

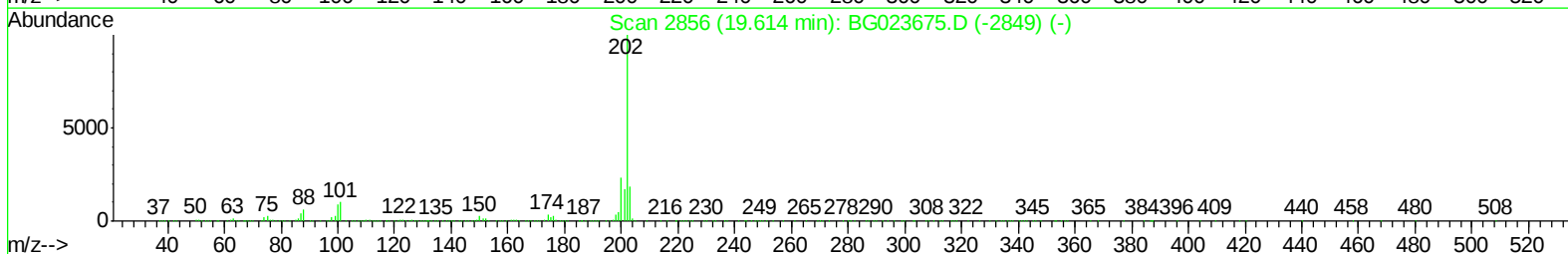
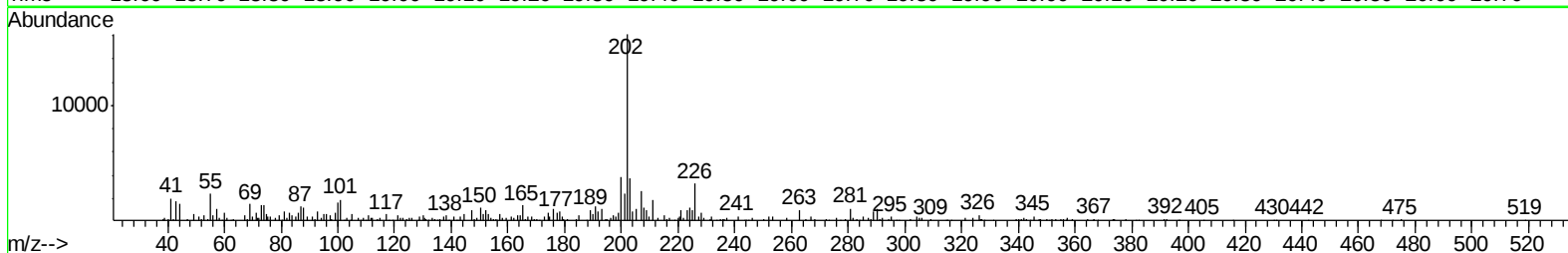
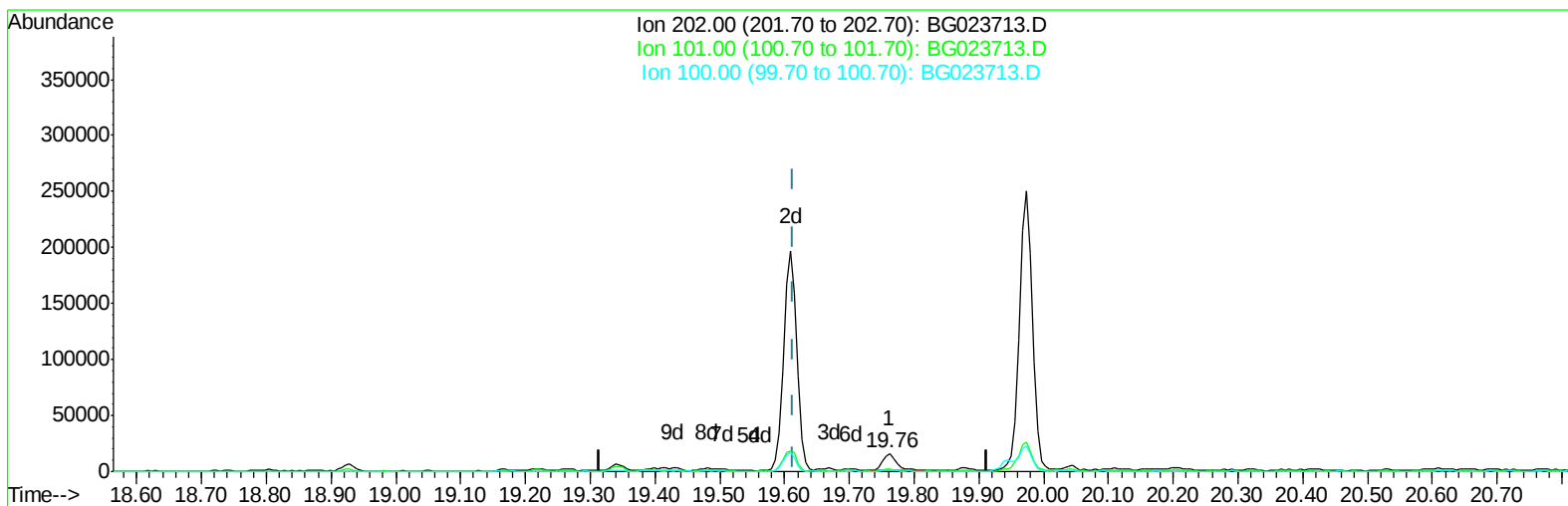
Data Path : Z:\HPCHEM1\BNA_G\Data\BG081316\
 Data File : BG023713.D
 Acq On : 14 Aug 2016 7:04
 Operator : UM/SJ
 Sample : H4388-08
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PR002-SS002-1218-01

Manual Integrations
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Quant Time: Aug 15 19:23:22 2016
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG080416.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Aug 15 18:52:38 2016
 Response via : Initial Calibration



TIC: BG023713.D

(74) Fluoranthene (C)

19.762min (+0.148) 0.58ng/ul

response 23286

Ion	Exp%	Act%
202.00	100	100
101.00	13.30	12.05
100.00	10.90	10.12
0.00	0.00	0.00

