

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE061419\
 Data File : BE100012.D
 Acq On : 14 Jun 2019 8:03
 Operator : JU/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Jun 15 00:58:56 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE061019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 10 14:57:20 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	0.392	2.0	124	0.00
3	n-Nitrosodimethylamine	0.400	0.238	40.5#	70	0.03
4 S	2-Fluorophenol	0.400	0.374	6.5	111	0.00
5 S	Phenol-d6	0.400	0.374	6.5	119	0.00
6	bis(2-Chloroethyl)ether	0.400	0.396	1.0	124	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	114	0.00
8 S	Nitrobenzene-d5	0.400	0.387	3.3	117	0.00
9	Naphthalene	0.400	0.415	-3.7	120	-0.01
10	Hexachlorobutadiene	0.400	0.466	-16.5	131	-0.01
11 SURR	2-Methylnaphthalene-d10	0.400	0.416	-4.0	121	0.00
12	2-Methylnaphthalene	0.400	0.406	-1.5	117	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	107	0.00
14 S	2,4,6-Tribromophenol	0.400	0.258	35.5#	76	0.00
15 S	2-Fluorobiphenyl	0.400	0.431	-7.7	122	0.00
16	Acenaphthylene	0.400	0.412	-3.0	114	-0.01
17	Acenaphthene	0.400	0.409	-2.2	114	0.00
18	Fluorene	0.400	0.396	1.0	112	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	96	0.00
20	4-Bromophenyl-phenylether	0.400	0.428	-7.0	112	0.00
21	Hexachlorobenzene	0.400	0.431	-7.7	110	0.00
22	Pentachlorophenol	0.400	0.449	-12.2	63	0.00
23	Phenanthrene	0.400	0.406	-1.5	105	0.00
24	Anthracene	0.400	0.368	8.0	94	0.00
25 SURR	Fluoranthene-d10	0.400	0.366	8.5	90	0.00
26	Fluoranthene	0.400	0.375	6.3	92	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	84	0.00
28	Pvrene	0.400	0.462	-15.5	91	0.00
29 S	Terphenyl-d14	0.400	0.428	-7.0	86	0.00
30	Benzo(a)anthracene	0.400	0.417	-4.2	84	0.01
31	Chrysene	0.400	0.464	-16.0	90	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.329	17.8	68	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.419	-4.7	83	0.00
34 I	Perylene-d12	0.400	0.400	0.0	82	0.00
35	Benzo(b)fluoranthene	0.400	0.400	0.0	86	0.00
36	Benzo(k)fluoranthene	0.400	0.432	-8.0	88	0.00
37 C	Benzo(a)pyrene	0.400	0.413	-3.2	87	0.00
38	Dibenzo(a,h)anthracene	0.400	0.370	7.5	79	0.00
39	Benzo(a,h,i)perylene	0.400	0.409	-2.2	86	0.00

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

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