

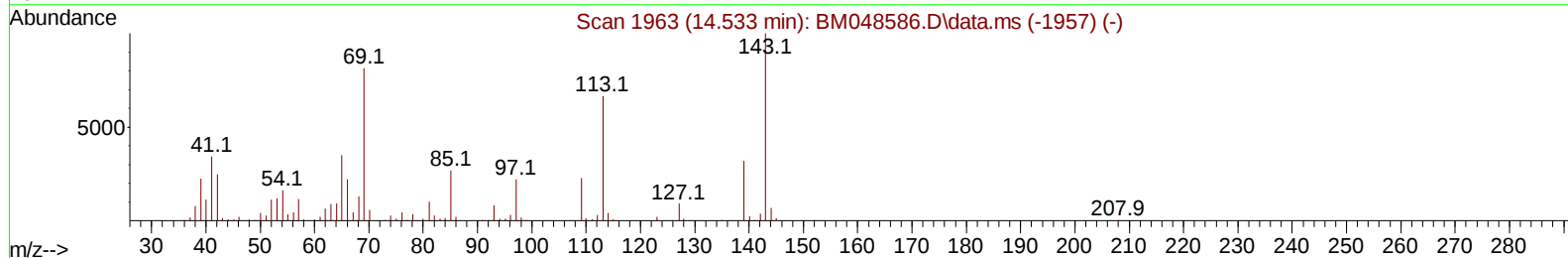
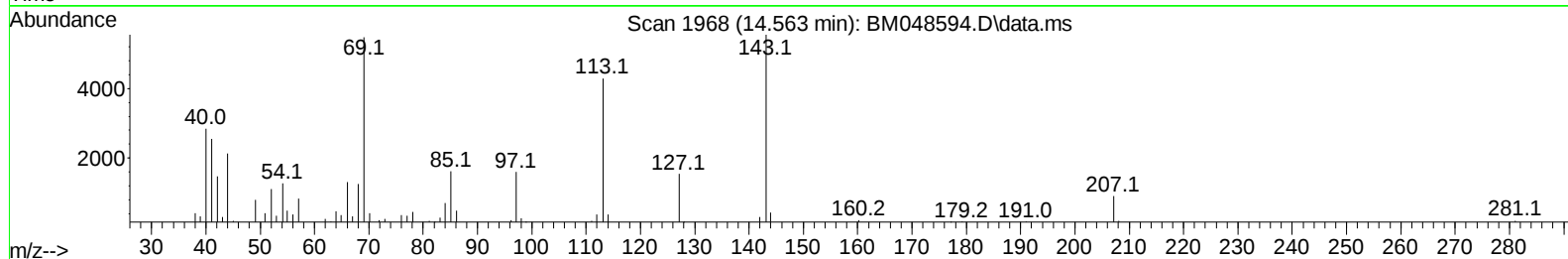
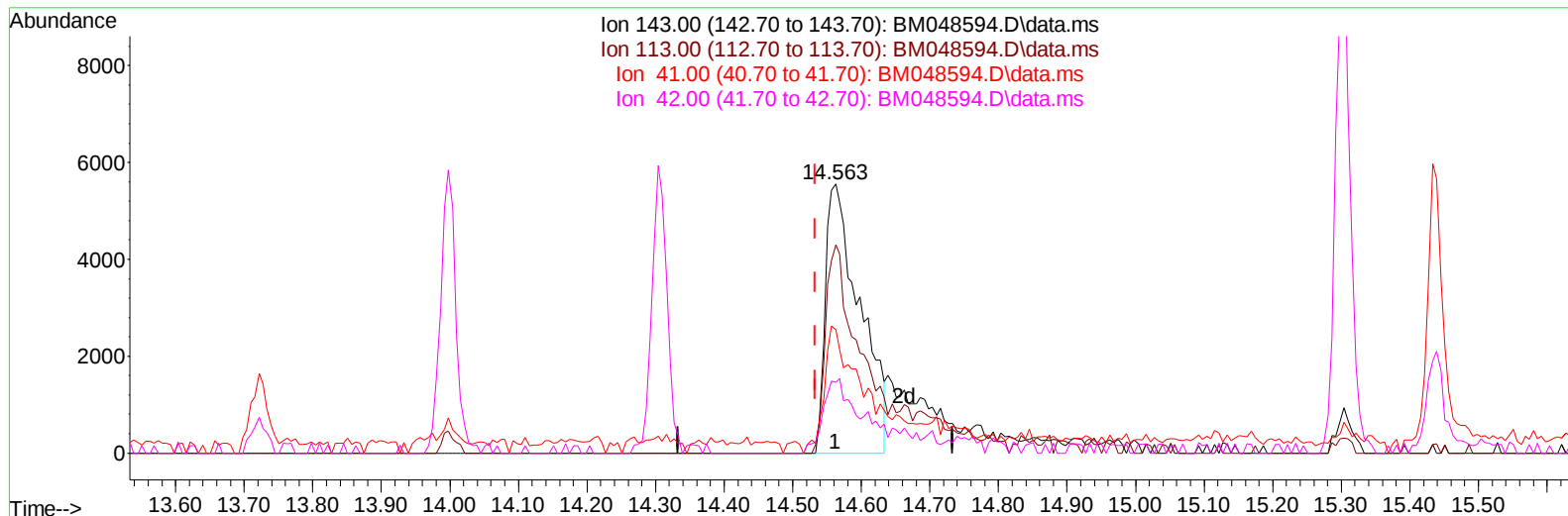
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111124\
Data File : BM048594.D
Acq On : 11 Nov 2024 21:01
Operator : RC/JU
Sample : P4714-12
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0AH3

