

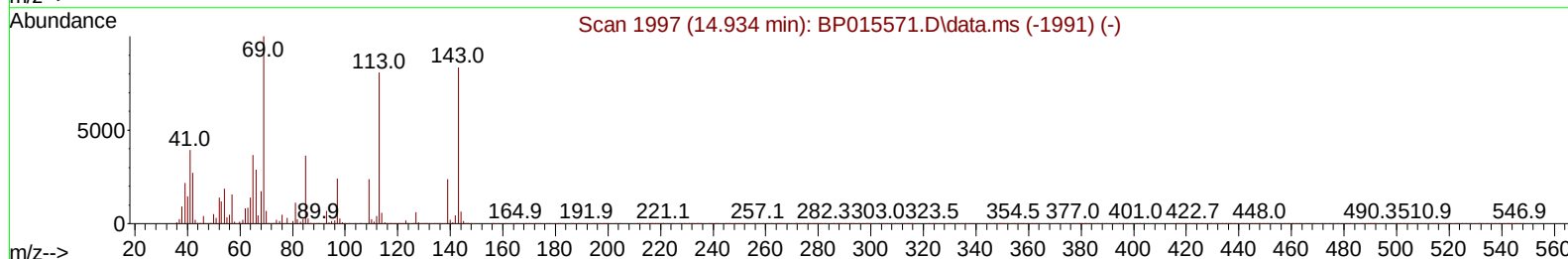
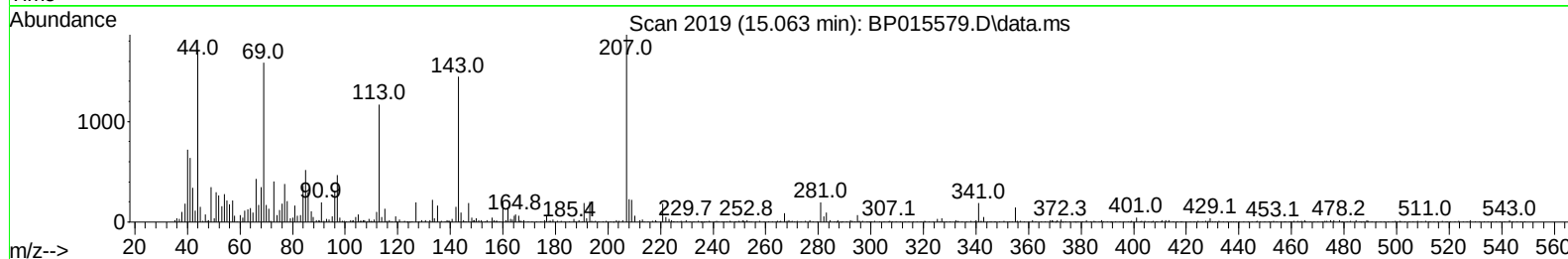
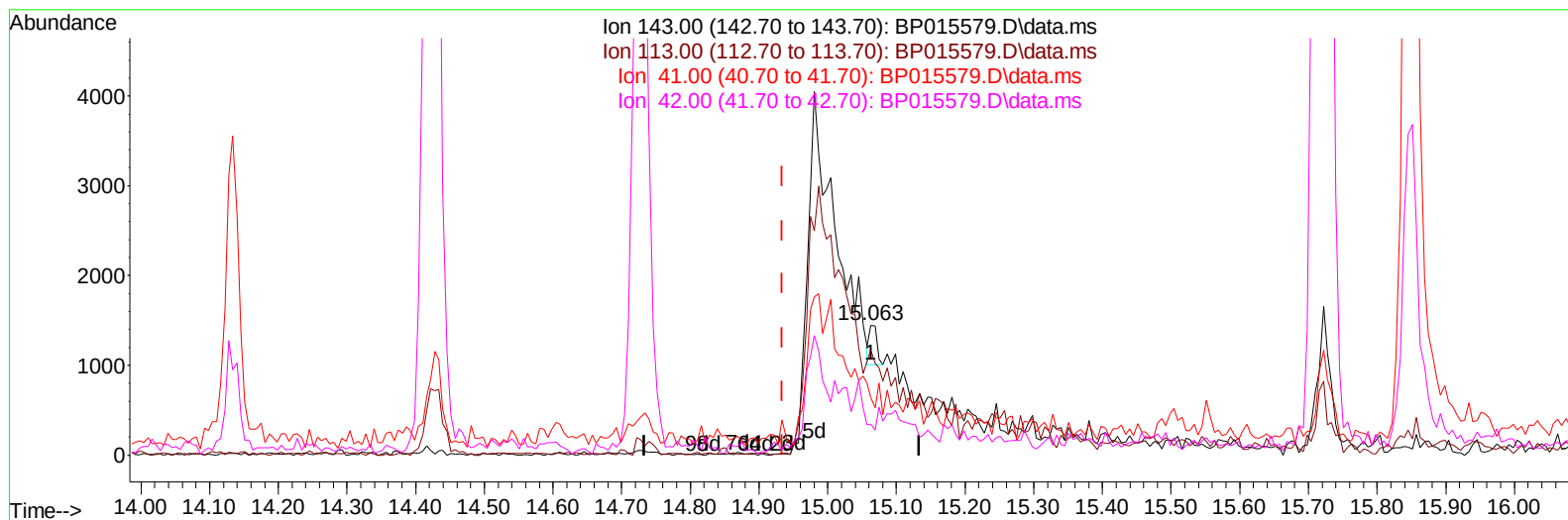
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061623\
 Data File : BP015579.D
 Acq On : 16 Jun 2023 14:06
 Operator : MA/JU
 Sample : 03220-09
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 CONR3

