

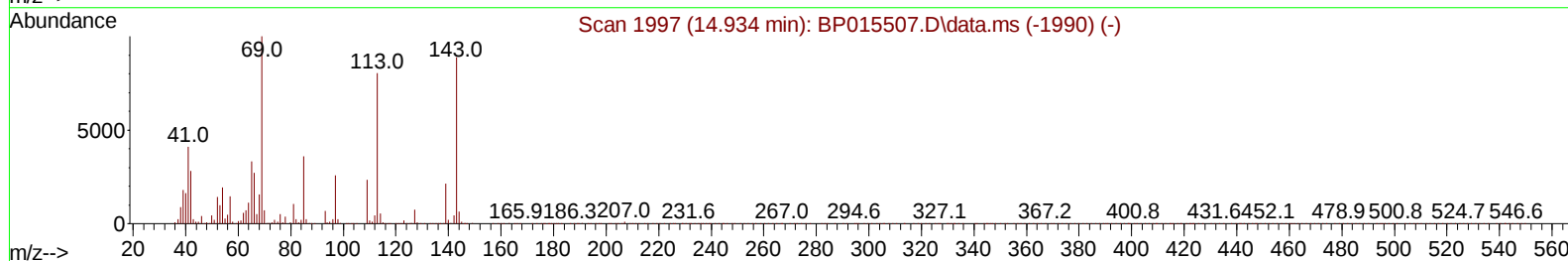
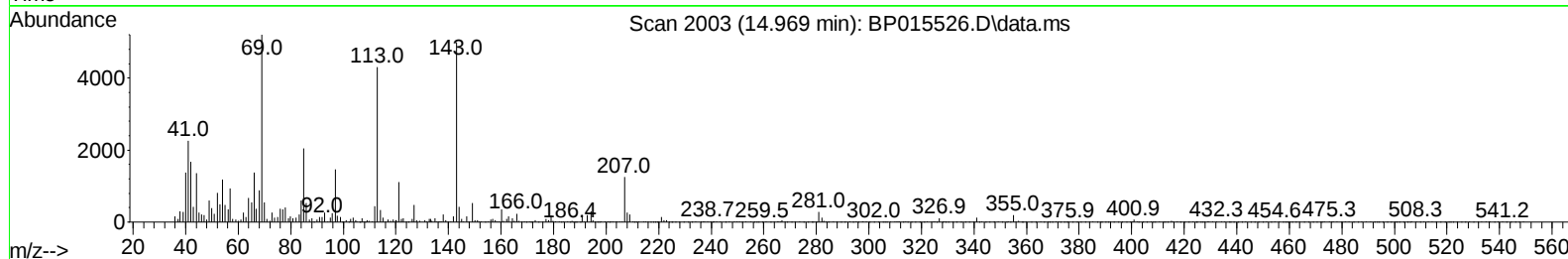
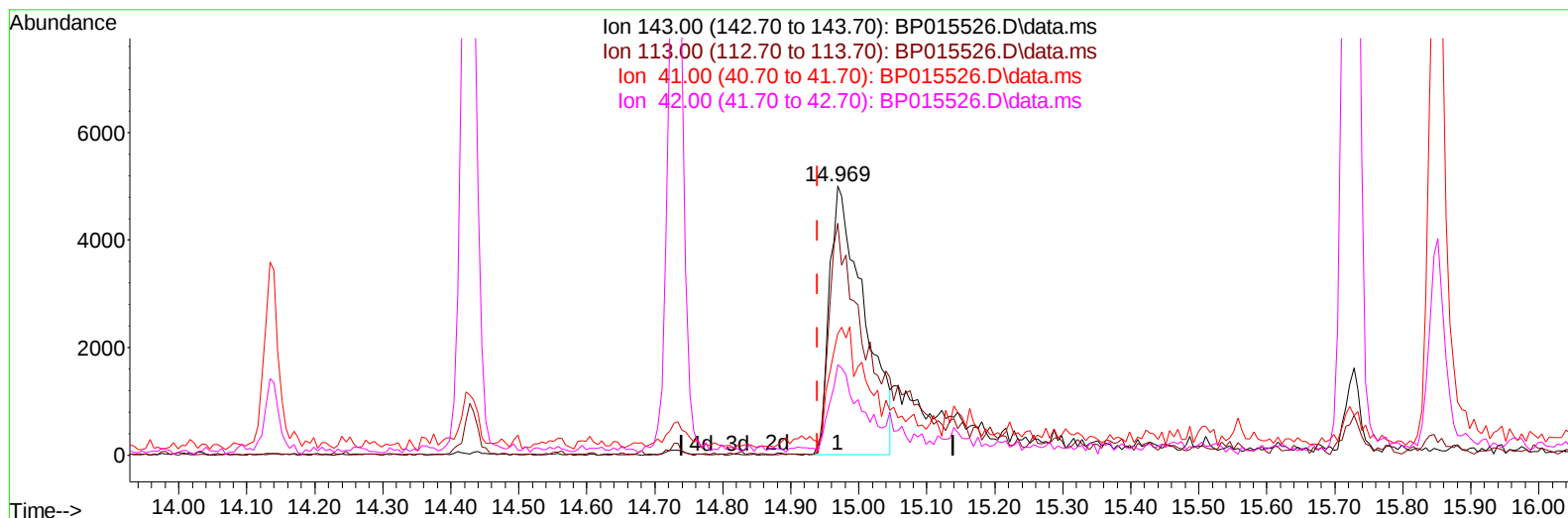
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061423\
Data File : BP015526.D
Acq On : 14 Jun 2023 20:00
Operator : MA/JU
Sample : 03097-18
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EZER0

