

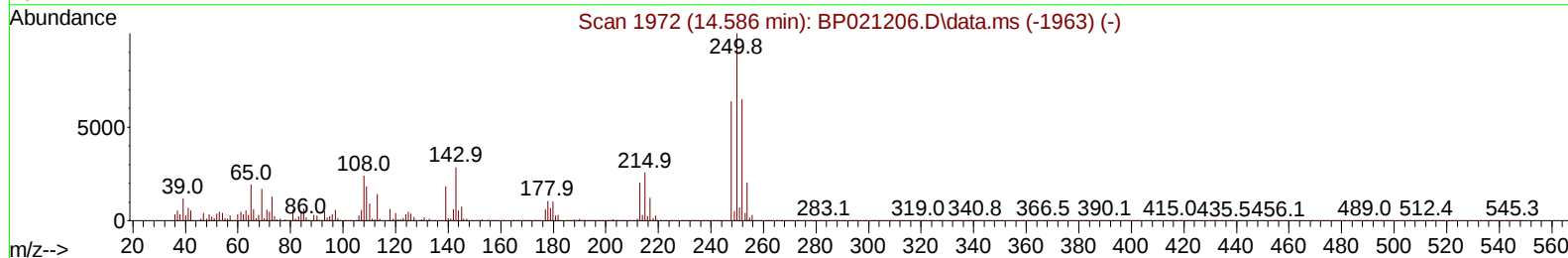
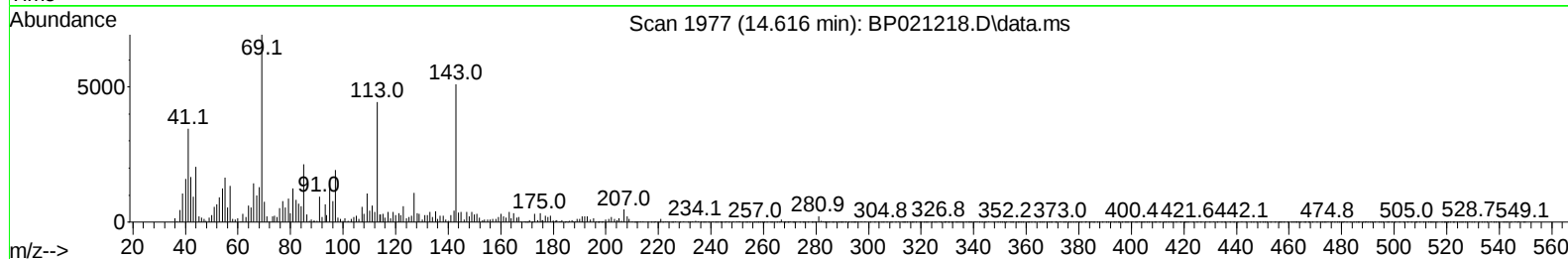
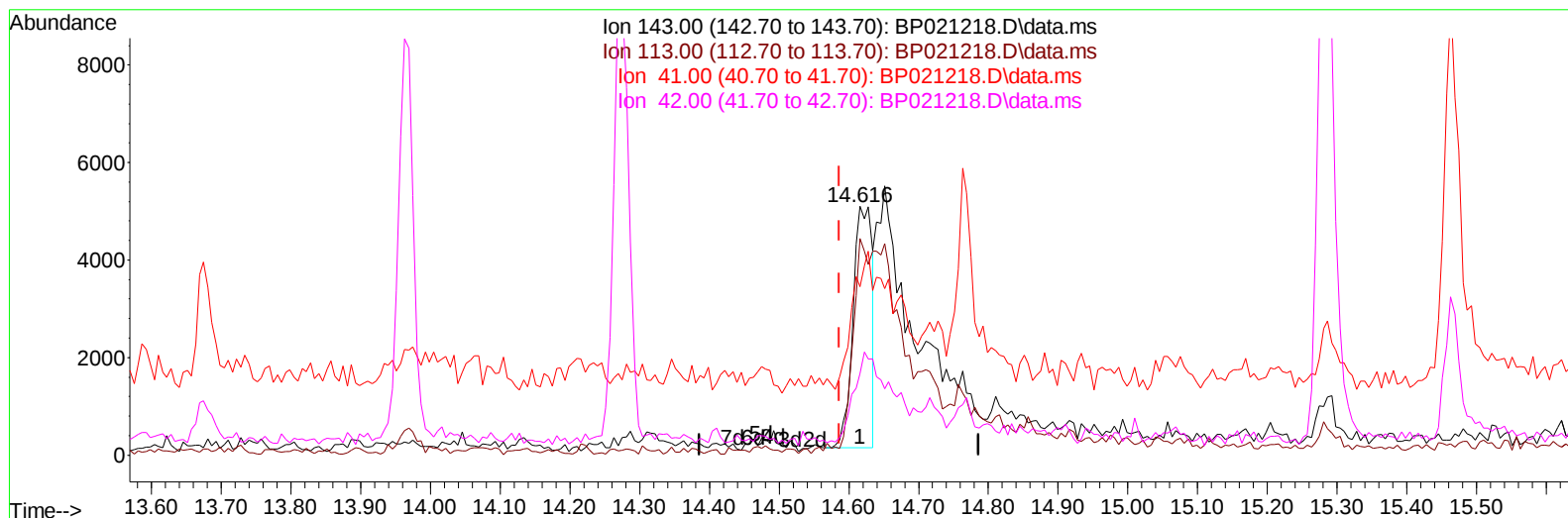
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072924\
Data File : BP021218.D
Acq On : 29 Jul 2024 21:29
Operator : MA/JU
Sample : P3280-07
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
YDZE1

