

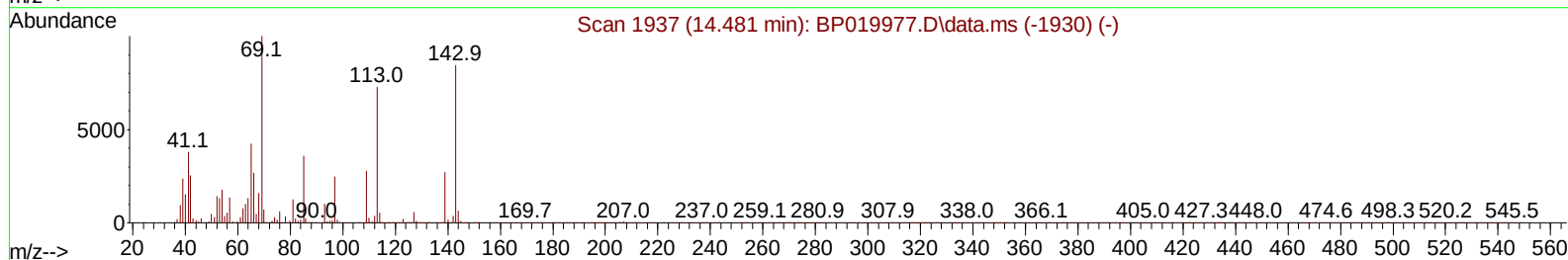
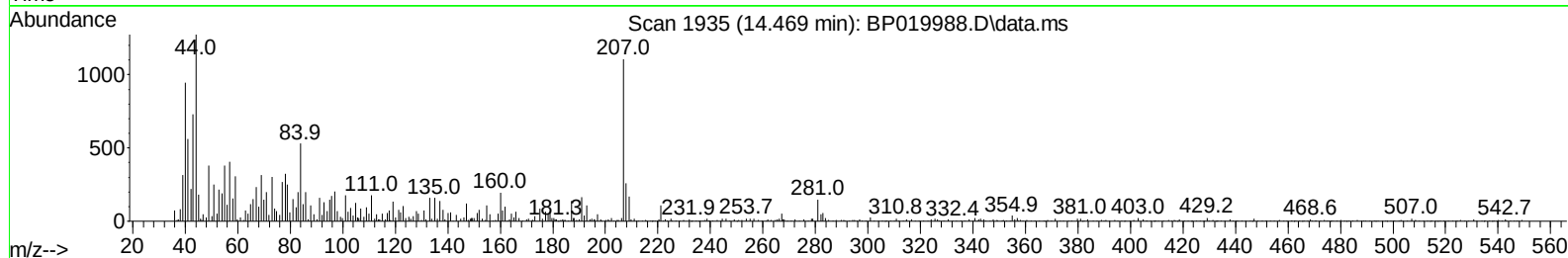
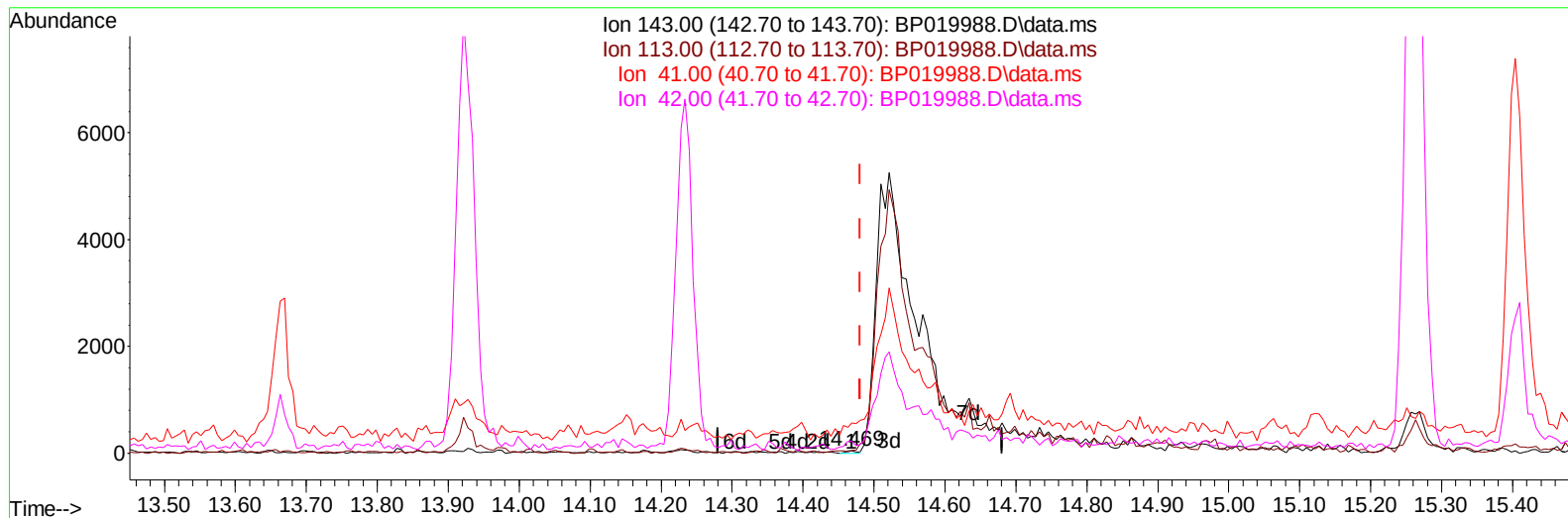
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032524\
Data File : BP019988.D
Acq On : 25 Mar 2024 19:35
Operator : MA/JU
Sample : P1813-13
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E0LY1

