

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE020718\
 Data File : BE095545.D
 Acq On : 7 Feb 2018 8:41
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Feb 08 01:26:14 2018
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE020618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 06 20:27:34 2018
 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	103	0.00
2	1,4-Dioxane	0.662	0.678	-2.4	105	0.00
3	n-Nitrosodimethylamine	0.833	0.803	3.6	98	0.00
4 S	2-Fluorophenol	1.207	1.204	0.2	101	0.00
5 S	Phenol-d6	1.671	1.633	2.3	99	-0.01
6	bis(2-Chloroethyl)ether	1.538	1.575	-2.4	102	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	103	0.00
8 S	Nitrobenzene-d5	0.424	0.411	3.1	99	0.00
9	Naphthalene	4.646	4.702	-1.2	101	0.00
10	Hexachlorobutadiene	0.249	0.241	3.2	90	0.00
11 SURR	2-Methylnaphthalene-d10	0.662	0.679	-2.6	103	0.00
12	2-Methylnaphthalene	0.789	0.808	-2.4	101	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	103	0.00
14 S	2,4,6-Tribromophenol	0.182	0.159	12.6	92	0.00
15 S	2-Fluorobiphenyl	1.725	1.776	-3.0	104	0.00
16	Acenaphthylene	2.282	2.212	3.1	100	0.00
17	Acenaphthene	1.418	1.396	1.6	101	0.00
18	Fluorene	1.782	1.760	1.2	100	-0.01
19 I	Phenanthrene-d10	1.000	1.000	0.0	100	0.00
20	4-Bromophenyl-phenylether	0.240	0.250	-4.2	101	0.00
21	Hexachlorobenzene	0.248	0.263	-6.0	104	0.00
22	Pentachlorophenol	0.069	0.055	20.3	95	0.00
23	Phenanthrene	1.228	1.275	-3.8	102	-0.01
24	Anthracene	1.147	1.133	1.2	100	-0.01
25 SURR	Fluoranthene-d10	5.268	5.314	-0.9	101	0.00
26	Fluoranthene	1.550	1.575	-1.6	103	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	101	0.00
28	Pyrene	1.680	1.545	8.0	97	0.00
29 S	Terphenyl-d14	0.852	0.857	-0.6	99	0.00
30	Benzo(a)anthracene	1.333	1.371	-2.9	103	0.00
31	Chrysene	1.314	1.356	-3.2	102	0.00
32	Bis(2-ethylhexyl)phthalate	3.072	2.671	13.1	92	0.00
33	Indeno(1,2,3-cd)pyrene	1.269	1.226	3.4	100	0.00
34 I	Perylene-d12	1.000	1.000	0.0	101	0.00
35	Benzo(b)fluoranthene	1.364	1.397	-2.4	102	0.00
36	Benzo(k)fluoranthene	1.402	1.425	-1.6	102	0.00
37 C	Benzo(a)pyrene	1.269	1.267	0.2	102	0.00
38	Dibenzo(a,h)anthracene	1.111	1.062	4.4	100	0.00
39	Benzo(g,h,i)perylene	1.193	1.159	2.8	98	0.00

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE020718\
Data File : BE095545.D
Acq On : 7 Feb 2018 8:41
Operator : SJ/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_E
LabSampleId :
SSTDCCC0.4

Quant Time: Feb 08 01:26:14 2018
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE020618.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Feb 06 20:27:34 2018
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			