

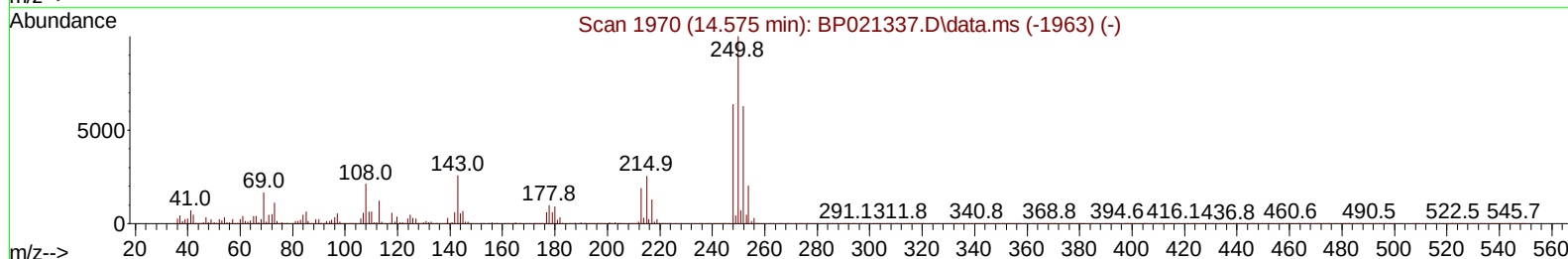
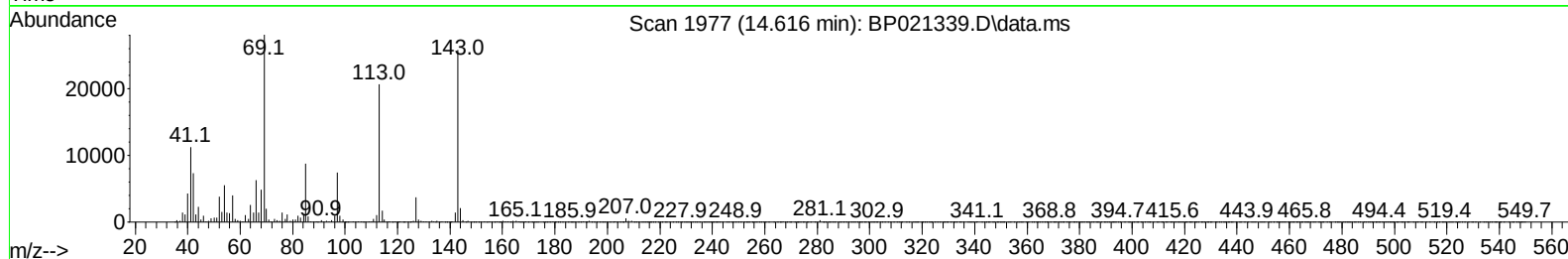
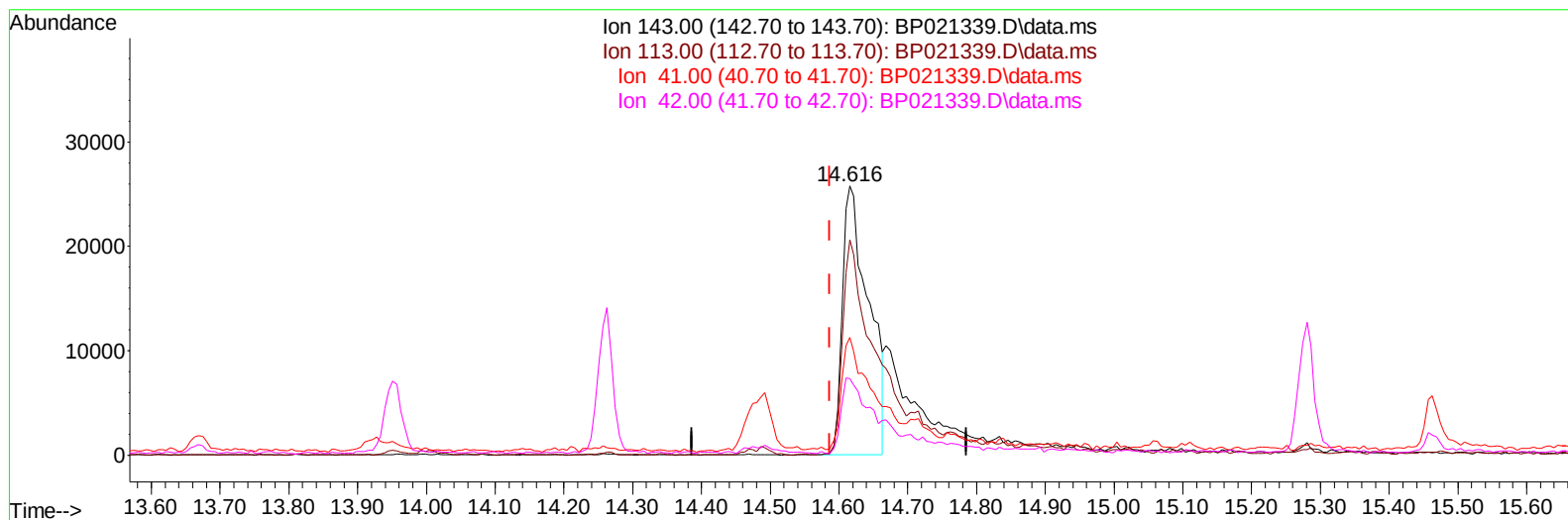
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080324\
Data File : BP021339.D
Acq On : 03 Aug 2024 10:56
Operator : CG/JU
Sample : P3278-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E1ZP6

