

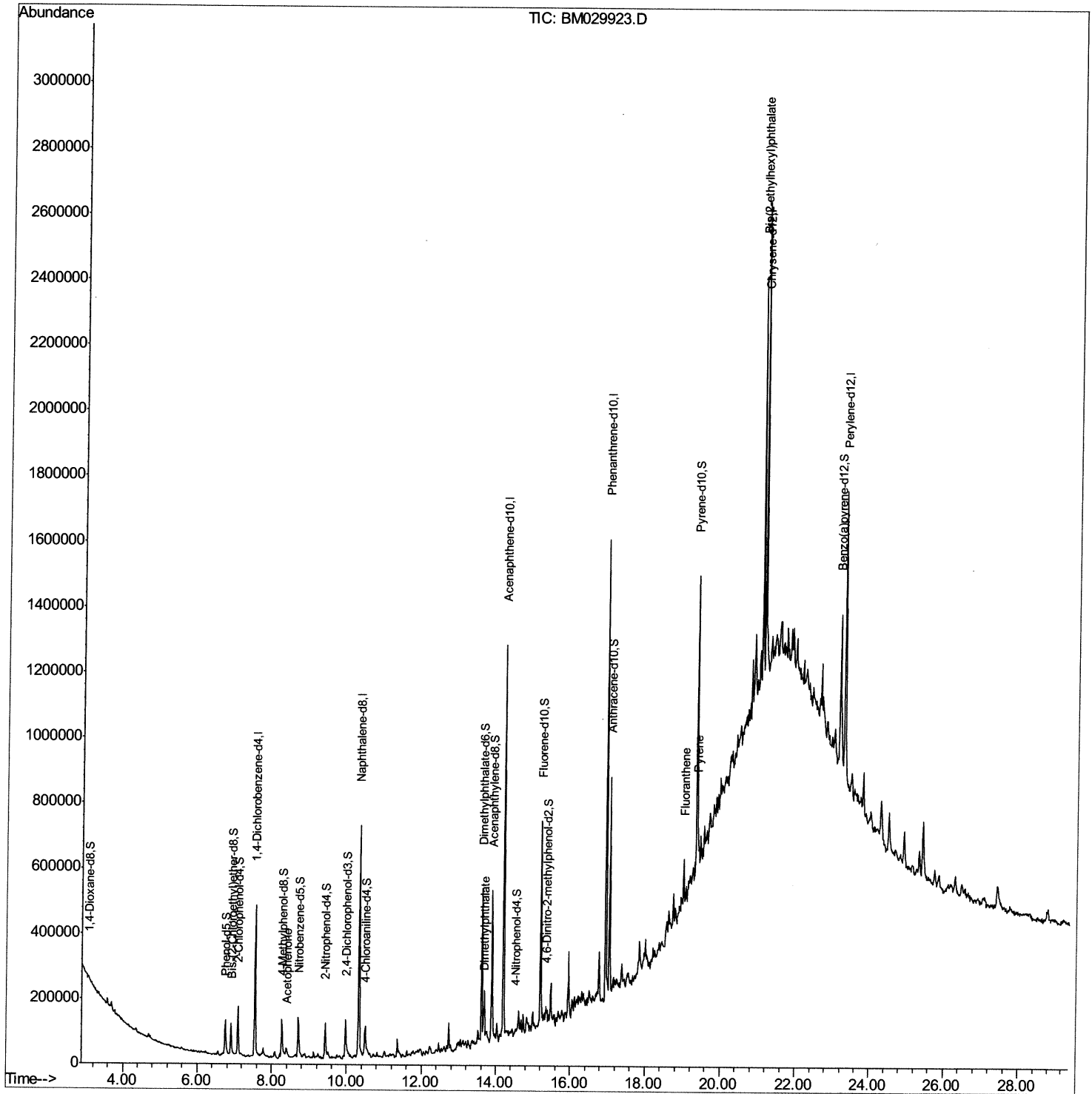
Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM051021\
Data File : BM029923.D
Acq On : 11 May 2021 02:12
Operator : CG/JU
Sample : M2177-19
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BG589

