

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE070119\
 Data File : BE100278.D
 Acq On : 1 Jul 2019 23:56
 Operator : JU/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Jul 02 02:01:51 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	164	0.00
2	1,4-Dioxane	0.400	0.321	19.8	128	0.01
3	n-Nitrosodimethylamine	0.400	0.431	-7.7	215	0.00
4 S	2-Fluorophenol	0.400	0.428	-7.0	175	0.01
5 S	Phenol-d6	0.400	0.438	-9.5	181	-0.01
6	bis(2-Chloroethyl)ether	0.400	0.447	-11.7	166	-0.01
7 I	Naphthalene-d8	0.400	0.400	0.0	160	-0.01
8 S	Nitrobenzene-d5	0.400	0.459	-14.8	186	-0.01
9	Naphthalene	0.400	0.419	-4.7	164	0.00
10	Hexachlorobutadiene	0.400	0.444	-11.0	172	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.434	-8.5	170	-0.01
12	2-Methylnaphthalene	0.400	0.413	-3.2	173	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	161	0.00
14 S	2,4,6-Tribromophenol	0.400	0.395	1.3	161	0.00
15 S	2-Fluorobiphenyl	0.400	0.428	-7.0	167	0.00
16	Acenaphthylene	0.400	0.428	-7.0	166	0.00
17	Acenaphthene	0.400	0.428	-7.0	163	0.00
18	Fluorene	0.400	0.402	-0.5	156	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	137	0.01
20	4-Bromophenyl-phenylether	0.400	0.448	-12.0	147	0.00
21	Hexachlorobenzene	0.400	0.455	-13.7	149	0.01
22	Pentachlorophenol	0.400	0.307	23.3	118	-0.03
23	Phenanthrene	0.400	0.431	-7.7	147	0.00
24	Anthracene	0.400	0.439	-9.7	153	0.00
25 SURR	Fluoranthene-d10	0.400	0.386	3.5	131	0.00
26	Fluoranthene	0.400	0.398	0.5	133	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	112	0.00
28	Pvrene	0.400	0.492	-23.0	135	0.00
29 S	Terphenyl-d14	0.400	0.494	-23.5	137	0.00
30	Benzo(a)anthracene	0.400	0.438	-9.5	120	0.01
31	Chrysene	0.400	0.460	-15.0	125	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.388	3.0	105	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.406	-1.5	105	0.02
34 I	Perylene-d12	0.400	0.400	0.0	101	0.01
35	Benzo(b)fluoranthene	0.400	0.437	-9.2	109	0.00
36	Benzo(k)fluoranthene	0.400	0.423	-5.7	103	0.00
37 C	Benzo(a)pyrene	0.400	0.432	-8.0	105	0.01
38	Dibenzo(a,h)anthracene	0.400	0.428	-7.0	108	0.01
39	Benzo(a,h,i)perylene	0.400	0.428	-7.0	104	0.01

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

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