Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN072320\
Data File : BN011224.D
Acq On : 22 Jul 2020 18:57
Operator : CG/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_N
LabSampleId : SSTDCCC0.4

Quant Time: Jul 23 03:55:41 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN071120.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jul 22 19:29:31 2020
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	87	0.00
2		1,4-Dioxane	0.400	0.345	13.8	70	0.00
3		n-Nitrosodimethylamine	0.400	0.365	8.8	84	0.00
4	S	2-Fluorophenol	0.400	0.366	8.5	77	0.00
5	S	Phenol-d6	0.400	0.421	-5.2	89	0.00
6		bis(2-Chloroethyl)ether	0.400	0.369	7.8	86	0.00
7	I	Naphthalene-d8	0.400	0.400	0.0	98	0.00
8	S	Nitrobenzene-d5	0.400	0.371	7.3	90	0.01
9		Naphthalene	0.400	0.377	5.8	93	0.00
10		Hexachlorobutadiene	0.400	0.378	5.5	89	0.00
11	SURR	2-Methylnaphthalene-d10	0.400	0.435	-8.7	101	0.00
12		2-Methylnaphthalene	0.400	0.406	-1.5	96	0.00
13	I	Acenaphthene-d10	0.400	0.400	0.0	116	0.00
14	S	2,4,6-Tribromophenol	0.400	0.377	5.8	107	0.00
15	S	2-Fluorobiphenyl	0.400	0.354	11.5	96	0.00
16		Acenaphthylene	0.400	0.360	10.0	108	0.00
17		Acenaphthene	0.400	0.365	8.8	106	0.00
18		Fluorene	0.400	0.376	6.0	96	0.00
19	I	Phenanthrene-d10	0.400	0.400	0.0	129	0.00
20		4,6-Dinitro-2-methylphenol	0.400	0.374	6.5	108	0.00
21		4-Bromophenyl-phenylether	0.400	0.361	9.8	116	0.00
22		Hexachlorobenzene	0.400	0.340	15.0	108	0.00
23		Atrazine	0.400	0.388	3.0	125	-0.01
24		Pentachlorophenol	0.400	0.352	12.0	122	0.00
25		Phenanthrene	0.400	0.374	6.5	121	0.00
26		Anthracene	0.400	0.360	10.0	117	0.00
27	SURR	Fluoranthene-d10	0.400	0.374	6.5	119	0.00
28		Fluoranthene	0.400	0.362	9.5	112	0.00
29	I	Chrysene-d12	0.400	0.400	0.0	106	-0.01
30		Pyrene	0.400	0.459	-14.8	107	0.00
31	S	Terphenyl-d14	0.400	0.447	-11.7	112	0.00
32		Benzo(a)anthracene	0.400	0.375	6.3	102	0.00
33		Chrysene	0.400	0.399	0.3	106	-0.01
34		Bis(2-ethylhexyl)phthalate	0.400	0.482	-20.5	129	0.00
35		Indeno(1,2,3-cd)pyrene	0.400	0.347	13.3	80	-0.01
36	I	Perylene-d12	0.400	0.400	0.0	98	-0.01
37		Benzo(b)fluoranthene	0.400	0.357	10.8	97	0.00
38		Benzo(k)fluoranthene	0.400	0.390	2.5	95	-0.01
39	C	Benzo(a)pyrene	0.400	0.348	13.0	84	-0.01
40		Dibenzo(a,h)anthracene	0.400	0.313	21.8	75	-0.02
41		Benzo(g,h,i)perylene	0.400	0.329	17.8	78	-0.01

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

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