

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE052819\
 Data File : BE099770.D
 Acq On : 28 May 2019 23:09
 Operator : JU/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: May 29 04:13:04 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	77	0.00
2	1,4-Dioxane	0.400	0.325	18.8	65	-0.02
3	n-Nitrosodimethylamine	0.400	0.360	10.0	79	-0.02
4 S	2-Fluorophenol	0.400	0.400	0.0	88	0.00
5 S	Phenol-d6	0.400	0.397	0.8	92	0.00
6	bis(2-Chloroethyl)ether	0.400	0.343	14.2	75	-0.02
7 I	Naphthalene-d8	0.400	0.400	0.0	79	-0.01
8 S	Nitrobenzene-d5	0.400	0.362	9.5	84	-0.01
9	Naphthalene	0.400	0.357	10.8	79	-0.01
10	Hexachlorobutadiene	0.400	0.354	11.5	79	-0.01
11 SURR	2-Methylnaphthalene-d10	0.400	0.365	8.8	81	-0.01
12	2-Methylnaphthalene	0.400	0.368	8.0	83	-0.01
13 I	Acenaphthene-d10	0.400	0.400	0.0	87	-0.01
14 S	2,4,6-Tribromophenol	0.400	0.454	-13.5	117	-0.01
15 S	2-Fluorobiphenyl	0.400	0.335	16.3	82	0.00
16	Acenaphthylene	0.400	0.366	8.5	92	-0.01
17	Acenaphthene	0.400	0.360	10.0	88	0.00
18	Fluorene	0.400	0.372	7.0	92	-0.01
19 I	Phenanthrene-d10	0.400	0.400	0.0	91	0.00
20	4-Bromophenyl-phenylether	0.400	0.341	14.8	89	-0.01
21	Hexachlorobenzene	0.400	0.337	15.8	87	-0.01
22	Pentachlorophenol	0.400	0.484	-21.0	129	-0.01
23	Phenanthrene	0.400	0.361	9.8	93	0.00
24	Anthracene	0.400	0.361	9.8	98	0.00
25 SURR	Fluoranthene-d10	0.400	0.358	10.5	93	0.00
26	Fluoranthene	0.400	0.360	10.0	93	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	92	-0.01
28	Pvrene	0.400	0.404	-1.0	95	-0.01
29 S	Terphenyl-d14	0.400	0.391	2.3	94	-0.01
30	Benzo(a)anthracene	0.400	0.403	-0.8	97	-0.01
31	Chrysene	0.400	0.393	1.8	92	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.522	-30.5#	134	-0.01
33	Indeno(1,2,3-cd)pyrene	0.400	0.381	4.8	93	-0.03
34 I	Perylene-d12	0.400	0.400	0.0	91	-0.02
35	Benzo(b)fluoranthene	0.400	0.368	8.0	94	-0.01
36	Benzo(k)fluoranthene	0.400	0.347	13.3	89	-0.02
37 C	Benzo(a)pyrene	0.400	0.365	8.8	95	-0.02
38	Dibenzo(a,h)anthracene	0.400	0.349	12.8	92	-0.03
39	Benzo(a,h,i)perylene	0.400	0.355	11.3	92	-0.02

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

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