

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE101717\
 Data File : BE094426.D
 Acq On : 17 Oct 2017 10:53
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Oct 17 20:21:52 2017
 Quant Method : Z:\HPCHEM1\BNA_E\Methods\8270-SIM-BE101617.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 16 14:54:40 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	96	-0.01
2	1,4-Dioxane	0.698	0.663	5.0	91	0.00
3	n-Nitrosodimethylamine	0.695	0.647	6.9	89	0.00
4 S	2-Fluorophenol	1.352	1.367	-1.1	96	0.00
5 S	Phenol-d6	2.108	2.047	2.9	90	-0.01
6	bis(2-Chloroethyl)ether	1.559	1.528	2.0	90	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	97	0.00
8 S	Nitrobenzene-d5	0.348	0.340	2.3	96	-0.02
9	Naphthalene	4.473	4.531	-1.3	98	-0.02
10	Hexachlorobutadiene	0.179	0.175	2.2	96	0.00
11 SURR	2-Methylnaphthalene-d10	0.617	0.632	-2.4	99	0.00
12	2-Methylnaphthalene	0.737	0.773	-4.9	103	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	105	0.00
14 S	2,4,6-Tribromophenol	0.152	0.149	2.0	97	0.00
15 S	2-Fluorobiphenyl	1.428	1.423	0.4	102	0.00
16	Acenaphthylene	2.169	2.125	2.0	102	0.00
17	Acenaphthene	1.366	1.362	0.3	105	0.00
18	Fluorene	1.761	1.752	0.5	105	-0.01
19 I	Phenanthrene-d10	1.000	1.000	0.0	98	-0.03
20	4-Bromophenyl-phenylether	0.200	0.206	-3.0	112	-0.03
21	Hexachlorobenzene	0.365	0.364	0.3	101	0.00
22	Pentachlorophenol	0.032	0.025	21.9	83	0.03
23	Phenanthrene	1.574	1.270	19.3	76	0.00
24	Anthracene	1.216	1.209	0.6	101	0.00
25 SURR	Fluoranthene-d10	6.441	6.507	-1.0	103	0.00
26	Fluoranthene	1.954	1.945	0.5	103	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	109	0.00
28	Pyrene	1.509	1.407	6.8	103	0.00
29 S	Terphenyl-d14	0.807	0.765	5.2	103	0.00
30	Benzo(a)anthracene	1.384	1.334	3.6	106	-0.01
31	Chrysene	1.332	1.274	4.4	103	-0.01
32	Bis(2-ethylhexyl)phthalate	3.984	3.650	8.4	104	0.00
33	Indeno(1,2,3-cd)pyrene	1.341	1.307	2.5	107	-0.01
34 I	Perylene-d12	1.000	1.000	0.0	108	0.00
35	Benzo(b)fluoranthene	1.357	1.300	4.2	104	0.00
36	Benzo(k)fluoranthene	1.337	1.327	0.7	106	0.00
37 C	Benzo(a)pyrene	1.275	1.235	3.1	105	0.00
38	Dibenzo(a,h)anthracene	1.207	1.184	1.9	105	0.00
39	Benzo(g,h,i)perylene	1.161	1.134	2.3	105	-0.02

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE101717\
Data File : BE094426.D
Acq On : 17 Oct 2017 10:53
Operator : SJ/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_E
LabSampleId :
SSTDCCC0.4

Quant Time: Oct 17 20:21:52 2017
Quant Method : Z:\HPCHEM1\BNA_E\Methods\8270-SIM-BE101617.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Oct 16 14:54:40 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	