

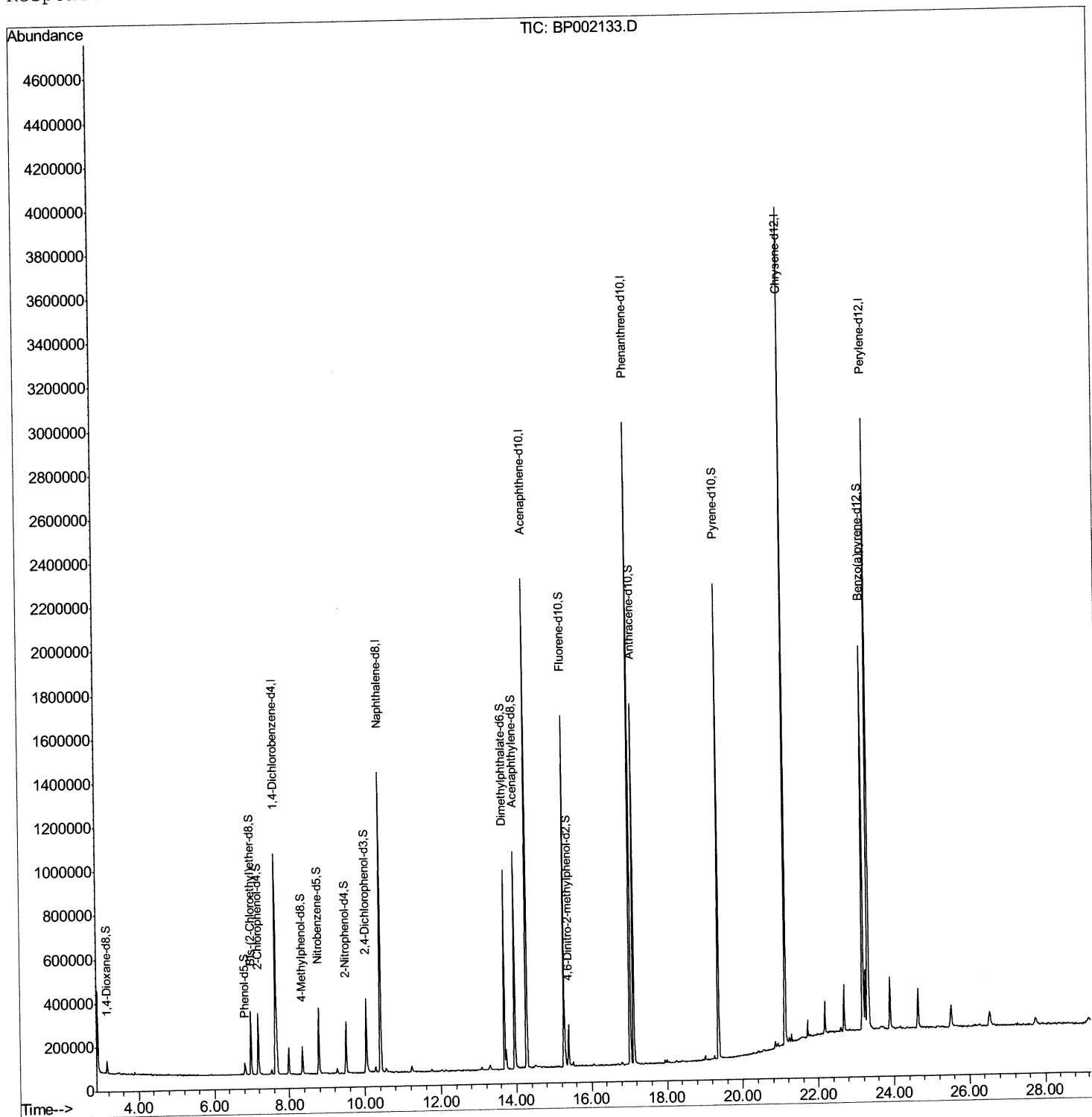
Data Path : Z:\svoasrv\HPCHEM1\BNA P\Data\BP030920\
Data File : BP002133.D
Acq On : 09 Mar 2020 22:20
Operator : CG/JU
Sample : L1834-13
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0DK6