Manual IntegrationsAPPROVED

Quant Time: Mar 25 23:12:51 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032224.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Mar 25 10:46:59 2024
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 03/26/2024
Supervised By :mohammad ahmed 03/27/2024



TIC: BP019988.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.469min (-0.012) 0.01 ng/u1

response 50

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	81.50	133.33#
41.00	40.90	1246.67#
42.00	28.50	493.33#

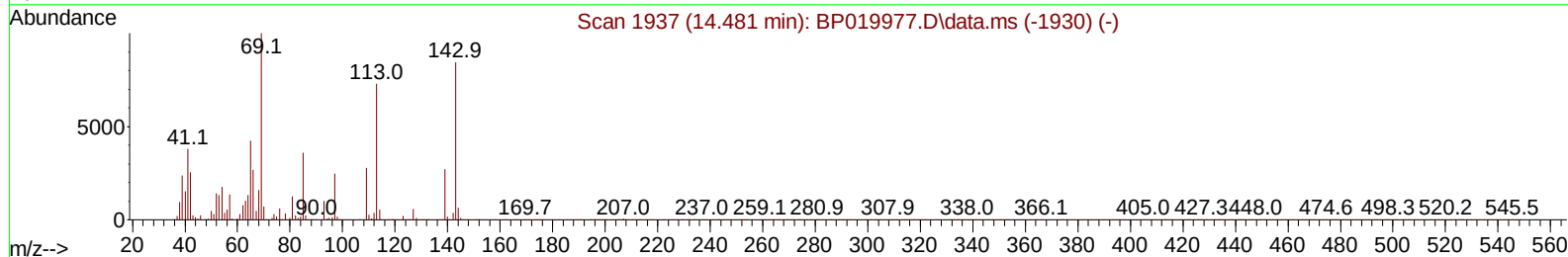
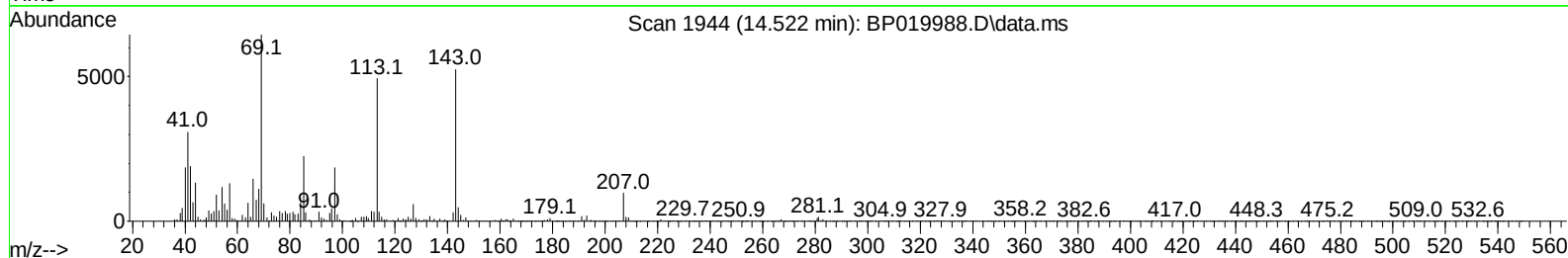
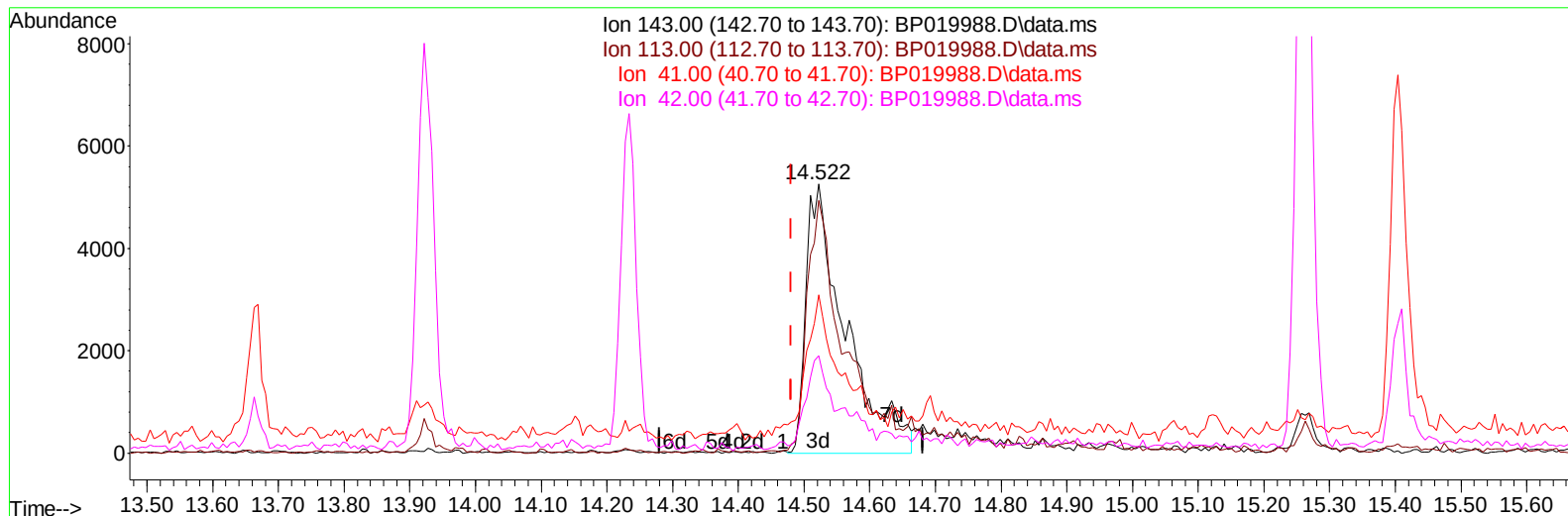
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TIC: BP019988.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.522min (+ 0.041) 5.96 ng/ul m