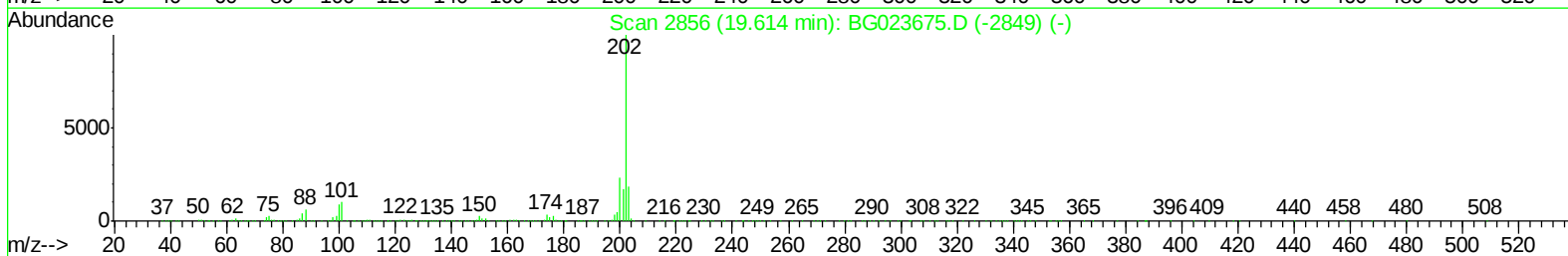
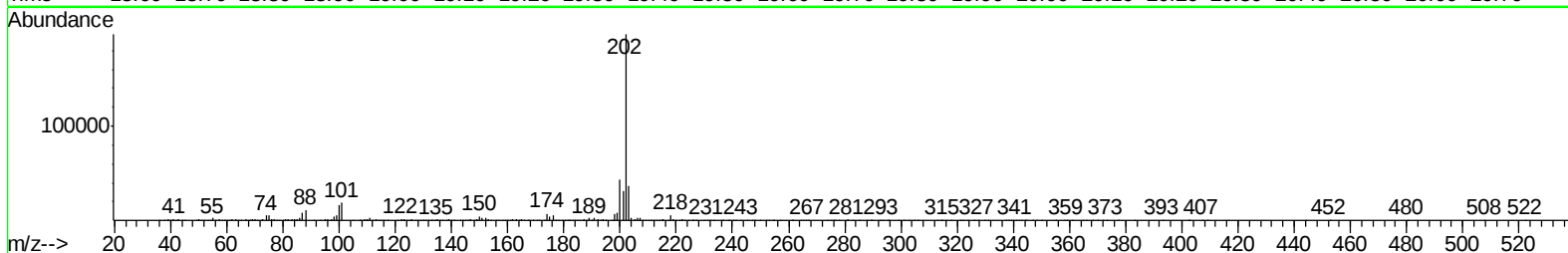
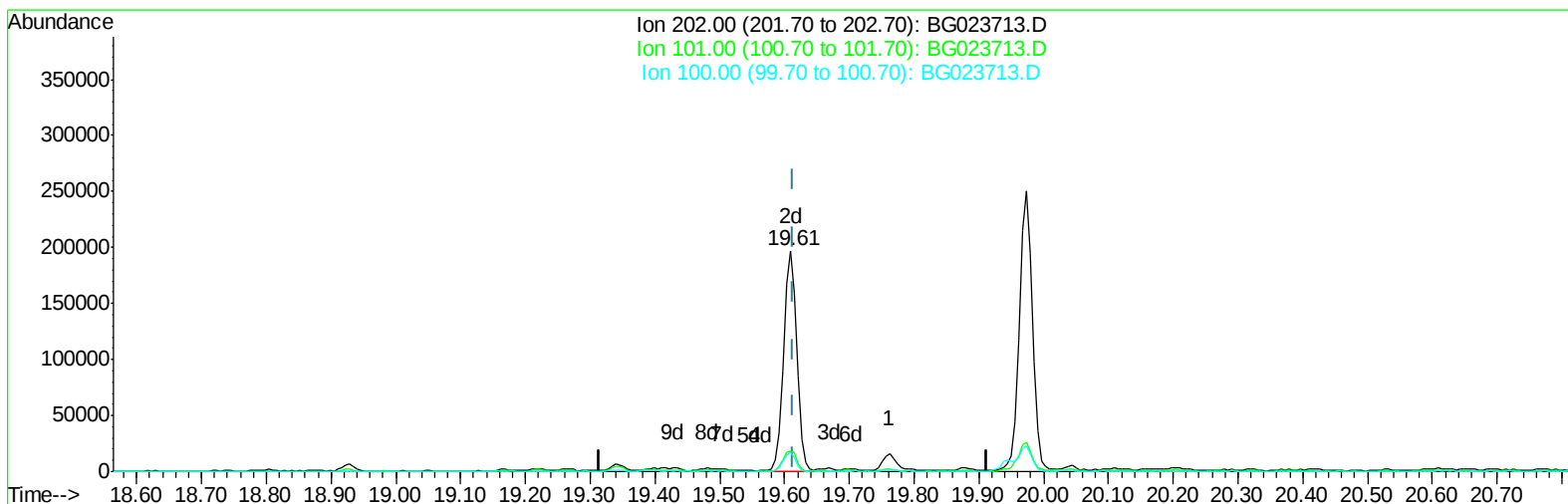
Data Path : Z:\HPCHEM1\BNA_G\Data\BG081316\
 Data File : BG023713.D
 Acq On : 14 Aug 2016 7:04
 Operator : UM/SJ
 Sample : H4388-08
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
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Manual Integrations
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 Quant Title : SVOA CALIBRATION
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TIC: BG023713.D

(74) Fluoranthene (C)

19.610min (-0.004) 6.90ng/ul m

response 277197

Ion	Exp%	Act%
202.00	100	100
101.00	13.30	9.48#
100.00	10.90	8.57#
0.00	0.00	0.00

