

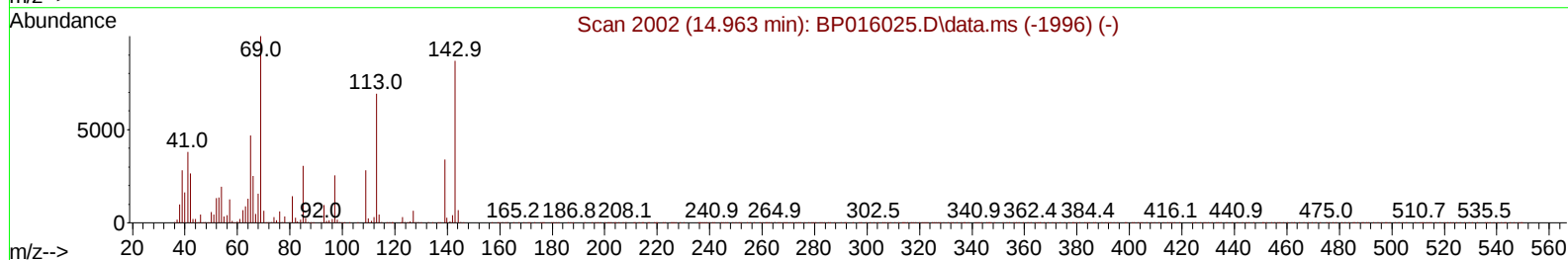
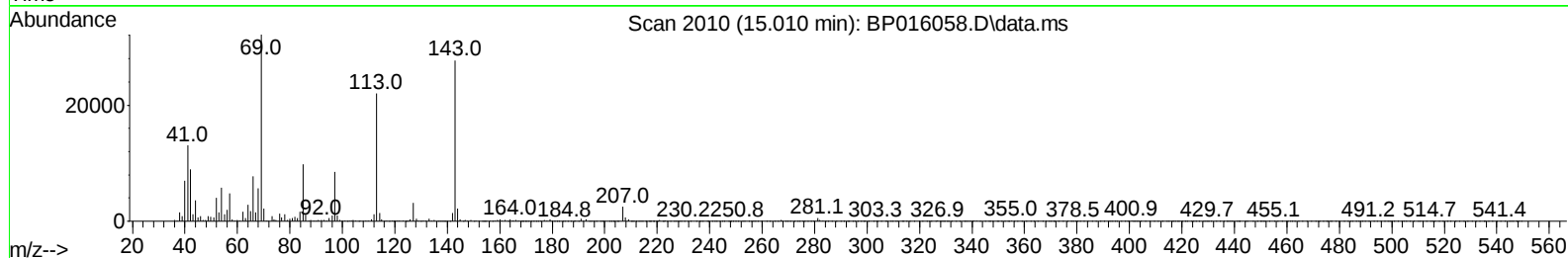
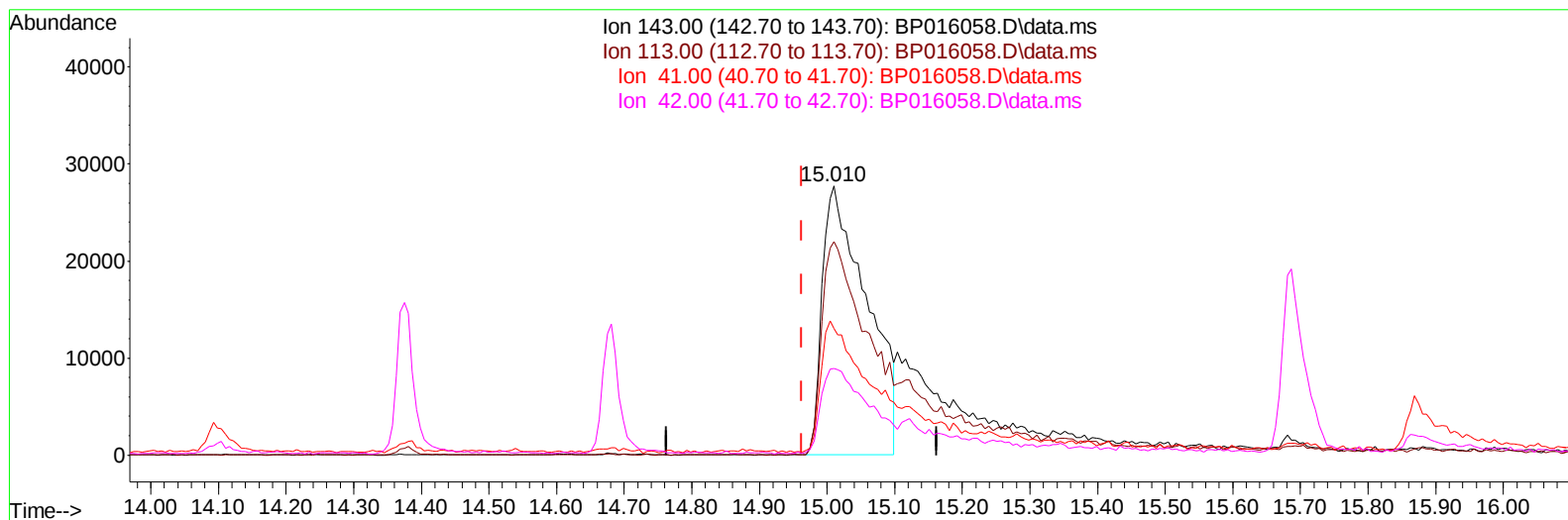
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070623\
Data File : BP016058.D
Acq On : 07 Jul 2023 06:33
Operator : MA/JU
Sample : PB153602BL
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SBLK602

