

Data Path : \\74.0.250.170\svoasrv\HPCHEM1\BNA E\Data\BE120117\
 Data File : BE094913.D
 Acq On : 1 Dec 2017 19:58
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Dec 02 00:37:09 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE113017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 01 13:33:28 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	AvqRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	86	0.00
2	1,4-Dioxane	0.419	0.417	0.5	80	0.00
3	n-Nitrosodimethylamine	0.533	0.510	4.3	83	-0.01
4 S	2-Fluorophenol	0.633	0.582	8.1	78	0.00
5 S	Phenol-d6	0.938	0.868	7.5	79	0.00
6	bis(2-Chloroethyl)ether	0.930	0.841	9.6	76	-0.01
7 I	Naphthalene-d8	1.000	1.000	0.0	80	0.00
8 S	Nitrobenzene-d5	0.237	0.293	-23.6	99	0.00
9	Naphthalene	4.690	4.667	0.5	79	0.00
10	Hexachlorobutadiene	0.198	0.199	-0.5	80	0.00
11 SURR	2-Methylnaphthalene-d10	0.638	0.649	-1.7	81	0.00
12	2-Methylnaphthalene	1.165	1.163	0.2	81	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	83	-0.02
14 S	2,4,6-Tribromophenol	0.086	0.098	-14.0	99	0.00
15 S	2-Fluorobiphenyl	1.713	1.709	0.2	82	0.00
16	Acenaphthylene	2.260	2.184	3.4	82	0.00
17	Acenaphthene	1.461	1.464	-0.2	85	-0.02
18	Fluorene	1.859	1.885	-1.4	83	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	86	0.00
20	4-Bromophenyl-phenylether	0.223	0.217	2.7	85	0.00
21	Hexachlorobenzene	0.218	0.221	-1.4	86	0.00
22	Pentachlorophenol	0.054	0.054	0.0	95	-0.01
23	Phenanthrene	1.275	1.231	3.5	82	0.00
24	Anthracene	1.351	1.263	6.5	83	0.00
25 SURR	Fluoranthene-d10	5.466	5.150	5.8	82	0.00
26	Fluoranthene	1.638	1.555	5.1	82	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	80	0.00
28	Pvrene	1.381	1.368	0.9	79	-0.02
29 S	Terphenyl-d14	0.739	0.743	-0.5	81	0.00
30	Benzo(a)anthracene	1.227	1.194	2.7	80	-0.02
31	Chrysene	1.355	1.356	-0.1	80	0.00
32	Bis(2-ethylhexyl)phthalate	2.033	1.640	19.3	75	0.00
33	Indeno(1,2,3-cd)pyrene	1.109	1.088	1.9	80	0.00
34 I	Perylene-d12	1.000	1.000	0.0	81	0.00
35	Benzo(b)fluoranthene	1.460	1.400	4.1	79	0.00
36	Benzo(k)fluoranthene	1.489	1.464	1.7	82	0.00
37 C	Benzo(a)pyrene	1.280	1.227	4.1	80	0.00
38	Dibenzo(a,h)anthracene	1.183	1.154	2.5	81	-0.02
39	Benzo(a,h,i)perylene	1.181	1.133	4.1	79	-0.02

Data Path : \\74.0.250.170\svoasrv\HPCHEM1\BNA E\Data\BE120117\
Data File : BE094913.D
Acq On : 1 Dec 2017 19:58
Operator : SJ/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_E
LabSampleId :
SSTDCCC0.4

Quant Time: Dec 02 00:37:09 2017
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE113017.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Dec 01 13:33:28 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)

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