

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE100819\
 Data File : BE100616.D
 Acq On : 8 Oct 2019 19:43
 Operator : JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Oct 09 01:46:28 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE100319.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 09 01:42:20 2019
 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	105	0.00
2	1,4-Dioxane	0.400	0.384	4.0	104	0.00
3	n-Nitrosodimethylamine	0.400	0.433	-8.2	137	-0.01
4 S	2-Fluorophenol	0.400	0.426	-6.5	119	0.00
5 S	Phenol-d6	0.400	0.440	-10.0	126	0.00
6	bis(2-Chloroethyl)ether	0.400	0.394	1.5	110	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	113	0.00
8 S	Nitrobenzene-d5	0.400	0.383	4.3	120	0.00
9	Naphthalene	0.400	0.393	1.8	112	0.00
10	Hexachlorobutadiene	0.400	0.381	4.8	107	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.397	0.8	114	-0.01
12	2-Methylnaphthalene	0.400	0.397	0.8	116	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	117	0.00
14 S	2,4,6-Tribromophenol	0.400	0.402	-0.5	131	0.00
15 S	2-Fluorobiphenyl	0.400	0.389	2.8	115	0.00
16	Acenaphthylene	0.400	0.381	4.8	119	0.00
17	Acenaphthene	0.400	0.385	3.8	117	0.00
18	Fluorene	0.400	0.389	2.8	121	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	116	0.00
20	4-Bromophenyl-phenylether	0.400	0.370	7.5	116	0.00
21	Hexachlorobenzene	0.400	0.384	4.0	114	0.00
22	Pentachlorophenol	0.400	0.552	-38.0#	147	-0.01
23	Phenanthrene	0.400	0.399	0.3	117	0.00
24	Anthracene	0.400	0.373	6.8	120	0.00
25 SURR	Fluoranthene-d10	0.400	0.379	5.3	120	0.00
26	Fluoranthene	0.400	0.391	2.3	120	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	111	0.00
28	Pvrene	0.400	0.443	-10.7	121	0.00
29 S	Terphenyl-d14	0.400	0.418	-4.5	121	0.00
30	Benzo(a)anthracene	0.400	0.380	5.0	110	0.00
31	Chrysene	0.400	0.391	2.3	109	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.371	7.3	119	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.389	2.8	108	0.00
34 I	Perylene-d12	0.400	0.400	0.0	109	0.00
35	Benzo(b)fluoranthene	0.400	0.357	10.8	108	0.00
36	Benzo(k)fluoranthene	0.400	0.368	8.0	107	0.00
37 C	Benzo(a)pyrene	0.400	0.371	7.3	110	0.00
38	Dibenzo(a,h)anthracene	0.400	0.366	8.5	108	0.00
39	Benzo(a,h,i)perylene	0.400	0.367	8.3	105	0.00

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Compound	Amount	Calc.	%Dev	Area%	Dev(min)

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