Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE072219\
Data File : BE100484.D
Acq On : 22 Jul 2019 17:21
Operator : HP/JU
Sample : SSTDICV0.4
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : ICVBE072219

Quant Time: Jul 22 18:02:51 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE072219.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jul 22 17:11:23 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	99	0.00
2		1,4-Dioxane	0.400	0.340	15.0	94	0.00
3		n-Nitrosodimethylamine	0.400	0.335	16.3	83	0.00
4	S	2-Fluorophenol	0.400	0.408	-2.0	103	0.00
5	S	Phenol-d6	0.400	0.394	1.5	99	0.00
6		bis(2-Chloroethyl)ether	0.400	0.412	-3.0	103	0.00
7	I	Naphthalene-d8	0.400	0.400	0.0	98	0.00
8	S	Nitrobenzene-d5	0.400	0.393	1.8	103	0.00
9		Naphthalene	0.400	0.396	1.0	98	0.00
10		Hexachlorobutadiene	0.400	0.391	2.3	96	0.00
11	SURR	2-Methylnaphthalene-d10	0.400	0.384	4.0	97	0.00
12		2-Methylnaphthalene	0.400	0.388	3.0	98	0.00
13	I	Acenaphthene-d10	0.400	0.400	0.0	100	0.00
14	S	2,4,6-Tribromophenol	0.400	0.378	5.5	109	0.00
15	S	2-Fluorobiphenyl	0.400	0.378	5.5	101	0.00
16		Acenaphthylene	0.400	0.373	6.8	100	0.00
17		Acenaphthene	0.400	0.376	6.0	100	0.00
18		Fluorene	0.400	0.381	4.8	100	0.00
19	I	Phenanthrene-d10	0.400	0.400	0.0	103	0.00
20		4-Bromophenyl-phenylether	0.400	0.377	5.8	101	0.00
21		Hexachlorobenzene	0.400	0.373	6.8	99	-0.01
22		Pentachlorophenol	0.400	0.432	-8.0	107	0.00
23		Phenanthrene	0.400	0.390	2.5	103	-0.01
24		Anthracene	0.400	0.378	5.5	102	0.00
25	SURR	Fluoranthene-d10	0.400	0.377	5.8	100	0.00
26		Fluoranthene	0.400	0.379	5.3	99	0.00
27	I	Chrysene-d12	0.400	0.400	0.0	97	0.00
28		Pyrene	0.400	0.403	-0.8	100	0.00
29	S	Terphenyl-d14	0.400	0.400	0.0	100	0.00
30		Benzo(a)anthracene	0.400	0.397	0.8	98	0.00
31		Chrysene	0.400	0.397	0.8	98	0.00
32		Bis(2-ethylhexyl)phthalate	0.400	0.392	2.0	95	0.00
33		Indeno(1,2,3-cd)pyrene	0.400	0.394	1.5	96	0.00
34	I	Perylene-d12	0.400	0.400	0.0	98	0.00
35		Benzo(b)fluoranthene	0.400	0.379	5.3	96	0.00
36		Benzo(k)fluoranthene	0.400	0.404	-1.0	102	0.00
37	C	Benzo(a)pyrene	0.400	0.391	2.3	99	0.00
38		Dibenzo(a,h)anthracene	0.400	0.387	3.3	100	0.00
39		Benzo(g,h,i)perylene	0.400	0.390	2.5	99	-0.01

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

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