Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN082120\
Data File : BN011597.D
Acq On : 22 Aug 2020 21:10
Operator : CG/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 42 Sample Multiplier: 1

Instrument : BNA_N
LabSampleId : SSTDCCC0.4

Quant Time: Aug 24 01:26:01 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN082120.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat Aug 22 04:52:07 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	198	0.00
2		1,4-Dioxane	0.400	0.360	10.0	179	0.00
3		n-Nitrosodimethylamine	0.400	0.386	3.5	204	-0.02
4	S	2-Fluorophenol	0.400	0.431	-7.7	244	0.00
5	S	Phenol-d6	0.400	0.466	-16.5	302	-0.02
6		bis(2-Chloroethyl)ether	0.400	0.402	-0.5	259	-0.02
7	I	Naphthalene-d8	0.400	0.400	0.0	227	-0.01
8	S	Nitrobenzene-d5	0.400	0.359	10.3	247	-0.01
9		Naphthalene	0.400	0.377	5.8	226	-0.03
10		Hexachlorobutadiene	0.400	0.339	15.3	192	-0.01
11	SURR	2-Methylnaphthalene-d10	0.400	0.307	23.3	196	-0.02
12		2-Methylnaphthalene	0.400	0.319	20.3	200	-0.02
13	I	Acenaphthene-d10	0.400	0.400	0.0	175	0.00
14	S	2,4,6-Tribromophenol	0.400	0.337	15.8	173	-0.01
15	S	2-Fluorobiphenyl	0.400	0.415	-3.7	191	-0.03
16		Acenaphthylene	0.400	0.392	2.0	191	-0.01
17		Acenaphthene	0.400	0.375	6.3	176	0.00
18		Fluorene	0.400	0.353	11.8	168	-0.01
19	I	Phenanthrene-d10	0.400	0.400	0.0	173	0.00
20		4,6-Dinitro-2-methylphenol	0.400	0.234	41.5#	56	-0.04
21		4-Bromophenyl-phenylether	0.400	0.170	57.5#	63	0.00
22		Hexachlorobenzene	0.400	0.359	10.3	163	-0.01
23		Atrazine	0.400	0.253	36.8#	177	-0.01
24		Pentachlorophenol	0.400	0.315	21.3	172	-0.02
25		Phenanthrene	0.400	0.353	11.8	170	-0.01
26		Anthracene	0.400	0.366	8.5	188	-0.01
27	SURR	Fluoranthene-d10	0.400	0.327	18.3	173	0.00
28		Fluoranthene	0.400	0.320	20.0	169	0.00
29	I	Chrysene-d12	0.400	0.400	0.0	222	-0.01
30		Pyrene	0.400	0.328	18.0	182	0.00
31	S	Terphenyl-d14	0.400	0.325	18.8	191	0.00
32		Benzo(a)anthracene	0.400	0.391	2.3	258	-0.01
33		Chrysene	0.400	0.370	7.5	214	0.00
34		Bis(2-ethylhexyl)phthalate	0.400	0.457	-14.2	281	-0.01
35		Indeno(1,2,3-cd)pyrene	0.400	0.497	-24.2	301	-0.02
36	I	Perylene-d12	0.400	0.400	0.0	247	-0.02
37		Benzo(b)fluoranthene	0.400	0.352	12.0	289	-0.01
38		Benzo(k)fluoranthene	0.400	0.314	21.5	237	-0.01
39	C	Benzo(a)pyrene	0.400	0.370	7.5	259	-0.01
40		Dibenzo(a,h)anthracene	0.400	0.437	-9.2	317	-0.03
41		Benzo(g,h,i)perylene	0.400	0.399	0.3	273	-0.02

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

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