

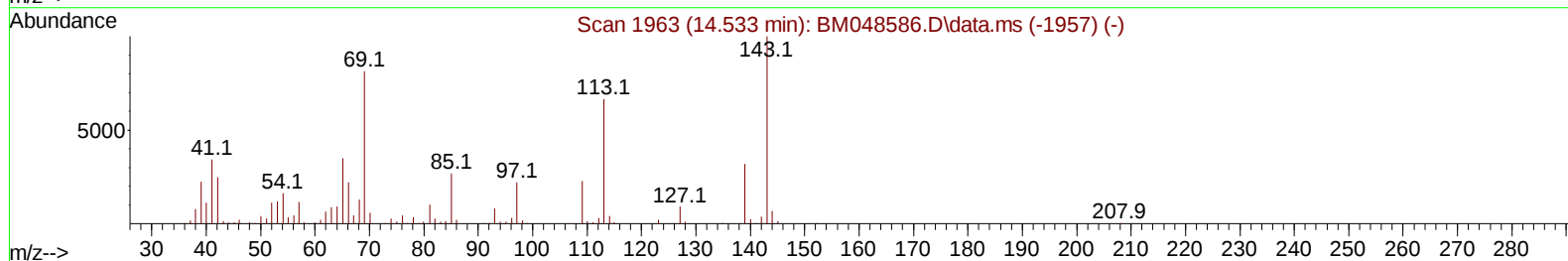
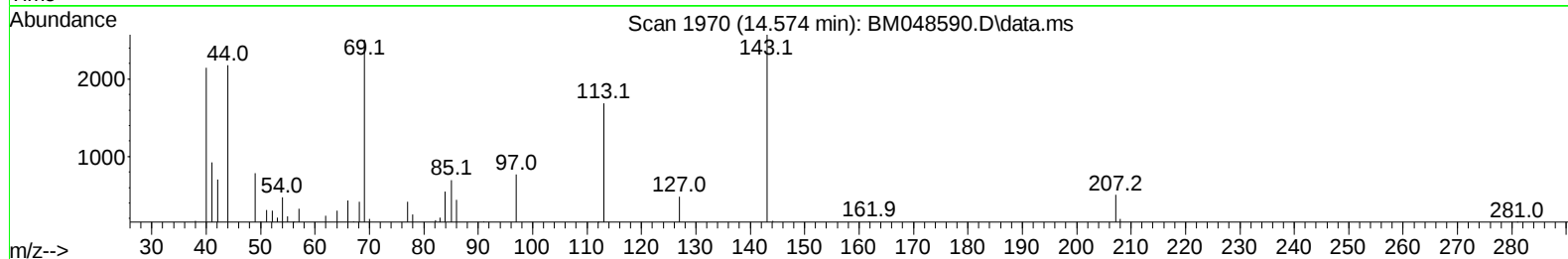
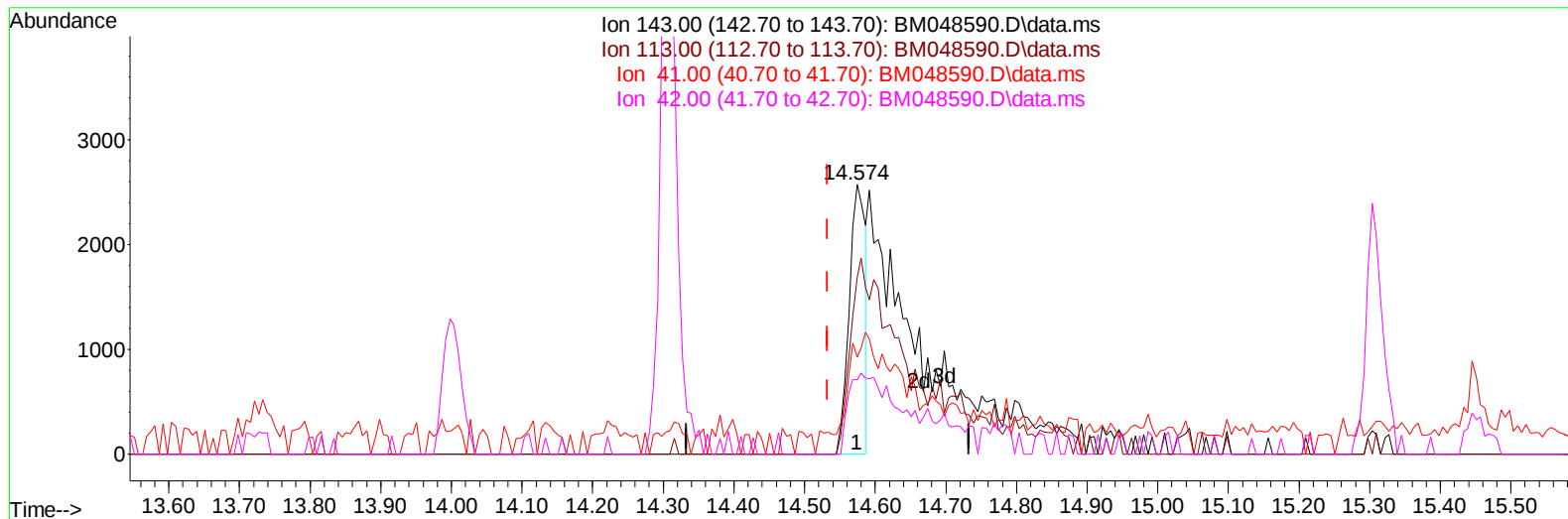
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111124\
Data File : BM048590.D
Acq On : 11 Nov 2024 18:23
Operator : RC/JU
Sample : P4585-05
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
A4ED9