Manual Integrations
APPROVED

mohammad
5/11/2021 2:33:14 PM

Quant Time: May 11 03:21:50 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM050721.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue May 11 01:19:00 2021
Response via : Initial Calibration



Quantitation Report (Qedit)

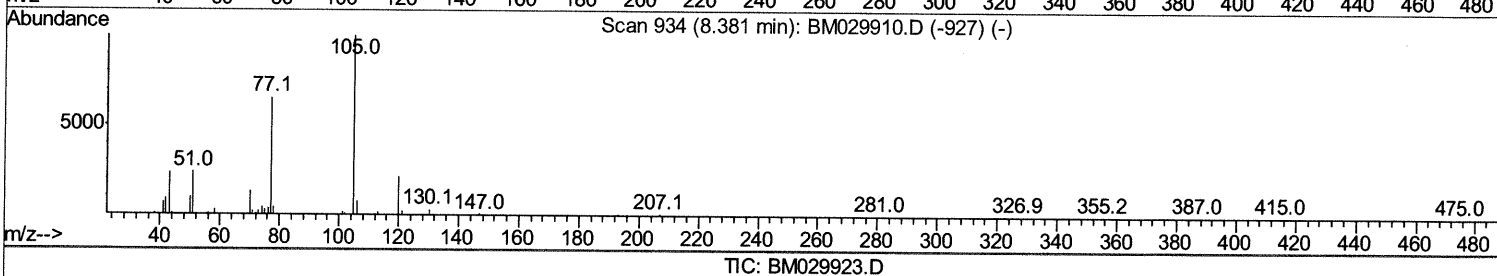
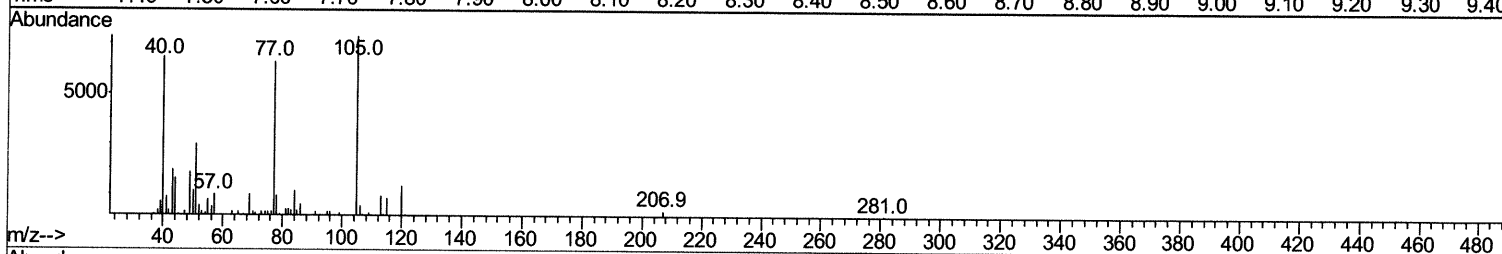
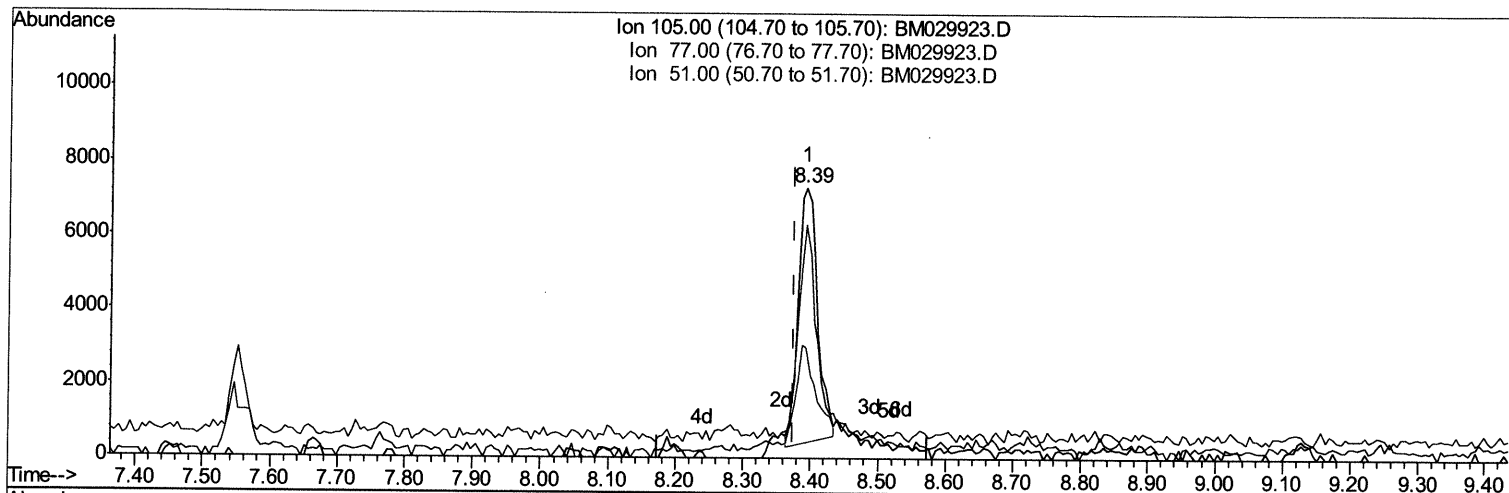
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TIC: BM029923.D

