

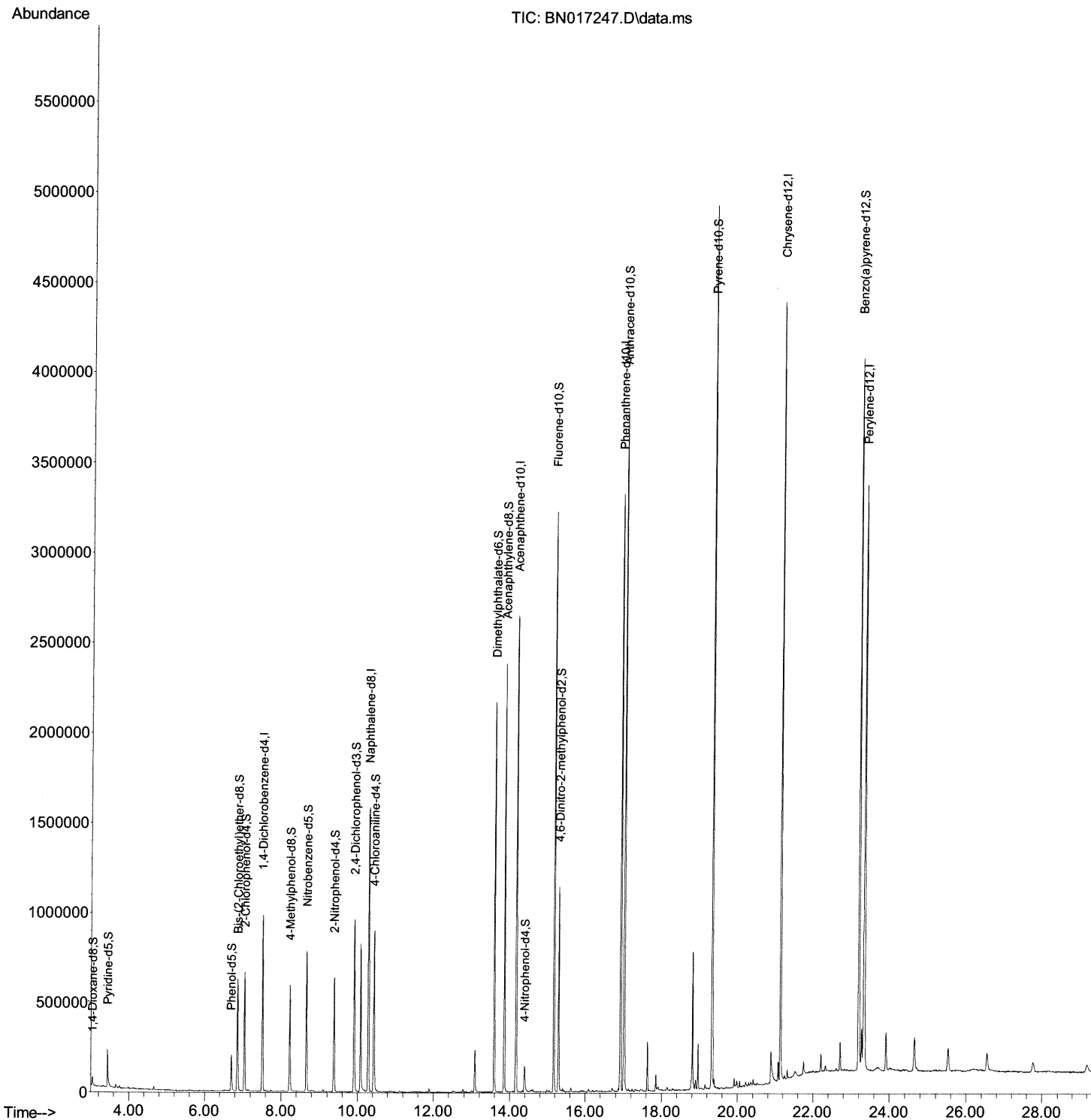
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN110221\
Data File : BN017247.D
Acq On : 02 Nov 2021 19:12
Operator : CG/JU
Sample : M4400-22
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BG356

Manual IntegrationsAPPROVED

Quant Time: Nov 02 21:36:49 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN110221.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Nov 02 15:59:34 2021
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 11/03/2021
Supervised By :mohammad ahmed 11/08/2021



