

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM122619\
Quantitation Report (QT Reviewed)

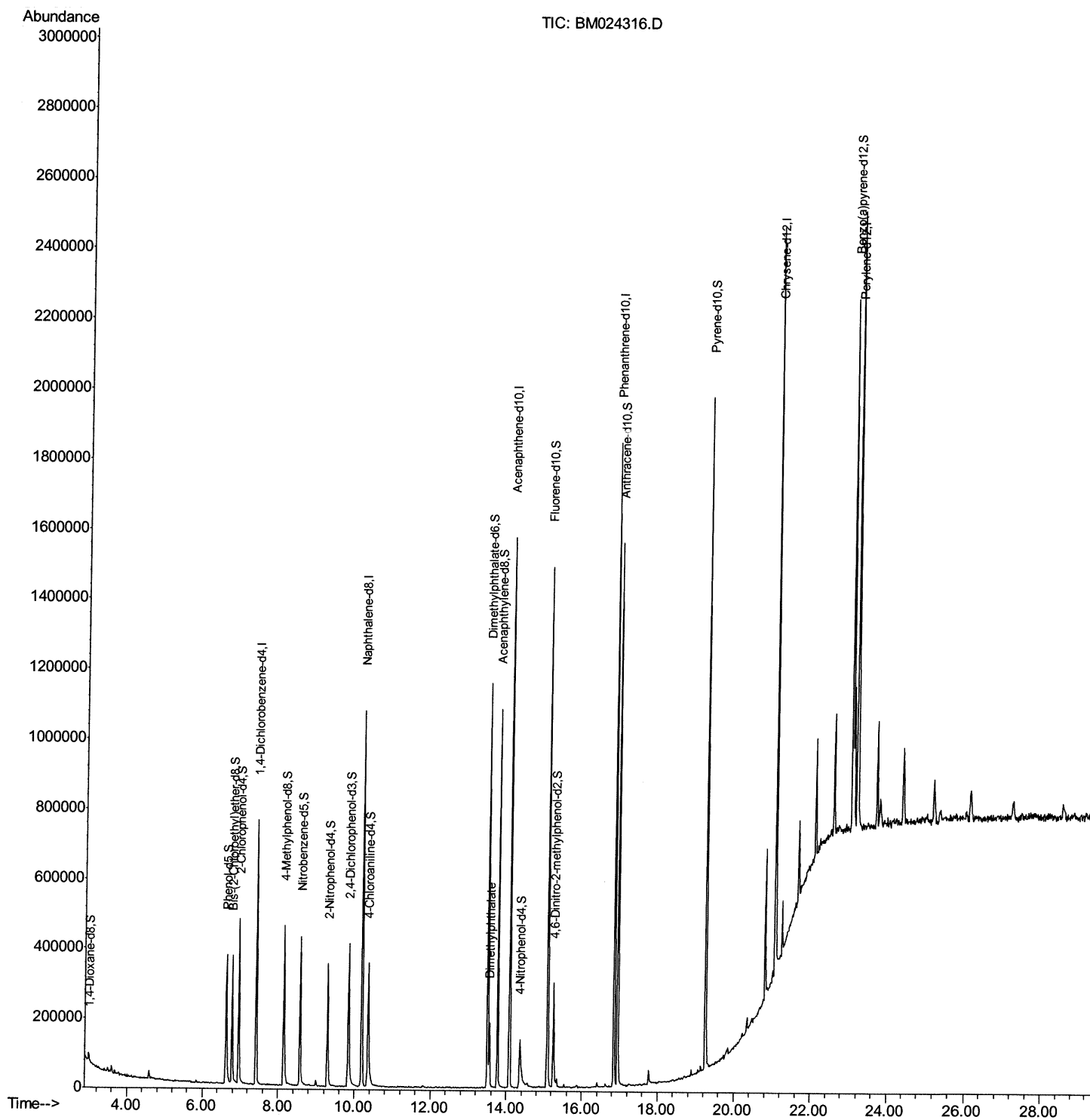
Data File : BM024316.D
Acq On : 27 Dec 2019 01:59
Operator : JU
Sample : K6279-09
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BFEY1

