

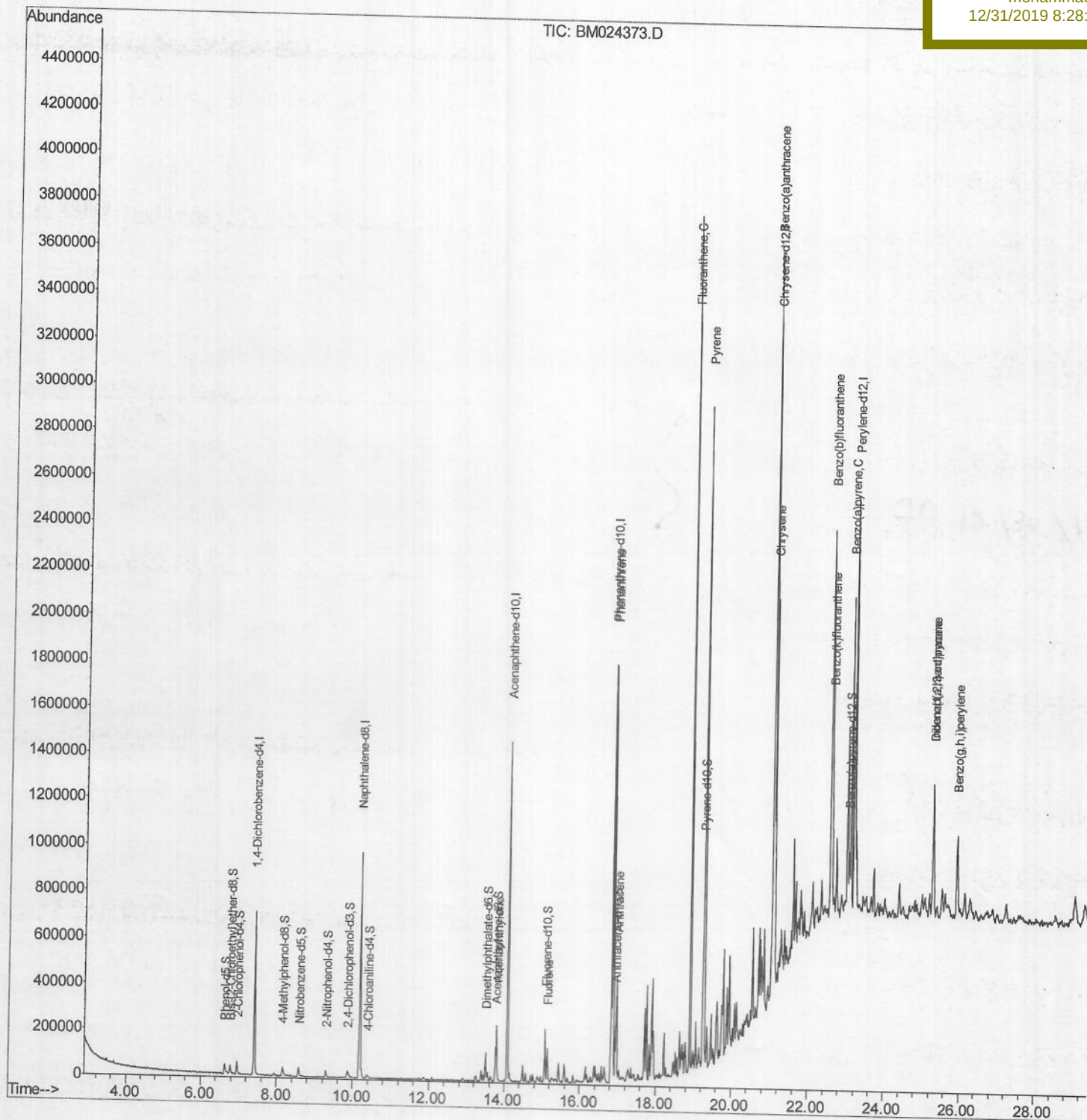
Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM123019\
Data File : BM024373.D
Acq On : 30 Dec 2019 15:10
Operator : JU
Sample : K6322-02DL 10X
Misc :
ALS Vial : 9 Sample Multiplier: 1

Quant Time: Dec 30 15:51:34 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM121719MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Dec 30 14:11:26 2019
Response via : Initial Calibration

