

Instrument :
BNA_M
ClientSampleId :
F9B07

Sohil
12/11/2017 6:16:24 PM

Abundance

TIC: BM013294.D

Time-->

1,4-Dioxane-d8,S

Phenol

Bis(2-chlorophenyl)ether-d8,S

2-Chlorophenol-d4,S

1,4-Dichlorobenzene-d4,l

4-Methylphenol-d8,S

Nitrobenzene-d5,S

2-Nitrophenol-d4,S

2,4-Dichlorophenol-d3,S

Naphthalene-d8,l

4-Chloroaniline-d4,S

Dimethylprthalate

Dimethylprthalate-d6,S

Acenaphthylene

Acenaphthylene-d8,S

Acenaphthene-d10,l

4-Nitrophenol-d4,S

4,6-Dinitro-2-methylphenol-d2,S

Fluorene-d10,S

Phenanthrene

Phenanthrene-d10,S

Fluoranthene,C

Pyrene

Pyrene-d10,S

Chrysene

Benzo(a)anthracene-d12,l

Benzo(b)fluoranthene

Benzo(a)pyrene-d12,S

Benzo(a)pyrene-d12,l

Dibenz(a,h)pyrene

Benzo(g,h,i)perylene

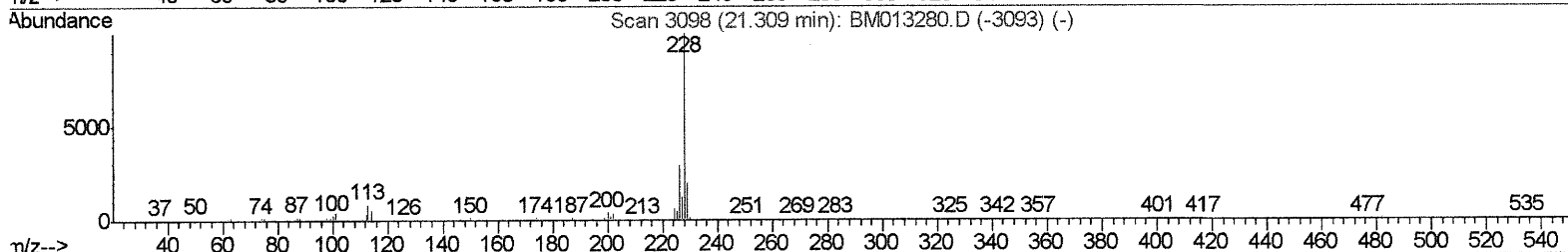
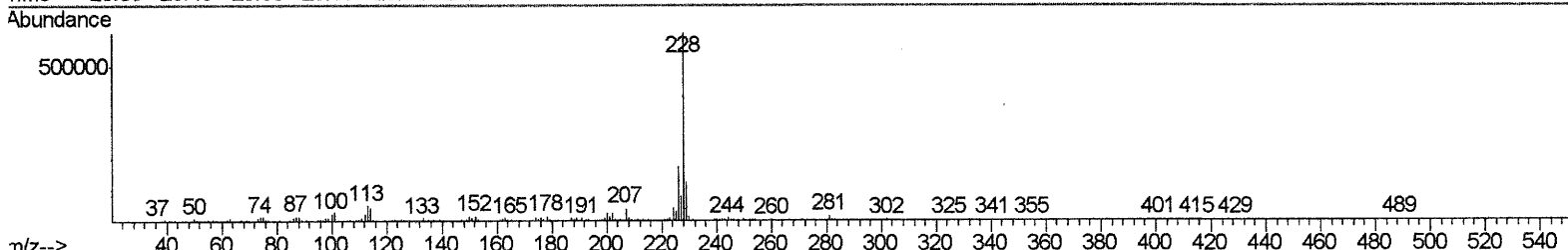
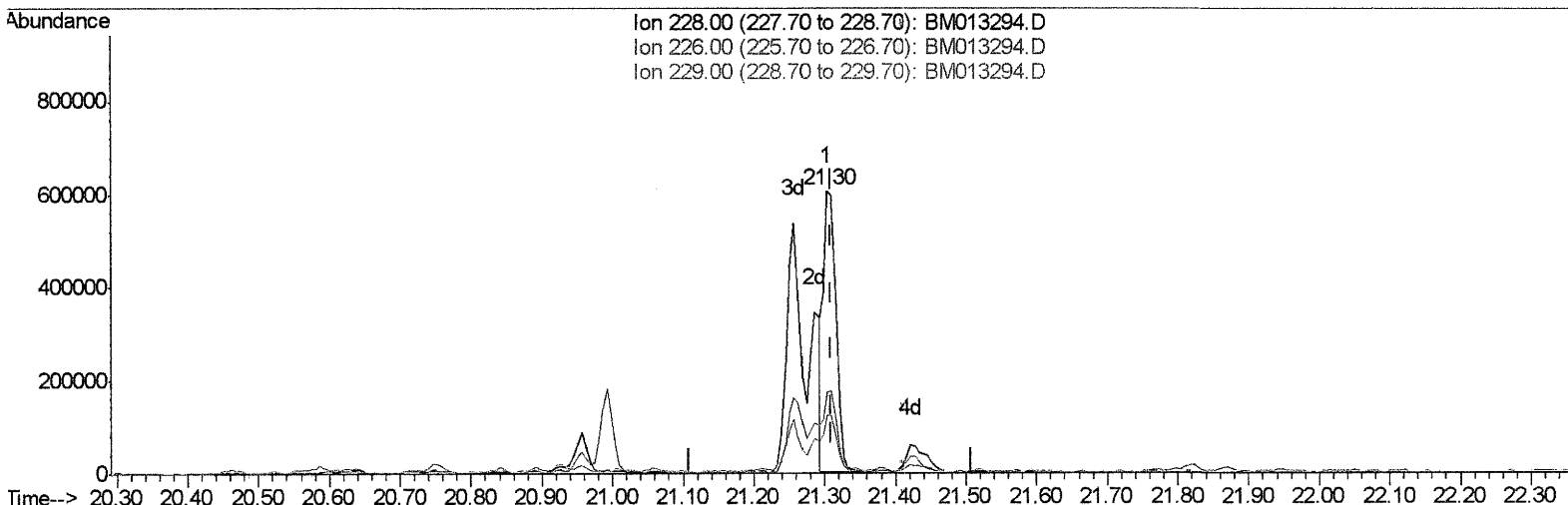
Data Path : Z:\HPCHEM1\BNA M\DATA\BM120817\
 Data File : BM013294.D
 Acq On : 09 Dec 2017 06:32
 Operator : SJ/JU
 Sample : I6758-06
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampled :
 F9B07

