

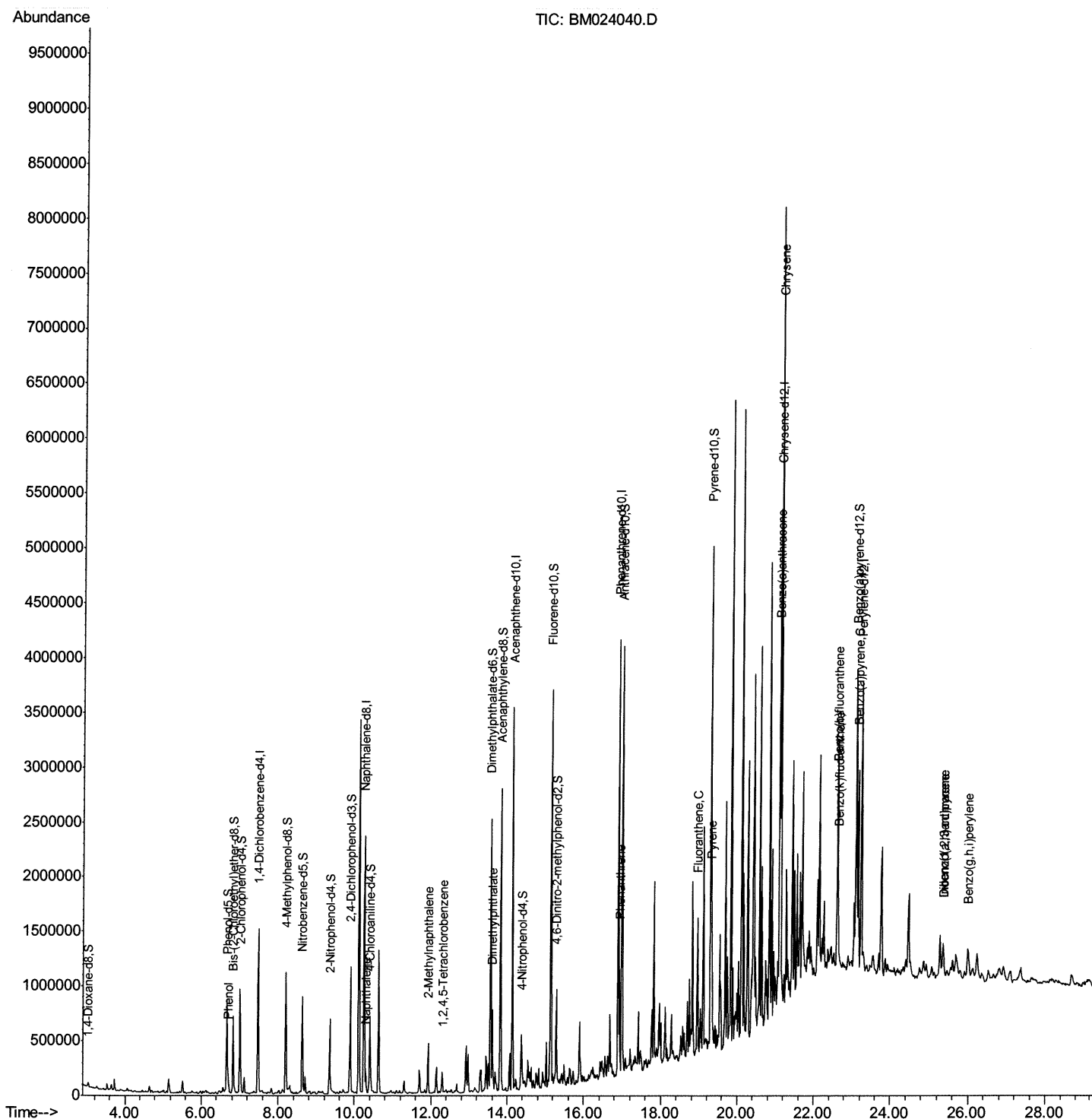
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121119\
Data File : BM024040.D
Acq On : 11 Dec 2019 19:04
Operator : JU
Sample : K6088-02
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0BM8

