

Data Path : Z:\svoasrv\HPCHEM1\BNA E\Data\BE100818\
 Data File : Be097800.d
 Acq On : 8 Oct 2018 19:19
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Oct 09 06:58:40 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE100318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 08 06:37:06 2018
 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	71	-0.01
2	1,4-Dioxane	0.400	0.379	5.3	72	0.00
3	n-Nitrosodimethylamine	0.400	0.358	10.5	67	0.00
4 S	2-Fluorophenol	0.400	0.376	6.0	70	-0.01
5 S	Phenol-d6	0.400	0.351	12.3	64	-0.01
6	bis(2-Chloroethyl)ether	0.400	0.384	4.0	71	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	75	0.00
8 S	Nitrobenzene-d5	0.400	0.622	-55.5#	124	-0.01
9	Naphthalene	0.400	0.384	4.0	73	-0.01
10	Hexachlorobutadiene	0.400	0.389	2.8	75	-0.01
11 SURR	2-Methylnaphthalene-d10	0.400	0.388	3.0	74	-0.02
12	2-Methylnaphthalene	0.400	0.384	4.0	75	-0.02
13 I	Acenaphthene-d10	0.400	0.400	0.0	80	-0.02
14 S	2,4,6-Tribromophenol	0.400	0.431	-7.7	107	-0.01
15 S	2-Fluorobiphenyl	0.400	0.359	10.3	76	-0.02
16	Acenaphthylene	0.400	0.347	13.3	73	0.00
17	Acenaphthene	0.400	0.372	7.0	79	0.00
18	Fluorene	0.400	0.360	10.0	76	-0.01
19 I	Phenanthrene-d10	0.400	0.400	0.0	78	-0.01
20	4-Bromophenyl-phenylether	0.400	0.378	5.5	79	-0.01
21	Hexachlorobenzene	0.400	0.368	8.0	75	0.00
22	Pentachlorophenol	0.400	0.451	-12.7	99	-0.01
23	Phenanthrene	0.400	0.372	7.0	76	0.00
24	Anthracene	0.400	0.349	12.8	74	-0.01
25 SURR	Fluoranthene-d10	0.400	0.364	9.0	76	-0.01
26	Fluoranthene	0.400	0.365	8.8	76	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	68	-0.01
28	Pvrene	0.400	0.404	-1.0	77	-0.01
29 S	Terphenyl-d14	0.400	0.401	-0.3	77	-0.01
30	Benzo(a)anthracene	0.400	0.363	9.3	69	-0.01
31	Chrysene	0.400	0.387	3.3	72	-0.01
32	Bis(2-ethylhexyl)phthalate	0.400	0.365	8.8	72	-0.01
33	Indeno(1,2,3-cd)pyrene	0.400	0.338	15.5	63	-0.03
34 I	Perylene-d12	0.400	0.400	0.0	68	-0.02
35	Benzo(b)fluoranthene	0.400	0.352	12.0	67	-0.02
36	Benzo(k)fluoranthene	0.400	0.390	2.5	74	-0.02
37 C	Benzo(a)pyrene	0.400	0.340	15.0	65	-0.02
38	Dibenzo(a,h)anthracene	0.400	0.346	13.5	66	-0.03
39	Benzo(a,h,i)perylene	0.400	0.360	10.0	67	-0.03

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

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