

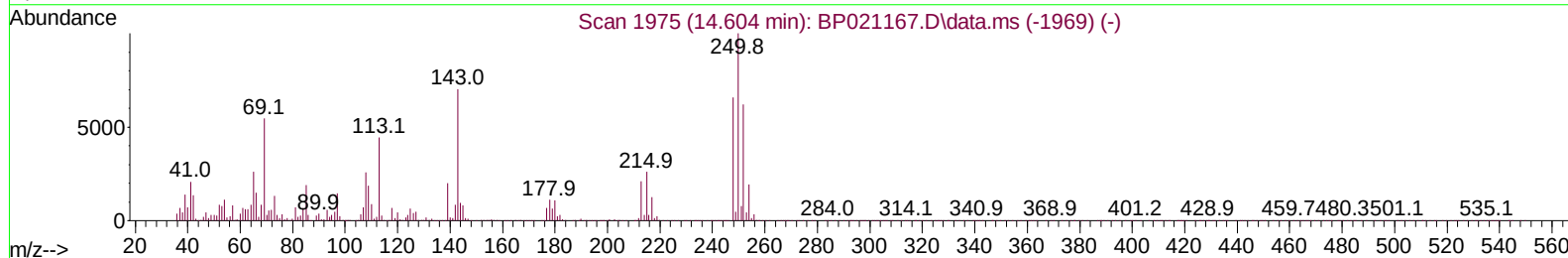
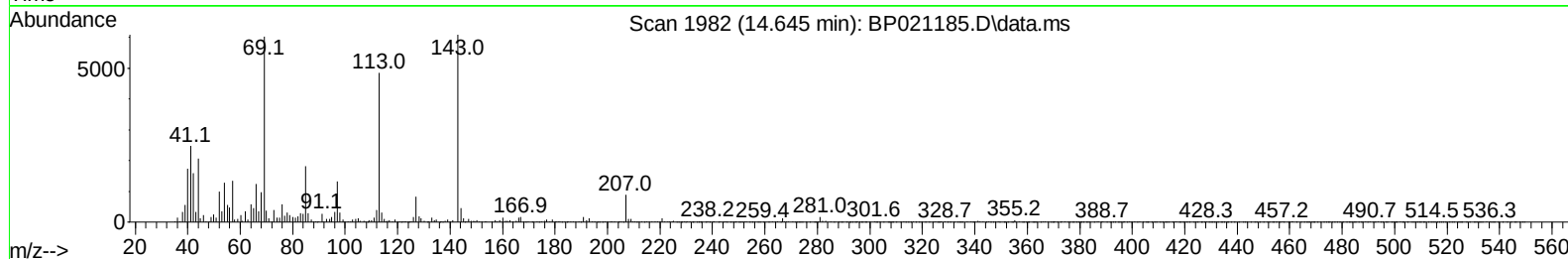
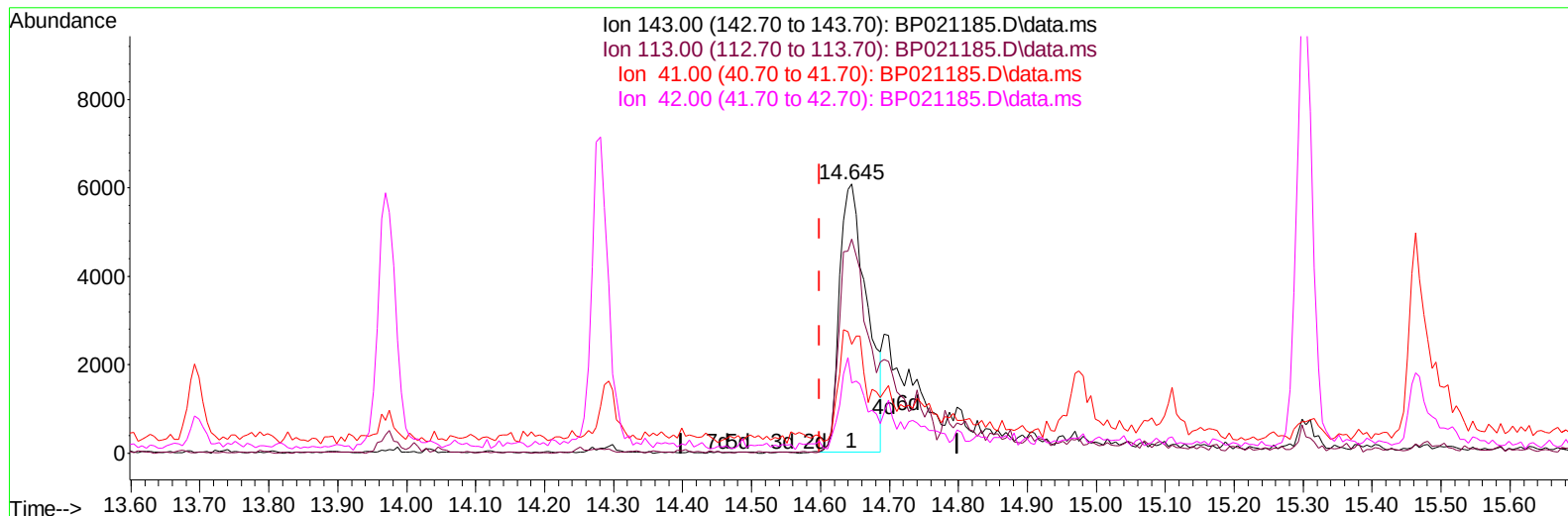
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072624\
Data File : BP021185.D
Acq On : 27 Jul 2024 02:09
Operator : MA/JU
Sample : P3301-02
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
F7E84

