Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN090619\
Data File : BN007557.D
Acq On : 07 Sep 2019 00:25
Operator : HP/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_N
LabSampleId : SSTDCCC0.4

Quant Time: Sep 07 00:58:48 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN082319.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Aug 29 12:18:45 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	158	-0.02
2		1,4-Dioxane	0.400	0.428	-7.0	163	0.00
3		n-Nitrosodimethylamine	0.400	0.250	37.5#	95	0.00
4	S	2-Fluorophenol	0.400	0.343	14.2	135	-0.02
5	S	Phenol-d6	0.400	0.319	20.3	133	0.00
6		bis(2-Chloroethyl)ether	0.400	0.376	6.0	150	0.00
7	I	Naphthalene-d8	0.400	0.400	0.0	157	-0.03
8	S	Nitrobenzene-d5	0.400	0.354	11.5	145	0.00
9		Naphthalene	0.400	0.397	0.8	151	-0.01
10		Hexachlorobutadiene	0.400	0.357	10.8	134	-0.01
11	SURR	2-Methylnaphthalene-d10	0.400	0.371	7.3	143	-0.01
12		2-Methylnaphthalene	0.400	0.379	5.3	144	0.00
13	I	Acenaphthene-d10	0.400	0.400	0.0	152	-0.02
14	S	2,4,6-Tribromophenol	0.400	0.293	26.8#	113	-0.02
15	S	2-Fluorobiphenyl	0.400	0.406	-1.5	150	0.00
16		Acenaphthylene	0.400	0.341	14.8	128	-0.02
17		Acenaphthene	0.400	0.397	0.8	147	0.00
18		Fluorene	0.400	0.387	3.3	144	-0.02
19	I	Phenanthrene-d10	0.400	0.400	0.0	150	0.00
20		4-Bromophenyl-phenylether	0.400	0.201	49.8#	87	0.00
21		Hexachlorobenzene	0.400	0.394	1.5	146	-0.02
22		Pentachlorophenol	0.400	0.218	45.5#	87	0.00
23		Phenanthrene	0.400	0.397	0.8	148	-0.02
24		Anthracene	0.400	0.342	14.5	127	0.00
25	SURR	Fluoranthene-d10	0.400	0.356	11.0	129	-0.01
26		Fluoranthene	0.400	0.366	8.5	133	-0.01
27	I	Chrysene-d12	0.400	0.400	0.0	144	-0.01
28		Pyrene	0.400	0.378	5.5	133	-0.01
29	S	Terphenyl-d14	0.400	0.372	7.0	130	-0.01
30		Benzo(a)anthracene	0.400	0.354	11.5	125	-0.01
31		Chrysene	0.400	0.401	-0.3	139	-0.01
32		Bis(2-ethylhexyl)phthalate	0.400	0.253	36.8#	89	-0.02
33		Indeno(1,2,3-cd)pyrene	0.400	0.393	1.8	137	-0.03
34	I	Perylene-d12	0.400	0.400	0.0	144	-0.02
35		Benzo(b)fluoranthene	0.400	0.366	8.5	128	-0.02
36		Benzo(k)fluoranthene	0.400	0.437	-9.2	151	-0.02
37	C	Benzo(a)pyrene	0.400	0.373	6.8	130	-0.02
38		Dibenzo(a,h)anthracene	0.400	0.381	4.8	132	-0.03
39		Benzo(g,h,i)perylene	0.400	0.391	2.3	136	-0.03

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

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