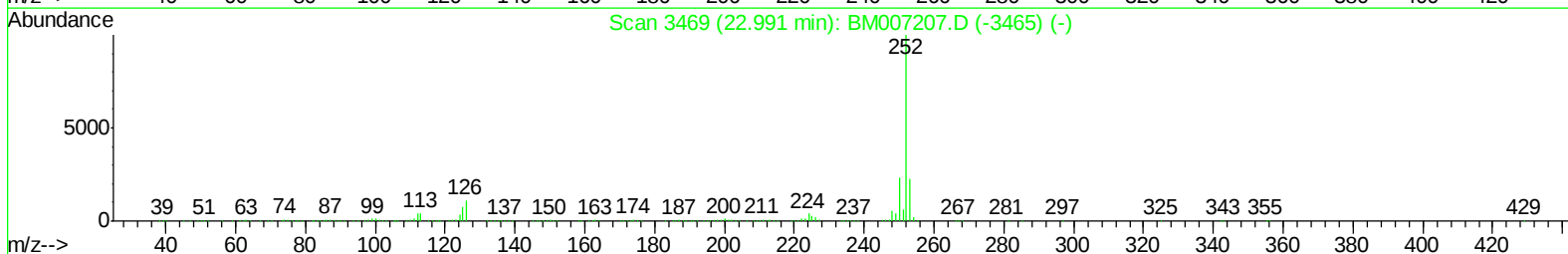
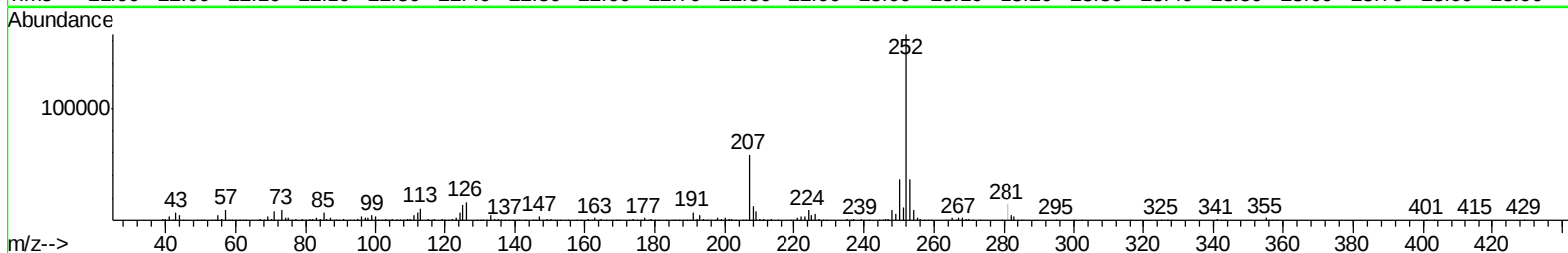
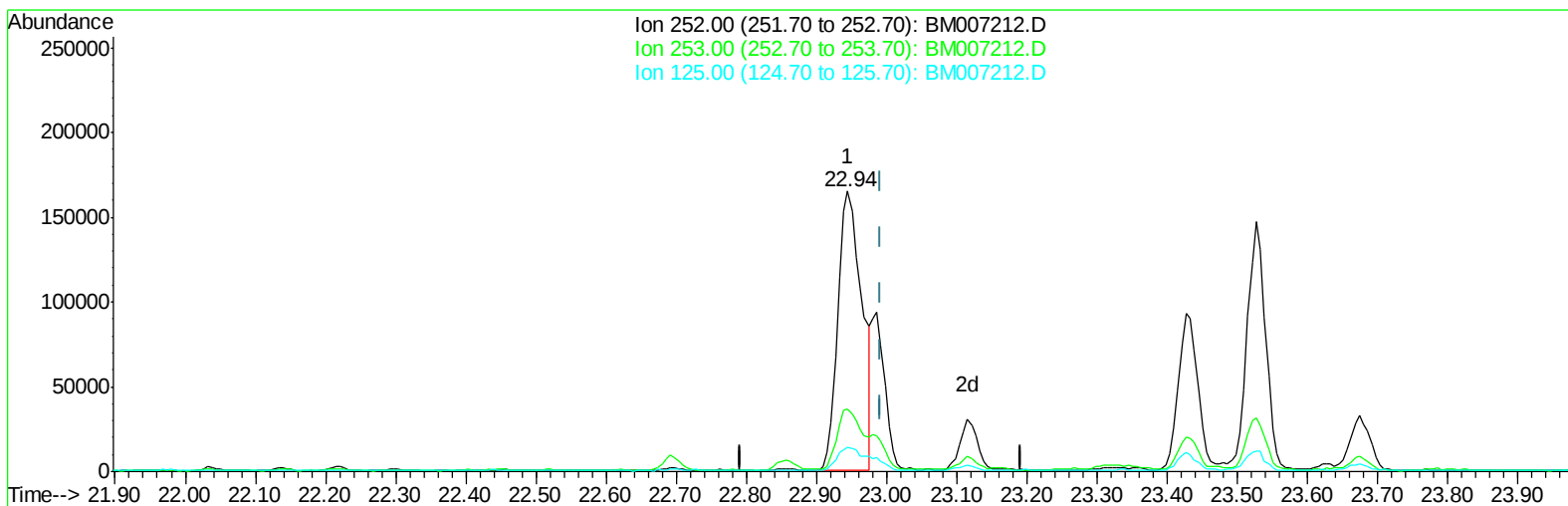


