

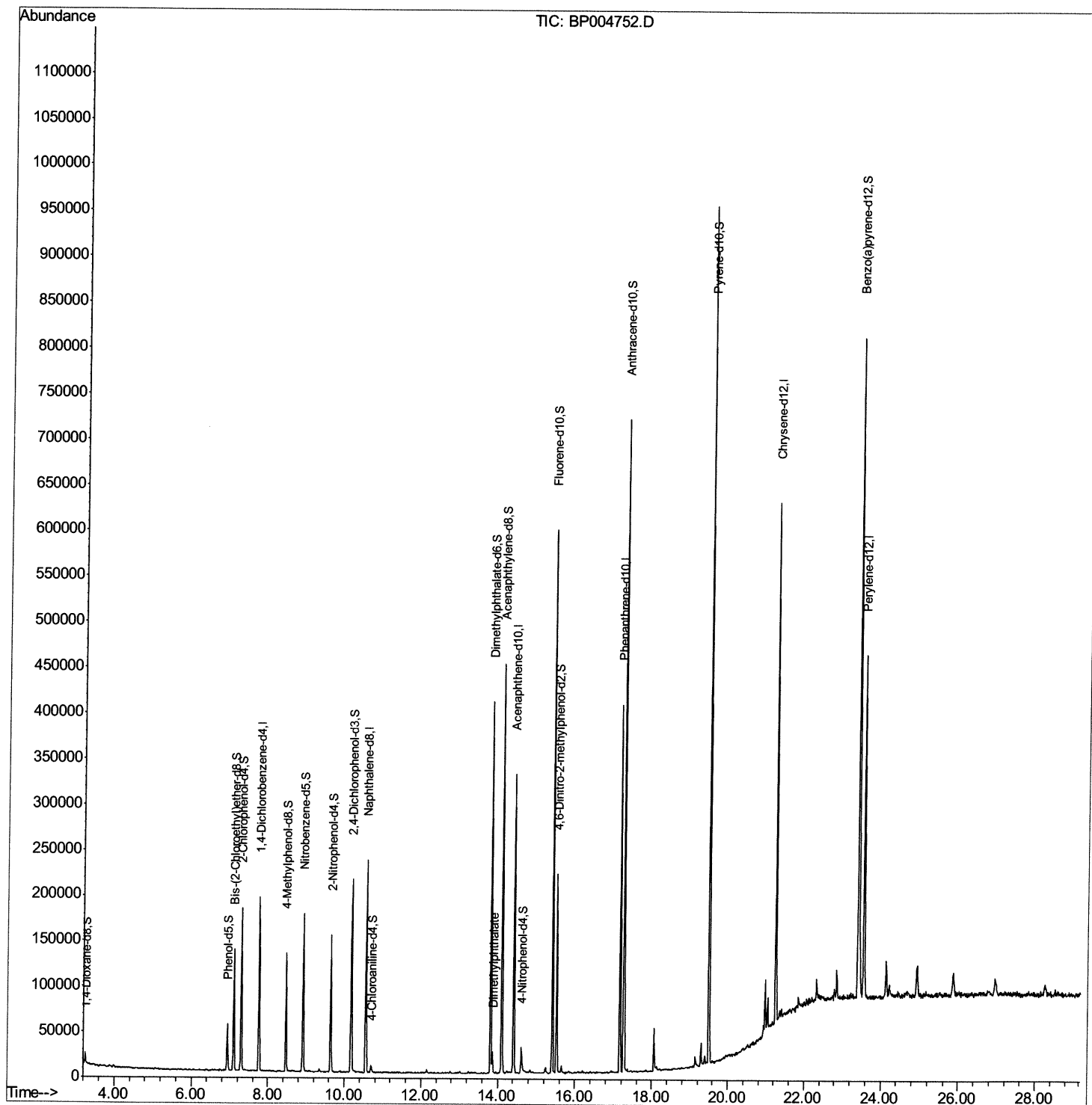
Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP021521\
Data File : BP004752.D
Acq On : 15 Feb 2021 18:24
Operator : CG/JU
Sample : M1349-15
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBG2

Manual Integrations
APPROVED

mohammad
2/16/2021 10:45:41 AM

Quant Time: Feb 16 00:26:13 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)