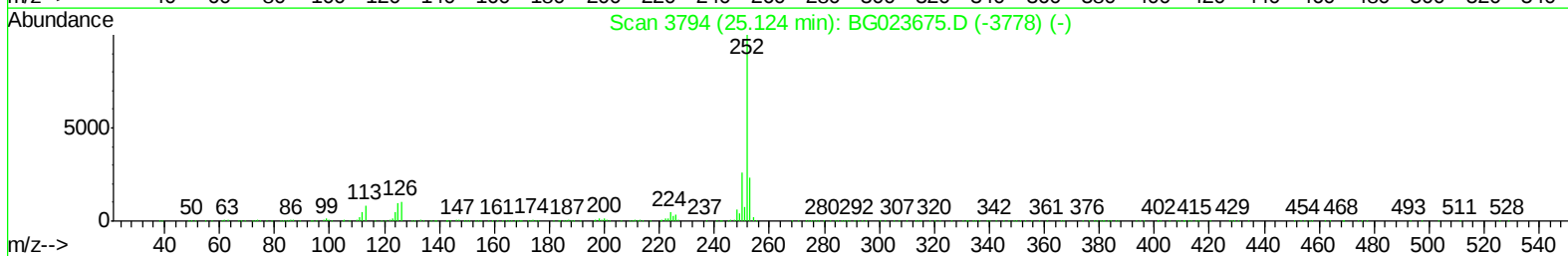
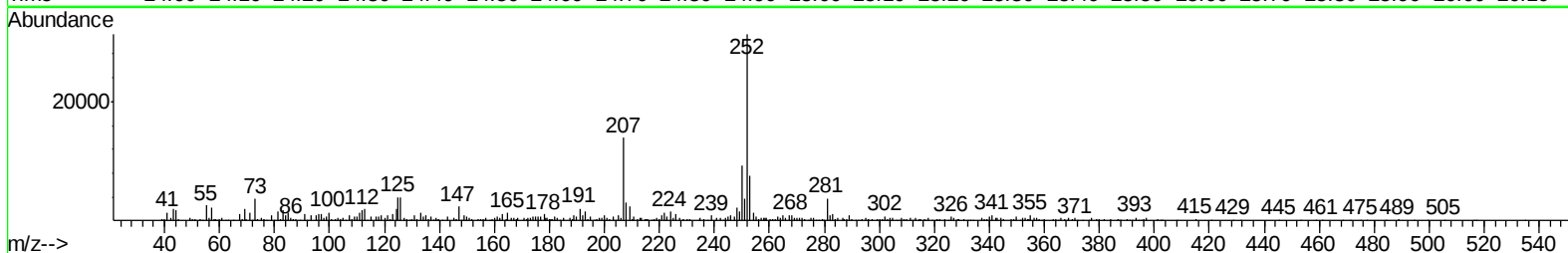
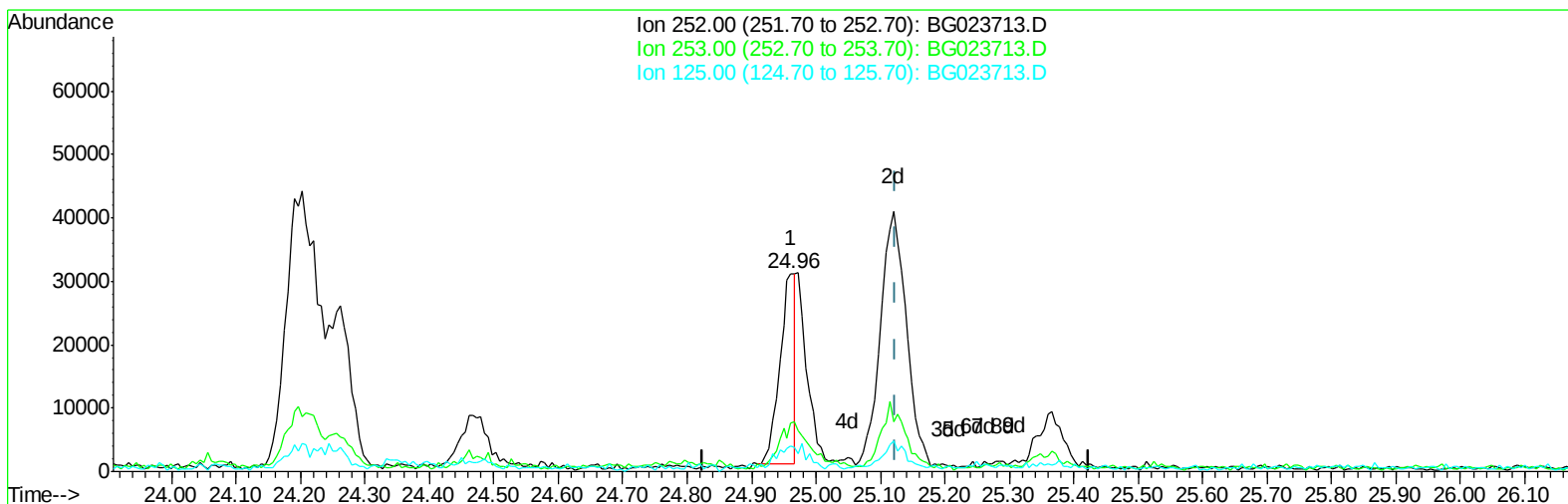
Data Path : Z:\HPCHEM1\BNA_G\Data\BG081316\
Data File : BG023713.D
Acq On : 14 Aug 2016 7:04
Operator : UM/SJ
Sample : H4388-08
Misc :
ALS Vial : 32 Sample Multiplier: 1

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ClientSampleId :
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Manual Integrations
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TIC: BG023713.D

(88) Benzo(a)pyrene (C)

24.961min (-0.163) 1.17ng/ul

response 52192

Ion	Exp%	Act%
252.00	100	100
253.00	23.70	24.49
125.00	15.00	12.90
0.00	0.00	0.00

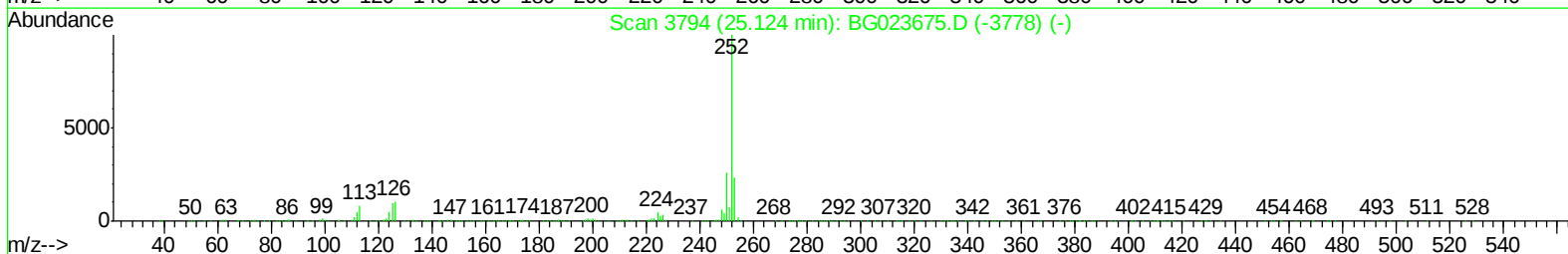
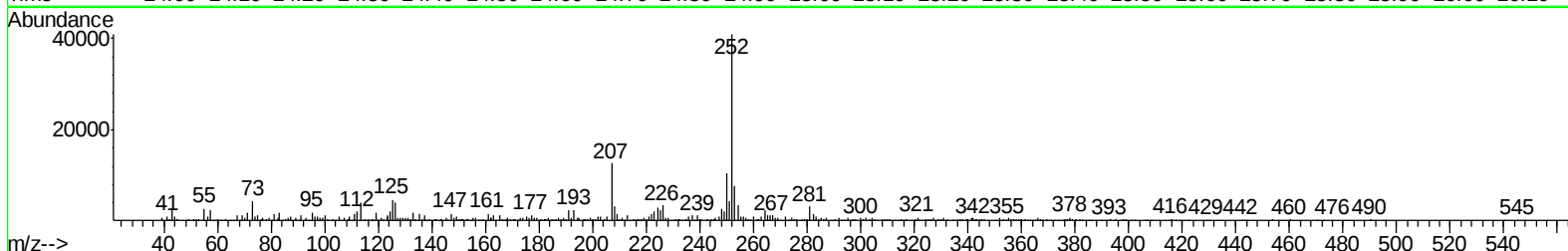
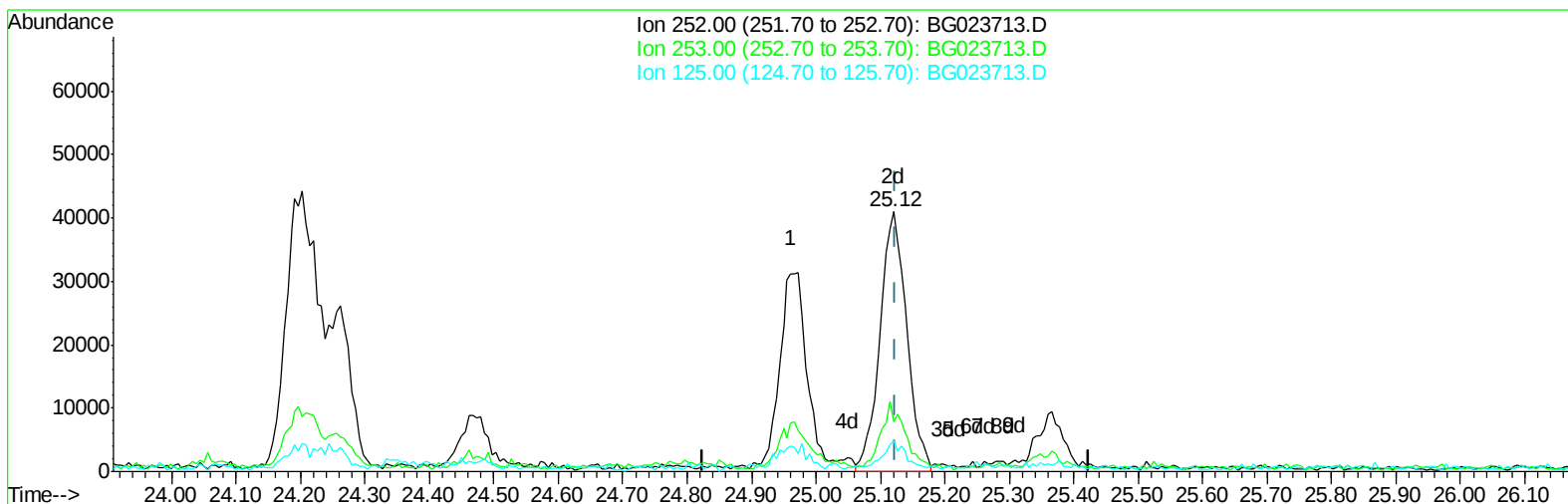
Data Path : Z:\HPCHEM1\BNA_G\Data\BG081316\
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TIC: BG023713.D

