Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN073018\
Data File : BN002188.D
Acq On : 30 Jul 2018 16:19
Operator : JU/SJ
Sample : SSTDICV0.4
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : ICVBN073018

Quant Time: Jul 30 16:49:10 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN073018.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jul 30 16:17:29 2018
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	85	0.00
2		1,4-Dioxane	0.400	0.388	3.0	92	0.00
3		n-Nitrosodimethylamine	0.400	0.374	6.5	84	0.00
4	S	2-Fluorophenol	0.400	0.389	2.8	87	0.00
5	S	Phenol-d6	0.400	0.389	2.8	89	0.00
6		bis(2-Chloroethyl)ether	0.400	0.391	2.3	84	0.00
7	I	Naphthalene-d8	0.400	0.400	0.0	88	0.00
8	S	Nitrobenzene-d5	0.400	0.399	0.3	89	0.00
9		Naphthalene	0.400	0.410	-2.5	92	0.00
10		Hexachlorobutadiene	0.400	0.410	-2.5	92	0.00
11	SURR	2-Methylnaphthalene-d10	0.400	0.404	-1.0	92	0.00
12		2-Methylnaphthalene	0.400	0.409	-2.2	91	0.00
13	I	Acenaphthene-d10	0.400	0.400	0.0	84	0.00
14	S	2,4,6-Tribromophenol	0.400	0.346	13.5	75	0.00
15	S	2-Fluorobiphenyl	0.400	0.406	-1.5	91	0.00
16		Acenaphthylene	0.400	0.392	2.0	87	0.01
17		Acenaphthene	0.400	0.402	-0.5	88	0.00
18		Fluorene	0.400	0.401	-0.3	85	0.00
19	I	Phenanthrene-d10	0.400	0.400	0.0	71	0.00
20		4-Bromophenyl-phenylether	0.400	0.411	-2.7	80	0.01
21		Hexachlorobenzene	0.400	0.424	-6.0	81	0.00
22		Pentachlorophenol	0.400	0.335	16.3	62	0.00
23		Phenanthrene	0.400	0.405	-1.3	75	0.00
24		Anthracene	0.400	0.383	4.3	72	0.00
25	SURR	Fluoranthene-d10	0.400	0.364	9.0	64	0.00
26		Fluoranthene	0.400	0.363	9.3	63	0.00
27	I	Chrysene-d12	0.400	0.400	0.0	52	0.00
28		Pyrene	0.400	0.441	-10.2	62	0.00
29	S	Terphenyl-d14	0.400	0.424	-6.0	60	0.00
30		Benzo(a)anthracene	0.400	0.392	2.0	54	0.00
31		Chrysene	0.400	0.412	-3.0	56	0.00
32		Bis(2-ethylhexyl)phthalate	0.400	0.355	11.3	47	-0.01
33		Indeno(1,2,3-cd)pyrene	0.400	0.448	-12.0	69	0.00
34	I	Perylene-d12	0.400	0.400	0.0	58	0.00
35		Benzo(b)fluoranthene	0.400	0.394	1.5	58	0.00
36		Benzo(k)fluoranthene	0.400	0.401	-0.3	61	0.00
37	C	Benzo(a)pyrene	0.400	0.406	-1.5	62	0.00
38		Dibenzo(a,h)anthracene	0.400	0.422	-5.5	70	0.00
39		Benzo(g,h,i)perylene	0.400	0.427	-6.7	70	0.00

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

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