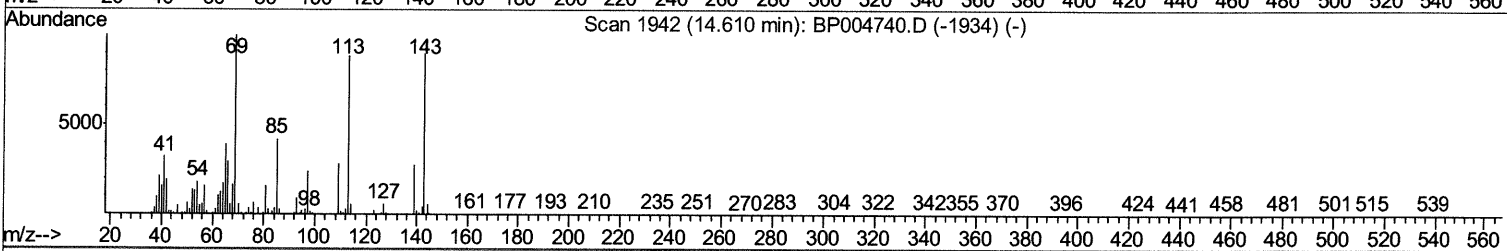
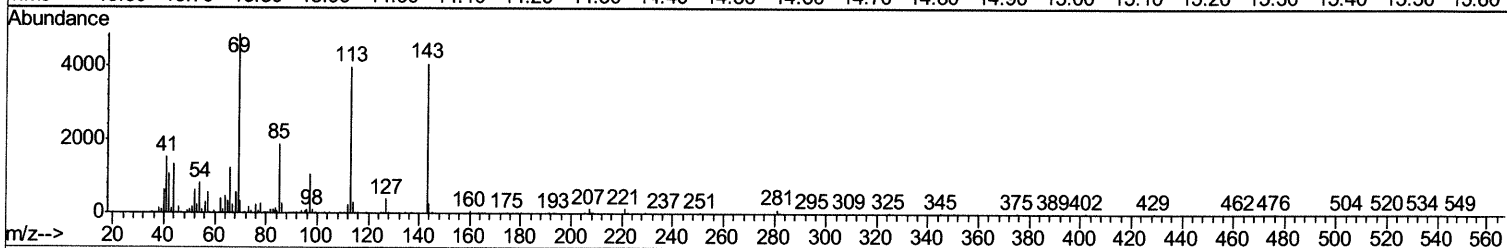
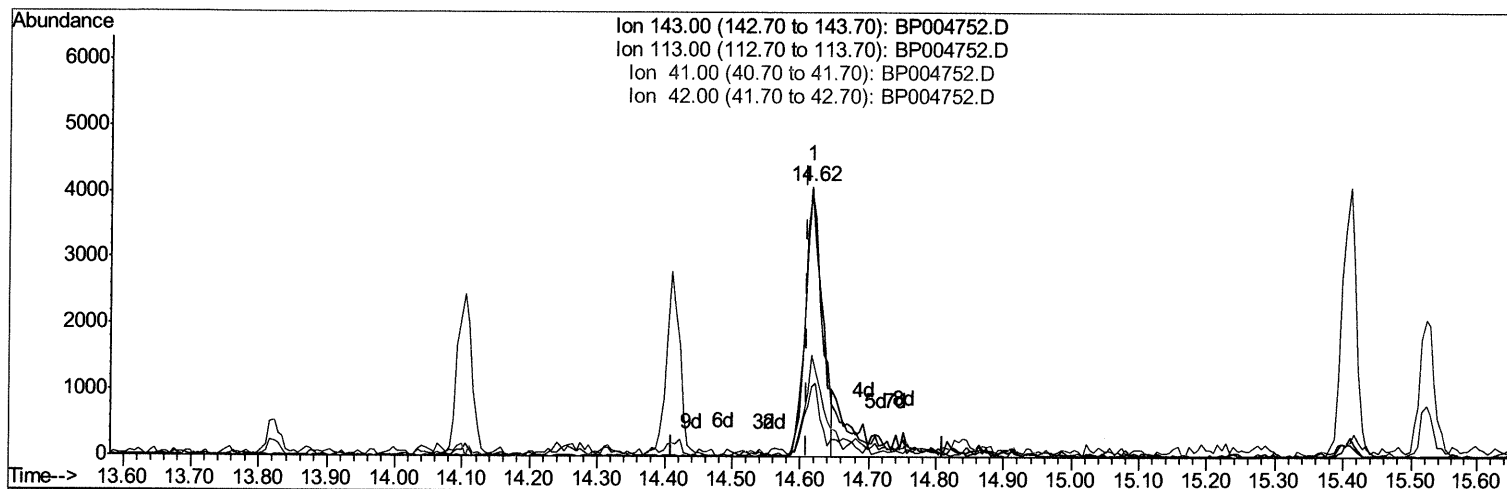
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TIC: BP004752.D

