

Data Path : Z:\svoasrv\HPCHEM1\BNA E\Data\BE100418\
 Data File : BE097726.D
 Acq On : 4 Oct 2018 15:31
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Oct 05 06:50:21 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE100318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 03 12:17:19 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	59	0.00
2	1,4-Dioxane	0.400	0.369	7.8	59	0.00
3	n-Nitrosodimethylamine	0.400	0.361	9.8	56	0.00
4 S	2-Fluorophenol	0.400	0.442	-10.5	69	0.00
5 S	Phenol-d6	0.400	0.489	-22.2	74	0.00
6	bis(2-Chloroethyl)ether	0.400	0.370	7.5	57	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	59	0.00
8 S	Nitrobenzene-d5	0.400	0.513	-28.2#	81	0.00
9	Naphthalene	0.400	0.386	3.5	58	0.00
10	Hexachlorobutadiene	0.400	0.388	3.0	60	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.368	8.0	56	0.00
12	2-Methylnaphthalene	0.400	0.377	5.8	58	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	60	0.00
14 S	2,4,6-Tribromophenol	0.400	0.426	-6.5	80	0.00
15 S	2-Fluorobiphenyl	0.400	0.377	5.8	60	0.00
16	Acenaphthylene	0.400	0.357	10.8	57	0.00
17	Acenaphthene	0.400	0.379	5.3	60	0.00
18	Fluorene	0.400	0.375	6.3	59	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	59	0.00
20	4-Bromophenyl-phenylether	0.400	0.379	5.3	60	0.00
21	Hexachlorobenzene	0.400	0.365	8.8	57	0.00
22	Pentachlorophenol	0.400	0.384	4.0	64	0.00
23	Phenanthrene	0.400	0.384	4.0	60	0.00
24	Anthracene	0.400	0.362	9.5	59	0.00
25 SURR	Fluoranthene-d10	0.400	0.376	6.0	60	0.00
26	Fluoranthene	0.400	0.371	7.3	59	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	56	0.00
28	Pvrene	0.400	0.385	3.8	60	0.00
29 S	Terphenyl-d14	0.400	0.378	5.5	59	0.00
30	Benzo(a)anthracene	0.400	0.377	5.8	59	0.00
31	Chrysene	0.400	0.391	2.3	60	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.367	8.3	60	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.358	10.5	55	0.00
34 I	Perylene-d12	0.400	0.400	0.0	58	0.00
35	Benzo(b)fluoranthene	0.400	0.352	12.0	57	0.00
36	Benzo(k)fluoranthene	0.400	0.389	2.8	63	0.00
37 C	Benzo(a)pyrene	0.400	0.370	7.5	60	0.00
38	Dibenzo(a,h)anthracene	0.400	0.349	12.8	56	0.00
39	Benzo(a,h,i)perylene	0.400	0.348	13.0	55	0.00

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

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