

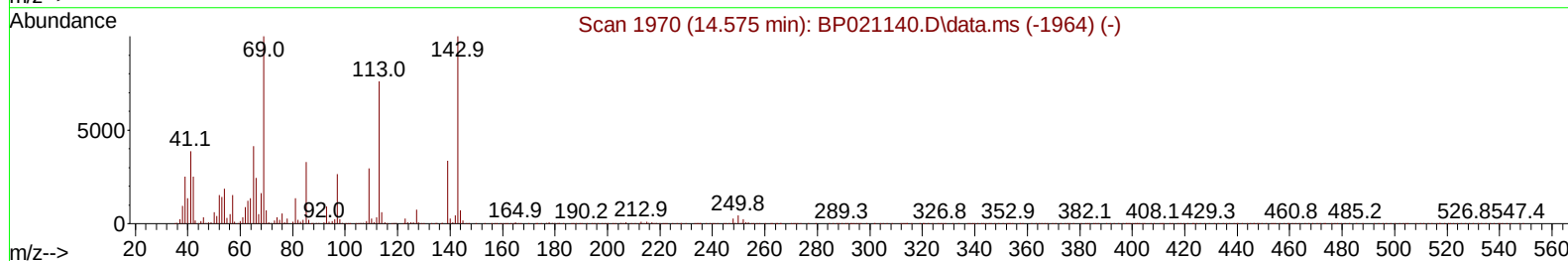
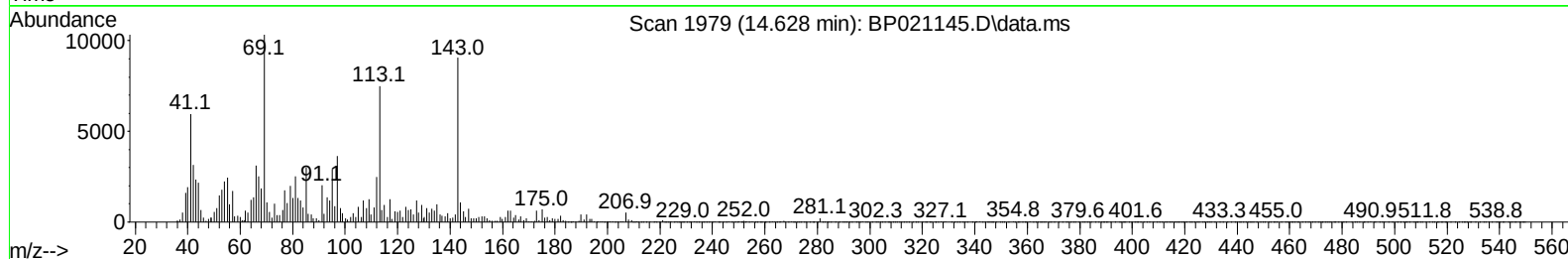
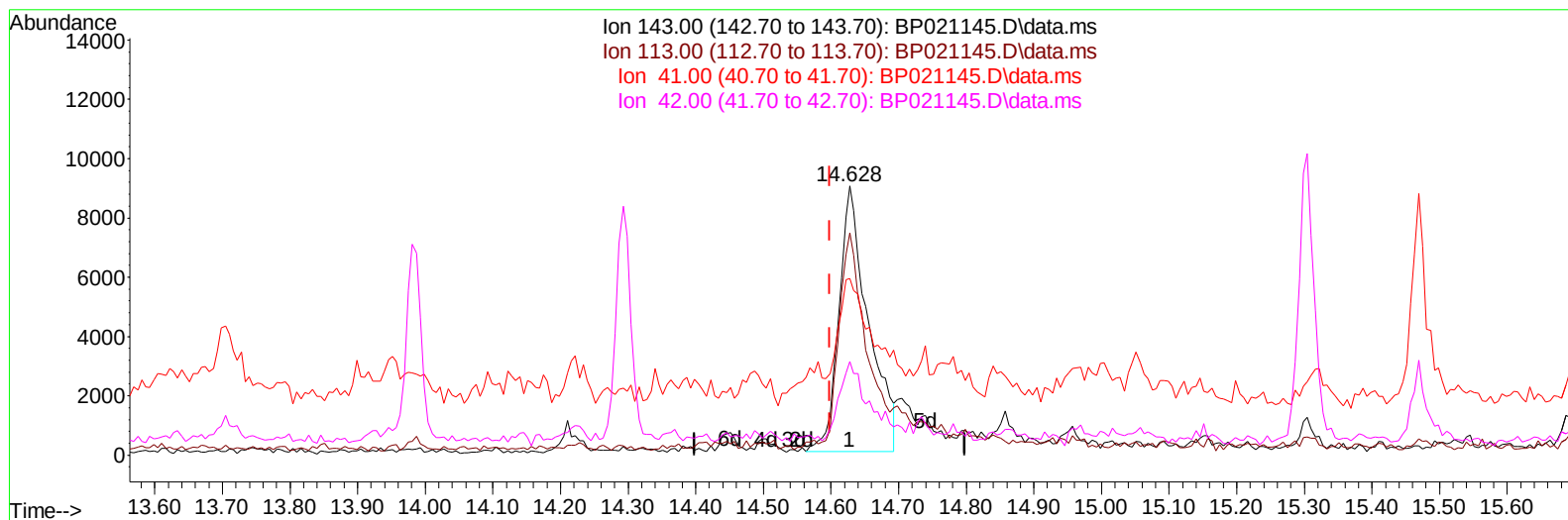
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072524\
Data File : BP021145.D
Acq On : 25 Jul 2024 13:06
Operator : MA/JU
Sample : P3244-13
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
YDZK7

