Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\Data\BE090115\
Data File : BE090677.D
Acq On : 1 Sep 2015 21:45
Operator : UM/IZ
Sample : SSTDICV001
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : ICVBE090115

Quant Time: Sep 02 11:46:04 2015
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE090115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Sep 02 11:40:50 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	87	0.00
2		1,4-Dioxane	1.000	0.915	8.5	76	0.00
3		n-Nitrosodimethylamine	1.000	0.846	15.4	72	0.00
4	S	2-Fluorophenol	1.000	0.860	14.0	73	0.00
5	S	Phenol-d6	1.000	0.792	20.8	73	0.01
6	I	Naphthalene-d8	5.000	5.000	0.0	88	0.00
7	S	Nitrobenzene-d5	1.000	0.815	18.5	71	0.00
8		Nitrobenzene	1.000	0.820	18.0	74	0.00
9		Naphthalene	1.000	0.881	11.9	77	0.00
10		Hexachlorobutadiene	1.000	0.920	8.0	78	0.00
11		2-Methylnaphthalene	1.000	0.850	15.0	77	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	88	0.00
13	S	2,4,6-Tribromophenol	1.000	0.771	22.9	67	0.02
14	S	2-Fluorobiphenyl	1.000	0.876	12.4	78	0.00
15		Acenaphthylene	1.000	0.844	15.6	77	0.00
16		Acenaphthene	1.000	0.901	9.9	78	0.00
17		Fluorene	1.000	0.879	12.1	78	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	89	0.00
19		4-Bromophenyl-phenylether	1.000	0.871	12.9	81	0.00
20		Hexachlorobenzene	1.000	0.936	6.4	82	0.00
21		Pentachlorophenol	1.000	0.902	9.8	62	0.02
22		Phenanthrene	1.000	0.880	12.0	79	0.00
23		Anthracene	1.000	0.836	16.4	78	0.00
24		Fluoranthene	1.000	0.794	20.6	76	0.00
25	I	Chrysene-d12	5.000	5.000	0.0	93	0.00
26		Pyrene	1.000	0.766	23.4	75	0.00
27	S	Terphenyl-d14	1.000	0.782	21.8	76	0.00
28		Benzo(a)anthracene	1.000	0.835	16.5	79	0.00
29		Chrysene	1.000	0.909	9.1	83	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	0.717	28.3#	72	0.00
31		Indeno(1,2,3-cd)pyrene	1.000	0.788	21.2	80	0.00
32	I	Perylene-d12	5.000	5.000	0.0	97	0.00
33		Benzo(b)fluoranthene	1.000	0.806	19.4	80	0.00
34		Benzo(k)fluoranthene	1.000	0.837	16.3	82	0.00
35	C	Benzo(a)pyrene	1.000	0.782	21.8#	81	0.00
36		Dibenzo(a,h)anthracene	1.000	0.806	19.4	80	0.00
37		Benzo(g,h,i)perylene	1.000	0.811	18.9	83	0.00

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