

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN022420\
Quantitation Report (QT Reviewed)

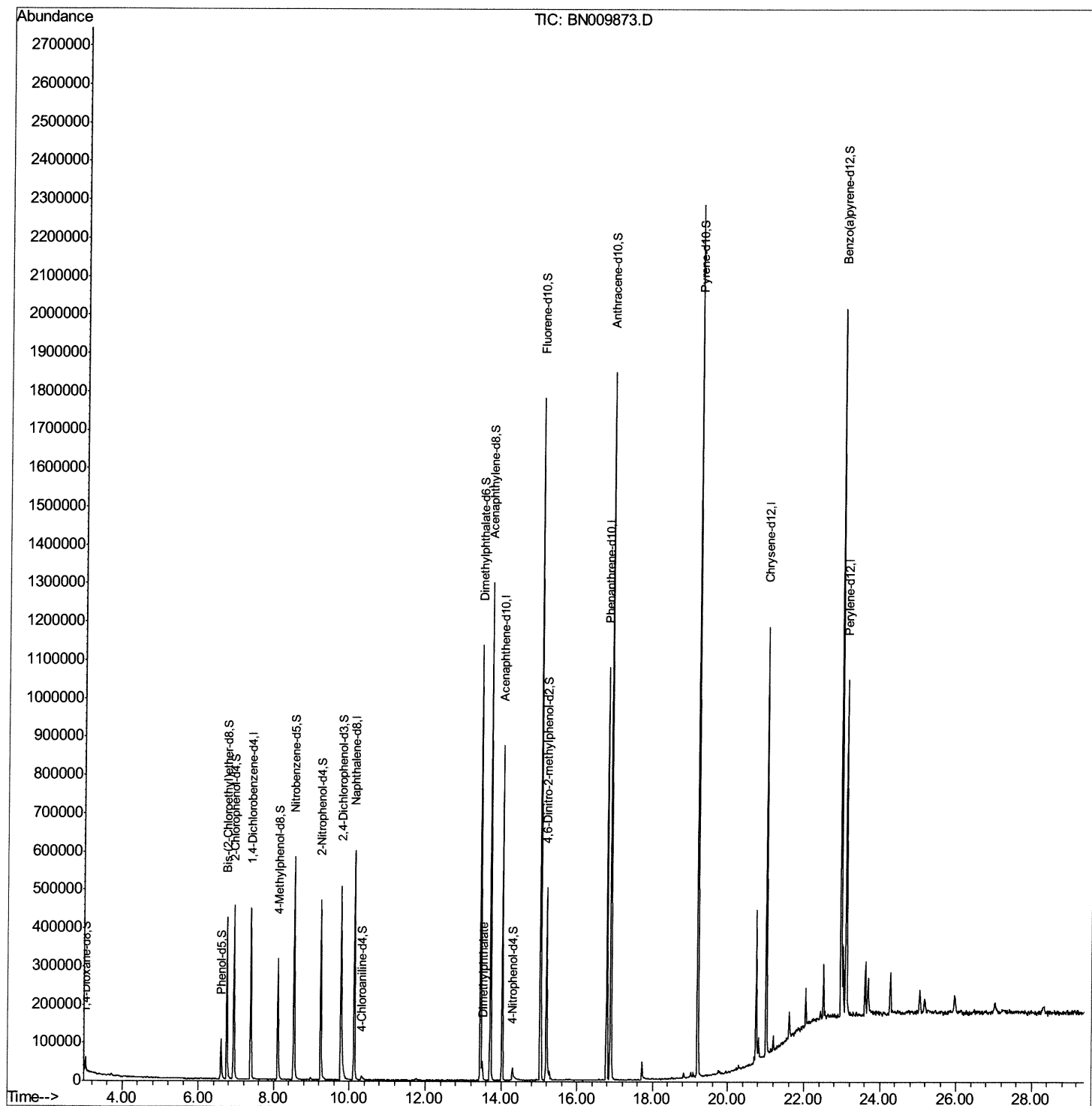
Data File : BN009873.D
Acq On : 24 Feb 2020 21:04
Operator : CG/JU
Sample : L1582-13
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DBDK5

