

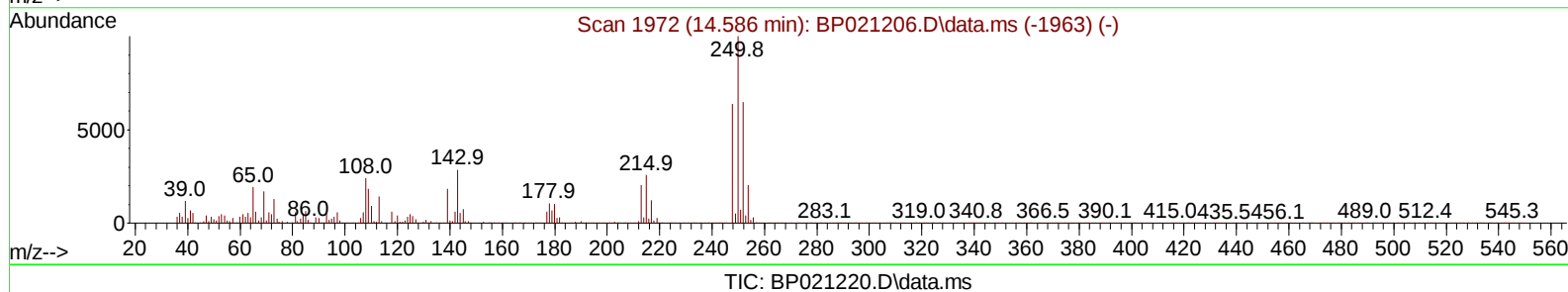
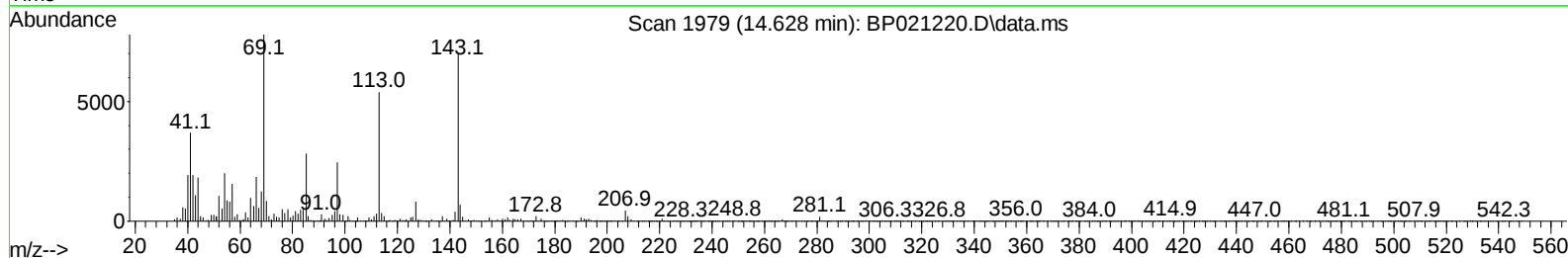
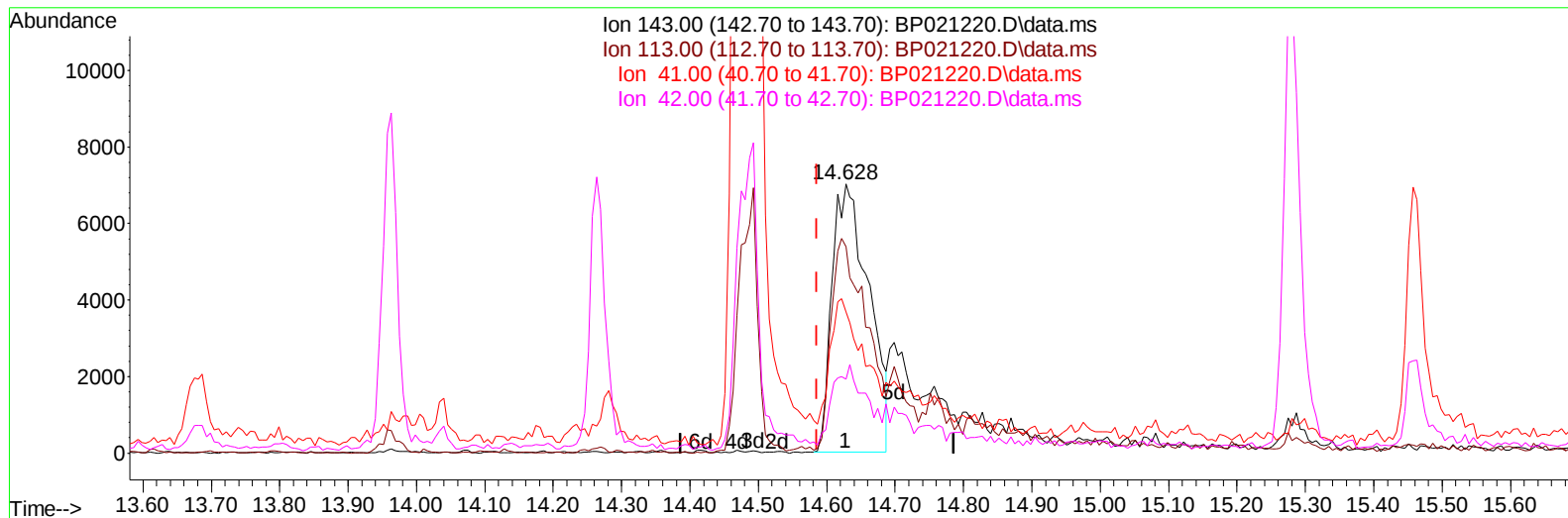
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072924\
Data File : BP021220.D
Acq On : 29 Jul 2024 22:57
Operator : MA/JU
Sample : P3280-13
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
YDZJ2

