

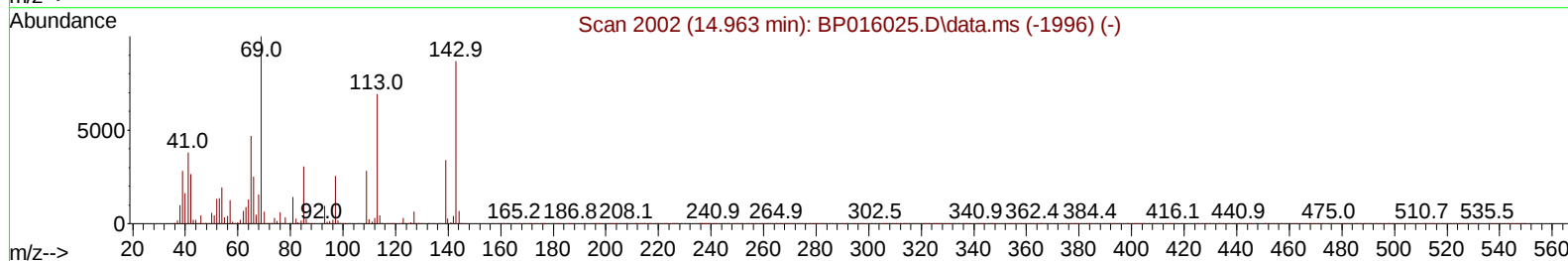
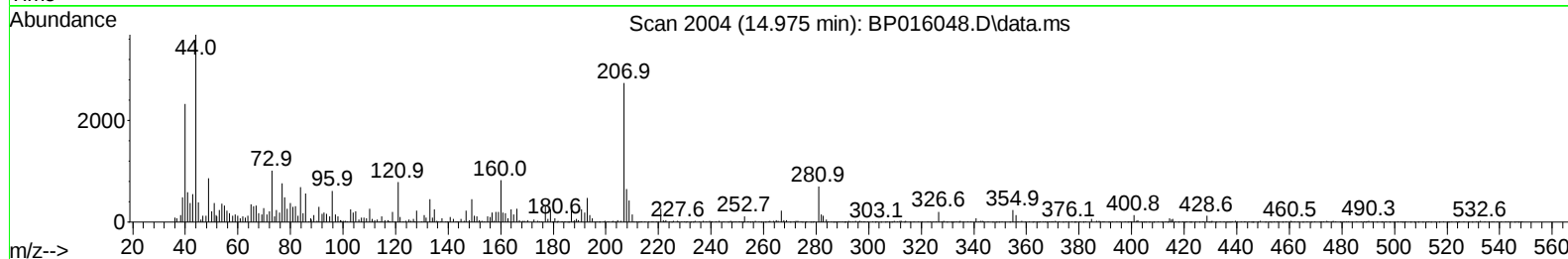
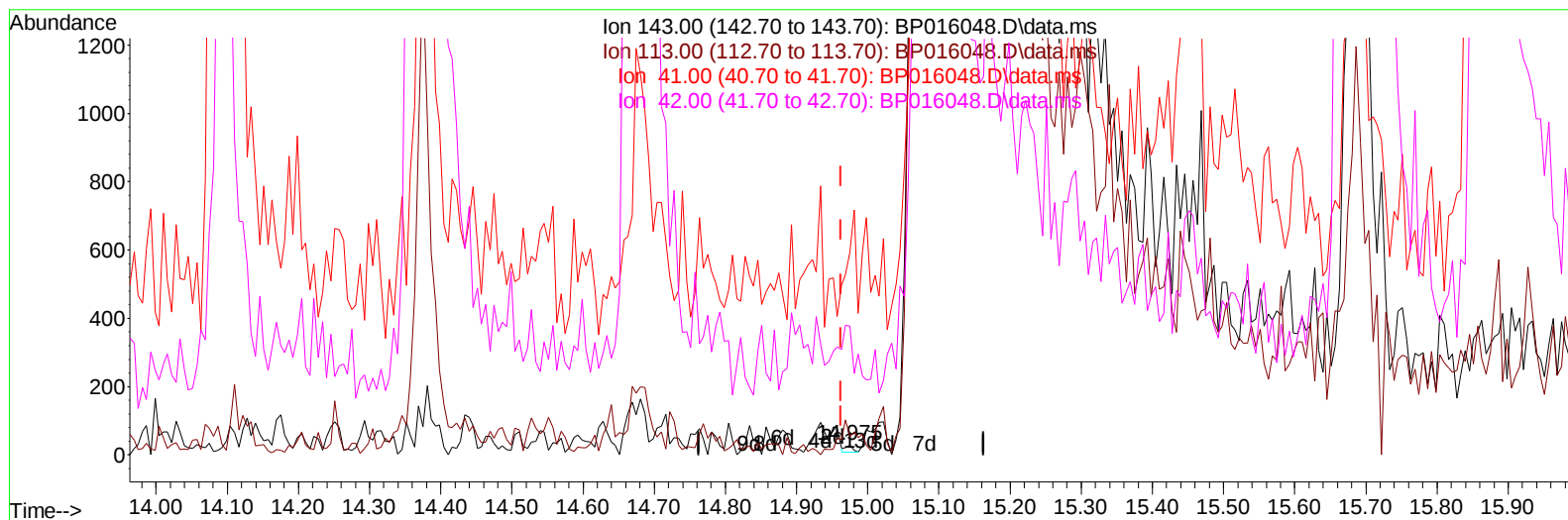
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070623\
Data File : BP016048.D
Acq On : 07 Jul 2023 00:19
Operator : MA/JU
Sample : 03283-08
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EXYX4