Manual IntegrationsAPPROVED

Quant Time: Jun 16 22:06:31 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jun 15 05:36:46 2023
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 06/19/2023
 Supervised By :mohammad ahmed 06/20/2023



TIC: BP015579.D\data.ms

(54) 4-Nitrophenol-d4 (S)

15.063min (+ 0.129) 0.07 ng/ul

response 330

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	86.60	80.84
41.00	42.20	44.12
42.00	29.40	23.58

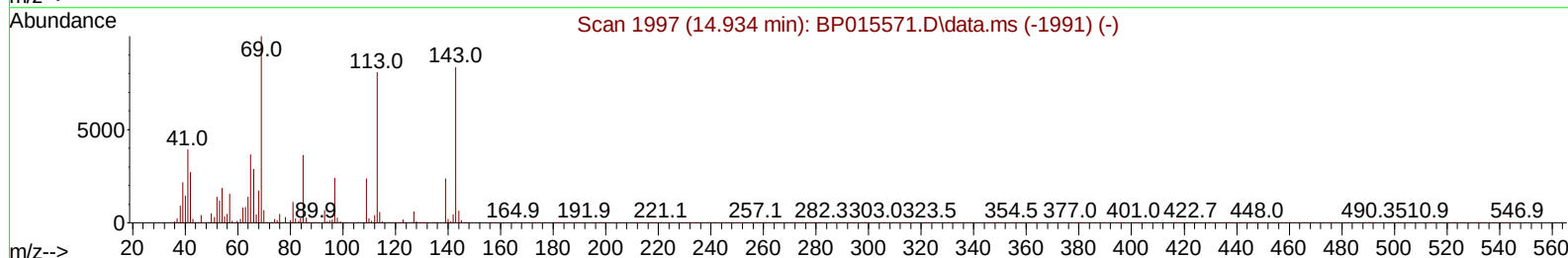
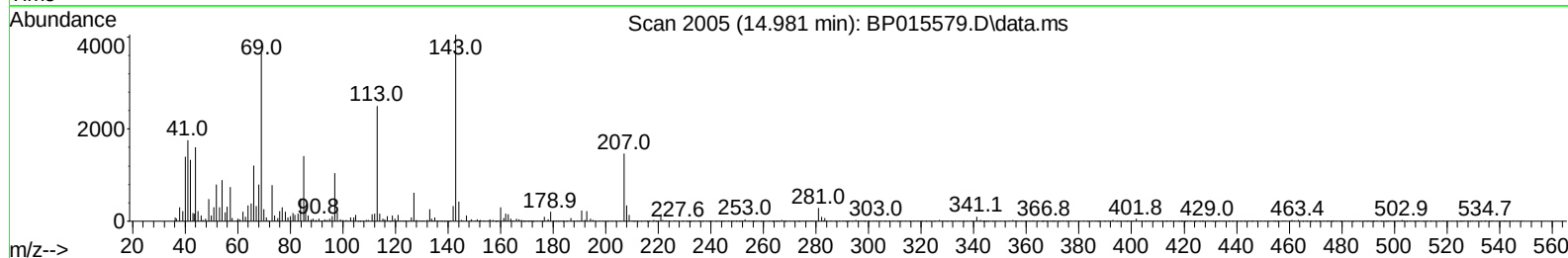
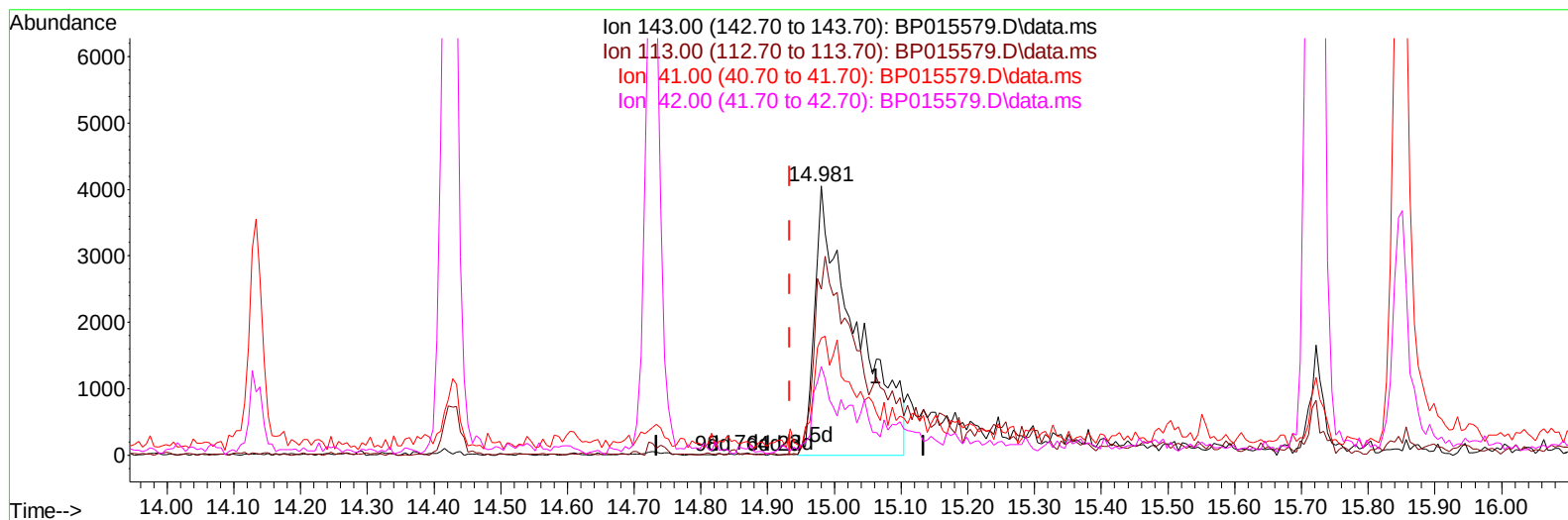
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(54) 4-Nitrophenol-d4 (S)

