

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE062719\
 Data File : BE100206.D
 Acq On : 27 Jun 2019 16:03
 Operator : JU/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Jun 27 19:17:25 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE062519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 27 19:12:41 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	156	0.00
2	1,4-Dioxane	0.400	0.388	3.0	140	0.00
3	n-Nitrosodimethylamine	0.400	0.437	-9.2	203	-0.01
4 S	2-Fluorophenol	0.400	0.486	-21.5	182	0.00
5 S	Phenol-d6	0.400	0.430	-7.5	177	-0.01
6	bis(2-Chloroethyl)ether	0.400	0.424	-6.0	180	-0.01
7 I	Naphthalene-d8	0.400	0.400	0.0	158	0.00
8 S	Nitrobenzene-d5	0.400	0.422	-5.5	167	-0.01
9	Naphthalene	0.400	0.420	-5.0	157	0.00
10	Hexachlorobutadiene	0.400	0.406	-1.5	150	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.412	-3.0	155	-0.01
12	2-Methylnaphthalene	0.400	0.423	-5.7	156	-0.01
13 I	Acenaphthene-d10	0.400	0.400	0.0	171	0.00
14 S	2,4,6-Tribromophenol	0.400	0.386	3.5	171	-0.01
15 S	2-Fluorobiphenyl	0.400	0.384	4.0	159	-0.01
16	Acenaphthylene	0.400	0.388	3.0	156	0.00
17	Acenaphthene	0.400	0.407	-1.7	162	0.00
18	Fluorene	0.400	0.414	-3.5	164	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	165	0.00
20	4-Bromophenyl-phenylether	0.400	0.401	-0.3	158	0.00
21	Hexachlorobenzene	0.400	0.412	-3.0	159	0.00
22	Pentachlorophenol	0.400	0.531	-32.8#	193	0.01
23	Phenanthrene	0.400	0.424	-6.0	163	0.00
24	Anthracene	0.400	0.436	-9.0	177	0.00
25 SURR	Fluoranthene-d10	0.400	0.391	2.3	155	0.00
26	Fluoranthene	0.400	0.396	1.0	158	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	154	0.00
28	Pvrene	0.400	0.465	-16.3	161	0.00
29 S	Terphenyl-d14	0.400	0.444	-11.0	155	0.00
30	Benzo(a)anthracene	0.400	0.444	-11.0	162	0.00
31	Chrysene	0.400	0.402	-0.5	149	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.600	-50.0#	247	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.315	21.3	132	0.00
34 I	Perylene-d12	0.400	0.400	0.0	140	0.00
35	Benzo(b)fluoranthene	0.400	0.425	-6.2	141	0.00
36	Benzo(k)fluoranthene	0.400	0.392	2.0	126	0.00
37 C	Benzo(a)pyrene	0.400	0.405	-1.3	135	0.00
38	Dibenzo(a,h)anthracene	0.400	0.360	10.0	130	0.00
39	Benzo(a,h,i)perylene	0.400	0.387	3.3	135	-0.01

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

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