

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032519\

Data File : BM019438.D

Acq On : 25 Mar 2019 19:14

Operator : JU/SJ

Sample : K2027-15

Misc :

ALS Vial : 18 Sample Multiplier: 1

Quant Time: Apr 05 07:30:28 2019

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM032219MA.M

Quant Title : SVOA CALIBRATION

QLast Update : Mon Mar 25 03:11:42 2019

Response via : Initial Calibration

Instrument :

BNA_M

ClientSampleId :

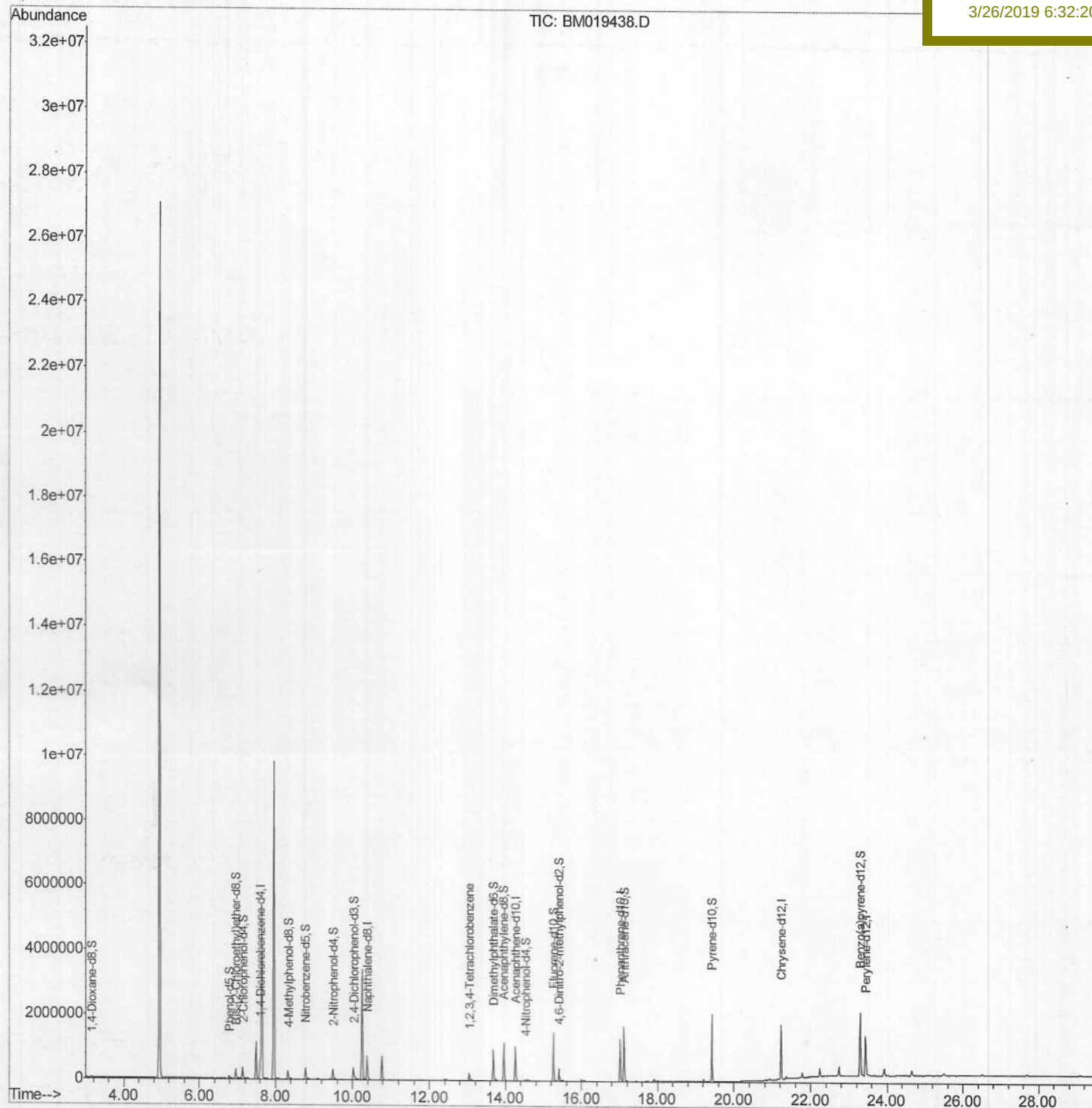
C09D1

Manual Integrations

APPROVED

Sohil

3/26/2019 6:32:20 PM



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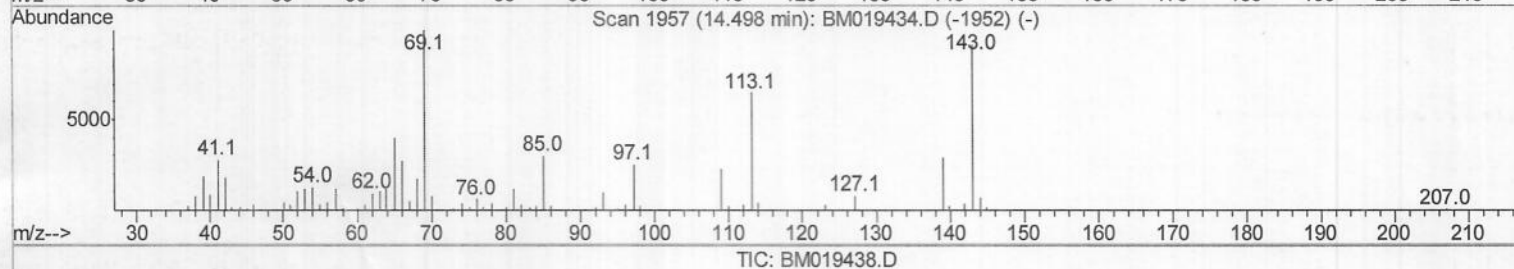
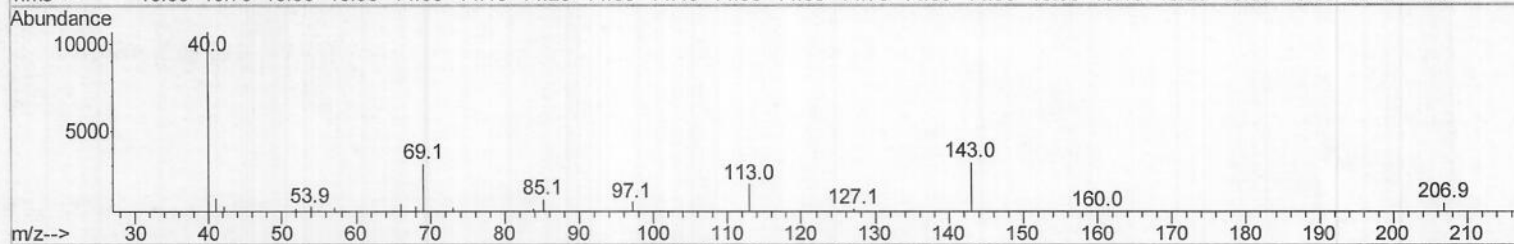
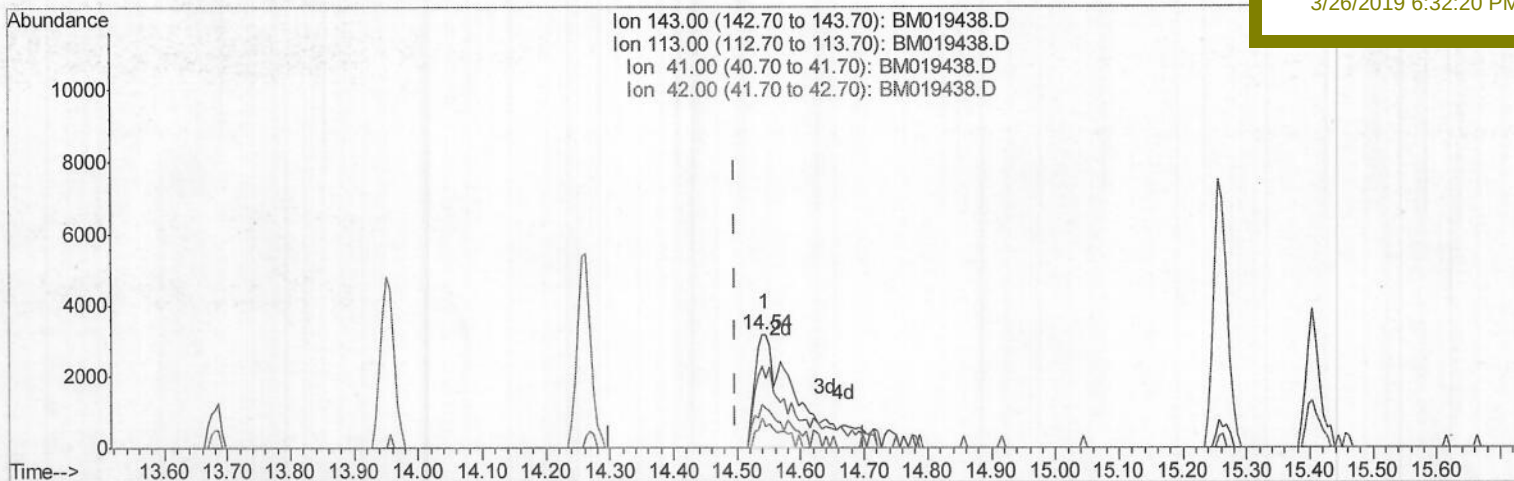
ClientSampleId :

