

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE062519\
 Data File : BE100185.D
 Acq On : 25 Jun 2019 18:32
 Operator : JU/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Jun 26 03:21:16 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\8270-SIM-BE062519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 25 15:40:27 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	103	0.00
2	1,4-Dioxane	0.400	0.420	-5.0	99	0.00
3	n-Nitrosodimethylamine	0.400	0.329	17.8	101	-0.03
4 S	2-Fluorophenol	0.400	0.405	-1.3	100	0.00
5 S	Phenol-d6	0.400	0.404	-1.0	110	0.00
6	bis(2-Chloroethyl)ether	0.400	0.296	26.0#	83	-0.03
7 I	Naphthalene-d8	0.400	0.400	0.0	105	0.00
8 S	Nitrobenzene-d5	0.400	0.369	7.8	97	0.00
9	Naphthalene	0.400	0.408	-2.0	102	0.01
10	Hexachlorobutadiene	0.400	0.408	-2.0	100	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.395	1.3	99	0.00
12	2-Methylnaphthalene	0.400	0.395	1.3	97	0.01
13 I	Acenaphthene-d10	0.400	0.400	0.0	97	0.01
14 S	2,4,6-Tribromophenol	0.400	0.412	-3.0	104	0.00
15 S	2-Fluorobiphenyl	0.400	0.427	-6.7	100	0.01
16	Acenaphthylene	0.400	0.415	-3.7	95	0.00
17	Acenaphthene	0.400	0.424	-6.0	95	0.00
18	Fluorene	0.400	0.428	-7.0	96	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	92	0.00
20	4-Bromophenyl-phenylether	0.400	0.427	-6.7	93	0.00
21	Hexachlorobenzene	0.400	0.436	-9.0	93	0.00
22	Pentachlorophenol	0.400	0.584	-46.0#	129	0.00
23	Phenanthrene	0.400	0.434	-8.5	93	0.00
24	Anthracene	0.400	0.415	-3.7	93	0.00
25 SURR	Fluoranthene-d10	0.400	0.423	-5.7	93	0.00
26	Fluoranthene	0.400	0.430	-7.5	95	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	102	0.00
28	Pyrene	0.400	0.422	-5.5	97	0.00
29 S	Terphenyl-d14	0.400	0.411	-2.7	95	0.00
30	Benzo(a)anthracene	0.400	0.427	-6.7	103	0.00
31	Chrysene	0.400	0.424	-6.0	105	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.371	7.3	102	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.465	-16.3	129	0.00
34 I	Perylene-d12	0.400	0.400	0.0	120	0.00
35	Benzo(b)fluoranthene	0.400	0.433	-8.2	123	0.00
36	Benzo(k)fluoranthene	0.400	0.402	-0.5	111	0.00
37 C	Benzo(a)pyrene	0.400	0.434	-8.5	124	0.00
38	Dibenzo(a,h)anthracene	0.400	0.419	-4.7	130	0.00
39	Benzo(g,h,i)perylene	0.400	0.416	-4.0	124	0.00

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

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