

Data Path : \\74.0.250.170\svoasrv\HPCHEM1\BNA E\Data\BE091317\
 Data File : BE093996.D
 Acq On : 13 Sep 2017 19:24
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Sep 14 03:55:23 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE090517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 05 13:34:56 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	Amount	Calc.	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	0.400	0.400	0.0	108	0.00
2	1,4-Dioxane	0.400	0.489	-22.2	133	0.00
3	n-Nitrosodimethylamine	0.400	0.416	-4.0	112	0.00
4 S	2-Fluorophenol	0.400	0.397	0.8	108	0.00
5 S	Phenol-d6	0.400	0.383	4.3	106	0.00
6	bis(2-Chloroethyl)ether	0.400	0.397	0.8	107	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	106	0.00
8 S	Nitrobenzene-d5	0.400	0.394	1.5	105	0.00
9	Naphthalene	0.400	0.403	-0.8	105	0.00
10	Hexachlorobutadiene	0.400	0.398	0.5	105	-0.02
11 SURR	2-Methylnaphthalene-d10	0.400	0.407	-1.7	105	0.00
12	2-Methylnaphthalene	0.400	0.401	-0.3	104	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	107	0.00
14 S	2,4,6-Tribromophenol	0.400	0.393	1.8	101	0.00
15 S	2-Fluorobiphenyl	0.400	0.398	0.5	104	0.00
16	Acenaphthylene	0.400	0.394	1.5	108	0.00
17	Acenaphthene	0.400	0.401	-0.3	105	0.00
18	Fluorene	0.400	0.401	-0.3	109	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	101	0.00
20	4-Bromophenyl-phenylether	0.400	0.408	-2.0	93	0.00
21	Hexachlorobenzene	0.400	0.329	17.8	78	0.00
22	Pentachlorophenol	0.400	0.442	-10.5	112	0.00
23	Phenanthrene	0.400	0.417	-4.2	96	0.00
24	Anthracene	0.400	0.400	0.0	100	0.00
25 SURR	Fluoranthene-d10	0.400	0.398	0.5	103	0.00
26	Fluoranthene	0.400	0.396	1.0	104	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	103	-0.01
28	Pvrene	0.400	0.400	0.0	104	0.00
29 S	Terphenyl-d14	0.400	0.397	0.8	102	0.00
30	Benzo(a)anthracene	0.400	0.368	8.0	94	0.00
31	Chrysene	0.400	0.409	-2.2	106	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.347	13.3	94	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.321	19.8	84	0.00
34 I	Perylene-d12	0.400	0.400	0.0	96	0.00
35	Benzo(b)fluoranthene	0.400	0.395	1.3	94	-0.01
36	Benzo(k)fluoranthene	0.400	0.432	-8.0	104	0.00
37 C	Benzo(a)pyrene	0.400	0.400	0.0	96	0.00
38	Dibenzo(a,h)anthracene	0.400	0.348	13.0	83	0.00
39	Benzo(a,h,i)perylene	0.400	0.361	9.8	87	0.00

Data Path : \\74.0.250.170\svoasrv\HPCHEM1\BNA E\Data\BE091317\
Data File : BE093996.D
Acq On : 13 Sep 2017 19:24
Operator : SJ/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_E
LabSampleId :
SSTDCCC0.4

Quant Time: Sep 14 03:55:23 2017
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE090517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Sep 05 13:34:56 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	Amount	Calc.	%Dev	Area%	Dev(min)

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