

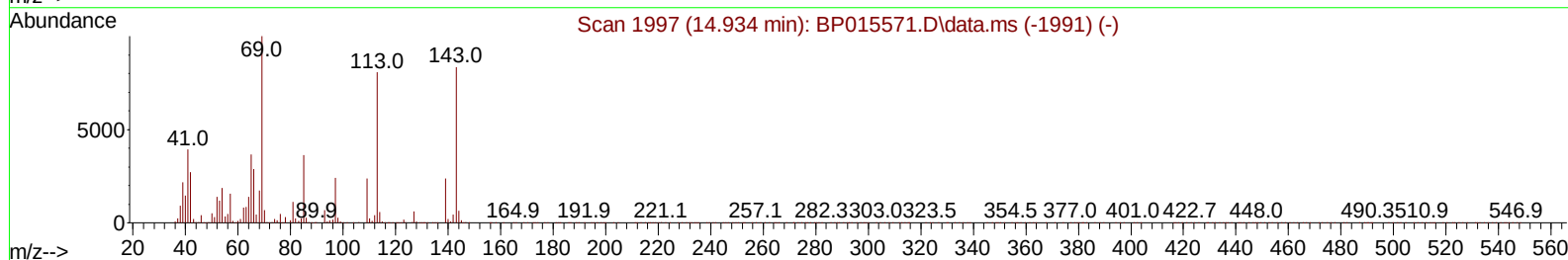
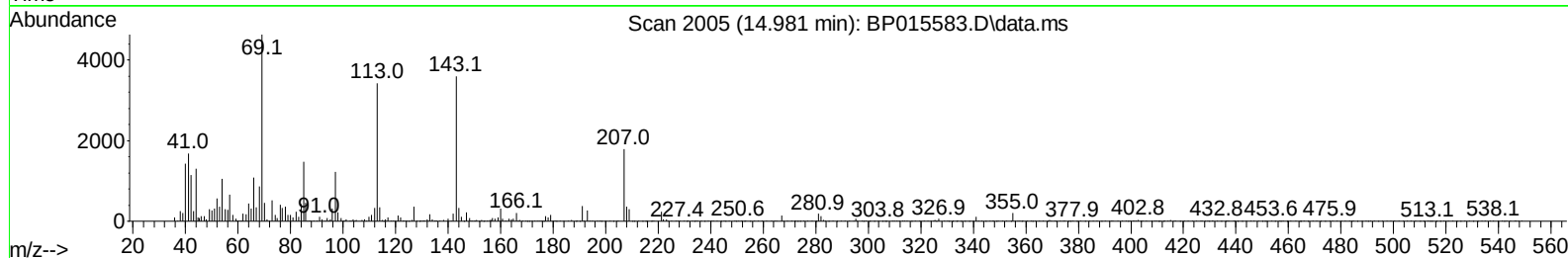
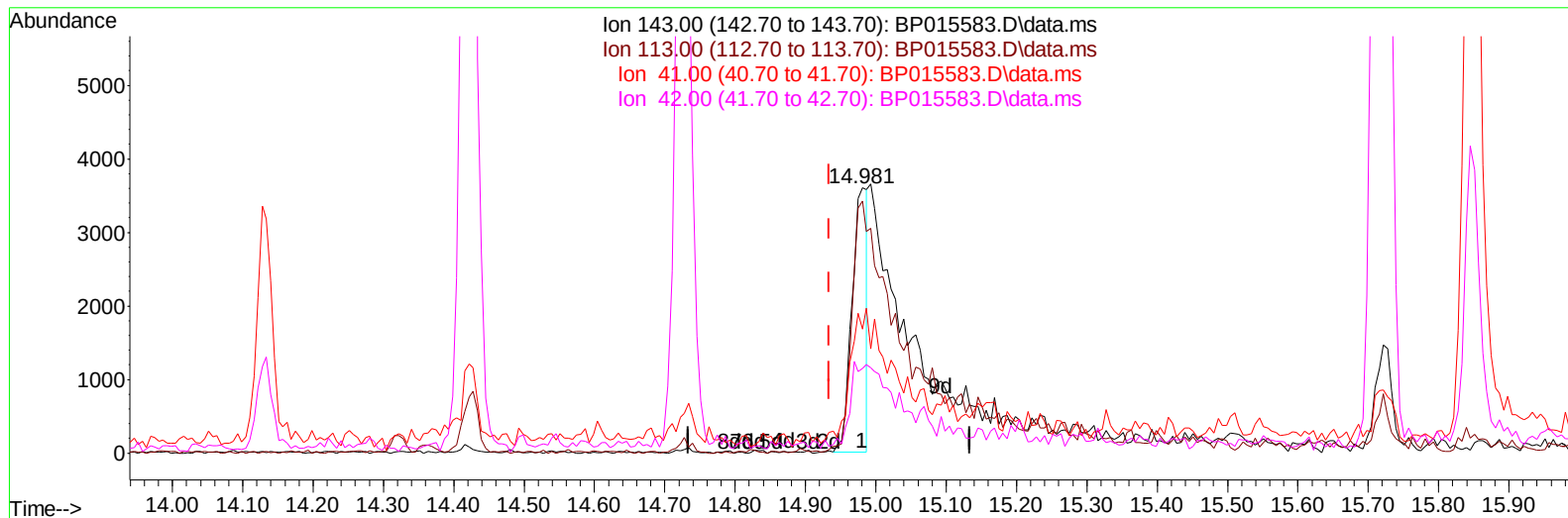
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061623\
 Data File : BP015583.D
 Acq On : 16 Jun 2023 16:21
 Operator : MA/JU
 Sample : 03220-01
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 CONK8

