

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE092319\
 Data File : BE100548.D
 Acq On : 23 Sep 2019 20:37
 Operator : HP/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Sep 24 05:50:17 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	93	0.00
2	1,4-Dioxane	0.577	0.554	4.0	88	0.00
3	n-Nitrosodimethylamine	0.339	0.319	5.9	115	0.00
4 S	2-Fluorophenol	0.936	1.052	-12.4	113	0.00
5 S	Phenol-d6	1.282	1.381	-7.7	110	0.00
6	bis(2-Chloroethyl)ether	1.297	1.276	1.6	96	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	96	-0.01
8 S	Nitrobenzene-d5	0.231	0.232	-0.4	107	0.00
9	Naphthalene	4.147	4.068	1.9	96	0.00
10	Hexachlorobutadiene	0.133	0.130	2.3	95	0.00
11 SURR	2-Methylnaphthalene-d10	0.598	0.591	1.2	100	0.00
12	2-Methylnaphthalene	0.660	0.644	2.4	99	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	97	0.00
14 S	2,4,6-Tribromophenol	0.065	0.073	-12.3	120	0.00
15 S	2-Fluorobiphenyl	1.428	1.408	1.4	98	0.00
16	Acenaphthylene	1.647	1.557	5.5	101	0.00
17	Acenaphthene	1.112	1.080	2.9	100	0.00
18	Fluorene	1.391	1.329	4.5	99	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	99	0.00
20	4-Bromophenyl-phenylether	0.188	0.177	5.9	97	0.00
21	Hexachlorobenzene	0.164	0.152	7.3	96	0.00
22	Pentachlorophenol	0.037	0.027	27.0#	122	-0.01
23	Phenanthrene	1.115	1.070	4.0	99	0.00
24	Anthracene	0.938	0.858	8.5	101	0.00
25 SURR	Fluoranthene-d10	4.758	4.463	6.2	98	0.00
26	Fluoranthene	1.344	1.260	6.3	98	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	93	0.00
28	Pvrene	1.151	1.181	-2.6	99	0.00
29 S	Terphenyl-d14	0.947	0.958	-1.2	98	0.00
30	Benzo(a)anthracene	1.068	1.102	-3.2	101	0.00
31	Chrysene	1.217	1.165	4.3	94	0.00
32	Bis(2-ethylhexyl)phthalate	1.151	1.420	-23.4	128	0.00
33	Indeno(1,2,3-cd)pyrene	1.353	1.335	1.3	95	0.00
34 I	Perylene-d12	1.000	1.000	0.0	96	0.00
35	Benzo(b)fluoranthene	1.082	0.996	7.9	97	0.00
36	Benzo(k)fluoranthene	1.252	1.184	5.4	97	0.00
37 C	Benzo(a)pyrene	1.010	0.938	7.1	95	0.00
38	Dibenzo(a,h)anthracene	1.052	0.967	8.1	95	0.00
39	Benzo(a,h,i)perylene	1.098	1.003	8.7	91	0.00

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE092319\
Data File : BE100548.D
Acq On : 23 Sep 2019 20:37
Operator : HP/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_E
LabSampleId :
SSTDCCC0.4

Quant Time: Sep 24 05:50:17 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			