

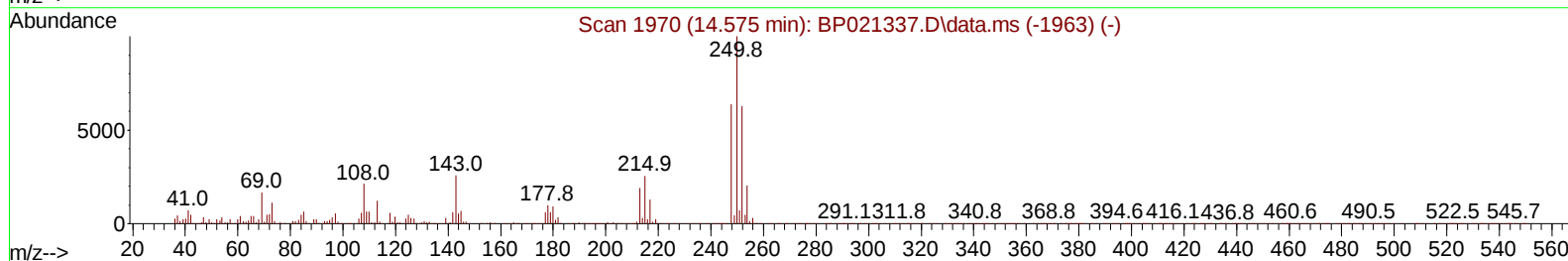
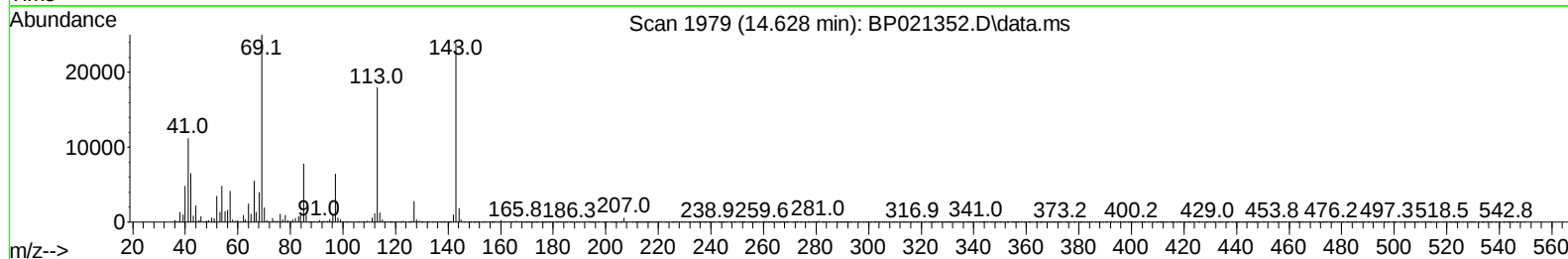
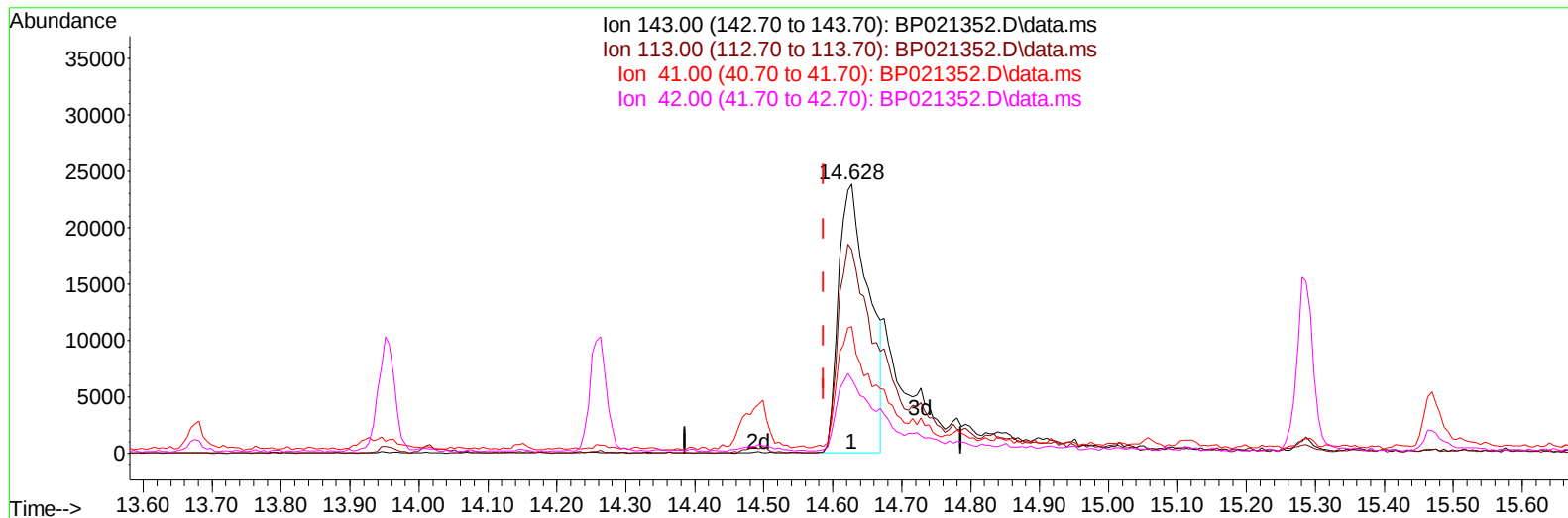
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080324\
Data File : BP021352.D
Acq On : 03 Aug 2024 21:26
Operator : CG/JU
Sample : P3278-08
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E1ZQ4