Manual Integrations
APPROVED

mohammad
12/27/2019 3:21:18 PM

Quant Time: Dec 27 06:49:19 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM121719MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Dec 26 03:43:38 2019
Response via : Initial Calibration



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM122619\
Quantitation Report (Qedit)

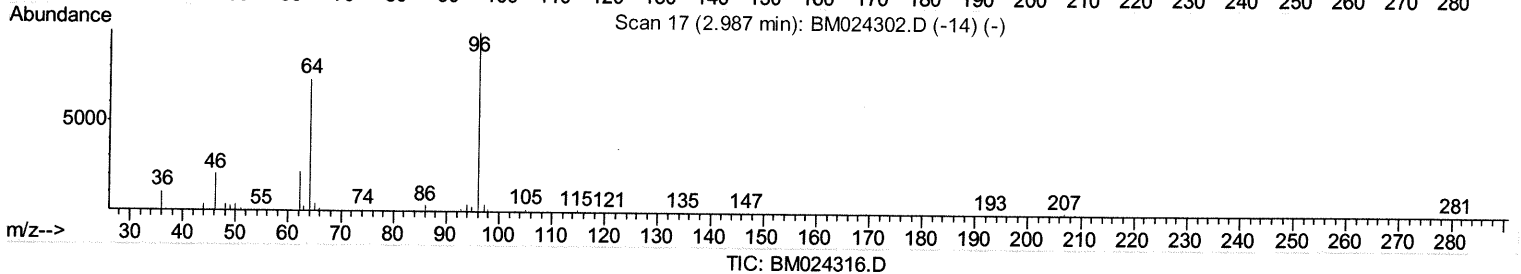
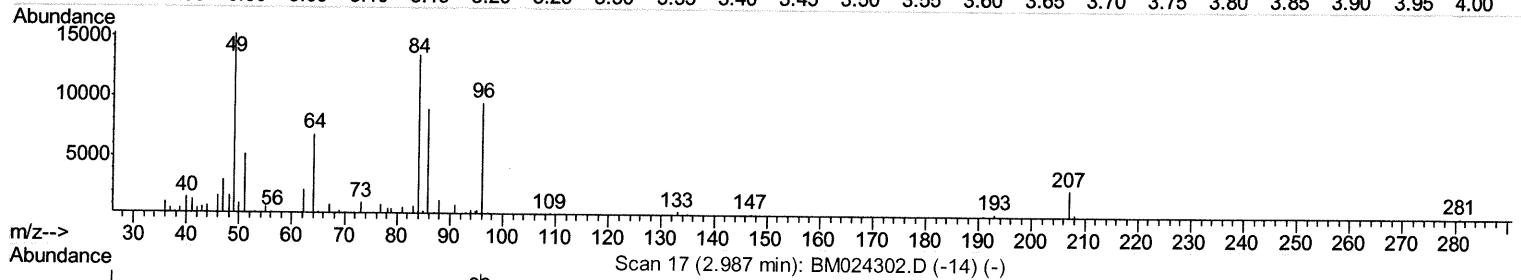
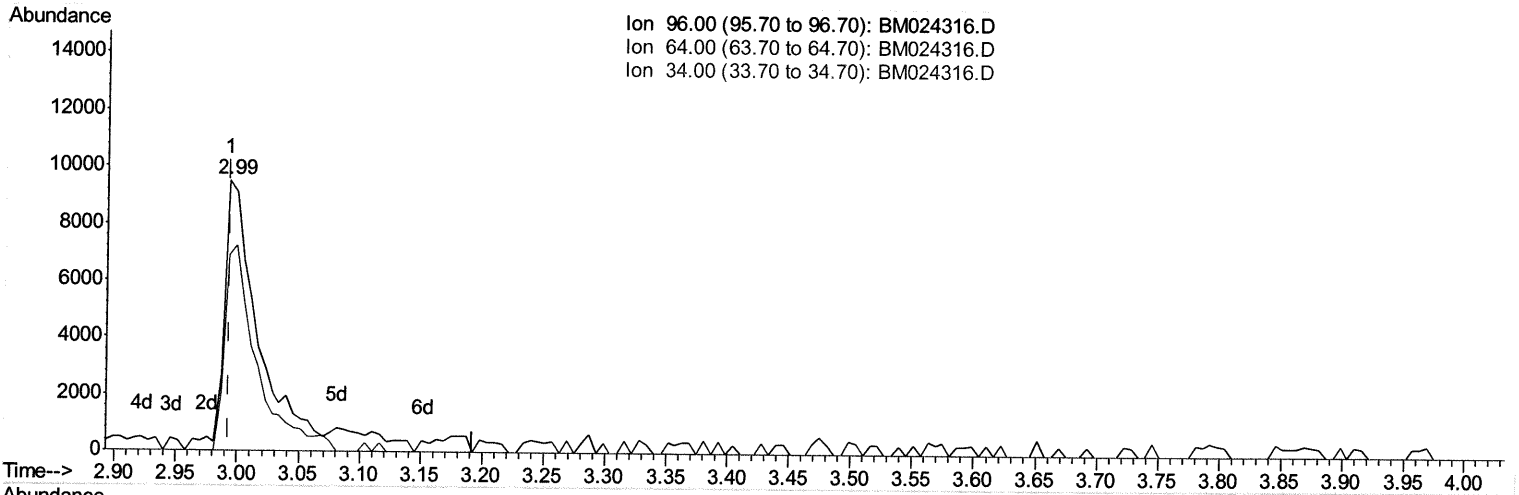
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(3) 1,4-Dioxane-d8 (S)