Manual Integrations
APPROVED

mohammad
3/10/2020 2:46:20 PM

Quant Time: Mar 09 23:38:49 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Mar 09 21:43:02 2020
Response via : Initial Calibration



Quantitation Report (Qedit)

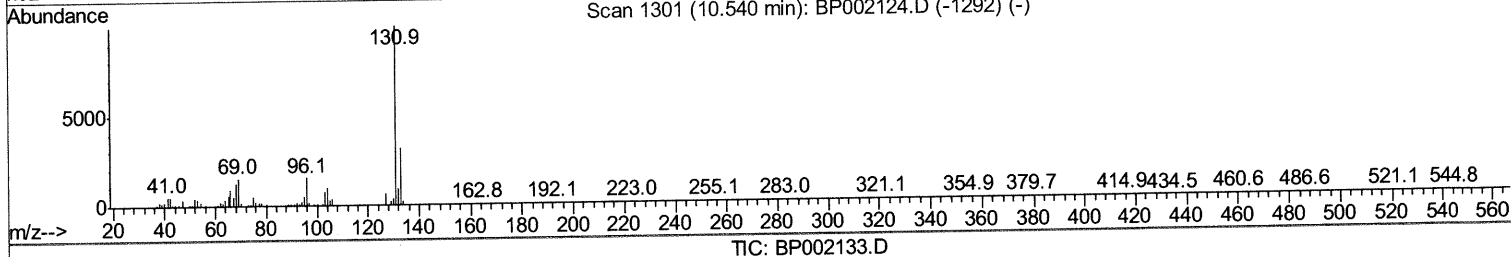
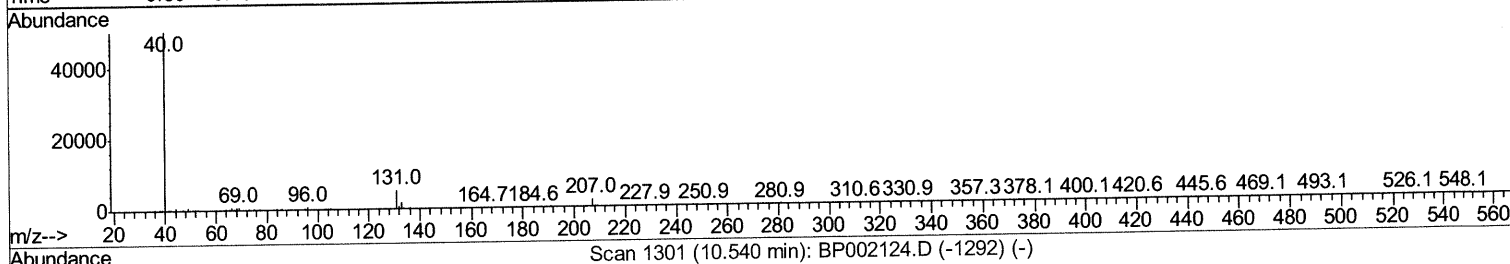
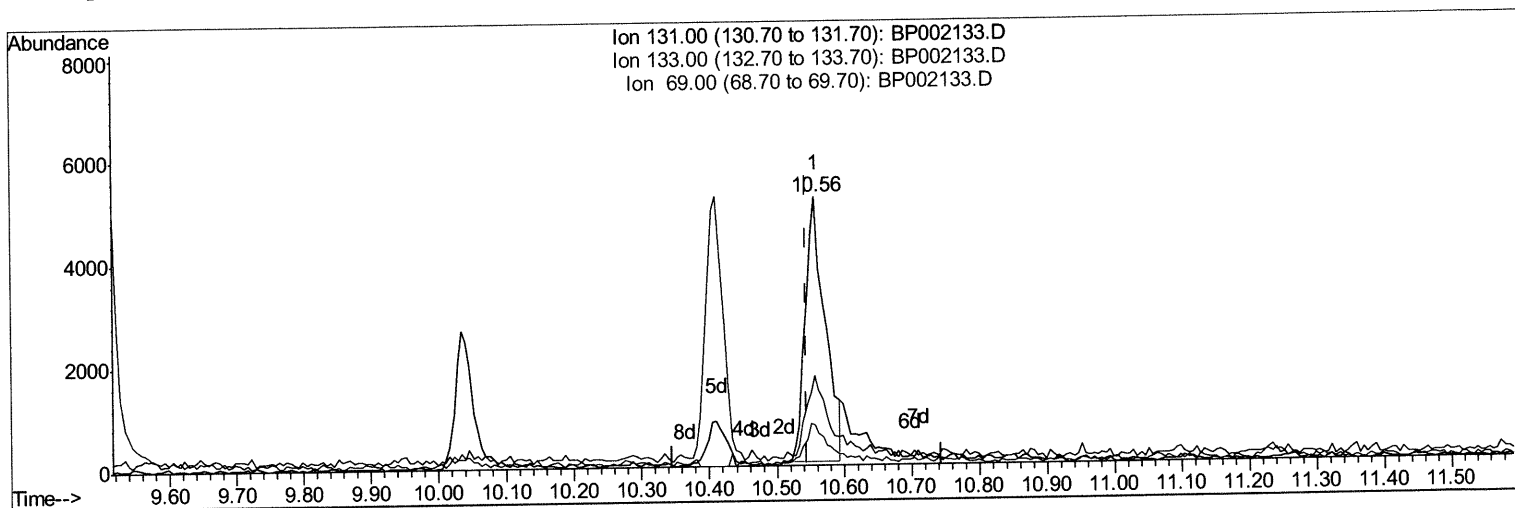
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(29) 4-Chloroaniline-d4 (S)