Manual IntegrationsAPPROVED

Quant Time: Jul 27 03:07:16 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 :Yogesh Patel 07/27/2024
Supervised By :mohammad ahmed 07/29/2024



TIC: BP021185.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.645min (+ 0.047) 4.13 ng/ul

response 16822

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	74.70	79.63
41.00	37.90	40.54
42.00	27.10	25.97

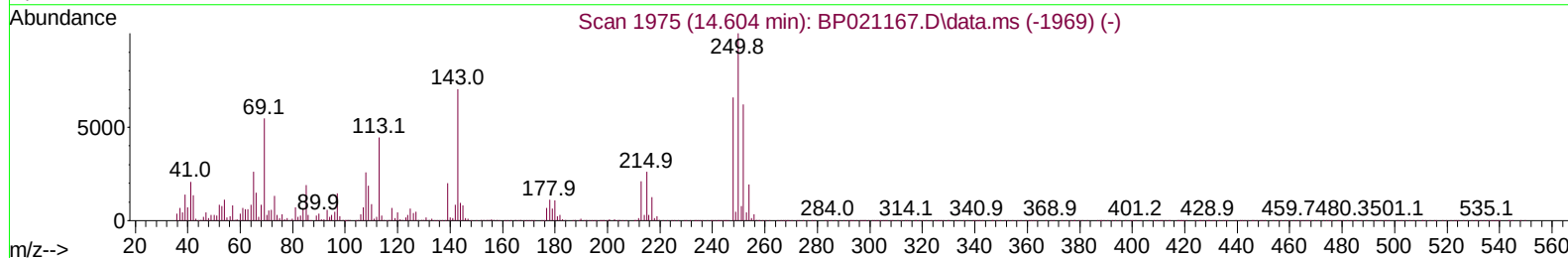
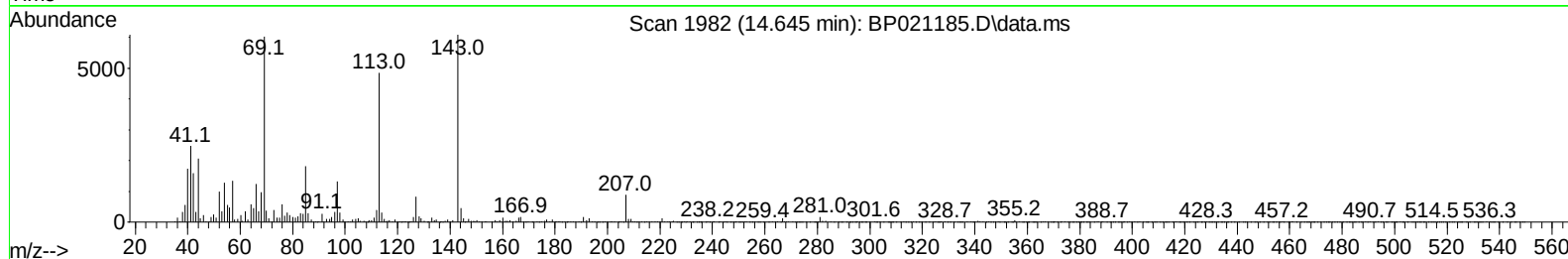
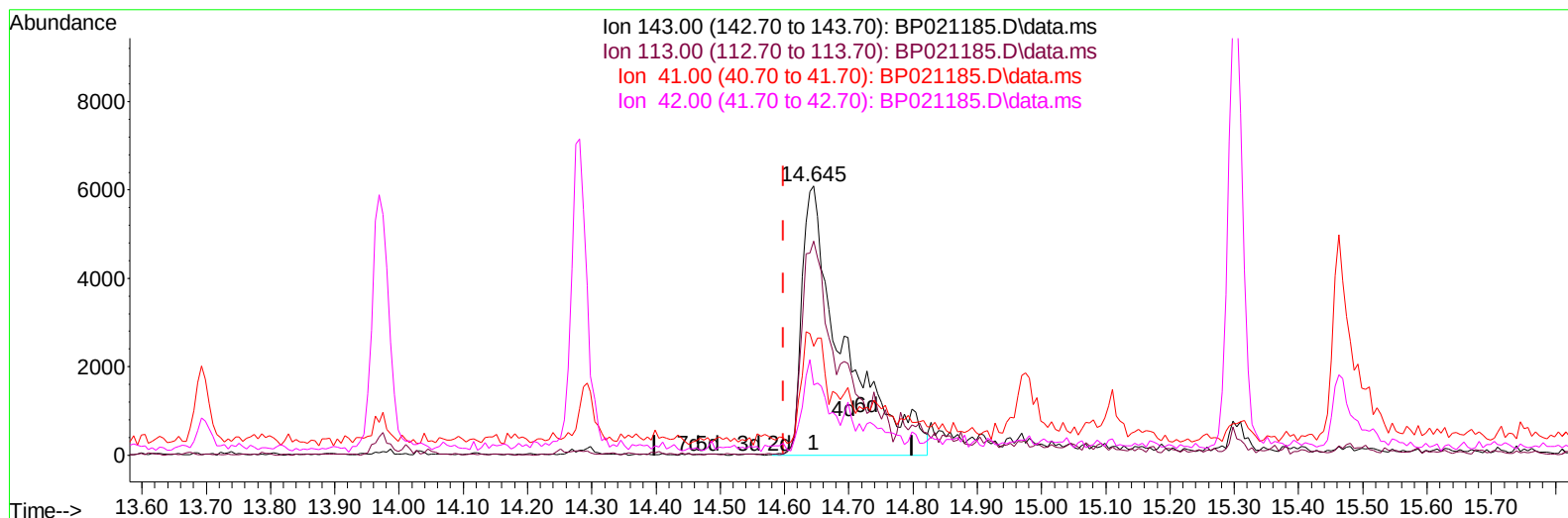
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TIC: BP021185.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.645min (+ 0.047) 6.71 ng/ul m

