

Data Path : Z:\svoasrv\HPCHEM1\BNA E\Data\BE092118\
 Data File : BE097626.D
 Acq On : 22 Sep 2018 00:03
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Sep 22 04:50:27 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE091918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Sep 20 10:54:43 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	93	0.00
2	1,4-Dioxane	0.400	0.354	11.5	80	0.00
3	n-Nitrosodimethylamine	0.400	0.390	2.5	104	0.01
4 S	2-Fluorophenol	0.400	0.334	16.5	85	0.00
5 S	Phenol-d6	0.400	0.374	6.5	93	0.00
6	bis(2-Chloroethyl)ether	0.400	0.333	16.8	85	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	97	0.00
8 S	Nitrobenzene-d5	0.400	0.330	17.5	91	0.01
9	Naphthalene	0.400	0.399	0.3	104	0.01
10	Hexachlorobutadiene	0.400	0.382	4.5	98	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.421	-5.2	112	0.00
12	2-Methylnaphthalene	0.400	0.412	-3.0	110	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	125	0.00
14 S	2,4,6-Tribromophenol	0.400	0.402	-0.5	144	0.01
15 S	2-Fluorobiphenyl	0.400	0.356	11.0	120	0.01
16	Acenaphthylene	0.400	0.382	4.5	134	0.00
17	Acenaphthene	0.400	0.387	3.3	129	0.00
18	Fluorene	0.400	0.408	-2.0	140	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	139	0.00
20	4-Bromophenyl-phenylether	0.400	0.385	3.8	146	0.00
21	Hexachlorobenzene	0.400	0.411	-2.7	152	0.00
22	Pentachlorophenol	0.400	0.300	25.0#	95	0.01
23	Phenanthrene	0.400	0.389	2.8	149	0.00
24	Anthracene	0.400	0.355	11.3	139	0.00
25 SURR	Fluoranthene-d10	0.400	0.319	20.3	121	0.00
26	Fluoranthene	0.400	0.326	18.5	123	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	96	0.00
28	Pvrene	0.400	0.450	-12.5	117	0.00
29 S	Terphenyl-d14	0.400	0.417	-4.2	109	0.00
30	Benzo(a)anthracene	0.400	0.352	12.0	94	0.00
31	Chrysene	0.400	0.412	-3.0	103	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.265	33.8#	70	-0.01
33	Indeno(1,2,3-cd)pyrene	0.400	0.391	2.3	96	0.00
34 I	Perylene-d12	0.400	0.400	0.0	90	0.00
35	Benzo(b)fluoranthene	0.400	0.368	8.0	92	0.00
36	Benzo(k)fluoranthene	0.400	0.408	-2.0	102	0.00
37 C	Benzo(a)pyrene	0.400	0.387	3.3	94	0.00
38	Dibenzo(a,h)anthracene	0.400	0.394	1.5	96	0.00
39	Benzo(a,h,i)perylene	0.400	0.392	2.0	95	0.01

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

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