Instrument :
BNA_M
ClientSampleId :
C0C30DL

Manual Integrations
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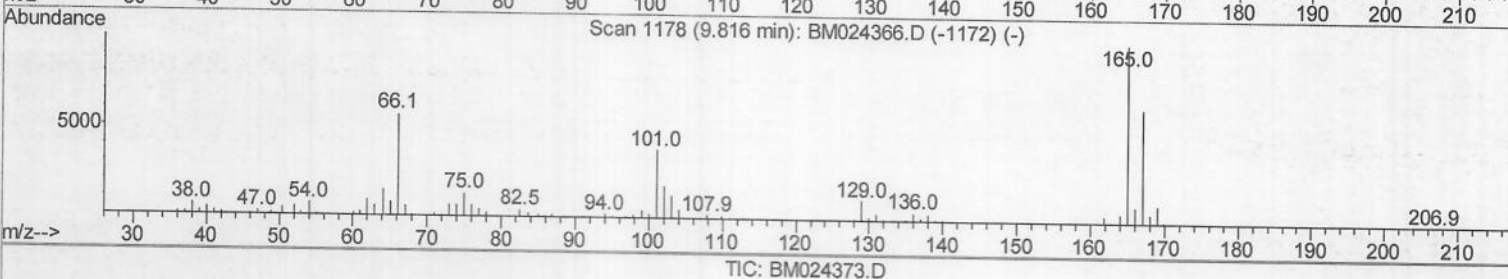
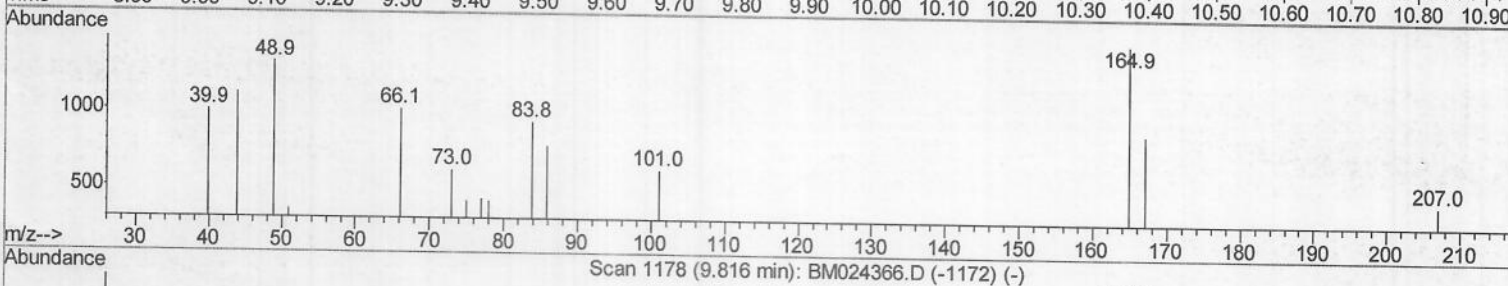
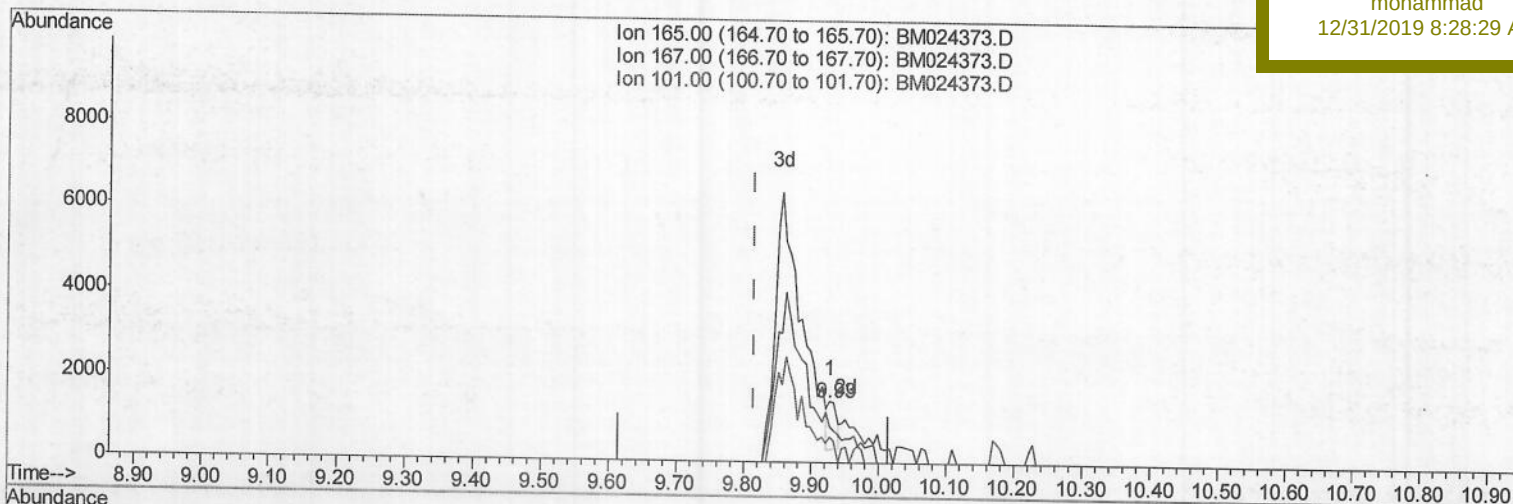
Quant Time: Dec 30 15:43:40 2019

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM121719MA.M

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(26) 2,4-Dichlorophenol-d3 (S)

9.928min (+0.112) 0.09ng/ul

response 1039

Ion	Exp%	Act%
165.00	100	100
167.00	64.20	59.69
101.00	37.00	42.27
0.00	0.00	0.00

Quantitation Report (Qedit)