Manual IntegrationsAPPROVED

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

Quant Time: Nov 11 22:55:15 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



TIC: BM048594.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.563min (+ 0.030) 4.39 ng/ul

response 19353

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	66.50	77.40
41.00	36.20	45.86#
42.00	24.50	26.40

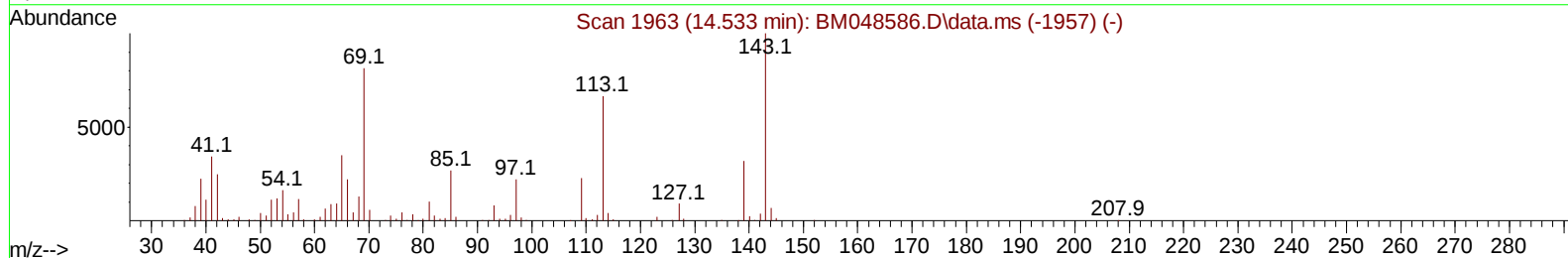
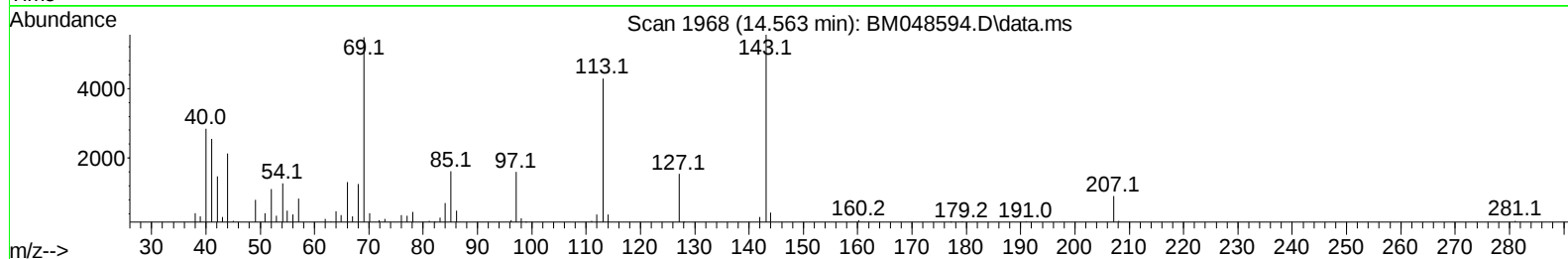
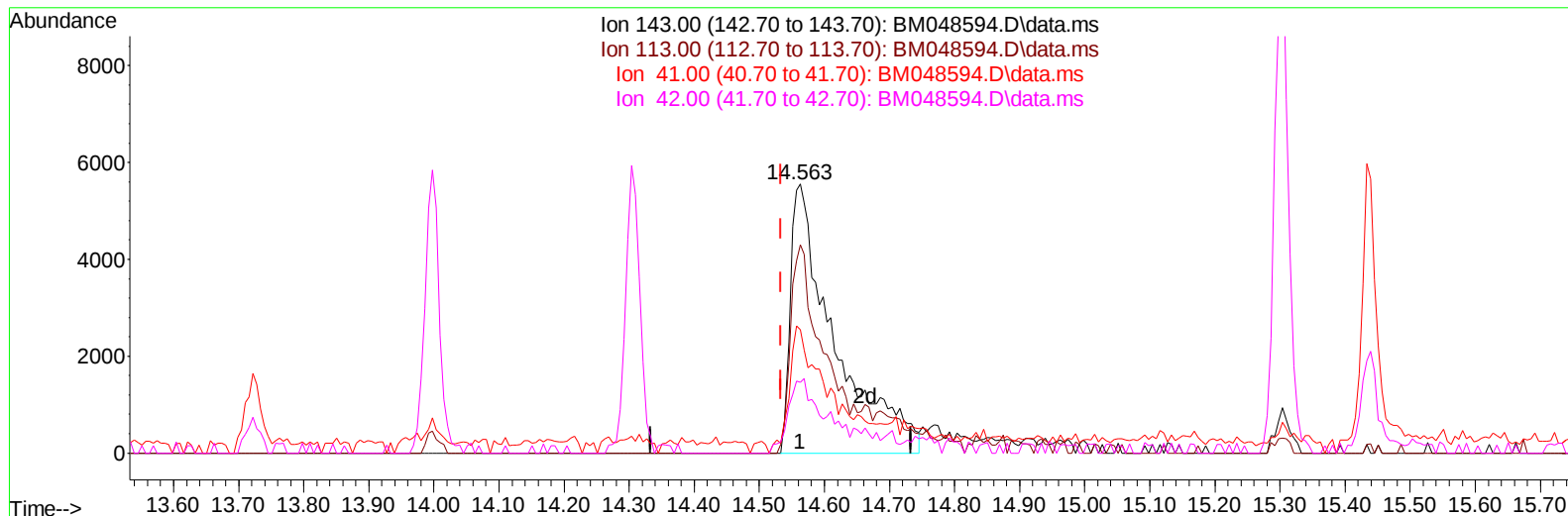
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TIC: BM048594.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.563min (+ 0.030) 5.85 ng/ul m

