

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE032519\
 Data File : BE099217.D
 Acq On : 26 Mar 2019 16:15
 Operator : JU/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Mar 26 17:03:30 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE032519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 26 07:02:06 2019
 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	113	-0.01
2	1,4-Dioxane	0.400	0.326	18.5	105	0.00
3	n-Nitrosodimethylamine	0.400	0.316	21.0	117	-0.01
4 S	2-Fluorophenol	0.400	0.343	14.2	107	0.00
5 S	Phenol-d6	0.400	0.344	14.0	109	0.00
6	bis(2-Chloroethyl)ether	0.400	0.351	12.3	113	-0.01
7 I	Naphthalene-d8	0.400	0.400	0.0	111	0.00
8 S	Nitrobenzene-d5	0.400	0.354	11.5	121	0.00
9	Naphthalene	0.400	0.347	13.3	109	0.00
10	Hexachlorobutadiene	0.400	0.356	11.0	114	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.344	14.0	107	0.00
12	2-Methylnaphthalene	0.400	0.350	12.5	109	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	111	-0.01
14 S	2,4,6-Tribromophenol	0.400	0.472	-18.0	166	0.00
15 S	2-Fluorobiphenyl	0.400	0.354	11.5	109	0.00
16	Acenaphthylene	0.400	0.348	13.0	114	0.00
17	Acenaphthene	0.400	0.353	11.8	112	0.00
18	Fluorene	0.400	0.345	13.8	111	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	113	-0.01
20	4-Bromophenyl-phenylether	0.400	0.323	19.3	108	-0.01
21	Hexachlorobenzene	0.400	0.333	16.8	112	-0.01
22	Pentachlorophenol	0.400	0.702	-75.5#	358	-0.01
23	Phenanthrene	0.400	0.338	15.5	111	0.00
24	Anthracene	0.400	0.334	16.5	115	0.00
25 SURR	Fluoranthene-d10	0.400	0.354	11.5	116	0.00
26	Fluoranthene	0.400	0.340	15.0	116	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	118	-0.01
28	Pvrene	0.400	0.343	14.2	117	0.00
29 S	Terphenyl-d14	0.400	0.336	16.0	110	0.00
30	Benzo(a)anthracene	0.400	0.342	14.5	121	0.00
31	Chrysene	0.400	0.341	14.8	117	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.355	11.3	149	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.310	22.5	115	-0.01
34 I	Perylene-d12	0.400	0.400	0.0	112	0.00
35	Benzo(b)fluoranthene	0.400	0.328	18.0	112	0.00
36	Benzo(k)fluoranthene	0.400	0.360	10.0	119	0.00
37 C	Benzo(a)pyrene	0.400	0.344	14.0	117	0.00
38	Dibenzo(a,h)anthracene	0.400	0.294	26.5#	107	-0.01
39	Benzo(a,h,i)perylene	0.400	0.300	25.0#	106	0.00

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

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