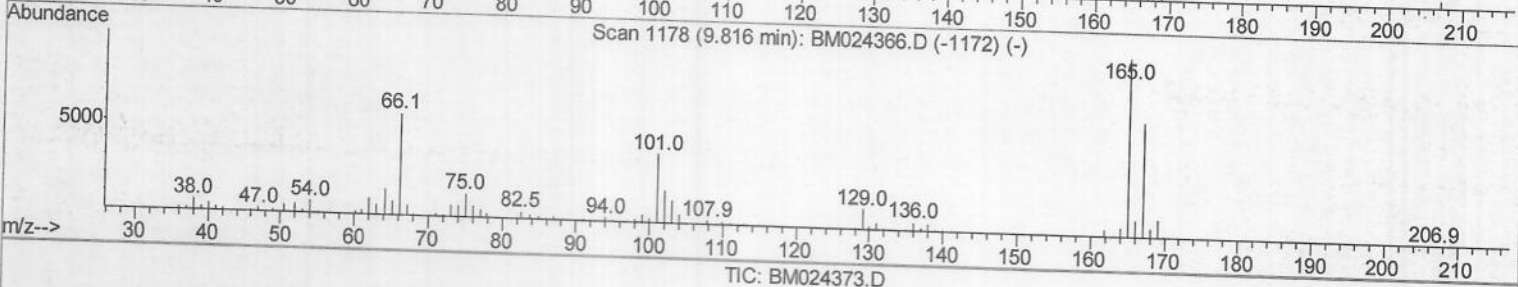
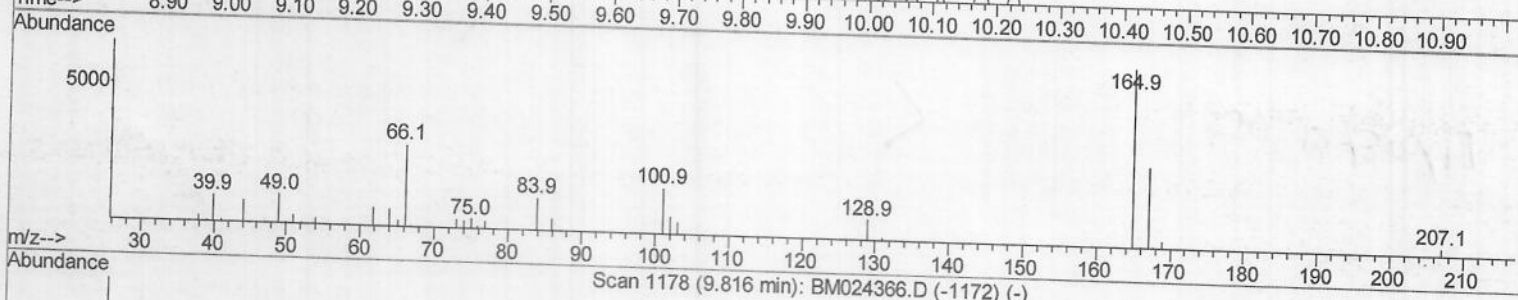
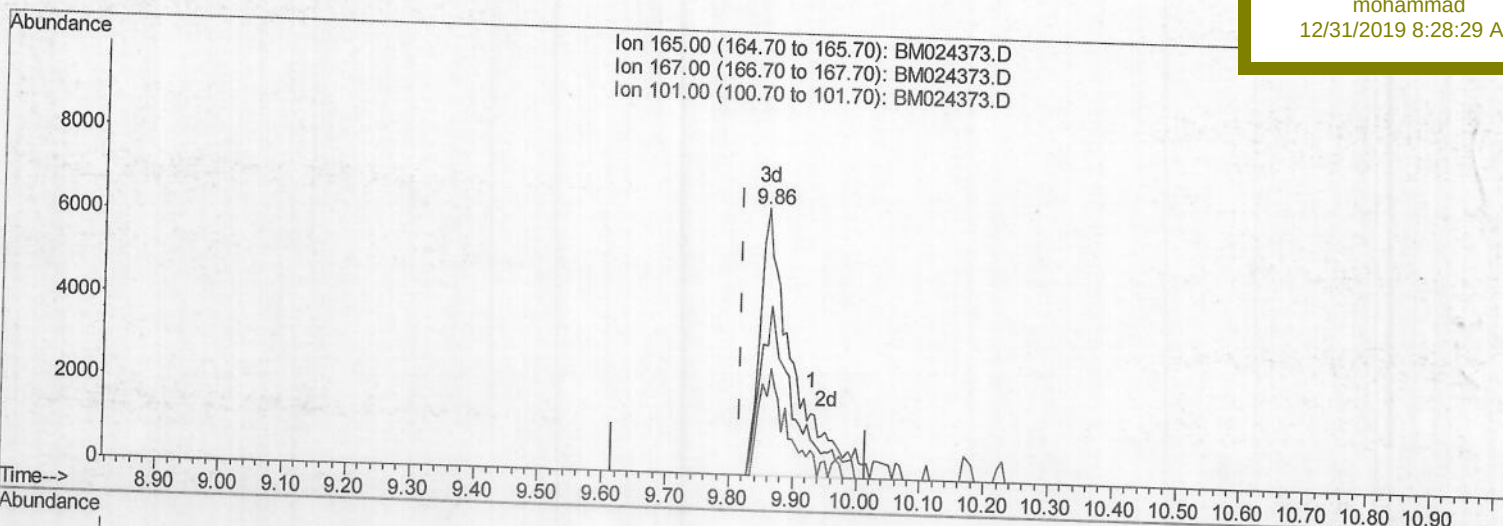
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(26) 2,4-Dichlorophenol-d3 (S)