10.557min (+0.012) 0.48ng/ul

response 10562

Ion	Exp%	Act%
131.00	100	100
133.00	32.40	33.54
69.00	15.50	15.11
0.00	0.00	0.00

Quantitation Report (Qedit)

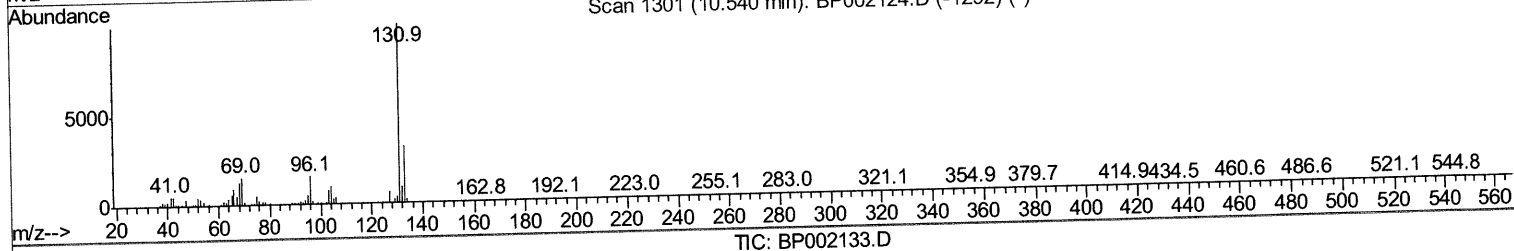
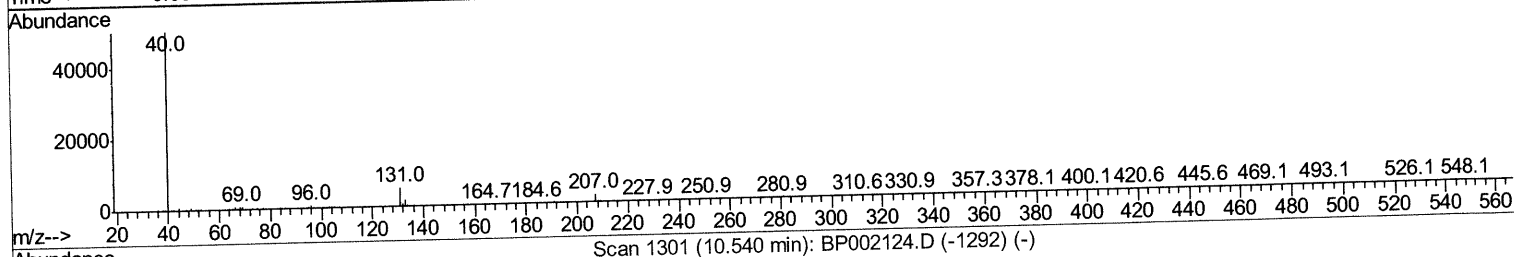
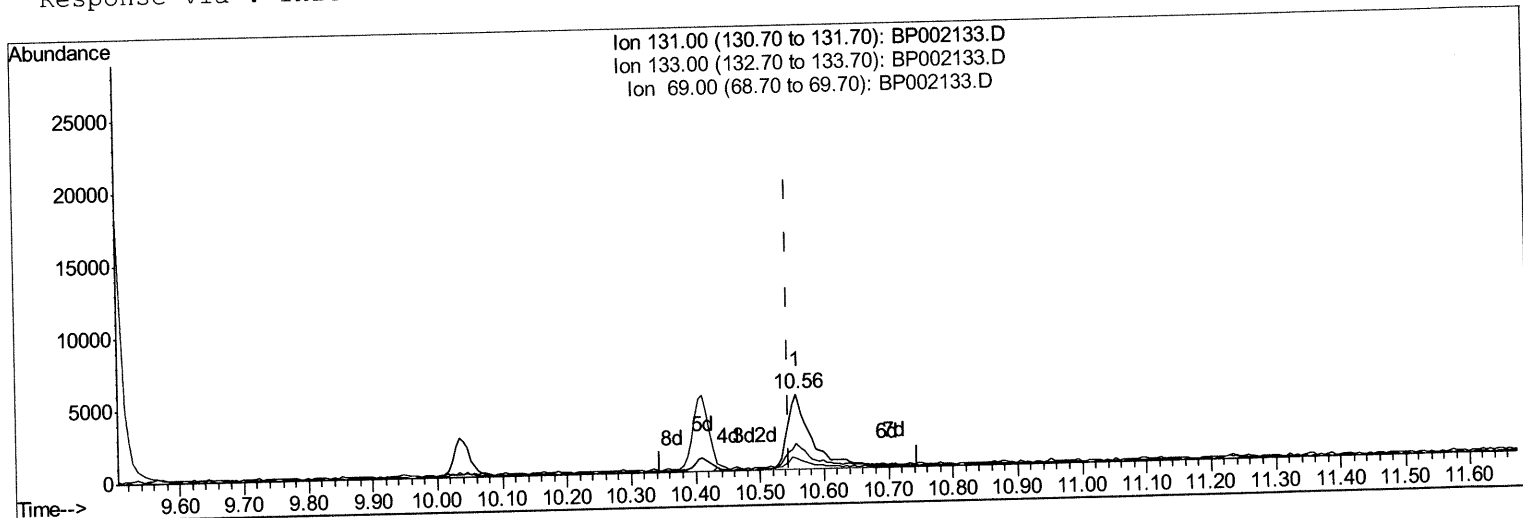
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Manual Integrations
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(29) 4-Chloroaniline-d4 (S)

10.557min (+0.012) 0.61ng/ul m

2003/12/20

response 13458

Ion	Exp%	Act%
131.00	100	100
133.00	32.40	33.54
69.00	15.50	15.11
0.00	0.00	0.00

Quantitation Report (Qedit)