Manual Integrations
APPROVED

mohammad
12/12/2019 11:25:52 AM

Quant Time: Dec 12 10:48:40 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM112719MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat Dec 07 04:26:57 2019
Response via : Initial Calibration



Quantitation Report (Qedit)

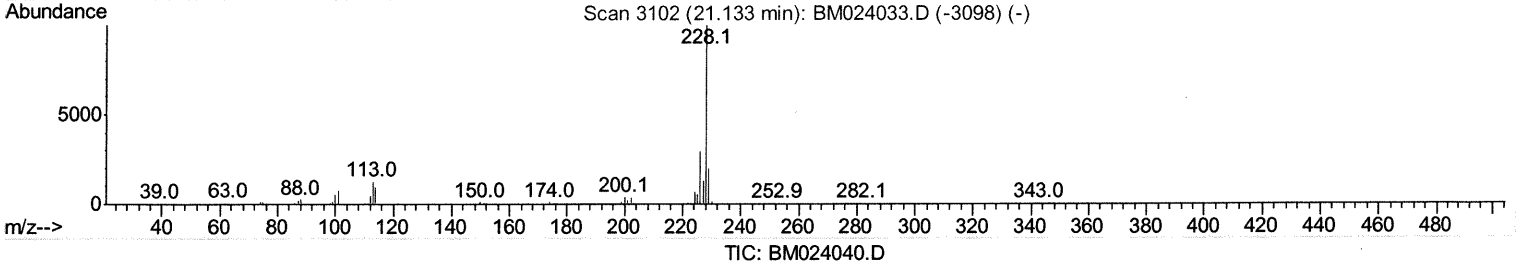
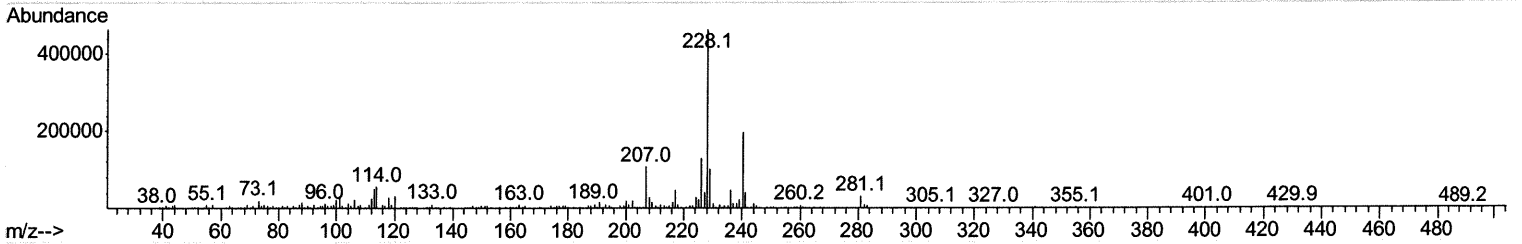
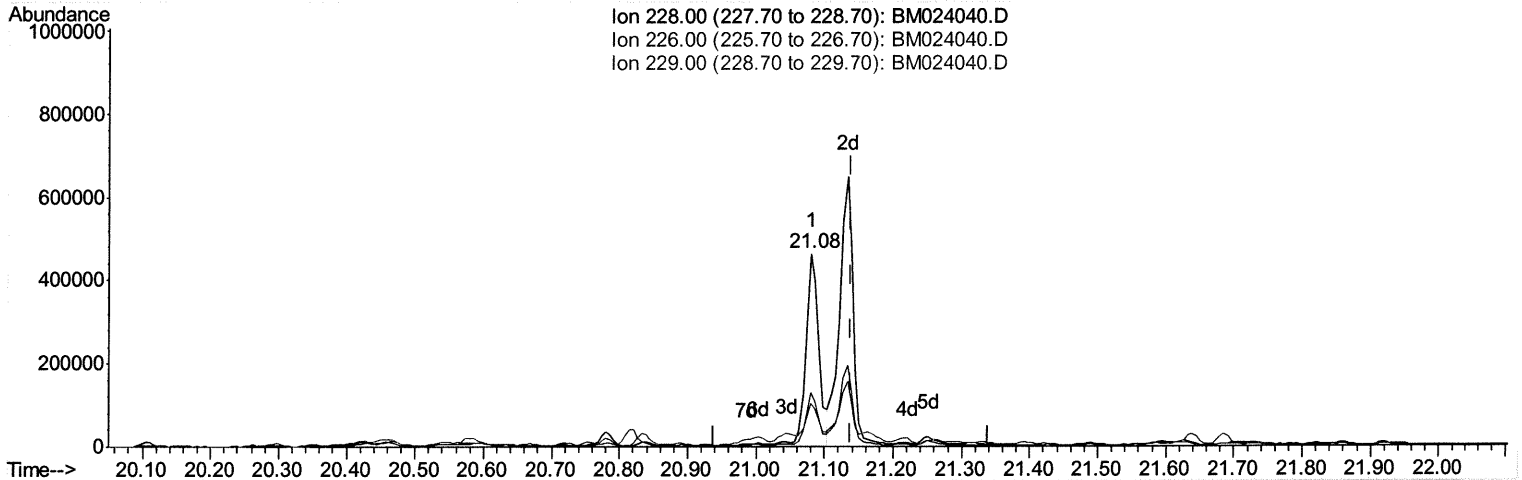
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(85) Chrysene

