Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN080720\
Data File : BN011392.D
Acq On : 07 Aug 2020 10:40
Operator : CG/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_N
LabSampleId : SSTDCCC0.4

Quant Time: Aug 07 11:52:46 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN072820.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jul 31 12:17:46 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	91	0.00
2		1,4-Dioxane	0.400	0.393	1.8	87	0.00
3		n-Nitrosodimethylamine	0.400	0.336	16.0	90	0.00
4	S	2-Fluorophenol	0.400	0.378	5.5	96	0.00
5	S	Phenol-d6	0.400	0.349	12.8	78	0.00
6		bis(2-Chloroethyl)ether	0.400	0.322	19.5	73	0.00
7	I	Naphthalene-d8	0.400	0.400	0.0	94	-0.01
8	S	Nitrobenzene-d5	0.400	0.333	16.8	85	0.00
9		Naphthalene	0.400	0.355	11.3	87	-0.01
10		Hexachlorobutadiene	0.400	0.391	2.3	95	-0.01
11	SURR	2-Methylnaphthalene-d10	0.400	0.336	16.0	85	-0.01
12		2-Methylnaphthalene	0.400	0.334	16.5	85	-0.01
13	I	Acenaphthene-d10	0.400	0.400	0.0	77	-0.01
14	S	2,4,6-Tribromophenol	0.400	0.319	20.3	73	0.00
15	S	2-Fluorobiphenyl	0.400	0.378	5.5	84	-0.01
16		Acenaphthylene	0.400	0.346	13.5	82	-0.01
17		Acenaphthene	0.400	0.357	10.8	76	-0.01
18		Fluorene	0.400	0.327	18.3	73	-0.01
19	I	Phenanthrene-d10	0.400	0.400	0.0	71	-0.01
20		4,6-Dinitro-2-methylphenol	0.400	0.331	17.3	55	0.00
21		4-Bromophenyl-phenylether	0.400	0.343	14.2	73	-0.01
22		Hexachlorobenzene	0.400	0.359	10.3	75	-0.01
23		Atrazine	0.400	0.341	14.8	76	-0.01
24		Pentachlorophenol	0.400	0.355	11.3	86	-0.01
25		Phenanthrene	0.400	0.352	12.0	68	-0.01
26		Anthracene	0.400	0.352	12.0	77	-0.01
27	SURR	Fluoranthene-d10	0.400	0.401	-0.3	89	0.00
28		Fluoranthene	0.400	0.399	0.3	90	0.00
29	I	Chrysene-d12	0.400	0.400	0.0	103	-0.01
30		Pyrene	0.400	0.331	17.3	91	0.00
31	S	Terphenyl-d14	0.400	0.321	19.8	89	0.00
32		Benzo(a)anthracene	0.400	0.368	8.0	97	-0.01
33		Chrysene	0.400	0.392	2.0	101	-0.01
34		Bis(2-ethylhexyl)phthalate	0.400	0.265	33.8#	69	-0.01
35		Indeno(1,2,3-cd)pyrene	0.400	0.480	-20.0	132	-0.02
36	I	Perylene-d12	0.400	0.400	0.0	128	-0.02
37		Benzo(b)fluoranthene	0.400	0.350	12.5	129	-0.01
38		Benzo(k)fluoranthene	0.400	0.333	16.8	114	-0.01
39	C	Benzo(a)pyrene	0.400	0.349	12.8	124	-0.02
40		Dibenzo(a,h)anthracene	0.400	0.376	6.0	139	-0.02
41		Benzo(g,h,i)perylene	0.400	0.376	6.0	132	-0.02

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

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