

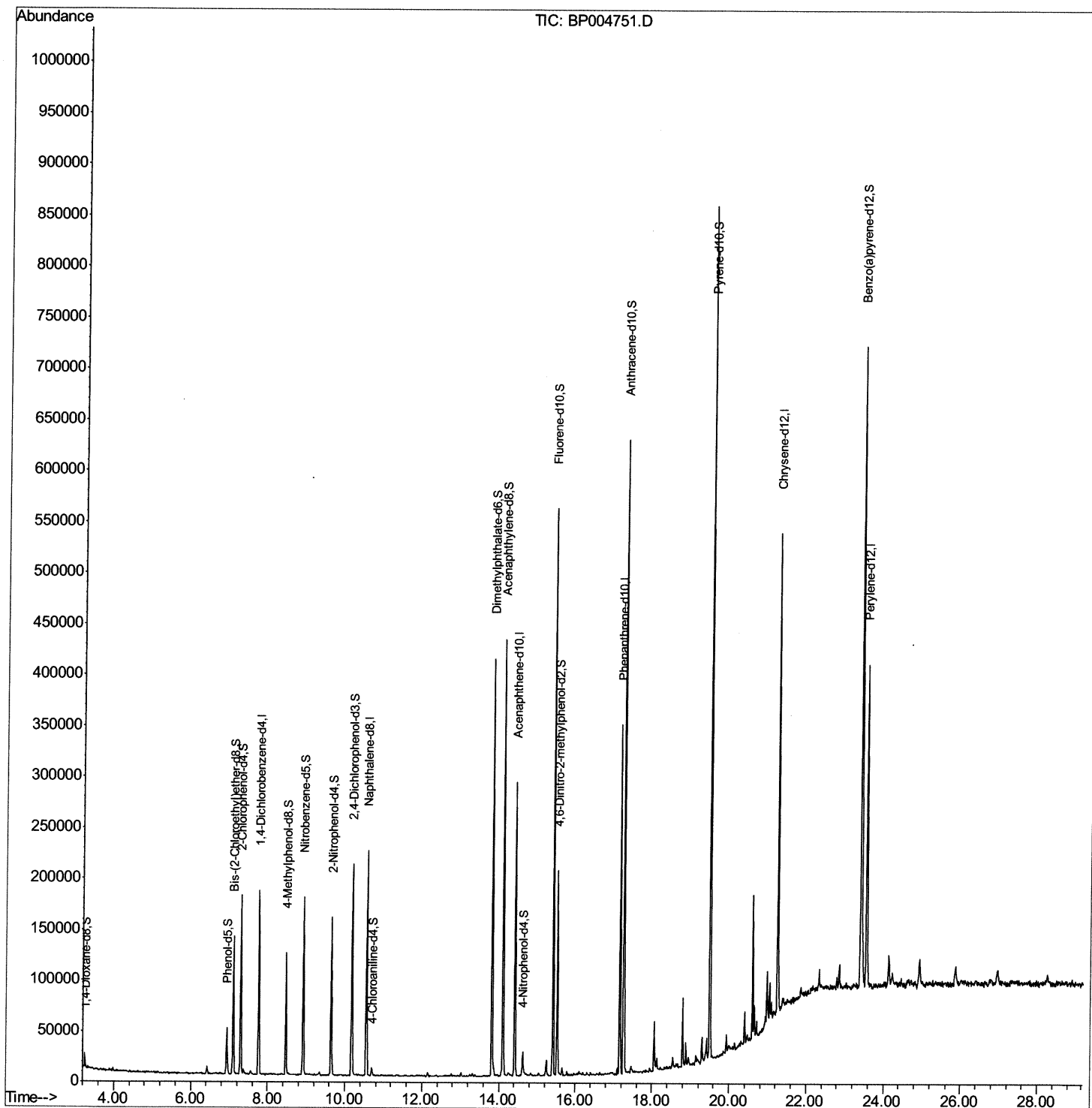
Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP021521\
Data File : BP004751.D
Acq On : 15 Feb 2021 17:51
Operator : CG/JU
Sample : M1349-17
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBGS9

Manual Integrations
APPROVED

mohammad
2/16/2021 10:45:38 AM

Quant Time: Feb 16 00:18:41 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP020821MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Feb 15 12:16:19 2021
Response via : Initial Calibration



Quantitation Report (Qedit)

