

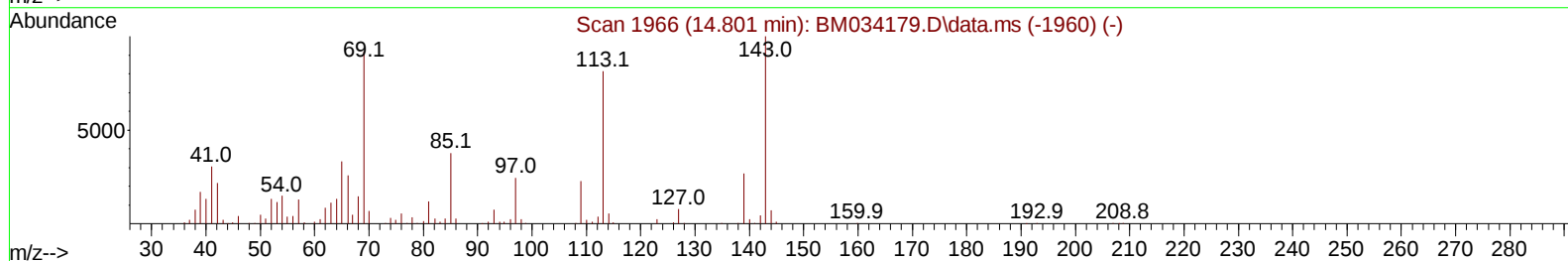
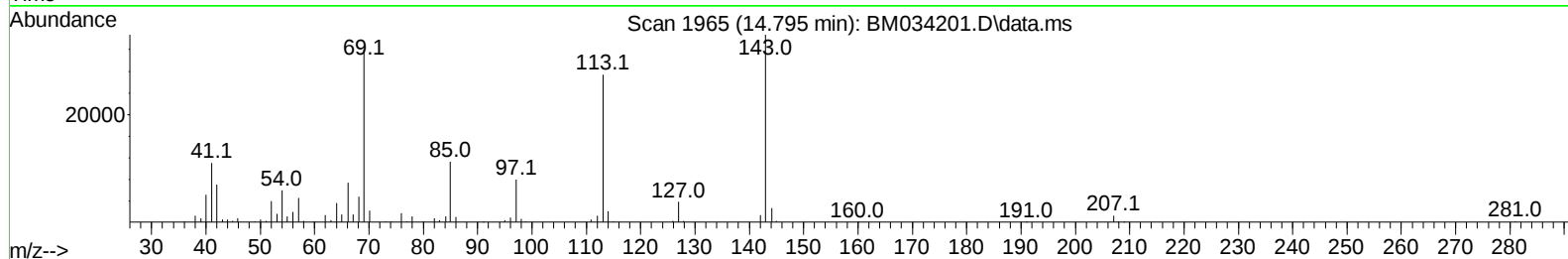
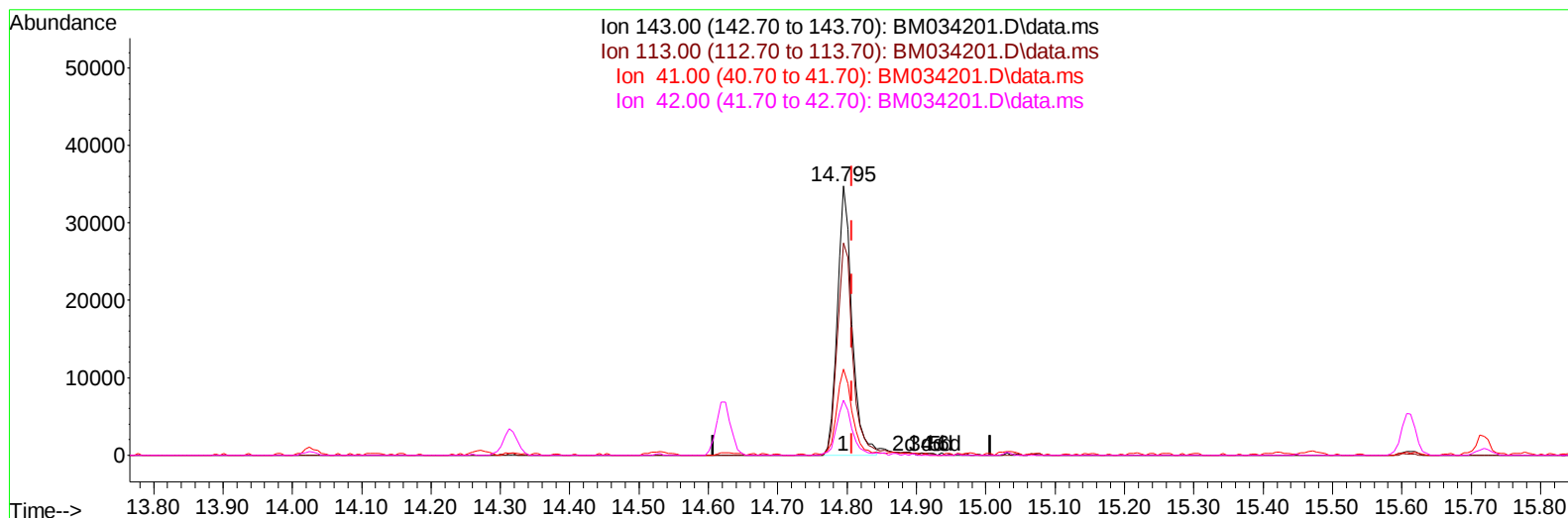
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030822\
Data File : BM034201.D
Acq On : 10 Mar 2022 00:48
Operator : CG/JU
Sample : N1775-17
Misc :
ALS Vial : 61 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DBSC0