Manual IntegrationsAPPROVED

Quant Time: Jul 07 07:07:36 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Jul 06 23:18:26 2023
Response via : Initial Calibration

Reviewed By :Yogesh Patel 07/07/2023
Supervised By :Sohil Jodhani 07/07/2023



TIC: BP016058.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.010min (+ 0.047) 21.85 ng/u1

response 126658

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	94.00	79.28
41.00	45.70	47.02
42.00	31.20	32.18

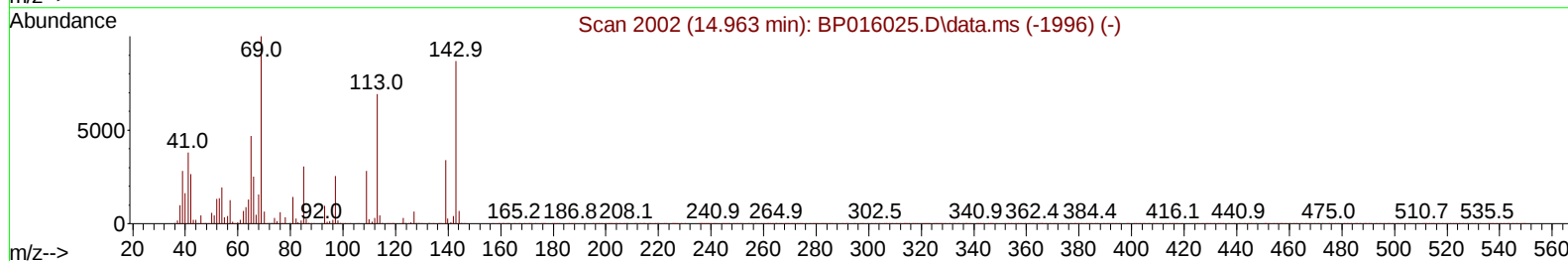
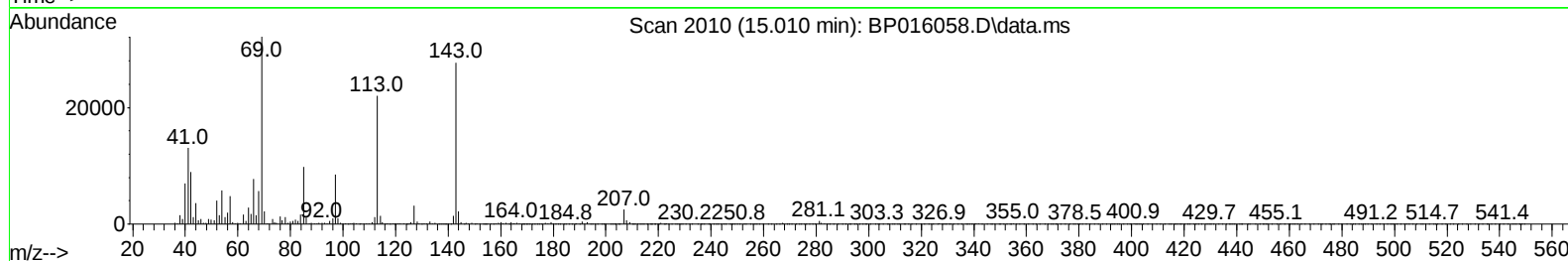
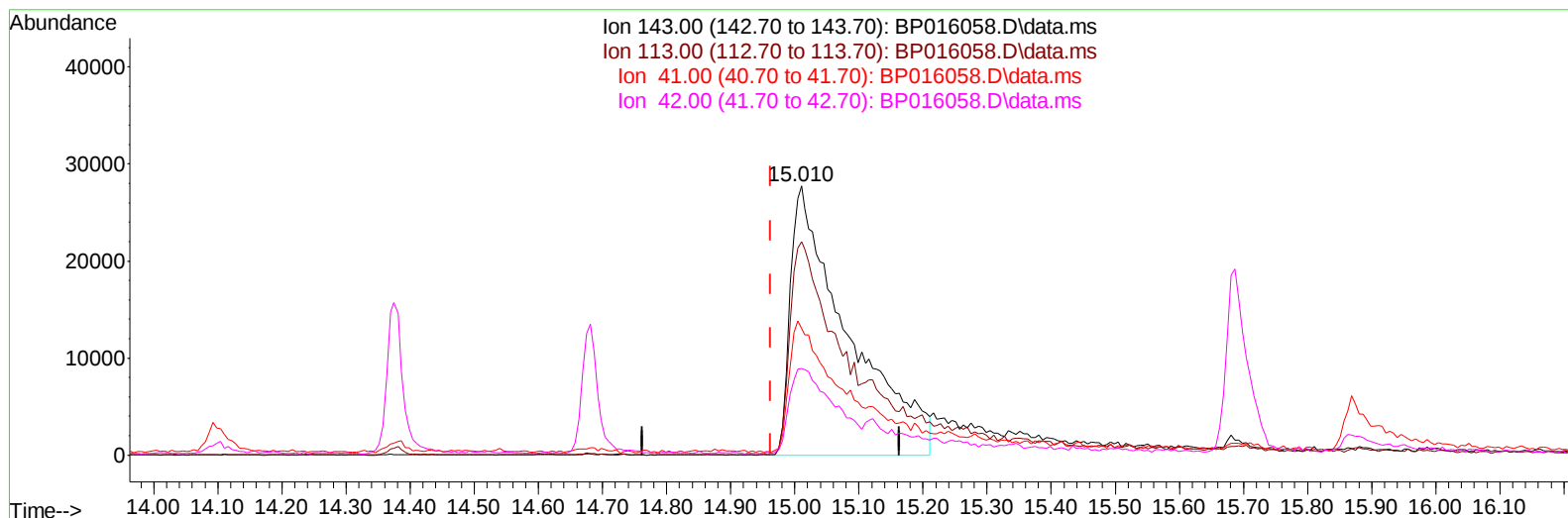
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TIC: BP016058.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.010min (+ 0.047) 29.90 ng/ul m

