

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE072618\
 Data File : BE097101.D
 Acq On : 26 Jul 2018 12:44
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Jul 27 03:35:03 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\8270-SIM-BE072418.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 24 13:24:10 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	76	0.00
2	1,4-Dioxane	0.400	0.457	-14.2	89	0.00
3	n-Nitrosodimethylamine	0.400	0.422	-5.5	86	0.00
4 S	2-Fluorophenol	0.400	0.409	-2.2	82	0.00
5 S	Phenol-d6	0.400	0.389	2.8	79	0.00
6	bis(2-Chloroethyl)ether	0.400	0.451	-12.7	92	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	73	0.00
8 S	Nitrobenzene-d5	0.400	0.448	-12.0	86	0.00
9	Naphthalene	0.400	0.462	-15.5	88	0.00
10	Hexachlorobutadiene	0.400	0.465	-16.3	88	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.443	-10.7	84	0.00
12	2-Methylnaphthalene	0.400	0.446	-11.5	85	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	74	0.00
14 S	2,4,6-Tribromophenol	0.400	0.380	5.0	77	0.00
15 S	2-Fluorobiphenyl	0.400	0.446	-11.5	84	0.00
16	Acenaphthylene	0.400	0.444	-11.0	84	0.00
17	Acenaphthene	0.400	0.433	-8.2	82	0.00
18	Fluorene	0.400	0.436	-9.0	84	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	74	0.00
20	4-Bromophenyl-phenylether	0.400	0.442	-10.5	86	0.00
21	Hexachlorobenzene	0.400	0.467	-16.8	88	0.00
22	Pentachlorophenol	0.400	0.380	5.0	74	0.00
23	Phenanthrene	0.400	0.447	-11.7	87	0.00
24	Anthracene	0.400	0.431	-7.7	86	0.00
25 SURR	Fluoranthene-d10	0.400	0.439	-9.7	89	0.00
26	Fluoranthene	0.400	0.447	-11.7	89	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	73	0.00
28	Pyrene	0.400	0.474	-18.5	90	0.00
29 S	Terphenyl-d14	0.400	0.454	-13.5	88	0.00
30	Benzo(a)anthracene	0.400	0.405	-1.3	82	0.00
31	Chrysene	0.400	0.465	-16.3	87	-0.01
32	Bis(2-ethylhexyl)phthalate	0.400	0.360	10.0	67	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.403	-0.8	79	0.00
34 I	Perylene-d12	0.400	0.400	0.0	75	0.00
35	Benzo(b)fluoranthene	0.400	0.398	0.5	79	0.00
36	Benzo(k)fluoranthene	0.400	0.433	-8.2	88	0.00
37 C	Benzo(a)pyrene	0.400	0.409	-2.2	82	0.00
38	Dibenzo(a,h)anthracene	0.400	0.388	3.0	77	0.00
39	Benzo(g,h,i)perylene	0.400	0.400	0.0	79	0.00

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