

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE061319\
 Data File : BE099991.D
 Acq On : 13 Jun 2019 15:59
 Operator : JU/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Jun 14 04:09:02 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	-0.01
2	1,4-Dioxane	0.400	0.343	14.2	108	0.00
3	n-Nitrosodimethylamine	0.400	0.413	-3.2	121	0.00
4 S	2-Fluorophenol	0.400	0.396	1.0	117	0.00
5 S	Phenol-d6	0.400	0.415	-3.7	132	0.00
6	bis(2-Chloroethyl)ether	0.400	0.397	0.8	124	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	115	-0.01
8 S	Nitrobenzene-d5	0.400	0.415	-3.7	127	0.00
9	Naphthalene	0.400	0.425	-6.2	123	-0.01
10	Hexachlorobutadiene	0.400	0.463	-15.8	130	-0.01
11 SURR	2-Methylnaphthalene-d10	0.400	0.437	-9.2	128	-0.01
12	2-Methylnaphthalene	0.400	0.429	-7.2	125	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	118	0.00
14 S	2,4,6-Tribromophenol	0.400	0.400	0.0	131	-0.01
15 S	2-Fluorobiphenyl	0.400	0.411	-2.7	128	-0.01
16	Acenaphthylene	0.400	0.407	-1.7	125	-0.01
17	Acenaphthene	0.400	0.396	1.0	122	0.00
18	Fluorene	0.400	0.415	-3.7	129	-0.01
19 I	Phenanthrene-d10	0.400	0.400	0.0	121	0.00
20	4-Bromophenyl-phenylether	0.400	0.398	0.5	131	-0.01
21	Hexachlorobenzene	0.400	0.427	-6.7	136	-0.01
22	Pentachlorophenol	0.400	0.609	-52.2#	163	-0.03
23	Phenanthrene	0.400	0.397	0.8	129	0.00
24	Anthracene	0.400	0.393	1.8	126	0.00
25 SURR	Fluoranthene-d10	0.400	0.377	5.8	117	0.00
26	Fluoranthene	0.400	0.379	5.3	117	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	107	0.00
28	Pvrene	0.400	0.465	-16.3	117	0.00
29 S	Terphenyl-d14	0.400	0.461	-15.3	118	0.00
30	Benzo(a)anthracene	0.400	0.440	-10.0	114	0.00
31	Chrysene	0.400	0.468	-17.0	115	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.436	-9.0	115	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.464	-16.0	118	-0.01
34 I	Perylene-d12	0.400	0.400	0.0	107	0.00
35	Benzo(b)fluoranthene	0.400	0.407	-1.7	114	0.00
36	Benzo(k)fluoranthene	0.400	0.409	-2.2	109	0.00
37 C	Benzo(a)pyrene	0.400	0.408	-2.0	112	0.00
38	Dibenzo(a,h)anthracene	0.400	0.421	-5.2	118	-0.01
39	Benzo(a,h,i)perylene	0.400	0.415	-3.7	113	0.00

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

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