

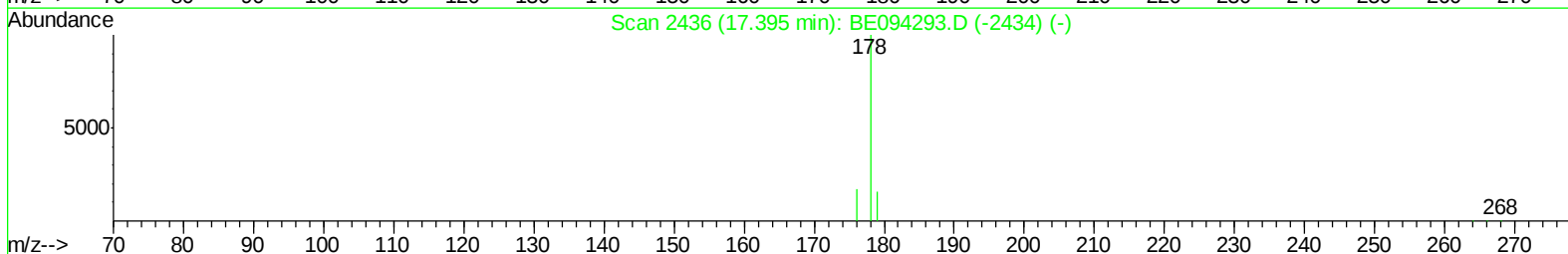
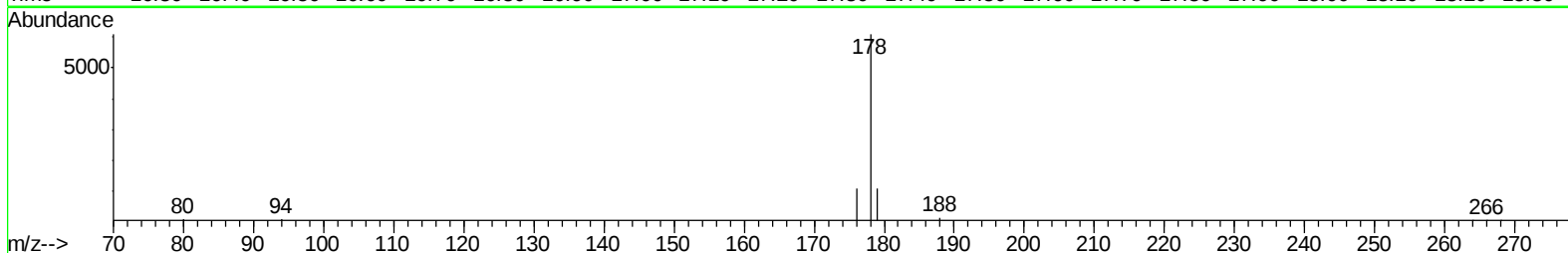
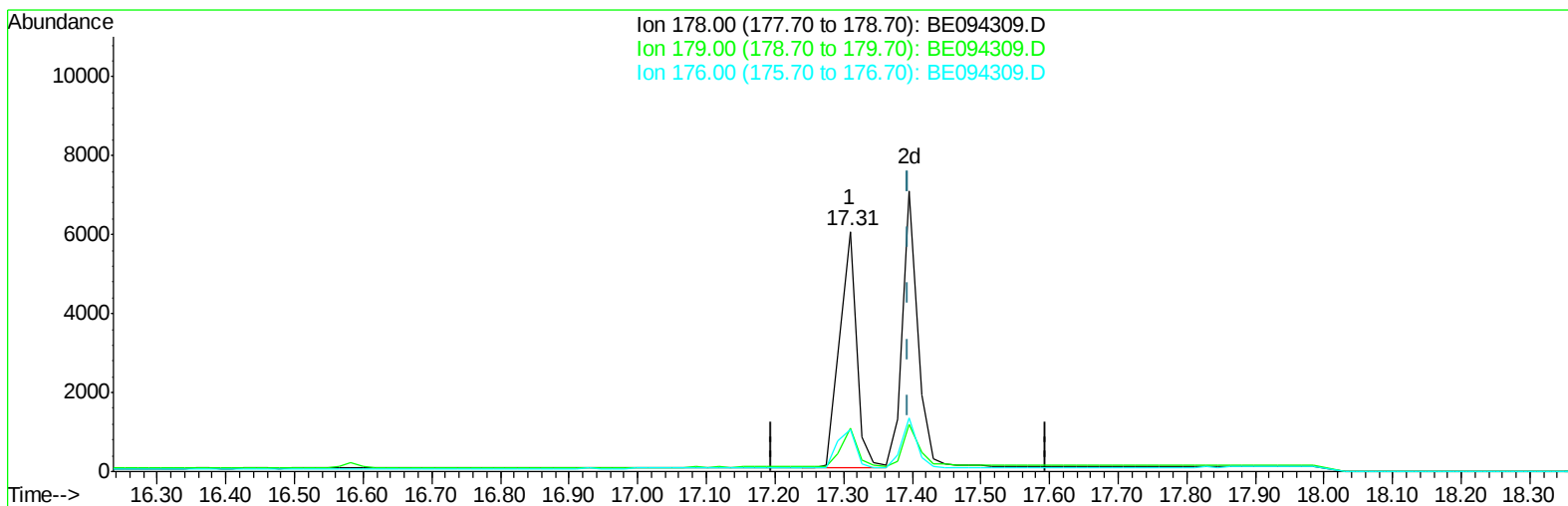
Data Path : Z:\HPCHEM1\BNA_E\Data\Be101117\
Data File : BE094309.D
Acq On : 12 Oct 2017 9:23
Operator : SJ/JU
Sample : SSTDCCC0.4EC
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_E
LabSampleId :
SSTD0.432

