

Data Path : Z:\svoasrv\HPCHEM1\BNA E\Data\BE091018\
 Data File : BE097445.D
 Acq On : 10 Sep 2018 13:37
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Sep 10 17:07:40 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	97	0.00
2	1,4-Dioxane	0.400	0.372	7.0	92	0.00
3	n-Nitrosodimethylamine	0.400	0.378	5.5	97	0.00
4 S	2-Fluorophenol	0.400	0.410	-2.5	111	-0.01
5 S	Phenol-d6	0.400	0.432	-8.0	116	-0.01
6	bis(2-Chloroethyl)ether	0.400	0.380	5.0	98	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	110	0.00
8 S	Nitrobenzene-d5	0.400	0.429	-7.2	132	0.00
9	Naphthalene	0.400	0.412	-3.0	120	-0.01
10	Hexachlorobutadiene	0.400	0.404	-1.0	117	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.457	-14.2	136	0.00
12	2-Methylnaphthalene	0.400	0.458	-14.5	135	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	139	-0.01
14 S	2,4,6-Tribromophenol	0.400	0.462	-15.5	194	0.00
15 S	2-Fluorobiphenyl	0.400	0.386	3.5	141	0.00
16	Acenaphthylene	0.400	0.420	-5.0	163	-0.01
17	Acenaphthene	0.400	0.412	-3.0	156	-0.01
18	Fluorene	0.400	0.429	-7.2	162	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	130	-0.01
20	4-Bromophenyl-phenylether	0.400	0.434	-8.5	158	0.00
21	Hexachlorobenzene	0.400	0.439	-9.7	155	-0.01
22	Pentachlorophenol	0.400	0.465	-16.3	160	-0.01
23	Phenanthrene	0.400	0.420	-5.0	149	0.00
24	Anthracene	0.400	0.437	-9.2	160	0.00
25 SURR	Fluoranthene-d10	0.400	0.383	4.3	137	0.00
26	Fluoranthene	0.400	0.377	5.8	136	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	106	0.00
28	Pvrene	0.400	0.470	-17.5	135	0.00
29 S	Terphenyl-d14	0.400	0.437	-9.2	125	0.00
30	Benzo(a)anthracene	0.400	0.421	-5.2	123	0.00
31	Chrysene	0.400	0.414	-3.5	118	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.501	-25.2#	154	-0.01
33	Indeno(1,2,3-cd)pyrene	0.400	0.372	7.0	111	0.00
34 I	Perylene-d12	0.400	0.400	0.0	108	0.00
35	Benzo(b)fluoranthene	0.400	0.411	-2.7	123	0.00
36	Benzo(k)fluoranthene	0.400	0.401	-0.3	118	0.00
37 C	Benzo(a)pyrene	0.400	0.403	-0.8	125	0.00
38	Dibenzo(a,h)anthracene	0.400	0.363	9.3	109	0.00
39	Benzo(a,h,i)perylene	0.400	0.341	14.8	104	0.00

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

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