

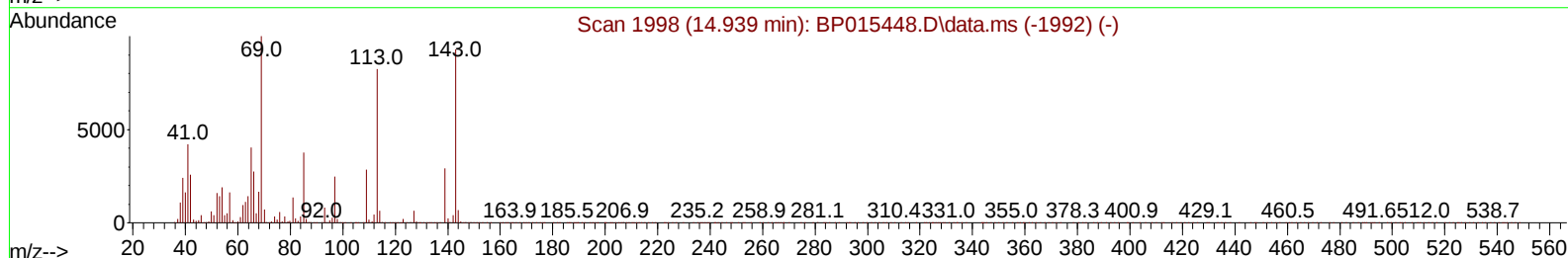
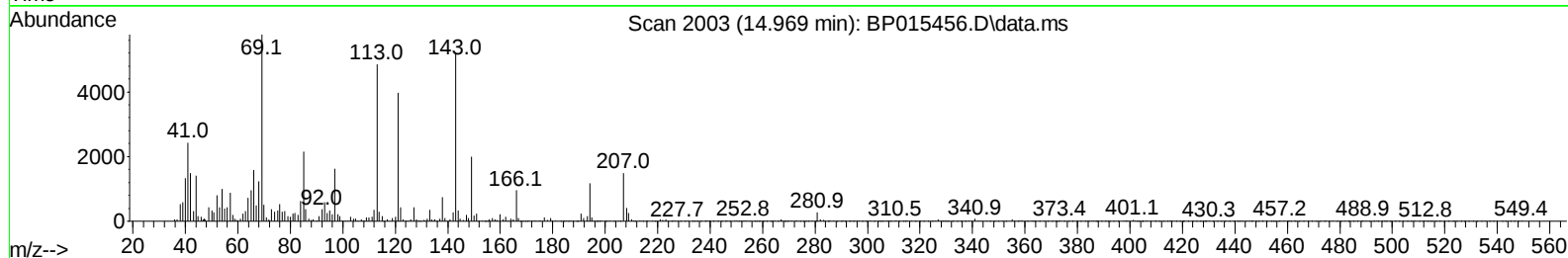
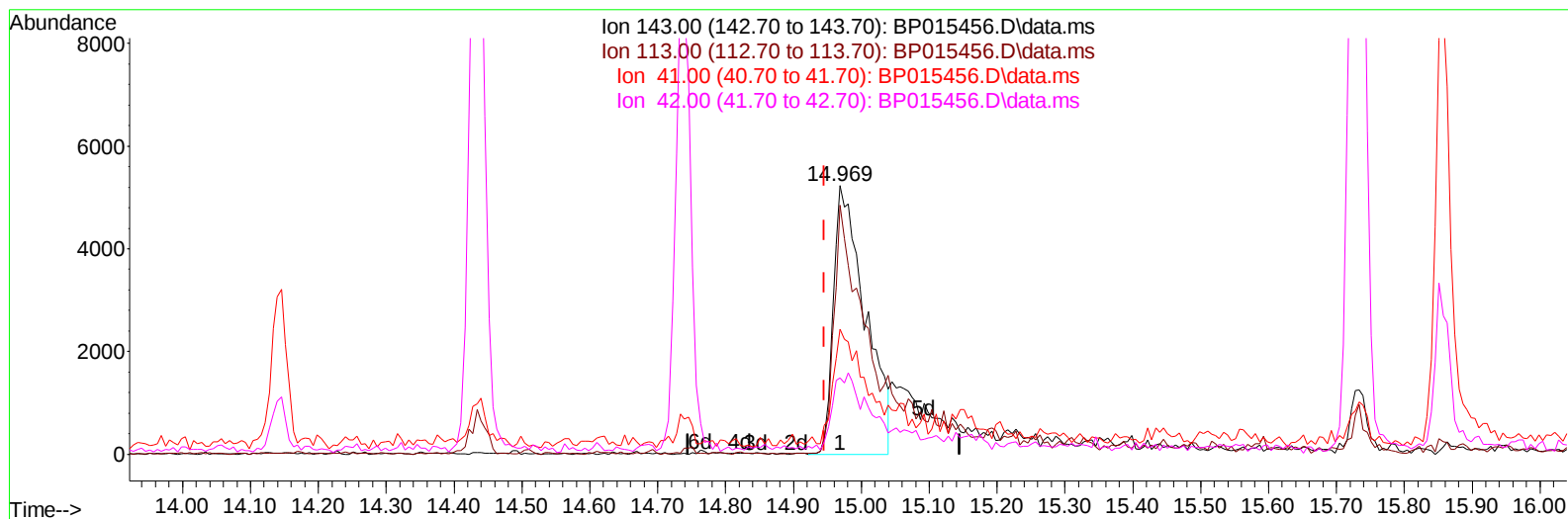
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP060923\
 Data File : BP015456.D
 Acq On : 09 Jun 2023 17:26
 Operator : MA/JU
 Sample : 03097-14
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EZEN7