2.993min (-0.000) 3.13ng/uL

response 14435

Ion	Exp%	Act%
96.00	100	100
64.00	78.60	72.39
34.00	0.00	0.00
0.00	0.00	0.00

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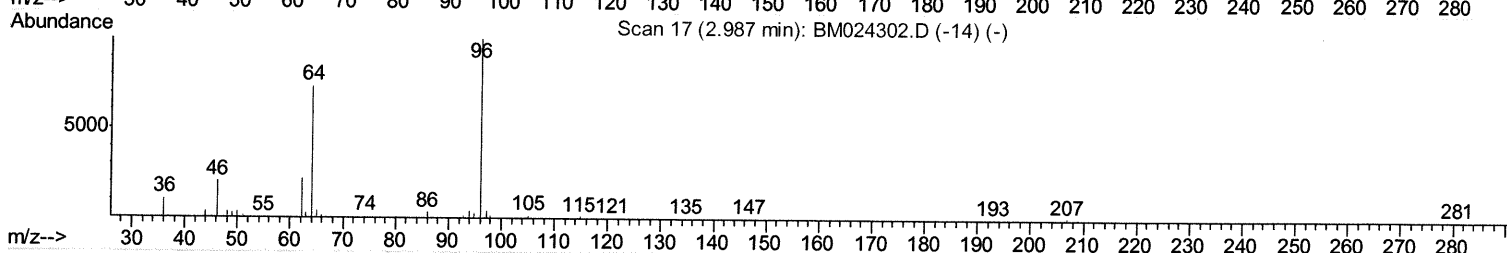
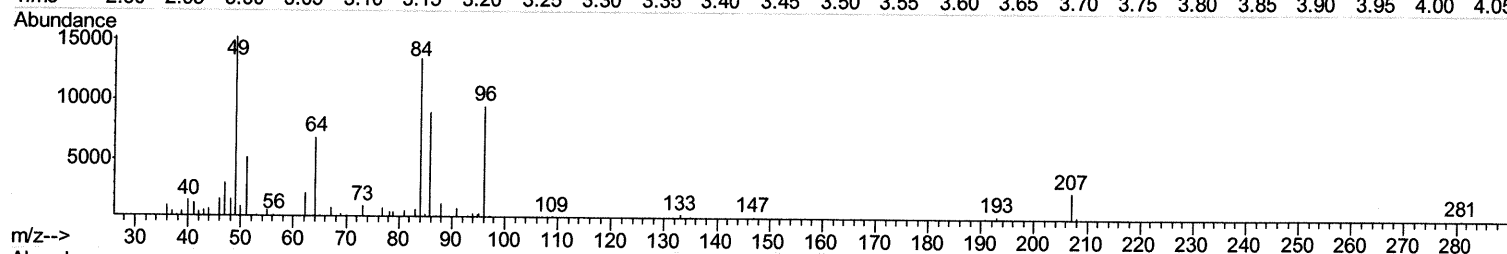
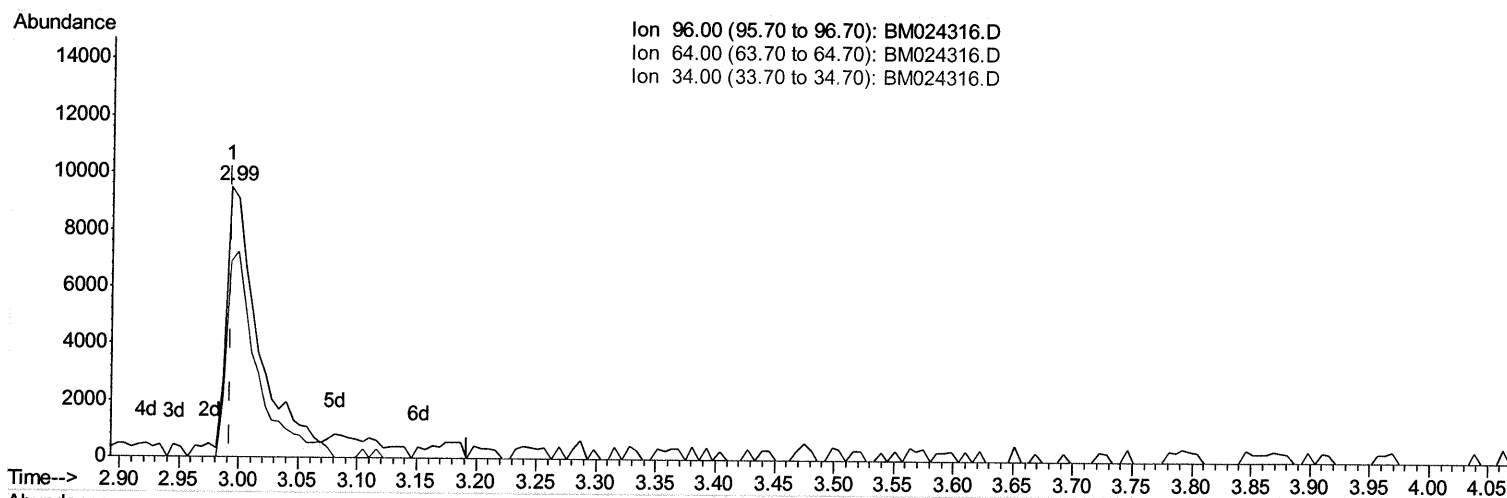
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TIC: BM024316.D

(3) 1,4-Dioxane-d8 (S)

2.993min (-0.000) 3.52ng/uL m *JU 12/27/19*

response 16210

Ion	Exp%	Act%
96.00	100	100
64.00	78.60	72.39
34.00	0.00	0.00
0.00	0.00	0.00

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Quantitation Report (Qedit)

