

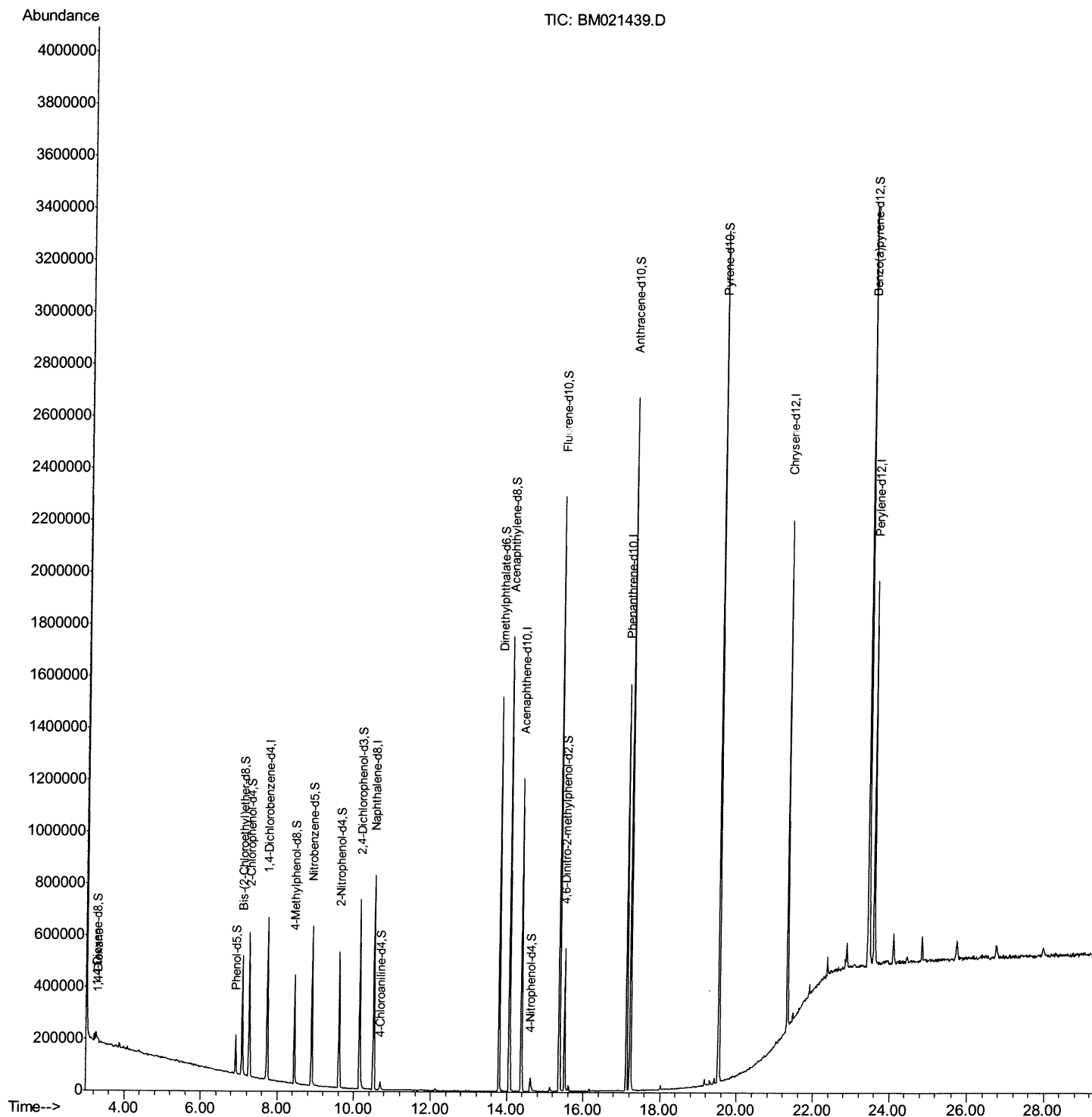
Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM071119\
Data File : BM021439.D
Acq On : 11 Jul 2019 22:00
Operator : HP/JU
Sample : K3762-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BF481

Manual Integrations
APPROVED

mohammad
7/12/2019 9:50:56 AM

Quant Time: Jul 12 06:03:10 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM070819MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Jul 08 17:45:45 2019
Response via : Initial Calibration



Quantitation Report (Qedit)

