Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\Data\BE121417\
Data File : BE094982.D
Acq On : 14 Dec 2017 18:25
Operator : SJ/JU
Sample : SSTDICV0.4
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : ICVBE121417

Quant Time: Dec 14 19:07:14 2017
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE121417.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Dec 14 18:42:53 2017
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	0.394	0.376	4.6	83	0.00
3		n-Nitrosodimethylamine	0.466	0.456	2.1	90	0.00
4	S	2-Fluorophenol	0.621	0.599	3.5	92	0.00
5	S	Phenol-d6	0.923	0.859	6.9	88	0.00
6		bis(2-Chloroethyl)ether	0.901	0.866	3.9	89	0.00
7	I	Naphthalene-d8	1.000	1.000	0.0	90	0.00
8	S	Nitrobenzene-d5	0.328	0.328	0.0	94	0.00
9		Naphthalene	4.697	4.744	-1.0	91	0.00
10		Hexachlorobutadiene	0.214	0.213	0.5	89	0.00
11	SURR	2-Methylnaphthalene-d10	0.655	0.665	-1.5	92	0.00
12		2-Methylnaphthalene	1.172	1.164	0.7	90	0.00
13	I	Acenaphthene-d10	1.000	1.000	0.0	91	0.00
14	S	2,4,6-Tribromophenol	0.093	0.086	7.5	91	0.00
15	S	2-Fluorobiphenyl	1.772	1.753	1.1	89	0.00
16		Acenaphthylene	2.138	2.060	3.6	92	0.00
17		Acenaphthene	1.418	1.398	1.4	91	0.00
18		Fluorene	1.762	1.713	2.8	91	-0.02
19	I	Phenanthrene-d10	1.000	1.000	0.0	90	0.00
20		4-Bromophenyl-phenylether	0.234	0.233	0.4	92	0.00
21		Hexachlorobenzene	0.239	0.243	-1.7	93	0.00
22		Pentachlorophenol	0.046	0.040	13.0	82	0.00
23		Phenanthrene	1.246	1.244	0.2	92	0.00
24		Anthracene	1.279	1.245	2.7	93	0.00
25	SURR	Fluoranthene-d10	5.189	5.023	3.2	91	0.00
26		Fluoranthene	1.538	1.503	2.3	91	0.00
27	I	Chrysene-d12	1.000	1.000	0.0	91	0.00
28		Pyrene	1.355	1.329	1.9	86	0.00
29	S	Terphenyl-d14	0.753	0.761	-1.1	91	0.00
30		Benzo(a)anthracene	1.169	1.122	4.0	89	0.00
31		Chrysene	1.316	1.327	-0.8	92	0.00
32		Bis(2-ethylhexyl)phthalate	1.489	1.232	17.3	84	0.00
33		Indeno(1,2,3-cd)pyrene	1.071	1.004	6.3	84	0.00
34	I	Perylene-d12	1.000	1.000	0.0	85	0.00
35		Benzo(b)fluoranthene	1.469	1.419	3.4	82	0.00
36		Benzo(k)fluoranthene	1.488	1.470	1.2	89	0.00
37	C	Benzo(a)pyrene	1.303	1.223	6.1	81	0.00
38		Dibenzo(a,h)anthracene	1.209	1.149	5.0	82	0.00
39		Benzo(g,h,i)perylene	1.203	1.192	0.9	86	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\Data\BE121417\
Data File : BE094982.D
Acq On : 14 Dec 2017 18:25
Operator : SJ/JU
Sample : SSTDICV0.4
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : ICVBE121417

Quant Time: Dec 14 19:07:14 2017
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE121417.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Dec 14 18:42:53 2017
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)

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