

Data Path : Z:\svoasrv\HPCHEM1\BNA E\Data\BE110719\
 Data File : BE100700.D
 Acq On : 7 Nov 2019 13:57
 Operator : JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Nov 08 11:18:09 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE101019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 23 16:49:33 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	71	0.00
2	1,4-Dioxane	0.400	0.362	9.5	65	0.00
3	n-Nitrosodimethylamine	0.400	0.360	10.0	70	0.00
4 S	2-Fluorophenol	0.400	0.395	1.3	75	0.00
5 S	Phenol-d6	0.400	0.438	-9.5	83	0.00
6	bis(2-Chloroethyl)ether	0.400	0.395	1.3	75	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	75	0.00
8 S	Nitrobenzene-d5	0.400	0.516	-29.0#	111	0.00
9	Naphthalene	0.400	0.394	1.5	76	0.00
10	Hexachlorobutadiene	0.400	0.399	0.3	78	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.387	3.3	76	0.00
12	2-Methylnaphthalene	0.400	0.391	2.3	77	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	81	0.00
14 S	2,4,6-Tribromophenol	0.400	0.507	-26.7#	117	0.00
15 S	2-Fluorobiphenyl	0.400	0.424	-6.0	88	0.00
16	Acenaphthylene	0.400	0.338	15.5	71	-0.01
17	Acenaphthene	0.400	0.379	5.3	80	0.00
18	Fluorene	0.400	0.385	3.8	81	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	91	-0.01
20	4-Bromophenyl-phenylether	0.400	0.356	11.0	84	0.00
21	Hexachlorobenzene	0.400	0.380	5.0	91	-0.01
22	Pentachlorophenol	0.400	0.303	24.3	77	0.01
23	Phenanthrene	0.400	0.402	-0.5	91	-0.01
24	Anthracene	0.400	0.339	15.3	80	0.00
25 SURR	Fluoranthene-d10	0.400	0.372	7.0	88	0.00
26	Fluoranthene	0.400	0.392	2.0	89	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	81	0.00
28	Pvrene	0.400	0.431	-7.7	88	0.00
29 S	Terphenyl-d14	0.400	0.465	-16.3	100	0.00
30	Benzo(a)anthracene	0.400	0.367	8.3	75	0.00
31	Chrysene	0.400	0.380	5.0	78	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.272	32.0#	55	-0.01
33	Indeno(1,2,3-cd)pyrene	0.400	0.339	15.3	71	-0.01
34 I	Perylene-d12	0.400	0.400	0.0	74	0.00
35	Benzo(b)fluoranthene	0.400	0.359	10.3	72	0.00
36	Benzo(k)fluoranthene	0.400	0.371	7.3	74	0.00
37 C	Benzo(a)pyrene	0.400	0.363	9.3	73	-0.01
38	Dibenzo(a,h)anthracene	0.400	0.346	13.5	71	-0.01
39	Benzo(a,h,i)perylene	0.400	0.370	7.5	74	-0.01

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