Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\Data\BE062315\
Data File : BE089931.D
Acq On : 24 Jun 2015 1:35
Operator : TP/IZ
Sample : SSTDICV001
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : SSTDICV001

Quant Time: Jun 24 09:30:59 2015
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\8270-SIM-BE062315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jun 24 09:11:03 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	83	0.00
2		1,4-Dioxane	1.000	1.130	-13.0	92	0.00
3		n-Nitrosodimethylamine	1.000	1.016	-1.6	89	0.00
4	S	2-Fluorophenol	1.000	0.715	28.5#	79	0.00
5	S	Phenol-d6	1.000	0.846	15.4	81	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	81	0.00
7	S	Nitrobenzene-d5	1.000	0.921	7.9	86	0.00
8		Nitrobenzene	1.000	0.915	8.5	87	0.00
9		Naphthalene	1.000	0.913	8.7	86	0.00
10		Hexachlorobutadiene	1.000	0.868	13.2	87	0.00
11		2-Methylnaphthalene	1.000	0.962	3.8	87	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	81	0.00
13	S	2,4,6-Tribromophenol	1.000	0.355	64.5#	76	0.00
14	S	2-Fluorobiphenyl	1.000	0.823	17.7	88	0.00
15		Acenaphthylene	1.000	2.756	-175.6#	85	0.00
16		Acenaphthene	1.000	0.853	14.7	83	0.00
17		Fluorene	1.000	1.302	-30.2#	85	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	80	0.00
19		4-Bromophenyl-phenylether	1.000	0.978	2.2	90	0.00
20		Hexachlorobenzene	1.000	3.787	-278.7#	89	0.00
21		Pentachlorophenol	1.000	0.456	54.4#	83	0.00
22		Phenanthrene	1.000	0.744	25.6#	87	0.00
23		Anthracene	1.000	1.502	-50.2#	87	0.00
24		Fluoranthene	1.000	1.221	-22.1	88	0.00
25	I	Chrysene-d12	5.000	5.000	0.0	83	0.00
26		Pyrene	1.000	0.742	25.8#	87	0.00
27	S	Terphenyl-d14	1.000	1.152	-15.2	91	0.00
28		Benzo(a)anthracene	1.000	0.783	21.7	87	0.00
29		Chrysene	1.000	0.830	17.0	90	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	0.121	87.9#	90	0.00
31		Indeno(1,2,3-cd)pyrene	1.000	0.630	37.0#	84	0.00
32	I	Perylene-d12	5.000	5.000	0.0	82	0.00
33		Benzo(b)fluoranthene	1.000	0.659	34.1#	85	0.00
34		Benzo(k)fluoranthene	1.000	0.839	16.1	88	0.00
35	C	Benzo(a)pyrene	1.000	0.685	31.5#	85	0.00
36		Dibenzo(a,h)anthracene	1.000	0.650	35.0#	86	0.00
37		Benzo(g,h,i)perylene	1.000	0.658	34.2#	84	0.00

(#) = Out of Range SPCC's out = 0 CCC's out = 1