response 25813

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	66.50	77.40
41.00	36.20	45.86#
42.00	24.50	26.40

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Manual IntegrationsAPPROVED

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

Quant Time: Nov 11 23:08:37 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	151869	20.000	ng/ul	0.00
20) Naphthalene-d8	10.445	136	611475	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.310	164	372135	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.051	188	784626	20.000	ng/ul	0.00
79) Chrysene-d12	21.280	240	760857	20.000	ng/ul	0.00
88) Perylene-d12	24.186	264	891932	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.122	96	15475	4.062	ng/uL	0.00
4) Pyridine-d5	3.546	84	100322	9.556	ng/ul	0.00
7) Phenol-d5	6.845	99	92009	7.959	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.993	67	196733	28.332	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.193	132	260959	27.250	ng/ul	0.00
15) 4-Methylphenol-d8	8.375	113	171150	18.828	ng/ul	0.00
21) Nitrobenzene-d5	8.816	128	131335	29.634	ng/ul	0.00
24) 2-Nitrophenol-d4	9.534	143	142712	30.274	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.081	165	241600	27.589	ng/ul	0.00
31) 4-Chloroaniline-d4	10.592	131	377514	30.350	ng/ul	0.00
46) Dimethylphthalate-d6	13.722	166	831246	34.992	ng/ul	0.00
49) Acenaphthylene-d8	13.998	160	871985	31.134	ng/ul	0.00
54) 4-Nitrophenol-d4	14.563	143	25813m	5.850	ng/ul	0.03
60) Fluorene-d10	15.304	176	727605	35.089	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.439	200	108948	27.602	ng/ul	0.00
73) Anthracene-d10	17.151	188	1270008	38.250	ng/ul	0.00
81) Pyrene-d10	19.451	212	1539655	40.322	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.991	264	1628108	39.068	ng/ul	0.00

Target Compounds Qvalue

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

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