(88) Benzo(a)pyrene (C)

25.120min (-0.004) 2.69ng/ul m

response 119637

Ion	Exp%	Act%
252.00	100	100
253.00	23.70	19.14
125.00	15.00	11.62#
0.00	0.00	0.00

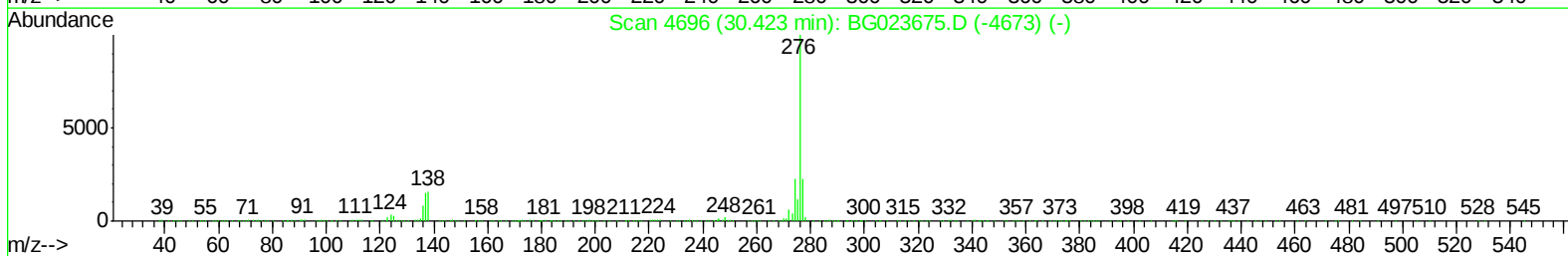
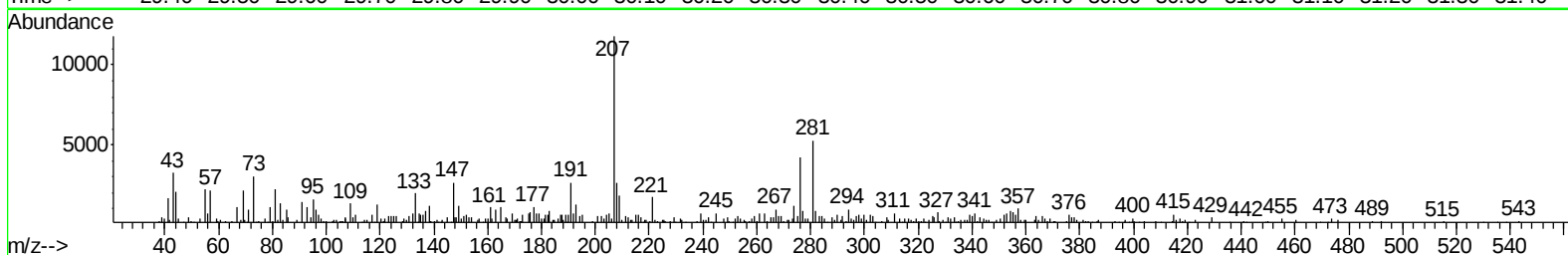
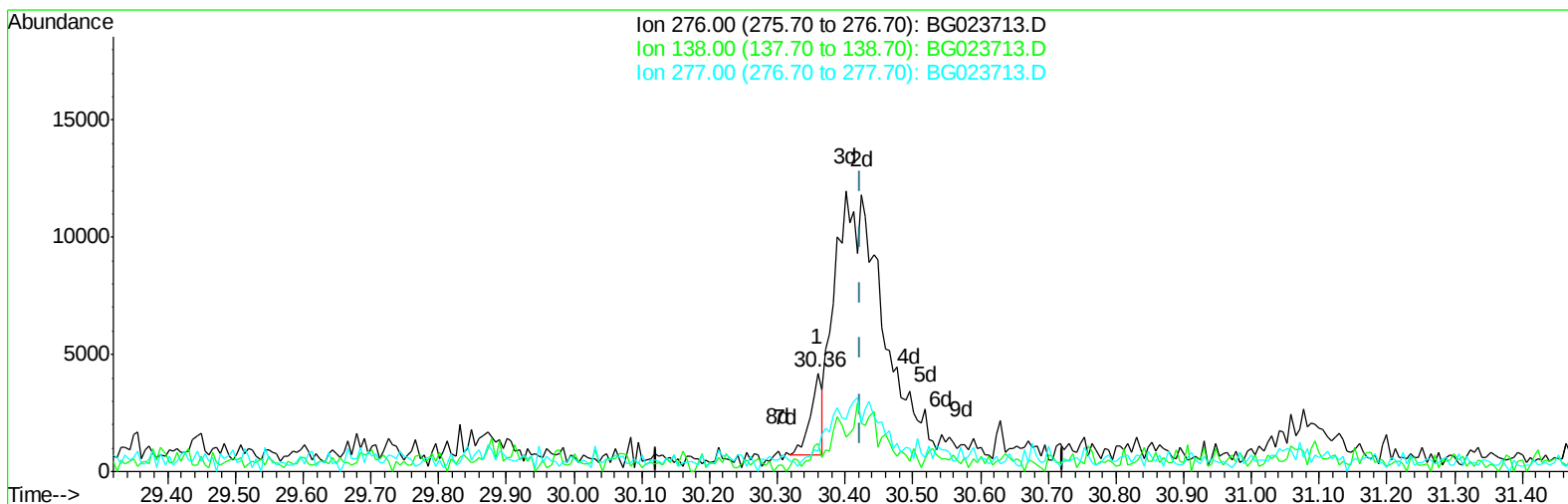
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Manual Integrations
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TIC: BG023713.D