response 21794

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	81.50	94.01
41.00	40.90	58.84#
42.00	28.50	36.10#

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Instrument :
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Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 03/26/2024
 Supervised By :mohammad ahmed 03/27/2024

Quant Time: Mar 25 23:58:32 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032224.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Mar 25 10:46:59 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.622	152	113511	20.000	ng/ul	0.00
20) Naphthalene-d8	10.369	136	494025	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.234	164	331849	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.063	188	644762	20.000	ng/ul	-0.01
79) Chrysene-d12	21.533	240	479262	20.000	ng/ul	-0.01
88) Perylene-d12	24.804	264	533407	20.000	ng/ul	-0.03
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.240	96	13441	5.188	ng/uL	0.00
4) Pyridine-d5	3.652	84	31555	4.034	ng/ul	0.01
7) Phenol-d5	6.811	99	71685	7.488	ng/ul	0.01
9) Bis-(2-Chloroethyl)eth...	6.975	67	250137	40.104	ng/ul	0.00
11) 2-Chlorophenol-d4	7.158	132	192446	27.991	ng/ul	0.00
15) 4-Methylphenol-d8	8.316	113	140953	19.030	ng/ul	0.01
21) Nitrobenzene-d5	8.769	128	143624	37.930	ng/ul	0.01
24) 2-Nitrophenol-d4	9.475	143	152731	36.050	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.004	165	247088	31.934	ng/ul	0.00
31) 4-Chloroaniline-d4	10.528	131	197134	19.544	ng/ul	0.01
46) Dimethylphthalate-d6	13.669	166	1017857	43.747	ng/ul	0.02
49) Acenaphthylene-d8	13.922	160	1102812	40.576	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.522	143	21794m	5.958	ng/ul	0.04
60) Fluorene-d10	15.263	176	847572	42.764	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.404	200	139200	36.067	ng/ul	-0.02
73) Anthracene-d10	17.169	188	1283137	44.958	ng/ul	0.00
81) Pyrene-d10	19.557	212	1320132	46.948	ng/ul	-0.01
92) Benzo(a)pyrene-d12	24.592	264	1213125	47.340	ng/ul	-0.02

Target Compounds Qvalue

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

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