Manual IntegrationsAPPROVED

Quant Time: Jul 29 22:48:06 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



TIC: BP021218.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.616min (+ 0.030) 1.89 ng/u1

response 9385

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	74.70	87.06
41.00	37.90	67.74#
42.00	27.10	32.79#

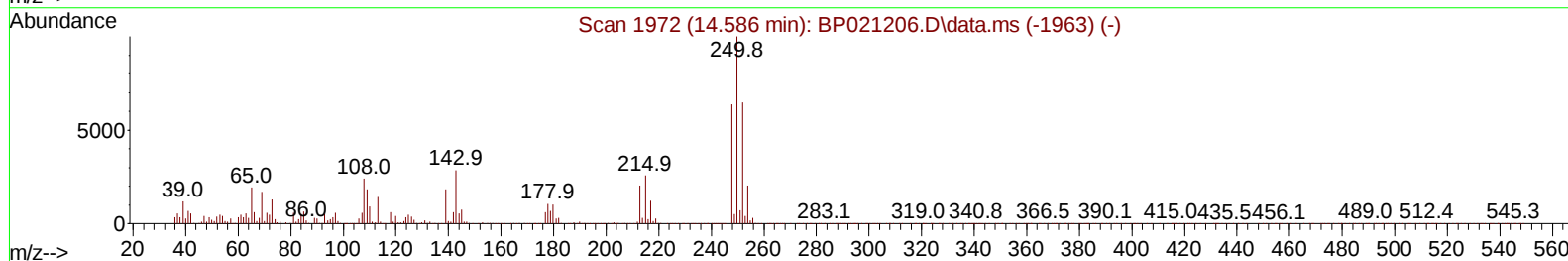
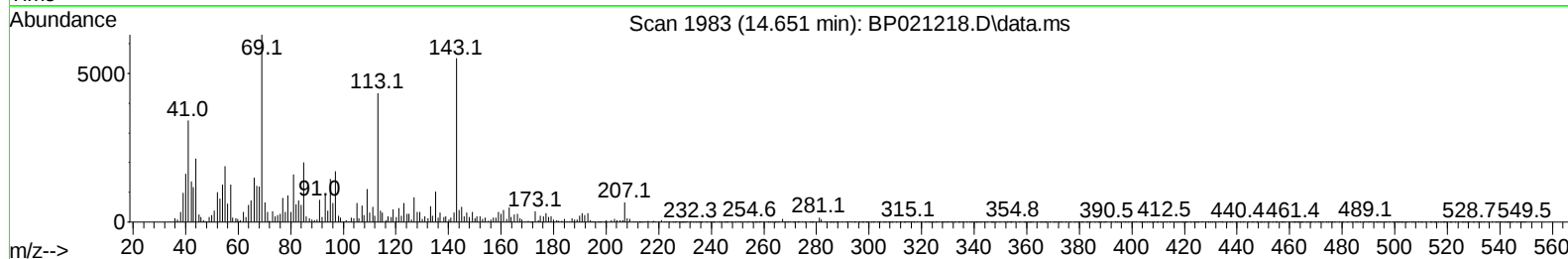
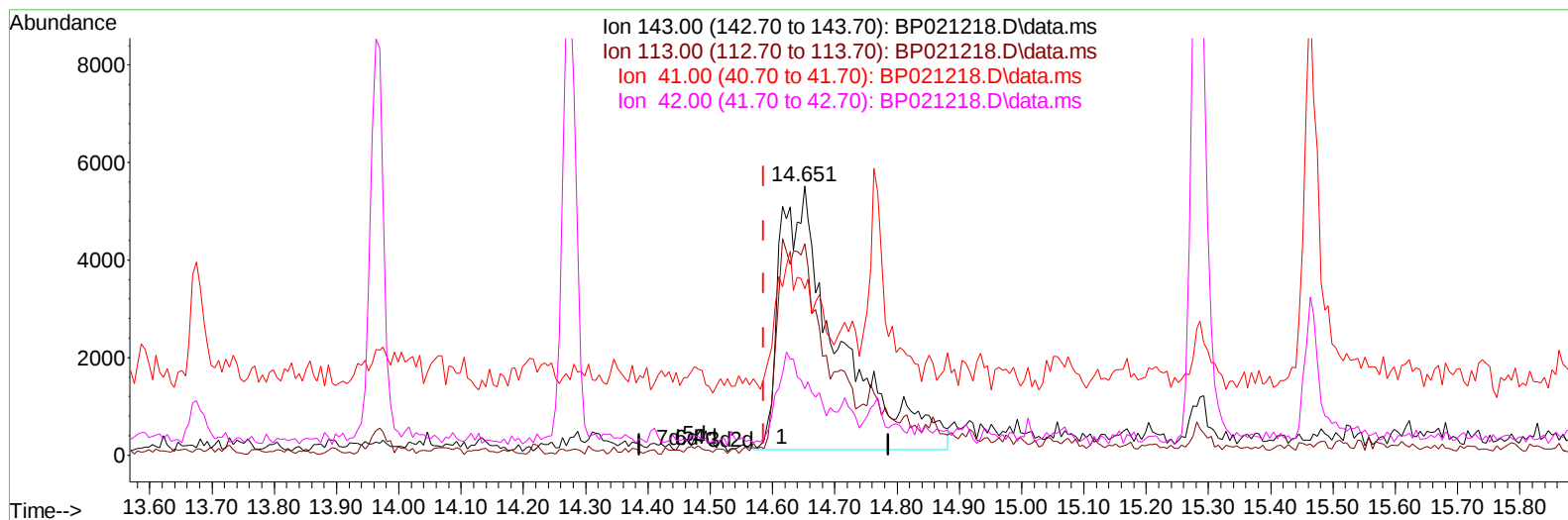
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(54) 4-Nitrophenol-d4 (S)