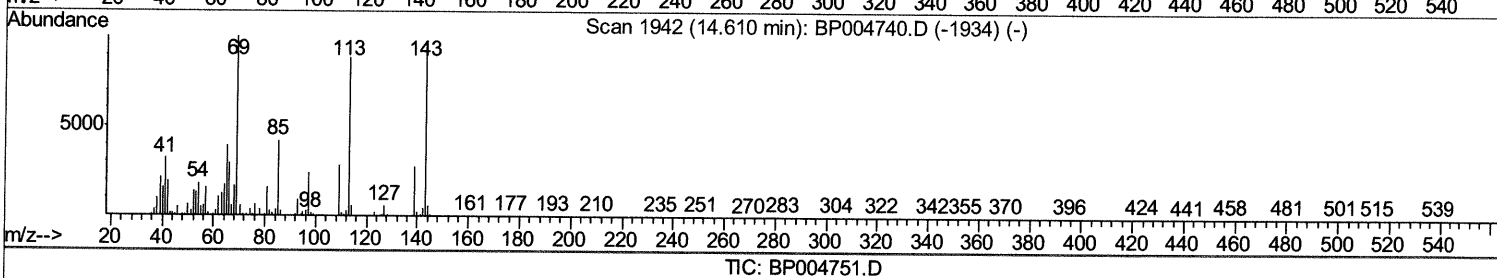
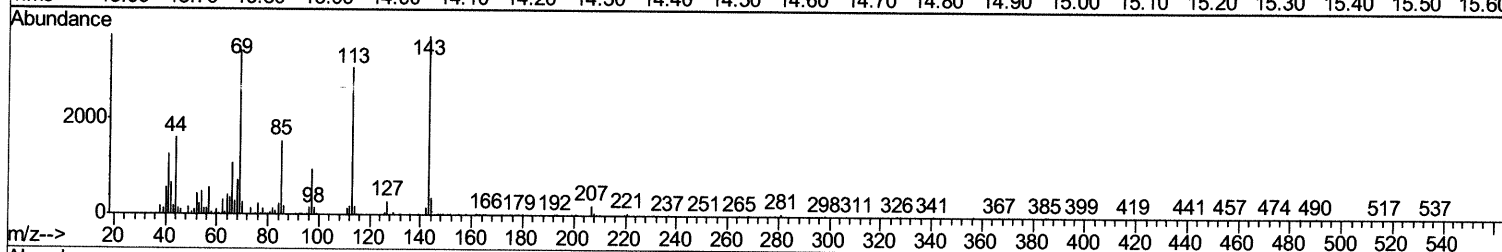
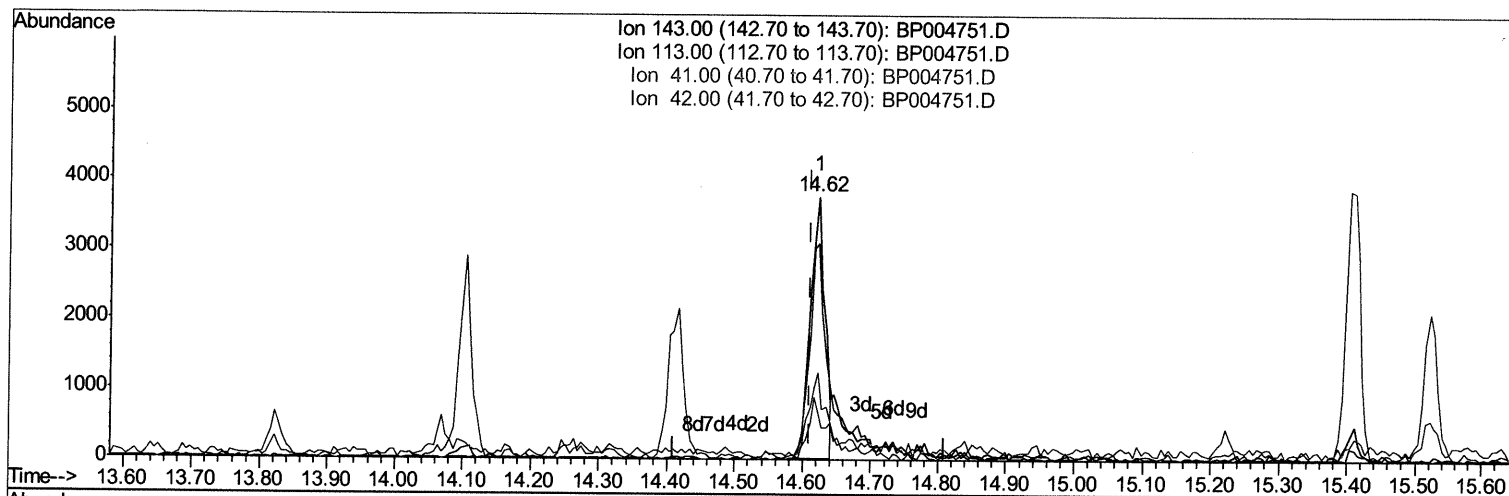
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mohammad
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Quant Time: Feb 16 00:05:19 2021
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Quant Title : SVOA CALIBRATION
QLast Update : Mon Feb 15 12:16:19 2021
Response via : Initial Calibration



