

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE092319\
 Data File : BE100535.D
 Acq On : 23 Sep 2019 10:34
 Operator : HP/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Sep 23 12:23:37 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE092019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 20 18:26:17 2019
 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	98	0.00
2	1,4-Dioxane	0.577	0.574	0.5	97	0.00
3	n-Nitrosodimethylamine	0.339	0.297	12.4	114	0.00
4 S	2-Fluorophenol	0.936	0.848	9.4	96	0.00
5 S	Phenol-d6	1.282	1.143	10.8	97	0.00
6	bis(2-Chloroethyl)ether	1.297	1.216	6.2	97	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	99	0.00
8 S	Nitrobenzene-d5	0.231	0.208	10.0	99	0.00
9	Naphthalene	4.147	4.041	2.6	99	0.00
10	Hexachlorobutadiene	0.133	0.130	2.3	99	0.00
11 SURR	2-Methylnaphthalene-d10	0.598	0.567	5.2	99	0.00
12	2-Methylnaphthalene	0.660	0.623	5.6	99	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	98	0.00
14 S	2,4,6-Tribromophenol	0.065	0.055	15.4	92	0.00
15 S	2-Fluorobiphenyl	1.428	1.409	1.3	99	0.00
16	Acenaphthylene	1.647	1.476	10.4	96	0.01
17	Acenaphthene	1.112	1.061	4.6	99	0.00
18	Fluorene	1.391	1.318	5.2	99	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	100	0.00
20	4-Bromophenyl-phenylether	0.188	0.179	4.8	99	0.01
21	Hexachlorobenzene	0.164	0.158	3.7	102	0.00
22	Pentachlorophenol	0.037	0.020	45.9#	89	0.00
23	Phenanthrene	1.115	1.058	5.1	99	0.00
24	Anthracene	0.938	0.834	11.1	99	0.00
25 SURR	Fluoranthene-d10	4.758	4.591	3.5	102	0.00
26	Fluoranthene	1.344	1.300	3.3	102	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	104	0.00
28	Pvrene	1.151	1.104	4.1	103	0.00
29 S	Terphenyl-d14	0.947	0.927	2.1	105	0.00
30	Benzo(a)anthracene	1.068	1.017	4.8	104	0.00
31	Chrysene	1.217	1.174	3.5	105	0.00
32	Bis(2-ethylhexyl)phthalate	1.151	1.038	9.8	104	0.00
33	Indeno(1,2,3-cd)pyrene	1.353	1.254	7.3	99	0.00
34 I	Perylene-d12	1.000	1.000	0.0	104	0.00
35	Benzo(b)fluoranthene	1.082	0.983	9.1	104	0.00
36	Benzo(k)fluoranthene	1.252	1.198	4.3	107	0.00
37 C	Benzo(a)pyrene	1.010	0.914	9.5	101	0.00
38	Dibenzo(a,h)anthracene	1.052	0.934	11.2	100	0.00
39	Benzo(a,h,i)perylene	1.098	1.010	8.0	99	0.00

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE092319\
Data File : BE100535.D
Acq On : 23 Sep 2019 10:34
Operator : HP/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_E
LabSampleId :
SSTDCCC0.4

Quant Time: Sep 23 12:23:37 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE092019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Sep 20 18:26:17 2019
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			