

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE070319\
 Data File : BE100323.D
 Acq On : 3 Jul 2019 11:46
 Operator : JU/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Jul 05 01:34:27 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	73	0.00
2	1,4-Dioxane	0.400	0.468	-17.0	83	-0.01
3	n-Nitrosodimethylamine	0.400	0.466	-16.5	104	0.00
4 S	2-Fluorophenol	0.400	0.185	53.8#	34	0.00
5 S	Phenol-d6	0.400	0.356	11.0	65	0.02
6	bis(2-Chloroethyl)ether	0.400	0.376	6.0	62	0.01
7 I	Naphthalene-d8	0.400	0.400	0.0	72	0.00
8 S	Nitrobenzene-d5	0.400	0.335	16.3	62	0.01
9	Naphthalene	0.400	0.428	-7.0	76	0.00
10	Hexachlorobutadiene	0.400	0.512	-28.0#	90	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.436	-9.0	78	0.00
12	2-Methylnaphthalene	0.400	0.402	-0.5	76	0.02
13 I	Acenaphthene-d10	0.400	0.400	0.0	82	0.00
14 S	2,4,6-Tribromophenol	0.400	0.092	77.0#	19	0.08
15 S	2-Fluorobiphenyl	0.400	0.414	-3.5	82	0.02
16	Acenaphthylene	0.400	0.423	-5.7	84	0.00
17	Acenaphthene	0.400	0.440	-10.0	85	0.00
18	Fluorene	0.400	0.433	-8.2	86	0.01
19 I	Phenanthrene-d10	0.400	0.400	0.0	81	0.01
20	4-Bromophenyl-phenylether	0.400	0.449	-12.2	87	0.00
21	Hexachlorobenzene	0.400	0.450	-12.5	87	0.00
22	Pentachlorophenol	0.400	0.002	99.5#	1	0.00
23	Phenanthrene	0.400	0.411	-2.7	83	0.00
24	Anthracene	0.400	0.374	6.5	77	0.00
25 SURR	Fluoranthene-d10	0.400	0.440	-10.0	88	0.00
26	Fluoranthene	0.400	0.441	-10.2	87	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	71	0.01
28	Pvrene	0.400	0.515	-28.7#	90	0.00
29 S	Terphenyl-d14	0.400	0.542	-35.5#	96	0.01
30	Benzo(a)anthracene	0.400	0.378	5.5	66	0.01
31	Chrysene	0.400	0.396	1.0	69	0.01
32	Bis(2-ethylhexyl)phthalate	0.400	0.459	-14.8	78	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.363	9.3	60	0.02
34 I	Perylene-d12	0.400	0.400	0.0	68	0.02
35	Benzo(b)fluoranthene	0.400	0.339	15.3	57	0.00
36	Benzo(k)fluoranthene	0.400	0.415	-3.7	68	0.01
37 C	Benzo(a)pyrene	0.400	0.406	-1.5	67	0.02
38	Dibenzo(a,h)anthracene	0.400	0.344	14.0	58	0.02
39	Benzo(a,h,i)perylene	0.400	0.415	-3.7	68	0.02

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

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