Manual IntegrationsAPPROVED

Quant Time: Jun 09 22:58:48 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Jun 07 23:46:15 2023
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 06/12/2023
 Supervised By :mohammad ahmed 06/13/2023



TIC: BP015456.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.969min (+ 0.023) 3.08 ng/u1

response 17015

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	86.60	92.88
41.00	42.20	46.52
42.00	29.40	28.54

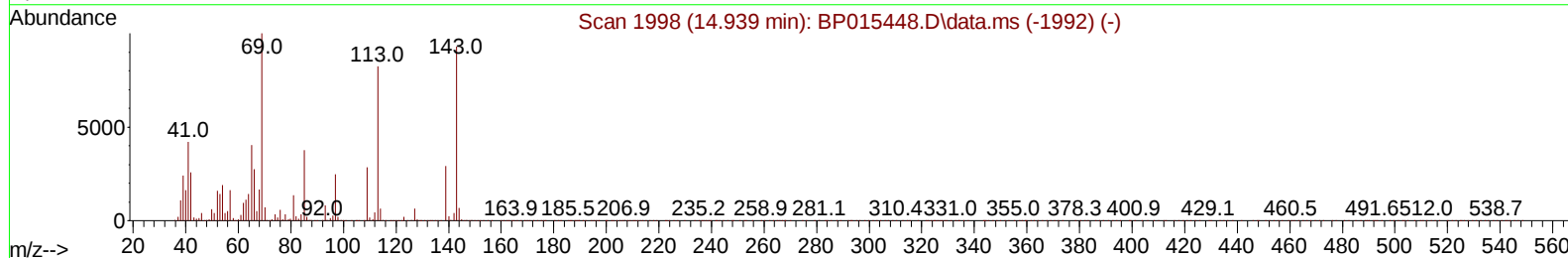
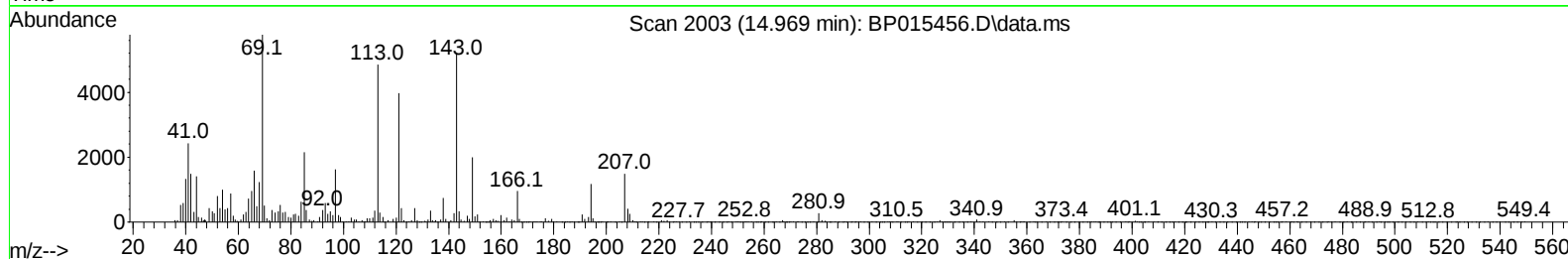
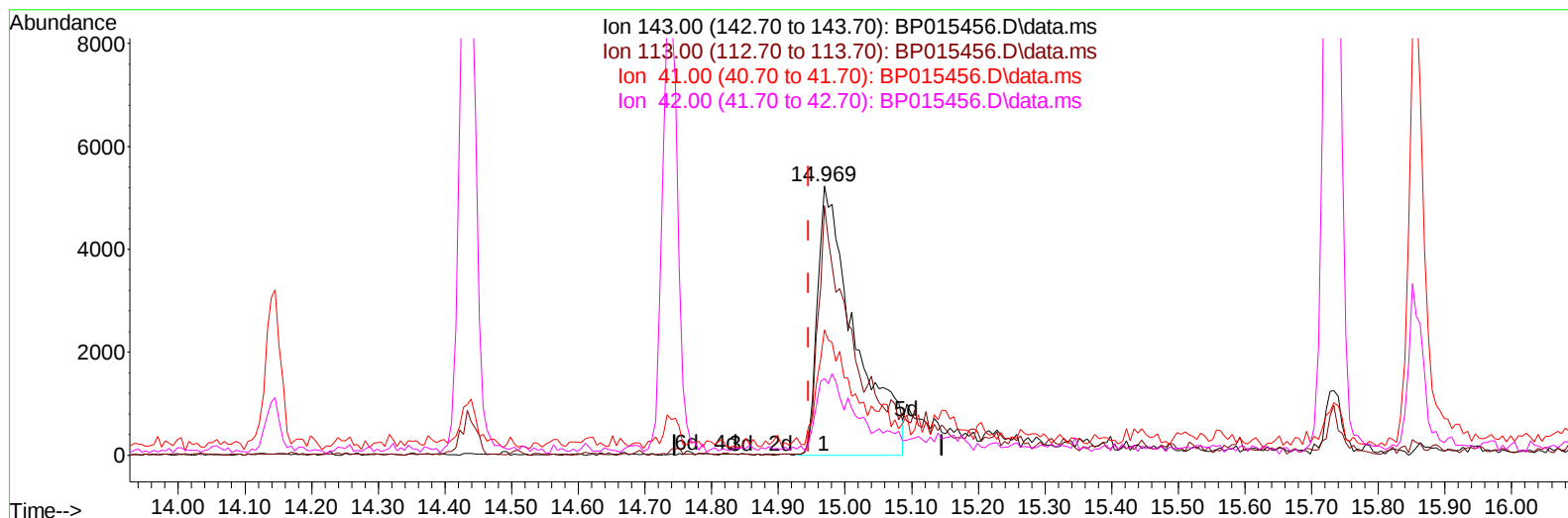
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TIC: BP015456.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.969min (+ 0.023) 3.68 ng/ul m

