

Data Path : Z:\HPCHEM1\BNA_E\Data\BE080517\
 Data File : BE093672.D
 Acq On : 5 Aug 2017 10:34
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Aug 07 18:08:47 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE080117.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 02 10:58:31 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	110	-0.01
2	1,4-Dioxane	0.400	0.403	-0.8	111	0.00
3	n-Nitrosodimethylamine	0.400	0.390	2.5	114	0.01
4 S	2-Fluorophenol	0.400	0.440	-10.0	126	0.00
5 S	Phenol-d6	0.400	0.435	-8.7	124	-0.01
6	bis(2-Chloroethyl)ether	0.400	0.389	2.8	110	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	112	0.00
8 S	Nitrobenzene-d5	0.400	0.430	-7.5	129	0.00
9	Naphthalene	0.400	0.393	1.8	111	0.00
10	Hexachlorobutadiene	0.400	0.394	1.5	113	-0.02
11 SURR	2-Methylnaphthalene-d10	0.400	0.401	-0.3	113	0.00
12	2-Methylnaphthalene	0.400	0.394	1.5	111	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	111	0.00
14 S	2,4,6-Tribromophenol	0.400	0.447	-11.7	149	0.00
15 S	2-Fluorobiphenyl	0.400	0.390	2.5	107	-0.02
16	Acenaphthylene	0.400	0.424	-6.0	122	0.00
17	Acenaphthene	0.400	0.400	0.0	112	0.00
18	Fluorene	0.400	0.389	2.8	106	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	116	0.00
20	4-Bromophenyl-phenylether	0.400	0.375	6.3	96	0.00
21	Hexachlorobenzene	0.400	0.427	-6.7	109	0.00
22	Pentachlorophenol	0.400	0.534	-33.5#	166	0.00
23	Phenanthrene	0.400	0.406	-1.5	115	0.00
24	Anthracene	0.400	0.432	-8.0	119	0.00
25 SURR	Fluoranthene-d10	0.400	0.425	-6.2	113	0.00
26	Fluoranthene	0.400	0.419	-4.7	112	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	114	-0.01
28	Pyrene	0.400	0.401	-0.3	115	0.00
29 S	Terphenyl-d14	0.400	0.391	2.3	110	0.00
30	Benzo(a)anthracene	0.400	0.444	-11.0	131	0.00
31	Chrysene	0.400	0.397	0.8	113	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.761	-90.3#	253	-0.01
33	Indeno(1,2,3-cd)pyrene	0.400	0.431	-7.7	122	0.00
34 I	Perylene-d12	0.400	0.400	0.0	124	0.00
35	Benzo(b)fluoranthene	0.400	0.383	4.3	122	0.00
36	Benzo(k)fluoranthene	0.400	0.365	8.8	117	0.00
37 C	Benzo(a)pyrene	0.400	0.396	1.0	127	0.00
38	Dibenzo(a,h)anthracene	0.400	0.386	3.5	122	0.00
39	Benzo(g,h,i)perylene	0.400	0.378	5.5	120	0.00

Data Path : Z:\HPCHEM1\BNA_E\Data\BE080517\
Data File : BE093672.D
Acq On : 5 Aug 2017 10:34
Operator : SJ/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_E
LabSampleId :
SSTDCCC0.4

Quant Time: Aug 07 18:08:47 2017
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE080117.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Aug 02 10:58:31 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				