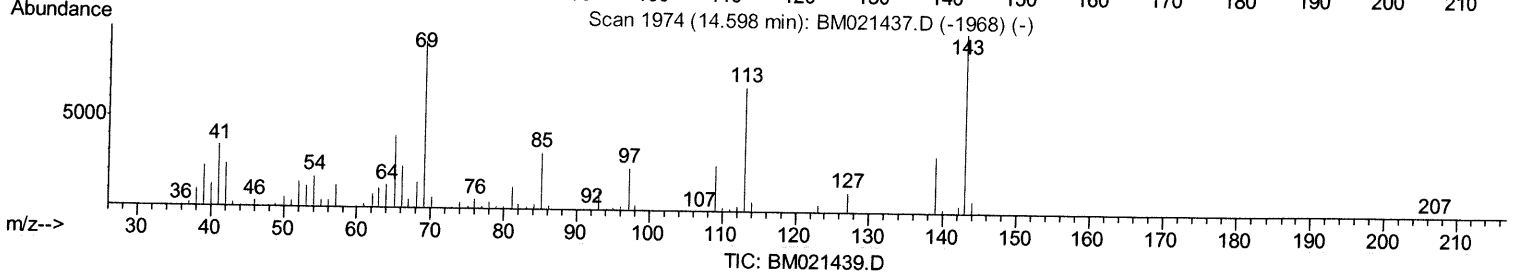
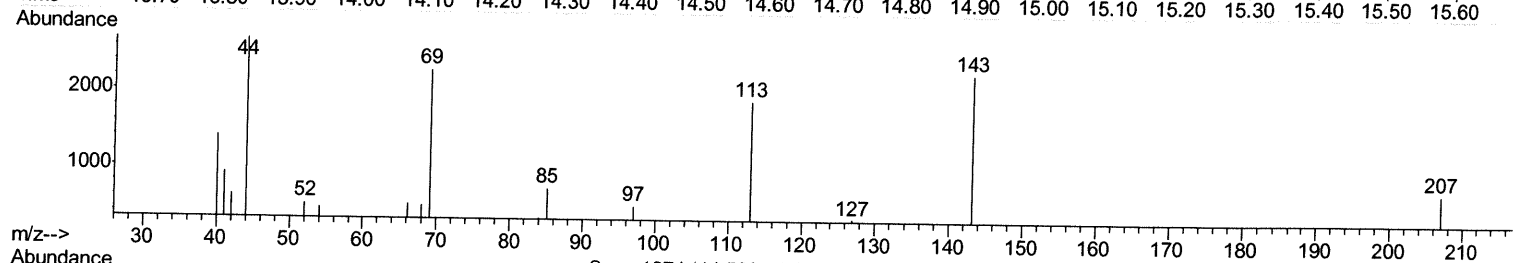
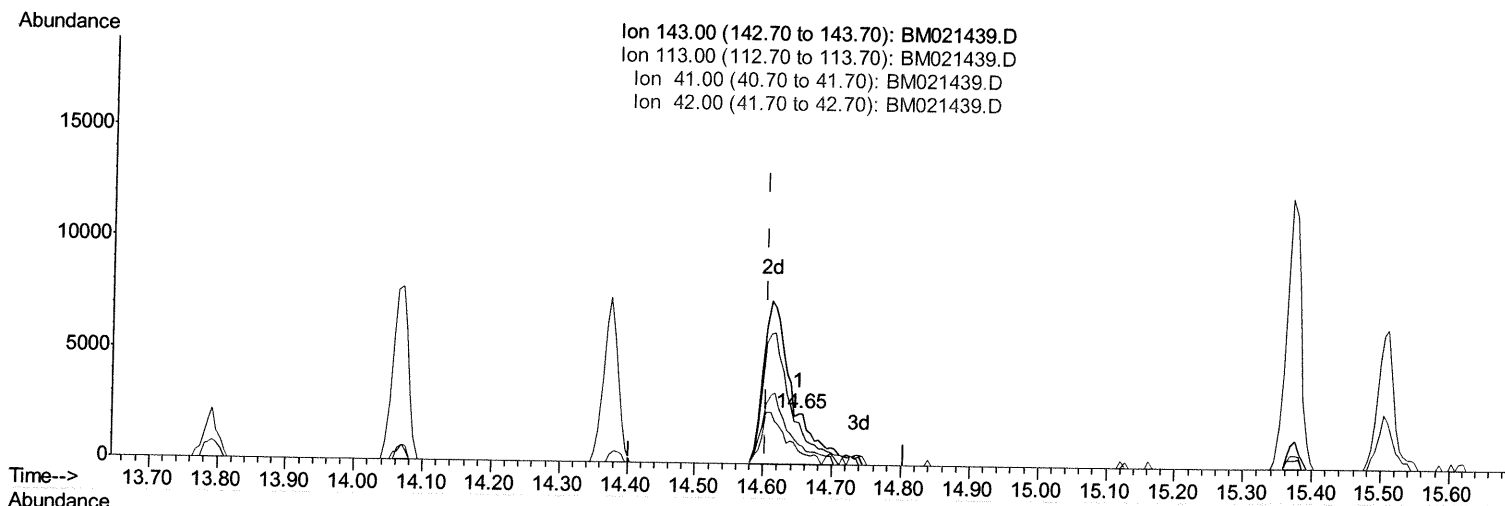
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ClientSampled :
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Response via : Initial Calibration



