

Data Path : \\74.0.250.170\svoasrv\HPCHEM1\BNA E\Data\BE110917\
 Data File : BE094793.D
 Acq On : 8 Nov 2017 21:17
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Nov 09 00:50:07 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	102	0.01
2	1,4-Dioxane	0.400	0.356	11.0	78	-0.01
3	n-Nitrosodimethylamine	0.400	0.291	27.3#	73	0.03
4 S	2-Fluorophenol	0.400	0.300	25.0#	76	0.05
5 S	Phenol-d6	0.400	0.368	8.0	96	0.06
6	bis(2-Chloroethyl)ether	0.400	0.351	12.3	90	0.01
7 I	Naphthalene-d8	0.400	0.400	0.0	108	0.00
8 S	Nitrobenzene-d5	0.400	0.326	18.5	92	0.05
9	Naphthalene	0.400	0.353	11.8	93	0.02
10	Hexachlorobutadiene	0.400	0.398	0.5	106	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.356	11.0	94	0.02
12	2-Methylnaphthalene	0.400	0.344	14.0	92	0.02
13 I	Acenaphthene-d10	0.400	0.400	0.0	96	0.02
14 S	2,4,6-Tribromophenol	0.400	0.363	9.3	91	0.03
15 S	2-Fluorobiphenyl	0.400	0.373	6.8	89	0.02
16	Acenaphthylene	0.400	0.367	8.3	89	0.00
17	Acenaphthene	0.400	0.365	8.8	87	0.00
18	Fluorene	0.400	0.355	11.3	86	0.02
19 I	Phenanthrene-d10	0.400	0.400	0.0	108	0.03
20	4-Bromophenyl-phenylether	0.400	0.296	26.0#	81	0.03
21	Hexachlorobenzene	0.400	0.482	-20.5	113	0.00
22	Pentachlorophenol	0.400	0.436	-9.0	118	0.03
23	Phenanthrene	0.400	0.478	-19.5	117	0.00
24	Anthracene	0.400	0.396	1.0	105	0.00
25 SURR	Fluoranthene-d10	0.400	0.348	13.0	89	0.01
26	Fluoranthene	0.400	0.345	13.8	88	0.01
27 I	Chrysene-d12	0.400	0.400	0.0	92	0.01
28	Pvrene	0.400	0.379	5.3	89	0.01
29 S	Terphenyl-d14	0.400	0.367	8.3	86	0.00
30	Benzo(a)anthracene	0.400	0.376	6.0	86	0.01
31	Chrysene	0.400	0.375	6.3	89	0.01
32	Bis(2-ethylhexyl)phthalate	0.400	0.366	8.5	92	-0.01
33	Indeno(1,2,3-cd)pyrene	0.400	0.266	33.5#	62	0.07
34 I	Perylene-d12	0.400	0.400	0.0	85	0.03
35	Benzo(b)fluoranthene	0.400	0.369	7.8	80	0.02
36	Benzo(k)fluoranthene	0.400	0.426	-6.5	91	0.02
37 C	Benzo(a)pyrene	0.400	0.386	3.5	82	0.03
38	Dibenzo(a,h)anthracene	0.400	0.317	20.8	69	0.05
39	Benzo(a,h,i)perylene	0.400	0.262	34.5#	56	0.09

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

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