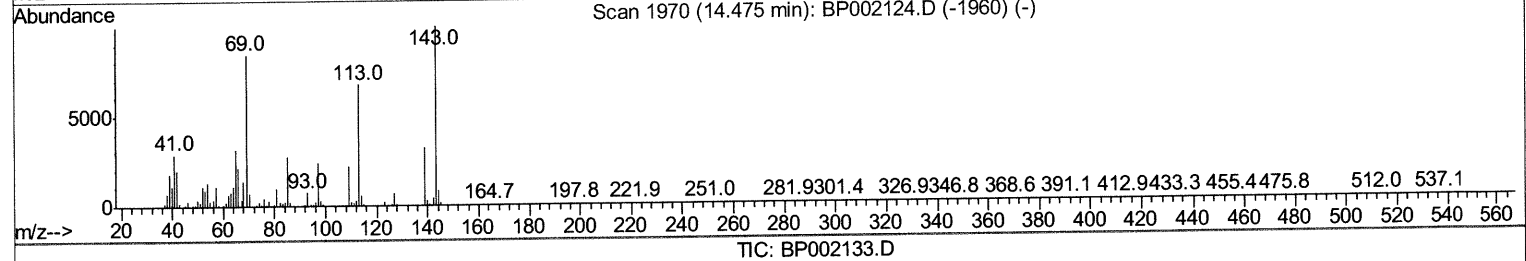
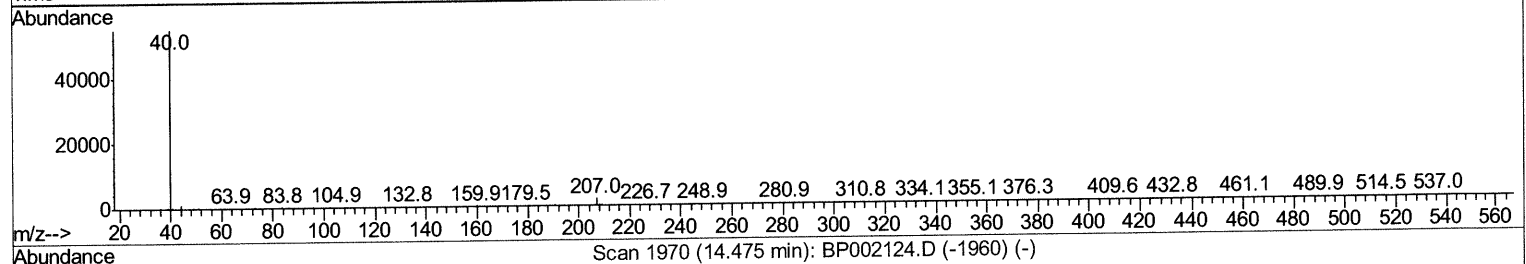
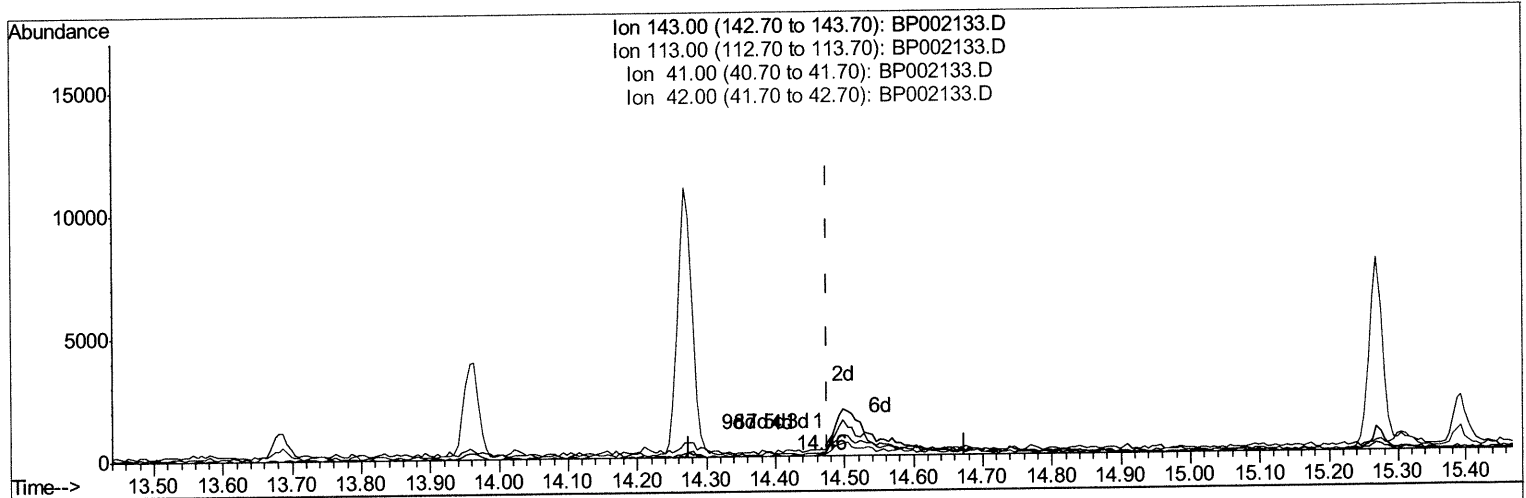
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(52) 4-Nitrophenol-d4 (S)
 14.463min (-0.012) 0.00ng/ul
 response 25