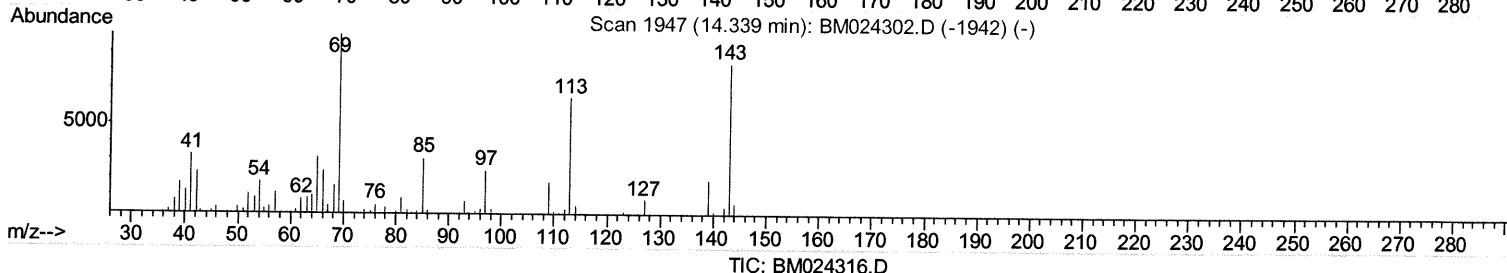
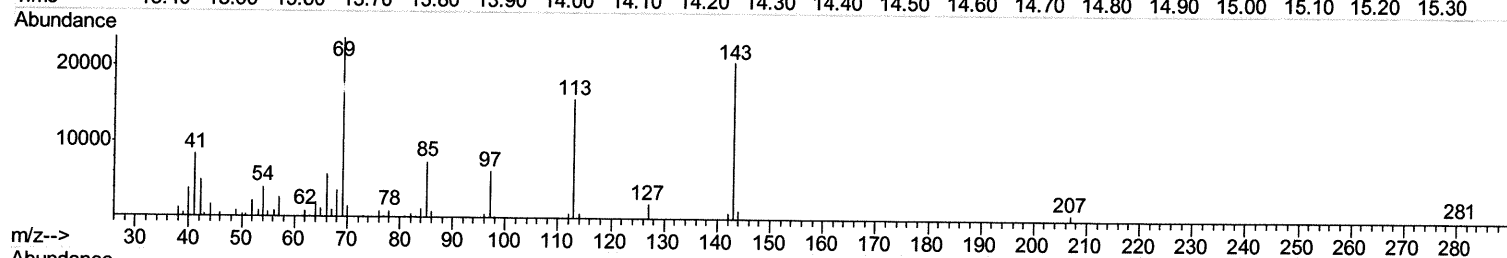
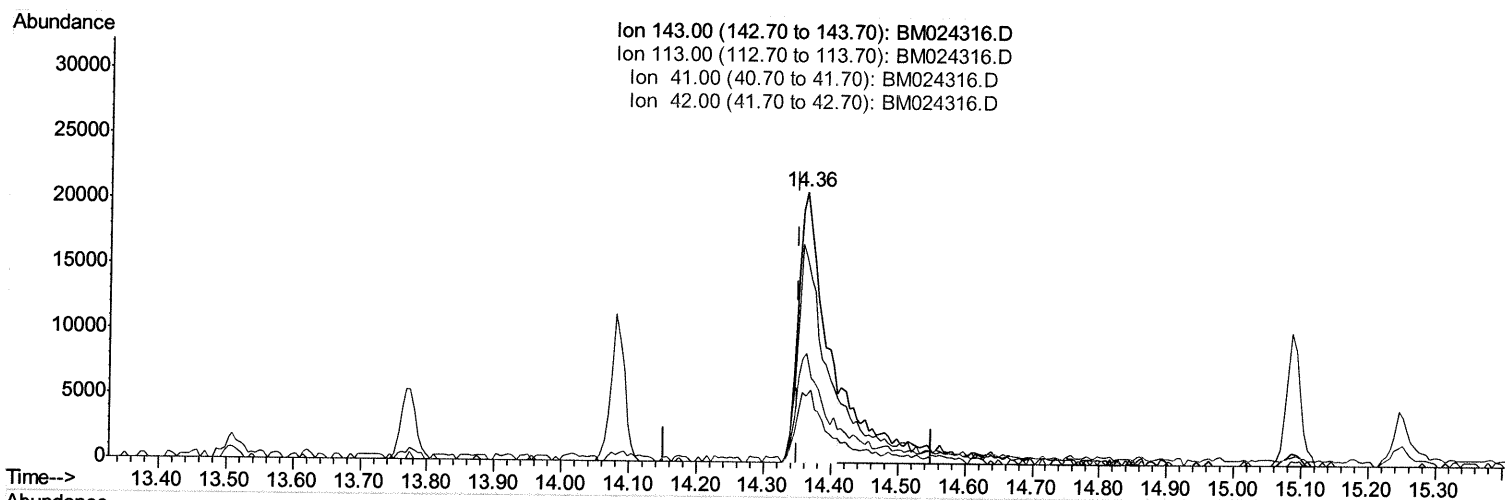
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(52) 4-Nitrophenol-d4 (S)

