

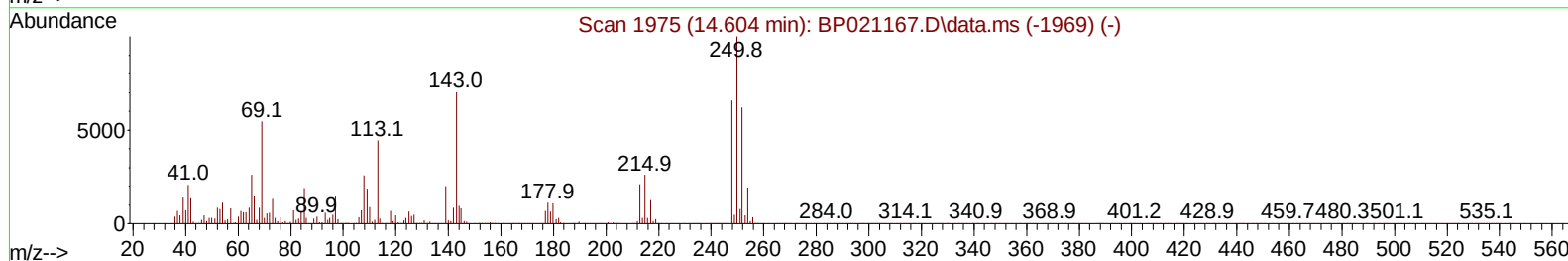
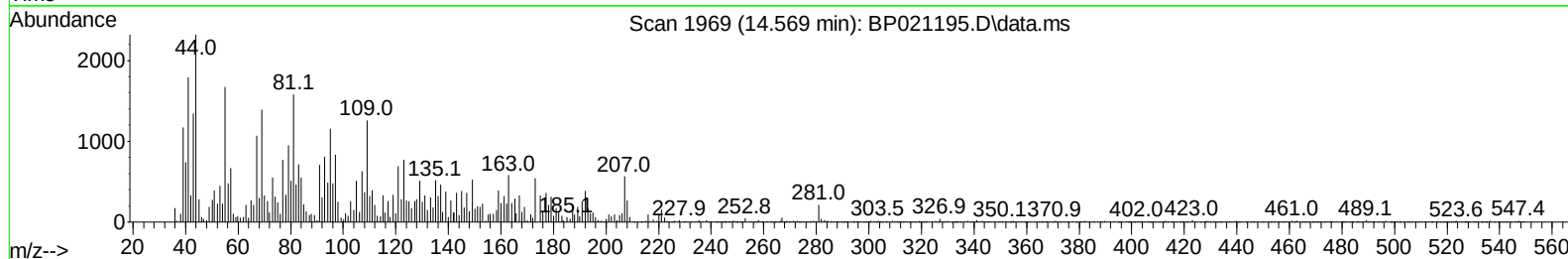
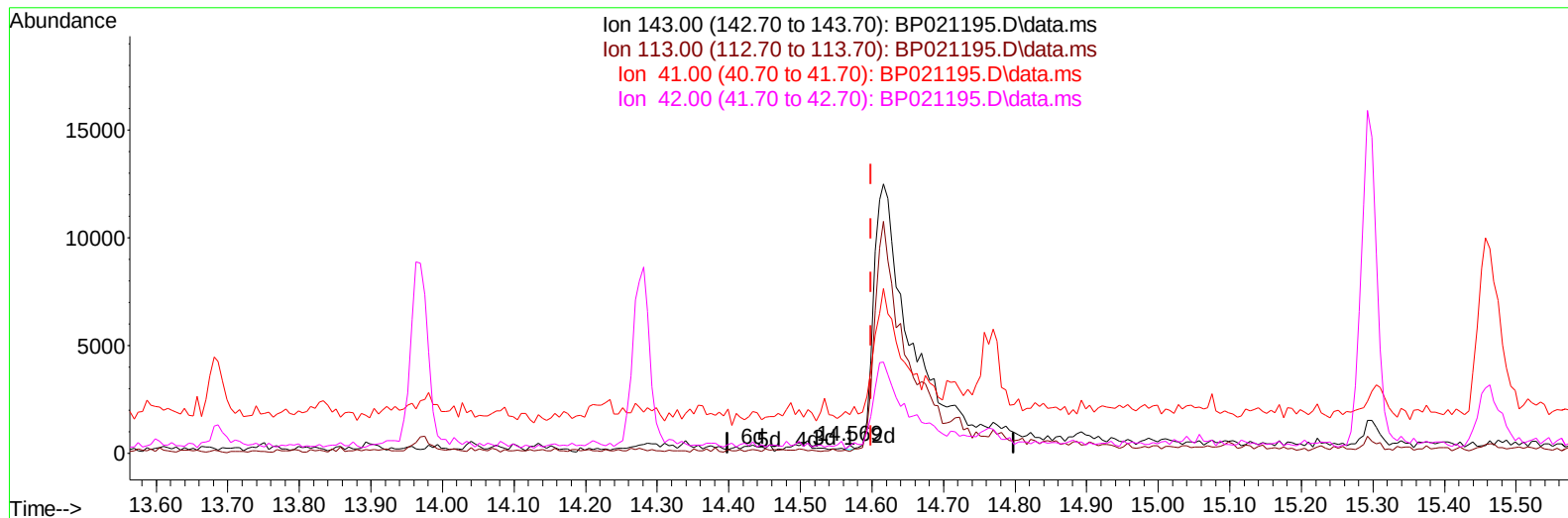
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072624\
Data File : BP021195.D
Acq On : 27 Jul 2024 10:09
Operator : MA/JU
Sample : P3280-07
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
YDZE1