9.857min (+0.041) 1.86ng/ul m

response 22685

Ion	Exp%	Act%
165.00	100	100
167.00	64.20	48.04#
101.00	37.00	29.09#
0.00	0.00	0.00

JU 12/30/19

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM123019\

Data File : BM024373.D

Acq On : 30 Dec 2019 15:10

Operator : JU

Sample : K6322-02DL 10X

Misc :

ALS Vial : 9 Sample Multiplier: 1

Quant Time: Dec 30 15:43:40 2019

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM121719MA.M

Quant Title : SVOA CALIBRATION

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Instrument :

BNA_M

Client Sampled :

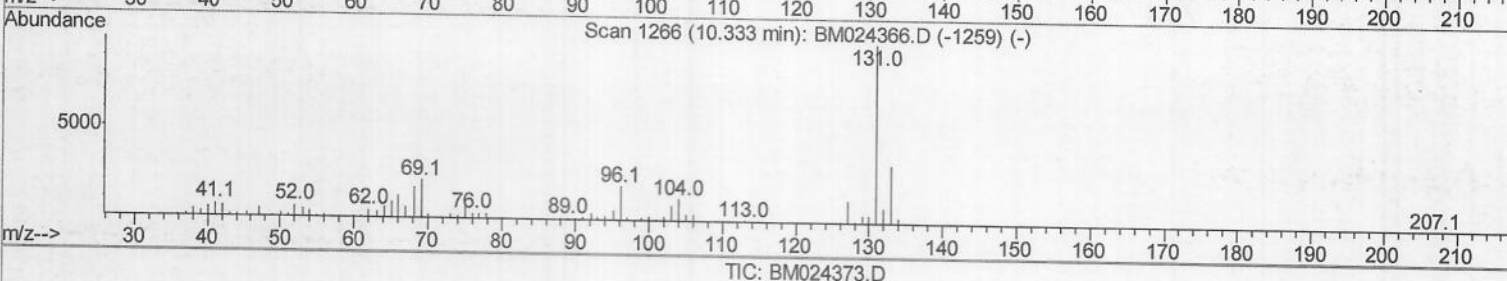
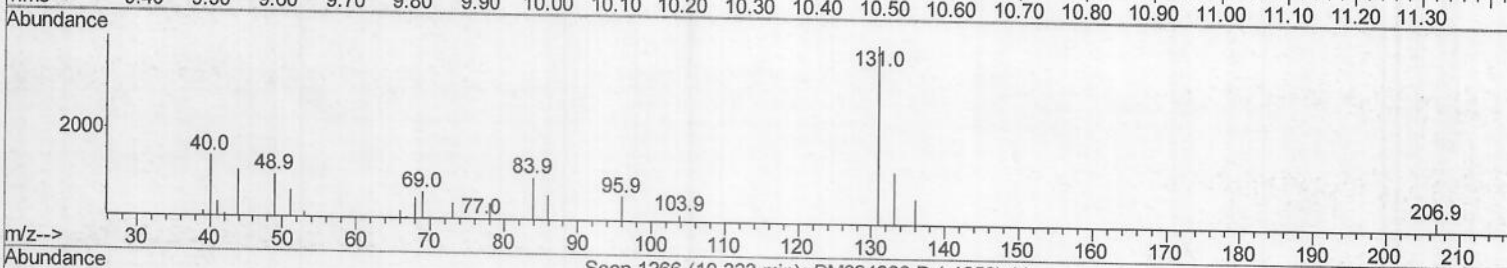
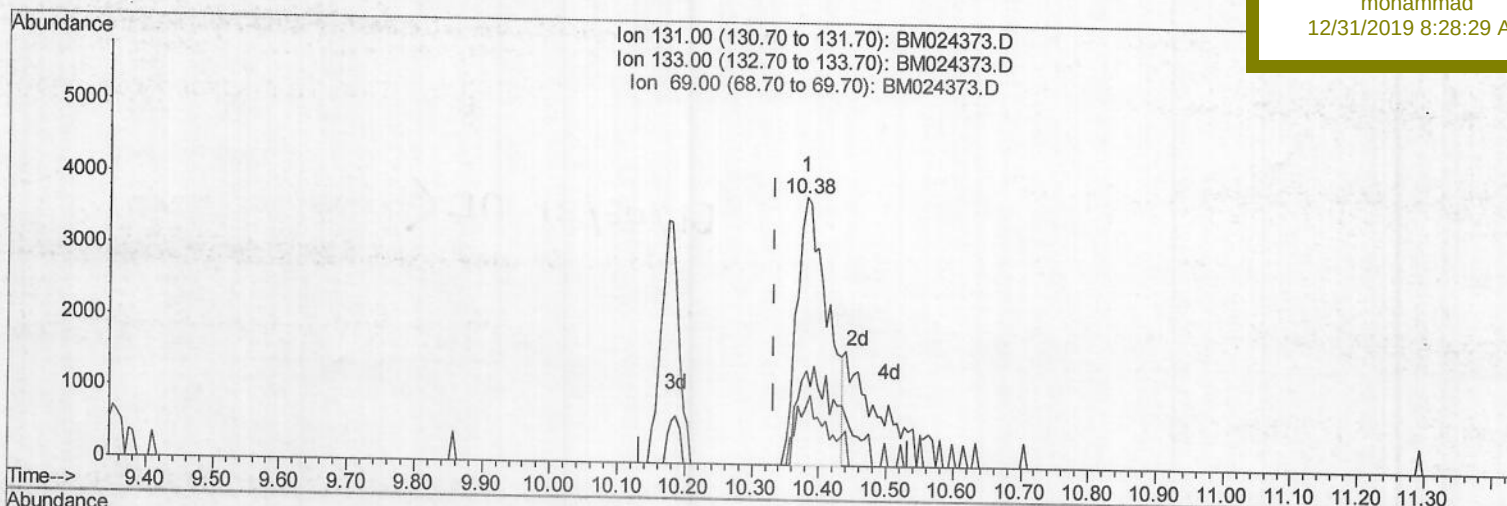
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(29) 4-Chloroaniline-d4 (S)