Manual IntegrationsAPPROVED

Quant Time: Jul 25 22:38:40 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Jul 25 00:15:24 2024
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 07/26/2024
Supervised By :mohammad ahmed 07/27/2024



TIC: BP021145.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.628min (+ 0.029) 6.01 ng/ul

response 27030

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	74.70	82.75
41.00	37.90	65.67#
42.00	27.10	34.91#

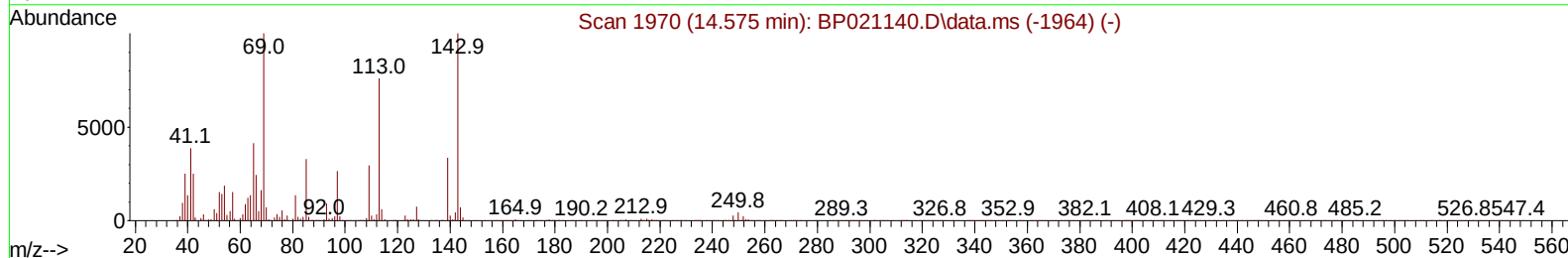
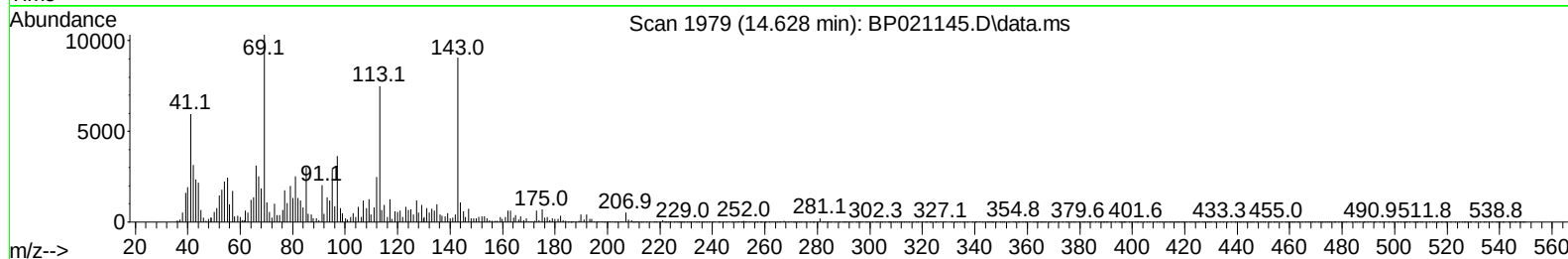
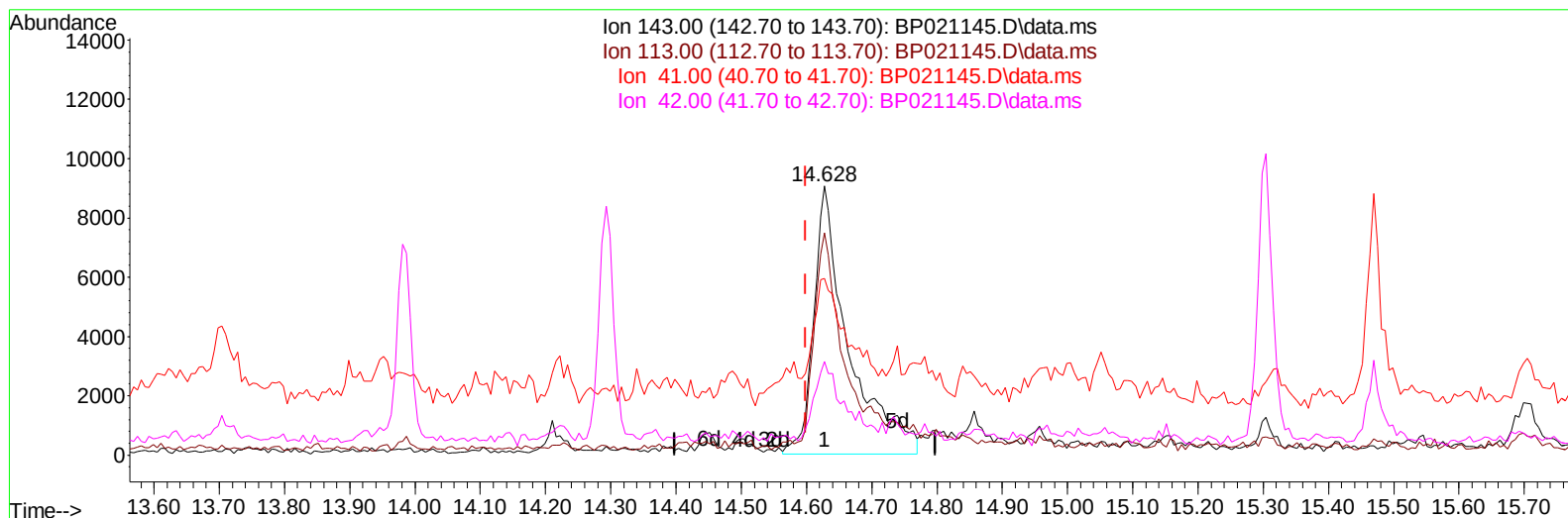
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(54) 4-Nitrophenol-d4 (S)

