

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062923\
Data File : BP015829.D
Acq On : 30 Jun 2023 01:59
Operator : MA/JU
Sample : 03220-10
Misc :
ALS Vial : 33 Sample Multiplier: 1

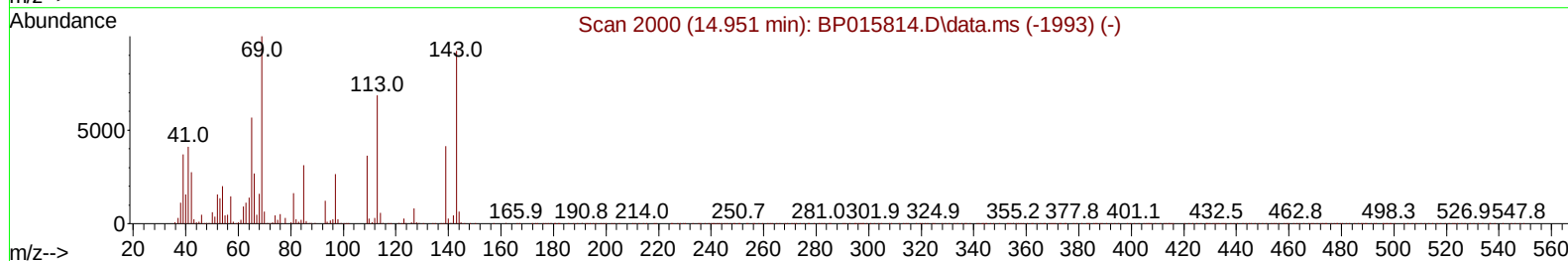
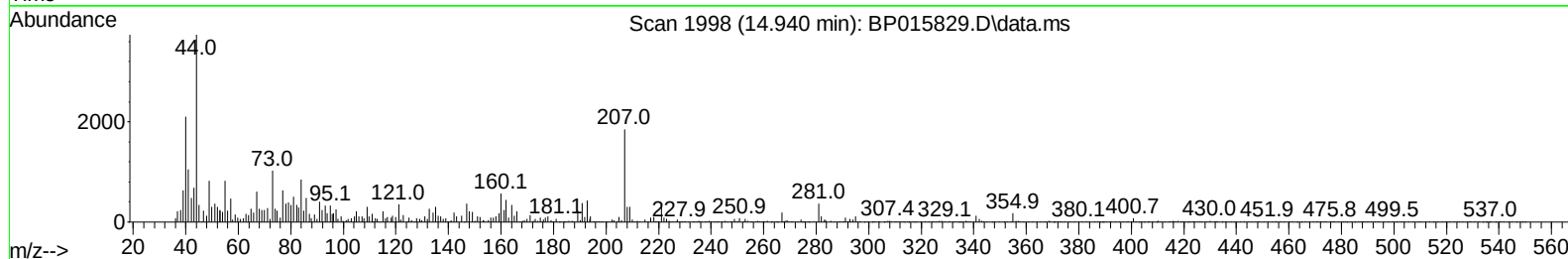
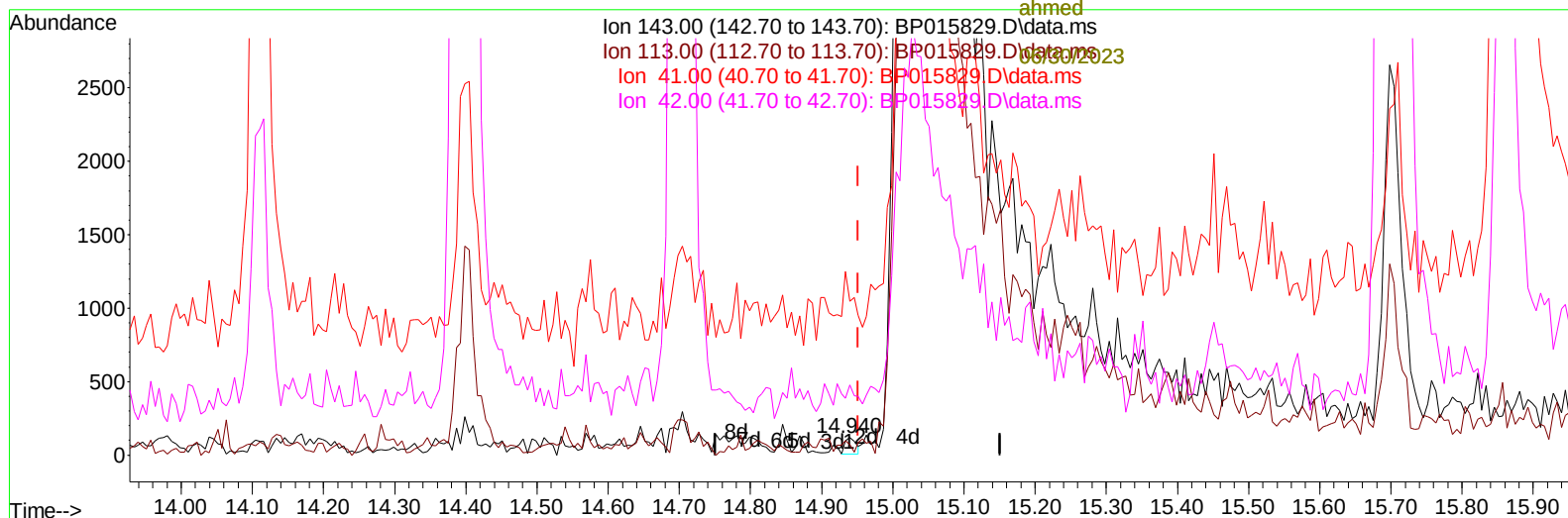
Instrument :
BNA_P
ClientSampleId :
CONR6