Quantitation Report (Qedit)

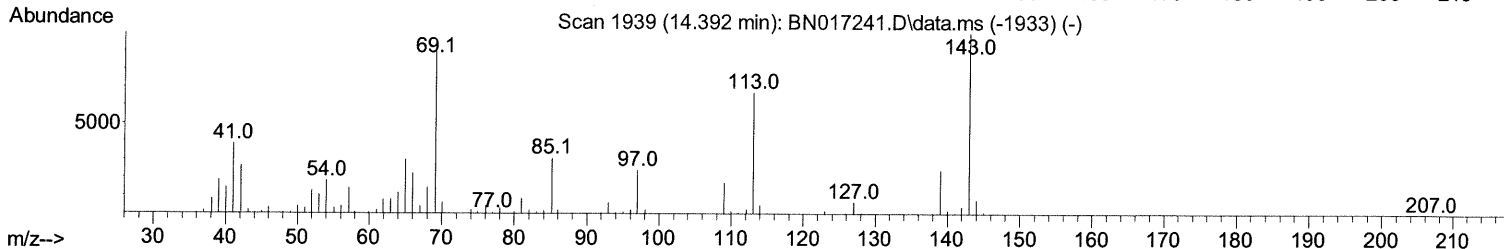
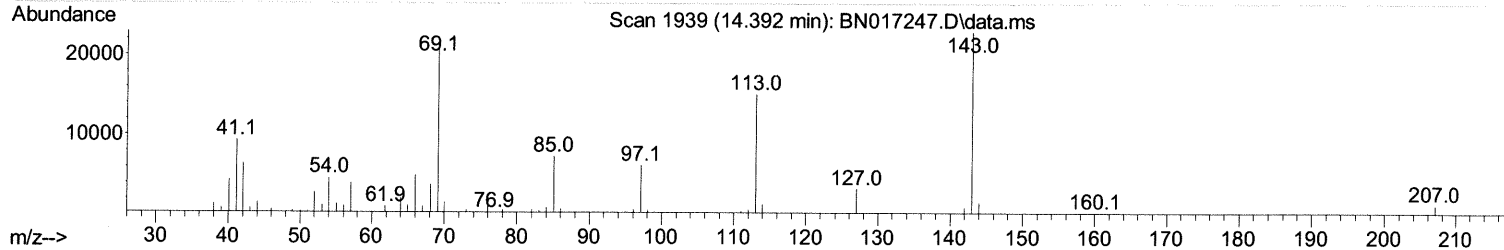
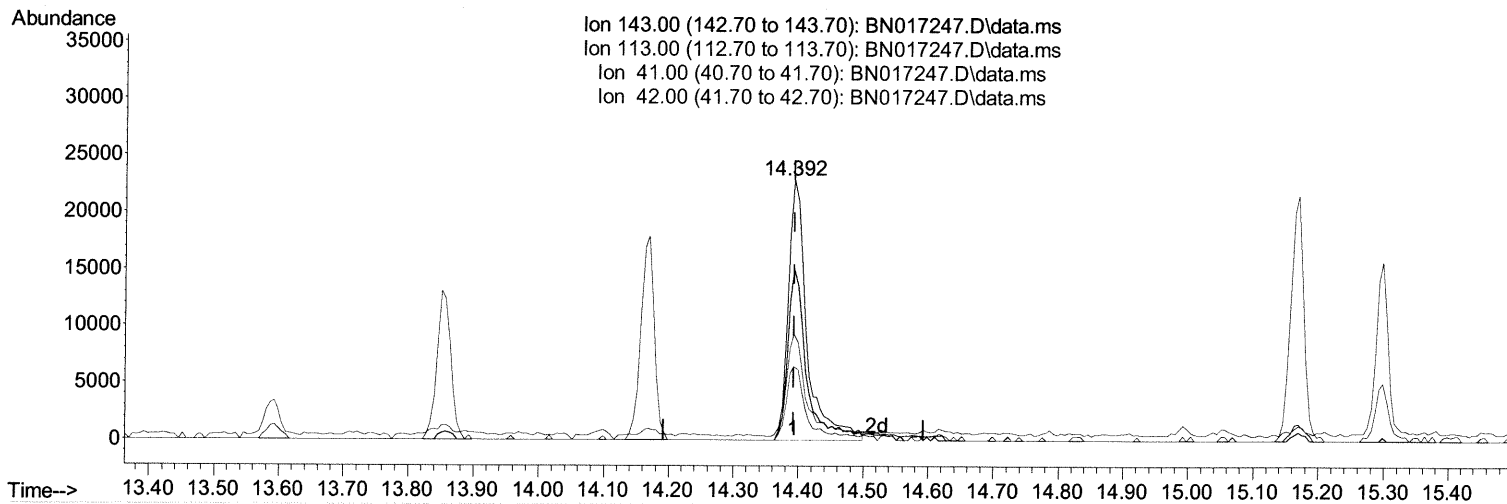
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TIC: BN017247.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.392min (-0.000) 3.65 ng/ul