(52) 4-Nitrophenol-d4 (S)

14.622min (+0.012) 4.67ng/ul

response 5642

Ion	Exp%	Act%
143.00	100	100
113.00	104.70	81.94#
41.00	37.50	33.30
42.00	24.30	18.14#

Quantitation Report (Qedit)

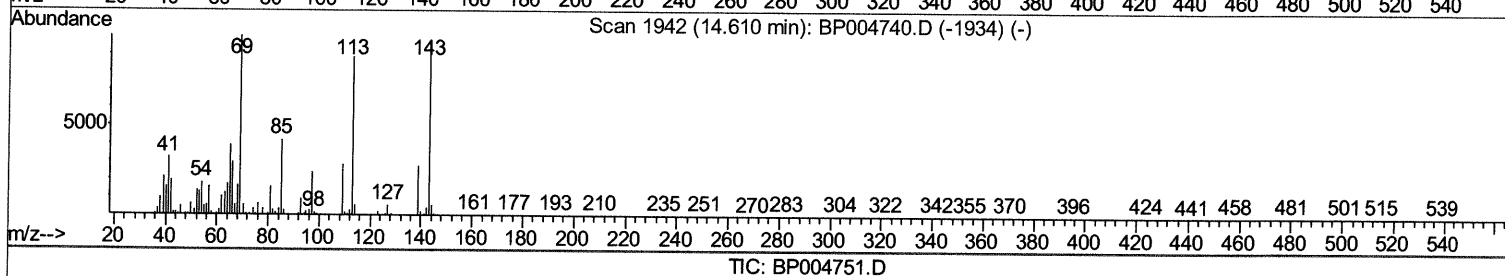
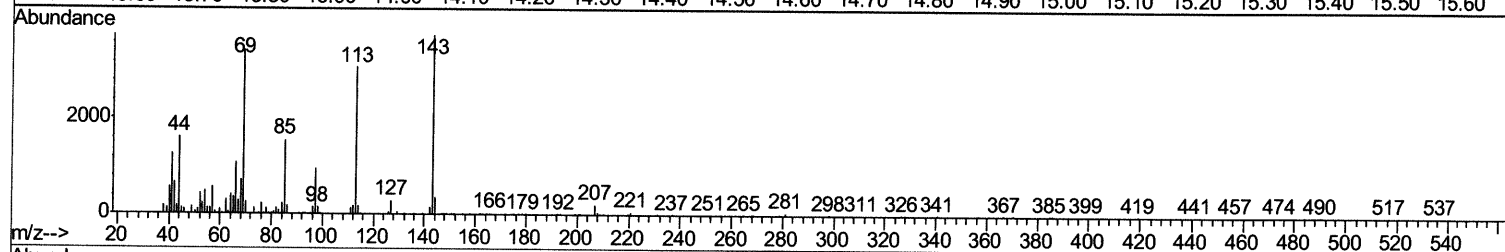
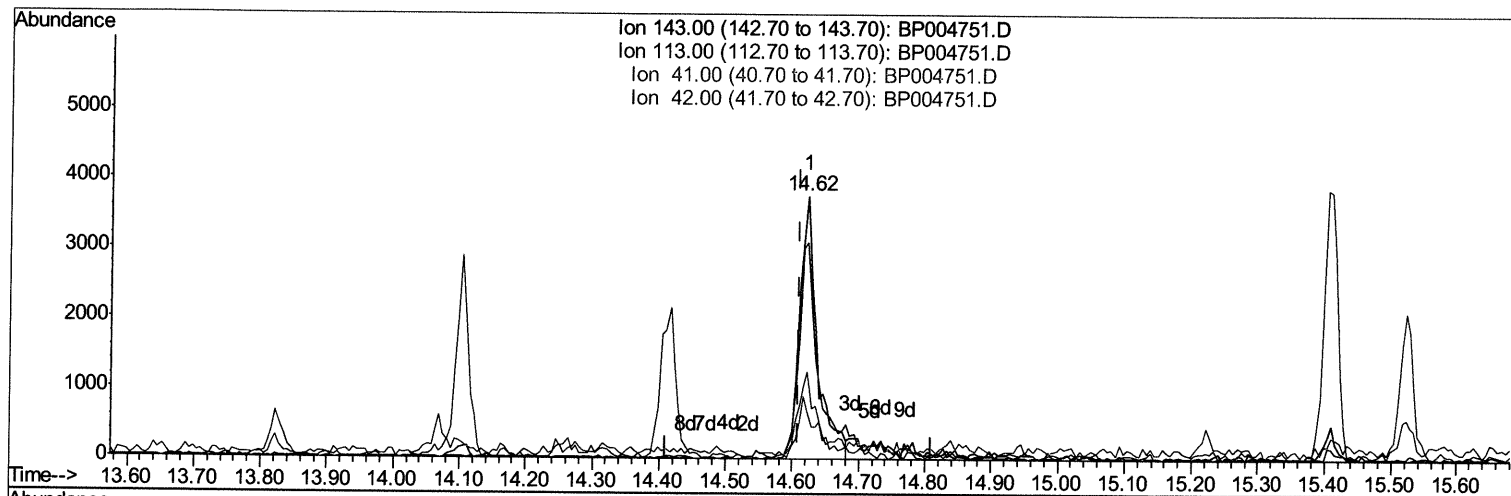
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Manual Integrations
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Quant Title : SVOA CALIBRATION
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Response via : Initial Calibration