(91) Benzo(g,h,i)perylene
30.360min (-0.063) 0.10ng/ul
response 4256

Ion	Exp%	Act%
276.00	100	100
138.00	27.60	28.55
277.00	21.20	20.65
0.00	0.00	0.00

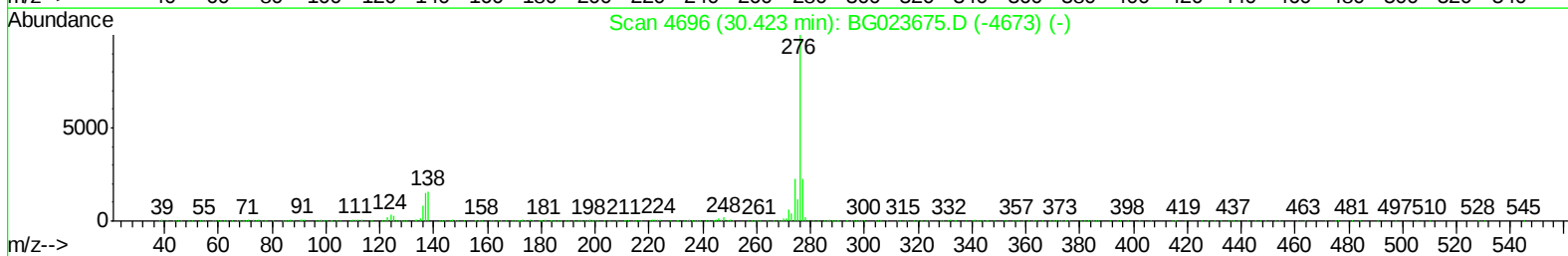
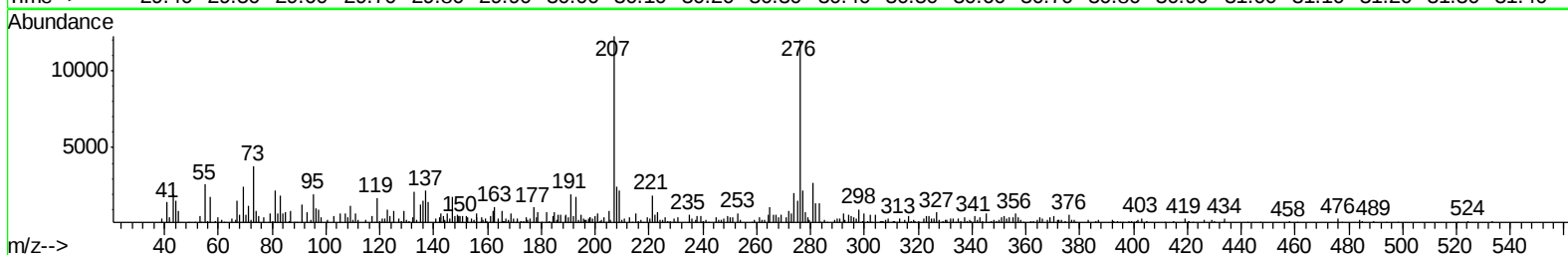
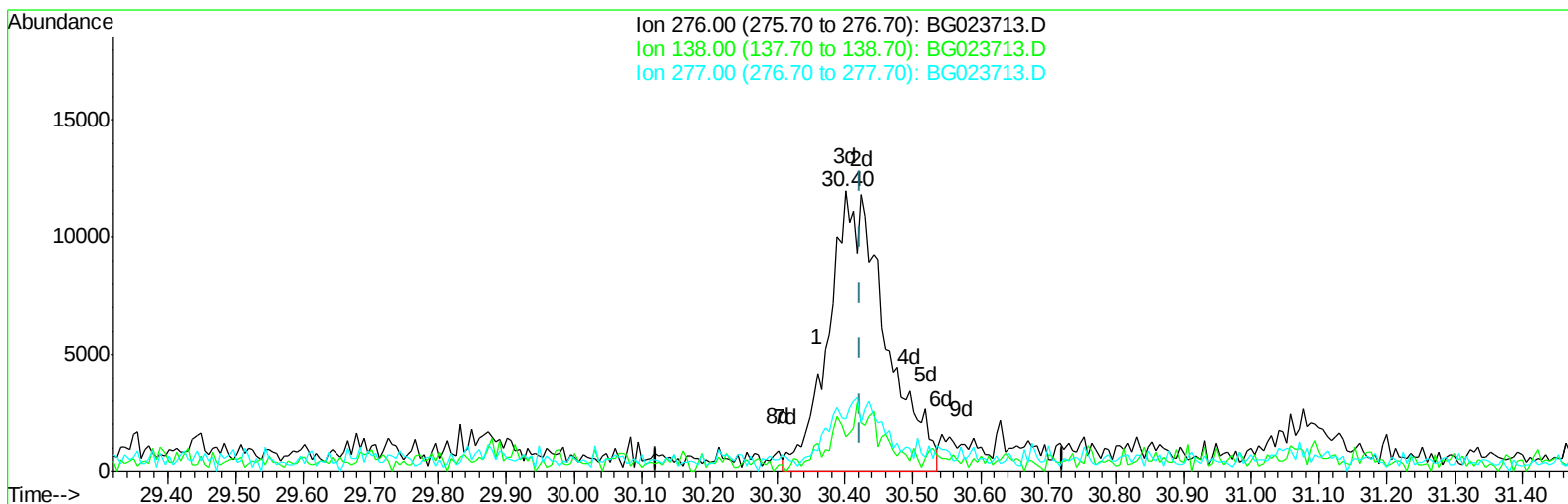
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Data File : BG023713.D
Acq On : 14 Aug 2016 7:04
Operator : UM/SJ
Sample : H4388-08
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
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Manual Integrations
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TIC: BG023713.D