14.651min (+ 0.065) 7.16 ng/ul m

response 35516

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	74.70	78.59
41.00	37.90	61.93#
42.00	27.10	24.69

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 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 07/30/2024
 Supervised By :mohammad ahmed 09/04/2024

Quant Time: Jul 29 22:59:04 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	7.628	152	203953	20.000	ng/ul	0.00
20) Naphthalene-d8	10.405	136	690725	20.000	ng/ul	0.01
38) Acenaphthene-d10	14.275	164	479404	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.087	188	980203	20.000	ng/ul	0.00
79) Chrysene-d12	21.533	240	1040171	20.000	ng/ul	0.00
88) Perylene-d12	24.833	264	1258826	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.176	96	20447	4.562	ng/uL	0.00
4) Pyridine-d5	3.599	84	107094	8.254	ng/ul	0.00
7) Phenol-d5	6.864	99	104233	6.269	ng/ul	0.02
9) Bis-(2-Chloroethyl)eth...	6.987	67	245693	24.382	ng/ul	0.00
11) 2-Chlorophenol-d4	7.187	132	303529	22.156	ng/ul	0.00
15) 4-Methylphenol-d8	8.363	113	195885	14.184	ng/ul	0.00
21) Nitrobenzene-d5	8.793	128	151282	28.198	ng/ul	0.00
24) 2-Nitrophenol-d4	9.510	143	172148	28.035	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.057	165	289201	26.046	ng/ul	0.00
31) 4-Chloroaniline-d4	10.563	131	291353	19.807	ng/ul	0.00
46) Dimethylphthalate-d6	13.675	166	994942	28.549	ng/ul	0.00
49) Acenaphthylene-d8	13.963	160	1049841	26.828	ng/ul	0.00
54) 4-Nitrophenol-d4	14.651	143	35516m	7.163	ng/ul	0.07
60) Fluorene-d10	15.287	176	827527	28.944	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.463	200	146758	24.745	ng/ul	0.00
73) Anthracene-d10	17.192	188	1414499	32.491	ng/ul	0.00
81) Pyrene-d10	19.575	212	1794144	27.899	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.604	264	2159211	34.778	ng/ul	0.00

Target Compounds Qvalue

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

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