14.628min (+ 0.029) 7.40 ng/ul m

response 33283

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	74.70	82.75
41.00	37.90	65.67#
42.00	27.10	34.91#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.658	152	189017	20.000	ng/ul	0.01
20) Naphthalene-d8	10.416	136	699232	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.293	164	434663	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.104	188	892272	20.000	ng/ul	0.00
79) Chrysene-d12	21.551	240	863269	20.000	ng/ul	0.00
88) Perylene-d12	24.857	264	1089558	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.181	96	16719	4.025	ng/uL	0.00
4) Pyridine-d5	3.611	84	61164	5.086	ng/ul	0.01
7) Phenol-d5	6.875	99	87270	5.663	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.005	67	191054	20.458	ng/ul	0.01
11) 2-Chlorophenol-d4	7.205	132	240653	18.954	ng/ul	0.00
15) 4-Methylphenol-d8	8.387	113	161689	12.633	ng/ul	0.00
21) Nitrobenzene-d5	8.816	128	122353	22.529	ng/ul	0.01
24) 2-Nitrophenol-d4	9.528	143	144236	23.204	ng/ul	0.01
28) 2,4-Dichlorophenol-d3	10.081	165	253831	22.582	ng/ul	0.00
31) 4-Chloroaniline-d4	10.587	131	147552	9.909	ng/ul	0.02
46) Dimethylphthalate-d6	13.704	166	804752	25.469	ng/ul	0.00
49) Acenaphthylene-d8	13.981	160	853813	24.065	ng/ul	0.00
54) 4-Nitrophenol-d4	14.628	143	33283m	7.403	ng/ul	0.03
60) Fluorene-d10	15.304	176	685306	26.437	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.469	200	121404	22.488	ng/ul	0.00
73) Anthracene-d10	17.210	188	1104187	27.862	ng/ul	0.00
81) Pyrene-d10	19.598	212	1350558	25.305	ng/ul	0.01
92) Benzo(a)pyrene-d12	24.633	264	1552876	28.898	ng/ul	-0.02

Target Compounds Qvalue

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

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