21.080min (-0.059) 5.52ng/ul

response 602690

Ion	Exp%	Act%
228.00	100	100
226.00	29.30	28.23
229.00	19.10	22.09
0.00	0.00	0.00

Quantitation Report (Qedit)

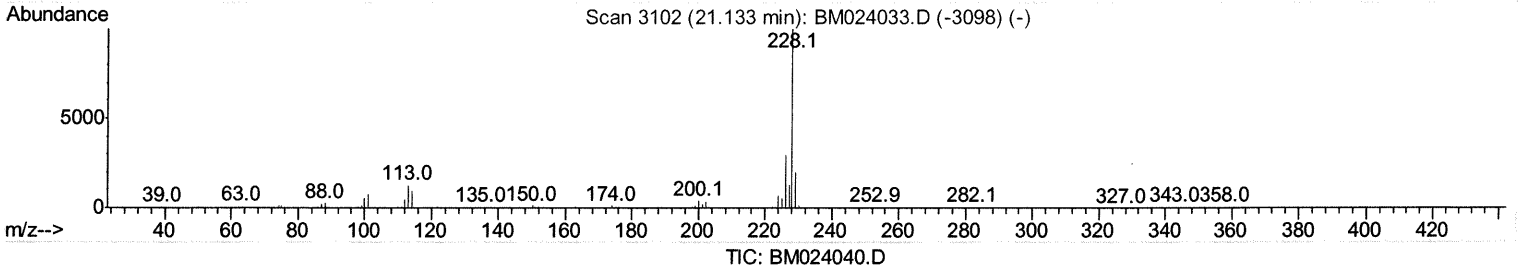
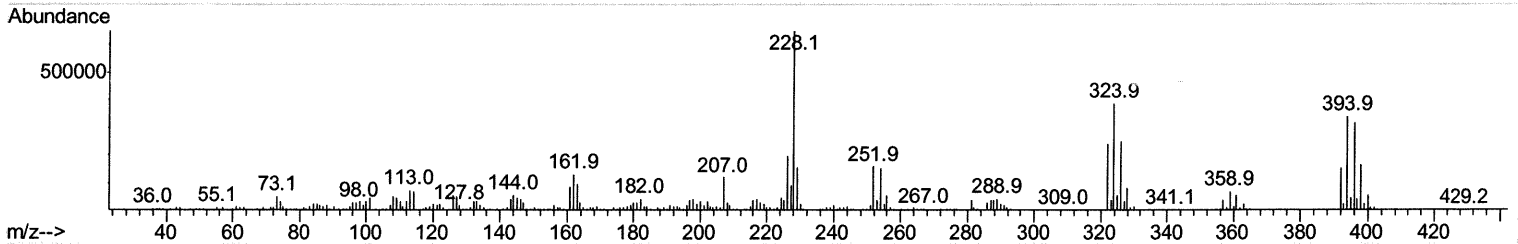
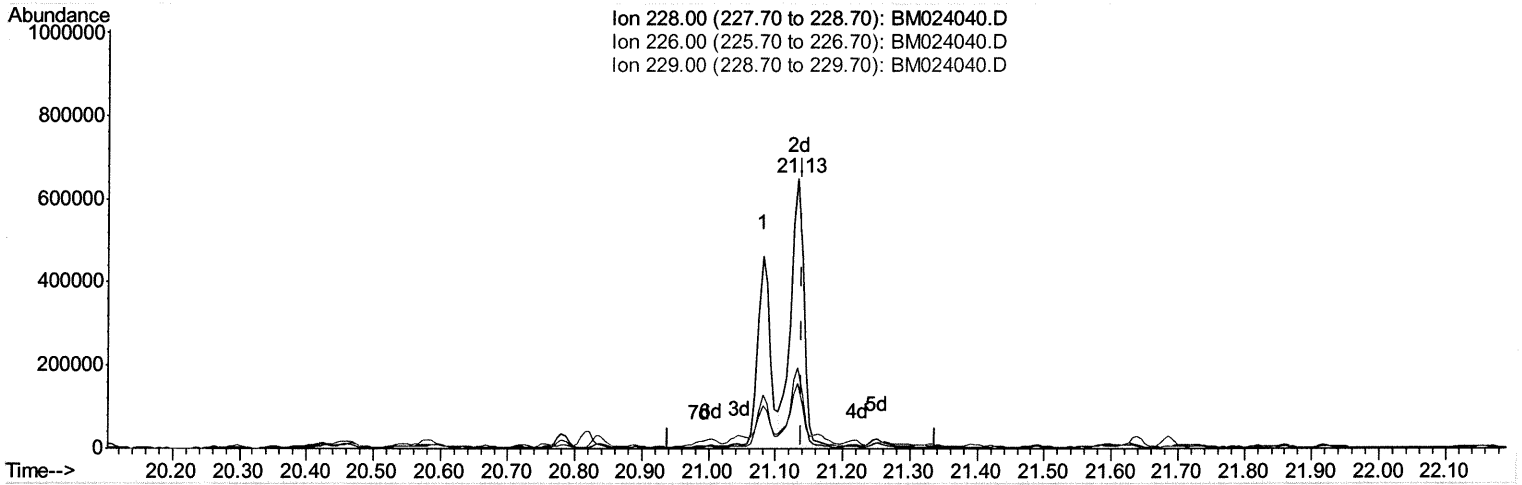
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(85) Chrysene

