Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\Data\BE092815\
Data File : BE090977.D
Acq On : 28 Sep 2015 21:57
Operator : UM/IZ
Sample : SSTDICV001
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : ICVBE092815

Quant Time: Sep 30 16:02:05 2015
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE092815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Sep 30 15:53:58 2015
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	5.000	5.000	0.0	204	0.00
2		1,4-Dioxane	1.000	0.706	29.4#	158	0.00
3		n-Nitrosodimethylamine	1.000	0.955	4.5	199	0.00
4	S	2-Fluorophenol	1.000	0.863	13.7	180	0.00
5	S	Phenol-d6	1.000	0.976	2.4	223	-0.03
6		bis(2-Chloroethyl)ether	1.000	0.912	8.8	166	-0.05
7	I	Naphthalene-d8	5.000	5.000	0.0	220	-0.01
8	S	Nitrobenzene-d5	1.000	0.976	2.4	234	-0.02
9		Nitrobenzene	1.000	1.087	-8.7	254	-0.02
10		Naphthalene	1.000	0.919	8.1	204	0.00
11		Hexachlorobutadiene	1.000	0.846	15.4	185	0.00
12		2-Methylnaphthalene	1.000	0.939	6.1	219	-0.03
13	I	Acenaphthene-d10	5.000	5.000	0.0	228	0.00
14	S	2,4,6-Tribromophenol	1.000	0.873	12.7	239	-0.02
15	S	2-Fluorobiphenyl	1.000	0.885	11.5	217	-0.01
16		Acenaphthylene	1.000	0.877	12.3	218	0.00
17		Acenaphthene	1.000	0.869	13.1	211	0.00
18		Fluorene	1.000	0.884	11.6	216	0.00
19	I	Phenanthrene-d10	5.000	5.000	0.0	221	0.00
20		4-Bromophenyl-phenylether	1.000	0.885	11.5	213	-0.02
21		Hexachlorobenzene	1.000	0.871	12.9	195	0.00
22		Pentachlorophenol	1.000	0.989	1.1	267	-0.02
23		Phenanthrene	1.000	0.932	6.8	216	-0.02
24		Anthracene	1.000	0.921	7.9	219	-0.02
25		Fluoranthene	1.000	0.944	5.6	229	-0.01
26	I	Chrysene-d12	5.000	5.000	0.0	221	0.00
27		Benzidine	1.000	0.201	79.9#	62	-0.01
28		Pyrene	1.000	0.998	0.2	240	-0.01
29	S	Terphenyl-d14	1.000	1.019	-1.9	246	-0.02
30		Benzo(a)anthracene	1.000	0.996	0.4	231	0.00
31		3,3'-Dichlorobenzidine	1.000	1.111	-11.1	284	0.00
32		Chrysene	1.000	0.871	12.9	201	0.00
33		Bis(2-ethylhexyl)phthalate	1.000	1.369	-36.9#	386	0.00
34		Indeno(1,2,3-cd)pyrene	1.000	1.089	-8.9	264	-0.01
35	I	Perylene-d12	5.000	5.000	0.0	233	0.00
36		Benzo(b)fluoranthene	1.000	0.966	3.4	251	0.00
37		Benzo(k)fluoranthene	1.000	0.833	16.7	207	0.00
38	C	Benzo(a)pyrene	1.000	0.942	5.8	254	0.00
39		Dibenzo(a,h)anthracene	1.000	1.021	-2.1	283	-0.01
40		Benzo(g,h,i)perylene	1.000	0.942	5.8	253	-0.02

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

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