

Data Path : Z:\HPCHEM1\BNA_E\Data\BE110217\
 Data File : BE094694.D
 Acq On : 2 Nov 2017 14:10
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Nov 02 18:01:54 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	114	0.00
2	1,4-Dioxane	0.400	0.354	11.5	87	-0.01
3	n-Nitrosodimethylamine	0.400	0.332	17.0	94	0.01
4 S	2-Fluorophenol	0.400	0.311	22.3	89	0.04
5 S	Phenol-d6	0.400	0.372	7.0	109	0.04
6	bis(2-Chloroethyl)ether	0.400	0.342	14.5	98	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	125	0.00
8 S	Nitrobenzene-d5	0.400	0.334	16.5	108	0.02
9	Naphthalene	0.400	0.344	14.0	104	0.02
10	Hexachlorobutadiene	0.400	0.368	8.0	113	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.340	15.0	103	0.00
12	2-Methylnaphthalene	0.400	0.333	16.8	103	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	107	0.00
14 S	2,4,6-Tribromophenol	0.400	0.320	20.0	89	0.03
15 S	2-Fluorobiphenyl	0.400	0.376	6.0	99	0.00
16	Acenaphthylene	0.400	0.348	13.0	94	0.00
17	Acenaphthene	0.400	0.364	9.0	96	0.00
18	Fluorene	0.400	0.350	12.5	94	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	101	0.03
20	4-Bromophenyl-phenylether	0.400	0.375	6.3	97	0.00
21	Hexachlorobenzene	0.400	0.544	-36.0#	120	0.00
22	Pentachlorophenol	0.400	0.350	12.5	89	0.07
23	Phenanthrene	0.400	0.422	-5.5	99	0.00
24	Anthracene	0.400	0.377	5.8	94	0.00
25 SURR	Fluoranthene-d10	0.400	0.417	-4.2	96	0.00
26	Fluoranthene	0.400	0.409	-2.2	95	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	104	0.00
28	Pyrene	0.400	0.364	9.0	97	0.00
29 S	Terphenyl-d14	0.400	0.352	12.0	94	0.00
30	Benzo(a)anthracene	0.400	0.354	11.5	92	0.01
31	Chrysene	0.400	0.368	8.0	99	0.01
32	Bis(2-ethylhexyl)phthalate	0.400	0.344	14.0	98	-0.01
33	Indeno(1,2,3-cd)pyrene	0.400	0.277	30.8#	74	0.04
34 I	Perylene-d12	0.400	0.400	0.0	100	0.01
35	Benzo(b)fluoranthene	0.400	0.351	12.3	90	0.00
36	Benzo(k)fluoranthene	0.400	0.402	-0.5	101	0.01
37 C	Benzo(a)pyrene	0.400	0.372	7.0	93	0.01
38	Dibenzo(a,h)anthracene	0.400	0.305	23.8	78	0.02
39	Benzo(g,h,i)perylene	0.400	0.267	33.3#	67	0.05

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

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