Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE050719\
Data File : BE099565.D
Acq On : 7 May 2019 14:30
Operator : JU/SJ
Sample : SSTDICV0.4
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : ICVBE050719

Quant Time: May 07 15:21:07 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE050719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue May 07 14:10:11 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	93	0.02
2		1,4-Dioxane	0.400	0.402	-0.5	113	0.00
3		n-Nitrosodimethylamine	0.400	0.373	6.8	96	0.00
4	S	2-Fluorophenol	0.400	0.406	-1.5	105	0.00
5	S	Phenol-d6	0.400	0.391	2.3	99	0.02
6		bis(2-Chloroethyl)ether	0.400	0.398	0.5	103	0.02
7	I	Naphthalene-d8	0.400	0.400	0.0	91	0.01
8	S	Nitrobenzene-d5	0.400	0.395	1.3	100	0.03
9		Naphthalene	0.400	0.413	-3.2	104	0.01
10		Hexachlorobutadiene	0.400	0.421	-5.2	106	0.01
11	SURR	2-Methylnaphthalene-d10	0.400	0.419	-4.7	106	0.03
12		2-Methylnaphthalene	0.400	0.407	-1.7	101	0.01
13	I	Acenaphthene-d10	0.400	0.400	0.0	89	0.01
14	S	2,4,6-Tribromophenol	0.400	0.375	6.3	98	0.01
15	S	2-Fluorobiphenyl	0.400	0.402	-0.5	103	0.01
16		Acenaphthylene	0.400	0.392	2.0	99	0.01
17		Acenaphthene	0.400	0.389	2.8	97	0.01
18		Fluorene	0.400	0.399	0.3	99	0.01
19	I	Phenanthrene-d10	0.400	0.400	0.0	91	0.01
20		4-Bromophenyl-phenylether	0.400	0.382	4.5	99	0.01
21		Hexachlorobenzene	0.400	0.399	0.3	101	0.00
22		Pentachlorophenol	0.400	0.348	13.0	99	0.01
23		Phenanthrene	0.400	0.407	-1.7	103	0.01
24		Anthracene	0.400	0.400	0.0	104	0.01
25	SURR	Fluoranthene-d10	0.400	0.398	0.5	102	0.00
26		Fluoranthene	0.400	0.398	0.5	101	0.00
27	I	Chrysene-d12	0.400	0.400	0.0	101	0.00
28		Pyrene	0.400	0.399	0.3	102	0.00
29	S	Terphenyl-d14	0.400	0.400	0.0	103	0.01
30		Benzo(a)anthracene	0.400	0.385	3.8	99	0.00
31		Chrysene	0.400	0.404	-1.0	102	0.01
32		Bis(2-ethylhexyl)phthalate	0.400	0.396	1.0	103	0.00
33		Indeno(1,2,3-cd)pyrene	0.400	0.390	2.5	100	0.01
34	I	Perylene-d12	0.400	0.400	0.0	91	0.01
35		Benzo(b)fluoranthene	0.400	0.407	-1.7	101	0.00
36		Benzo(k)fluoranthene	0.400	0.415	-3.7	104	0.00
37	C	Benzo(a)pyrene	0.400	0.404	-1.0	101	0.00
38		Dibenzo(a,h)anthracene	0.400	0.403	-0.8	101	0.02
39		Benzo(g,h,i)perylene	0.400	0.404	-1.0	100	0.01

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

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