Manual IntegrationsAPPROVED

Quant Time: Aug 04 23:20:46 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Jul 29 12:46:18 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 08/05/2024
Supervised By :mohammad ahmed 08/05/2024



TIC: BP021339.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.616min (+ 0.030) 9.31 ng/ul

response 68015

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	74.70	80.07
41.00	37.90	43.64
42.00	27.10	28.32

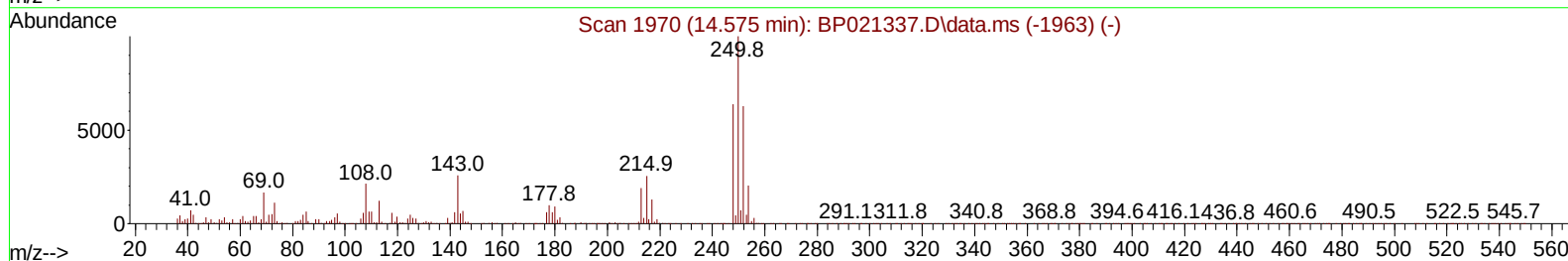
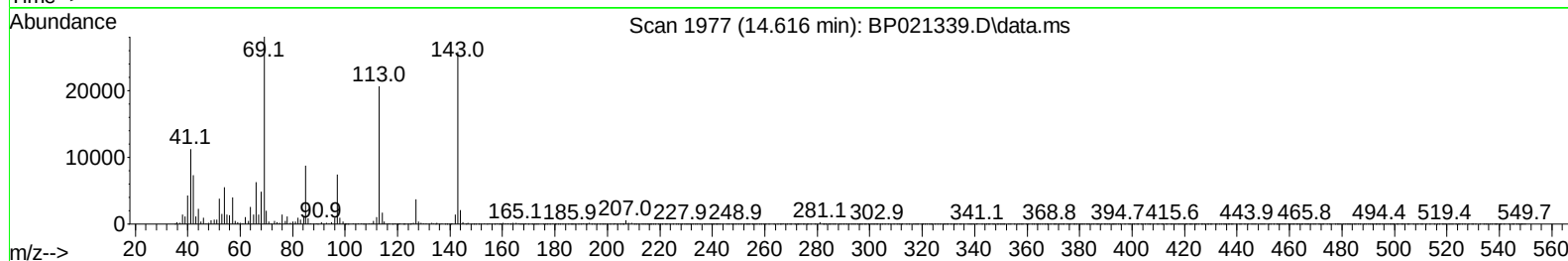
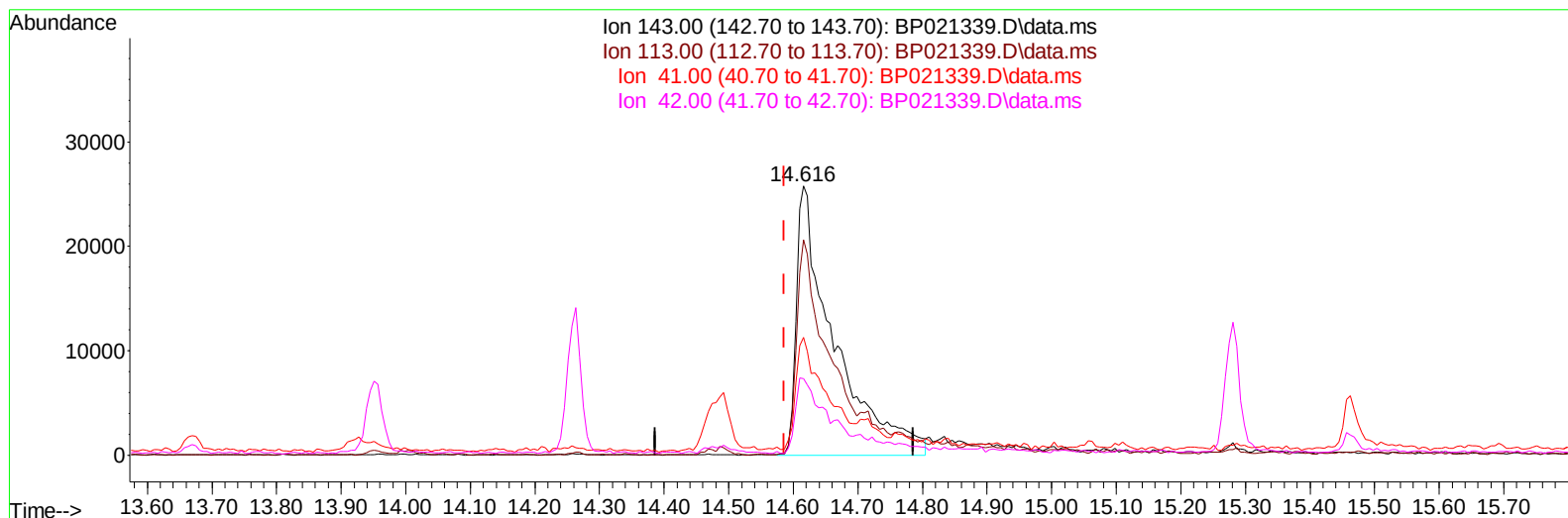
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(54) 4-Nitrophenol-d4 (S)

