

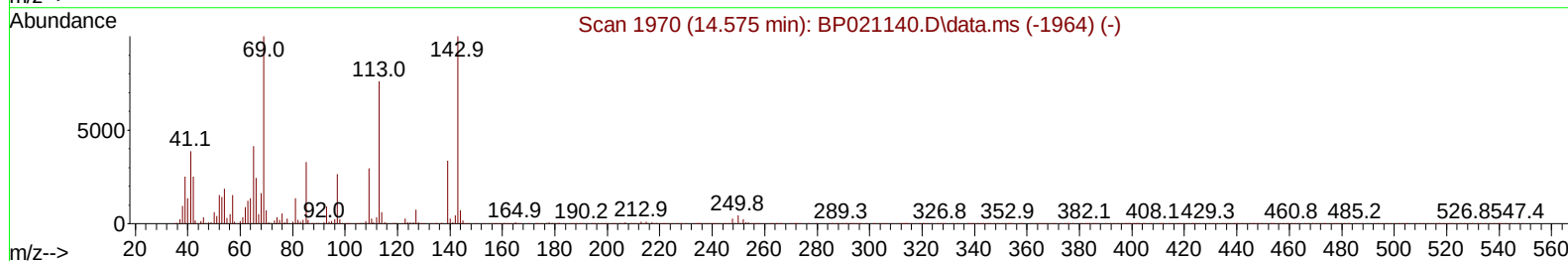
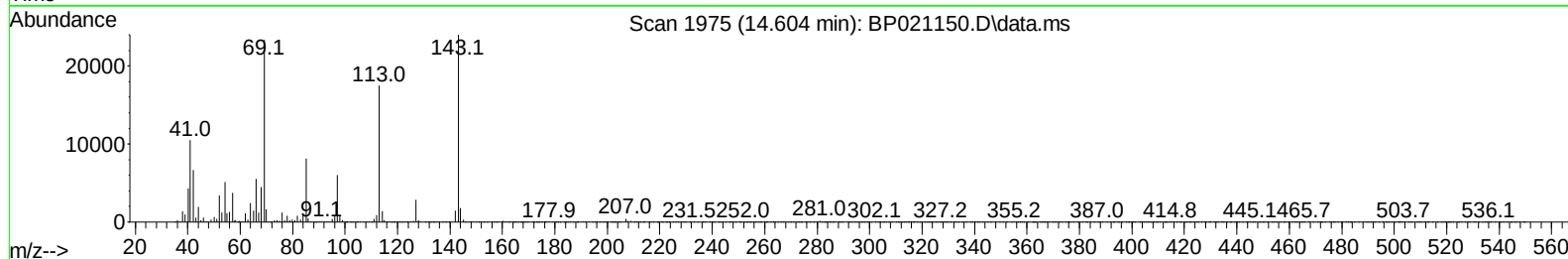
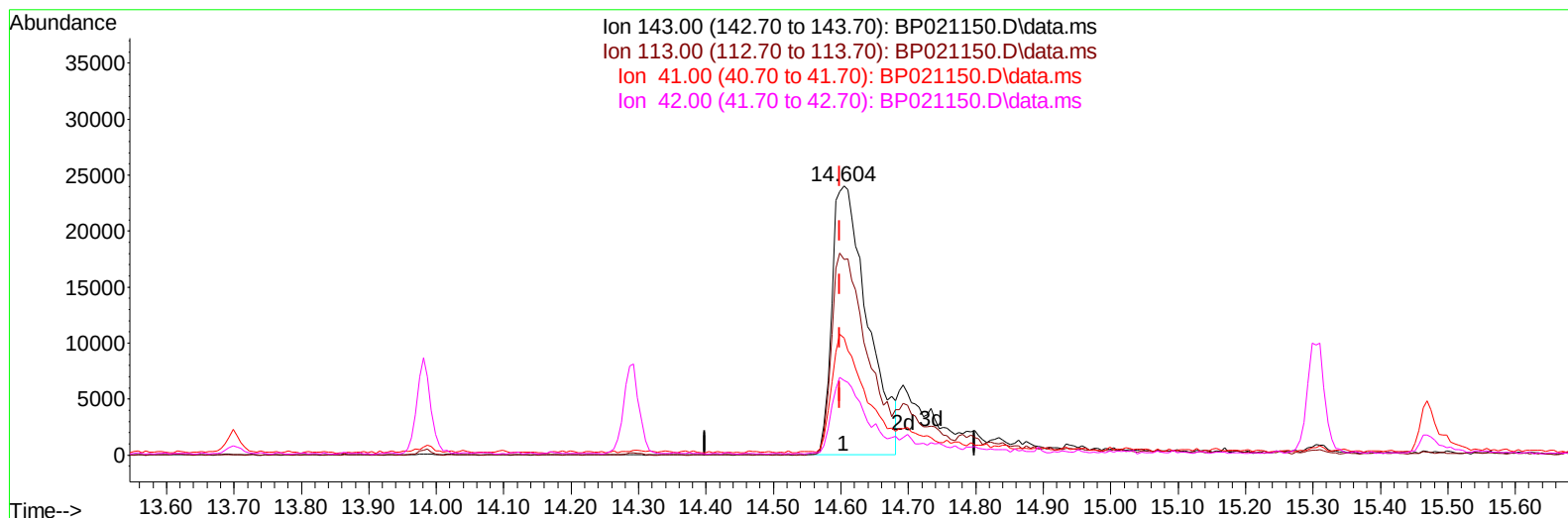
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072524\
 Data File : BP021150.D
 Acq On : 25 Jul 2024 16:39
 Operator : MA/JU
 Sample : PB162155BL
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK155