Manual IntegrationsAPPROVED

Quant Time: Aug 04 23:24:13 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: BP021352.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.628min (+ 0.041) 12.08 ng/u1

response 72132

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	74.70	75.53
41.00	37.90	47.04#
42.00	27.10	27.42

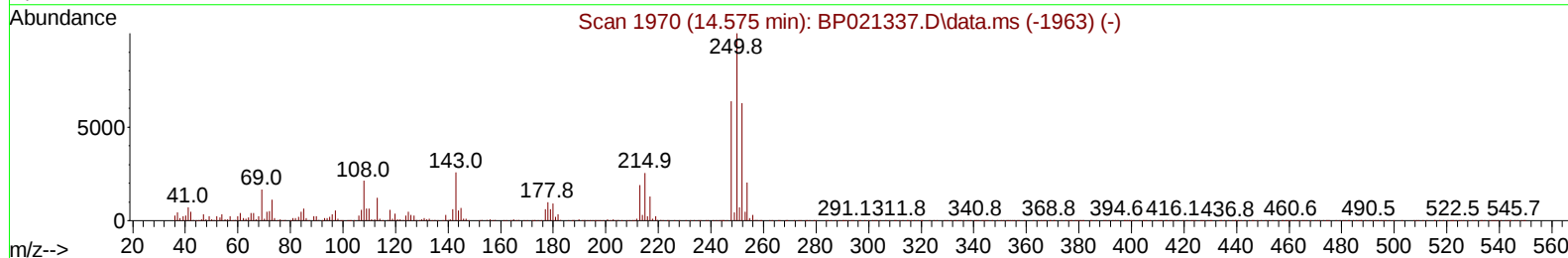
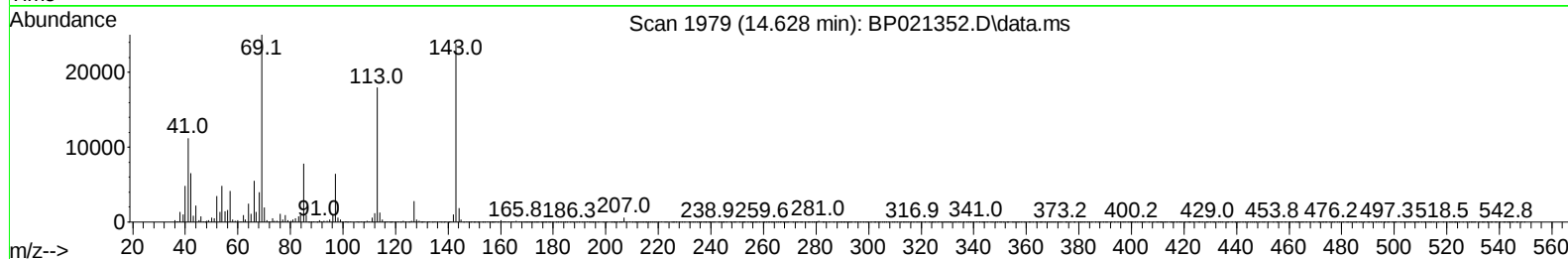
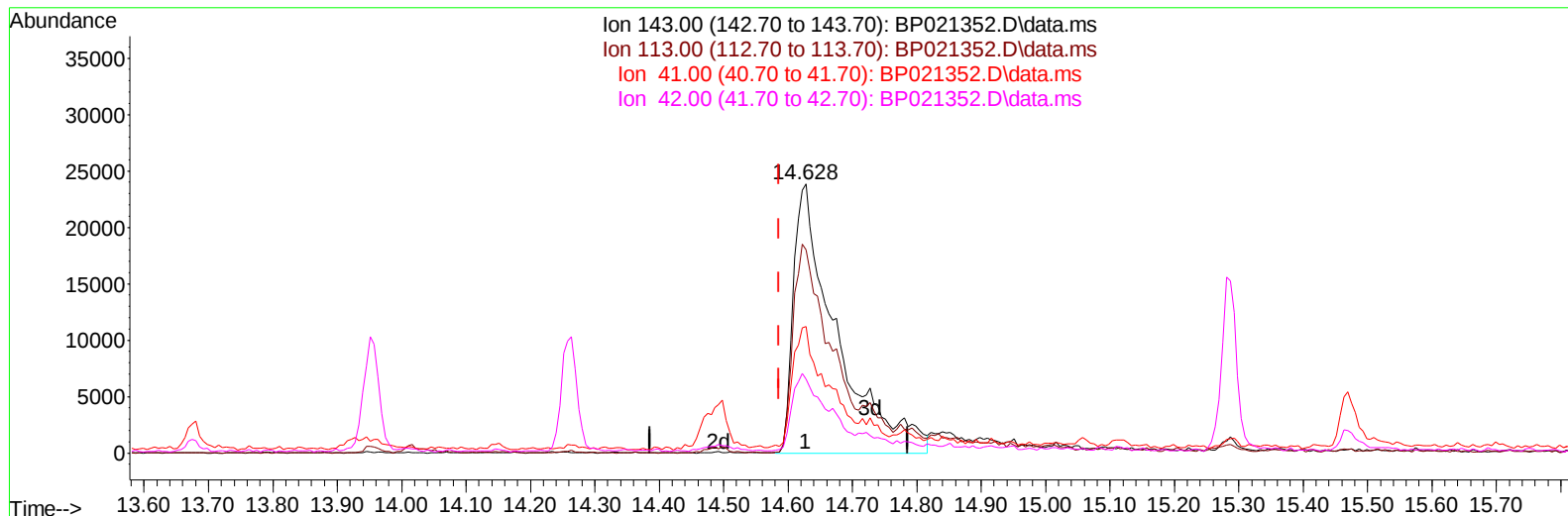
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(54) 4-Nitrophenol-d4 (S)

