

Data Path : Z:\svoasrv\HPCHEM1\BNA E\Data\BE102418\
 Data File : BE098047.D
 Acq On : 24 Oct 2018 20:26
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Oct 25 04:03:57 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE102418.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 24 13:49:35 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	88	0.00
2	1,4-Dioxane	0.400	0.397	0.8	93	0.00
3	n-Nitrosodimethylamine	0.400	0.367	8.3	85	0.00
4 S	2-Fluorophenol	0.400	0.368	8.0	84	0.00
5 S	Phenol-d6	0.400	0.376	6.0	85	0.00
6	bis(2-Chloroethyl)ether	0.400	0.332	17.0	75	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	89	0.00
8 S	Nitrobenzene-d5	0.400	0.371	7.3	86	0.00
9	Naphthalene	0.400	0.369	7.8	85	0.00
10	Hexachlorobutadiene	0.400	0.387	3.3	91	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.368	8.0	87	0.00
12	2-Methylnaphthalene	0.400	0.369	7.8	84	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	91	0.00
14 S	2,4,6-Tribromophenol	0.400	0.358	10.5	88	0.00
15 S	2-Fluorobiphenyl	0.400	0.355	11.3	87	0.00
16	Acenaphthylene	0.400	0.371	7.3	89	-0.01
17	Acenaphthene	0.400	0.370	7.5	87	0.00
18	Fluorene	0.400	0.367	8.3	88	-0.01
19 I	Phenanthrene-d10	0.400	0.400	0.0	95	0.00
20	4-Bromophenyl-phenylether	0.400	0.361	9.8	87	0.00
21	Hexachlorobenzene	0.400	0.365	8.8	90	-0.01
22	Pentachlorophenol	0.400	0.372	7.0	98	0.00
23	Phenanthrene	0.400	0.359	10.3	88	-0.01
24	Anthracene	0.400	0.354	11.5	89	0.00
25 SURR	Fluoranthene-d10	0.400	0.356	11.0	88	0.00
26	Fluoranthene	0.400	0.353	11.8	87	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	90	0.00
28	Pvrene	0.400	0.353	11.8	86	0.00
29 S	Terphenyl-d14	0.400	0.353	11.8	88	0.00
30	Benzo(a)anthracene	0.400	0.365	8.8	89	-0.01
31	Chrysene	0.400	0.354	11.5	86	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.341	14.8	83	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.360	10.0	87	-0.01
34 I	Perylene-d12	0.400	0.400	0.0	92	-0.01
35	Benzo(b)fluoranthene	0.400	0.360	10.0	87	-0.01
36	Benzo(k)fluoranthene	0.400	0.354	11.5	87	0.00
37 C	Benzo(a)pyrene	0.400	0.356	11.0	88	-0.01
38	Dibenzo(a,h)anthracene	0.400	0.356	11.0	86	0.00
39	Benzo(a,h,i)perylene	0.400	0.356	11.0	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				