(16) Acetophenone

8.393min (+0.018) 0.88ng/ul

response 13754

Ion	Exp%	Act%
105.00	100	100
77.00	80.30	86.24
51.00	30.40	40.85#
0.00	0.00	0.00

Quantitation Report (Qedit)

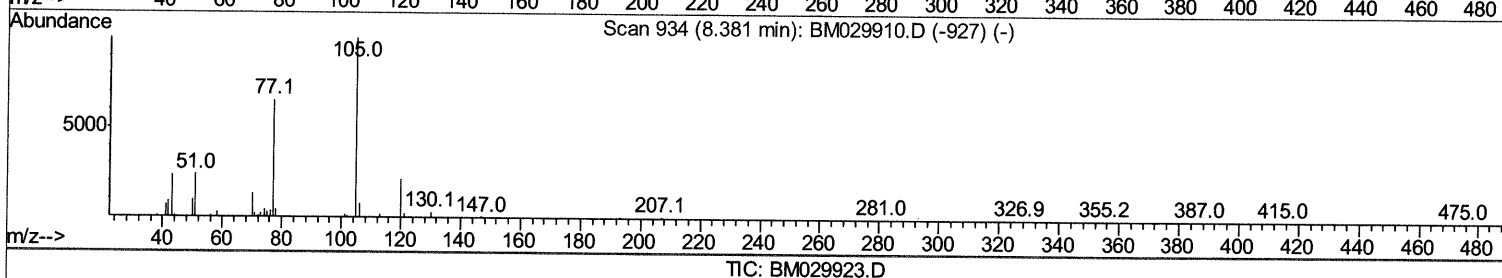
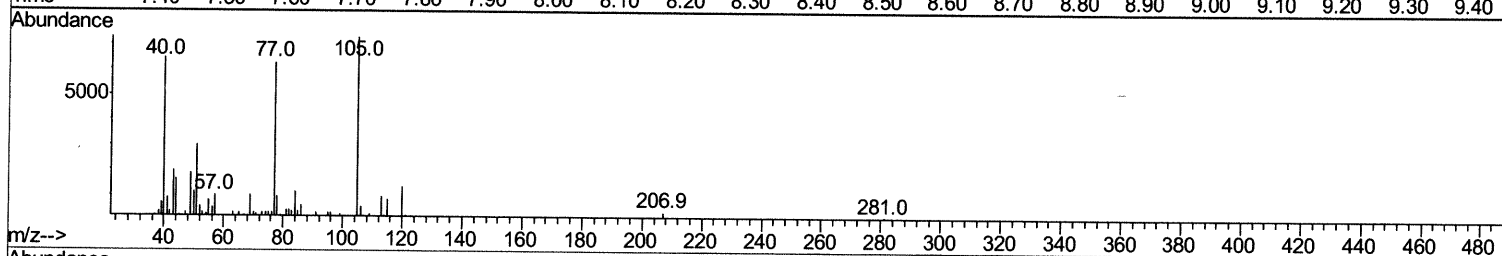