Manual IntegrationsAPPROVED

Quant Time: Nov 11 22:53:48 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM110724.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Nov 07 22:50:20 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 11/12/2024
Supervised By :mohammad ahmed 11/13/2024



TIC: BM048590.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.574min (+ 0.041) 1.06 ng/u1

response 4091

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	66.50	65.73
41.00	36.20	35.90
42.00	24.50	27.38

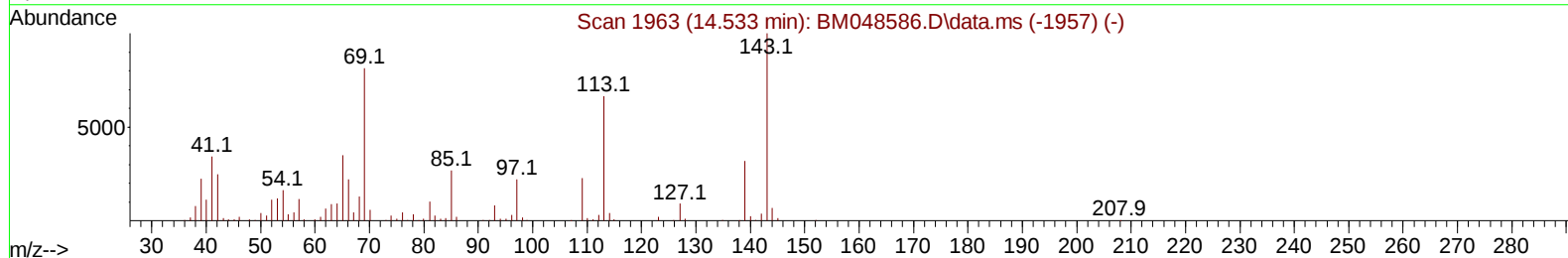
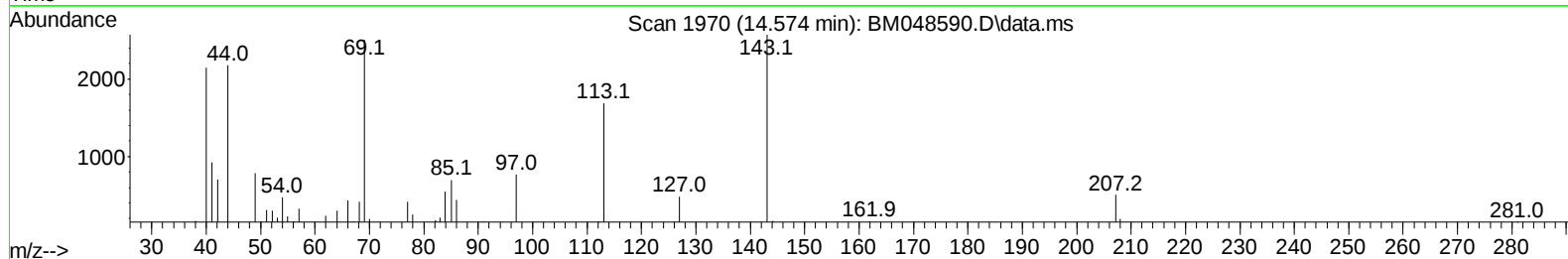
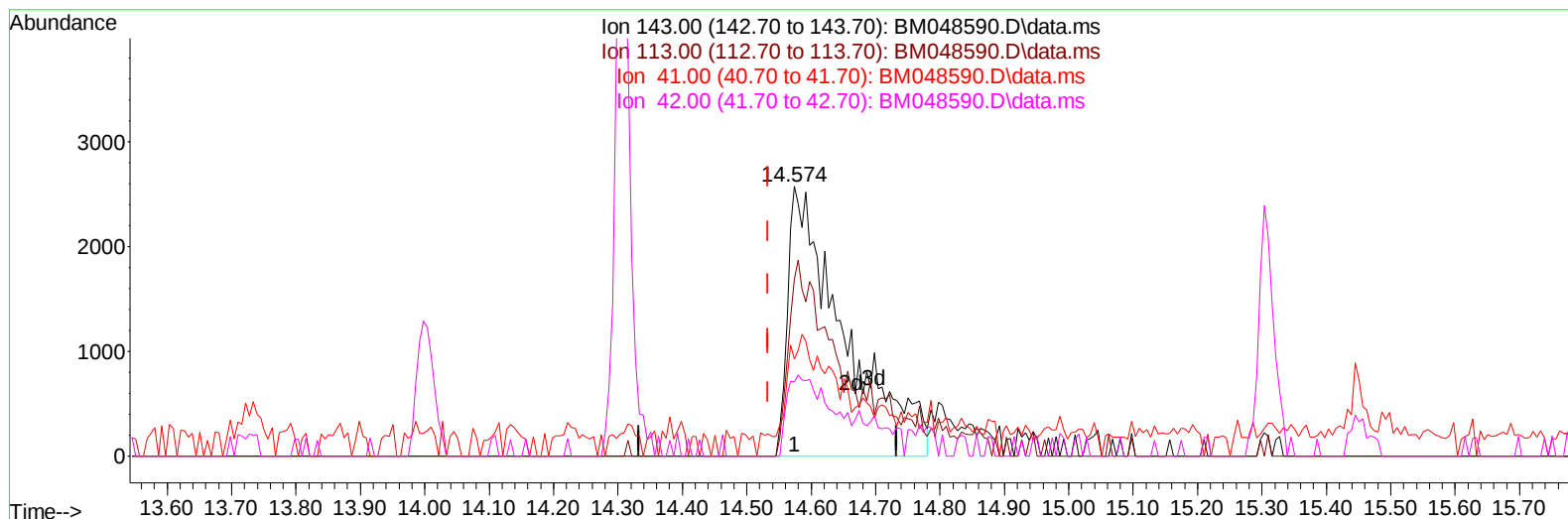
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TIC: BM048590.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.574min (+ 0.041) 4.01 ng/ul m