14.981min (+ 0.047) 3.67 ng/ul m

response 17097

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	86.60	61.73#
41.00	42.20	43.51
42.00	29.40	32.96

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.087	152	97312	20.000	ng/ul	0.00
20) Naphthalene-d8	10.910	136	459016	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.728	164	329928	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.486	188	778747	20.000	ng/ul	0.00
79) Chrysene-d12	21.568	240	798234	20.000	ng/ul	0.00
88) Perylene-d12	24.115	264	863753	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.405	96	13228	5.034	ng/uL	0.00
4) Pyridine-d5	3.858	84	72192	10.576	ng/ul	0.01
7) Phenol-d5	7.246	99	62403	6.961	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.405	67	226545	42.313	ng/ul	0.00
11) 2-Chlorophenol-d4	7.610	132	204400	31.447	ng/ul	0.00
15) 4-Methylphenol-d8	8.799	113	140587	19.107	ng/ul	0.00
21) Nitrobenzene-d5	9.263	128	150239	41.690	ng/ul	0.00
24) 2-Nitrophenol-d4	9.993	143	164979	40.088	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.534	165	267309	35.607	ng/ul	0.00
31) 4-Chloroaniline-d4	11.057	131	322560	29.934	ng/ul	0.00
46) Dimethylphthalate-d6	14.134	166	1234312	47.443	ng/ul	0.00
49) Acenaphthylene-d8	14.422	160	1283762	45.126	ng/ul	0.00
54) 4-Nitrophenol-d4	14.981	143	17097m	3.670	ng/ul	0.05
60) Fluorene-d10	15.722	176	1031455	47.014	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.851	200	181673	36.825	ng/ul	0.00
73) Anthracene-d10	17.586	188	1714035	48.375	ng/ul	0.00
81) Pyrene-d10	19.810	212	2135257	49.979	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.951	264	2112616	48.746	ng/ul	0.00

Target Compounds Qvalue

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

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