Manual Integrations
APPROVED

mohammad
2/25/2020 3:40:04 PM

Quant Time: Feb 25 02:58:41 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN021220MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Feb 25 02:39:29 2020
Response via : Initial Calibration



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN022420\
Quantitation Report (Qedit)

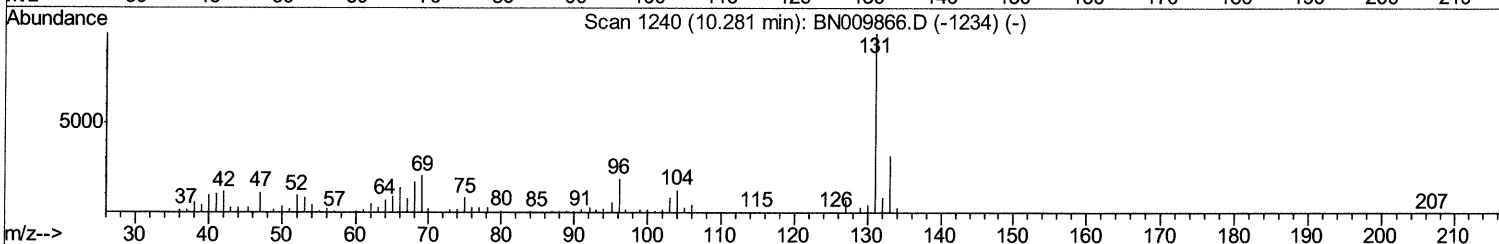
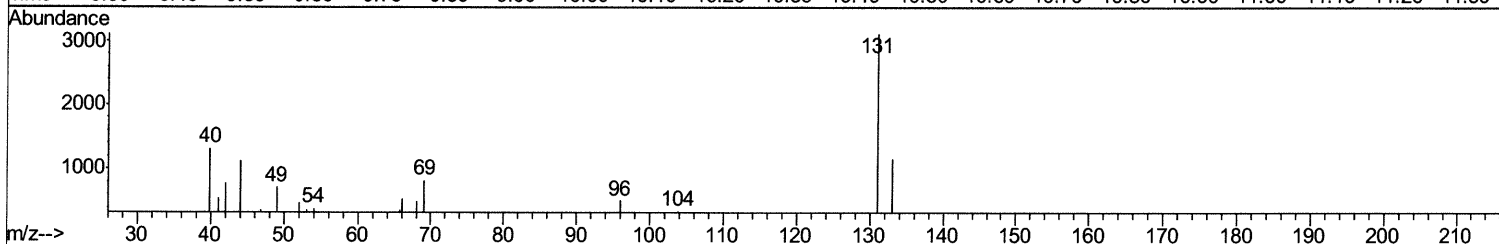
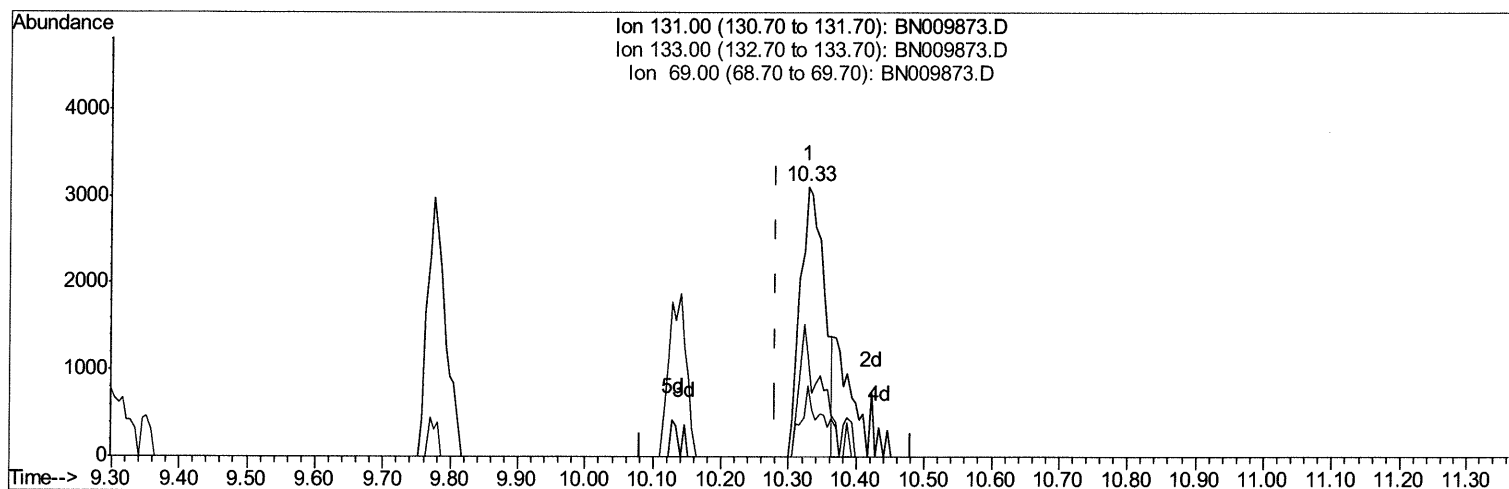
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Quant Title : SVOA CALIBRATION
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TIC: BN009873.D

