

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE061719\
 Data File : BE100069.D
 Acq On : 17 Jun 2019 19:25
 Operator : JU/SJ
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDC00.4

Quant Time: Jun 18 03:44:44 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE061719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 17 17:23:36 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	0.698	0.628	10.0	84	0.01
3	n-Nitrosodimethylamine	0.289	0.157	45.7#	69	0.01
4 S	2-Fluorophenol	1.083	0.975	10.0	89	0.00
5 S	Phenol-d6	1.306	1.109	15.1	83	0.01
6	bis(2-Chloroethyl)ether	1.225	1.034	15.6	93	0.01
7 I	Naphthalene-d8	1.000	1.000	0.0	96	0.00
8 S	Nitrobenzene-d5	0.396	0.351	11.4	94	0.01
9	Naphthalene	4.423	4.494	-1.6	97	0.00
10	Hexachlorobutadiene	0.380	0.403	-6.1	101	0.00
11 SURR	2-Methylnaphthalene-d10	0.723	0.731	-1.1	100	0.00
12	2-Methylnaphthalene	0.693	0.714	-3.0	102	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	97	0.00
14 S	2,4,6-Tribromophenol	0.245	0.209	14.7	96	0.00
15 S	2-Fluorobiphenyl	1.551	1.530	1.4	94	0.00
16	Acenaphthylene	2.004	2.025	-1.0	100	0.00
17	Acenaphthene	1.084	1.087	-0.3	97	-0.01
18	Fluorene	1.614	1.682	-4.2	102	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	100	0.01
20	4-Bromophenyl-phenylether	0.241	0.235	2.5	101	0.00
21	Hexachlorobenzene	0.247	0.246	0.4	102	0.00
22	Pentachlorophenol	0.077	0.057	26.0#	88	0.01
23	Phenanthrene	0.975	0.933	4.3	99	0.00
24	Anthracene	0.857	0.805	6.1	101	0.00
25 SURR	Fluoranthene-d10	5.626	5.302	5.8	98	0.00
26	Fluoranthene	1.604	1.529	4.7	99	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	94	0.00
28	Pvrene	0.995	1.149	-15.5	98	0.00
29 S	Terphenyl-d14	0.674	0.736	-9.2	94	0.01
30	Benzo(a)anthracene	1.152	1.230	-6.8	94	0.00
31	Chrysene	1.299	1.504	-15.8	97	0.00
32	Bis(2-ethylhexyl)phthalate	1.604	1.403	12.5	81	-0.01
33	Indeno(1,2,3-cd)pyrene	1.535	1.698	-10.6	93	0.00
34 I	Perylene-d12	1.000	1.000	0.0	92	0.00
35	Benzo(b)fluoranthene	1.153	1.176	-2.0	92	0.00
36	Benzo(k)fluoranthene	1.249	1.273	-1.9	91	0.00
37 C	Benzo(a)pyrene	1.144	1.202	-5.1	96	0.00
38	Dibenzo(a,h)anthracene	1.116	1.132	-1.4	92	0.01
39	Benzo(a,h,i)perylene	1.181	1.213	-2.7	92	0.00

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE061719\
Data File : BE100069.D
Acq On : 17 Jun 2019 19:25
Operator : JU/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_E
LabSampleId :
SSTDCCC0.4

Quant Time: Jun 18 03:44:44 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE061719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jun 17 17:23:36 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			