response 15478

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	66.50	65.73
41.00	36.20	35.90
42.00	24.50	27.38

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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 11/12/2024
 Supervised By :mohammad ahmed 11/13/2024

Quant Time: Nov 11 23:02:38 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM110724.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Nov 07 22:50:20 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.657	152	129989	20.000	ng/ul	0.00
20) Naphthalene-d8	10.445	136	526247	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.310	164	325600	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.057	188	681807	20.000	ng/ul	0.00
79) Chrysene-d12	21.280	240	688904	20.000	ng/ul	0.00
88) Perylene-d12	24.191	264	812531	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.128	96	2883	0.884	ng/uL	0.00
4) Pyridine-d5	3.569	84	15424	1.716	ng/ul	0.02
7) Phenol-d5	6.845	99	70272	7.102	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.998	67	44092	7.418	ng/ul	0.00
11) 2-Chlorophenol-d4	7.198	132	62770	7.658	ng/ul	0.00
15) 4-Methylphenol-d8	8.387	113	47394	6.091	ng/ul	0.00
21) Nitrobenzene-d5	8.822	128	25898	6.790	ng/ul	0.00
24) 2-Nitrophenol-d4	9.545	143	27683	6.824	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.092	165	53602	7.112	ng/ul	0.01
31) 4-Chloroaniline-d4	10.610	131	58020	5.420	ng/ul	0.01
46) Dimethylphthalate-d6	13.727	166	183530	8.830	ng/ul	0.00
49) Acenaphthylene-d8	14.004	160	200131	8.167	ng/ul	0.00
54) 4-Nitrophenol-d4	14.574	143	15478m	4.009	ng/ul	0.04
60) Fluorene-d10	15.310	176	160892	8.868	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.451	200	11648	3.396	ng/ul	0.01
73) Anthracene-d10	17.157	188	240177	8.324	ng/ul	0.00
81) Pyrene-d10	19.451	212	279911	8.096	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.991	264	254865	6.713	ng/ul	0.00
Target Compounds						
					Qvalue	
16) Acetophenone	8.492	105	49523	4.083	ng/ul	99
80) Fluoranthene	19.115	202	85585	2.133	ng/ul	98
82) Pyrene	19.480	202	47534	1.093	ng/ul	99
86) Bis(2-ethylhexyl)phtha...	21.186	149	146424	6.007	ng/ul	100
87) Chrysene	21.321	228	52725	1.263	ng/ul	96
90) Benzo(b)fluoranthene	23.274	252	62872	1.406	ng/ul	98

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

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