Manual IntegrationsAPPROVED

Quant Time: Jul 29 23:57:15 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 07/30/2024
Supervised By :mohammad ahmed 09/04/2024



(54) 4-Nitrophenol-d4 (S)

14.628min (+ 0.042) 6.81 ng/u1

response 26122

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	74.70	76.72
41.00	37.90	52.57#
42.00	27.10	27.40

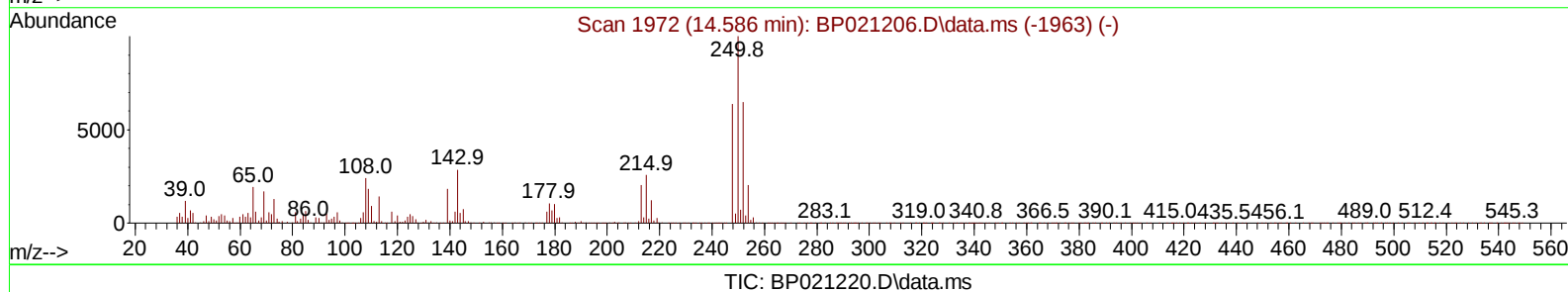
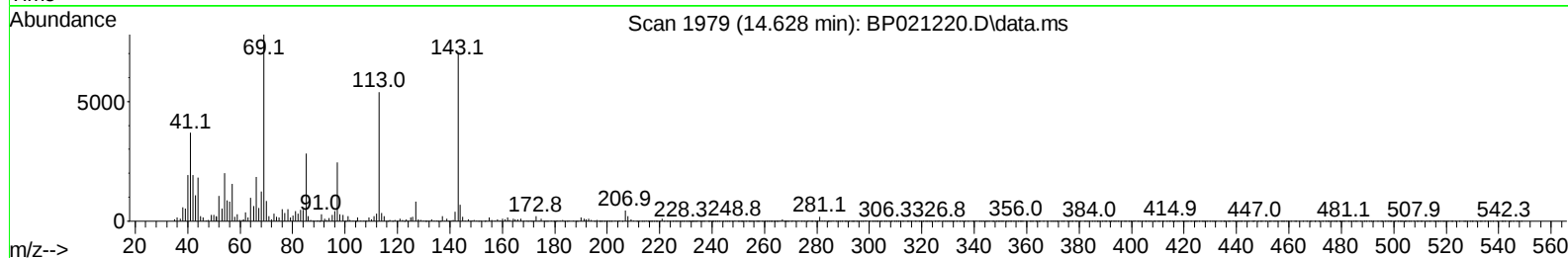
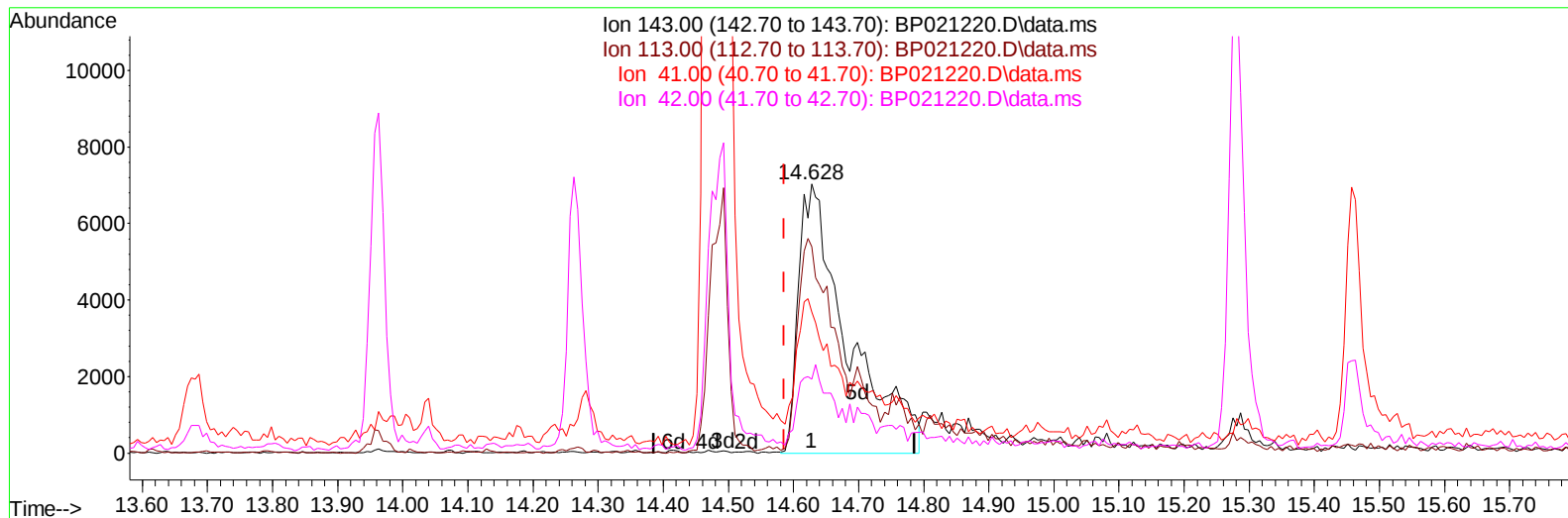
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(54) 4-Nitrophenol-d4 (S)