Manual IntegrationsAPPROVED

Quant Time: Jun 30 02:40:27 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Jun 29 02:30:23 2023
Response via : Initial Calibration

Reviewed By :Yogesh
Patel

06/30/2023
Supervised By :mohammad
ahmed



TIC: BP015829.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.940min (-0.012) 0.02 ng/u1

response 120

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	94.00	58.82#
41.00	45.70	879.83#
42.00	31.20	401.68#

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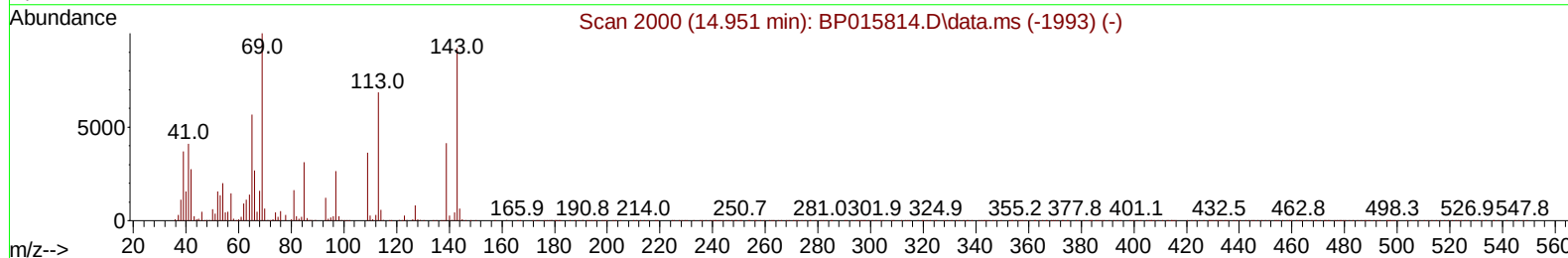
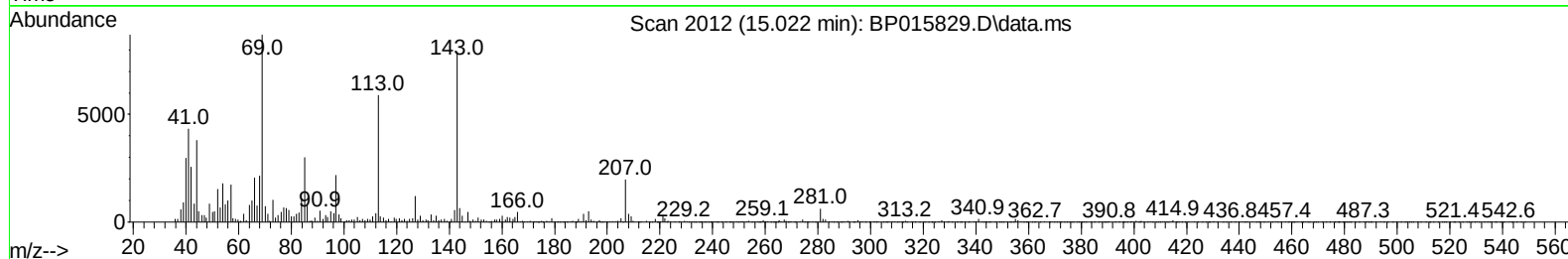
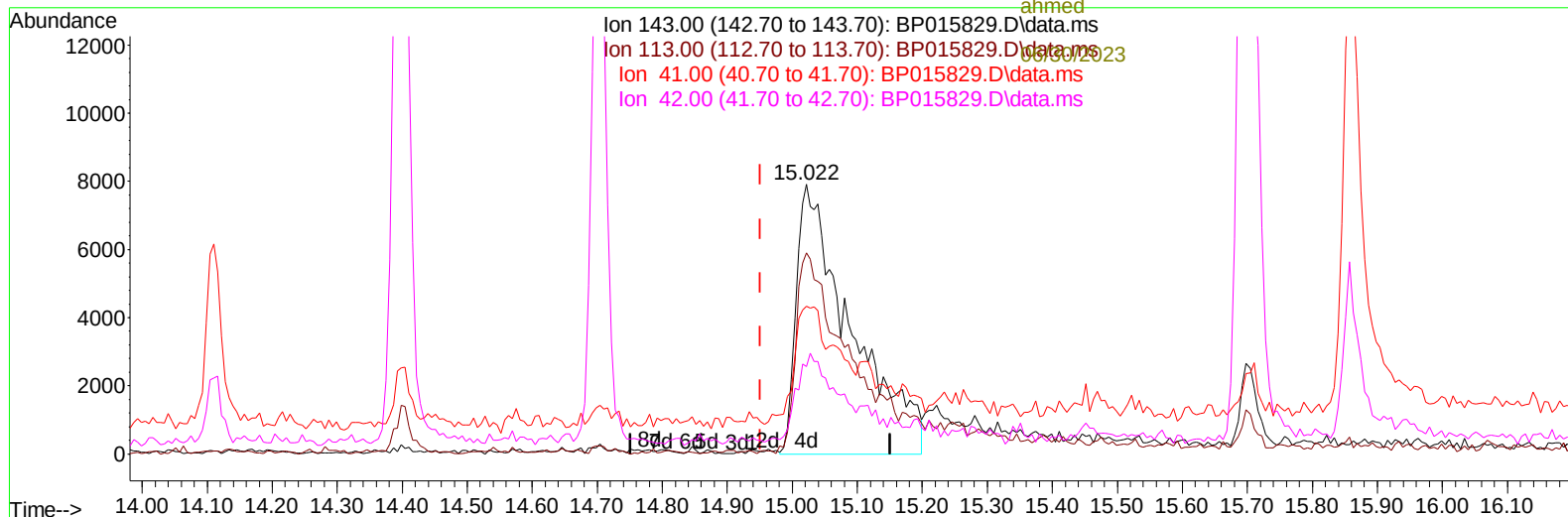
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(54) 4-Nitrophenol-d4 (S)