Manual IntegrationsAPPROVED

Quant Time: Jul 07 02:17:13 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 :mohammad ahmed 07/07/2023



TIC: BP016048.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.975min (+ 0.012) 0.00 ng/u1

response

14

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	94.00	189.66#
41.00	45.70	2041.38#
42.00	31.20	1303.45#

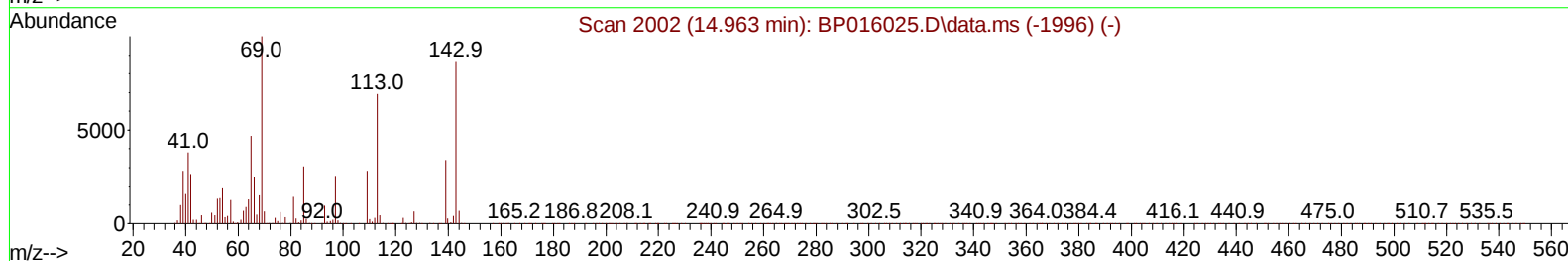
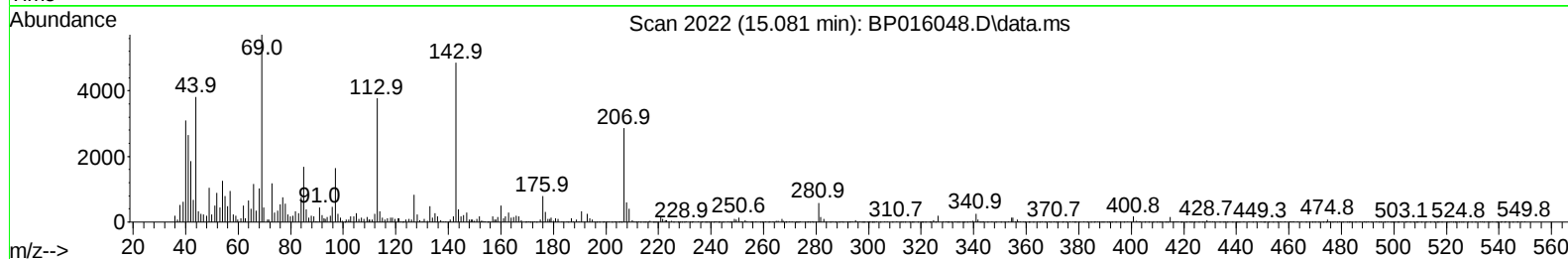
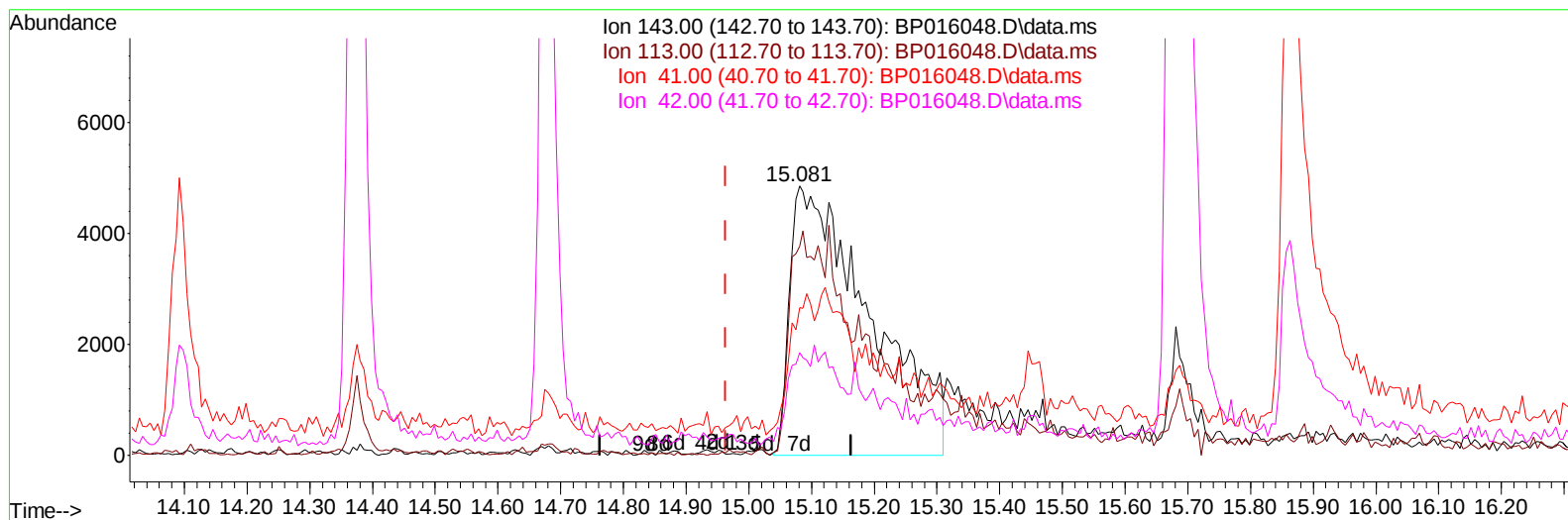
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TIC: BP016048.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.081min (+ 0.118) 6.30 ng/ul m