21.133min (-0.006) 8.00ng/ul m JU 12/13/19

response 873340

Ion	Exp%	Act%
228.00	100	100
226.00	29.30	29.97
229.00	19.10	23.95#
0.00	0.00	0.00

Quantitation Report (Qedit)

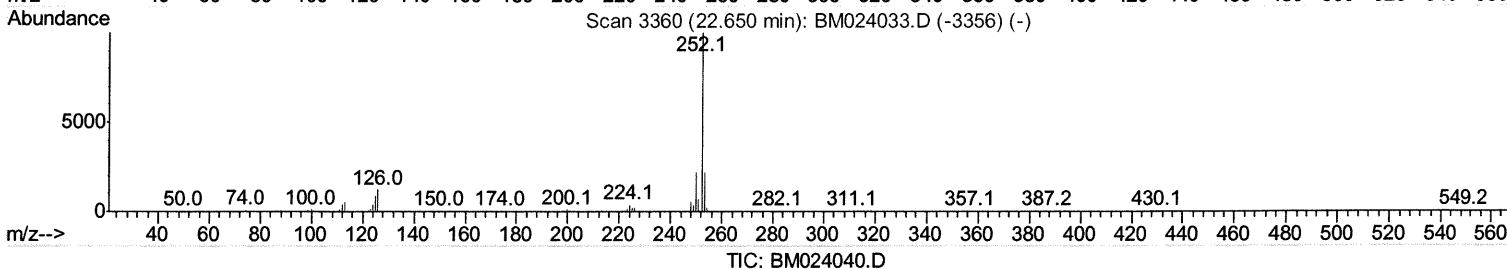
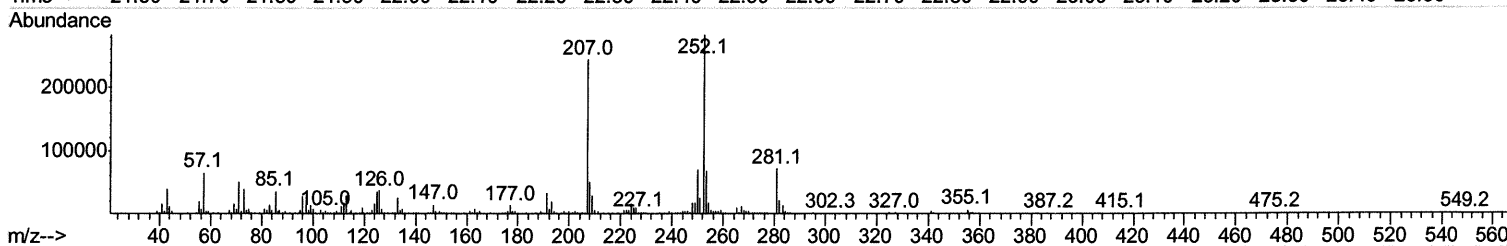
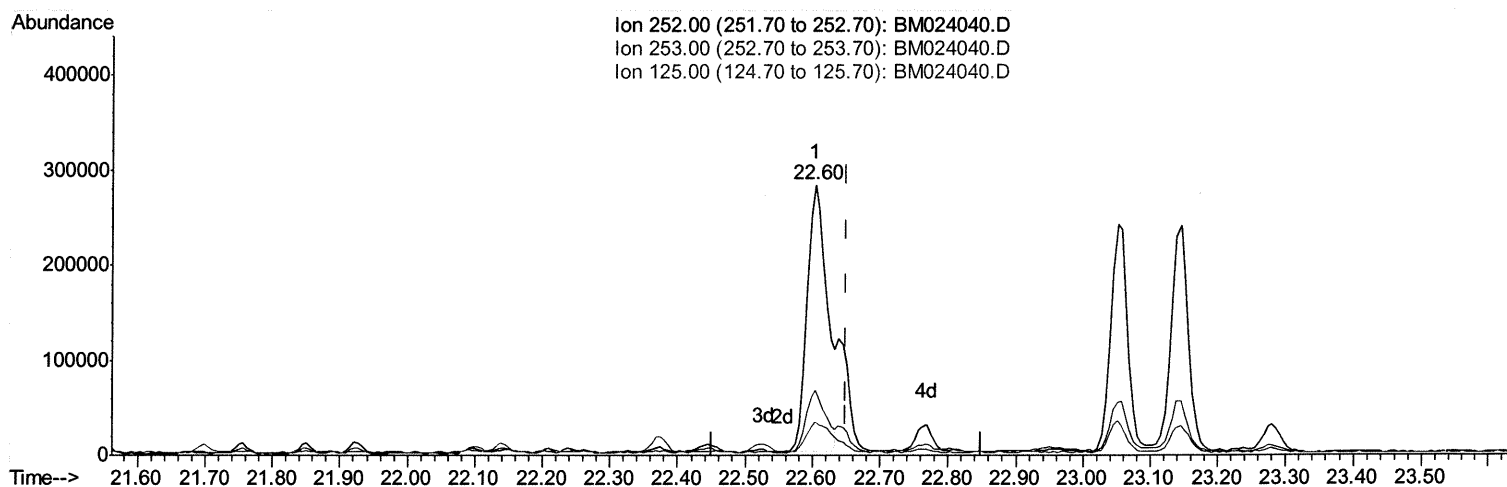
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Instrument :
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Manual Integrations
APPROVED

mohammad
12/12/2019 11:25:52 AM

Quant Time: Dec 12 03:38:36 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM112719MA.M
Quant Title : SVOA CALIBRATION
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(89) Benzo(k)fluoranthene

22.603min (-0.047) 5.32ng/ul

response 580213

Ion	Exp%	Act%
252.00	100	100
253.00	21.50	24.22
125.00	8.70	12.23#
0.00	0.00	0.00

Quantitation Report (Qedit)