Manual IntegrationsAPPROVED

Quant Time: Mar 10 01:22:34 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM030822.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Mar 08 13:09:16 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 03/10/2022
Supervised By :Yogesh Patel 03/11/2022



TIC: BM034201.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.795min (-0.012) 9.03 ng/u1

response 50532

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	105.00	78.87#
41.00	57.20	31.94#
42.00	39.50	20.43#

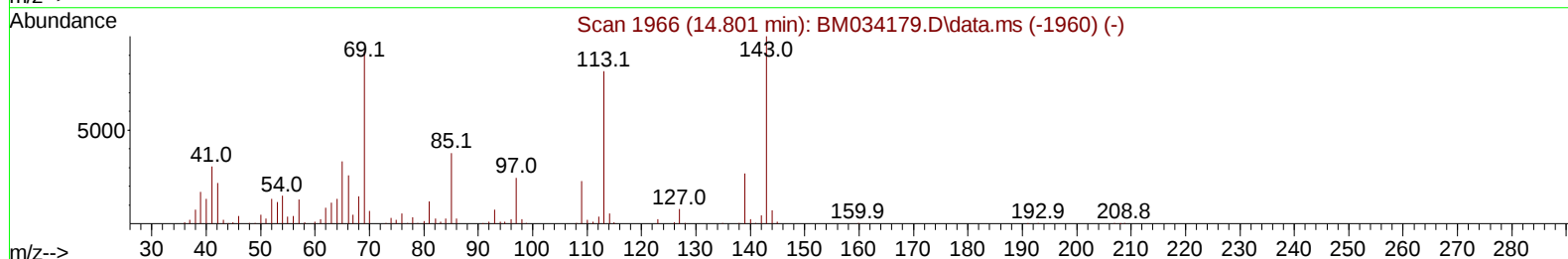
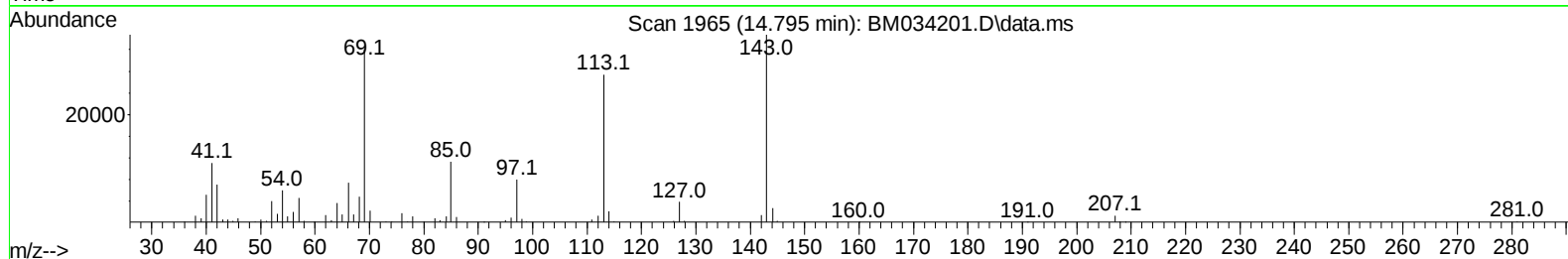
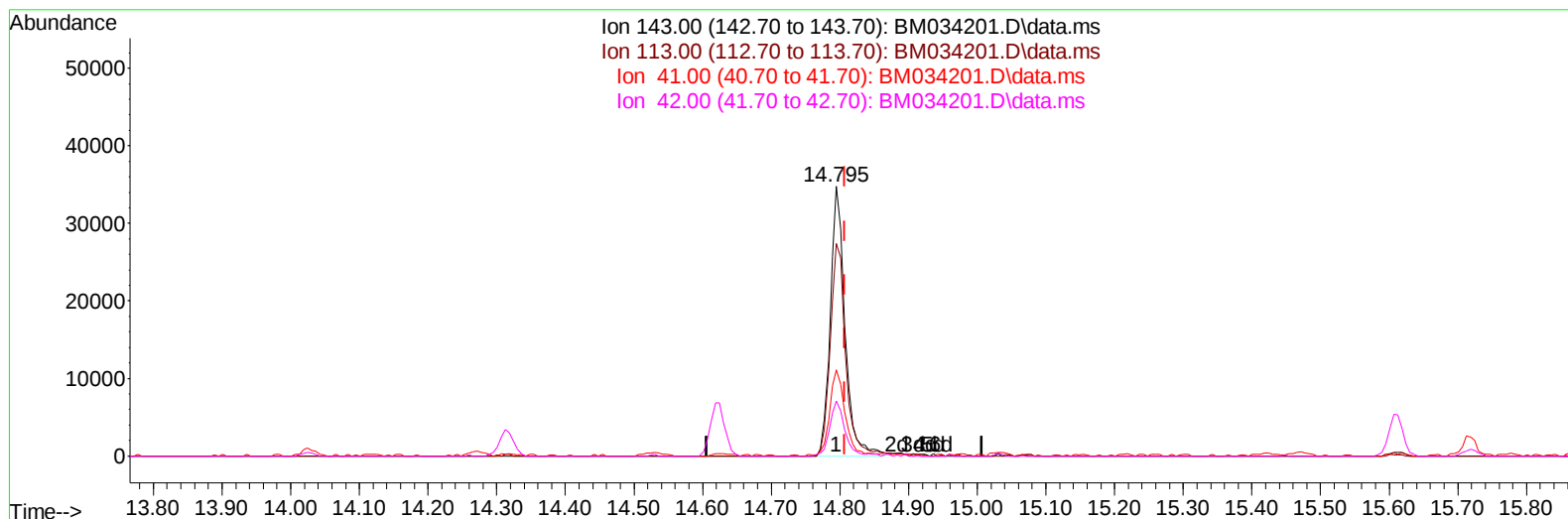
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TIC: BM034201.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.795min (-0.012) 9.18 ng/ul m