10.381min (+0.047) 0.80ng/ul

response 11915

Ion	Exp%	Act%
131.00	100	100
133.00	32.50	35.23
69.00	19.00	21.83
0.00	0.00	0.00

Quantitation Report (Qedit)

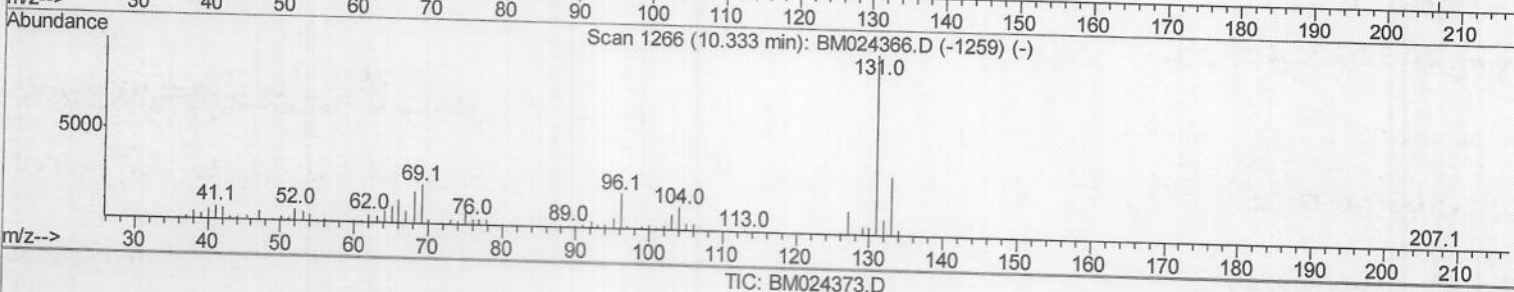
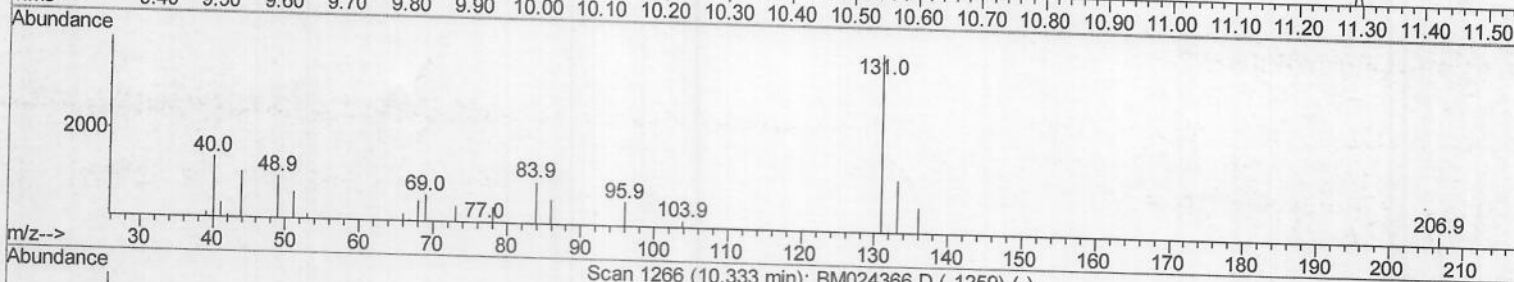
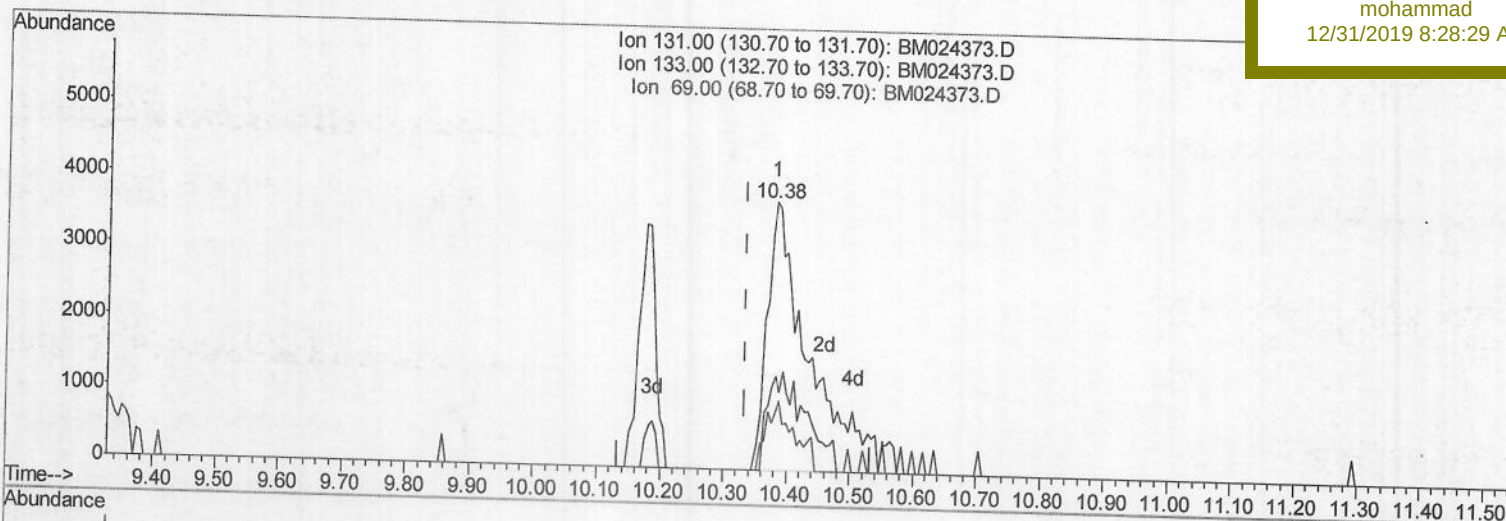
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(29) 4-Chloroaniline-d4 (S)