response 173353

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	94.00	79.28
41.00	45.70	47.02
42.00	31.20	32.18

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 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 07/07/2023
 Supervised By :Sohil Jodhani 07/07/2023

Quant Time: Jul 07 07:08:42 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jul 06 23:18:26 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.028	152	198473	20.000	ng/ul	0.00
20) Naphthalene-d8	10.857	136	871007	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.681	164	568424	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.445	188	1245344	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.557	240	1112630	20.000	ng/ul	0.02
88) Perylene-d12	24.086	264	1122275	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.352	96	38038	6.895	ng/uL	0.00
4) Pyridine-d5	3.799	84	461845	33.213	ng/ul	0.00
7) Phenol-d5	7.222	99	566109	33.337	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.352	67	406084	38.652	ng/ul	0.00
11) 2-Chlorophenol-d4	7.563	132	470186	36.679	ng/ul	0.00
15) 4-Methylphenol-d8	8.775	113	487621	36.348	ng/ul	0.00
21) Nitrobenzene-d5	9.228	128	233435	35.069	ng/ul	0.01
24) 2-Nitrophenol-d4	9.957	143	259984	33.464	ng/ul	0.01
28) 2,4-Dichlorophenol-d3	10.540	165	356802	25.772	ng/ul	0.04
31) 4-Chloroaniline-d4	11.051	131	605454	34.044	ng/ul	0.03
46) Dimethylphthalate-d6	14.098	166	1641370	37.359	ng/ul	0.01
49) Acenaphthylene-d8	14.375	160	1772701	35.893	ng/ul	0.00
54) 4-Nitrophenol-d4	15.010	143	173353m	29.900	ng/ul	0.05
60) Fluorene-d10	15.687	176	1332851	36.844	ng/ul	0.01
65) 4,6-Dinitro-2-methylph...	15.869	200	170345	22.242	ng/ul	0.02
73) Anthracene-d10	17.551	188	2120173	37.271	ng/ul	0.00
81) Pyrene-d10	19.780	212	2534942	34.891	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.916	264	2097684	37.053	ng/ul	0.00

Target Compounds Qvalue

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

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