

Data Path : Z:\HPCHEM1\BNA_E\Data\BE111617\
 Data File : BE094823.D
 Acq On : 15 Nov 2017 20:12
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Nov 16 02:54:17 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE102317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 23 22:40:47 2017
 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	66	0.01
2	1,4-Dioxane	0.400	0.482	-20.5	69	-0.02
3	n-Nitrosodimethylamine	0.400	0.376	6.0	62	0.00
4 S	2-Fluorophenol	0.400	0.333	16.8	55	0.04
5 S	Phenol-d6	0.400	0.384	4.0	65	0.07
6	bis(2-Chloroethyl)ether	0.400	0.325	18.8	54	0.03
7 I	Naphthalene-d8	0.400	0.400	0.0	67	0.02
8 S	Nitrobenzene-d5	0.400	0.334	16.5	59	0.07
9	Naphthalene	0.400	0.355	11.3	58	0.03
10	Hexachlorobutadiene	0.400	0.404	-1.0	67	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.366	8.5	60	0.03
12	2-Methylnaphthalene	0.400	0.359	10.3	60	0.03
13 I	Acenaphthene-d10	0.400	0.400	0.0	71	0.02
14 S	2,4,6-Tribromophenol	0.400	0.513	-28.2#	95	0.03
15 S	2-Fluorobiphenyl	0.400	0.346	13.5	61	0.03
16	Acenaphthylene	0.400	0.361	9.8	65	0.00
17	Acenaphthene	0.400	0.372	7.0	65	0.00
18	Fluorene	0.400	0.393	1.8	70	0.02
19 I	Phenanthrene-d10	0.400	0.400	0.0	97	0.03
20	4-Bromophenyl-phenylether	0.400	0.301	24.8	74	0.03
21	Hexachlorobenzene	0.400	0.474	-18.5	100	0.00
22	Pentachlorophenol	0.400	0.339	15.3	82	0.07
23	Phenanthrene	0.400	0.497	-24.2	108	0.00
24	Anthracene	0.400	0.438	-9.5	104	0.00
25 SURR	Fluoranthene-d10	0.400	0.422	-5.5	93	0.01
26	Fluoranthene	0.400	0.419	-4.7	92	0.01
27 I	Chrysene-d12	0.400	0.400	0.0	102	0.01
28	Pyrene	0.400	0.358	10.5	93	0.01
29 S	Terphenyl-d14	0.400	0.356	11.0	92	0.00
30	Benzo(a)anthracene	0.400	0.362	9.5	92	0.01
31	Chrysene	0.400	0.400	0.0	105	0.01
32	Bis(2-ethylhexyl)phthalate	0.400	0.347	13.3	96	-0.01
33	Indeno(1,2,3-cd)pyrene	0.400	0.391	2.3	101	0.07
34 I	Perylene-d12	0.400	0.400	0.0	101	0.03
35	Benzo(b)fluoranthene	0.400	0.364	9.0	95	0.02
36	Benzo(k)fluoranthene	0.400	0.399	0.3	102	0.02
37 C	Benzo(a)pyrene	0.400	0.382	4.5	97	0.04
38	Dibenzo(a,h)anthracene	0.400	0.392	2.0	101	0.05
39	Benzo(g,h,i)perylene	0.400	0.414	-3.5	106	0.07

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

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