14.363min (+0.012) 8.08ng/ul

response 53284

Ion	Exp%	Act%
143.00	100	100
113.00	73.60	75.90
41.00	33.90	40.42
42.00	22.80	24.48

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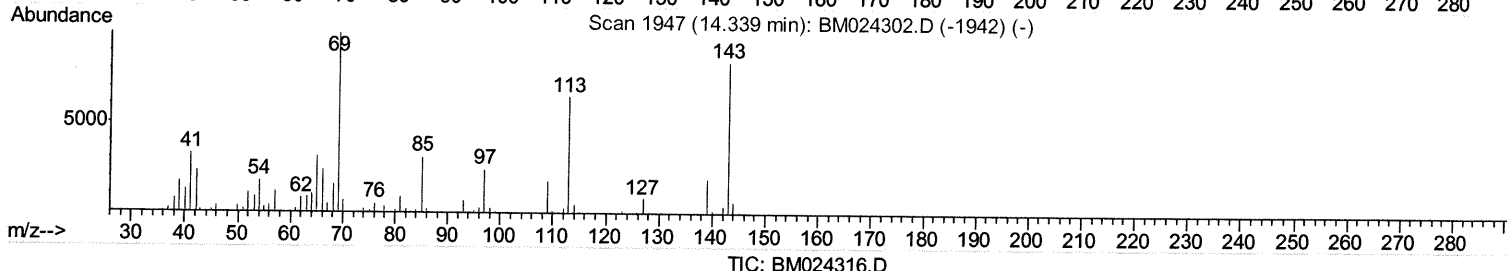
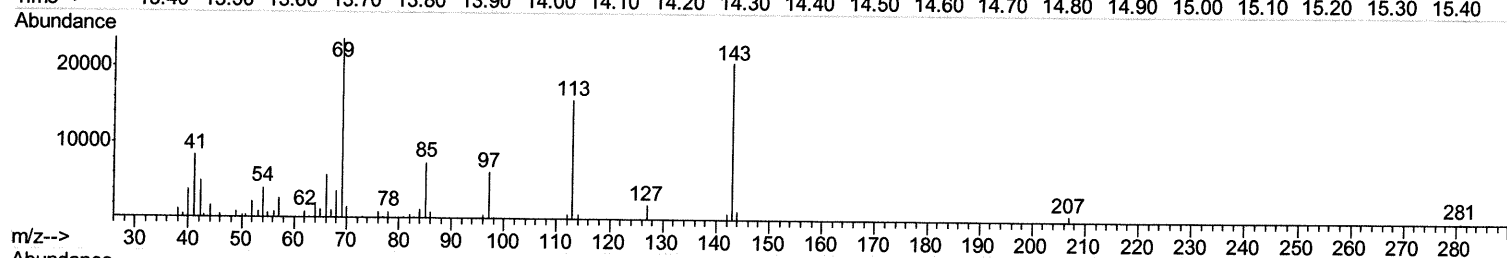
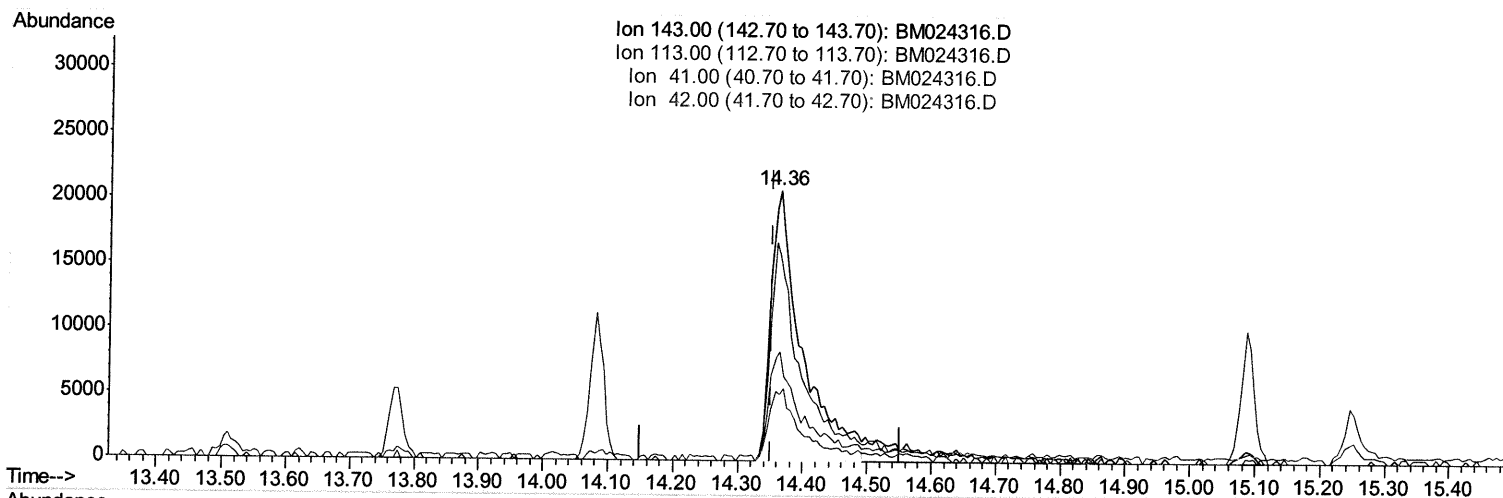
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Instrument :
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Manual Integrations
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(52) 4-Nitrophenol-d4 (S)