Manual IntegrationsAPPROVED

Quant Time: Jun 16 22:07:15 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: BP015583.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.981min (+ 0.047) 0.95 ng/ul

response 5479

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	86.60	95.09
41.00	42.20	46.78
42.00	29.40	31.96

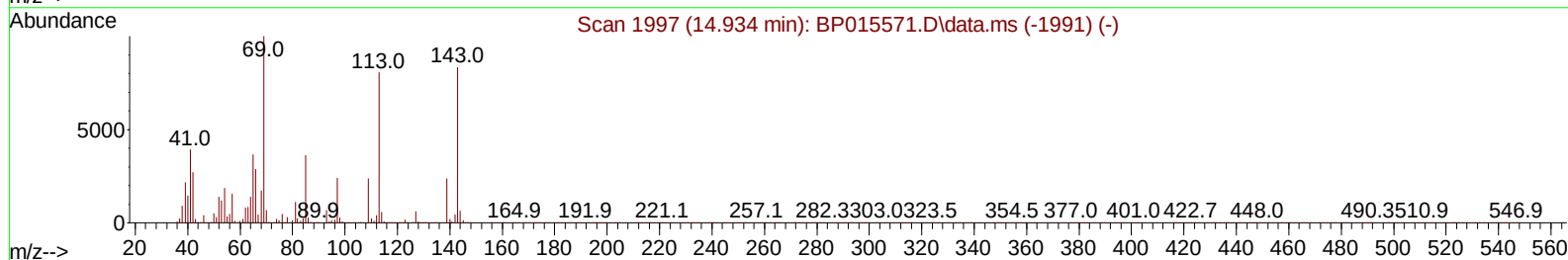
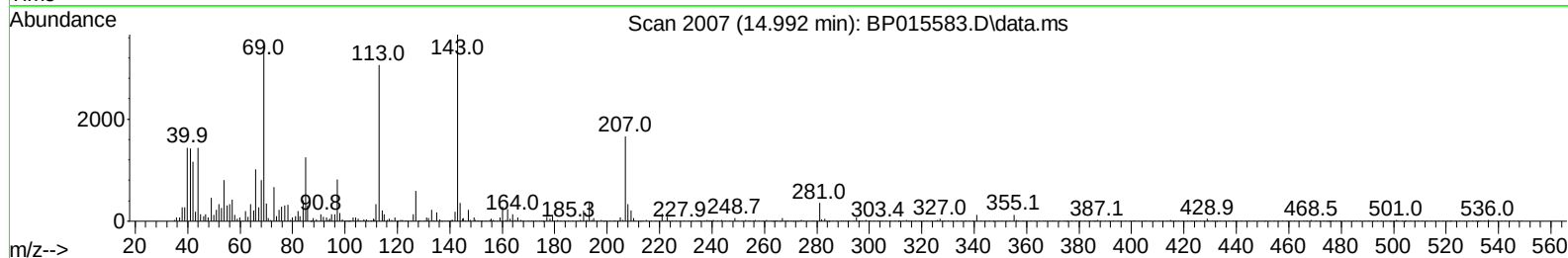
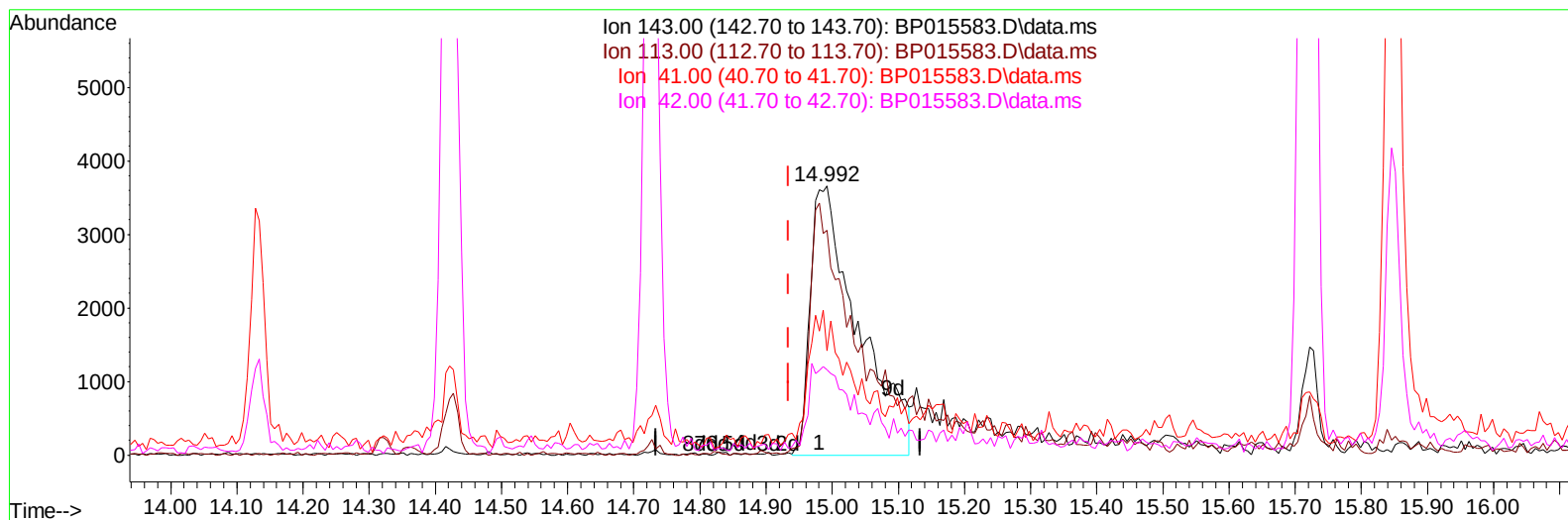
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(54) 4-Nitrophenol-d4 (S)