Manual Integrations
APPROVED

Sohil
10/13/2017 12:06:22 PM

Quant Time: Oct 12 10:27:20 2017
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE092017.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Oct 12 08:55:52 2017
Response via : Initial Calibration



TIC: BE094309.D

(13) Anthracene

17.309min (-0.086) 0.41ng/ul

response 10197

Ion	Exp%	Act%
178.00	100	100
179.00	22.60	18.33
176.00	18.40	17.74
0.00	0.00	0.00

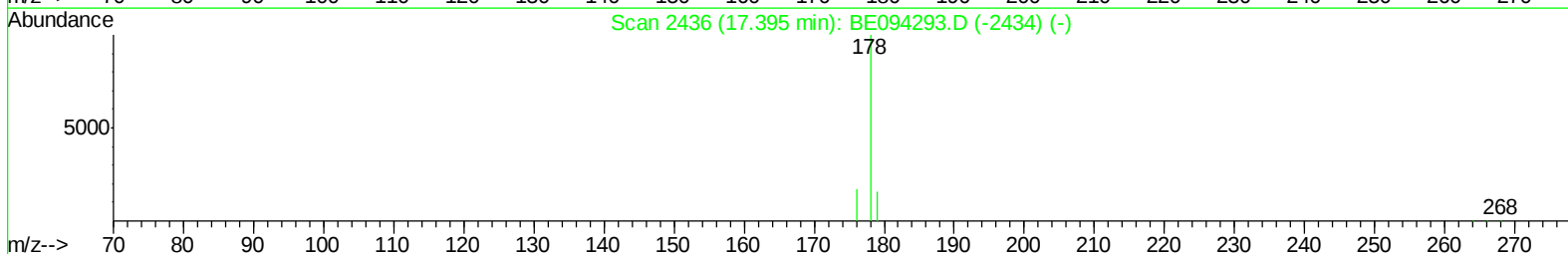
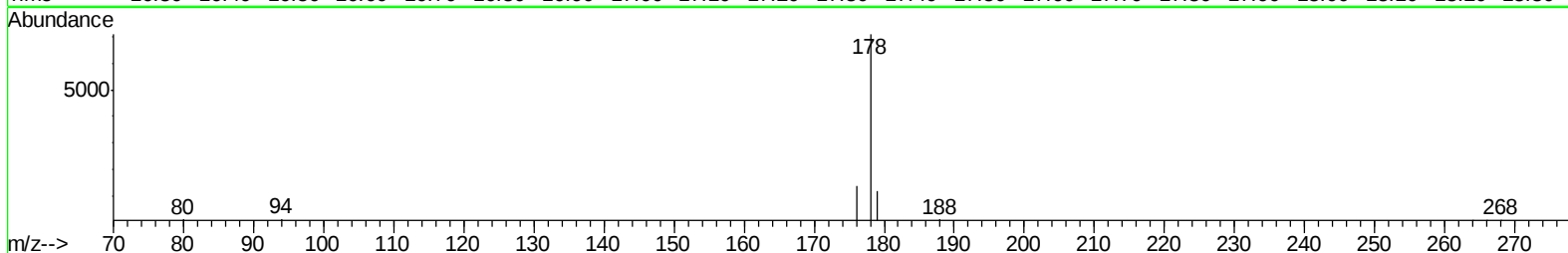
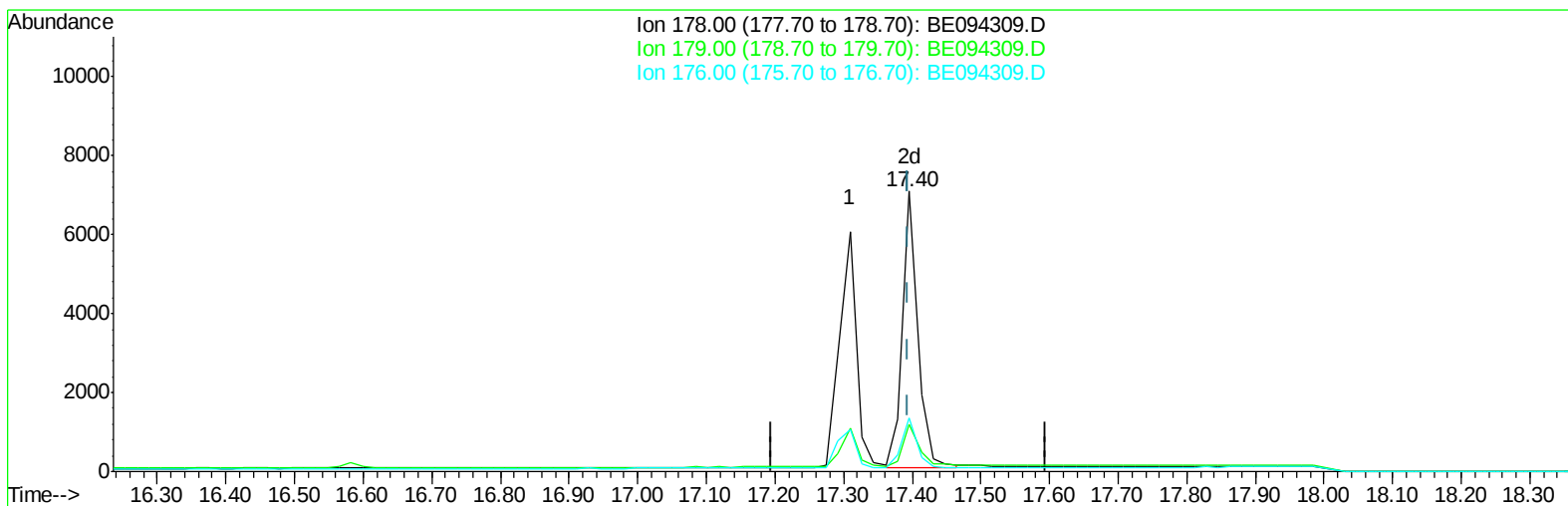
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Acq On : 12 Oct 2017 9:23
Operator : SJ/JU
Sample : SSTDCCC0.4EC
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_E
LabSampleId :
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Manual Integrations
APPROVED

