

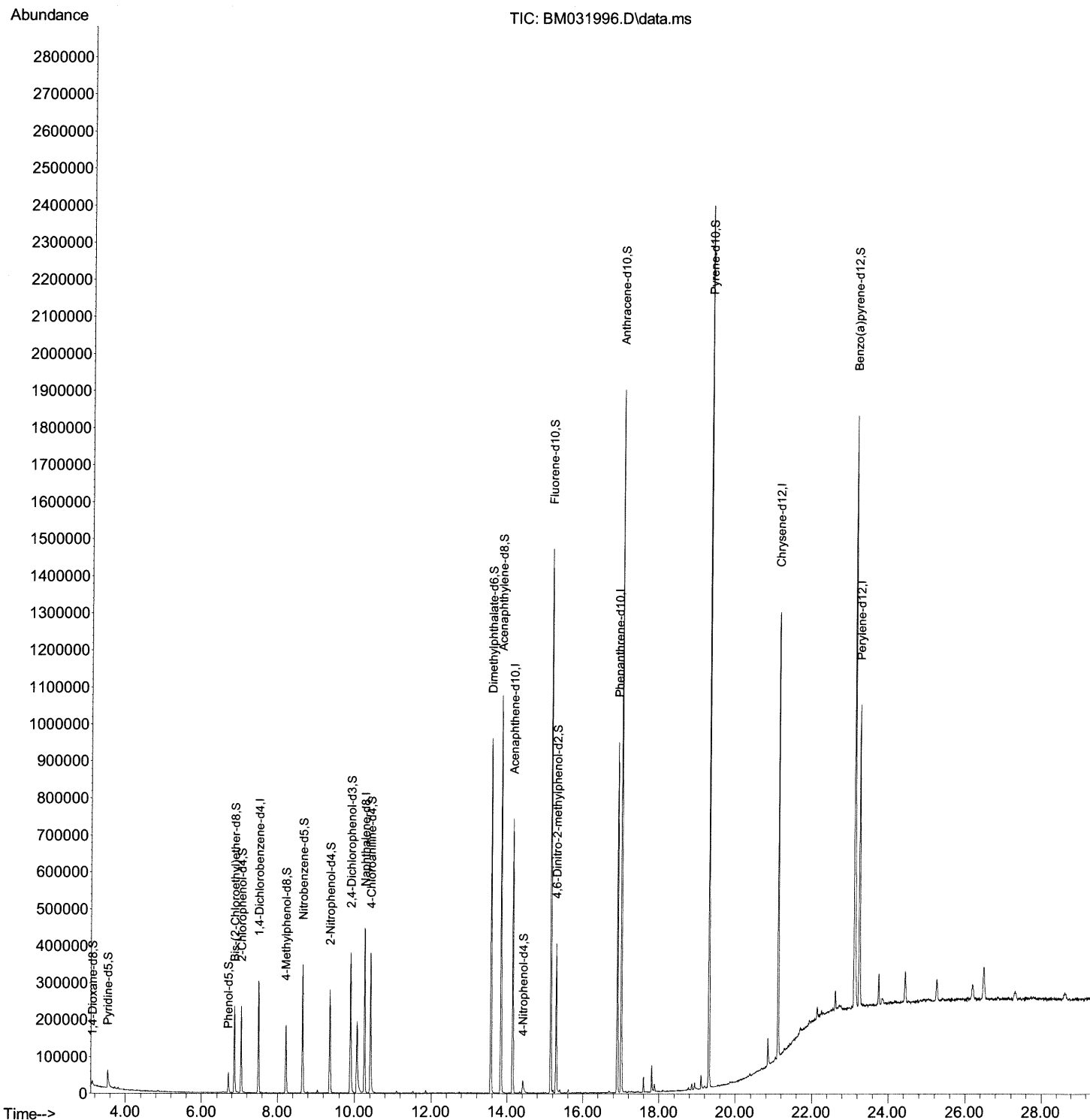
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM083021\
Data File : BM031996.D
Acq On : 31 Aug 2021 22:07
Operator : CG/JU
Sample : M3517-06
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0AB9

Manual Integrations
APPROVED

mohammad
9/1/2021 4:03:34 PM

Quant Time: Sep 01 01:05:03 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM082821.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Aug 30 09:21:30 2021
Response via : Initial Calibration



