

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE062819\
 Data File : BE100238.D
 Acq On : 28 Jun 2019 17:14
 Operator : JU/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Jun 29 00:27:32 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	80	0.00
2	1,4-Dioxane	0.400	0.421	-5.2	81	0.00
3	n-Nitrosodimethylamine	0.400	0.440	-10.0	106	0.01
4 S	2-Fluorophenol	0.400	0.397	0.8	79	0.00
5 S	Phenol-d6	0.400	0.343	14.2	69	0.00
6	bis(2-Chloroethyl)ether	0.400	0.439	-9.7	79	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	75	0.00
8 S	Nitrobenzene-d5	0.400	0.402	-0.5	76	0.00
9	Naphthalene	0.400	0.421	-5.2	77	0.00
10	Hexachlorobutadiene	0.400	0.460	-15.0	83	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.427	-6.7	78	0.00
12	2-Methylnaphthalene	0.400	0.388	3.0	76	0.01
13 I	Acenaphthene-d10	0.400	0.400	0.0	79	0.01
14 S	2,4,6-Tribromophenol	0.400	0.373	6.8	75	0.00
15 S	2-Fluorobiphenyl	0.400	0.415	-3.7	80	0.01
16	Acenaphthylene	0.400	0.426	-6.5	82	0.00
17	Acenaphthene	0.400	0.418	-4.5	79	0.00
18	Fluorene	0.400	0.418	-4.5	80	0.01
19 I	Phenanthrene-d10	0.400	0.400	0.0	77	0.01
20	4-Bromophenyl-phenylether	0.400	0.433	-8.2	80	0.00
21	Hexachlorobenzene	0.400	0.439	-9.7	81	0.00
22	Pentachlorophenol	0.400	0.274	31.5#	59	0.01
23	Phenanthrene	0.400	0.415	-3.7	79	0.00
24	Anthracene	0.400	0.409	-2.2	80	0.00
25 SURR	Fluoranthene-d10	0.400	0.414	-3.5	79	0.00
26	Fluoranthene	0.400	0.425	-6.2	80	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	80	0.00
28	Pvrene	0.400	0.395	1.3	77	0.00
29 S	Terphenyl-d14	0.400	0.393	1.8	78	0.00
30	Benzo(a)anthracene	0.400	0.394	1.5	77	0.01
31	Chrysene	0.400	0.428	-7.0	83	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.297	25.8#	60	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.413	-3.2	76	0.00
34 I	Perylene-d12	0.400	0.400	0.0	74	0.00
35	Benzo(b)fluoranthene	0.400	0.414	-3.5	76	0.00
36	Benzo(k)fluoranthene	0.400	0.427	-6.7	76	0.00
37 C	Benzo(a)pyrene	0.400	0.439	-9.7	78	0.00
38	Dibenzo(a,h)anthracene	0.400	0.418	-4.5	77	0.00
39	Benzo(a,h,i)perylene	0.400	0.432	-8.0	77	0.00

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

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