15.022min (+ 0.071) 5.92 ng/ul m

response 45917

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	94.00	74.57#
41.00	45.70	54.81
42.00	31.20	32.52

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
-----06/30/2023						
Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.052	152	278680	20.000	ng/ul	0.00
20) Naphthalene-d8	10.881	136	1245665	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.704	164	760057	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.463	188	1669050	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.557	240	1263039	20.000	ng/ul	# 0.00
88) Perylene-d12	24.098	264	1249058	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.375	96	32169	4.153	ng/uL	0.00
4) Pyridine-d5	3.834	84	157736	8.079	ng/ul	0.01
7) Phenol-d5	7.246	99	160158	6.717	ng/ul	0.02
9) Bis-(2-Chloroethyl)eth...	7.375	67	601129	40.749	ng/ul	0.00
11) 2-Chlorophenol-d4	7.587	132	530031	29.447	ng/ul	0.00
15) 4-Methylphenol-d8	8.793	113	342025	18.157	ng/ul	0.00
21) Nitrobenzene-d5	9.240	128	372279	39.106	ng/ul	0.00
24) 2-Nitrophenol-d4	9.969	143	395690	35.613	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.522	165	621407	31.385	ng/ul	0.00
31) 4-Chloroaniline-d4	11.040	131	893999	35.149	ng/ul	0.00
46) Dimethylphthalate-d6	14.110	166	2410725	41.036	ng/ul	0.00
49) Acenaphthylene-d8	14.398	160	2735271	41.419	ng/ul	0.00
54) 4-Nitrophenol-d4	15.022	143	45917m	5.923	ng/ul	0.07
60) Fluorene-d10	15.698	176	2075701	42.911	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.857	200	278345	27.117	ng/ul	0.00
73) Anthracene-d10	17.569	188	3393545	44.512	ng/ul	0.00
81) Pyrene-d10	19.792	212	3716716	45.065	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.933	264	2827862	44.881	ng/ul	0.00

Target Compounds Qvalue

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

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