(91) Benzo(g,h,i)perylene

30.401min (-0.022) 1.62ng/ul m

response 69950

Ion	Exp%	Act%
276.00	100	100
138.00	27.60	12.33#
277.00	21.20	18.63
0.00	0.00	0.00

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 Operator : UM/SJ
 Sample : H4388-08
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
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Manual Integrations
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Quant Time: Aug 15 19:25:22 2016
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG080416.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Aug 15 18:52:38 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.11	152	124674	20.00	ng/ul	0.00
18) Naphthalene-d8	10.94	136	552982	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.79	164	354065	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.55	188	758158	20.00	ng/ul	0.00
75) Chrysene-d12	21.88	240	818205	20.00	ng/ul	0.00
83) Perylene-d12	25.29	264	809010	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.48	96	7572	2.58	ng/uL	0.00
5) Phenol-d5	7.27	99	205210	16.34	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.43	67	142746	17.74	ng/ul	0.00
9) 2-Chlorophenol-d4	7.64	132	150896	17.61	ng/ul	0.00
13) 4-Methylphenol-d8	8.83	113	158026	16.12	ng/ul	0.00
19) Nitrobenzene-d5	9.35	128	590	0.14	ng/ul	0.06
22) 2-Nitrophenol-d4	10.02	143	96038	19.18	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.57	165	166086	17.59	ng/ul	0.00
29) 4-Chloroaniline-d4	11.09	131	170749	15.41	ng/ul	0.00
43) Dimethylphthalate-d6	14.18	166	556847	19.21	ng/ul	0.00
46) Acenaphthylene-d8	14.48	160	639695	19.61	ng/ul	0.00
51) 4-Nitrophenol-d4	15.01	143	77420	15.64	ng/ul	0.00
57) Fluorene-d10	15.79	176	483524	18.05	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.91	200	77201	15.87	ng/ul	0.00
70) Anthracene-d10	17.65	188	690570	20.22	ng/ul	0.00
76) Pyrene-d10	19.94	212	765475	18.48	ng/ul	0.00
87) Benzo(a)pyrene-d12	25.05	264	726204	19.83	ng/ul	0.00

Target Compounds

						Qvalue
44) Dimethylphthalate	14.23	163	249041	8.68	ng/ul	97
69) Phenanthrene	17.59	178	197617	5.03	ng/ul	99
74) Fluoranthene	19.61	202	277197m	6.90	ng/ul	
77) Pyrene	19.97	202	348355	6.89	ng/ul#	89
80) Benzo(a)anthracene	21.85	228	129502	2.81	ng/ul	98
81) Bis(2-ethylhexyl)phthalate	21.73	149	31966	1.24	ng/ul#	98
82) Chrysene	21.92	228	164124	3.92	ng/ul	97
85) Benzo(b)fluoranthene	24.20	252	151087	3.28	ng/ul#	93
88) Benzo(a)pyrene	25.12	252	119637m	2.69	ng/ul	
91) Benzo(g,h,i)perylene	30.40	276	69950m	1.62	ng/ul	