Quantitation Report (Qedit)

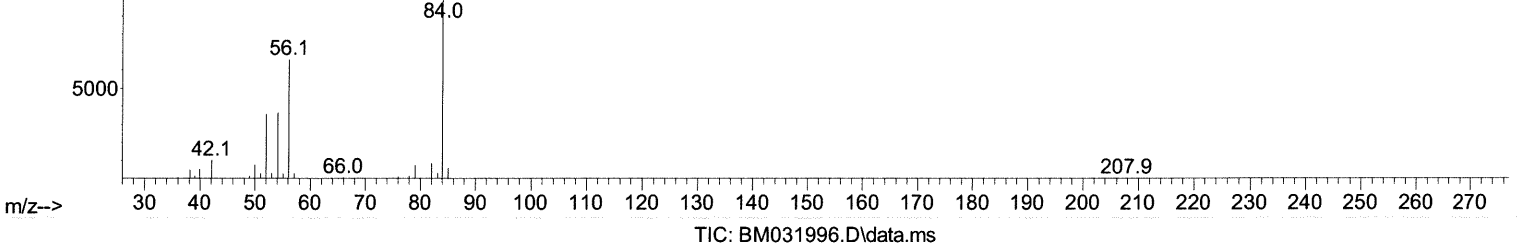
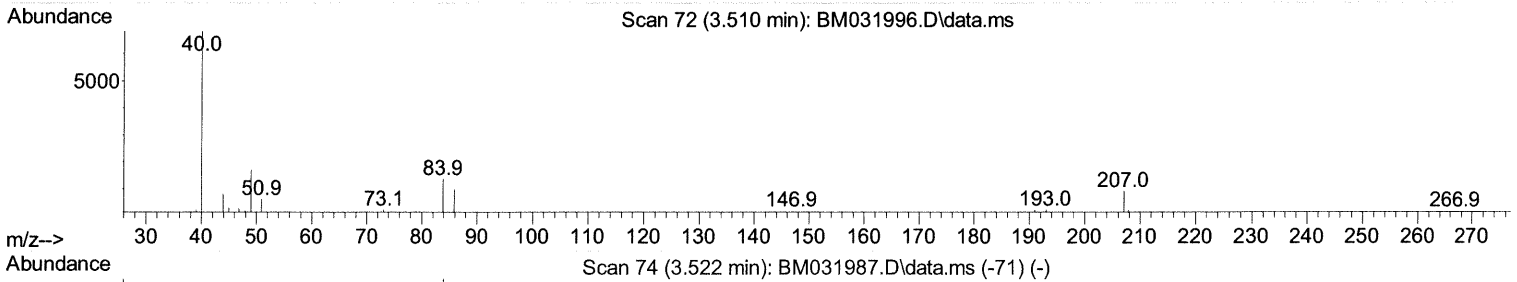
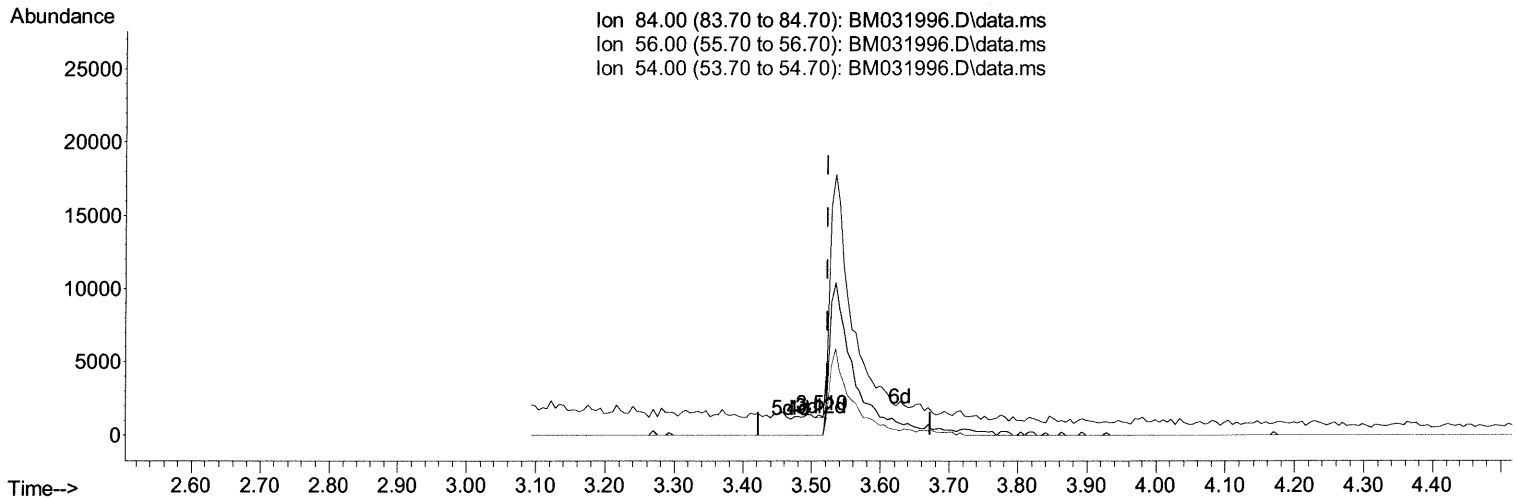
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(4) Pyridine-d5 (S)