10.381min (+0.047) 1.16ng/ul m

response 17116

JU 12/30/19

Ion	Exp%	Act%
131.00	100	100
133.00	32.50	35.23
69.00	19.00	21.83
0.00	0.00	0.00

Quantitation Report (Qedit)

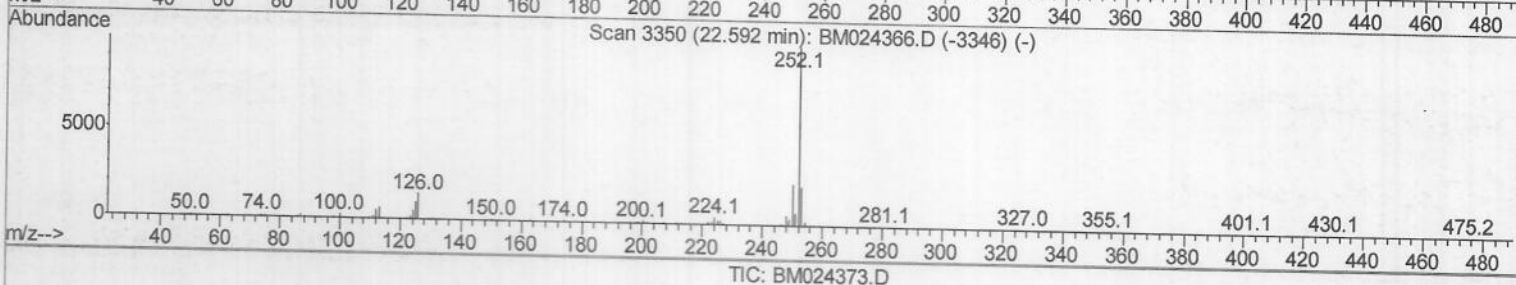
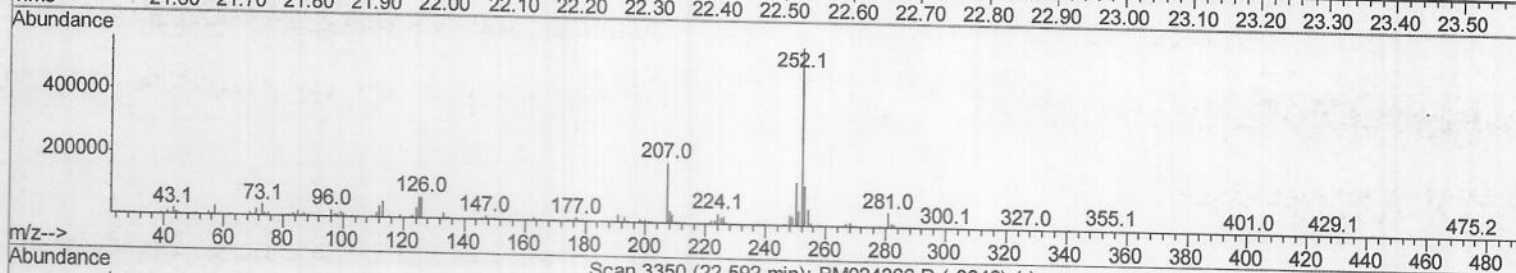
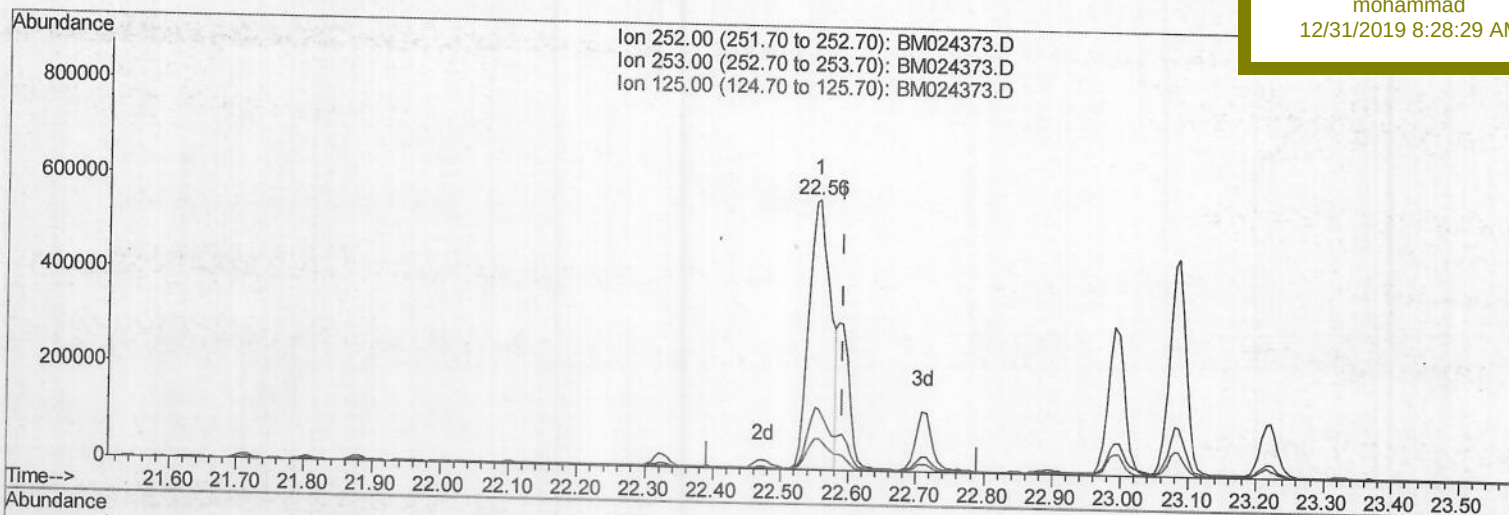
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 Data File : BM024373.D
 Acq On : 30 Dec 2019 15:10
 Operator : JU
 Sample : K6322-02DL 10X
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Quant Time: Dec 30 15:43:40 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM121719MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Dec 30 14:11:26 2019
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Manual Integrations
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TIC: BM024373.D

