Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_N\DATA\BN022718\
Data File : BN000187.D
Acq On : 27 Feb 2018 12:42
Operator : JU/SJ
Sample : SSTDICV0.4
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : ICVBE022718

Quant Time: Feb 27 15:14:39 2018
Quant Method : Z:\HPCHEM1\BNA_N\METHODS\8270-SIM-BN022718.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Feb 27 15:07:57 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	Amount	Calc.	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	0.400	0.400	0.0	82	0.00
2		1,4-Dioxane	0.400	0.392	2.0	85	0.00
3		n-Nitrosodimethylamine	0.400	0.376	6.0	78	0.00
4	S	2-Fluorophenol	0.400	0.392	2.0	79	0.00
5	S	Phenol-d6	0.400	0.386	3.5	80	0.00
6		bis(2-Chloroethyl)ether	0.400	0.387	3.3	77	0.00
7	I	Naphthalene-d8	0.400	0.400	0.0	77	0.00
8	S	Nitrobenzene-d5	0.400	0.391	2.3	75	0.00
9		Naphthalene	0.400	0.401	-0.3	76	0.00
10		Hexachlorobutadiene	0.400	0.407	-1.7	76	0.00
11	SURR	2-Methylnaphthalene-d10	0.400	0.398	0.5	76	0.00
12		2-Methylnaphthalene	0.400	0.395	1.3	76	0.00
13	I	Acenaphthene-d10	0.400	0.400	0.0	76	0.01
14	S	2,4,6-Tribromophenol	0.400	0.399	0.3	77	0.00
15	S	2-Fluorobiphenyl	0.400	0.410	-2.5	76	0.00
16		Acenaphthylene	0.400	0.378	5.5	74	0.00
17		Acenaphthene	0.400	0.400	0.0	76	0.00
18		Fluorene	0.400	0.393	1.8	76	0.01
19	I	Phenanthrene-d10	0.400	0.400	0.0	77	0.00
20		4-Bromophenyl-phenylether	0.400	0.390	2.5	76	0.00
21		Hexachlorobenzene	0.400	0.400	0.0	77	0.01
22		Pentachlorophenol	0.400	0.380	5.0	74	0.00
23		Phenanthrene	0.400	0.395	1.3	76	0.00
24		Anthracene	0.400	0.378	5.5	75	0.00
25	SURR	Fluoranthene-d10	0.400	0.387	3.3	76	0.00
26		Fluoranthene	0.400	0.382	4.5	75	0.00
27	I	Chrysene-d12	0.400	0.400	0.0	75	0.00
28		Pyrene	0.400	0.402	-0.5	75	0.00
29	S	Terphenyl-d14	0.400	0.411	-2.7	76	0.01
30		Benzo(a)anthracene	0.400	0.382	4.5	72	0.00
31		Chrysene	0.400	0.413	-3.2	76	0.01
32		Bis(2-ethylhexyl)phthalate	0.400	0.375	6.3	77	0.01
33		Indeno(1,2,3-cd)pyrene	0.400	0.435	-8.7	78	0.01
34	I	Perylene-d12	0.400	0.400	0.0	75	0.00
35		Benzo(b)fluoranthene	0.400	0.405	-1.3	77	0.00
36		Benzo(k)fluoranthene	0.400	0.391	2.3	73	0.01
37	C	Benzo(a)pyrene	0.400	0.396	1.0	75	0.00
38		Dibenzo(a,h)anthracene	0.400	0.429	-7.2	82	0.01
39		Benzo(g,h,i)perylene	0.400	0.397	0.8	74	0.01

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

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