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

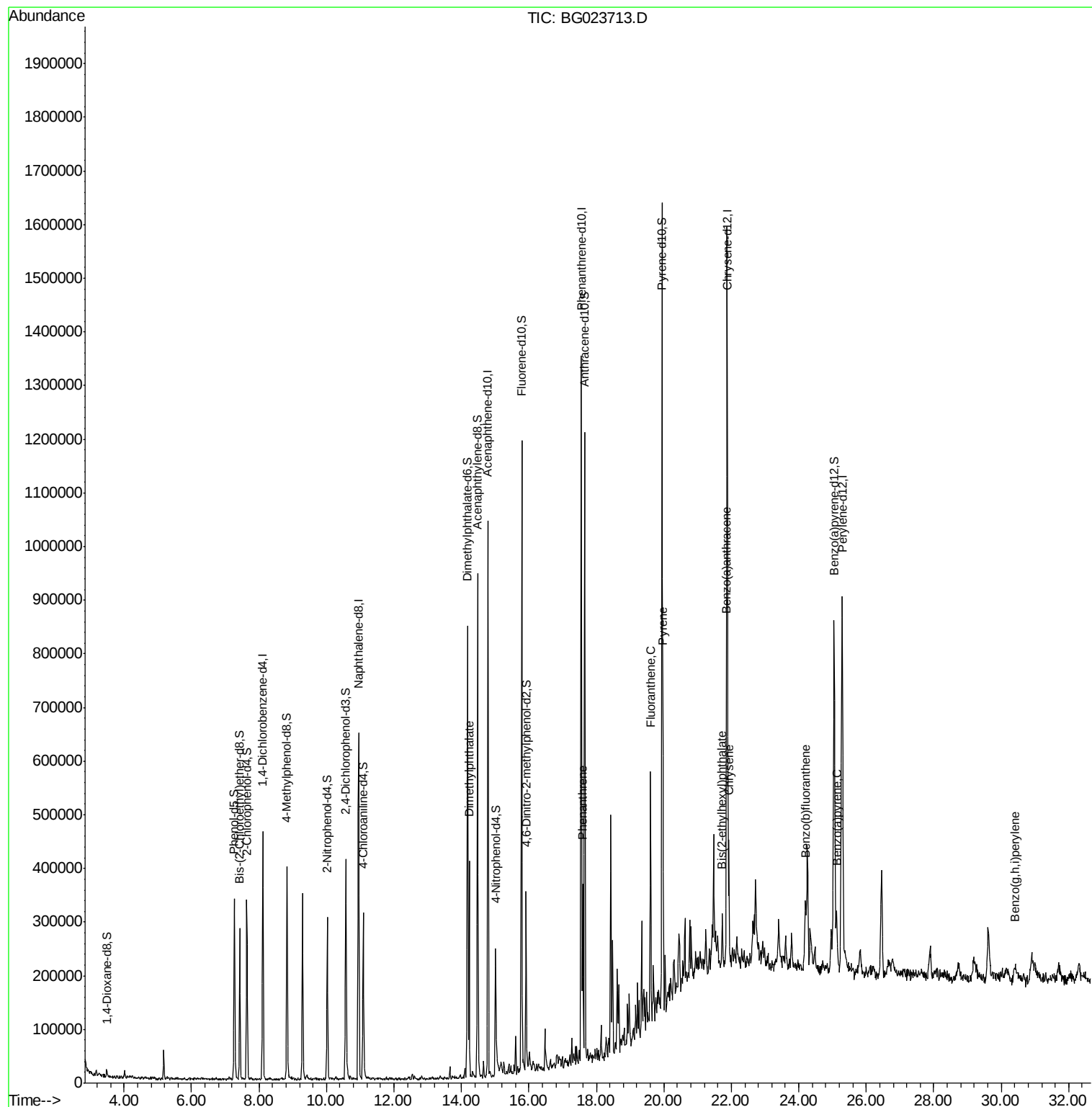
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Misc :
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Manual Integrations
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Quant Time: Aug 15 19:25:22 2016
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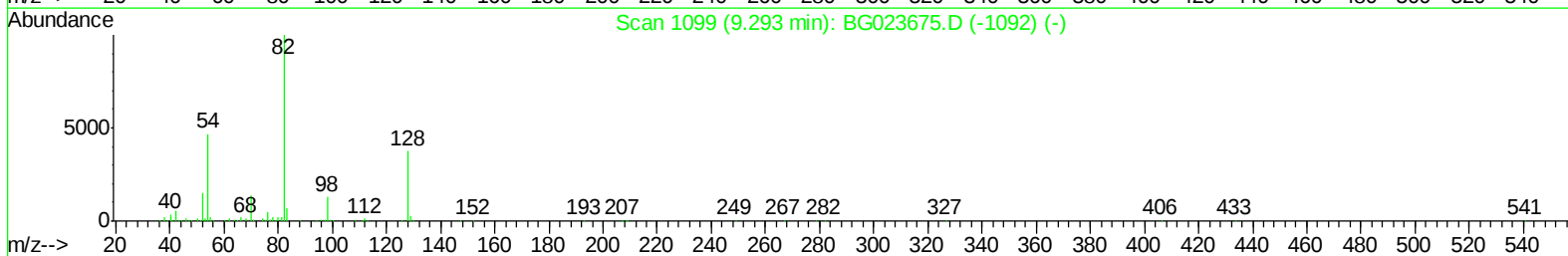
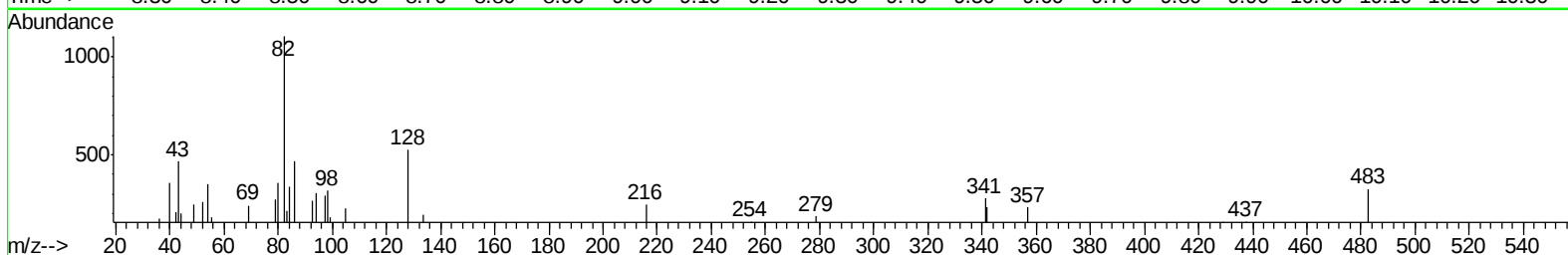
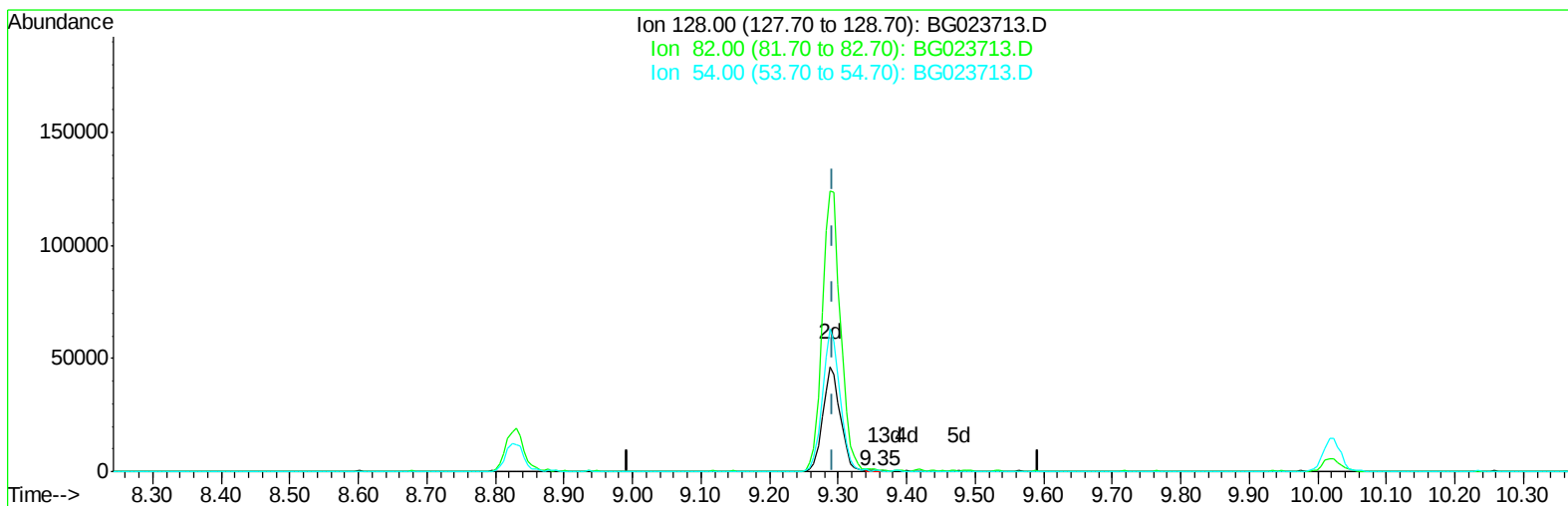
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TIC: BG023713.D

(19) Nitrobenzene-d5 (S)

9.353min (+0.060) 0.14ng/ul

response 590

Ion	Exp%	Act%
128.00	100	100
82.00	219.50	210.46
54.00	78.50	66.54
0.00	0.00	0.00