response 45966

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	67.10	65.94
41.00	38.90	40.67
42.00	26.50	28.25

Quantitation Report (Qedit)

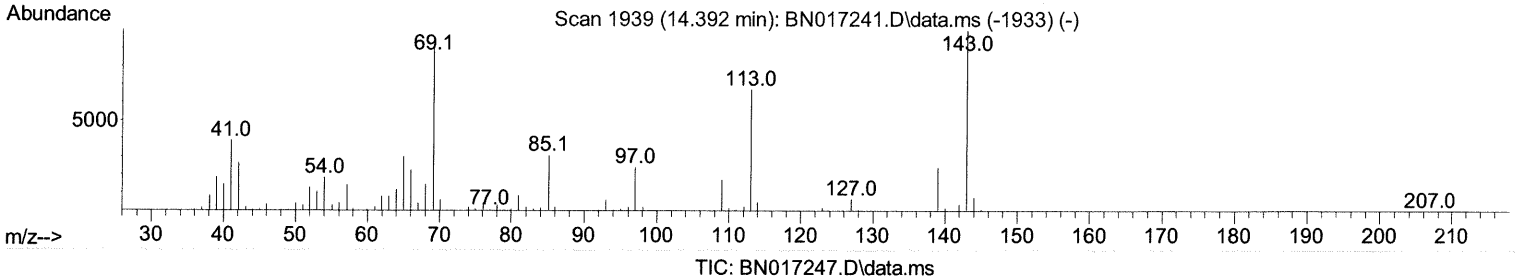
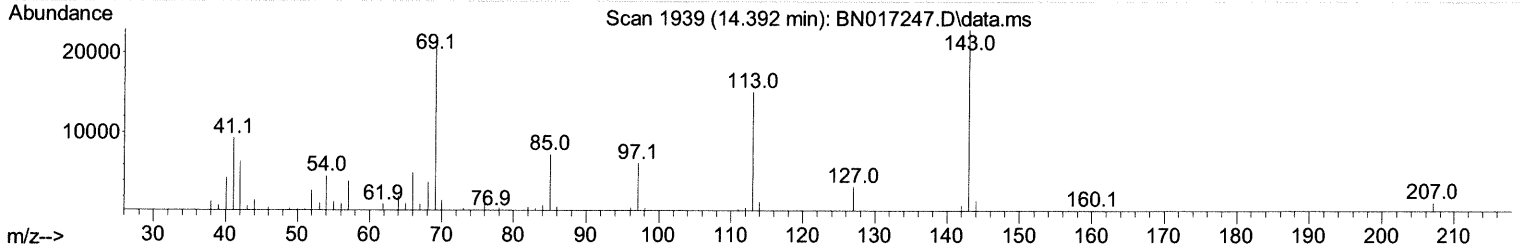
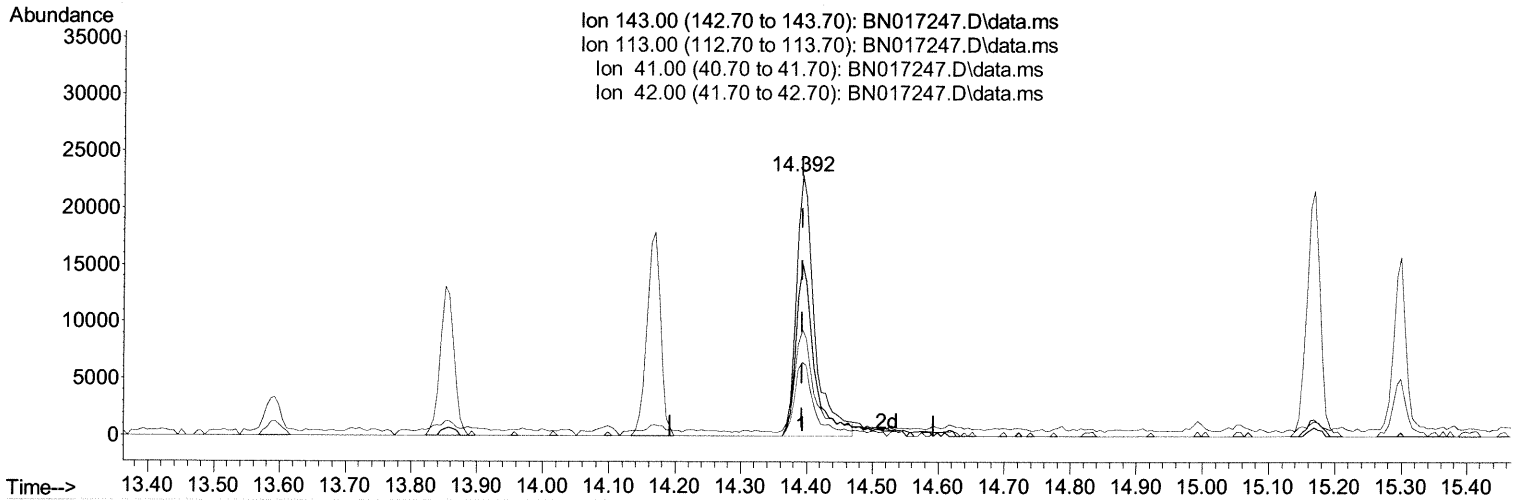
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 Supervised By :mohammad ahmed 11/08/2021