3.510min (-0.012) 0.02 ng/ul

response 86

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	73.70	0.00#
54.00	35.80	0.00#
0.00	0.00	0.00

Quantitation Report (Qedit)

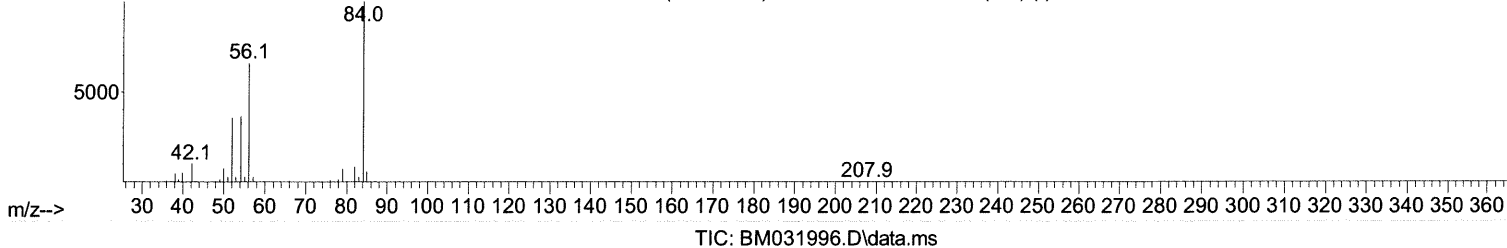
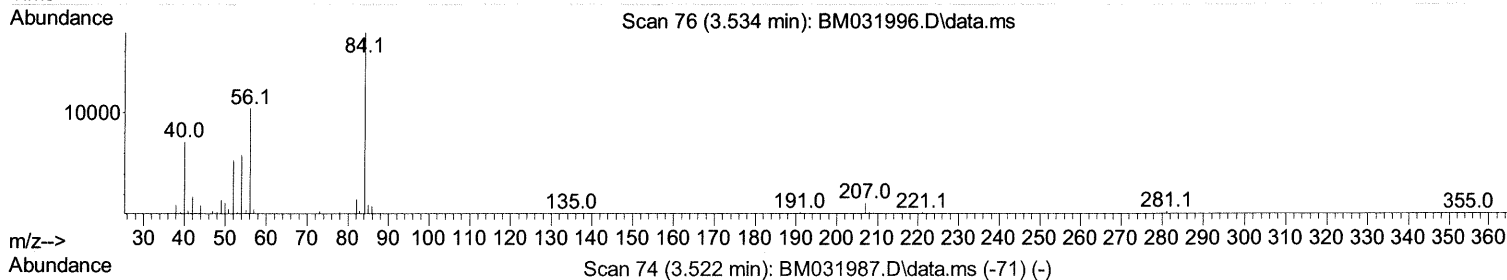
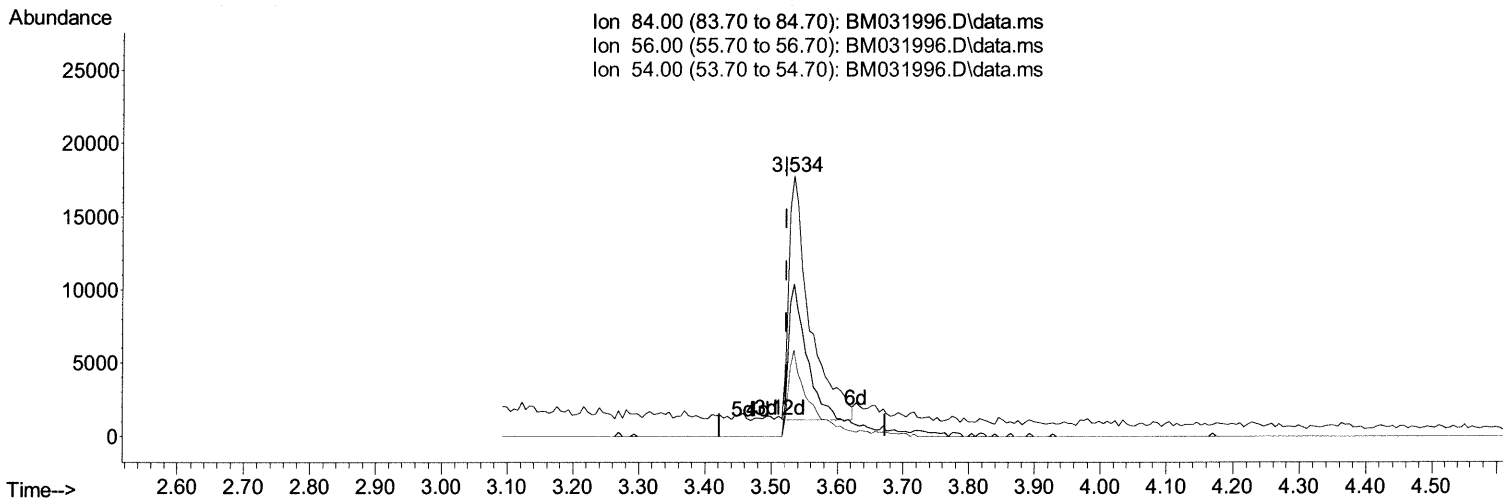
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Manual Integrations
 APPROVED