(29) 4-Chloroaniline-d4 (S)
10.328min (+0.047) 1.06ng/ui
response 7750

Ion	Exp%	Act%
131.00	100	100
133.00	33.00	37.28
69.00	22.10	26.37
0.00	0.00	0.00

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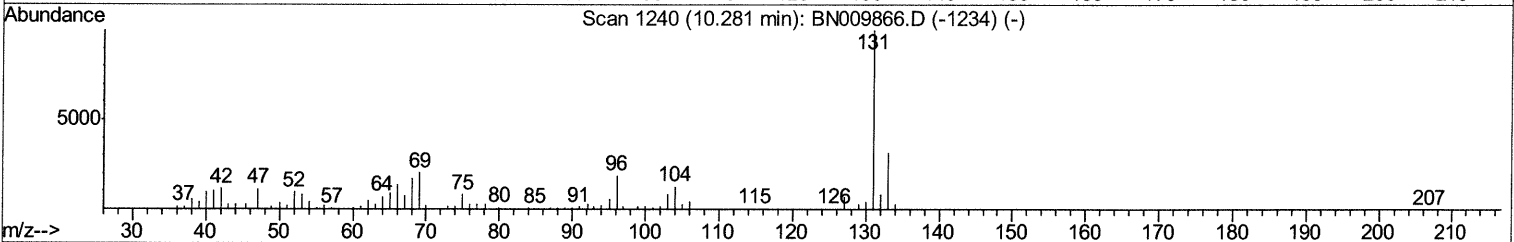
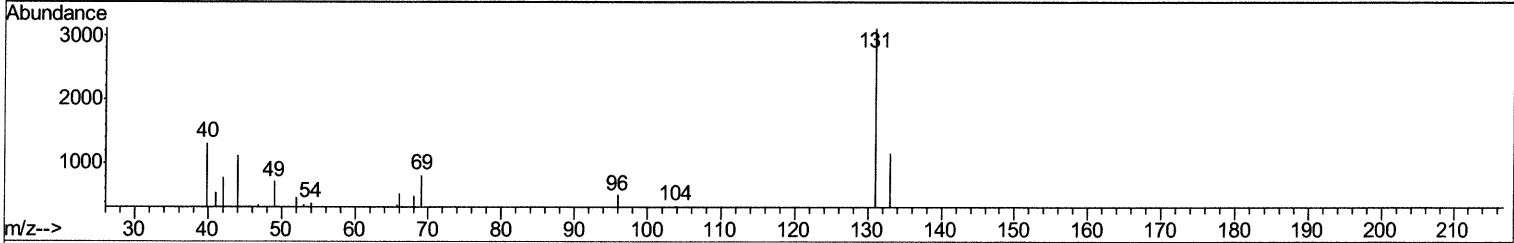
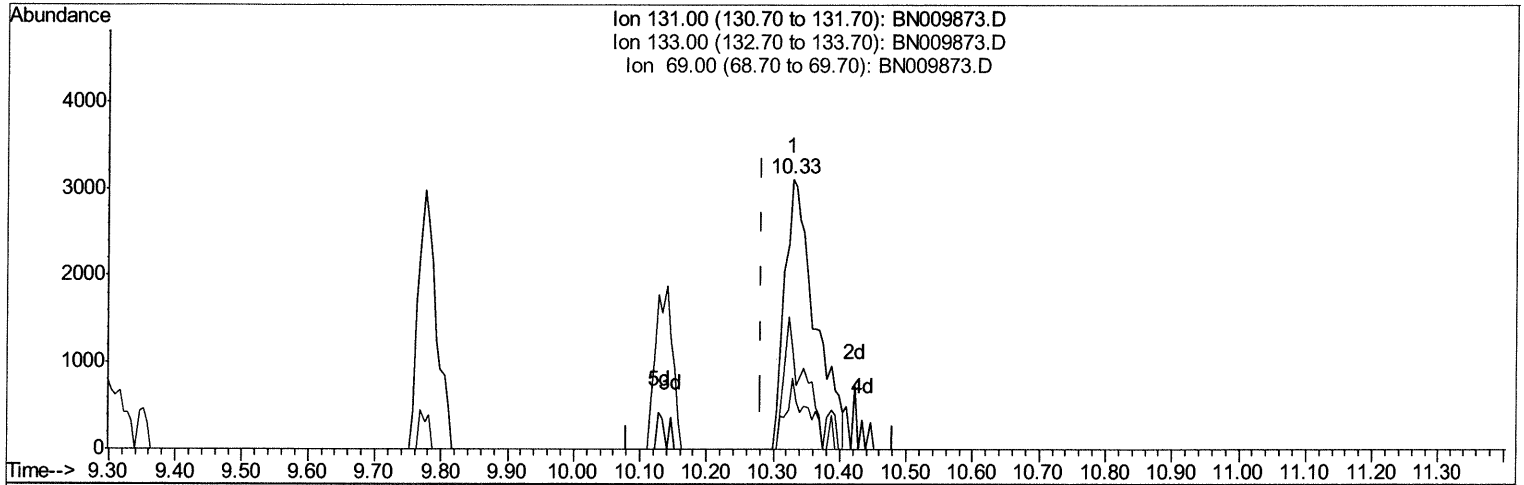
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TIC: BN009873.D

(29) 4-Chloroaniline-d4 (S)

