

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE052919\
 Data File : BE099772.D
 Acq On : 29 May 2019 12:32
 Operator : JU/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: May 29 17:25:15 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE052019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon May 20 19:29:40 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	87	0.00
2	1,4-Dioxane	0.400	0.371	7.3	84	-0.02
3	n-Nitrosodimethylamine	0.400	0.356	11.0	88	-0.02
4 S	2-Fluorophenol	0.400	0.368	8.0	91	0.00
5 S	Phenol-d6	0.400	0.359	10.3	94	0.00
6	bis(2-Chloroethyl)ether	0.400	0.346	13.5	86	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	85	0.00
8 S	Nitrobenzene-d5	0.400	0.384	4.0	95	-0.01
9	Naphthalene	0.400	0.365	8.8	86	-0.01
10	Hexachlorobutadiene	0.400	0.366	8.5	87	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.360	10.0	85	-0.02
12	2-Methylnaphthalene	0.400	0.357	10.8	87	-0.02
13 I	Acenaphthene-d10	0.400	0.400	0.0	86	0.00
14 S	2,4,6-Tribromophenol	0.400	0.401	-0.3	96	-0.01
15 S	2-Fluorobiphenyl	0.400	0.353	11.8	85	0.00
16	Acenaphthylene	0.400	0.359	10.3	90	-0.02
17	Acenaphthene	0.400	0.351	12.3	85	0.00
18	Fluorene	0.400	0.352	12.0	86	-0.01
19 I	Phenanthrene-d10	0.400	0.400	0.0	83	0.00
20	4-Bromophenyl-phenylether	0.400	0.344	14.0	82	-0.01
21	Hexachlorobenzene	0.400	0.340	15.0	80	-0.01
22	Pentachlorophenol	0.400	0.416	-4.0	90	-0.01
23	Phenanthrene	0.400	0.353	11.8	83	0.00
24	Anthracene	0.400	0.344	14.0	85	0.00
25 SURR	Fluoranthene-d10	0.400	0.352	12.0	84	-0.01
26	Fluoranthene	0.400	0.355	11.3	84	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	88	0.00
28	Pvrene	0.400	0.386	3.5	86	-0.01
29 S	Terphenyl-d14	0.400	0.366	8.5	83	-0.01
30	Benzo(a)anthracene	0.400	0.408	-2.0	94	-0.01
31	Chrysene	0.400	0.376	6.0	83	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.448	-12.0	109	-0.01
33	Indeno(1,2,3-cd)pyrene	0.400	0.386	3.5	90	-0.02
34 I	Perylene-d12	0.400	0.400	0.0	87	-0.01
35	Benzo(b)fluoranthene	0.400	0.374	6.5	91	-0.01
36	Benzo(k)fluoranthene	0.400	0.349	12.8	85	-0.01
37 C	Benzo(a)pyrene	0.400	0.359	10.3	89	-0.01
38	Dibenzo(a,h)anthracene	0.400	0.357	10.8	90	-0.02
39	Benzo(a,h,i)perylene	0.400	0.361	9.8	89	-0.01

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

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