TIC: BP004751.D

(52) 4-Nitrophenol-d4 (S)

14.622min (+0.012) 5.76ng/ul m 02/17/21 JU

response 6951

Ion	Exp%	Act%
143.00	100	100
113.00	104.70	81.94#
41.00	37.50	33.30
42.00	24.30	18.14#

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP021521\
 Data File : BP004751.D
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 Operator : CG/JU
 Sample : M1349-17
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBGS9

Manual Integrations
 APPROVED

mohammad
 2/16/2021 10:45:38 AM

Quant Time: Feb 16 00:18:41 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP020821MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Feb 15 12:16:19 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
1) 1,4-Dichlorobenzene-d4	7.77	152	44527	20.00	ng/ul	0.00
18) Naphthalene-d8	10.56	136	165860	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.41	164	95179	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.17	188	195070	20.00	ng/ul	0.00
78) Chrysene-d12	21.25	240	217041	20.00	ng/ul	0.00
86) Perylene-d12	23.56	264	229002	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.25	96	6239	5.42	ng/uL	0.00
5) Phenol-d5	6.93	99	24015	6.72	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.10	67	55096	29.68	ng/ul	0.00
9) 2-Chlorophenol-d4	7.30	132	70404	26.00	ng/ul	0.00
13) 4-Methylphenol-d8	8.48	113	41584	14.79	ng/ul	0.00
19) Nitrobenzene-d5	8.92	128	37429	32.69	ng/ul	0.00
22) 2-Nitrophenol-d4	9.65	143	44277	35.23	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.18	165	70782	27.98	ng/ul	0.00
29) 4-Chloroaniline-d4	10.70	131	3922	1.39	ng/ul	0.00
44) Dimethylphthalate-d6	13.82	166	244874	35.44	ng/ul	0.00
47) Acenaphthylene-d8	14.10	160	282651	33.63	ng/ul	0.00
52) 4-Nitrophenol-d4	14.62	143	6951m >	5.76	ng/ul >	0.01
58) Fluorene-d10	15.41	176	206554	34.89	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.52	200	40065	33.98	ng/ul	0.00
71) Anthracene-d10	17.27	188	341691	39.72	ng/ul	0.00
79) Pyrene-d10	19.50	212	401940	39.56	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.41	264	471466	40.70	ng/ul	0.00

02/17/2021

Target Compounds Ovalue

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