Manual Integrations
 APPROVED

Sohil
 12/11/2017 6:16:24 PM

Quant Time: Dec 11 00:50:01 2017
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM113017.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Dec 08 23:45:50 2017
 Response via : Initial Calibration



TIC: BM013294.D

(82) Chrysene

21.303min (-0.006) 8.12ng/ui

response 755615

Ion	Exp%	Act%
228.00	100	100
226.00	29.30	28.72
229.00	19.40	20.37
0.00	0.00	0.00

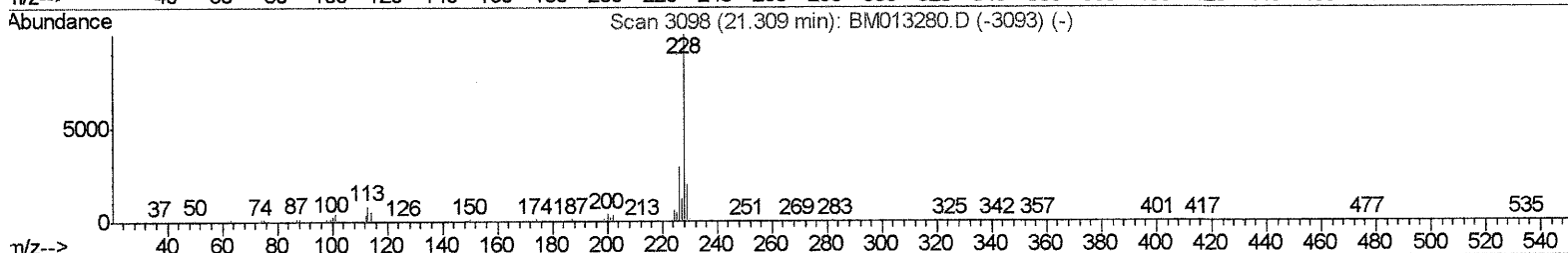
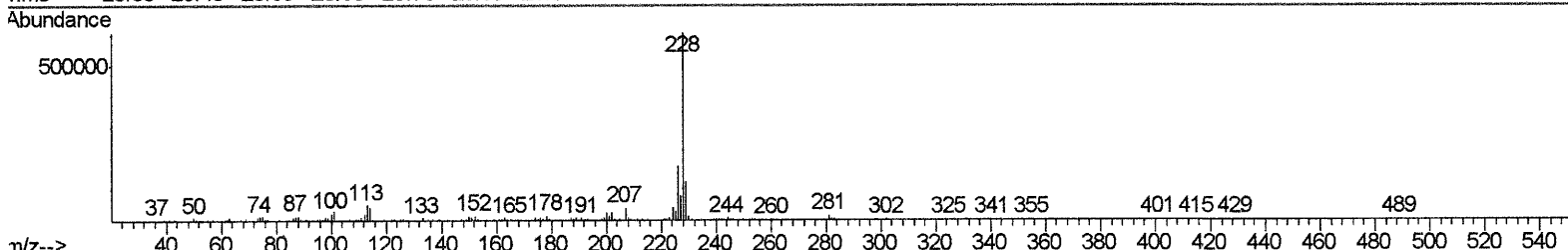
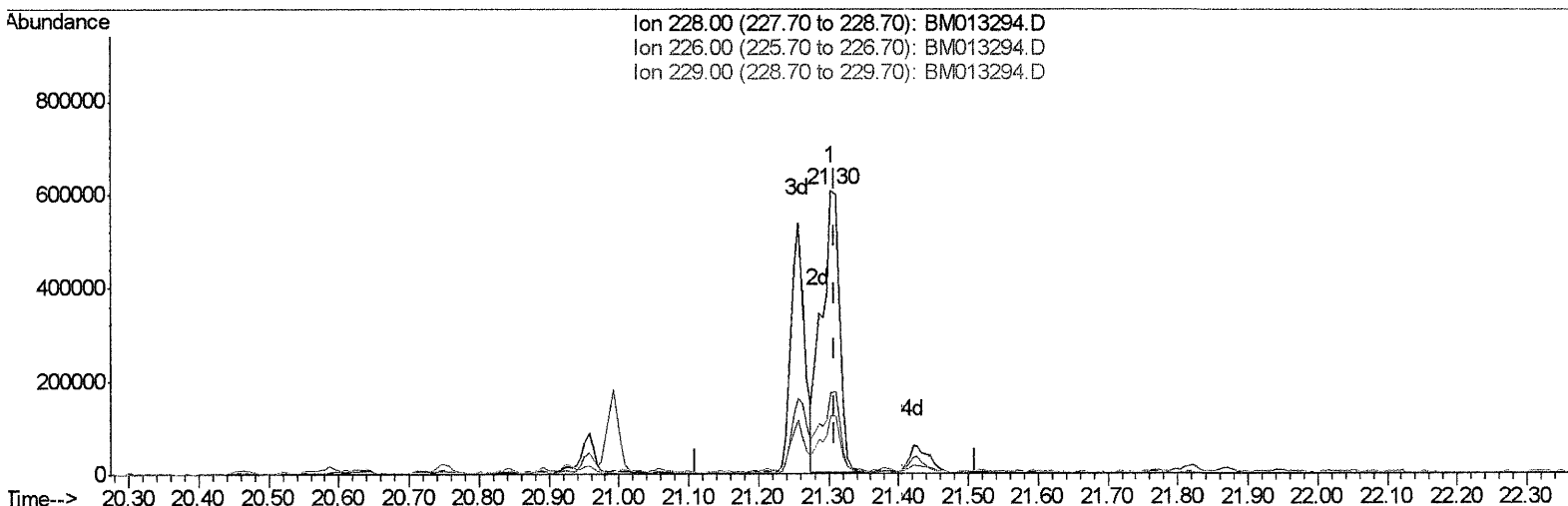
Data Path : Z:\HPCHEM1\BNA M\DATA\BM120817\
 Data File : BM013294.D
 Acq On : 09 Dec 2017 06:32
 Operator : SJ/JU
 Sample : I6758-06
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
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 ClientSampled :
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Manual Integrations
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 Response via : Initial Calibration



TIC: BM013294.D

(82) Chrysene

21.303min (-0.006) 11.58ng/ul m> SJ 12/11/17

response 1077553

Ion	Exp%	Act%
228.00	100	100
226.00	29.30	28.72
229.00	19.40	20.37
0.00	0.00	0.00