C09D1

Manual Integrations
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Sohil

3/26/2019 6:32:20 PM



(52) 4-Nitrophenol-d4 (S)

14.545min (+0.047) 3.54ng/ul m

response 15259

Ion Exp% Act%

143.00 100 100

113.00 77.10 60.99#

41.00 32.60 34.76

42.00 22.30 18.07

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Misc :

ALS Vial : 18 Sample Multiplier: 1

Quant Time: Apr 05 07:37:17 2019

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM032219MA.M

Quant Title : SVOA CALIBRATION

QLast Update : Fri Apr 05 07:28:31 2019

Response via : Initial Calibration

Instrument :

BNA_M

ClientSampleId :

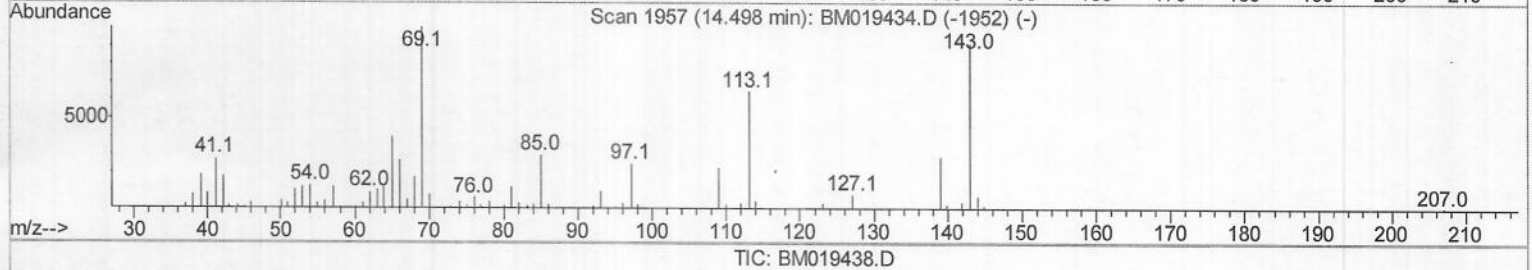
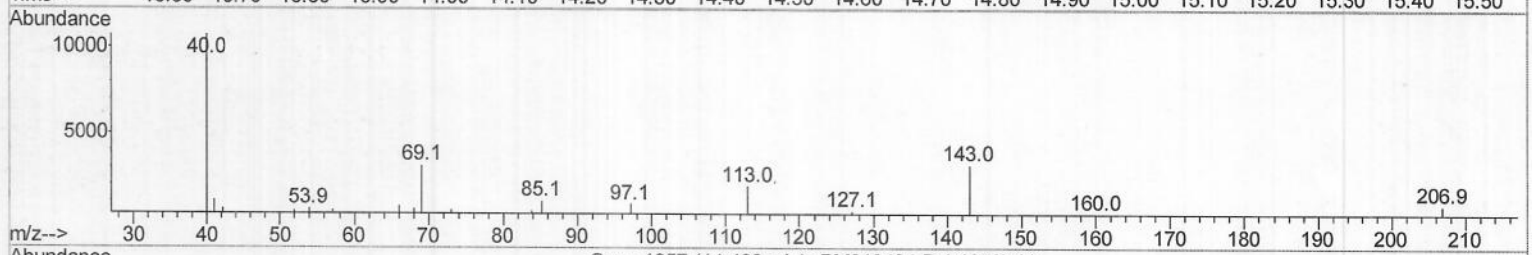
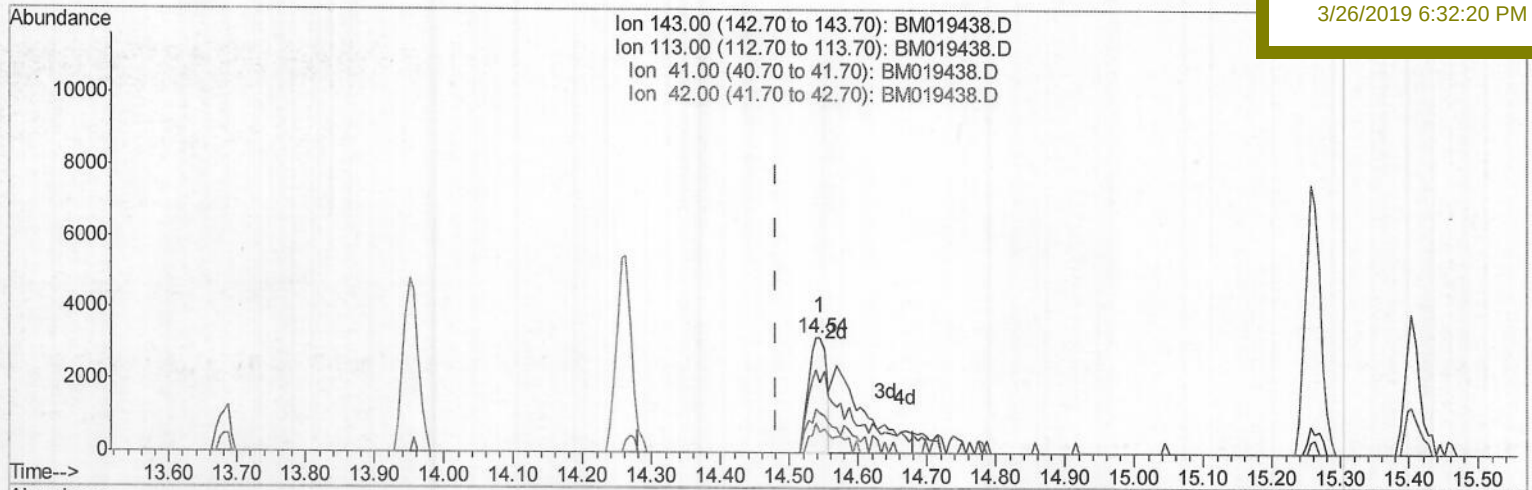
C09D1