(52) 4-Nitrophenol-d4 (S)

14.616min (+0.006) 5.06ng/ul

response 6967

Ion	Exp%	Act%
143.00	100	100
113.00	104.70	97.70
41.00	37.50	37.57
42.00	24.30	26.06

Quantitation Report (Qedit)

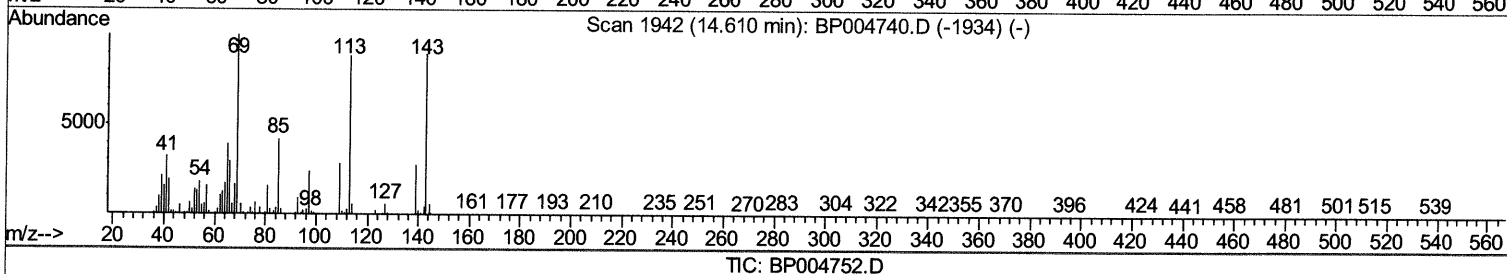
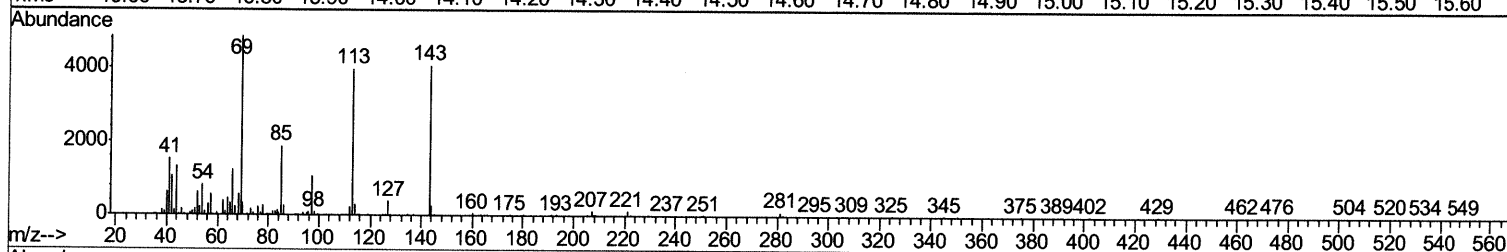
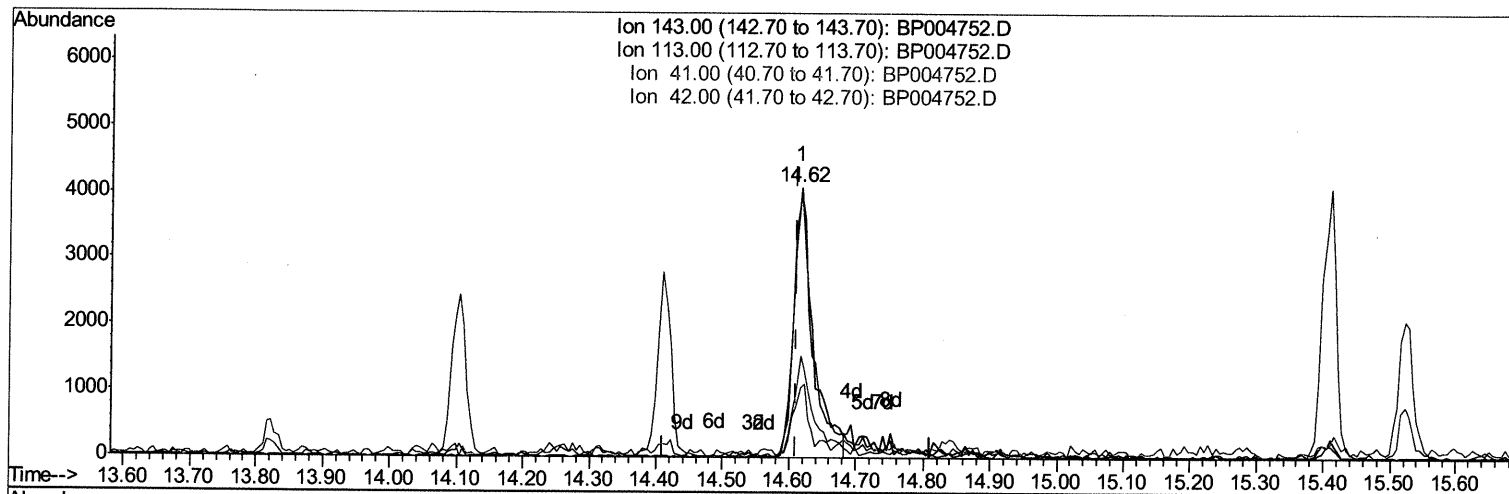
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Instrument :
BNA_P
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DBGT2