response 51406

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	105.00	78.87#
41.00	57.20	31.94#
42.00	39.50	20.43#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.990	152	147740	20.000	ng/ul	0.00
20) Naphthalene-d8	10.801	136	594434	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.624	164	396532	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.360	188	752785	20.000	ng/ul	#-0.01
79) Chrysene-d12	21.530	240	642195	20.000	ng/ul	#-0.01
88) Perylene-d12	23.971	264	688776	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.372	96	8121	2.528	ng/uL	0.00
4) Pyridine-d5	3.796	84	53086	5.677	ng/ul	0.00
7) Phenol-d5	7.137	99	137246	11.856	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.307	67	80978	12.996	ng/ul	0.00
11) 2-Chlorophenol-d4	7.513	132	120149	12.521	ng/ul	0.00
15) 4-Methylphenol-d8	8.690	113	117431	12.107	ng/ul	0.00
21) Nitrobenzene-d5	9.142	128	58982	12.806	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.872	143	63681	12.730	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.413	165	124224	12.042	ng/ul	0.00
31) 4-Chloroaniline-d4	10.925	131	155012	10.998	ng/ul	-0.01
46) Dimethylphthalate-d6	14.024	166	444427	14.770	ng/ul	-0.01
49) Acenaphthylene-d8	14.313	160	498843	13.257	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.795	143	51406m	9.181	ng/ul	-0.01
60) Fluorene-d10	15.613	176	370742	13.844	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.719	200	52003	10.705	ng/ul	0.00
73) Anthracene-d10	17.460	188	583582	15.738	ng/ul	0.00
81) Pyrene-d10	19.748	212	648876	18.382	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.812	264	609103	16.671	ng/ul	-0.02

Target Compounds Qvalue

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

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