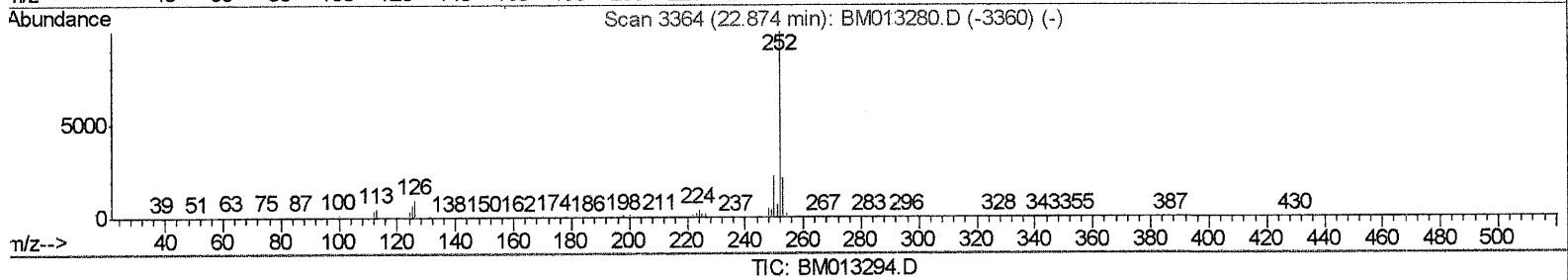
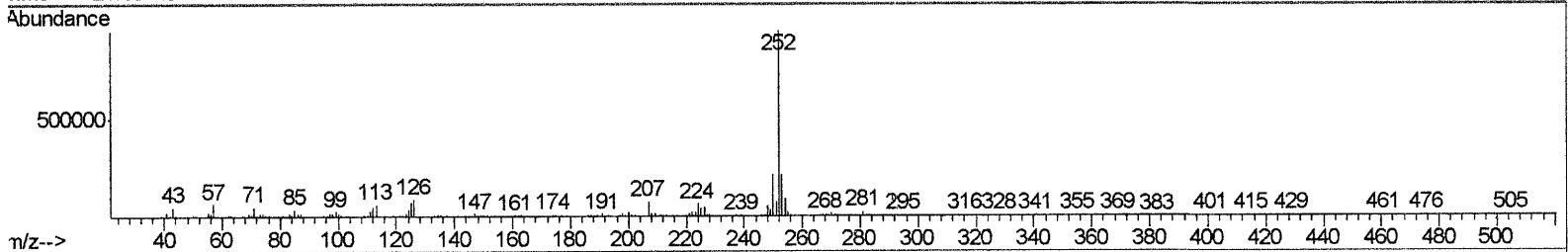
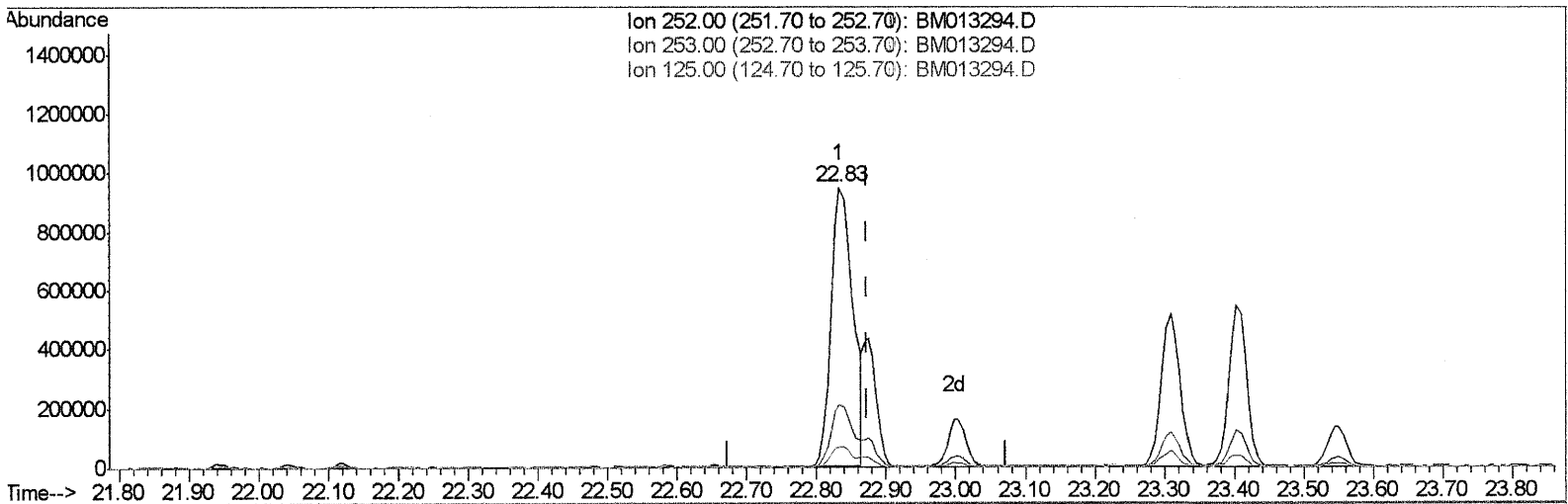
Data Path : Z:\HPCHEM1\BNA M\DATA\BM120817\
 Data File : BM013294.D
 Acq On : 09 Dec 2017 06:32
 Operator : SJ/JU
 Sample : I6758-06
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
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Quant Time: Dec 11 00:50:01 2017
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(86) Benzo(k)fluoranthene
 22.833min (-0.041) 23.65ng/ul
 response 2050925