(51) 4-Nitrophenol-d4 (S)

14.651min (+0.047) 0.43ng/ul

response 2360

Ion	Exp%	Act%
143.00	100	100
113.00	74.00	83.55
41.00	34.20	40.21
42.00	23.00	27.28

Quantitation Report (Qedit)

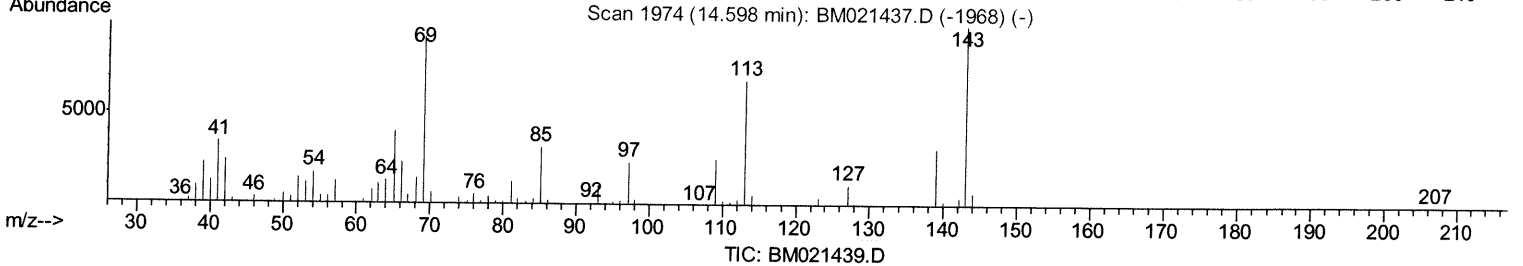
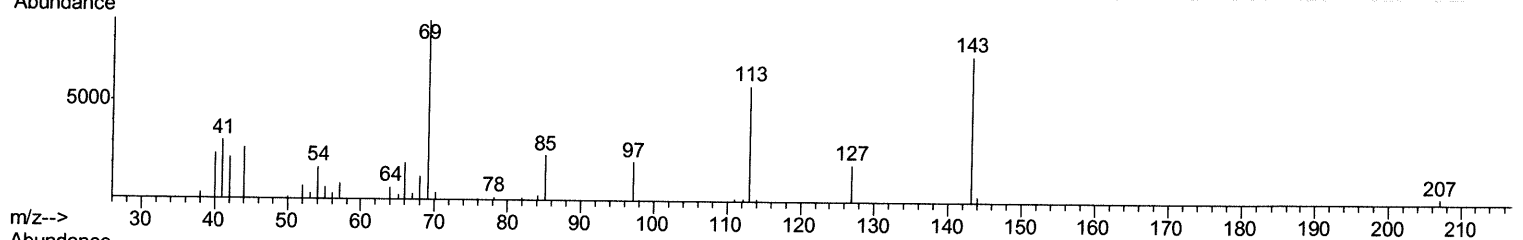
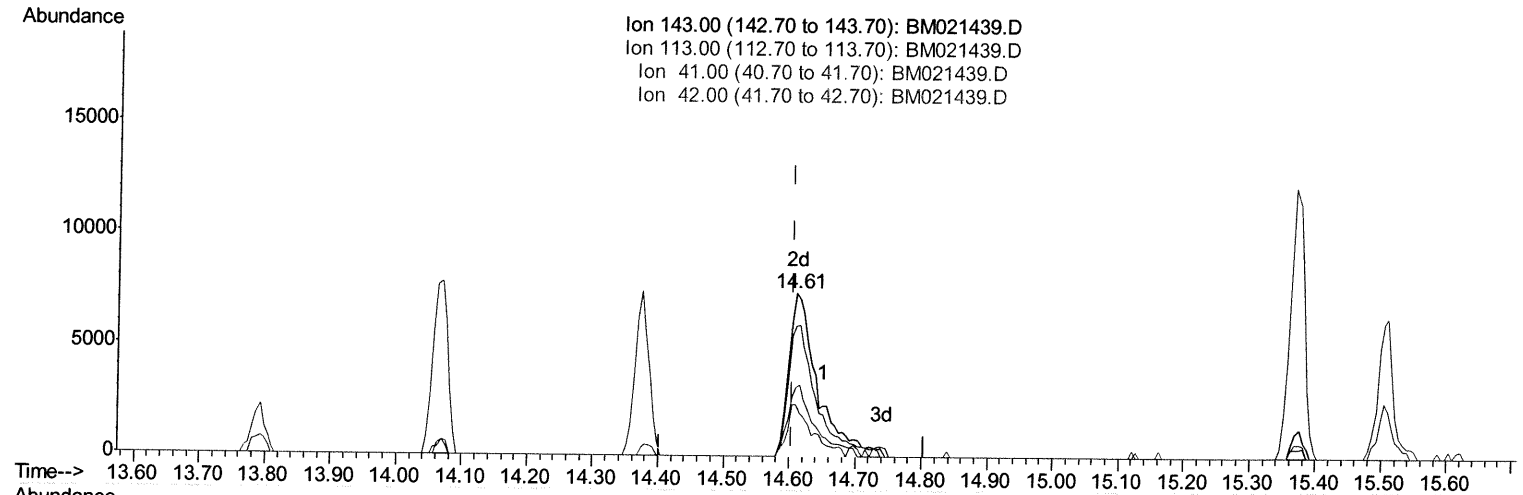
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Manual Integrations
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 QLast Update : Mon Jul 08 17:45:45 2019
 Response via : Initial Calibration