Manual IntegrationsAPPROVED

Quant Time: Jun 14 23:11:31 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Jun 09 23:33:38 2023
Response via : Initial Calibration

Reviewed By :Yogesh Patel 06/15/2023
Supervised By :mohammad ahmed 06/16/2023



TIC: BP015526.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.969min (+ 0.029) 2.82 ng/u1

response 17631

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	86.60	86.15
41.00	42.20	45.14
42.00	29.40	33.65

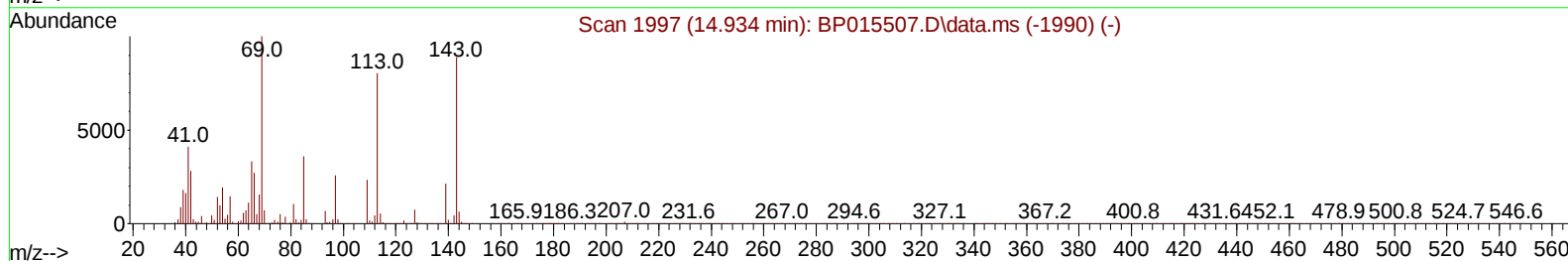
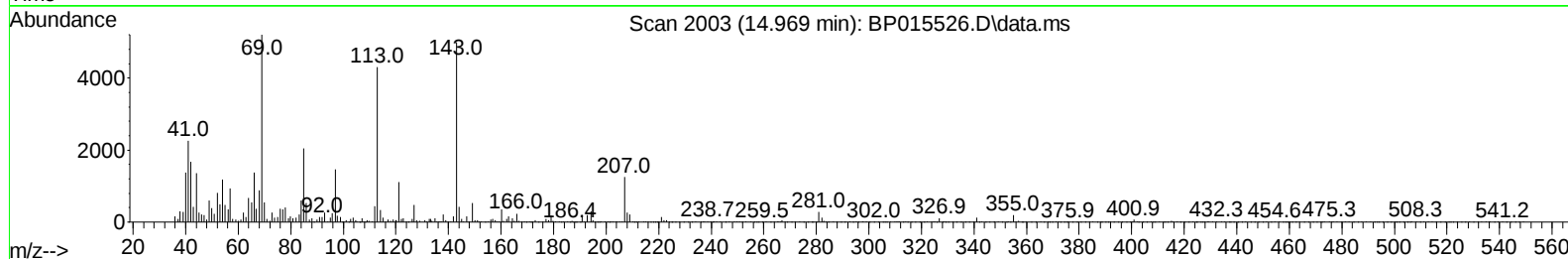
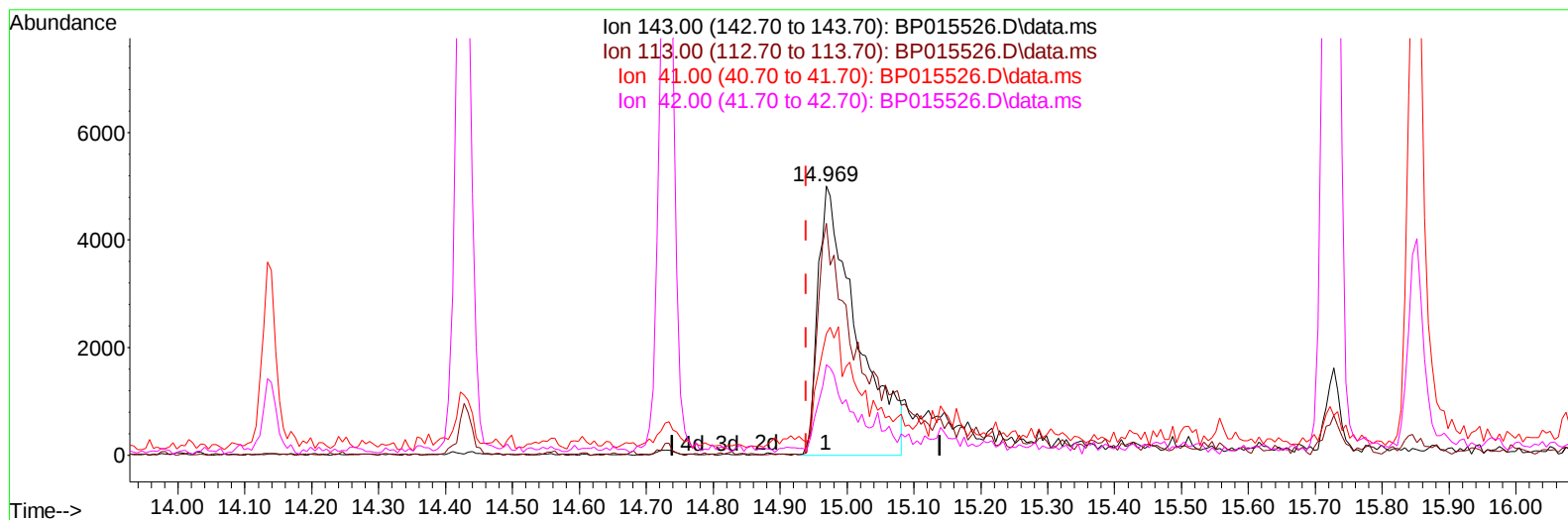
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TIC: BP015526.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.969min (+ 0.029) 3.21 ng/ul m

