

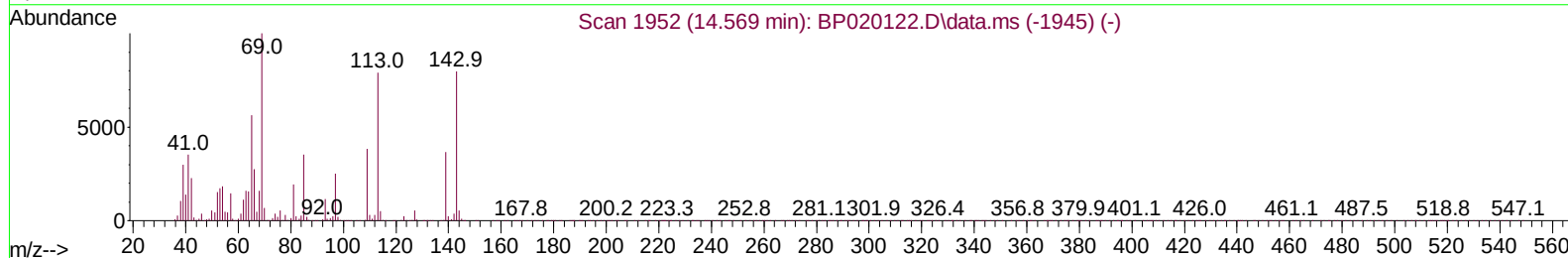
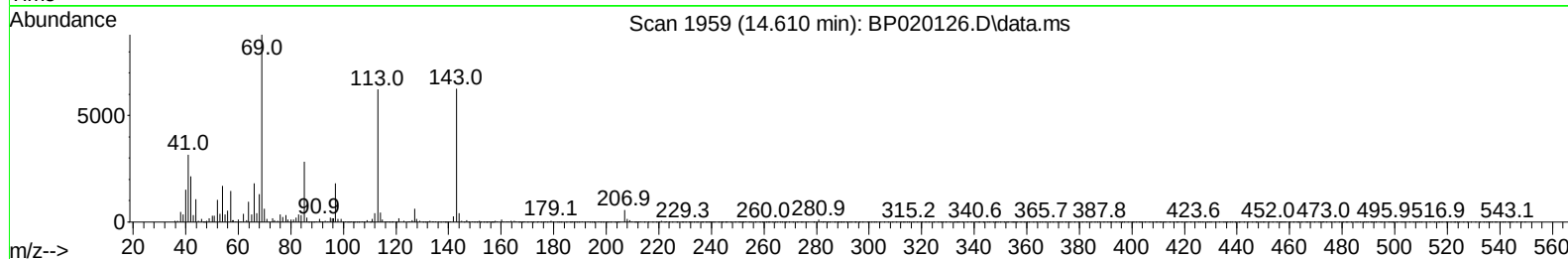
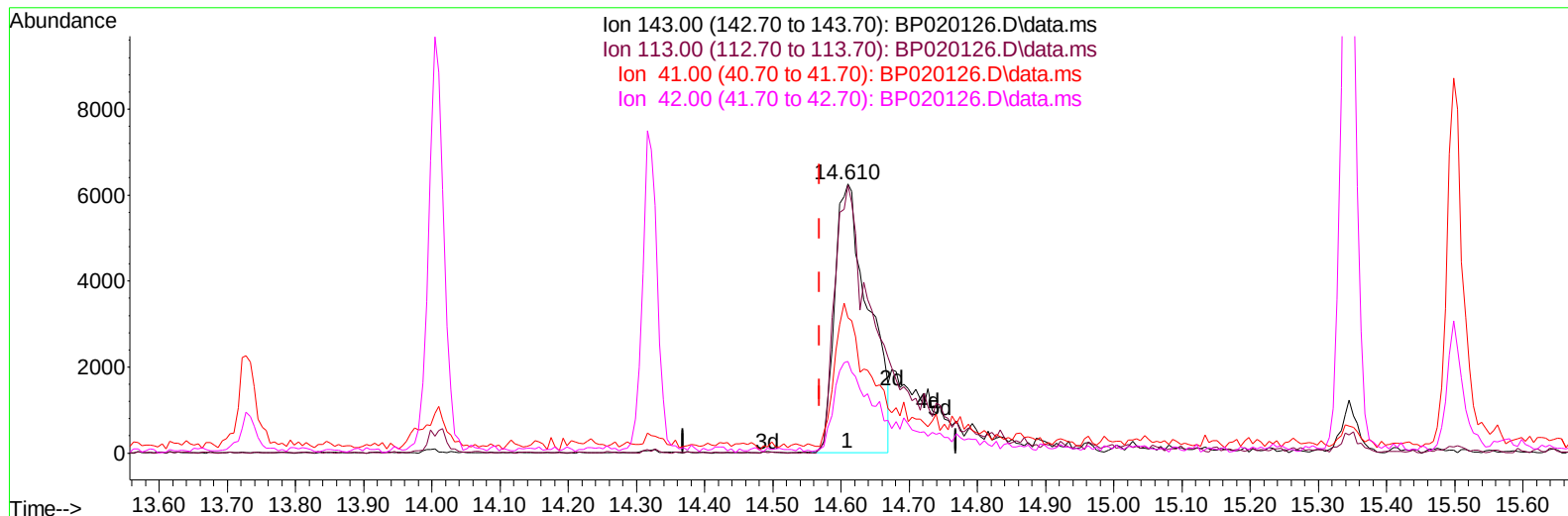
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP043024\
Data File : BP020126.D
Acq On : 30 Apr 2024 19:36
Operator : MA/JU
Sample : P2220-01 10X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E0808

