

Data Path : Z:\svoasrv\HPCHEM1\BNA E\Data\BE082318\
 Data File : BE097357.D
 Acq On : 23 Aug 2018 14:25
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Aug 23 15:42:58 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE082218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 22 18:07:26 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	96	0.00
2	1,4-Dioxane	0.400	0.440	-10.0	109	0.00
3	n-Nitrosodimethylamine	0.400	0.389	2.8	99	0.00
4 S	2-Fluorophenol	0.400	0.356	11.0	96	-0.01
5 S	Phenol-d6	0.400	0.355	11.3	95	-0.01
6	bis(2-Chloroethyl)ether	0.400	0.381	4.8	98	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	91	0.00
8 S	Nitrobenzene-d5	0.400	0.405	-1.3	104	0.00
9	Naphthalene	0.400	0.417	-4.2	101	0.00
10	Hexachlorobutadiene	0.400	0.423	-5.7	102	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.409	-2.2	101	0.00
12	2-Methylnaphthalene	0.400	0.412	-3.0	101	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	90	-0.01
14 S	2,4,6-Tribromophenol	0.400	0.362	9.5	98	0.00
15 S	2-Fluorobiphenyl	0.400	0.434	-8.5	102	0.00
16	Acenaphthylene	0.400	0.407	-1.7	102	0.00
17	Acenaphthene	0.400	0.425	-6.2	104	0.00
18	Fluorene	0.400	0.421	-5.2	103	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	90	0.00
20	4-Bromophenyl-phenylether	0.400	0.424	-6.0	107	0.00
21	Hexachlorobenzene	0.400	0.424	-6.0	104	0.00
22	Pentachlorophenol	0.400	0.413	-3.2	89	0.00
23	Phenanthrene	0.400	0.423	-5.7	104	0.00
24	Anthracene	0.400	0.410	-2.5	104	0.00
25 SURR	Fluoranthene-d10	0.400	0.395	1.3	98	0.00
26	Fluoranthene	0.400	0.399	0.3	100	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	89	0.00
28	Pvrene	0.400	0.408	-2.0	98	0.00
29 S	Terphenyl-d14	0.400	0.402	-0.5	97	0.00
30	Benzo(a)anthracene	0.400	0.388	3.0	95	0.00
31	Chrysene	0.400	0.418	-4.5	99	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.389	2.8	94	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.350	12.5	89	0.00
34 I	Perylene-d12	0.400	0.400	0.0	91	0.00
35	Benzo(b)fluoranthene	0.400	0.375	6.3	94	0.00
36	Benzo(k)fluoranthene	0.400	0.419	-4.7	104	0.00
37 C	Benzo(a)pyrene	0.400	0.361	9.8	94	0.00
38	Dibenzo(a,h)anthracene	0.400	0.362	9.5	92	0.00
39	Benzo(a,h,i)perylene	0.400	0.338	15.5	87	0.00

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

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