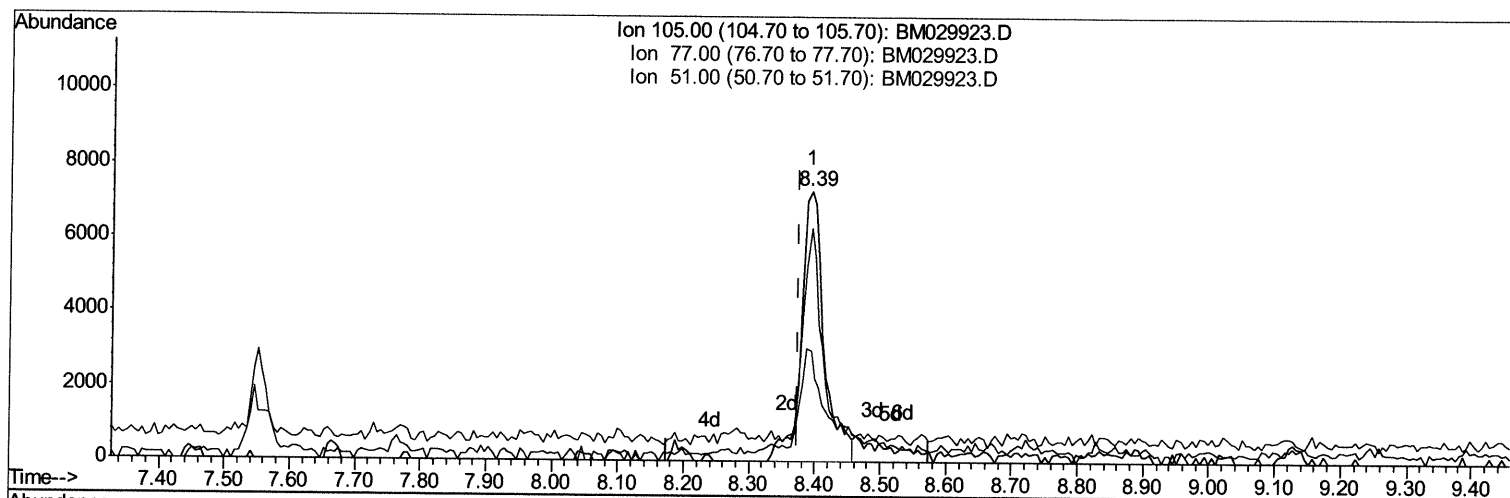
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(16) Acetophenone

8.393min (+0.018) 1.14ng/ul m 05/12/2021

response 17872

Ion	Exp%	Act%
105.00	100	100
77.00	80.30	86.24
51.00	30.40	40.85#
0.00	0.00	0.00

Quantitation Report (Qedit)

