

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE062019\
 Data File : BE100101.D
 Acq On : 20 Jun 2019 10:11
 Operator : JU/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Jun 21 01:53:01 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE061919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 19 15:07:24 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	88	-0.01
2	1,4-Dioxane	0.400	0.386	3.5	97	-0.01
3	n-Nitrosodimethylamine	0.400	0.560	-40.0#	292	-0.02
4 S	2-Fluorophenol	0.400	0.402	-0.5	94	-0.01
5 S	Phenol-d6	0.400	0.428	-7.0	94	0.00
6	bis(2-Chloroethyl)ether	0.400	0.455	-13.7	110	-0.01
7 I	Naphthalene-d8	0.400	0.400	0.0	95	0.00
8 S	Nitrobenzene-d5	0.400	0.372	7.0	88	0.01
9	Naphthalene	0.400	0.405	-1.3	97	0.00
10	Hexachlorobutadiene	0.400	0.384	4.0	94	-0.01
11 SURR	2-Methylnaphthalene-d10	0.400	0.407	-1.7	101	0.00
12	2-Methylnaphthalene	0.400	0.403	-0.8	100	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	112	0.00
14 S	2,4,6-Tribromophenol	0.400	0.419	-4.7	132	0.00
15 S	2-Fluorobiphenyl	0.400	0.414	-3.5	107	0.00
16	Acenaphthylene	0.400	0.408	-2.0	113	-0.01
17	Acenaphthene	0.400	0.421	-5.2	115	0.00
18	Fluorene	0.400	0.414	-3.5	114	-0.01
19 I	Phenanthrene-d10	0.400	0.400	0.0	120	0.00
20	4-Bromophenyl-phenylether	0.400	0.406	-1.5	127	0.00
21	Hexachlorobenzene	0.400	0.375	6.3	113	-0.01
22	Pentachlorophenol	0.400	0.502	-25.5#	174	-0.01
23	Phenanthrene	0.400	0.387	3.3	118	0.00
24	Anthracene	0.400	0.392	2.0	127	0.00
25 SURR	Fluoranthene-d10	0.400	0.416	-4.0	127	0.00
26	Fluoranthene	0.400	0.412	-3.0	126	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	127	0.00
28	Pvrene	0.400	0.409	-2.2	131	0.00
29 S	Terphenyl-d14	0.400	0.415	-3.7	132	-0.01
30	Benzo(a)anthracene	0.400	0.410	-2.5	142	0.00
31	Chrysene	0.400	0.402	-0.5	124	-0.01
32	Bis(2-ethylhexyl)phthalate	0.400	0.540	-35.0#	199	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.425	-6.2	130	-0.01
34 I	Perylene-d12	0.400	0.400	0.0	137	-0.01
35	Benzo(b)fluoranthene	0.400	0.364	9.0	126	0.00
36	Benzo(k)fluoranthene	0.400	0.382	4.5	135	-0.01
37 C	Benzo(a)pyrene	0.400	0.394	1.5	135	-0.01
38	Dibenzo(a,h)anthracene	0.400	0.376	6.0	133	0.00
39	Benzo(a,h,i)perylene	0.400	0.372	7.0	129	-0.02

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

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