response 27292

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	74.70	79.63
41.00	37.90	40.54
42.00	27.10	25.97

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 ALS Vial : 20 Sample Multiplier: 1

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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 07/27/2024
 Supervised By :mohammad ahmed 07/29/2024

Quant Time: Jul 27 03:09:08 2024
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 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jul 25 00:15:24 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.646	152	175167	20.000	ng/ul	0.00
20) Naphthalene-d8	10.404	136	666697	20.000	ng/ul	-0.02
38) Acenaphthene-d10	14.281	164	393506	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.104	188	774464	20.000	ng/ul	0.00
79) Chrysene-d12	21.551	240	782848	20.000	ng/ul	0.00
88) Perylene-d12	24.863	264	991551	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.175	96	17757	4.613	ng/uL	0.00
4) Pyridine-d5	3.605	84	98563	8.844	ng/ul	0.00
7) Phenol-d5	6.869	99	100245	7.020	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.993	67	213947	24.721	ng/ul	0.00
11) 2-Chlorophenol-d4	7.193	132	272206	23.135	ng/ul	-0.01
15) 4-Methylphenol-d8	8.381	113	180161	15.189	ng/ul	0.00
21) Nitrobenzene-d5	8.805	128	135920	26.248	ng/ul	0.00
24) 2-Nitrophenol-d4	9.516	143	156020	26.324	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.063	165	258350	24.106	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.575	131	336865	23.727	ng/ul	0.00
46) Dimethylphthalate-d6	13.698	166	815515	28.509	ng/ul	0.00
49) Acenaphthylene-d8	13.969	160	881712	27.450	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.645	143	27292m	6.706	ng/ul	0.05
60) Fluorene-d10	15.304	176	702440	29.932	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.463	200	113358	24.191	ng/ul	0.00
73) Anthracene-d10	17.210	188	1167212	33.933	ng/ul	0.00
81) Pyrene-d10	19.592	212	1401508	28.957	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.633	264	1760289	35.995	ng/ul	-0.02

Target Compounds Qvalue

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

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