

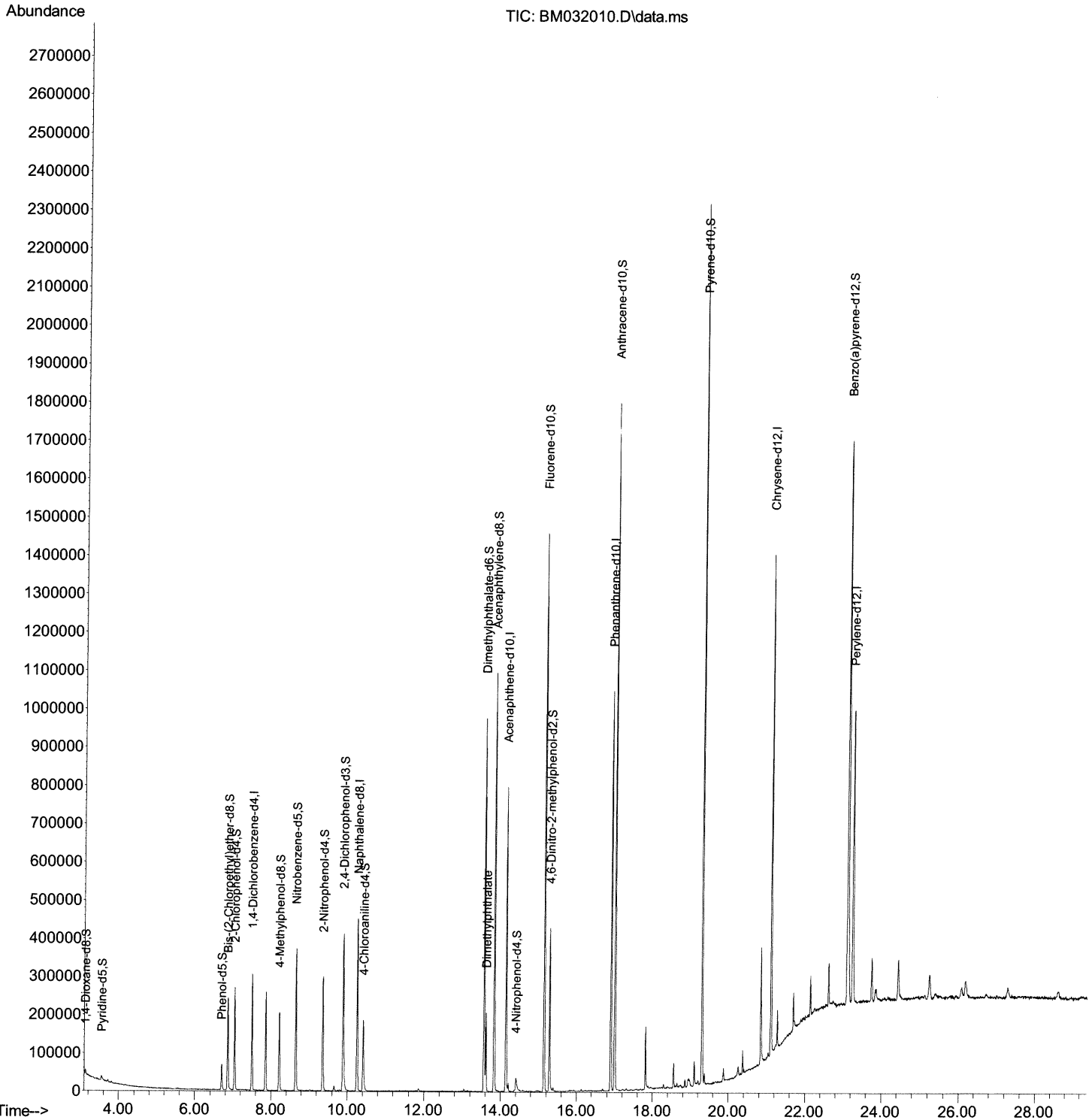
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM083021\
Data File : BM032010.D
Acq On : 01 Sep 2021 07:21
Operator : CG/JU
Sample : M3494-06
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
YAS57

Manual Integrations
APPROVED

mohammad
9/1/2021 4:04:12 PM

Quant Time: Sep 01 08:05:12 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)