10.328min (+0.047) 1.35ng/ul m *Tu 02/28/20*

response 9900

Ion	Exp%	Act%
131.00	100	100
133.00	33.00	37.28
69.00	22.10	26.37
0.00	0.00	0.00

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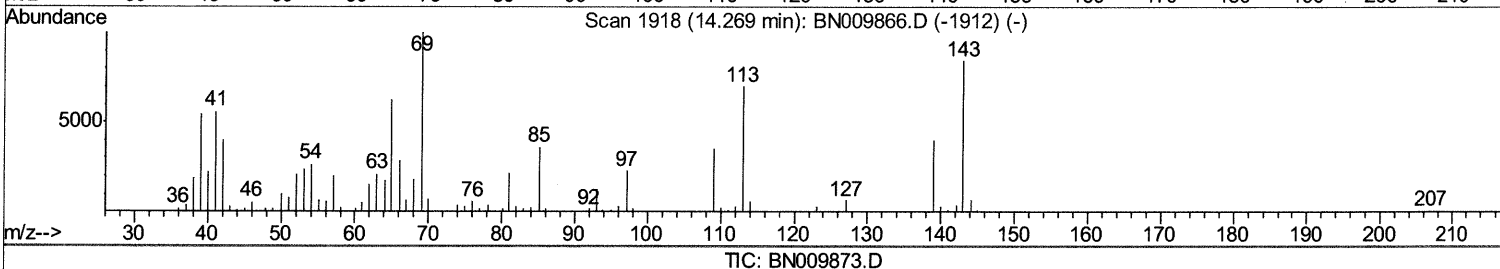
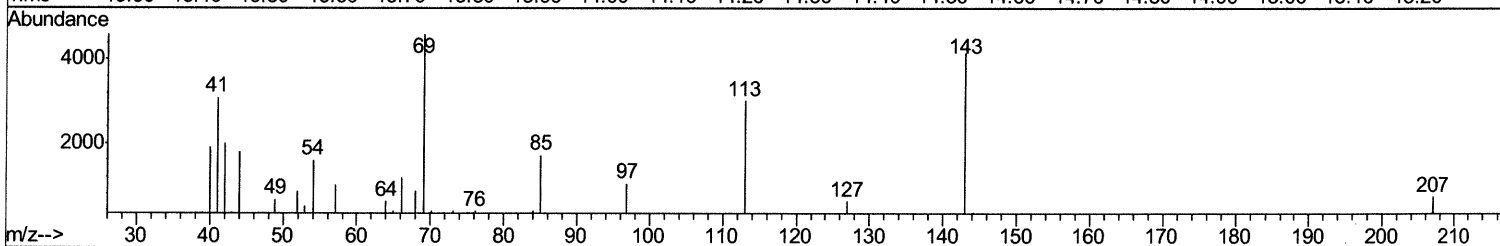
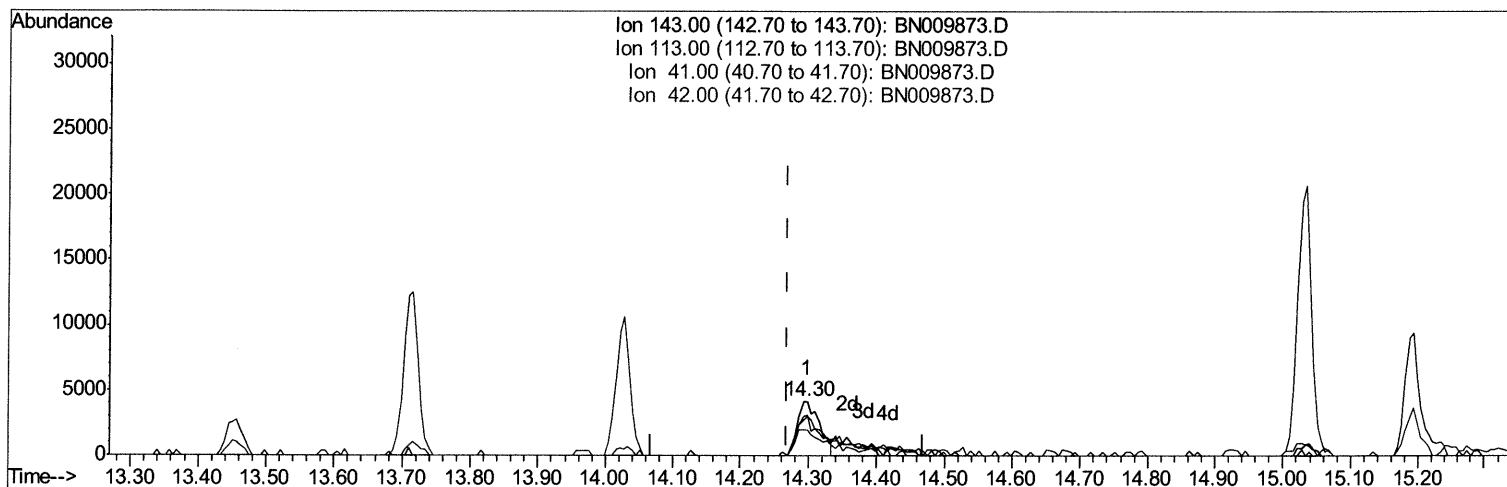
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(52) 4-Nitrophenol-d4 (S)