Manual IntegrationsAPPROVED

Quant Time: Jul 29 00:41:36 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 07/29/2024
Supervised By :mohammad ahmed 07/29/2024



TIC: BP021195.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.569min (-0.029) 0.02 ng/u1

response 94

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	74.70	20.72#
41.00	37.90	495.58#
42.00	27.10	90.88#

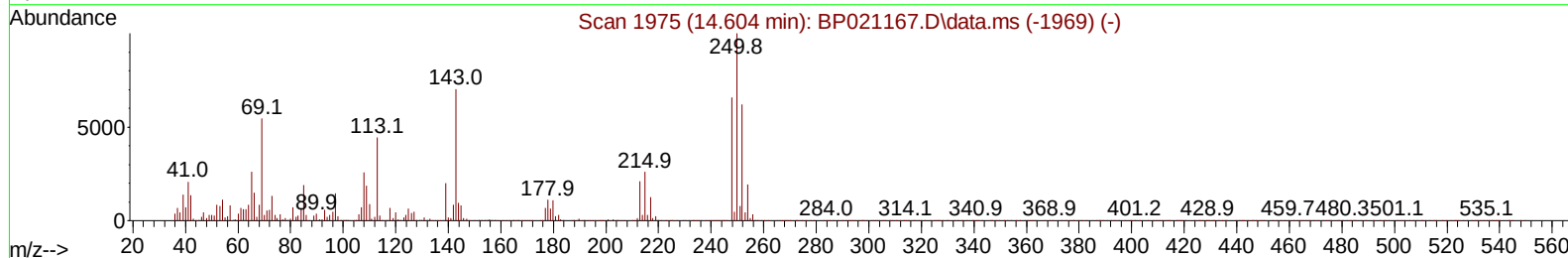
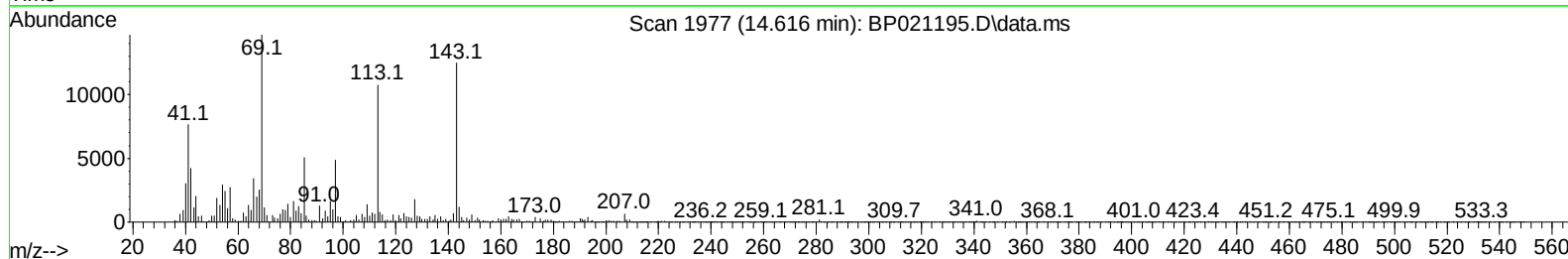
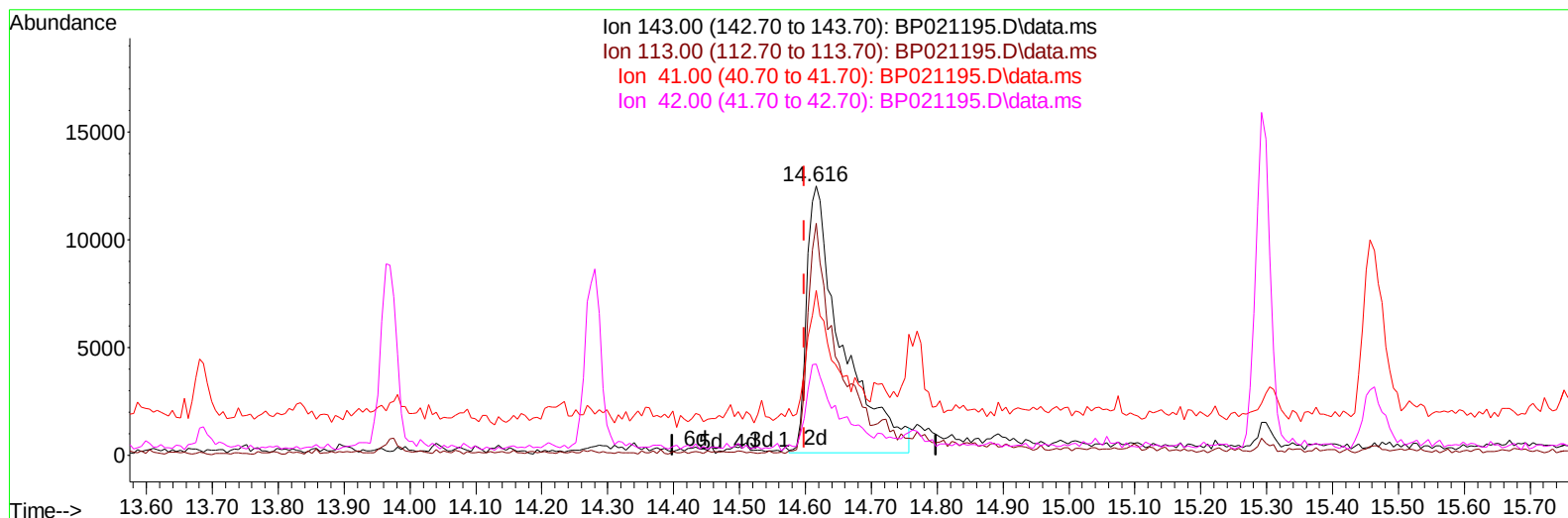
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(54) 4-Nitrophenol-d4 (S)

