Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE020618\
Data File : BE095538.D
Acq On : 6 Feb 2018 20:27
Operator : SJ/JU
Sample : SSTDICV0.4
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : ICVBE020618

Quant Time: Feb 06 21:03:20 2018
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE020618.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Feb 06 20:27:34 2018
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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	100	0.00
2		1,4-Dioxane	0.662	0.658	0.6	99	0.00
3		n-Nitrosodimethylamine	0.833	0.818	1.8	97	0.00
4	S	2-Fluorophenol	1.207	1.179	2.3	96	0.00
5	S	Phenol-d6	1.671	1.590	4.8	94	0.00
6		bis(2-Chloroethyl)ether	1.538	1.546	-0.5	98	0.00
7	I	Naphthalene-d8	1.000	1.000	0.0	98	0.00
8	S	Nitrobenzene-d5	0.424	0.412	2.8	94	0.00
9		Naphthalene	4.646	4.738	-2.0	97	0.00
10		Hexachlorobutadiene	0.249	0.270	-8.4	96	0.00
11	SURR	2-Methylnaphthalene-d10	0.662	0.672	-1.5	97	0.00
12		2-Methylnaphthalene	0.789	0.801	-1.5	96	0.00
13	I	Acenaphthene-d10	1.000	1.000	0.0	99	0.00
14	S	2,4,6-Tribromophenol	0.182	0.155	14.8	86	0.00
15	S	2-Fluorobiphenyl	1.725	1.733	-0.5	98	0.00
16		Acenaphthylene	2.282	2.195	3.8	95	0.00
17		Acenaphthene	1.418	1.391	1.9	97	0.00
18		Fluorene	1.782	1.740	2.4	95	-0.01
19	I	Phenanthrene-d10	1.000	1.000	0.0	99	0.00
20		4-Bromophenyl-phenylether	0.240	0.229	4.6	92	0.00
21		Hexachlorobenzene	0.248	0.249	-0.4	98	0.00
22		Pentachlorophenol	0.069	0.052	24.6	90	0.00
23		Phenanthrene	1.228	1.220	0.7	97	-0.01
24		Anthracene	1.147	1.106	3.6	96	-0.01
25	SURR	Fluoranthene-d10	5.268	5.139	2.4	97	0.00
26		Fluoranthene	1.550	1.529	1.4	99	0.00
27	I	Chrysene-d12	1.000	1.000	0.0	101	0.01
28		Pyrene	1.680	1.641	2.3	104	0.00
29	S	Terphenyl-d14	0.852	0.838	1.6	98	0.00
30		Benzo(a)anthracene	1.333	1.268	4.9	96	0.00
31		Chrysene	1.314	1.283	2.4	97	0.00
32		Bis(2-ethylhexyl)phthalate	3.072	2.527	17.7	88	0.00
33		Indeno(1,2,3-cd)pyrene	1.269	1.190	6.2	98	0.00
34	I	Perylene-d12	1.000	1.000	0.0	97	0.00
35		Benzo(b)fluoranthene	1.364	1.377	-1.0	97	0.00
36		Benzo(k)fluoranthene	1.402	1.423	-1.5	98	0.00
37	C	Benzo(a)pyrene	1.269	1.250	1.5	96	0.00
38		Dibenzo(a,h)anthracene	1.111	1.090	1.9	98	0.00
39		Benzo(g,h,i)perylene	1.193	1.167	2.2	95	0.00

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Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)

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