(89) Benzo(k)fluoranthene

22.556min (-0.035) 18.33ng/ul

response 1241603

Ion	Exp%	Act%
252.00	100	100
253.00	22.30	21.89
125.00	9.90	11.47
0.00	0.00	0.00

Quantitation Report (Qedit)

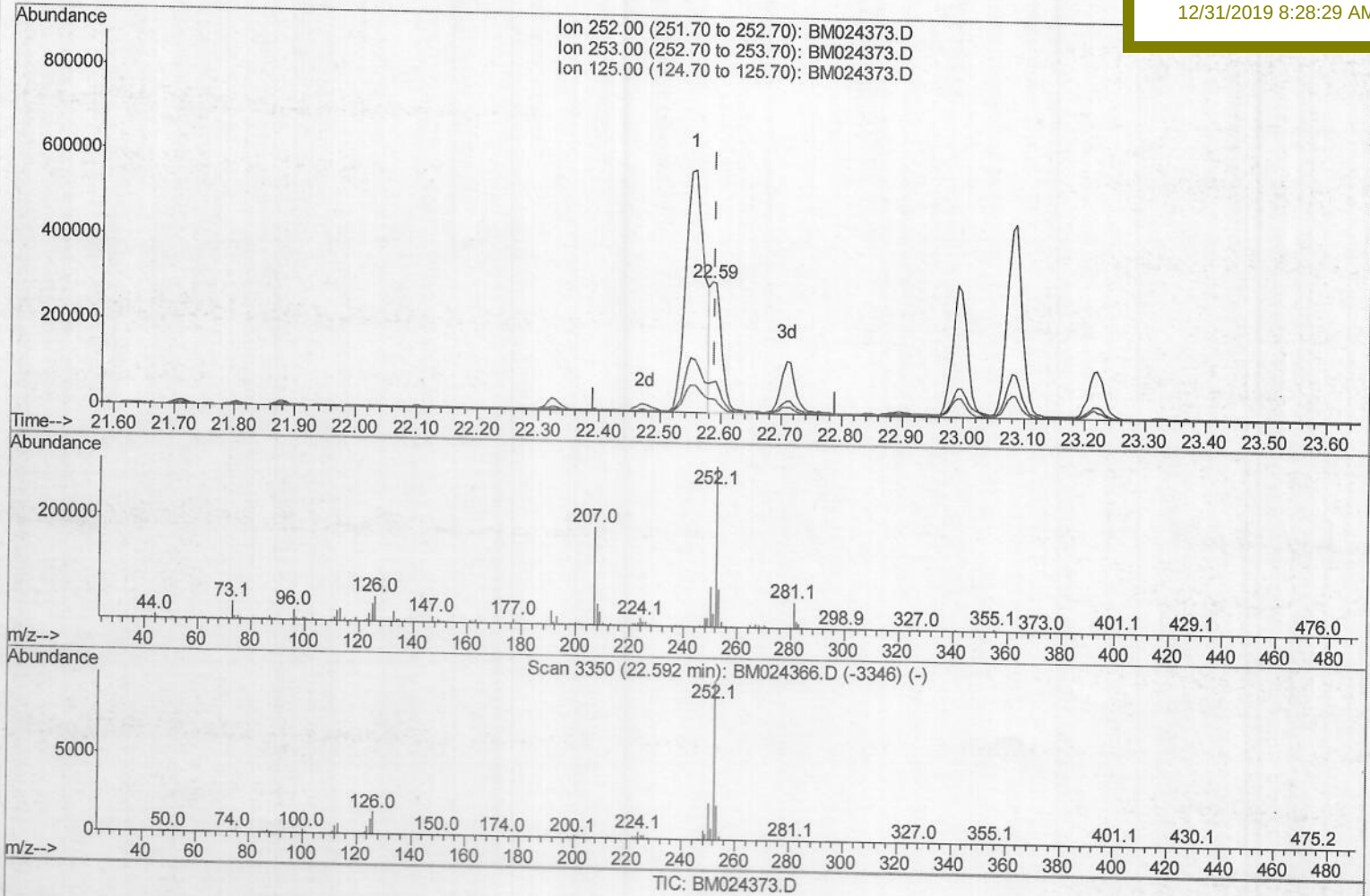
Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM123019\
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 Misc :
 ALS Vial : 9 Sample Multiplier: 1