Manual IntegrationsAPPROVED

Quant Time: Jul 25 22:39:58 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: BP021150.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.604min (+ 0.006) 17.58 ng/ul

response 87522

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	74.70	72.78
41.00	37.90	43.60
42.00	27.10	27.66

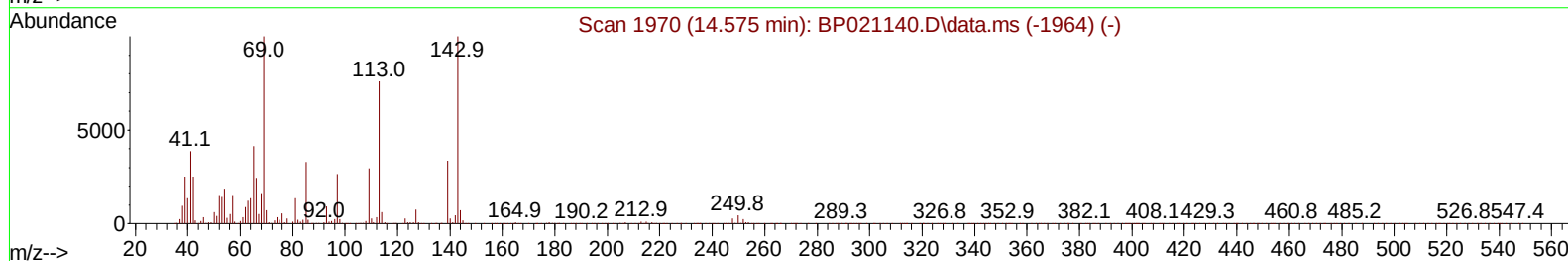
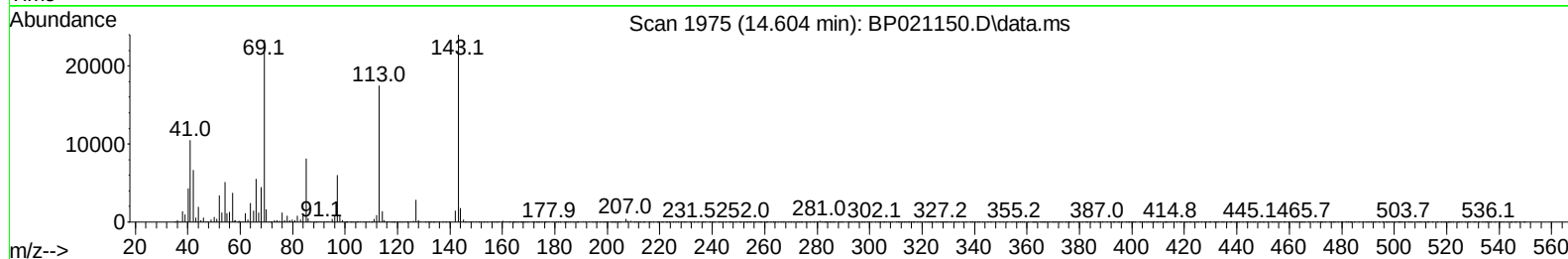
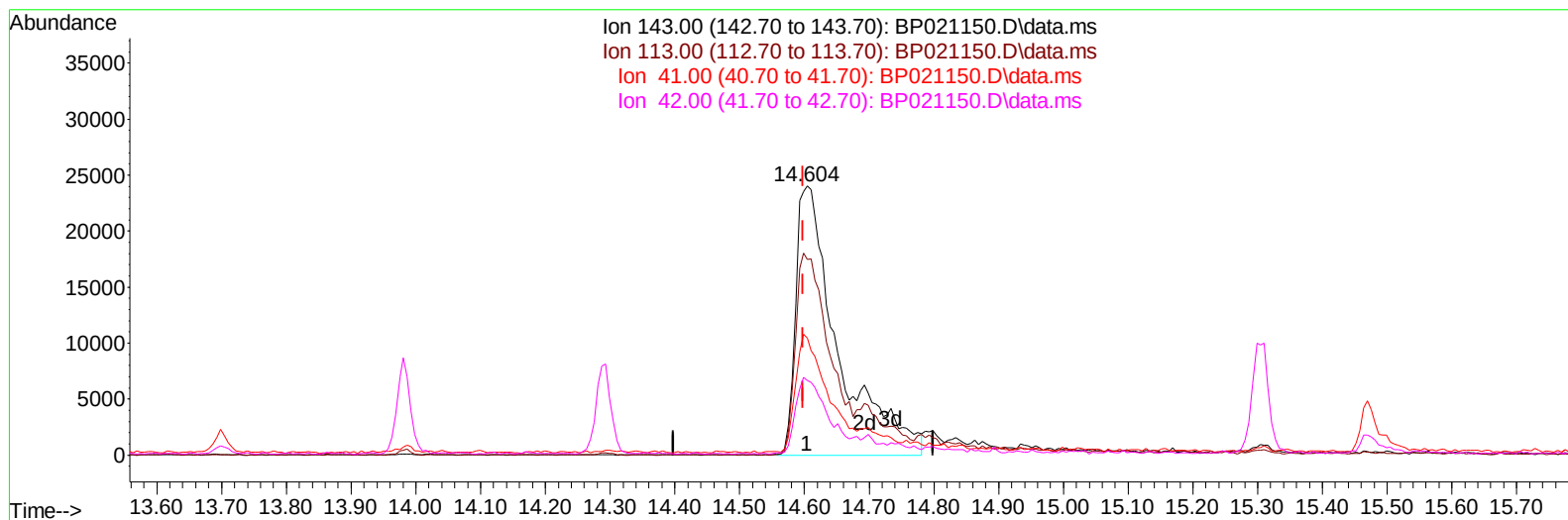
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TIC: BP021150.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.604min (+ 0.006) 21.90 ng/ul m