Ion	Exp%	Act%
252.00	100	100
253.00	21.40	22.41
125.00	4.30	7.48#
0.00	0.00	0.00

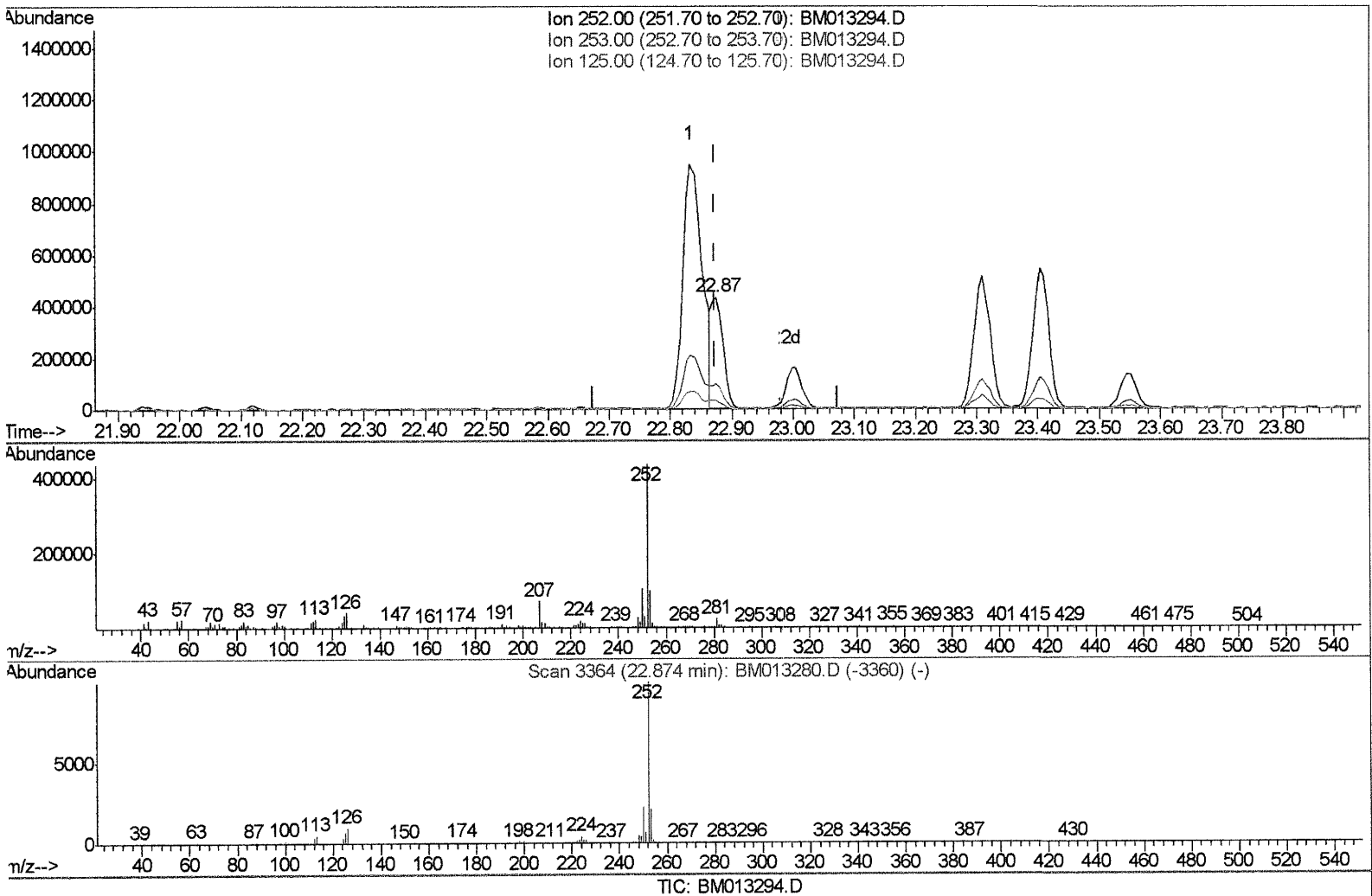
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 Data File : BM013294.D
 Acq On : 09 Dec 2017 06:32
 Operator : SJ/JU
 Sample : I6758-06
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampled :
 F9B07

Manual Integrations
 APPROVED

Sohil
 12/11/2017 6:16:24 PM

Quant Time: Dec 11 00:50:01 2017
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM113017.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Dec 08 23:45:50 2017
 Response via : Initial Calibration