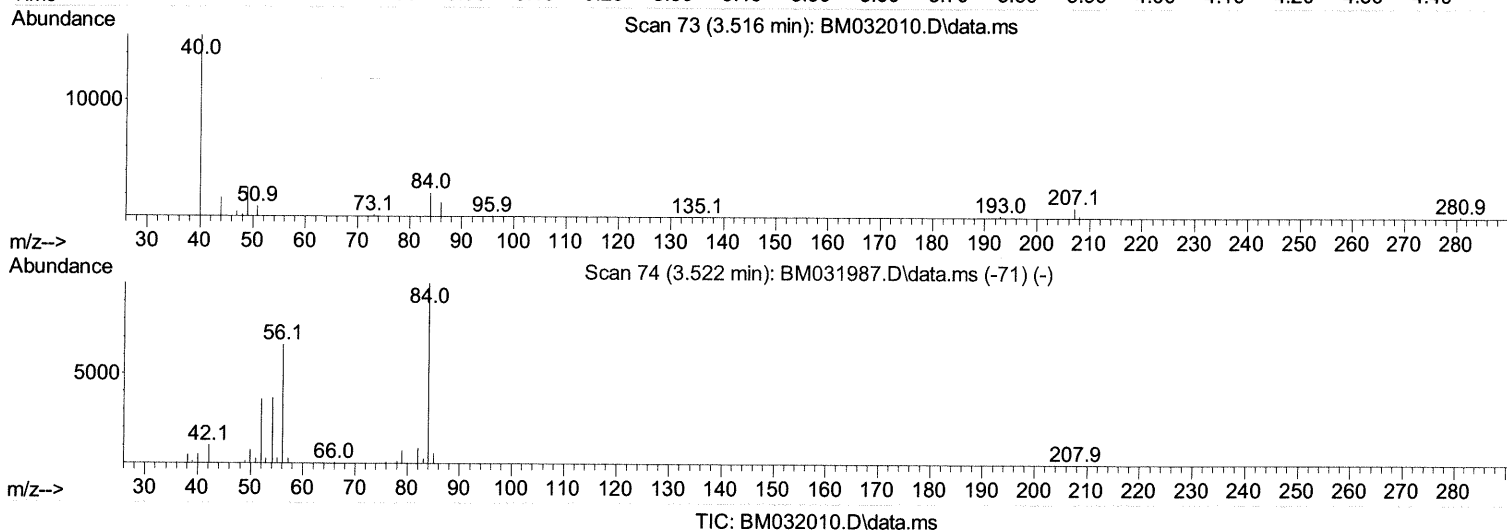
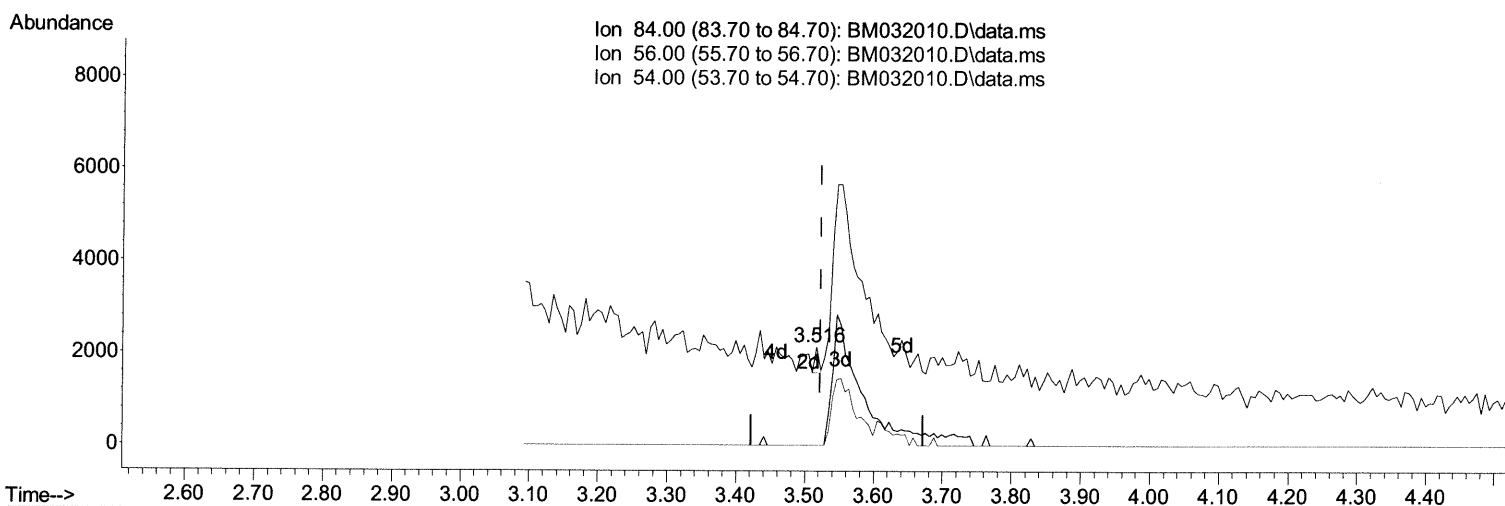
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(4) Pyridine-d5 (S)