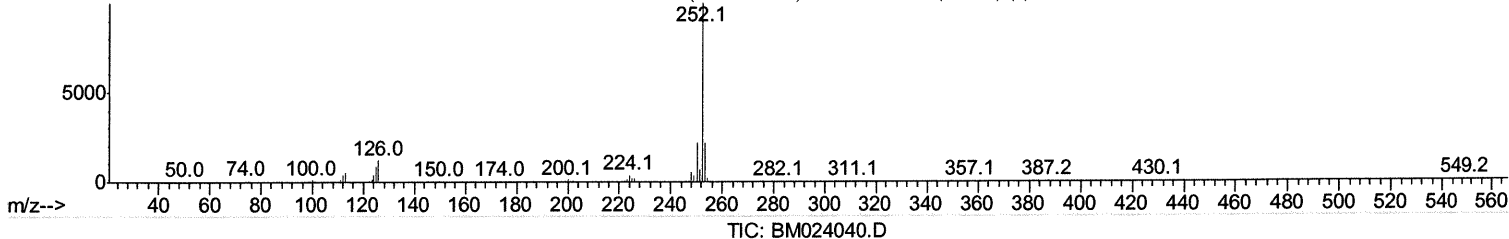
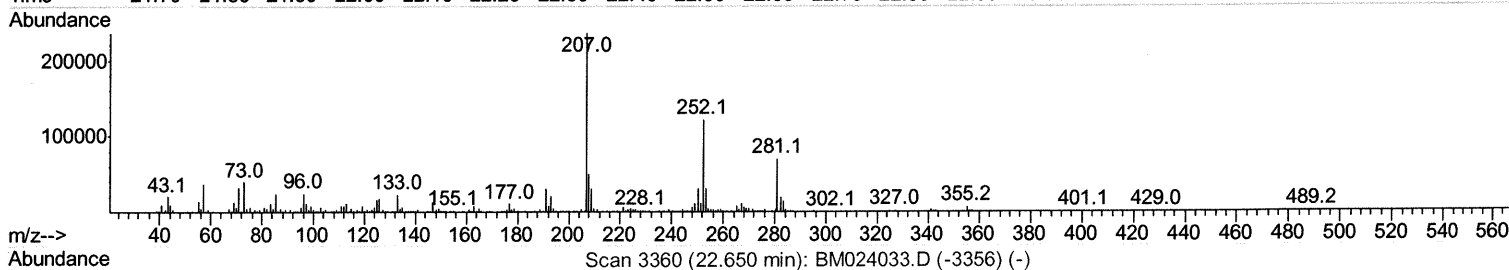
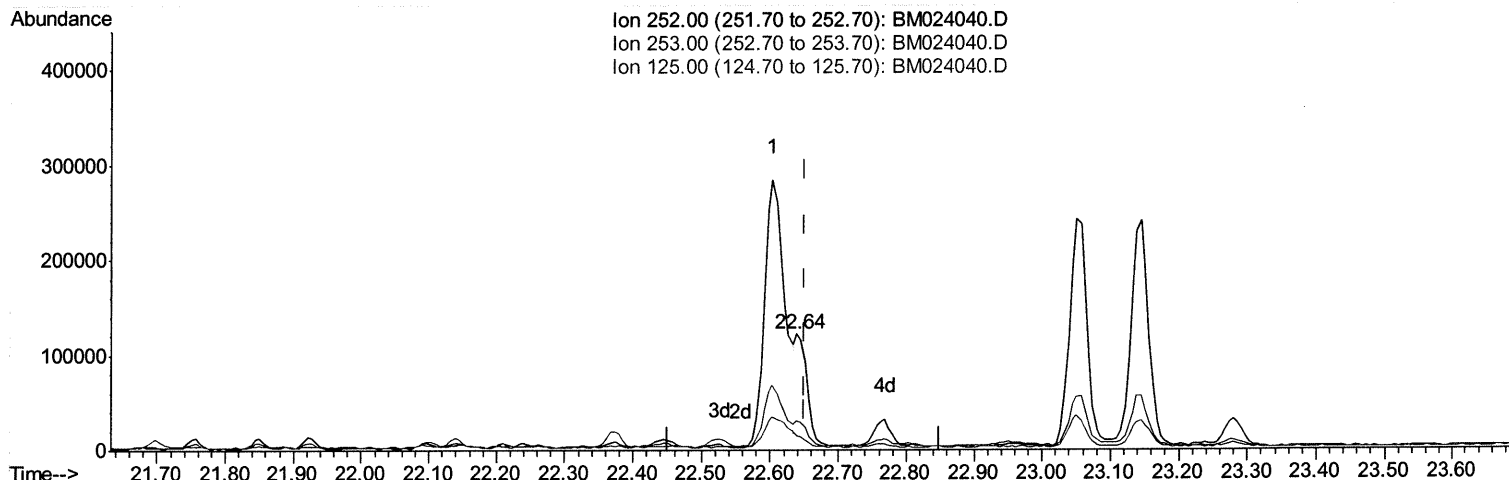
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Manual Integrations
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(89) Benzo(k)fluoranthene