Manual Integrations

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Sohil

3/26/2019 6:32:20 PM



(52) 4-Nitrophenol-d4 (S)

14.545min (+0.065) 1.35ng/ul

response 5820

Ion	Exp%	Act%
143.00	100	100
113.00	77.10	60.99#
41.00	32.60	34.76
42.00	22.30	18.07

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Misc :

ALS Vial : 18 Sample Multiplier: 1

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Quant Title : SVOA CALIBRATION

QLast Update : Mon Mar 25 03:11:42 2019

Response via : Initial Calibration

Instrument :

BNA_M

Client Sampled :

C09D1

Manual Integrations

APPROVED

Sohil

3/26/2019 6:32:20 PM

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.60	152	139109	20.00	ng/ul	0.00
18) Naphthalene-d8	10.39	136	589194	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.26	164	341263	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.02	188	753846	20.00	ng/ul	0.00
78) Chrysene-d12	21.23	240	789450	20.00	ng/ul	0.00
86) Perylene-d12	23.44	264	862473	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.15	96	11435	3.22	ng/uL	0.00
5) Phenol-d5	6.79	99	62320	5.03	ng/ul	0.00
7) Bis-(2-Chloroethyl) ether-d	6.96	67	157520	18.99	ng/ul	0.00
9) 2-Chlorophenol-d4	7.14	132	174653	18.60	ng/ul	0.00
13) 4-Methylphenol-d8	8.32	113	112031	12.05	ng/ul	0.00
19) Nitrobenzene-d5	8.77	128	95146	22.64	ng/ul	0.00
22) 2-Nitrophenol-d4	9.49	143	109005	23.30	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.02	165	189232	21.44	ng/ul	0.00
29) 4-Chloroaniline-d4	0.00	131	0d	0.00	ng/ul	0.00
44) Dimethylphthalate-d6	13.68	166	667253	24.23	ng/ul	0.00
47) Acenaphthylene-d8	13.95	160	841874	24.45	ng/ul	0.00
52) 4-Nitrophenol-d4	14.54	143	15259m	3.54	ng/ul	0.05
58) Fluorene-d10	15.26	176	597605	25.19	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.40	200	89776	18.01	ng/ul	0.00
71) Anthracene-d10	17.12	188	937159	25.91	ng/ul	0.00
79) Pyrene-d10	19.43	212	1022558	28.04	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.30	264	1113069	24.34	ng/ul	0.00

Target Compounds

Target Compounds	R.T.	QIon	Response	Conc	Units	Ovalue
73) 1,2,3,4-Tetrachlorobenzene	13.05	216	66141	6.052	ng/uL	98

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

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03/28/19