Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE090116\
Data File : BE092333.D
Acq On : 1 Sep 2016 17:47
Operator : UM/SJ
Sample : SSTDICV0.4
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : ICVBE090116

Quant Time: Sep 02 10:55:34 2016
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE090116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Sep 02 03:36:16 2016
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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	92	0.00
2		1,4-Dioxane	0.717	0.800	-11.6	106	0.00
3		n-Nitrosodimethylamine	0.744	0.715	3.9	94	0.00
4	S	2-Fluorophenol	1.330	1.289	3.1	98	0.00
5	S	Phenol-d6	1.842	1.548	16.0	98	0.00
6		bis(2-Chloroethyl)ether	1.424	1.284	9.8	101	0.02
7	I	Naphthalene-d8	1.000	1.000	0.0	91	0.00
8	S	Nitrobenzene-d5	0.365	0.300	17.8	97	0.00
9		Naphthalene	1.106	1.172	-6.0	104	0.00
10		Hexachlorobutadiene	0.175	0.193	-10.3	108	0.00
11		2-Methylnaphthalene	0.728	0.750	-3.0	103	0.01
12	I	Acenaphthene-d10	1.000	1.000	0.0	89	-0.02
13	S	2,4,6-Tribromophenol	0.196	0.171	12.8	91	0.00
14	S	2-Fluorobiphenyl	1.526	1.566	-2.6	99	0.00
15		Acenaphthylene	8.673	9.155	-5.6	103	0.00
16		Acenaphthene	1.394	1.500	-7.6	102	0.00
17		Fluorene	1.772	1.896	-7.0	102	0.00
18	I	Phenanthrene-d10	1.000	1.000	0.0	91	0.00
19		Hexachlorobenzene	0.220	0.236	-7.3	104	0.00
20		Pentachlorophenol	0.070	0.053	24.3	89	0.00
21		Phenanthrene	1.215	1.282	-5.5	100	0.00
22		Anthracene	1.181	1.195	-1.2	100	0.00
23		Fluoranthene	6.088	6.294	-3.4	98	0.00
24	I	Chrysene-d12	1.000	1.000	0.0	92	0.00
25		Pyrene	4.561	4.563	-0.0	97	0.00
26	S	Terphenyl-d14	0.689	0.668	3.0	95	0.00
27		Benzo(a)anthracene	5.548	5.794	-4.4	102	0.00
28		Chrysene	4.695	5.242	-11.7	102	0.00
29		Bis(2-ethylhexyl)phthalate	0.534	0.481	9.9	87	0.00
30		Indeno(1,2,3-cd)pyrene	5.479	5.746	-4.9	100	0.00
31	I	Perylene-d12	1.000	1.000	0.0	89	0.00
32		Benzo(b)fluoranthene	1.350	1.327	1.7	103	0.00
33		Benzo(k)fluoranthene	1.317	1.462	-11.0	100	-0.01
34	C	Benzo(a)pyrene	1.282	1.329	-3.7	101	0.00
35		Dibenzo(a,h)anthracene	1.211	1.235	-2.0	100	-0.02
36		Benzo(g,h,i)perylene	5.006	5.210	-4.1	100	0.00

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