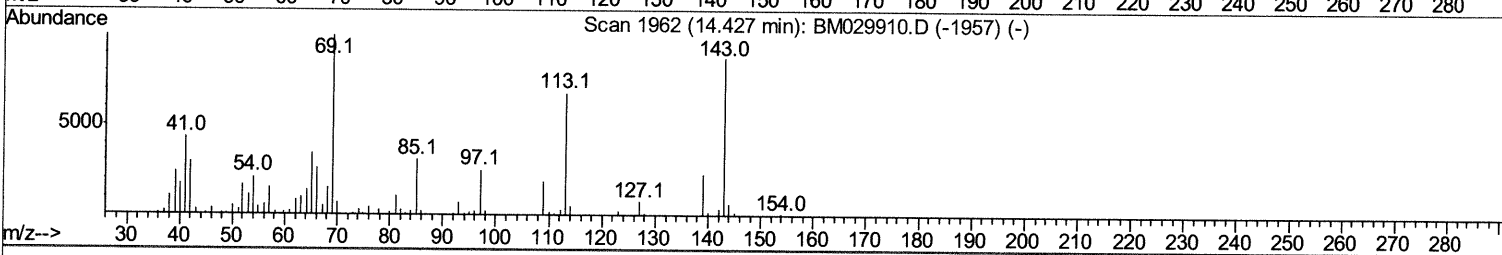
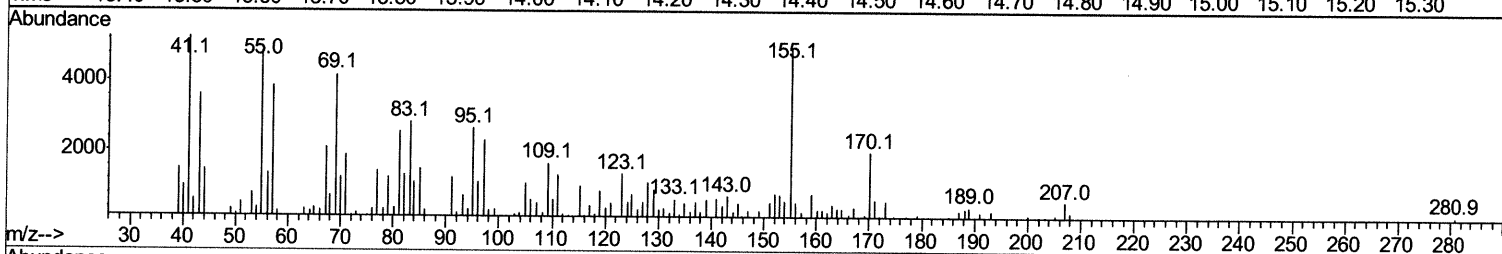
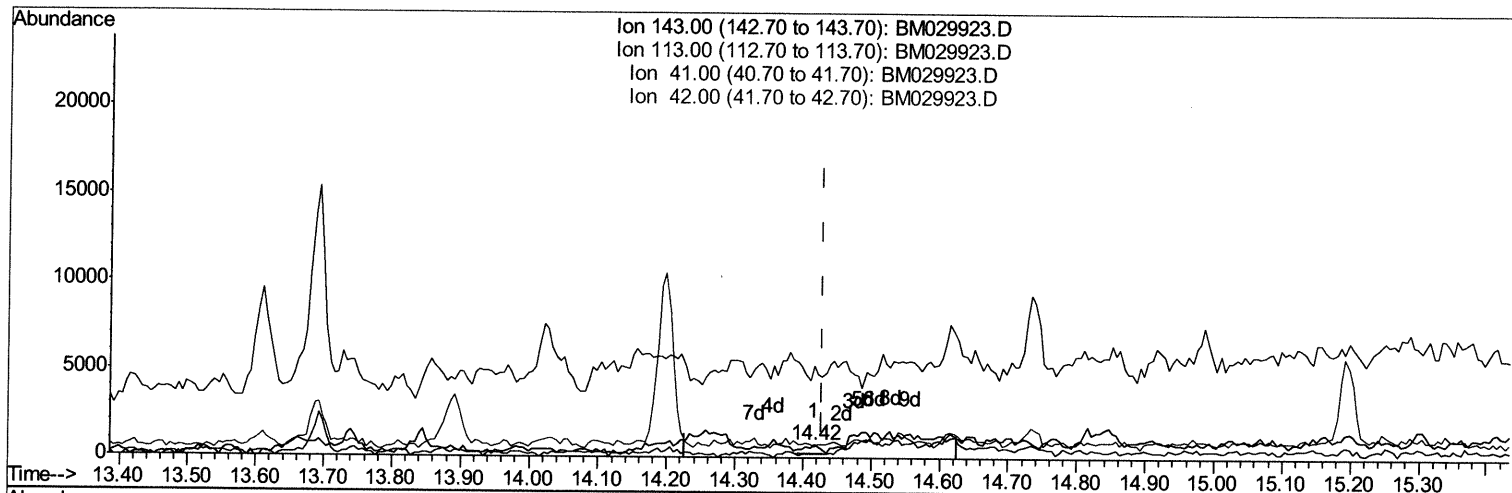
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TIC: BM029923.D

(54) 4-Nitrophenol-d4 (S)

