

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE092619\
 Data File : BE100565.D
 Acq On : 26 Sep 2019 12:43
 Operator : JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Sep 26 15:24:28 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE092019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 20 18:26:17 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	95	0.00
2	1,4-Dioxane	0.400	0.414	-3.5	97	0.03
3	n-Nitrosodimethylamine	0.400	0.463	-15.8	145	0.02
4 S	2-Fluorophenol	0.400	0.553	-38.3#	142	0.02
5 S	Phenol-d6	0.400	0.509	-27.2#	134	0.00
6	bis(2-Chloroethyl)ether	0.400	0.400	0.0	100	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	94	0.00
8 S	Nitrobenzene-d5	0.400	0.524	-31.0#	137	0.00
9	Naphthalene	0.400	0.397	0.8	96	0.00
10	Hexachlorobutadiene	0.400	0.390	2.5	93	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.387	3.3	96	0.00
12	2-Methylnaphthalene	0.400	0.399	0.3	99	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	89	0.00
14 S	2,4,6-Tribromophenol	0.400	0.654	-63.5#	161	0.00
15 S	2-Fluorobiphenyl	0.400	0.404	-1.0	92	0.00
16	Acenaphthylene	0.400	0.442	-10.5	108	0.01
17	Acenaphthene	0.400	0.411	-2.7	97	0.00
18	Fluorene	0.400	0.417	-4.2	99	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	84	0.00
20	4-Bromophenyl-phenylether	0.400	0.414	-3.5	91	0.00
21	Hexachlorobenzene	0.400	0.384	4.0	85	0.01
22	Pentachlorophenol	0.400	0.795	-98.8#	242	-0.01
23	Phenanthrene	0.400	0.409	-2.2	90	0.00
24	Anthracene	0.400	0.438	-9.5	103	0.00
25 SURR	Fluoranthene-d10	0.400	0.385	3.8	86	0.00
26	Fluoranthene	0.400	0.391	2.3	87	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	79	0.00
28	Pvrene	0.400	0.445	-11.2	91	0.00
29 S	Terphenyl-d14	0.400	0.420	-5.0	86	0.00
30	Benzo(a)anthracene	0.400	0.458	-14.5	95	0.01
31	Chrysene	0.400	0.394	1.5	82	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	1.185	-196.3#	260	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.438	-9.5	89	0.00
34 I	Perylene-d12	0.400	0.400	0.0	83	0.00
35	Benzo(b)fluoranthene	0.400	0.403	-0.8	93	0.00
36	Benzo(k)fluoranthene	0.400	0.359	10.3	80	0.00
37 C	Benzo(a)pyrene	0.400	0.408	-2.0	91	0.00
38	Dibenzo(a,h)anthracene	0.400	0.399	0.3	90	0.00
39	Benzo(a,h,i)perylene	0.400	0.381	4.8	83	0.01

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

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