Manual IntegrationsAPPROVED

Quant Time: May 01 01:14:48 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042424.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed May 01 00:58:38 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 05/03/2024
Supervised By :mohammad ahmed 05/03/2024



TIC: BP020126.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.610min (+ 0.041) 5.77 ng/u1

response 21499

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	101.00	99.50
41.00	44.80	50.42
42.00	29.30	34.14

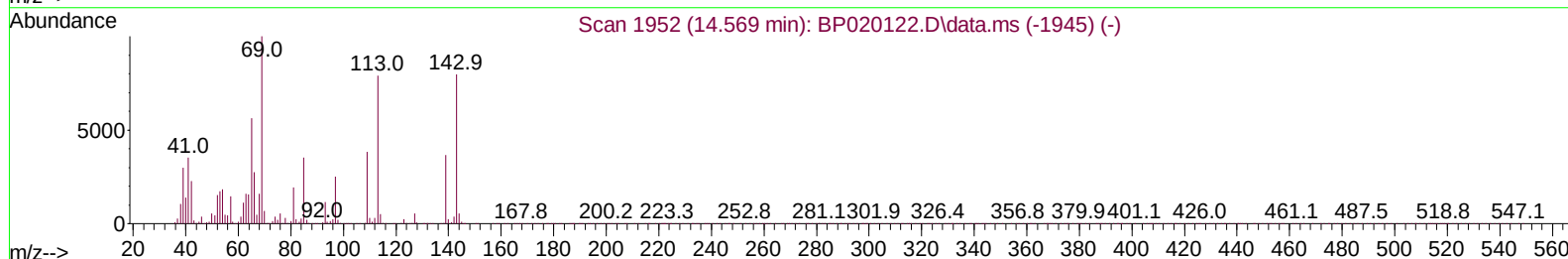
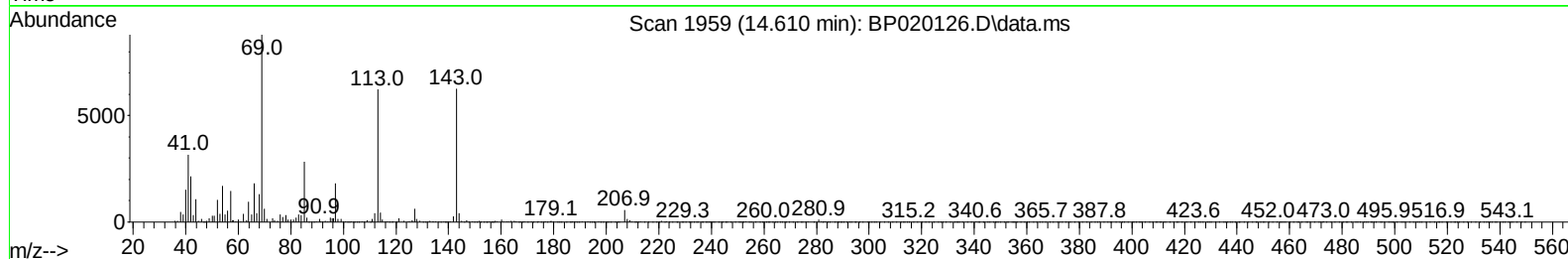
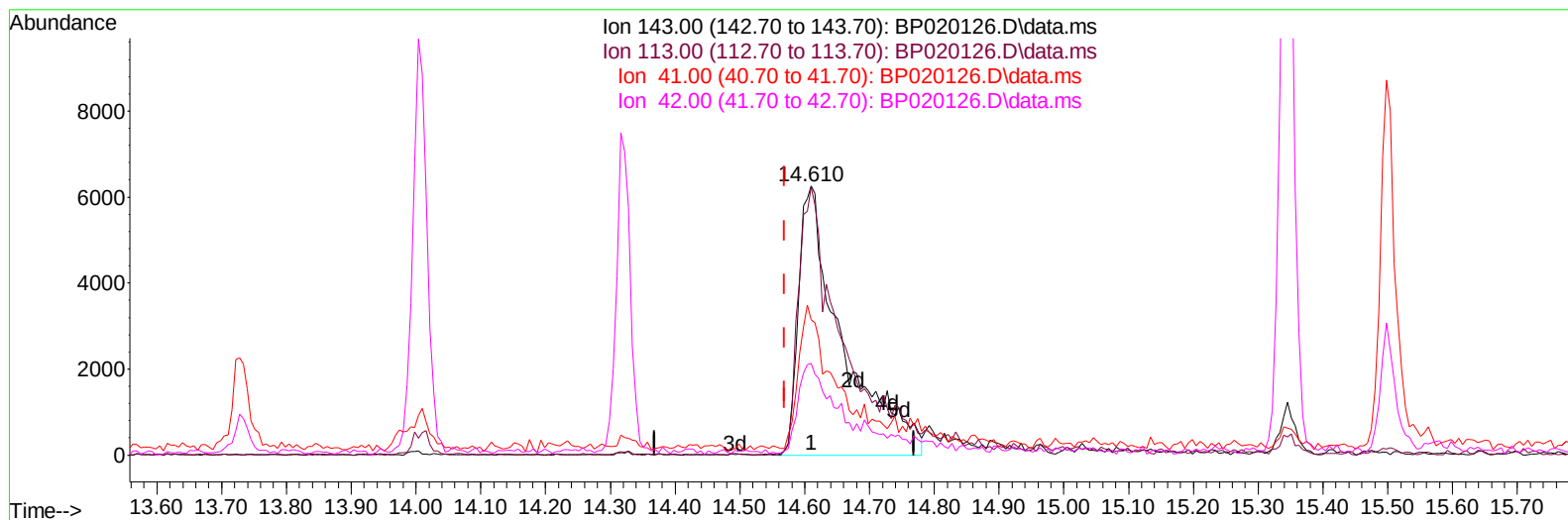
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TIC: BP020126.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.610min (+ 0.041) 7.89 ng/ul m