(86) Benzo(k)fluoranthene

22.874min (+0.000) 6.69ng/ul m> SJ 12/11/17

response 579760

Ion	Exp%	Act%
252.00	100	100
253.00	21.40	22.62
125.00	4.30	7.69#
0.00	0.00	0.00

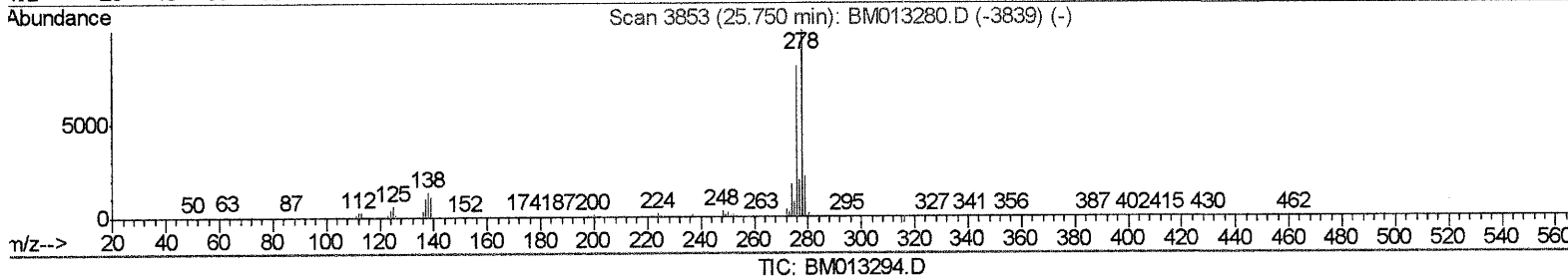
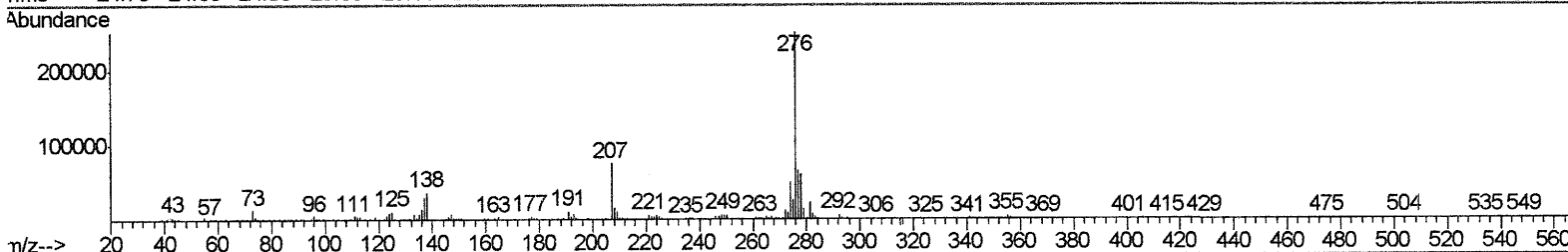
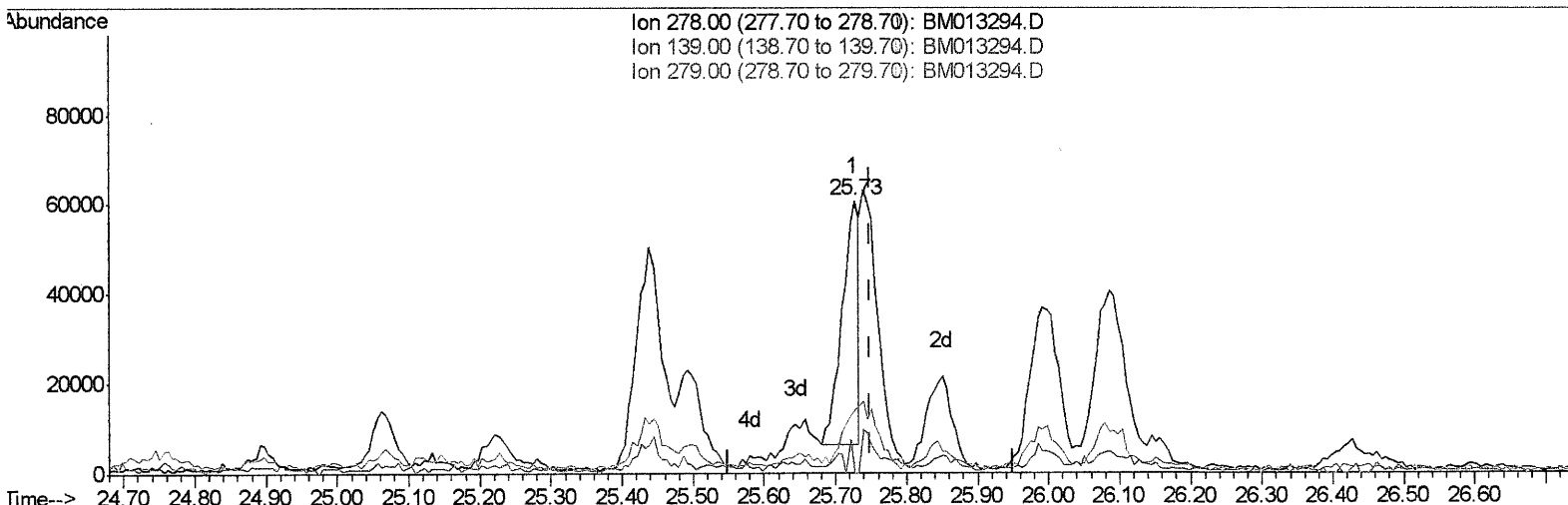
Data Path : Z:\HPCHEM1\BNA M\DATA\BM120817\
 Data File : BM013294.D
 Acq On : 09 Dec 2017 06:32
 Operator : SJ/JU
 Sample : I6758-06
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampled :
 F9B07

Manual Integrations
 APPROVED

Sohil
 12/11/2017 6:16:24 PM

Quant Time: Dec 11 00:50:01 2017
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM113017.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Dec 08 23:45:50 2017
 Response via : Initial Calibration



