

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE010920\
 Data File : BE101010.D
 Acq On : 9 Jan 2020 12:54
 Operator : JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Jan 09 18:11:44 2020
 Quant Method : Z:\svoasrv\HPCHEM1\BNA E\Methods\8270-SIM-BE010820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jan 08 18:36:25 2020
 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	101	0.00
2	1,4-Dioxane	0.400	0.363	9.3	96	0.00
3	n-Nitrosodimethylamine	0.400	0.361	9.8	105	0.00
4 S	2-Fluorophenol	0.400	0.373	6.8	95	0.00
5 S	Phenol-d6	0.400	0.347	13.3	98	0.00
6	bis(2-Chloroethyl)ether	0.400	0.338	15.5	93	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	103	0.01
8 S	Nitrobenzene-d5	0.400	0.334	16.5	96	0.00
9	Naphthalene	0.400	0.411	-2.7	103	0.00
10	Hexachlorobutadiene	0.400	0.427	-6.7	106	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.406	-1.5	104	0.00
12	2-Methylnaphthalene	0.400	0.396	1.0	103	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	99	0.00
14 S	2,4,6-Tribromophenol	0.400	0.350	12.5	100	0.00
15 S	2-Fluorobiphenyl	0.400	0.413	-3.2	101	0.00
16	Acenaphthylene	0.400	0.359	10.3	99	0.00
17	Acenaphthene	0.400	0.399	0.3	102	0.00
18	Fluorene	0.400	0.393	1.8	100	-0.01
19 I	Phenanthrene-d10	0.400	0.400	0.0	98	0.00
20	4-Bromophenyl-phenylether	0.400	0.385	3.8	99	0.00
21	Hexachlorobenzene	0.400	0.414	-3.5	100	0.00
22	Pentachlorophenol	0.400	0.424	-6.0	79	0.00
23	Phenanthrene	0.400	0.394	1.5	98	0.00
24	Anthracene	0.400	0.380	5.0	98	0.00
25 SURR	Fluoranthene-d10	0.400	0.382	4.5	97	0.00
26	Fluoranthene	0.400	0.380	5.0	99	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	95	0.00
28	Pvrene	0.400	0.417	-4.2	96	0.00
29 S	Terphenyl-d14	0.400	0.406	-1.5	95	0.00
30	Benzo(a)anthracene	0.400	0.323	19.3	80	0.00
31	Chrysene	0.400	0.462	-15.5	107	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.328	18.0	89	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.427	-6.7	104	0.00
34 I	Perylene-d12	0.400	0.400	0.0	107	0.00
35	Benzo(b)fluoranthene	0.400	0.332	17.0	95	0.00
36	Benzo(k)fluoranthene	0.400	0.341	14.8	91	0.00
37 C	Benzo(a)pyrene	0.400	0.372	7.0	103	0.00
38	Dibenzo(a,h)anthracene	0.400	0.369	7.8	107	0.00
39	Benzo(a,h,i)perylene	0.400	0.363	9.3	94	-0.02

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

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