Ion	Exp%	Act%
143.00	100	100
113.00	65.90	21.88#
41.00	29.10	709.38#
42.00	20.10	265.63#

Quantitation Report (Qedit)

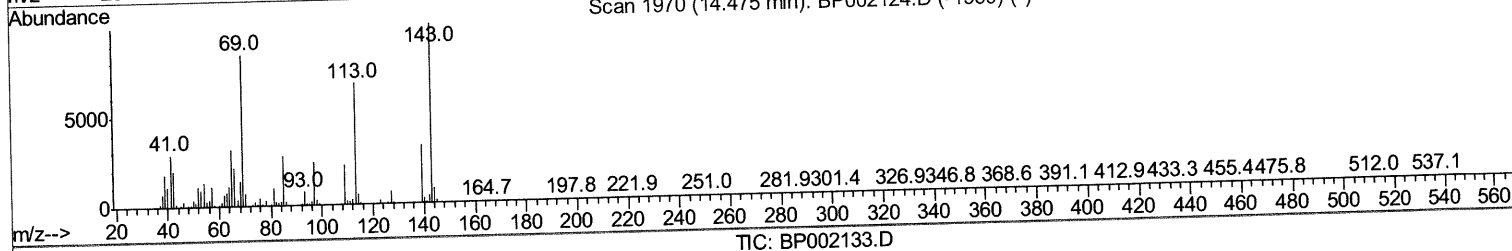
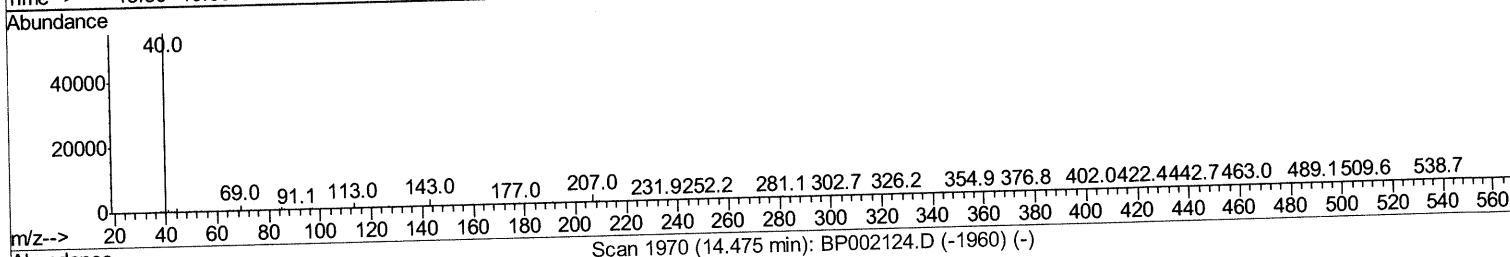
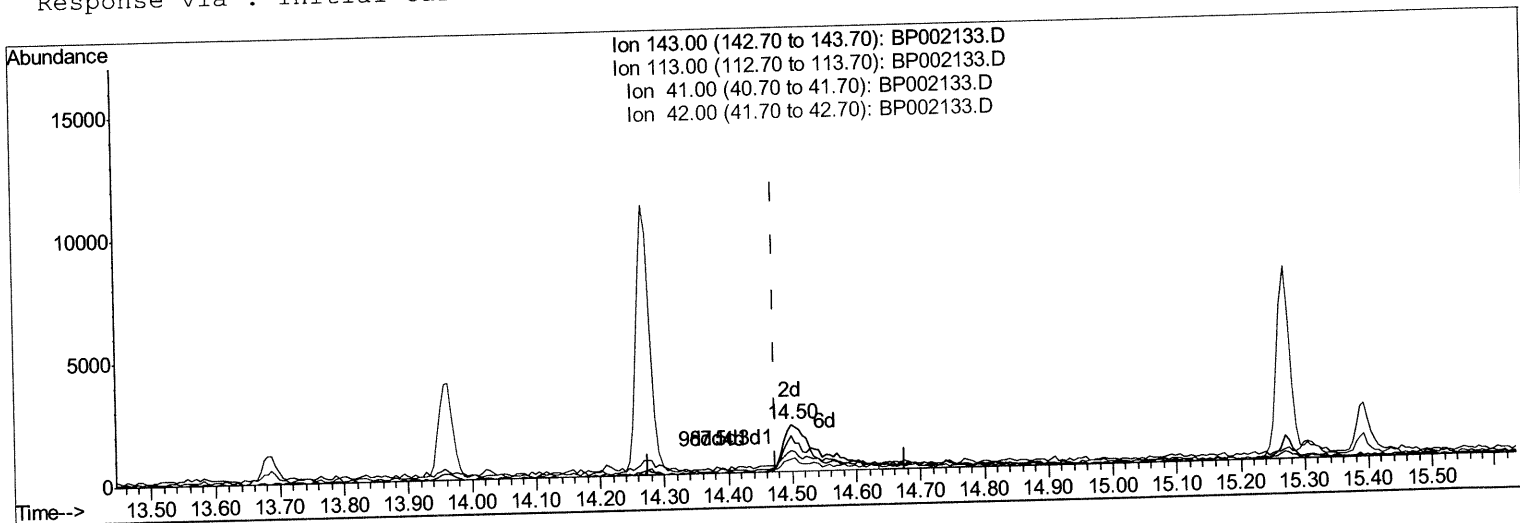
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Instrument :
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Manual Integrations
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(52) 4-Nitrophenol-d4 (S)