(90) Dibenzo(a,h)anthracene

25.727min (-0.023) 1.31ng/ul

response 91736

Ion	Exp%	Act%
278.00	100	100
139.00	7.30	0.00#
279.00	23.20	23.47
0.00	0.00	0.00

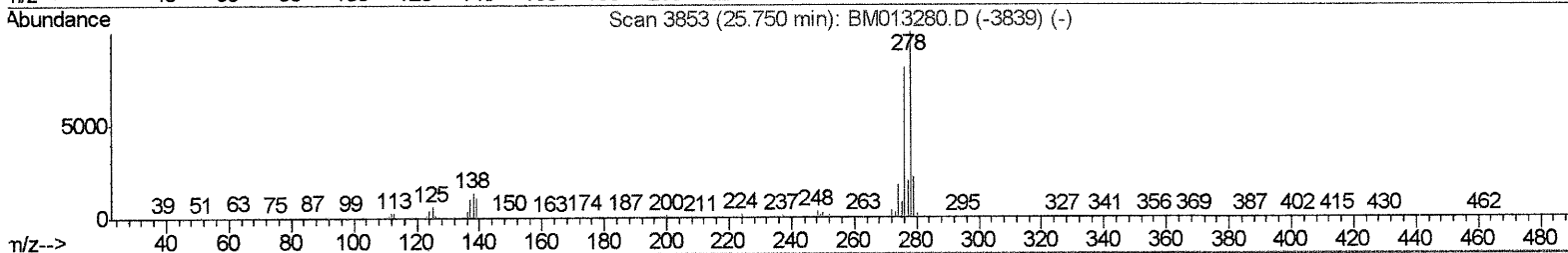
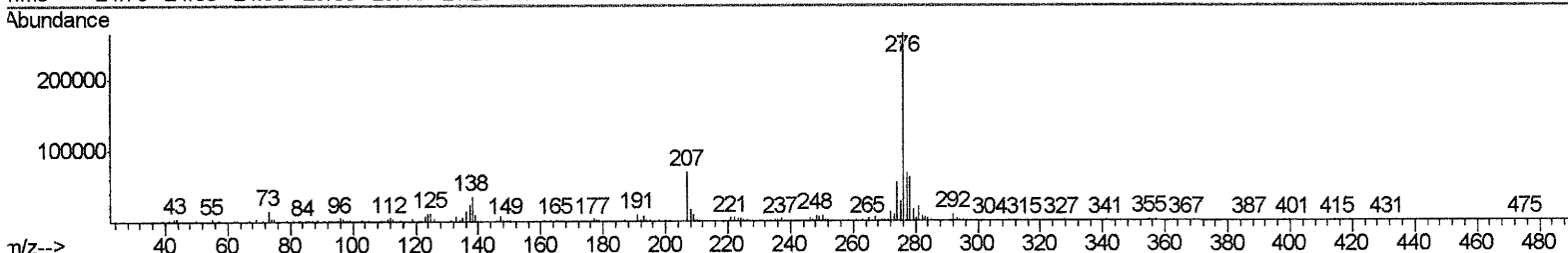
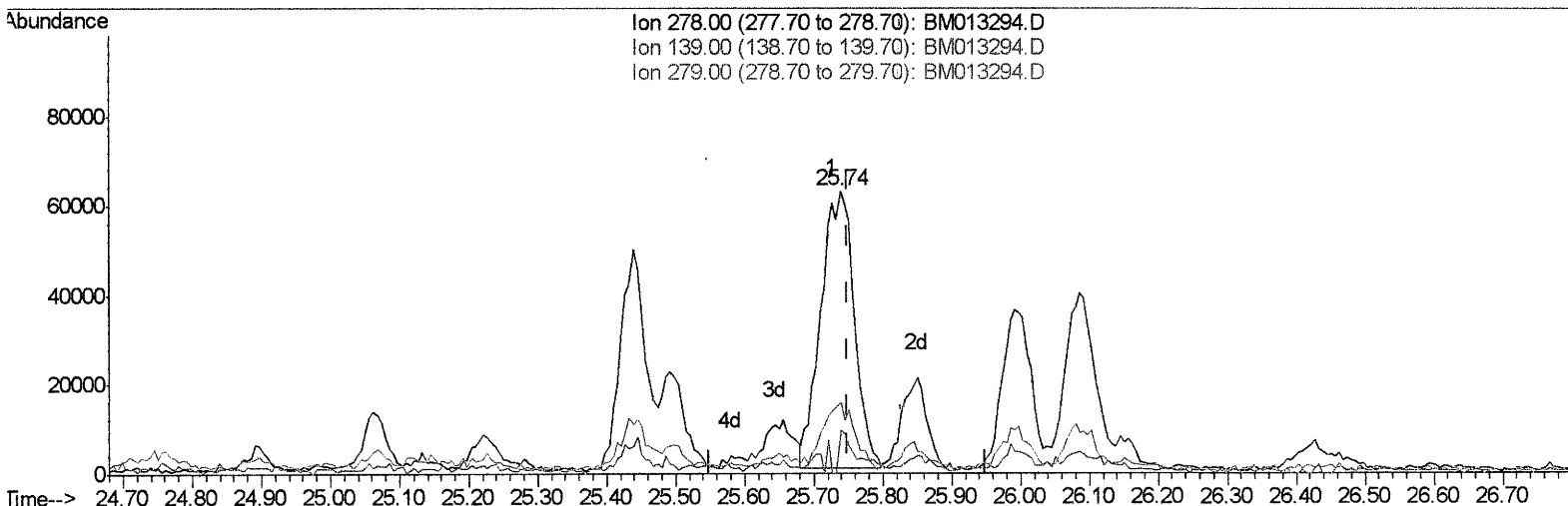
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 Data File : BM013294.D
 Acq On : 09 Dec 2017 06:32
 Operator : SJ/JU
 Sample : I6758-06
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampled :
 F9B07