response 29378

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	101.00	99.50
41.00	44.80	50.42
42.00	29.30	34.14

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 QLast Update : Wed May 01 00:58:38 2024
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Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)

Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.634	152		104129	20.000	ng/ul	0.00
20) Naphthalene-d8	10.410	136		438907	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.322	164		320807	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.163	188		720314	20.000	ng/ul	0.00
79) Chrysene-d12	21.651	240		569762	20.000	ng/ul	-0.02
88) Perylene-d12	25.027	264		556069	20.000	ng/ul	-0.02
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.222	96		6577	2.638	ng/uL	0.00
4) Pyridine-d5	3.628	84		71610	9.749	ng/ul	0.01
7) Phenol-d5	6.822	99		81268	9.204	ng/ul	0.01
9) Bis-(2-Chloroethyl)eth...	6.987	67		226481	39.816	ng/ul	0.00
11) 2-Chlorophenol-d4	7.169	132		187066	30.175	ng/ul	0.00
15) 4-Methylphenol-d8	8.340	113		143838	21.397	ng/ul	0.00
21) Nitrobenzene-d5	8.793	128		126075	37.405	ng/ul	0.00
24) 2-Nitrophenol-d4	9.510	143		136303	36.439	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.040	165		227955	32.382	ng/ul	0.00
31) 4-Chloroaniline-d4	10.569	131		287708	31.341	ng/ul	0.01
46) Dimethylphthalate-d6	13.734	166		945096	41.247	ng/ul	0.00
49) Acenaphthylene-d8	14.010	160		1012782	38.780	ng/ul	0.00
54) 4-Nitrophenol-d4	14.610	143		29378m	7.888	ng/ul	0.04
60) Fluorene-d10	15.345	176		796899	39.983	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.498	200		155590	33.782	ng/ul	0.00
73) Anthracene-d10	17.275	188		1333630	41.700	ng/ul	0.01
81) Pyrene-d10	19.663	212		1521218	50.185	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.798	264		1166379	42.989	ng/ul	-0.01

Target Compounds Qvalue

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

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