14.298min (+0.029) 2.59ng/ul

response 9376

Ion	Exp%	Act%
143.00	100	100
113.00	86.40	72.67
41.00	66.00	74.26
42.00	43.20	47.83

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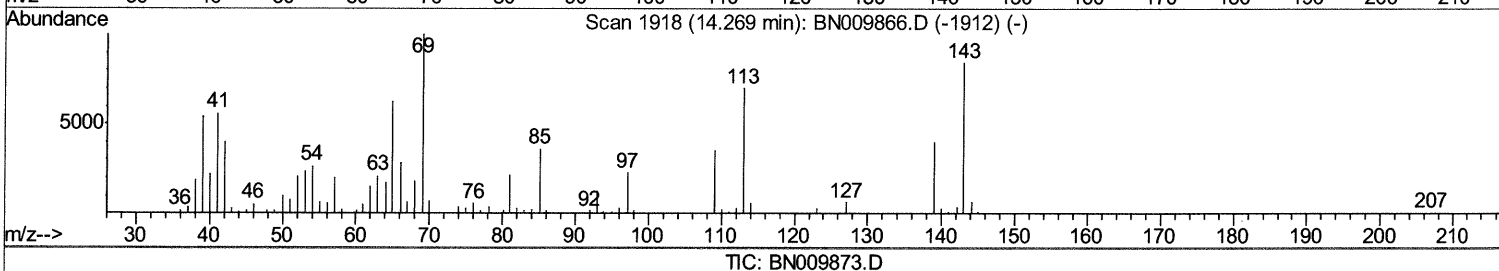
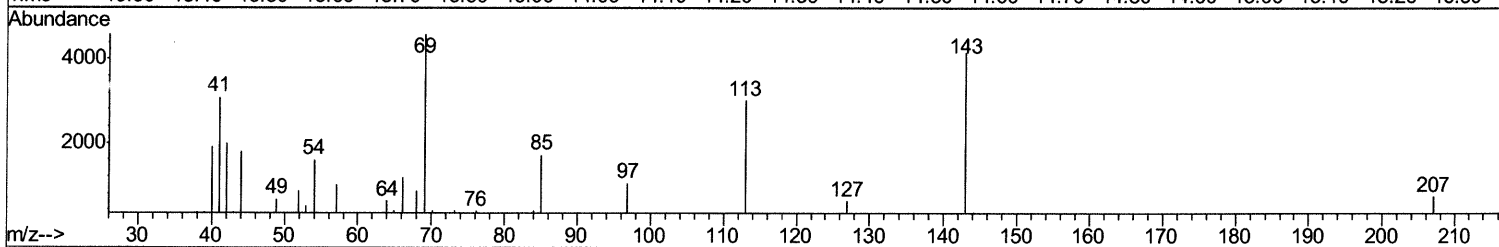
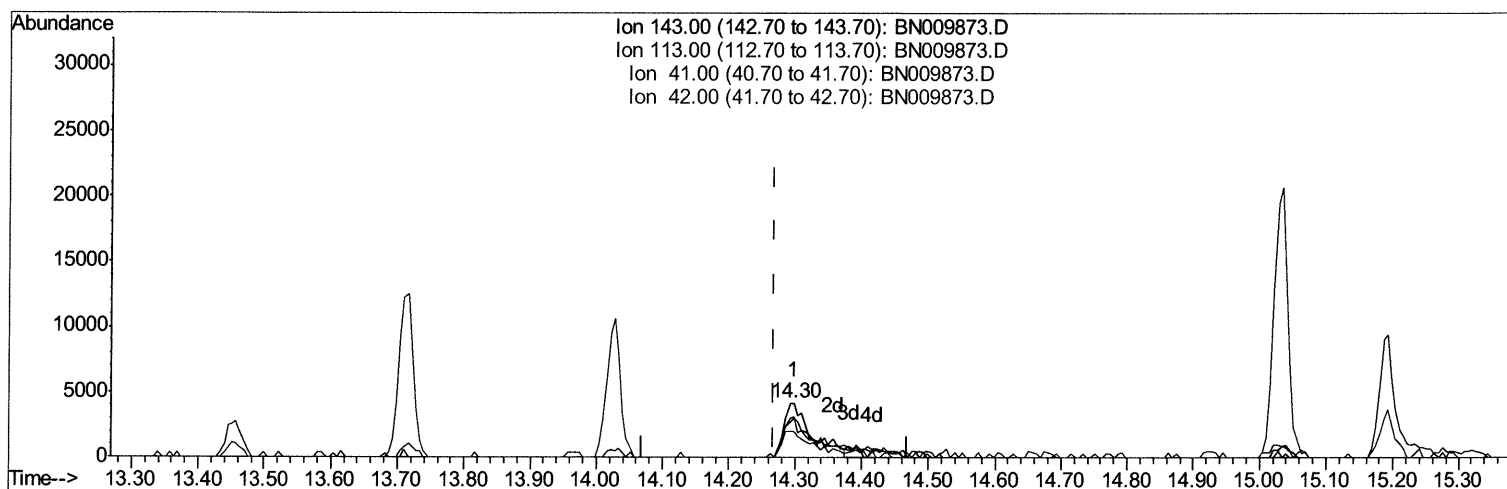
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ALS Vial : 12 Sample Multiplier: 1

