Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN031320\
Data File : BN010212.D
Acq On : 13 Mar 2020 22:10
Operator : CG/JU
Sample : SSTDICC5.0
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId :

Quant Time: Mar 16 02:04:22 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN031320.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Mar 16 01:51:14 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	404	0.00
2		1,4-Dioxane	0.400	4.467	-1016.7#	4536	0.00
3		n-Nitrosodimethylamine	0.400	2.750	-587.5#	3154	-0.02
4	S	2-Fluorophenol	0.400	4.701	-1075.2#	4929	0.00
5	S	Phenol-d6	0.400	4.937	-1134.3#	5249	0.00
6		bis(2-Chloroethyl)ether	0.400	4.373	-993.3#	4590	0.00
7	I	Naphthalene-d8	0.400	0.400	0.0	406	0.00
8	S	Nitrobenzene-d5	0.400	3.841	-860.3#	4142	0.00
9		Naphthalene	0.400	4.618	-1054.5#	4821	0.00
10		Hexachlorobutadiene	0.400	4.729	-1082.2#	4781	0.00
11	SURR	2-Methylnaphthalene-d10	0.400	4.376	-994.0#	4596	0.00
12		2-Methylnaphthalene	0.400	4.859	-1114.8#	5116	0.00
13	I	Acenaphthene-d10	0.400	0.400	0.0	419	0.00
14	S	2,4,6-Tribromophenol	0.400	5.200	-1200.0#	6095	-0.01
15	S	2-Fluorobiphenyl	0.400	4.788	-1097.0#	5224	0.00
16		Acenaphthylene	0.400	5.518	-1279.5#	6439	0.00
17		Acenaphthene	0.400	4.820	-1105.0#	5267	0.00
18		Fluorene	0.400	4.826	-1106.5#	5211	0.00
19	I	Phenanthrene-d10	0.400	0.400	0.0	423	0.00
20		4-Bromophenyl-phenylether	0.400	4.242	-960.5#	4731	0.00
21		Hexachlorobenzene	0.400	4.551	-1037.8#	5048	0.00
22		Pentachlorophenol	0.400	6.474	-1518.5#	9191	0.00
23		Phenanthrene	0.400	4.505	-1026.2#	5155	0.00
24		Anthracene	0.400	5.351	-1237.7#	6293	0.00
25	SURR	Fluoranthene-d10	0.400	4.490	-1022.5#	5094	0.00
26		Fluoranthene	0.400	4.687	-1071.8#	5351	0.00
27	I	Chrysene-d12	0.400	0.400	0.0	478	0.00
28		Pyrene	0.400	4.449	-1012.2#	5482	0.00
29	S	Terphenyl-d14	0.400	4.798	-1099.5#	5875	0.00
30		Benzo(a)anthracene	0.400	5.316	-1229.0#	7152	0.00
31		Chrysene	0.400	4.400	-1000.0#	5393	0.00
32		Bis(2-ethylhexyl)phthalate	0.400	5.087	-1171.7#	6761	0.00
33		Indeno(1,2,3-cd)pyrene	0.400	5.075	-1168.7#	6075	0.00
34	I	Perylene-d12	0.400	0.400	0.0	516	0.00
35		Benzo(b)fluoranthene	0.400	4.305	-976.2#	6067	0.00
36		Benzo(k)fluoranthene	0.400	3.925	-881.3#	5322	0.00
37	C	Benzo(a)pyrene	0.400	4.452	-1013.0#	5899	0.00
38		Dibenzo(a,h)anthracene	0.400	4.513	-1028.3#	6193	0.00
39		Benzo(g,h,i)perylene	0.400	4.233	-958.3#	5783	0.00

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

(#) = Out of Range SPCC's out = 0 CCC's out = 1