

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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	66	0.01
2	1,4-Dioxane	0.677	0.816	-20.5	69	-0.02
3	n-Nitrosodimethylamine	0.648	0.609	6.0	62	0.00
4 S	2-Fluorophenol	1.370	1.140	16.8	55	0.04
5 S	Phenol-d6	2.069	1.983	4.2	65	0.07
6	bis(2-Chloroethyl)ether	1.600	1.300	18.8	54	0.03
7 I	Naphthalene-d8	1.000	1.000	0.0	67	0.02
8 S	Nitrobenzene-d5	0.319	0.266	16.6	59	0.07
9	Naphthalene	4.527	4.023	11.1	58	0.03
10	Hexachlorobutadiene	0.164	0.165	-0.6	67	0.00
11 SURR	2-Methylnaphthalene-d10	0.642	0.587	8.6	60	0.03
12	2-Methylnaphthalene	0.768	0.690	10.2	60	0.03
13 I	Acenaphthene-d10	1.000	1.000	0.0	71	0.02
14 S	2,4,6-Tribromophenol	0.110	0.142	-29.1#	95	0.03
15 S	2-Fluorobiphenyl	1.410	1.218	13.6	61	0.03
16	Acenaphthylene	2.112	1.905	9.8	65	0.00
17	Acenaphthene	1.356	1.262	6.9	65	0.00
18	Fluorene	1.738	1.709	1.7	70	0.02
19 I	Phenanthrene-d10	1.000	1.000	0.0	97	0.03
20	4-Bromophenyl-phenylether	0.218	0.164	24.8	74	0.03
21	Hexachlorobenzene	0.191	0.226	-18.3	100	0.00
22	Pentachlorophenol	0.024	0.021	12.5	82	0.07
23	Phenanthrene	1.141	1.286	-12.7	108	0.00
24	Anthracene	1.176	1.287	-9.4	104	0.00
25 SURR	Fluoranthene-d10	4.676	5.140	-9.9	93	0.01
26	Fluoranthene	1.420	1.556	-9.6	92	0.01
27 I	Chrysene-d12	1.000	1.000	0.0	102	0.01
28	Pyrene	1.399	1.254	10.4	93	0.01
29 S	Terphenyl-d14	0.754	0.671	11.0	92	0.00
30	Benzo(a)anthracene	1.302	1.177	9.6	92	0.01
31	Chrysene	1.322	1.323	-0.1	105	0.01
32	Bis(2-ethylhexyl)phthalate	3.193	2.766	13.4	96	-0.01
33	Indeno(1,2,3-cd)pyrene	1.222	1.194	2.3	101	0.07
34 I	Perylene-d12	1.000	1.000	0.0	101	0.03
35	Benzo(b)fluoranthene	1.294	1.178	9.0	95	0.02
36	Benzo(k)fluoranthene	1.379	1.376	0.2	102	0.02
37 C	Benzo(a)pyrene	1.255	1.198	4.5	97	0.04
38	Dibenzo(a,h)anthracene	1.127	1.103	2.1	101	0.05
39	Benzo(g,h,i)perylene	1.055	1.091	-3.4	106	0.07

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