-8.5	338	0.00
33		Indeno(1,2,3-cd)pyrene	0.400	0.369	7.8	277	0.03
34	I	Perylene-d12	0.400	0.400	0.0	307	0.02
35		Benzo(b)fluoranthene	0.400	0.428	-7.0	324	0.00
36		Benzo(k)fluoranthene	0.400	0.398	0.5	294	0.01
37	C	Benzo(a)pyrene	0.400	0.421	-5.2	310	0.01
38		Dibenzo(a,h)anthracene	0.400	0.397	0.8	302	0.02
39		Benzo(g,h,i)perylene	0.400	0.350	12.5	258	0.02

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

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