Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE062915\
Data File : BE089990.D
Acq On : 29 Jun 2015 19:08
Operator : TP/IZ
Sample : SSTDICV001
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : ICVBE062915

Quant Time: Jun 30 07:02:35 2015
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\8270-SIM-BE062915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jun 30 07:01:01 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	130	0.00
2		1,4-Dioxane	1.000	0.841	15.9	114	0.00
3		n-Nitrosodimethylamine	1.000	0.839	16.1	114	0.00
4	S	2-Fluorophenol	1.000	0.836	16.4	109	0.00
5	S	Phenol-d6	1.000	0.861	13.9	114	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	134	0.00
7	S	Nitrobenzene-d5	1.000	0.840	16.0	114	0.00
8		Nitrobenzene	1.000	0.842	15.8	116	0.00
9		Naphthalene	1.000	0.911	8.9	124	0.00
10		Hexachlorobutadiene	1.000	0.926	7.4	126	0.00
11		2-Methylnaphthalene	1.000	0.896	10.4	122	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	135	0.00
13	S	2,4,6-Tribromophenol	1.000	0.776	22.4	107	0.00
14	S	2-Fluorobiphenyl	1.000	0.883	11.7	123	0.00
15		Acenaphthylene	1.000	0.902	9.8	126	0.00
16		Acenaphthene	1.000	0.942	5.8	129	0.00
17		Fluorene	1.000	0.922	7.8	128	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	134	0.00
19		4-Bromophenyl-phenylether	1.000	0.889	11.1	122	0.00
20		Hexachlorobenzene	1.000	0.928	7.2	130	0.00
21		Pentachlorophenol	1.000	0.754	24.6	96	0.00
22		Phenanthrene	1.000	0.921	7.9	127	0.00
23		Anthracene	1.000	0.923	7.7	126	0.00
24		Fluoranthene	1.000	0.922	7.8	126	0.00
25	I	Chrysene-d12	5.000	5.000	0.0	137	0.00
26		Pyrene	1.000	0.915	8.5	128	0.00
27	S	Terphenyl-d14	1.000	0.884	11.6	121	0.00
28		Benzo(a)anthracene	1.000	0.917	8.3	128	0.00
29		Chrysene	1.000	0.895	10.5	124	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	0.706	29.4#	92	0.00
31		Indeno(1,2,3-cd)pyrene	1.000	0.912	8.8	127	0.00
32	I	Perylene-d12	5.000	5.000	0.0	137	0.00
33		Benzo(b)fluoranthene	1.000	0.948	5.2	131	0.00
34		Benzo(k)fluoranthene	1.000	0.912	8.8	126	0.00
35	C	Benzo(a)pyrene	1.000	0.908	9.2	125	0.00
36		Dibenzo(a,h)anthracene	1.000	0.897	10.3	125	0.00
37		Benzo(g,h,i)perylene	1.000	0.898	10.2	125	0.00

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