Sohil
10/13/2017 12:06:22 PM

Quant Time: Oct 12 10:27:20 2017
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TIC: BE094309.D

(13) Anthracene

17.396min (+0.001) 0.43ng/ul m

response 10868

Ion	Exp%	Act%
178.00	100	100
179.00	22.60	16.72#
176.00	18.40	18.99
0.00	0.00	0.00

Data Path : Z:\HPCHEM1\BNA_E\Data\Be101117\
 Data File : BE094309.D
 Acq On : 12 Oct 2017 9:23
 Operator : SJ/JU
 Sample : SSTDC00.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTD0.432

Manual Integrations
 APPROVED

Sohil
 10/13/2017 12:06:22 PM

Quant Time: Oct 12 20:23:07 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE092017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 12 08:55:52 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.92	152	1760	0.40	ng/ul	0.00
2) Naphthalene-d8	10.71	136	8770	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.53	164	4536	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.27	188	9899	0.40	ng/ul	0.02
16) Chrysene-d12	21.43	240	9994	0.40	ng/ul	0.00
20) Perylene-d12	23.96	264	9791	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	5411	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.29	212	12270	0.37	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.76	128	8427	0.392	ng/ul#	93
5) 2-Methylnaphthalene	12.36	142	5719	0.410	ng/ul	98
7) Acenaphthylene	14.25	152	8543	0.483	ng/ul	97
8) Acenaphthene	14.59	153	5885	0.404	ng/ul	97
9) Fluorene	15.58	166	6623	0.388	ng/ul	94
11) Pentachlorophenol	16.96	266	249	0.214	ng/ul#	81
12) Phenanthrene	17.31	178	10248	0.396	ng/ul#	94
13) Anthracene	17.40	178	10868m	0.432	ng/ul	
15) Fluoranthene	19.32	202	12633	0.372	ng/ul	84
17) Pyrene	19.68	202	13171	0.474	ng/ul#	83
18) Benzo(a)anthracene	21.41	228	12180	0.459	ng/ul#	86
19) Chrysene	21.47	228	11341	0.370	ng/ul#	81
21) Benzo(b)fluoranthene	23.18	252	11653	0.438	ng/ul#	73
22) Benzo(k)fluoranthene	23.23	252	11060	0.360	ng/ul#	69
23) Benzo(a)pyrene	23.85	252	10913	0.395	ng/ul#	73
24) Indeno(1,2,3-cd)pyrene	26.65	276	10271	0.363	ng/ul#	80
25) Dibenzo(a,h)anthracene	26.66	278	8944	0.377	ng/ul#	59
26) Benzo(g,h,i)perylene	27.48	276	7742	0.317	ng/ul#	57

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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