

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE051019\
 Data File : BE099603.D
 Acq On : 10 May 2019 17:42
 Operator : JU/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: May 11 01:01:41 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE050919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu May 09 13:32:58 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	119	0.00
2	1,4-Dioxane	0.400	0.356	11.0	124	0.01
3	n-Nitrosodimethylamine	0.400	0.332	17.0	118	0.00
4 S	2-Fluorophenol	0.400	0.404	-1.0	139	0.00
5 S	Phenol-d6	0.400	0.401	-0.3	136	0.00
6	bis(2-Chloroethyl)ether	0.400	0.417	-4.2	141	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	118	0.00
8 S	Nitrobenzene-d5	0.400	0.412	-3.0	138	0.00
9	Naphthalene	0.400	0.417	-4.2	138	0.00
10	Hexachlorobutadiene	0.400	0.419	-4.7	137	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.405	-1.3	131	0.00
12	2-Methylnaphthalene	0.400	0.412	-3.0	136	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	119	0.00
14 S	2,4,6-Tribromophenol	0.400	0.409	-2.2	140	0.00
15 S	2-Fluorobiphenyl	0.400	0.409	-2.2	136	0.00
16	Acenaphthylene	0.400	0.398	0.5	134	0.00
17	Acenaphthene	0.400	0.409	-2.2	135	0.00
18	Fluorene	0.400	0.413	-3.2	139	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	118	0.00
20	4-Bromophenyl-phenylether	0.400	0.409	-2.2	138	0.00
21	Hexachlorobenzene	0.400	0.426	-6.5	138	0.00
22	Pentachlorophenol	0.400	0.511	-27.7#	146	0.00
23	Phenanthrene	0.400	0.424	-6.0	140	0.00
24	Anthracene	0.400	0.409	-2.2	140	0.00
25 SURR	Fluoranthene-d10	0.400	0.402	-0.5	133	0.00
26	Fluoranthene	0.400	0.420	-5.0	138	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	118	0.00
28	Pvrene	0.400	0.458	-14.5	137	0.00
29 S	Terphenyl-d14	0.400	0.467	-16.8	140	0.00
30	Benzo(a)anthracene	0.400	0.460	-15.0	137	0.00
31	Chrysene	0.400	0.448	-12.0	133	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.424	-6.0	131	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.449	-12.2	138	0.00
34 I	Perylene-d12	0.400	0.400	0.0	119	0.00
35	Benzo(b)fluoranthene	0.400	0.423	-5.7	138	0.00
36	Benzo(k)fluoranthene	0.400	0.429	-7.2	138	0.00
37 C	Benzo(a)pyrene	0.400	0.425	-6.2	139	0.00
38	Dibenzo(a,h)anthracene	0.400	0.419	-4.7	138	0.00
39	Benzo(a,h,i)perylene	0.400	0.413	-3.2	135	0.00

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

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