response 20292

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	86.60	92.88
41.00	42.20	46.52
42.00	29.40	28.54

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 Operator : MA/JU
 Sample : 03097-14
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EZEN7

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 06/12/2023
 Supervised By :mohammad ahmed 06/13/2023

Quant Time: Jun 09 23:25:33 2023
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 Quant Title : SVOA CALIBRATION
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.093	152	124438	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.922	136	571397	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.739	164	390883	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.492	188	921529	20.000	ng/ul	-0.01
79) Chrysene-d12	21.574	240	937096	20.000	ng/ul	-0.01
88) Perylene-d12	24.127	264	970098	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.411	96	7859	2.339	ng/uL	0.00
4) Pyridine-d5	3.858	84	109146	12.504	ng/ul	0.00
7) Phenol-d5	7.252	99	79730	6.955	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.416	67	242453	35.412	ng/ul	0.00
11) 2-Chlorophenol-d4	7.622	132	231844	27.894	ng/ul	0.00
15) 4-Methylphenol-d8	8.810	113	172735	18.359	ng/ul	0.00
21) Nitrobenzene-d5	9.275	128	160579	35.795	ng/ul	0.00
24) 2-Nitrophenol-d4	9.999	143	177597	34.666	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.540	165	302695	32.390	ng/ul	-0.01
31) 4-Chloroaniline-d4	11.063	131	394950	29.443	ng/ul	0.00
46) Dimethylphthalate-d6	14.145	166	1228242	39.848	ng/ul	0.00
49) Acenaphthylene-d8	14.434	160	1323806	39.277	ng/ul	0.00
54) 4-Nitrophenol-d4	14.969	143	20292m	3.676	ng/ul	0.02
60) Fluorene-d10	15.734	176	1033388	39.757	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.857	200	152100	26.054	ng/ul	0.00
73) Anthracene-d10	17.592	188	1705839	40.685	ng/ul	0.00
81) Pyrene-d10	19.816	212	2117723	42.224	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.963	264	2017275	41.443	ng/ul	-0.01

Target Compounds Qvalue

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

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