14.628min (+ 0.042) 9.67 ng/ul m

response 37073

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	74.70	76.72
41.00	37.90	52.57#
42.00	27.10	27.40

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.634	152	130166	20.000	ng/ul	0.00
20) Naphthalene-d8	10.399	136	584025	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.263	164	370750	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.087	188	802932	20.000	ng/ul	0.00
79) Chrysene-d12	21.539	240	749026	20.000	ng/ul	0.00
88) Perylene-d12	24.827	264	982265	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.176	96	21109	7.380	ng/uL	0.00
4) Pyridine-d5	3.599	84	133368	16.105	ng/ul	0.00
7) Phenol-d5	6.858	99	185637	17.493	ng/ul	0.01
9) Bis-(2-Chloroethyl)eth...	6.987	67	246751	38.368	ng/ul	0.00
11) 2-Chlorophenol-d4	7.187	132	269338	30.805	ng/ul	0.00
15) 4-Methylphenol-d8	8.369	113	275602	31.269	ng/ul	0.00
21) Nitrobenzene-d5	8.793	128	156426	34.484	ng/ul	0.00
24) 2-Nitrophenol-d4	9.505	143	176380	33.972	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.057	165	292565	31.163	ng/ul	0.00
31) 4-Chloroaniline-d4	10.557	131	331044	26.617	ng/ul	0.00
46) Dimethylphthalate-d6	13.675	166	990273	36.743	ng/ul	0.00
49) Acenaphthylene-d8	13.963	160	1069568	35.343	ng/ul	0.00
54) 4-Nitrophenol-d4	14.628	143	37073m	9.668	ng/ul	0.04
60) Fluorene-d10	15.281	176	821091	37.135	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.463	200	143888	29.618	ng/ul	0.00
73) Anthracene-d10	17.187	188	1400997	39.285	ng/ul	0.00
81) Pyrene-d10	19.575	212	1560344	33.695	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.598	264	2013509	41.563	ng/ul	0.00

Target Compounds Qvalue

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

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