Manual Integrations
 APPROVED

Sohil
 12/11/2017 6:16:24 PM

Quant Time: Dec 11 00:50:01 2017
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM113017.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Dec 08 23:45:50 2017
 Response via : Initial Calibration



TIC: BM013294.D

(90) Dibenzo(a,h)anthracene

25.738min (-0.012) 3.01ng/ul m> 50 12/11/17

response 210228

Ion	Exp%	Act%
278.00	100	100
139.00	7.30	15.34#
279.00	23.20	25.69
0.00	0.00	0.00

Data Path : Z:\HPCHEM1\BNA M\DATA\BM120817\
 Data File : BM013294.D
 Acq On : 09 Dec 2017 06:32
 Operator : SJ/JU
 Sample : I6758-06
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampled :
 F9B07

Manual Integrations
 APPROVED

Sohil
 12/11/2017 6:16:24 PM

Quant Time: Dec 11 02:03:40 2017
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM113017.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Dec 08 23:45:50 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.69	152	233887	20.00	ng/ul	0.00
18) Naphthalene-d8	10.47	136	1008004	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.33	164	616416	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.08	188	1431248	20.00	ng/ul	0.00
75) Chrysene-d12	21.27	240	1580758	20.00	ng/ul	0.00
83) Perylene-d12	23.50	264	1366876	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.23	96	11165	2.36	ng/uL	0.00
5) Phenol-d5	6.87	99	350145	17.30	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.03	67	237360	20.57	ng/ul	0.00
9) 2-Chlorophenol-d4	7.23	132	307375	19.44	ng/ul	0.00
13) 4-Methylphenol-d8	8.40	113	306015	17.70	ng/ul	0.00
19) Nitrobenzene-d5	8.85	128	177912	22.26	ng/ul	0.00
22) 2-Nitrophenol-d4	9.56	143	178010	20.81	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.10	165	329505	18.69	ng/ul	0.00
29) 4-Chloroaniline-d4	10.61	131	150842	9.31	ng/ul	0.00
43) Dimethylphthalate-d6	13.75	166	899698	16.40	ng/ul	0.00
46) Acenaphthylene-d8	14.02	160	1472259	21.75	ng/ul	0.00
51) 4-Nitrophenol-d4	14.55	143	63983	6.71	ng/ul	0.00
57) Fluorene-d10	15.33	176	1024432	21.72	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.45	200	67533	7.50	ng/ul	0.00
70) Anthracene-d10	17.18	188	1565797	21.84	ng/ul	0.00
76) Pyrene-d10	19.48	212	1716281	21.79	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.36	264	1420693	20.41	ng/ul	0.00

Target Compounds

					Ovalue	
6) Phenol	6.90	94	29983	1.499	ng/ul#	77
44) Dimethylphthalate	13.79	163	88079	1.679	ng/ul	96
47) Acenaphthylene	14.05	152	242492	3.871	ng/ul	97
71) Anthracene	17.22	178	212432	2.556	ng/ul	99
74) Fluoranthene	19.14	202	1286175	12.907	ng/ul#	94
77) Pyrene	19.51	202	1751460	18.238	ng/ul#	92
80) Benzo(a)anthracene	21.26	228	723512	7.214	ng/ul	94
82) Chrysene	21.30	228	1077553m)	11.581	ng/ul	94
85) Benzo(b)fluoranthene	22.83	252	2050925	22.258	ng/ul#	98
86) Benzo(k)fluoranthene	22.87	252	579760m)	6.686	ng/ul	98
88) Benzo(a)pyrene	23.40	252	967616	11.690	ng/ul#	99
89) Indeno(1,2,3-cd)pyrene	25.73	276	740221	9.013	ng/ul#	95
90) Dibenzo(a,h)anthracene	25.74	278	210228m)	3.006	ng/ul	95
91) Benzo(a,h,i)perylene	26.42	276	580289	8.958	ng/ul#	95

SD 12/11/17
 SD 12/11/17
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(#) = qualifier out of range (m) = manual integration (+) = signals summed