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

14.298min (+0.029) 3.28ng/ul m *JU 02/28/20*

response 11842

Ion	Exp%	Act%
143.00	100	100
113.00	86.40	72.67
41.00	66.00	74.26
42.00	43.20	47.83

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.39	152	102014	20.00	ng/ul	0.00
18) Naphthalene-d8	10.13	136	423972	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.03	164	265050	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.79	188	564376	20.00	ng/ul	0.00
78) Chrysene-d12	21.00	240	500260	20.00	ng/ul	0.00
86) Perylene-d12	23.10	264	565603	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.03	96	14544	5.86	ng/uL	0.00
5) Phenol-d5	6.60	99	54081	6.23	ng/ul	0.02
7) Bis-(2-Chloroethyl)ether-d	6.75	67	200113	33.42	ng/ul	0.00
9) 2-Chlorophenol-d4	6.93	132	169932	26.01	ng/ul	0.00
13) 4-Methylphenol-d8	8.10	113	104266	15.11	ng/ul	0.00
19) Nitrobenzene-d5	8.53	128	102889	33.17	ng/ul	0.01
22) 2-Nitrophenol-d4	9.25	143	115487	34.03	ng/ul	0.01
26) 2,4-Dichlorophenol-d3	9.77	165	184837	28.33	ng/ul	0.00
29) 4-Chloroaniline-d4	10.33	131	9900m>	1.35	ng/ul>	0.05 Ju02/28/20
44) Dimethylphthalate-d6	13.46	166	678607	34.28	ng/ul	0.00
47) Acenaphthylene-d8	13.71	160	818509	35.10	ng/ul	0.00
52) 4-Nitrophenol-d4	14.30	143	11842m>	3.28	ng/ul>	0.03 Ju02/28/20
58) Fluorene-d10	15.03	176	624107	36.51	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.19	200	86478	24.67	ng/ul	0.00
71) Anthracene-d10	16.89	188	971299	37.54	ng/ul	0.00
79) Pyrene-d10	19.19	212	1086730	41.56	ng/ul	0.00
90) Benzo(a)pyrene-d12	22.97	264	1084623	38.40	ng/ul	0.00

Target Compounds				Qvalue
45) Dimethylphthalate	13.50	163	24699	1.196 ng/ul 98

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