

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE061119\
 Data File : BE099944.D
 Acq On : 12 Jun 2019 3:09
 Operator : JU/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Jun 12 05:29:11 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	102	-0.02
2	1,4-Dioxane	0.400	0.359	10.3	97	0.00
3	n-Nitrosodimethylamine	0.400	0.217	45.8#	55	0.02
4 S	2-Fluorophenol	0.400	0.364	9.0	93	0.00
5 S	Phenol-d6	0.400	0.386	3.5	106	0.00
6	bis(2-Chloroethyl)ether	0.400	0.344	14.0	92	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	108	-0.01
8 S	Nitrobenzene-d5	0.400	0.344	14.0	99	0.00
9	Naphthalene	0.400	0.403	-0.8	111	-0.03
10	Hexachlorobutadiene	0.400	0.426	-6.5	113	-0.03
11 SURR	2-Methylnaphthalene-d10	0.400	0.406	-1.5	113	-0.01
12	2-Methylnaphthalene	0.400	0.400	0.0	110	-0.01
13 I	Acenaphthene-d10	0.400	0.400	0.0	122	-0.03
14 S	2,4,6-Tribromophenol	0.400	0.270	32.5#	91	-0.03
15 S	2-Fluorobiphenyl	0.400	0.368	8.0	119	-0.03
16	Acenaphthylene	0.400	0.384	4.0	121	-0.04
17	Acenaphthene	0.400	0.381	4.8	121	-0.03
18	Fluorene	0.400	0.406	-1.5	131	-0.04
19 I	Phenanthrene-d10	0.400	0.400	0.0	126	-0.04
20	4-Bromophenyl-phenylether	0.400	0.387	3.3	133	-0.04
21	Hexachlorobenzene	0.400	0.397	0.8	132	-0.04
22	Pentachlorophenol	0.400	0.392	2.0	51	0.00
23	Phenanthrene	0.400	0.375	6.3	127	-0.03
24	Anthracene	0.400	0.330	17.5	110	-0.03
25 SURR	Fluoranthene-d10	0.400	0.339	15.3	110	-0.04
26	Fluoranthene	0.400	0.351	12.3	113	-0.04
27 I	Chrysene-d12	0.400	0.400	0.0	99	-0.05
28	Pvrene	0.400	0.473	-18.2	110	-0.04
29 S	Terphenyl-d14	0.400	0.431	-7.7	102	-0.03
30	Benzo(a)anthracene	0.400	0.366	8.5	88	-0.04
31	Chrysene	0.400	0.457	-14.2	105	-0.04
32	Bis(2-ethylhexyl)phthalate	0.400	0.336	16.0	82	-0.04
33	Indeno(1,2,3-cd)pyrene	0.400	0.413	-3.2	97	-0.11
34 I	Perylene-d12	0.400	0.400	0.0	99	-0.07
35	Benzo(b)fluoranthene	0.400	0.394	1.5	102	-0.06
36	Benzo(k)fluoranthene	0.400	0.416	-4.0	103	-0.06
37 C	Benzo(a)pyrene	0.400	0.396	1.0	101	-0.07
38	Dibenzo(a,h)anthracene	0.400	0.344	14.0	89	-0.09
39	Benzo(a,h,i)perylene	0.400	0.391	2.3	99	-0.12

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

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