14.628min (+ 0.041) 18.52 ng/ul m

response 110580

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	74.70	75.53
41.00	37.90	47.04#
42.00	27.10	27.42

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Manual IntegrationsAPPROVED

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

Quant Time: Aug 05 00:16:02 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

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.616	152	242399	20.000	ng/ul	-0.02
20) Naphthalene-d8	10.381	136	965170	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.257	164	577352	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.086	188	1057433	20.000	ng/ul	0.00
79) Chrysene-d12	21.521	240	952905	20.000	ng/ul	-0.01
88) Perylene-d12	24.804	264	1249294	20.000	ng/ul	-0.03
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.164	96	24170	4.537	ng/uL	-0.01
4) Pyridine-d5	3.587	84	329524	21.368	ng/ul	0.00
7) Phenol-d5	6.840	99	489368	24.763	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.969	67	300913	25.126	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.169	132	408678	25.100	ng/ul	-0.01
15) 4-Methylphenol-d8	8.352	113	416231	25.359	ng/ul	-0.01
21) Nitrobenzene-d5	8.781	128	195035	26.016	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.493	143	227791	26.548	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.046	165	398264	25.669	ng/ul	0.00
31) 4-Chloroaniline-d4	10.551	131	524328	25.510	ng/ul	0.00
46) Dimethylphthalate-d6	13.681	166	1250403	29.792	ng/ul	0.00
49) Acenaphthylene-d8	13.951	160	1410507	29.930	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.628	143	110580m	18.518	ng/ul	0.04
60) Fluorene-d10	15.286	176	1035448	30.072	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.469	200	107699	16.833	ng/ul	0.00
73) Anthracene-d10	17.186	188	1406085	29.938	ng/ul	0.00
81) Pyrene-d10	19.569	212	1577864	26.783	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.586	264	2023187	32.836	ng/ul	-0.01

Target Compounds Qvalue

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

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