(54) 4-Nitrophenol-d4 (S)

14.392min (-0.000) 3.52 ng/ul m 11/04/2021

response 44382

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	67.10	65.94
41.00	38.90	40.67
42.00	26.50	28.25

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN110221\
 Data File : BN017247.D
 Acq On : 02 Nov 2021 19:12
 Operator : CG/JU
 Sample : M4400-22
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BG356

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 11/03/2021
 Supervised By :mohammad ahmed 11/08/2021

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.510	152	253644	20.000	ng/ul	0.00
20) Naphthalene-d8	10.287	136	1290012	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.169	164	866327	20.000	ng/ul	0.00
64) Phenanthrene-d10	16.922	188	1848999	20.000	ng/ul	0.00
79) Chrysene-d12	21.133	240	2007217	20.000	ng/ul	0.00
88) Perylene-d12	23.327	264	2173033	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.034	96	25386	3.956	ng/uL	0.00
4) Pyridine-d5	3.434	84	124204	6.936	ng/ul	0.00
7) Phenol-d5	6.699	99	111089	4.745	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.863	67	275808	19.736	ng/ul	0.00
11) 2-Chlorophenol-d4	7.046	132	305296	16.409	ng/ul	0.00
15) 4-Methylphenol-d8	8.228	113	213514	11.061	ng/ul	0.00
21) Nitrobenzene-d5	8.663	128	185513	18.534	ng/ul	0.00
24) 2-Nitrophenol-d4	9.381	143	204632	18.330	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.916	165	349732	17.423	ng/ul	0.00
31) 4-Chloroaniline-d4	10.434	131	473135	15.728	ng/ul	0.00
46) Dimethylphthalate-d6	13.592	166	1397519	21.757	ng/ul	0.00
49) Acenaphthylene-d8	13.857	160	1624092	20.102	ng/ul	0.00
54) 4-Nitrophenol-d4	14.392	143	44382m	3.520	ng/ul >	0.00 11/04/21 JU
60) Fluorene-d10	15.169	176	1201143	21.913	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.298	200	221380	18.718	ng/ul	0.00
73) Anthracene-d10	17.022	188	2011010	23.047	ng/ul	0.00
81) Pyrene-d10	19.327	212	2552165	23.497	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.192	264	2592145	22.080	ng/ul	0.00

Target Compounds Qvalue

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