response 43389

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	94.00	77.59
41.00	45.70	54.78
42.00	31.20	38.20#

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 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EXYX4

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 07/07/2023
 Supervised By :mohammad ahmed 07/07/2023

Quant Time: Jul 07 02:18:43 2023
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 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	8.034	152	228005	20.000	ng/ul	0.00
20) Naphthalene-d8	10.857	136	1003566	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.681	164	674955	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.445	188	1546418	20.000	ng/ul	0.00
79) Chrysene-d12	21.551	240	1504408	20.000	ng/ul	0.01
88) Perylene-d12	24.086	264	1568614	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.352	96	32394	5.112	ng/uL	0.00
4) Pyridine-d5	3.817	84	109717	6.868	ng/ul	0.01
7) Phenol-d5	7.252	99	131043	6.717	ng/ul	0.04
9) Bis-(2-Chloroethyl)eth...	7.358	67	514342	42.615	ng/ul	0.00
11) 2-Chlorophenol-d4	7.575	132	460918	31.299	ng/ul	0.01
15) 4-Methylphenol-d8	8.787	113	303777	19.711	ng/ul	0.02
21) Nitrobenzene-d5	9.228	128	314015	40.943	ng/ul	0.01
24) 2-Nitrophenol-d4	9.957	143	345309	38.576	ng/ul	0.01
28) 2,4-Dichlorophenol-d3	10.540	165	469227	29.416	ng/ul	0.04
31) 4-Chloroaniline-d4	11.051	131	551822	26.930	ng/ul	0.03
46) Dimethylphthalate-d6	14.092	166	2481431	47.565	ng/ul	0.00
49) Acenaphthylene-d8	14.375	160	2590432	44.171	ng/ul	0.00
54) 4-Nitrophenol-d4	15.081	143	43389m	6.302	ng/ul	0.12
60) Fluorene-d10	15.681	176	2065427	48.083	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.863	200	354590	37.285	ng/ul	0.01
73) Anthracene-d10	17.551	188	3469549	49.118	ng/ul	0.00
81) Pyrene-d10	19.775	212	4239068	43.152	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.910	264	3807201	48.114	ng/ul	0.00
Target Compounds						
2) 1,4-Dioxane	3.387	88	45462	6.666	ng/uL	Qvalue 97

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

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