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 9/1/2021 4:03:34 PM

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(4) Pyridine-d5 (S)

3.534min (+ 0.012) 7.49 ng/ul m

09/03/2021 JU

response 36633

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	73.70	58.55#
54.00	35.80	33.07
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM083021\
 Data File : BM031996.D
 Acq On : 31 Aug 2021 22:07
 Operator : CG/JU
 Sample : M3517-06
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AB9

Manual Integrations
 APPROVED

mohammad
 9/1/2021 4:03:34 PM

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.504	152	79421	20.000	ng/ul	0.00
20) Naphthalene-d8	10.263	136	352184	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.145	164	259226	20.000	ng/ul	0.00
64) Phenanthrene-d10	16.904	188	560512	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.109	240	574019	20.000	ng/ul	0.00
88) Perylene-d12	23.244	264	527830	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.134	96	7841	4.555	ng/ul	0.00
4) Pyridine-d5	3.534	84	36633m	7.494	ng/ul	> 0.01
7) Phenol-d5	6.698	99	30013	4.557	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.863	67	104175	27.339	ng/ul	0.00
11) 2-Chlorophenol-d4	7.045	132	103617	20.446	ng/ul	0.00
15) 4-Methylphenol-d8	8.216	113	65930	11.668	ng/ul	0.00
21) Nitrobenzene-d5	8.657	128	77421	30.266	ng/ul	0.00
24) 2-Nitrophenol-d4	9.363	143	81688	26.984	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.898	165	144753	23.832	ng/ul	0.00
31) 4-Chloroaniline-d4	10.416	131	196107	22.889	ng/ul	0.00
46) Dimethylphthalate-d6	13.568	166	642593	32.760	ng/ul	0.00
49) Acenaphthylene-d8	13.833	160	716407	30.840	ng/ul	0.00
54) 4-Nitrophenol-d4	14.410	143	13099	3.689	ng/ul	0.02
60) Fluorene-d10	15.145	176	576272	33.421	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.292	200	86450	22.446	ng/ul	0.00
73) Anthracene-d10	17.003	188	1045526	39.299	ng/ul	0.00
81) Pyrene-d10	19.309	212	1225105	45.566	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.115	264	1073415	38.695	ng/ul	0.00

Target Compounds Qvalue

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