Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE050919\
Data File : BE099589.D
Acq On : 9 May 2019 13:41
Operator : JU/SJ
Sample : SSTDICV0.4
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : ICVBE050919

Quant Time: May 09 16:06:11 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE050919.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu May 09 13:32:58 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	80	0.01
2		1,4-Dioxane	0.400	0.436	-9.0	103	0.00
3		n-Nitrosodimethylamine	0.400	0.386	3.5	93	0.00
4	S	2-Fluorophenol	0.400	0.404	-1.0	94	0.00
5	S	Phenol-d6	0.400	0.383	4.3	88	0.01
6		bis(2-Chloroethyl)ether	0.400	0.418	-4.5	96	0.01
7	I	Naphthalene-d8	0.400	0.400	0.0	80	0.01
8	S	Nitrobenzene-d5	0.400	0.386	3.5	87	0.01
9		Naphthalene	0.400	0.405	-1.3	91	0.00
10		Hexachlorobutadiene	0.400	0.398	0.5	88	0.01
11	SURR	2-Methylnaphthalene-d10	0.400	0.401	-0.3	88	0.01
12		2-Methylnaphthalene	0.400	0.406	-1.5	90	0.01
13	I	Acenaphthene-d10	0.400	0.400	0.0	78	0.00
14	S	2,4,6-Tribromophenol	0.400	0.409	-2.2	92	0.01
15	S	2-Fluorobiphenyl	0.400	0.410	-2.5	89	0.00
16		Acenaphthylene	0.400	0.397	0.8	87	0.01
17		Acenaphthene	0.400	0.402	-0.5	87	0.00
18		Fluorene	0.400	0.404	-1.0	89	0.01
19	I	Phenanthrene-d10	0.400	0.400	0.0	78	0.00
20		4-Bromophenyl-phenylether	0.400	0.386	3.5	87	0.01
21		Hexachlorobenzene	0.400	0.406	-1.5	88	0.01
22		Pentachlorophenol	0.400	0.467	-16.8	81	0.01
23		Phenanthrene	0.400	0.405	-1.3	89	0.00
24		Anthracene	0.400	0.392	2.0	89	0.00
25	SURR	Fluoranthene-d10	0.400	0.404	-1.0	89	0.00
26		Fluoranthene	0.400	0.408	-2.0	89	0.00
27	I	Chrysene-d12	0.400	0.400	0.0	85	0.00
28		Pyrene	0.400	0.408	-2.0	87	0.00
29	S	Terphenyl-d14	0.400	0.405	-1.3	87	0.00
30		Benzo(a)anthracene	0.400	0.401	-0.3	86	0.01
31		Chrysene	0.400	0.421	-5.2	90	0.00
32		Bis(2-ethylhexyl)phthalate	0.400	0.395	1.3	87	0.01
33		Indeno(1,2,3-cd)pyrene	0.400	0.420	-5.0	92	0.01
34	I	Perylene-d12	0.400	0.400	0.0	81	0.00
35		Benzo(b)fluoranthene	0.400	0.409	-2.2	91	0.00
36		Benzo(k)fluoranthene	0.400	0.394	1.5	87	0.00
37	C	Benzo(a)pyrene	0.400	0.400	0.0	90	0.00
38		Dibenzo(a,h)anthracene	0.400	0.402	-0.5	91	0.01
39		Benzo(g,h,i)perylene	0.400	0.408	-2.0	91	0.01

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

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