3.516min (-0.006) 0.04 ng/ul

response 216

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)

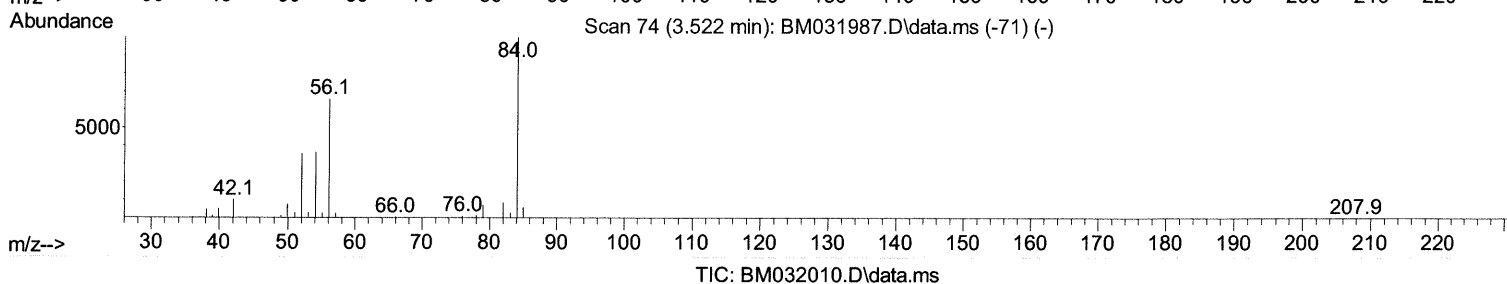
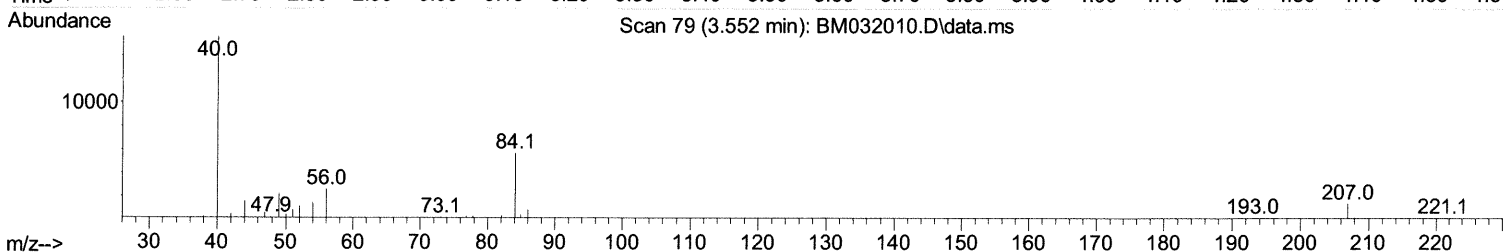
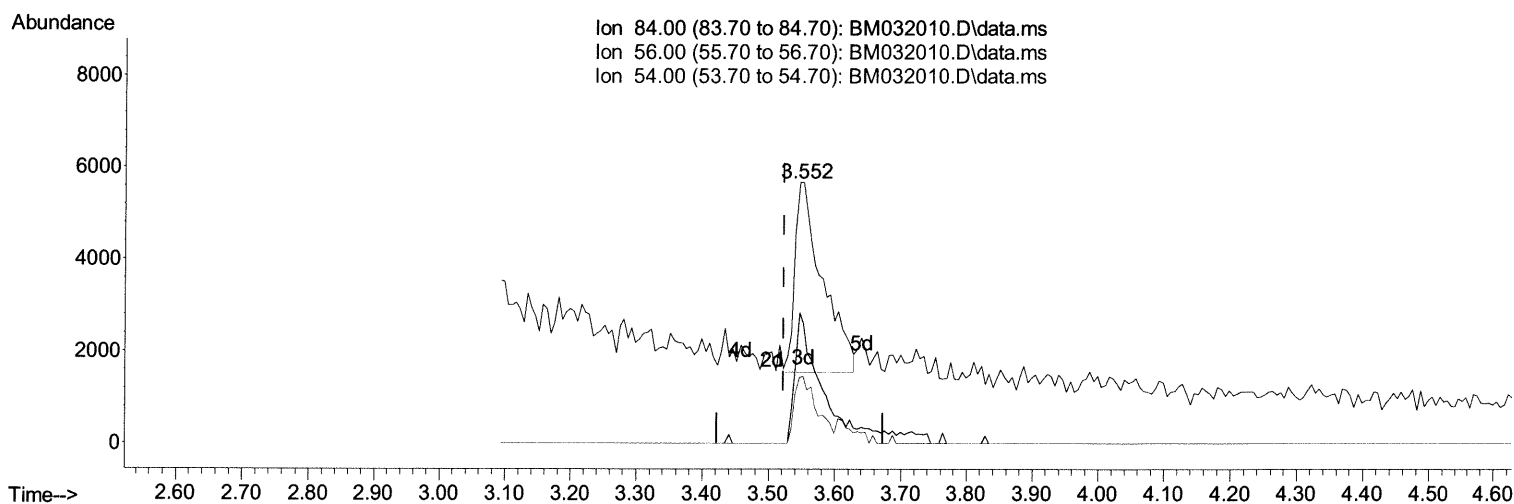
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM083021\
Data File : BM032010.D
Acq On : 01 Sep 2021 07:21
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Sample : M3494-06
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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Manual Integrations
APPROVED