Data Path : Z:\HPCHEM1\BNA_M\DATA\BM081916\
Data File : BM007212.D
Acq On : 18 Aug 2016 20:03
Operator : UM/SJ
Sample : H4445-06
Misc :
ALS Vial : 13 Sample Multiplier: 1

Quant Time: Aug 19 05:14:42 2016
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM081116.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Aug 19 04:48:28 2016
Response via : Initial Calibration



TIC: BM007212.D

(86) Benzo(k)fluoranthene

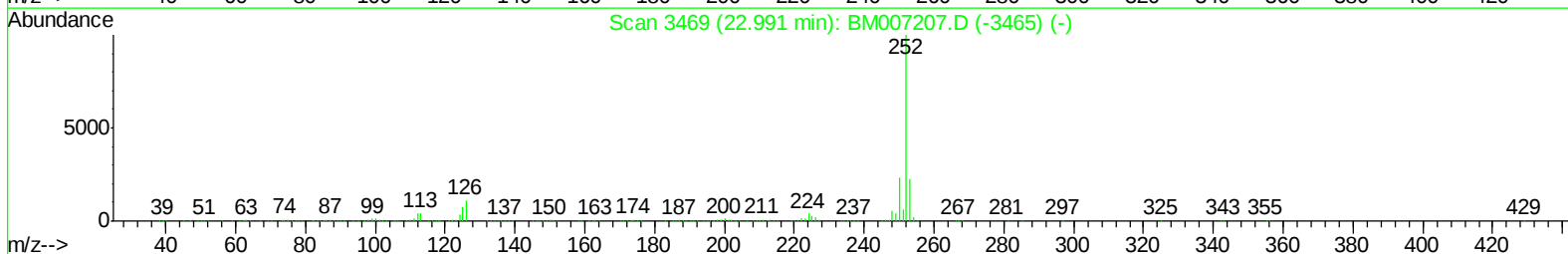
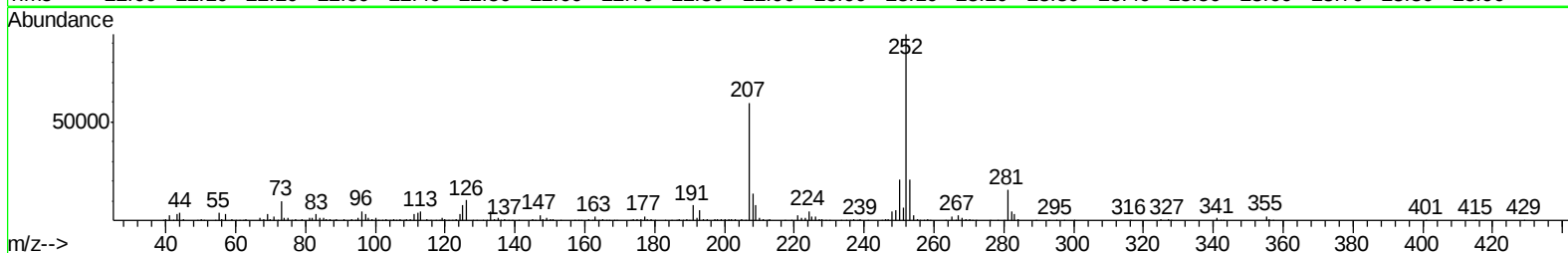
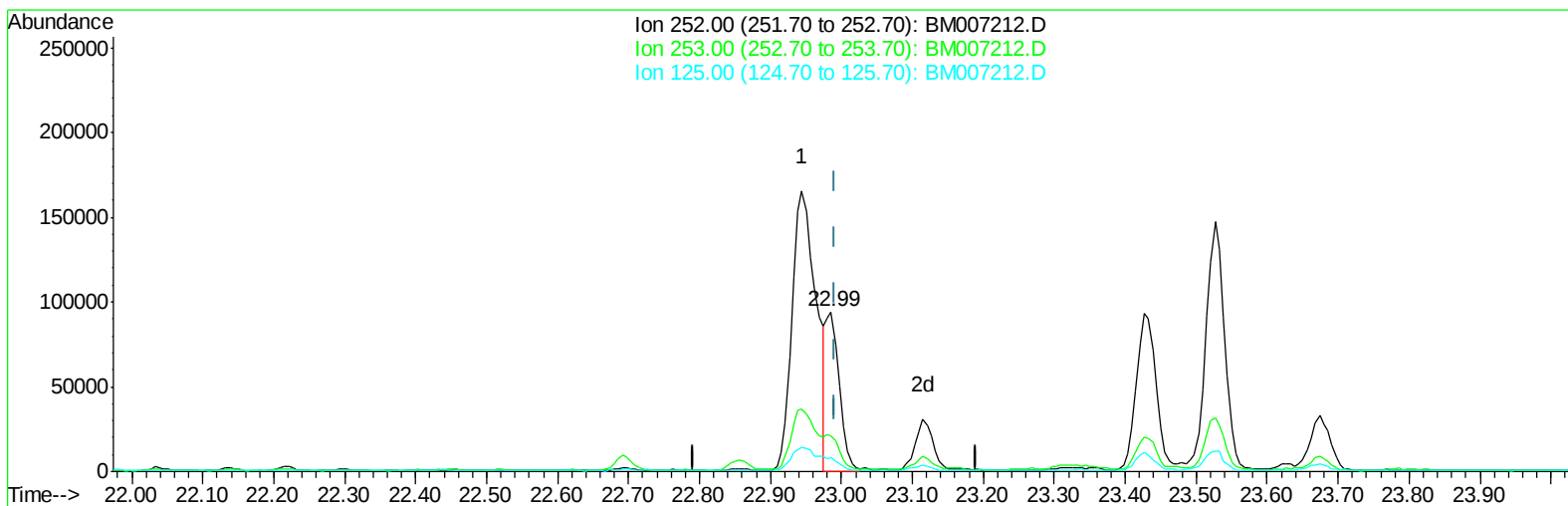
22.944min (-0.047) 3.65ng/ul

