Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE082415\
Data File : BE090608.D
Acq On : 24 Aug 2015 19:23
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : SSTDCCC001EC

Quant Time: Aug 24 23:01:04 2015
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE081415.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Aug 14 16:52:42 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	95	0.00
2		1,4-Dioxane	1.000	1.290	-29.0#	104	0.00
3		n-Nitrosodimethylamine	1.000	1.046	-4.6	100	0.00
4	S	2-Fluorophenol	1.000	0.753	24.7	71	0.00
5	S	Phenol-d6	1.000	0.676	32.4#	67	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	87	0.00
7	S	Nitrobenzene-d5	1.000	0.902	9.8	82	0.00
8		Nitrobenzene	1.000	0.873	12.7	81	0.00
9		Naphthalene	1.000	0.977	2.3	85	0.00
10		Hexachlorobutadiene	1.000	1.031	-3.1	89	0.00
11		2-Methylnaphthalene	1.000	0.957	4.3	85	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	95	0.00
13	S	2,4,6-Tribromophenol	1.000	0.842	15.8	88	0.00
14	S	2-Fluorobiphenyl	1.000	0.938	6.2	89	0.00
15		Acenaphthylene	1.000	0.935	6.5	91	0.00
16		Acenaphthene	1.000	0.963	3.7	93	0.00
17		Fluorene	1.000	1.022	-2.2	98	0.02
18	I	Phenanthrene-d10	5.000	5.000	0.0	110	0.00
19		4-Bromophenyl-phenylether	1.000	0.926	7.4	105	0.00
20		Hexachlorobenzene	1.000	0.969	3.1	107	0.00
21		Pentachlorophenol	1.000	0.813	18.7	109	0.00
22		Phenanthrene	1.000	0.950	5.0	106	0.00
23		Anthracene	1.000	0.932	6.8	106	0.00
24		Fluoranthene	1.000	0.987	1.3	111	0.00
25	I	Chrysene-d12	5.000	5.000	0.0	125	0.00
26		Pyrene	1.000	0.906	9.4	116	0.00
27	S	Terphenyl-d14	1.000	0.904	9.6	115	0.00
28		Benzo(a)anthracene	1.000	0.920	8.0	121	0.00
29		Chrysene	1.000	0.998	0.2	126	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	0.649	35.1#	86	0.00
31		Indeno(1,2,3-cd)pyrene	1.000	0.888	11.2	120	0.01
32	I	Perylene-d12	5.000	5.000	0.0	122	0.00
33		Benzo(b)fluoranthene	1.000	0.939	6.1	117	0.00
34		Benzo(k)fluoranthene	1.000	1.035	-3.5	125	0.00
35	C	Benzo(a)pyrene	1.000	0.943	5.7	118	0.00
36		Dibenzo(a,h)anthracene	1.000	0.912	8.8	117	0.00
37		Benzo(g,h,i)perylene	1.000	0.927	7.3	115	0.00

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