22.639min (-0.012) 1.32ng/ul m

JU 12/13/19

response 143985

Ion	Exp%	Act%
252.00	100	100
253.00	21.50	25.03
125.00	8.70	13.10#
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	7.47	152	402583	20.00	ng/ul	0.00
18) Naphthalene-d8	10.25	136	1724802	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.13	164	1099557	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.89	188	2283782	20.00	ng/ul	0.00
78) Chrysene-d12	21.10	240	1739739	20.00	ng/ul	0.00
86) Perylene-d12	23.23	264	1858747	20.00	ng/ul	-0.01

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.03	96	24943	2.76	ng/uL	0.00
5) Phenol-d5	6.67	99	531215	15.06	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.83	67	346543	16.89	ng/ul	0.00
9) 2-Chlorophenol-d4	7.01	132	419549	15.38	ng/ul	0.00
13) 4-Methylphenol-d8	8.19	113	429114	15.41	ng/ul	0.00
19) Nitrobenzene-d5	8.63	128	207389	15.49	ng/ul	0.00
22) 2-Nitrophenol-d4	9.35	143	223180	14.67	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.88	165	429576	14.91	ng/ul	0.00
29) 4-Chloroaniline-d4	10.40	131	420166	13.29	ng/ul	0.00
44) Dimethylphthalate-d6	13.56	166	1546905	18.11	ng/ul	0.00
47) Acenaphthylene-d8	13.82	160	1756730	17.54	ng/ul	0.00
52) 4-Nitrophenol-d4	14.37	143	141904	9.88	ng/ul	0.00
58) Fluorene-d10	15.13	176	1375837	19.14	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.28	200	181072	12.63	ng/ul	0.00
71) Anthracene-d10	16.99	188	2087536	19.14	ng/ul	0.00
79) Pyrene-d10	19.29	212	2123746	23.62	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.10	264	1901895	19.36	ng/ul	-0.01

Target Compounds

					Qvalue	
6) Phenol	6.69	94	59824	1.648	ng/ul	94
28) Naphthalene	10.29	128	162879	1.791	ng/ul	98
34) 2-Methylnaphthalene	11.92	142	220236	3.336	ng/ul	97
37) 1,2,4,5-Tetrachlorobenzene	12.30	216	57989	1.691	ng/ul	98
45) Dimethylphthalate	13.60	163	456236	5.455	ng/ul	100
70) Phenanthrene	16.93	178	574602	4.454	ng/ul	100
77) Fluoranthene	18.96	202	608295	4.397	ng/ul	99
80) Pyrene	19.32	202	738144	6.297	ng/ul	100
83) Benzo(a)anthracene	21.08	228	586307	5.042	ng/ul	96
85) Chrysene	21.13	228	873340m	7.998	ng/ul	> 94 12/13/19
88) Benzo(b)fluoranthene	22.60	252	580213	5.048	ng/ul#	94
89) Benzo(k)fluoranthene	22.64	252	143985m	1.320	ng/ul	> 94 12/13/19
91) Benzo(a)pyrene	23.14	252	422543	3.888	ng/ul#	94
92) Indeno(1,2,3-cd)pyrene	25.35	276	221743	1.643	ng/ul	96
93) Dibenzo(a,h)anthracene	25.36	278	169888	1.499	ng/ul#	77
94) Benzo(g,h,i)perylene	25.99	276	279276	2.471	ng/ul	99

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