Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE060519\
Data File : BE099805.D
Acq On : 5 Jun 2019 14:53
Operator : JU/SJ
Sample : SSTCICV0.4
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : ICVBE060519

Quant Time: Jun 05 15:30:27 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE060519.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jun 05 14:35:58 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	140	0.01
2		1,4-Dioxane	0.400	0.434	-8.5	178	0.01
3		n-Nitrosodimethylamine	0.400	0.407	-1.7	175	0.02
4	S	2-Fluorophenol	0.400	0.448	-12.0	183	0.02
5	S	Phenol-d6	0.400	0.432	-8.0	170	0.02
6		bis(2-Chloroethyl)ether	0.400	0.423	-5.7	174	0.01
7	I	Naphthalene-d8	0.400	0.400	0.0	140	0.01
8	S	Nitrobenzene-d5	0.400	0.421	-5.2	167	0.02
9		Naphthalene	0.400	0.417	-4.2	162	0.02
10		Hexachlorobutadiene	0.400	0.430	-7.5	167	0.02
11	SURR	2-Methylnaphthalene-d10	0.400	0.403	-0.8	157	0.01
12		2-Methylnaphthalene	0.400	0.409	-2.2	163	0.01
13	I	Acenaphthene-d10	0.400	0.400	0.0	126	0.01
14	S	2,4,6-Tribromophenol	0.400	0.423	-5.7	141	0.00
15	S	2-Fluorobiphenyl	0.400	0.452	-13.0	166	0.01
16		Acenaphthylene	0.400	0.429	-7.2	154	0.01
17		Acenaphthene	0.400	0.426	-6.5	152	0.01
18		Fluorene	0.400	0.416	-4.0	144	0.01
19	I	Phenanthrene-d10	0.400	0.400	0.0	110	0.00
20		4-Bromophenyl-phenylether	0.400	0.413	-3.2	139	0.01
21		Hexachlorobenzene	0.400	0.410	-2.5	132	0.01
22		Pentachlorophenol	0.400	0.419	-4.7	99	0.01
23		Phenanthrene	0.400	0.410	-2.5	128	0.00
24		Anthracene	0.400	0.387	3.3	128	0.00
25	SURR	Fluoranthene-d10	0.400	0.352	12.0	112	0.00
26		Fluoranthene	0.400	0.359	10.3	112	0.00
27	I	Chrysene-d12	0.400	0.400	0.0	96	0.00
28		Pyrene	0.400	0.450	-12.5	109	0.00
29	S	Terphenyl-d14	0.400	0.451	-12.7	111	0.00
30		Benzo(a)anthracene	0.400	0.407	-1.7	102	0.01
31		Chrysene	0.400	0.413	-3.2	100	0.00
32		Bis(2-ethylhexyl)phthalate	0.400	0.386	3.5	96	0.00
33		Indeno(1,2,3-cd)pyrene	0.400	0.407	-1.7	99	0.02
34	I	Perylene-d12	0.400	0.400	0.0	87	0.01
35		Benzo(b)fluoranthene	0.400	0.415	-3.7	101	0.00
36		Benzo(k)fluoranthene	0.400	0.416	-4.0	101	0.00
37	C	Benzo(a)pyrene	0.400	0.414	-3.5	100	0.00
38		Dibenzo(a,h)anthracene	0.400	0.396	1.0	98	0.02
39		Benzo(g,h,i)perylene	0.400	0.400	0.0	99	0.02

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

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