14.498min (+0.024) 0.66ng/ul m

Ju 03/12/20

response 7698

Ion	Exp%	Act%
143.00	100	100
113.00	65.90	76.30
41.00	29.10	45.32#
42.00	20.10	25.10#

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	296887	20.00	ng/ul	0.00
18) Naphthalene-d8	10.40	136	1212128	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.27	164	802743	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.02	188	1809916	20.00	ng/ul	0.00
78) Chrysene-d12	21.12	240	1765760	20.00	ng/ul	0.00
86) Perylene-d12	23.32	264	2046115	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.16	96	32996	4.78	ng/uL	0.00
5) Phenol-d5	6.82	99	35787	1.58	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.98	67	125609	10.47	ng/ul	0.00
9) 2-Chlorophenol-d4	7.18	132	140396	7.17	ng/ul	0.00
13) 4-Methylphenol-d8	8.35	113	55279	2.97	ng/ul	0.00
19) Nitrobenzene-d5	8.78	128	84493	9.07	ng/ul	0.00
22) 2-Nitrophenol-d4	9.50	143	86635	8.27	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.03	165	157475	7.80	ng/ul	0.00
29) 4-Chloroaniline-d4	10.56	131	13458m	0.61	ng/ul	0.01
44) Dimethylphthalate-d6	13.69	166	659841	10.86	ng/ul	0.00
47) Acenaphthylene-d8	13.96	160	744193	10.24	ng/ul	0.00
52) 4-Nitrophenol-d4	14.50	143	7698m	0.66	ng/ul	0.02
58) Fluorene-d10	15.27	176	633968	12.06	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.39	200	49616	4.97	ng/ul	0.00
71) Anthracene-d10	17.12	188	974139	11.98	ng/ul	0.00
79) Pyrene-d10	19.37	212	1129827	12.22	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.18	264	1281345	12.24	ng/ul	0.00

Jul 3/12/20

Jul 03/12/20

Target Compounds Ovalue

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