Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE061219\
Data File : BE099946.D
Acq On : 12 Jun 2019 8:32
Operator : JU/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Jun 13 04:24:54 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE061019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jun 10 14:57:20 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	127	-0.01
2		1,4-Dioxane	0.400	0.422	-5.5	142	0.00
3		n-Nitrosodimethylamine	0.400	0.255	36.3#	80	0.03
4	S	2-Fluorophenol	0.400	0.336	16.0	107	0.00
5	S	Phenol-d6	0.400	0.365	8.8	124	0.00
6		bis(2-Chloroethyl)ether	0.400	0.383	4.3	128	0.00
7	I	Naphthalene-d8	0.400	0.400	0.0	127	-0.01
8	S	Nitrobenzene-d5	0.400	0.373	6.8	126	-0.01
9		Naphthalene	0.400	0.410	-2.5	132	-0.03
10		Hexachlorobutadiene	0.400	0.450	-12.5	140	-0.03
11	SURR	2-Methylnaphthalene-d10	0.400	0.414	-3.5	134	-0.03
12		2-Methylnaphthalene	0.400	0.404	-1.0	130	-0.01
13	I	Acenaphthene-d10	0.400	0.400	0.0	123	-0.03
14	S	2,4,6-Tribromophenol	0.400	0.254	36.5#	87	-0.03
15	S	2-Fluorobiphenyl	0.400	0.411	-2.7	134	-0.03
16		Acenaphthylene	0.400	0.405	-1.3	129	-0.03
17		Acenaphthene	0.400	0.411	-2.7	132	-0.03
18		Fluorene	0.400	0.406	-1.5	132	-0.03
19	I	Phenanthrene-d10	0.400	0.400	0.0	119	-0.03
20		4-Bromophenyl-phenylether	0.400	0.415	-3.7	134	-0.03
21		Hexachlorobenzene	0.400	0.422	-5.5	132	-0.04
22		Pentachlorophenol	0.400	0.348	13.0	25	0.03
23		Phenanthrene	0.400	0.388	3.0	123	-0.03
24		Anthracene	0.400	0.350	12.5	110	-0.03
25	SURR	Fluoranthene-d10	0.400	0.348	13.0	106	-0.04
26		Fluoranthene	0.400	0.352	12.0	106	-0.04
27	I	Chrysene-d12	0.400	0.400	0.0	87	-0.04
28		Pyrene	0.400	0.501	-25.2#	103	-0.03
29	S	Terphenyl-d14	0.400	0.448	-12.0	94	-0.03
30		Benzo(a)anthracene	0.400	0.403	-0.8	85	-0.04
31		Chrysene	0.400	0.452	-13.0	91	-0.04
32		Bis(2-ethylhexyl)phthalate	0.400	0.325	18.8	70	-0.04
33		Indeno(1,2,3-cd)pyrene	0.400	0.385	3.8	80	-0.10
34	I	Perylene-d12	0.400	0.400	0.0	85	-0.07
35		Benzo(b)fluoranthene	0.400	0.366	8.5	81	-0.06
36		Benzo(k)fluoranthene	0.400	0.431	-7.7	92	-0.06
37	C	Benzo(a)pyrene	0.400	0.395	1.3	87	-0.07
38		Dibenzo(a,h)anthracene	0.400	0.322	19.5	72	-0.09
39		Benzo(g,h,i)perylene	0.400	0.385	3.8	84	-0.11

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

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