14.616min (+ 0.018) 9.35 ng/ul m

response 45980

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	74.70	86.25
41.00	37.90	61.22#
42.00	27.10	33.94#

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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 07/29/2024
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Quant Time: Jul 29 01:11:25 2024
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 QLast Update : Thu Jul 25 00:15:24 2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.652	152	189059	20.000	ng/ul	0.00
20) Naphthalene-d8	10.410	136	748239	20.000	ng/ul	-0.02
38) Acenaphthene-d10	14.281	164	475479	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.098	188	1034737	20.000	ng/ul	0.00
79) Chrysene-d12	21.539	240	1060535	20.000	ng/ul	-0.02
88) Perylene-d12	24.851	264	1253805	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.181	96	19020	4.578	ng/uL	0.00
4) Pyridine-d5	3.605	84	120581	10.025	ng/ul	0.00
7) Phenol-d5	6.875	99	120533	7.820	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.993	67	262207	28.071	ng/ul	0.00
11) 2-Chlorophenol-d4	7.205	132	334059	26.305	ng/ul	0.00
15) 4-Methylphenol-d8	8.375	113	229669	17.941	ng/ul	0.00
21) Nitrobenzene-d5	8.799	128	174567	30.037	ng/ul	0.00
24) 2-Nitrophenol-d4	9.510	143	199925	30.056	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.069	165	343806	28.584	ng/ul	0.00
31) 4-Chloroaniline-d4	10.569	131	380415	23.874	ng/ul	0.00
46) Dimethylphthalate-d6	13.687	166	1199781	34.711	ng/ul	-0.01
49) Acenaphthylene-d8	13.969	160	1239905	31.947	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.616	143	45980m	9.350	ng/ul	0.02
60) Fluorene-d10	15.293	176	996681	35.148	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.457	200	200534	32.031	ng/ul	-0.01
73) Anthracene-d10	17.198	188	1725496	37.545	ng/ul	0.00
81) Pyrene-d10	19.581	212	2200511	33.561	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.621	264	2488066	40.236	ng/ul	-0.03

Target Compounds Qvalue

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

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