

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE070119\
 Data File : BE100259.D
 Acq On : 1 Jul 2019 11:15
 Operator : JU/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Jul 01 14:50:24 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE062819.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 28 14:06:53 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	117	0.00
2	1,4-Dioxane	0.400	0.451	-12.7	128	0.00
3	n-Nitrosodimethylamine	0.400	0.593	-48.2#	210	-0.01
4 S	2-Fluorophenol	0.400	0.500	-25.0	145	0.00
5 S	Phenol-d6	0.400	0.464	-16.0	136	-0.01
6	bis(2-Chloroethyl)ether	0.400	0.426	-6.5	112	-0.01
7 I	Naphthalene-d8	0.400	0.400	0.0	109	-0.01
8 S	Nitrobenzene-d5	0.400	0.443	-10.7	122	-0.01
9	Naphthalene	0.400	0.420	-5.0	112	0.00
10	Hexachlorobutadiene	0.400	0.422	-5.5	111	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.424	-6.0	113	-0.01
12	2-Methylnaphthalene	0.400	0.410	-2.5	117	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	117	0.00
14 S	2,4,6-Tribromophenol	0.400	0.609	-52.2#	180	-0.01
15 S	2-Fluorobiphenyl	0.400	0.420	-5.0	119	0.00
16	Acenaphthylene	0.400	0.432	-8.0	121	0.00
17	Acenaphthene	0.400	0.438	-9.5	121	0.00
18	Fluorene	0.400	0.443	-10.7	125	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	125	0.01
20	4-Bromophenyl-phenylether	0.400	0.422	-5.5	126	-0.01
21	Hexachlorobenzene	0.400	0.399	0.3	119	0.01
22	Pentachlorophenol	0.400	0.330	17.5	115	-0.01
23	Phenanthrene	0.400	0.423	-5.7	131	0.00
24	Anthracene	0.400	0.458	-14.5	144	0.00
25 SURR	Fluoranthene-d10	0.400	0.418	-4.5	129	0.00
26	Fluoranthene	0.400	0.418	-4.5	127	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	116	0.00
28	Pvrene	0.400	0.453	-13.2	129	0.00
29 S	Terphenyl-d14	0.400	0.464	-16.0	134	0.00
30	Benzo(a)anthracene	0.400	0.436	-9.0	124	0.01
31	Chrysene	0.400	0.441	-10.2	125	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.415	-3.7	116	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.382	4.5	103	0.02
34 I	Perylene-d12	0.400	0.400	0.0	105	0.01
35	Benzo(b)fluoranthene	0.400	0.432	-8.0	112	0.00
36	Benzo(k)fluoranthene	0.400	0.419	-4.7	106	0.00
37 C	Benzo(a)pyrene	0.400	0.428	-7.0	108	0.01
38	Dibenzo(a,h)anthracene	0.400	0.415	-3.7	108	0.02
39	Benzo(a,h,i)perylene	0.400	0.410	-2.5	103	0.02

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

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