(51) 4-Nitrophenol-d4 (S)

14.610min (+0.006) 3.99ng/ul m

JU
 07/12/19

response 21850

Ion	Exp%	Act%
143.00	100	100
113.00	74.00	79.89
41.00	34.20	42.41#
42.00	23.00	31.50#

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM071119\
 Data File : BM021439.D
 Acq On : 11 Jul 2019 22:00
 Operator : HP/JU
 Sample : K3762-01
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
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Manual Integrations
 APPROVED

mohammad
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Quant Time: Jul 12 06:03:10 2019
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 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Jul 08 17:45:45 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.73	152	174333	20.00	ng/ul	0.00
18) Naphthalene-d8	10.52	136	682521	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.37	164	394776	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.13	188	906987	20.00	ng/ul	0.00
77) Chrysene-d12	21.32	240	908255	20.00	ng/ul	0.00
85) Perylene-d12	23.58	264	1076844	20.00	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.26	96	19993	5.65	ng/uL	0.00
5) Phenol-d5	6.91	99	87542	6.46	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.07	67	210335	26.17	ng/ul	0.00
9) 2-Chlorophenol-d4	7.27	132	260565	22.33	ng/ul	0.00
13) 4-Methylphenol-d8	8.44	113	163051	15.10	ng/ul	0.00
19) Nitrobenzene-d5	8.89	128	153254	27.26	ng/ul	0.00
22) 2-Nitrophenol-d4	9.61	143	171282	26.94	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.15	165	312028	24.61	ng/ul	0.00
29) 4-Chloroaniline-d4	10.69	131	23781	2.07	ng/ul	0.01
43) Dimethylphthalate-d6	13.79	166	1051535	30.41	ng/ul	0.00
46) Acenaphthylene-d8	14.07	160	1184778	27.29	ng/ul	0.00
51) 4-Nitrophenol-d4	14.61	143	21850m>	3.99	ng/ul	0.00
57) Fluorene-d10	15.37	176	933173	30.24	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.51	200	128580	21.09	ng/ul	0.00
70) Anthracene-d10	17.23	188	1560677	33.13	ng/ul	0.00
78) Pyrene-d10	19.52	212	1771478	35.54	ng/ul	0.00
89) Benzo(a)pyrene-d12	23.43	264	2123558	34.29	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	3.30	88	5161	1.387	ng/uL	96

JU 07/12/19

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