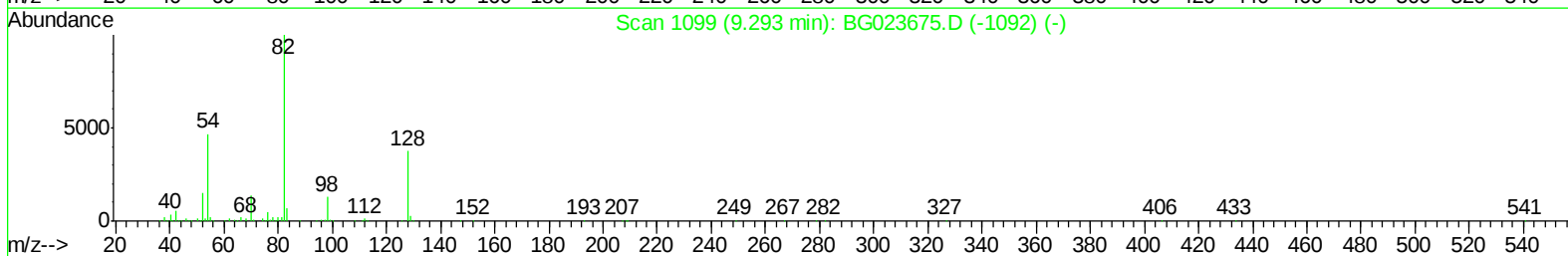
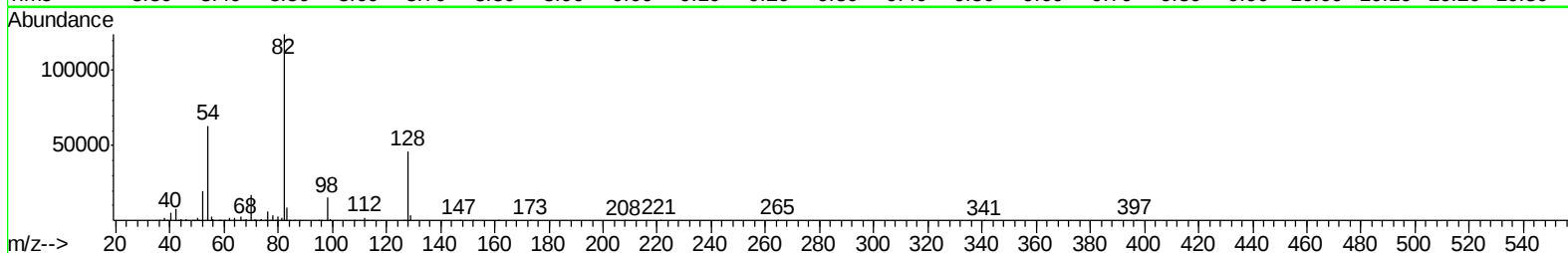
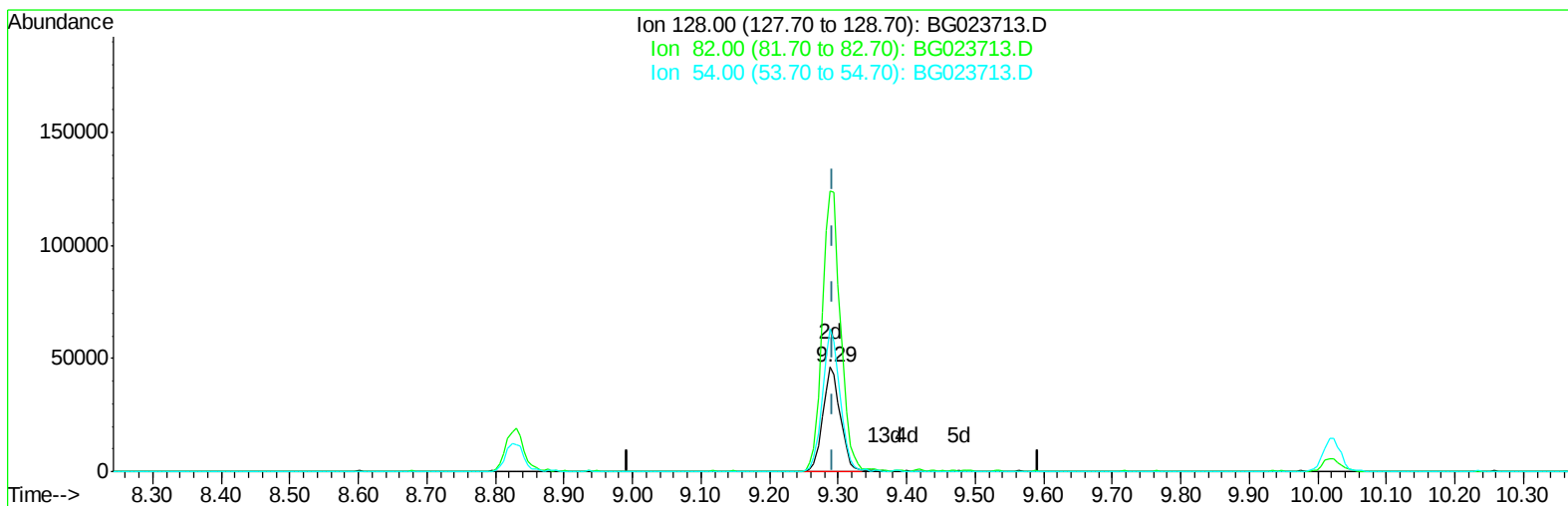
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TIC: BG023713.D

(19) Nitrobenzene-d5 (S)

9.288min (-0.004) 18.76ng/ul m

response 81783

Ion	Exp%	Act%
128.00	100	100
82.00	219.50	268.94#
54.00	78.50	136.82#
0.00	0.00	0.00

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.11	152	124674	20.00	ng/ul	0.00
18) Naphthalene-d8	10.94	136	552982	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.79	164	354065	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.55	188	758158	20.00	ng/ul	0.00
75) Chrysene-d12	21.88	240	818205	20.00	ng/ul	0.00
83) Perylene-d12	25.29	264	809010	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.48	96	7572	2.58	ng/uL	0.00
5) Phenol-d5	7.27	99	205210	16.34	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.43	67	142746	17.74	ng/ul	0.00
9) 2-Chlorophenol-d4	7.64	132	150896	17.61	ng/ul	0.00
13) 4-Methylphenol-d8	8.83	113	158026	16.12	ng/ul	0.00
19) Nitrobenzene-d5	9.29	128	81783m	18.76	ng/ul	0.00
22) 2-Nitrophenol-d4	10.02	143	96038	19.18	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.57	165	166086	17.59	ng/ul	0.00
29) 4-Chloroaniline-d4	11.09	131	170749	15.41	ng/ul	0.00
43) Dimethylphthalate-d6	14.18	166	556847	19.21	ng/ul	0.00
46) Acenaphthylene-d8	14.48	160	639695	19.61	ng/ul	0.00
51) 4-Nitrophenol-d4	15.01	143	77420	15.64	ng/ul	0.00
57) Fluorene-d10	15.79	176	483524	18.05	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.91	200	77201	15.87	ng/ul	0.00
70) Anthracene-d10	17.65	188	690570	20.22	ng/ul	0.00
76) Pyrene-d10	19.94	212	765475	18.48	ng/ul	0.00
87) Benzo(a)pyrene-d12	25.05	264	726204	19.83	ng/ul	0.00

Target Compounds

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44) Dimethylphthalate	14.23	163	249041	8.68	ng/ul	97
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74) Fluoranthene	19.61	202	277197m	6.90	ng/ul	
77) Pyrene	19.97	202	348355	6.89	ng/ul#	89
80) Benzo(a)anthracene	21.85	228	129502	2.81	ng/ul	98
81) Bis(2-ethylhexyl)phthalate	21.73	149	31966	1.24	ng/ul#	98
82) Chrysene	21.92	228	164124	3.92	ng/ul	97
85) Benzo(b)fluoranthene	24.20	252	151087	3.28	ng/ul#	93
88) Benzo(a)pyrene	25.12	252	119637m	2.69	ng/ul	
91) Benzo(g,h,i)perylene	30.40	276	69950m	1.62	ng/ul	

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

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