Manual Integrations
APPROVED

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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)

14.616min (+0.006) 5.91ng/ul m 02/17/21 JU

response 8152

Ion	Exp%	Act%
143.00	100	100
113.00	104.70	97.70
41.00	37.50	37.57
42.00	24.30	26.06

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP021521\
 Data File : BP004752.D
 Acq On : 15 Feb 2021 18:24
 Operator : CG/JU
 Sample : M1349-15
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBG2

Manual Integrations
 APPROVED

mohammad
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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.76	152	48045	20.00	ng/ul	0.00
18) Naphthalene-d8	10.56	136	180605	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.41	164	108636	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.17	188	229249	20.00	ng/ul	0.00
78) Chrysene-d12	21.25	240	258091	20.00	ng/ul	0.00
86) Perylene-d12	23.56	264	264612	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.25	96	6559	5.28	ng/uL	0.00
5) Phenol-d5	6.93	99	26250	6.81	ng/ul	0.00
7) Bis-(2-Chloroethyl) ether-d	7.10	67	53911	26.92	ng/ul	0.00
9) 2-Chlorophenol-d4	7.30	132	70503	24.13	ng/ul	0.00
13) 4-Methylphenol-d8	8.47	113	45182	14.89	ng/ul	0.00
19) Nitrobenzene-d5	8.92	128	36813	29.52	ng/ul	0.00
22) 2-Nitrophenol-d4	9.64	143	42678	31.19	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.18	165	71617	26.00	ng/ul	0.00
29) 4-Chloroaniline-d4	10.70	131	4268	1.39	ng/ul	0.00
44) Dimethylphthalate-d6	13.82	166	243151	30.83	ng/ul	0.00
47) Acenaphthylene-d8	14.10	160	294854	30.73	ng/ul	0.00
52) 4-Nitrophenol-d4	14.62	143	8152m>	5.91	ng/ul>	0.00
58) Fluorene-d10	15.41	176	218310	32.31	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.52	200	45575	32.89	ng/ul	0.00
71) Anthracene-d10	17.27	188	385889	38.17	ng/ul	0.00
79) Pyrene-d10	19.50	212	460099	38.08	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.41	264	532399	39.77	ng/ul	0.00

02/17/21J4

Target Compounds	R.T.	QIon	Response	Conc	Units	Ovalue
45) Dimethylphthalate	13.87	163	12355	1.519	ng/ul#	96

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