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Instrument :
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 C0C30DL

Manual Integrations
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(89) Benzo(k)fluoranthene

22.586min (-0.006) 5.67ng/ul m > JU 12/30/19
 response 384041

Ion	Exp%	Act%
252.00	100	100
253.00	22.30	23.58
125.00	9.90	11.33
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	7.41	152	188702	20.00	ng/ul	0.00
18) Naphthalene-d8	10.17	136	793858	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.07	164	461312	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.83	188	960733	20.00	ng/ul	0.00
78) Chrysene-d12	21.05	240	965816	20.00	ng/ul	0.00
86) Perylene-d12	23.17	264	1153848	20.00	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.01	96	2438	0.57	ng/uL	0.03
5) Phenol-d5	6.63	99	29333	1.64	ng/ul	0.02
7) Bis-(2-Chloroethyl)ether-d	6.77	67	23307	2.18	ng/ul	0.01
9) 2-Chlorophenol-d4	6.96	132	27298	2.14	ng/ul	0.00
13) 4-Methylphenol-d8	8.16	113	20999	1.44	ng/ul	0.02
19) Nitrobenzene-d5	8.58	128	11397	2.00	ng/ul	0.02
22) 2-Nitrophenol-d4	9.29	143	11044	1.76	ng/ul	0.01
26) 2,4-Dichlorophenol-d3	9.86	165	22685m	1.86	ng/ul	0.04
29) 4-Chloroaniline-d4	10.38	131	17116m	1.16	ng/ul	0.05
44) Dimethylphthalate-d6	13.50	166	88183	2.59	ng/ul	0.00
47) Acenaphthylene-d8	13.76	160	97540	2.39	ng/ul	0.00
52) 4-Nitrophenol-d4	14.43	143	364	0.06	ng/ul	0.09
58) Fluorene-d10	15.08	176	79278	2.65	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.27	200	2467	0.42	ng/ul	0.03
71) Anthracene-d10	16.93	188	124200	2.84	ng/ul	0.00
79) Pyrene-d10	19.24	212	138814	3.02	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.04	264	162184	2.83	ng/ul	0.00
Target Compounds						
48) Acenaphthylene	13.79	152	166709	3.842	ng/ul	98
59) Fluorene	15.14	166	61087	1.744	ng/ul	96
70) Phenanthrene	16.87	178	980373	18.227	ng/ul	99
72) Anthracene	16.97	178	273431	5.057	ng/ul	99
77) Fluoranthene	18.91	202	2068196	35.280	ng/ul	99
80) Pyrene	19.27	202	1632205	25.887	ng/ul	99
83) Benzo(a)anthracene	21.04	228	1045026	16.983	ng/ul	99
85) Chrysene	21.09	228	900644	14.932	ng/ul	97
88) Benzo(b)fluoranthene	22.56	252	1241603	17.984	ng/ul	98
89) Benzo(k)fluoranthene	22.59	252	384041m	5.669	ng/ul	98
91) Benzo(a)pyrene	23.09	252	765096	12.250	ng/ul	98
92) Indeno(1,2,3-cd)pyrene	25.26	276	476133	6.349	ng/ul	97
93) Dibenzo(a,h)anthracene	25.26	278	146318	2.327	ng/ul#	87
94) Benzo(g,h,i)perylene	25.90	276	390020	6.369	ng/ul	98

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

JU 12/30/19

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