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Quant Title : SVOA CALIBRATION
QLast Update : Mon Aug 30 09:21:30 2021
Response via : Initial Calibration



(4) Pyridine-d5 (S)

3.552min (+ 0.029) 2.41 ng/ul m 09/09/21 JU

response 11902

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	73.70	45.97#
54.00	35.80	25.60#
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM083021\
 Data File : BM032010.D
 Acq On : 01 Sep 2021 07:21
 Operator : CG/JU
 Sample : M3494-06
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 YAS57

Manual Integrations
 APPROVED

mohammad
 9/1/2021 4:04:12 PM

Quant Time: Sep 01 08:05:12 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

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.498	152	80124	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.257	136	351414	20.000	ng/ul	#-0.01
38) Acenaphthene-d10	14.139	164	263654	20.000	ng/ul	-0.01
64) Phenanthrene-d10	16.898	188	580638	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.103	240	588994	20.000	ng/ul	0.00
88) Perylene-d12	23.244	264	524670	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.128	96	8584	4.943	ng/ul	0.00
4) Pyridine-d5	3.552	84	11902m>	2.414	ng/ul>	0.03
7) Phenol-d5	6.698	99	36257	5.457	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.857	67	107613	27.993	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.040	132	114667	22.427	ng/ul	-0.01
15) 4-Methylphenol-d8	8.216	113	75578	13.258	ng/ul	0.00
21) Nitrobenzene-d5	8.651	128	78630	30.806	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.363	143	86772	28.726	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.892	165	158468	26.147	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.416	131	102023	11.934	ng/ul	0.00
46) Dimethylphthalate-d6	13.569	166	621975	31.177	ng/ul	0.00
49) Acenaphthylene-d8	13.827	160	751349	31.801	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.404	143	13384	3.706	ng/ul	0.02
60) Fluorene-d10	15.145	176	556420	31.728	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.292	200	92657	23.224	ng/ul	0.00
73) Anthracene-d10	16.998	188	965028	35.016	ng/ul	0.00
81) Pyrene-d10	19.303	212	1157715	41.965	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.109	264	974174	35.329	ng/ul	0.00
Target Compounds						
47) Dimethylphthalate	13.616	163	123603	6.310	ng/ul	99

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