response 109011

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	74.70	72.78
41.00	37.90	43.60
42.00	27.10	27.66

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Quant Time: Jul 25 23:07:52 2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.652	152	206867	20.000	ng/ul	0.00
20) Naphthalene-d8	10.416	136	785556	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.293	164	481256	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.104	188	928786	20.000	ng/ul	0.00
79) Chrysene-d12	21.563	240	725927	20.000	ng/ul	0.00
88) Perylene-d12	24.863	264	962172	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.181	96	24041	5.288	ng/uL	0.00
4) Pyridine-d5	3.599	84	316046	24.014	ng/ul	0.00
7) Phenol-d5	6.864	99	405017	24.015	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.999	67	245609	24.030	ng/ul	0.00
11) 2-Chlorophenol-d4	7.199	132	341007	24.541	ng/ul	0.00
15) 4-Methylphenol-d8	8.381	113	330319	23.582	ng/ul	0.00
21) Nitrobenzene-d5	8.811	128	161695	26.501	ng/ul	0.00
24) 2-Nitrophenol-d4	9.522	143	186116	26.651	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.069	165	301049	23.840	ng/ul	0.00
31) 4-Chloroaniline-d4	10.581	131	433303	25.901	ng/ul	0.01
46) Dimethylphthalate-d6	13.699	166	917952	26.239	ng/ul	0.00
49) Acenaphthylene-d8	13.981	160	1050206	26.734	ng/ul	0.00
54) 4-Nitrophenol-d4	14.604	143	109011m	21.900	ng/ul	0.00
60) Fluorene-d10	15.304	176	776730	27.063	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.469	200	107515	19.132	ng/ul	0.00
73) Anthracene-d10	17.210	188	1166604	28.280	ng/ul	0.00
81) Pyrene-d10	19.598	212	1241369	27.660	ng/ul	0.01
92) Benzo(a)pyrene-d12	24.645	264	1336827	28.171	ng/ul	0.00

Target Compounds Qvalue

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

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