14.992min (+ 0.059) 3.20 ng/ul m

response 18369

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	86.60	83.68
41.00	42.20	38.83
42.00	29.40	31.83

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.081	152	124120	20.000	ng/ul	0.00
20) Naphthalene-d8	10.910	136	555326	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.728	164	406910	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.486	188	979615	20.000	ng/ul	0.00
79) Chrysene-d12	21.568	240	901500	20.000	ng/ul	0.00
88) Perylene-d12	24.110	264	972797	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.405	96	17764	5.300	ng/uL	0.00
4) Pyridine-d5	3.852	84	85943	9.871	ng/ul	0.00
7) Phenol-d5	7.246	99	68682	6.006	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.405	67	237506	34.779	ng/ul	0.00
11) 2-Chlorophenol-d4	7.610	132	215891	26.041	ng/ul	0.00
15) 4-Methylphenol-d8	8.799	113	144910	15.441	ng/ul	0.00
21) Nitrobenzene-d5	9.263	128	149391	34.265	ng/ul	0.00
24) 2-Nitrophenol-d4	9.987	143	164815	33.103	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.534	165	259024	28.519	ng/ul	0.00
31) 4-Chloroaniline-d4	11.051	131	366795	28.136	ng/ul	0.00
46) Dimethylphthalate-d6	14.134	166	1287751	40.133	ng/ul	0.00
49) Acenaphthylene-d8	14.422	160	1290956	36.794	ng/ul	0.00
54) 4-Nitrophenol-d4	14.992	143	18369m	3.197	ng/ul	0.06
60) Fluorene-d10	15.722	176	1072992	39.655	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.845	200	200286	32.273	ng/ul	0.00
73) Anthracene-d10	17.586	188	1823300	40.908	ng/ul	0.00
81) Pyrene-d10	19.810	212	2206193	45.725	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.945	264	2059331	42.190	ng/ul	0.00

Target Compounds Qvalue

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

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