14.416min (-0.011) 0.13ng/ul

response 661

Ion	Exp%	Act%
143.00	100	100
113.00	71.10	27.24#
41.00	44.70	704.41#
42.00	28.30	87.58#

Quantitation Report (Qedit)

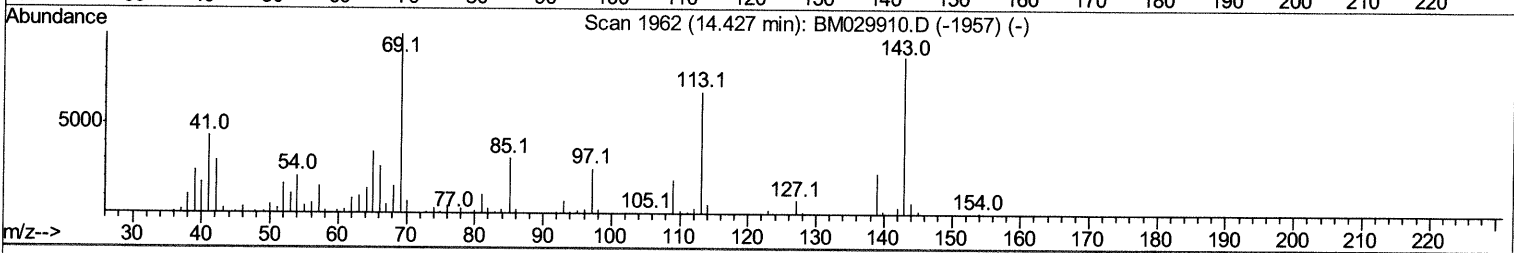
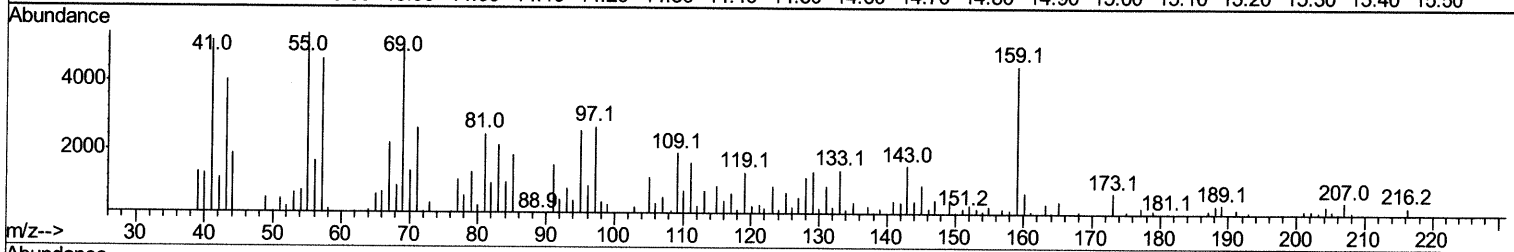
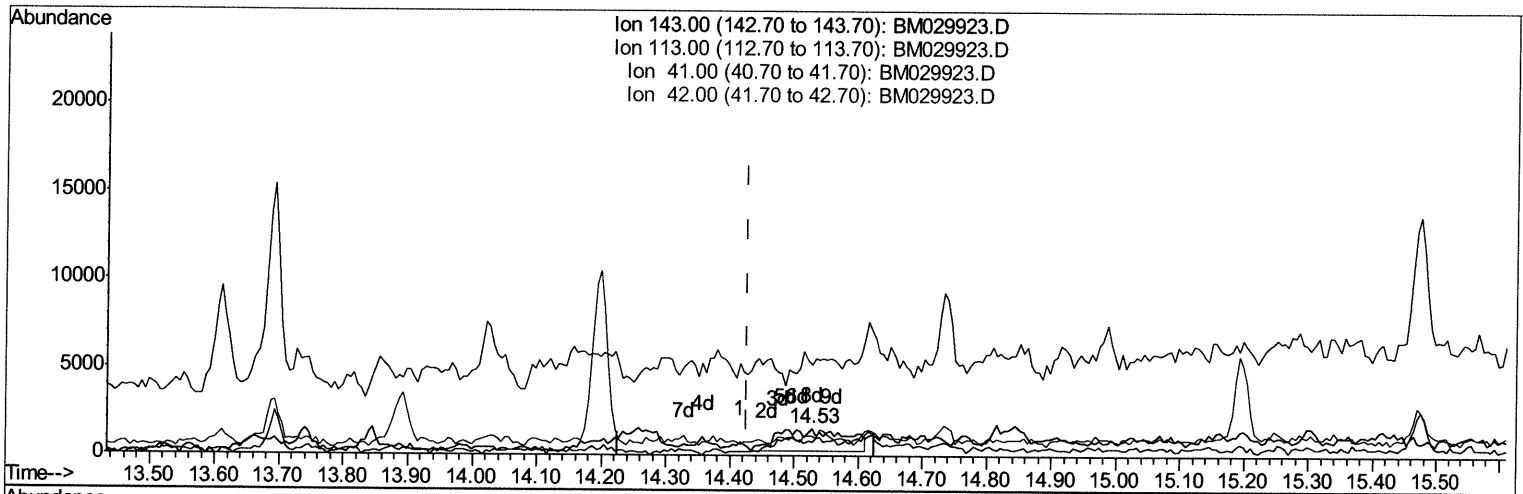
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 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampled :
 BG589

