

Data Path : Z:\HPCHEM1\BNA E\Data\BE032618\
 Data File : BE095848.D
 Acq On : 26 Mar 2018 20:25
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Mar 27 05:42:48 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE032318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 23 12:41:11 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	104	0.00
2	1,4-Dioxane	0.400	0.391	2.3	107	0.00
3	n-Nitrosodimethylamine	0.400	0.365	8.8	92	0.00
4 S	2-Fluorophenol	0.400	0.386	3.5	96	0.00
5 S	Phenol-d6	0.400	0.394	1.5	98	0.00
6	bis(2-Chloroethyl)ether	0.400	0.392	2.0	98	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	101	0.00
8 S	Nitrobenzene-d5	0.400	0.398	0.5	100	0.00
9	Naphthalene	0.400	0.410	-2.5	100	-0.01
10	Hexachlorobutadiene	0.400	0.385	3.8	93	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.401	-0.3	98	-0.01
12	2-Methylnaphthalene	0.400	0.412	-3.0	102	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	101	-0.01
14 S	2,4,6-Tribromophenol	0.400	0.374	6.5	98	-0.01
15 S	2-Fluorobiphenyl	0.400	0.404	-1.0	100	-0.01
16	Acenaphthylene	0.400	0.405	-1.3	102	0.00
17	Acenaphthene	0.400	0.396	1.0	101	0.00
18	Fluorene	0.400	0.410	-2.5	102	-0.01
19 I	Phenanthrene-d10	0.400	0.400	0.0	102	-0.01
20	4-Bromophenyl-phenylether	0.400	0.402	-0.5	101	0.00
21	Hexachlorobenzene	0.400	0.401	-0.3	101	0.00
22	Pentachlorophenol	0.400	0.416	-4.0	114	-0.01
23	Phenanthrene	0.400	0.421	-5.2	104	0.00
24	Anthracene	0.400	0.405	-1.3	102	0.00
25 SURR	Fluoranthene-d10	0.400	0.404	-1.0	101	0.00
26	Fluoranthene	0.400	0.411	-2.7	104	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	102	-0.01
28	Pvrene	0.400	0.380	5.0	88	-0.01
29 S	Terphenyl-d14	0.400	0.399	0.3	100	0.00
30	Benzo(a)anthracene	0.400	0.413	-3.2	103	-0.01
31	Chrysene	0.400	0.417	-4.2	104	-0.01
32	Bis(2-ethylhexyl)phthalate	0.400	0.363	9.3	98	-0.01
33	Indeno(1,2,3-cd)pyrene	0.400	0.406	-1.5	102	-0.03
34 I	Perylene-d12	0.400	0.400	0.0	101	-0.02
35	Benzo(b)fluoranthene	0.400	0.414	-3.5	103	-0.01
36	Benzo(k)fluoranthene	0.400	0.415	-3.7	104	-0.01
37 C	Benzo(a)pyrene	0.400	0.412	-3.0	102	-0.02
38	Dibenzo(a,h)anthracene	0.400	0.412	-3.0	101	-0.02
39	Benzo(a,h,i)perylene	0.400	0.413	-3.2	102	-0.03

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

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