14.616min (+ 0.030) 14.19 ng/ul m

response 103679

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	74.70	80.07
41.00	37.90	43.64
42.00	27.10	28.32

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.616	152	261634	20.000	ng/ul	-0.02
20) Naphthalene-d8	10.387	136	1063487	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.263	164	706184	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.086	188	1504101	20.000	ng/ul	0.00
79) Chrysene-d12	21.521	240	1045520	20.000	ng/ul	-0.01
88) Perylene-d12	24.815	264	1190947	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.169	96	17953	3.123	ng/uL	0.00
4) Pyridine-d5	3.610	84	30331	1.822	ng/ul	0.02
7) Phenol-d5	6.851	99	341224	15.997	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.975	67	214869	16.622	ng/ul	0.00
11) 2-Chlorophenol-d4	7.181	132	291518	16.588	ng/ul	0.00
15) 4-Methylphenol-d8	8.357	113	275994	15.579	ng/ul	0.00
21) Nitrobenzene-d5	8.781	128	138080	16.716	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.492	143	159755	16.898	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.057	165	275457	16.113	ng/ul	0.00
31) 4-Chloroaniline-d4	10.557	131	290624	12.832	ng/ul	0.00
46) Dimethylphthalate-d6	13.669	166	939496	18.301	ng/ul	-0.01
49) Acenaphthylene-d8	13.951	160	1029307	17.857	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.616	143	103679m	14.195	ng/ul	0.03
60) Fluorene-d10	15.280	176	811148	19.260	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.463	200	105433	11.585	ng/ul	0.00
73) Anthracene-d10	17.186	188	1223907	18.321	ng/ul	0.00
81) Pyrene-d10	19.568	212	1266991	19.601	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.586	264	842379	14.341	ng/ul	-0.01

Target Compounds Qvalue

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

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