response 20061

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	86.60	86.15
41.00	42.20	45.14
42.00	29.40	33.65

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061423\
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 Operator : MA/JU
 Sample : 03097-18
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EZERO

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 06/15/2023
 Supervised By :mohammad ahmed 06/16/2023

Quant Time: Jun 14 23:56:51 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Jun 09 23:33:38 2023
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.087	152	137318	20.000	ng/ul	0.00
20) Naphthalene-d8	10.916	136	649451	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.734	164	442570	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.487	188	1059442	20.000	ng/ul	0.00
79) Chrysene-d12	21.569	240	1092276	20.000	ng/ul	0.00
88) Perylene-d12	24.116	264	1145927	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.411	96	18268	4.926	ng/uL	0.00
4) Pyridine-d5	3.858	84	71477	7.421	ng/ul	0.01
7) Phenol-d5	7.246	99	70831	5.599	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.411	67	233575	30.916	ng/ul	0.00
11) 2-Chlorophenol-d4	7.617	132	219510	23.933	ng/ul	0.00
15) 4-Methylphenol-d8	8.805	113	154505	14.881	ng/ul	0.00
21) Nitrobenzene-d5	9.269	128	155278	30.454	ng/ul	0.00
24) 2-Nitrophenol-d4	9.993	143	171776	29.500	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.534	165	282401	26.587	ng/ul	0.00
31) 4-Chloroaniline-d4	11.057	131	286766	18.809	ng/ul	0.00
46) Dimethylphthalate-d6	14.140	166	1204091	34.502	ng/ul	0.00
49) Acenaphthylene-d8	14.428	160	1264383	33.133	ng/ul	0.00
54) 4-Nitrophenol-d4	14.969	143	20061m	3.210	ng/ul	0.03
60) Fluorene-d10	15.728	176	1019462	34.641	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.851	200	198046	29.508	ng/ul	0.00
73) Anthracene-d10	17.587	188	1715977	35.599	ng/ul	0.00
81) Pyrene-d10	19.810	212	2239940	38.316	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.951	264	2213289	38.493	ng/ul	0.00

Target Compounds Qvalue

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

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