response 388010

Ion	Exp%	Act%
252.00	100	100
253.00	22.00	22.38
125.00	7.10	8.58#
0.00	0.00	0.00

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TIC: BM007212.D

(86) Benzo(k)fluoranthene

22.986min (-0.006) 1.18ng/ul m

response 125542

Ion	Exp%	Act%
252.00	100	100
253.00	22.00	22.50
125.00	7.10	8.81#
0.00	0.00	0.00

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM081916\
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 Operator : UM/SJ
 Sample : H4445-06
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Quant Time: Aug 19 05:16:10 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM081116.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Aug 19 04:48:28 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.80	152	242685	20.00	ng/ul	0.00
18) Naphthalene-d8	10.59	136	981630	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.43	164	586100	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.17	188	1403906	20.00	ng/ul	0.00
75) Chrysene-d12	21.36	240	1816727	20.00	ng/ul	0.00
83) Perylene-d12	23.63	264	1880240	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.32	96	20018	3.60	ng/uL	0.00
5) Phenol-d5	6.97	99	362954	18.46	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.14	67	231400	20.53	ng/ul	0.00
9) 2-Chlorophenol-d4	7.33	132	310135	19.81	ng/ul	0.00
13) 4-Methylphenol-d8	8.50	113	247494	15.30	ng/ul	0.00
19) Nitrobenzene-d5	8.96	128	151479	21.25	ng/ul	0.00
22) 2-Nitrophenol-d4	9.67	143	164449	21.59	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.20	165	313681	19.66	ng/ul	0.00
29) 4-Chloroaniline-d4	10.72	131	378172	19.92	ng/ul	0.00
43) Dimethylphthalate-d6	13.84	166	1018266	21.03	ng/ul	0.00
46) Acenaphthylene-d8	14.12	160	1246557	21.43	ng/ul	0.00
51) 4-Nitrophenol-d4	14.63	143	131896	15.20	ng/ul	0.00
57) Fluorene-d10	15.43	176	909474	21.48	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.54	200	107076	13.80	ng/ul	0.00
70) Anthracene-d10	17.27	188	1452364	22.33	ng/ul	0.00
76) Pyrene-d10	19.57	212	1799452	23.47	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.48	264	2000197	23.26	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
44) Dimethylphthalate	13.89	163	317361	6.56	ng/ul	98
69) Phenanthrene	17.22	178	294570	3.90	ng/ul	99
74) Fluoranthene	19.23	202	510660	5.58	ng/ul	99
77) Pyrene	19.60	202	453522	4.52	ng/ul	98
80) Benzo(a)anthracene	21.34	228	323798	3.02	ng/ul	98
82) Chrysene	21.39	228	290929	2.97	ng/ul	97
85) Benzo(b)fluoranthene	22.94	252	388010	3.43	ng/ul	98
86) Benzo(k)fluoranthene	22.99	252	125542m	1.18	ng/ul	
88) Benzo(a)pyrene	23.53	252	265628	2.47	ng/ul	99
89) Indeno(1,2,3-cd)pyrene	25.91	276	167126	1.27	ng/ul	97
91) Benzo(g,h,i)perylene	26.61	276	139110	1.25	ng/ul	96

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

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