14.363min (+0.012) 10.49ng/ul m JU 12/27/19

response 69190

Ion	Exp%	Act%
143.00	100	100
113.00	73.60	75.90
41.00	33.90	40.42
42.00	22.80	24.48

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.42	152	203714	20.00	ng/ul	0.00
18) Naphthalene-d8	10.19	136	851924	20.00	ng/ul	-0.01
36) Acenaphthene-d10	14.09	164	496675	20.00	ng/ul	-0.01
62) Phenanthrene-d10	16.84	188	1008254	20.00	ng/ul	-0.01
78) Chrysene-d12	21.06	240	928296	20.00	ng/ul	-0.01
86) Perylene-d12	23.18	264	1134282	20.00	ng/ul	-0.02

System Monitoring Compounds

3) 1,4-Dioxane-d8	2.99	96	16210m >	3.52	ng/uL >	0.00	5412/27/19
5) Phenol-d5	6.63	99	260069	13.48	ng/ul	0.00	
7) Bis-(2-Chloroethyl) ether-d	6.77	67	204984	17.72	ng/ul	-0.01	
9) 2-Chlorophenol-d4	6.96	132	213970	15.55	ng/ul	0.00	
13) 4-Methylphenol-d8	8.15	113	197984	12.56	ng/ul	0.00	
19) Nitrobenzene-d5	8.58	128	108663	17.75	ng/ul	0.00	
22) 2-Nitrophenol-d4	9.29	143	114007	16.94	ng/ul	0.00	
26) 2,4-Dichlorophenol-d3	9.83	165	182502	13.95	ng/ul	0.00	
29) 4-Chloroaniline-d4	10.35	131	248491	15.64	ng/ul	0.00	
44) Dimethylphthalate-d6	13.51	166	752447	20.49	ng/ul	-0.01	
47) Acenaphthylene-d8	13.77	160	715803	16.26	ng/ul	0.00	
52) 4-Nitrophenol-d4	14.36	143	69190m >	10.49	ng/ul >	0.01	5412/27/19
58) Fluorene-d10	15.09	176	575012	17.86	ng/ul	-0.01	
63) 4,6-Dinitro-2-methylphenol	15.25	200	68102	11.07	ng/ul	0.00	
71) Anthracene-d10	16.94	188	866462	18.89	ng/ul	-0.01	
79) Pyrene-d10	19.25	212	917618	20.75	ng/ul	-0.01	
90) Benzo(a)pyrene-d12	23.05	264	1055896	18.74	ng/ul	-0.02	

Target Compounds					Qvalue
45) Dimethylphthalate	13.56	163	115644	3.030 ng/ul	99

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