Manual Integrations
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TIC: BM029923.D

(54) 4-Nitrophenol-d4 (S)

14.528min (+0.100) 1.82ng/ul m 05/12/2021

response 8994

Ion	Exp%	Act%
143.00	100	100
113.00	71.10	51.00#
41.00	44.70	333.14#
42.00	28.30	73.95#

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 Misc :
 ALS Vial : 22 Sample Multiplier: 1

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.55	152	126892	20.00	ng/ul	0.00
20) Naphthalene-d8	10.33	136	588199	20.00	ng/ul	0.00
38) Acenaphthene-d10	14.20	164	392172	20.00	ng/ul	0.00
64) Phenanthrene-d10	16.95	188	796827	20.00	ng/ul	0.00
79) Chrysene-d12	21.13	240	687788	20.00	ng/ul	0.00
88) Perylene-d12	23.29	264	585551	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.08	96	5881	1.85	ng/uL	0.01
4) Pyridine-d5	0.00	84	0d	0.00	ng/ul	
7) Phenol-d5	6.74	99	78303	6.64	ng/ul	0.00
9) Bis-(2-Chloroethyl)ether-d	6.90	67	54659	7.47	ng/ul	0.00
11) 2-Chlorophenol-d4	7.09	132	73360	8.05	ng/ul	0.00
15) 4-Methylphenol-d8	8.28	113	60325	6.18	ng/ul	0.00
21) Nitrobenzene-d5	8.72	128	32057	7.91	ng/ul	0.00
24) 2-Nitrophenol-d4	9.43	143	35972	7.86	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.97	165	62740	6.84	ng/ul	0.01
31) 4-Chloroaniline-d4	10.49	131	54038	3.87	ng/ul	0.01
46) Dimethylphthalate-d6	13.62	166	275615	9.07	ng/ul	0.00
49) Acenaphthylene-d8	13.89	160	325954	8.46	ng/ul	0.00
54) 4-Nitrophenol-d4	14.53	143	8994m >	1.82	ng/ul >	0.1005/12.1217u
60) Fluorene-d10	15.20	176	245989	9.53	ng/ul	0.00
65) 4,6-Dinitro-2-methylphenol	15.34	200	8338	1.61	ng/ul	0.00
73) Anthracene-d10	17.05	188	371591	9.26	ng/ul	0.00
81) Pyrene-d10	19.34	212	423671	12.01	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.14	264	315253	9.50	ng/ul	0.00

Target Compounds

					Ovalue
16) Acetophenone	8.39	105	17872m	1.141	ng/ul
47) Dimethylphthalate	13.67	163	37465	1.231	ng/ul 99
80) Fluoranthene	19.00	202	75292	1.753	ng/ul 97
82) Pyrene	19.37	202	81812	1.810	ng/ul 99
86) Bis(2-ethylhexyl)phthalate	21.06	149	400508	14.464	ng/ul 100

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