

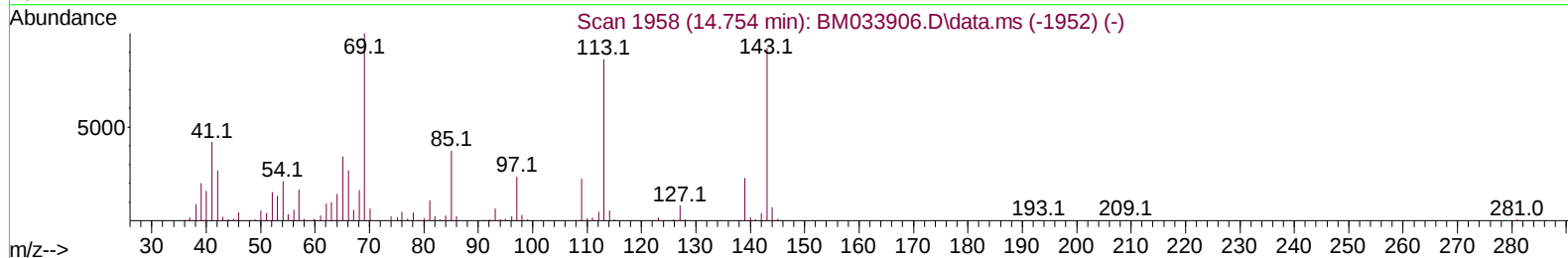
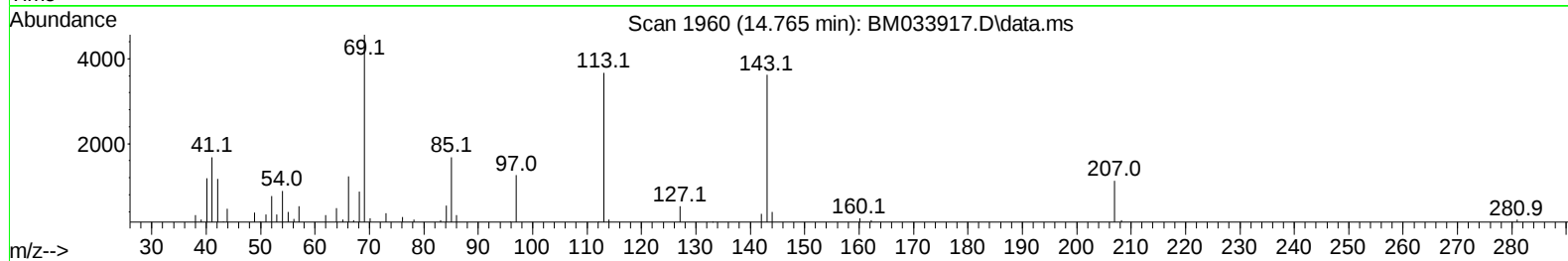
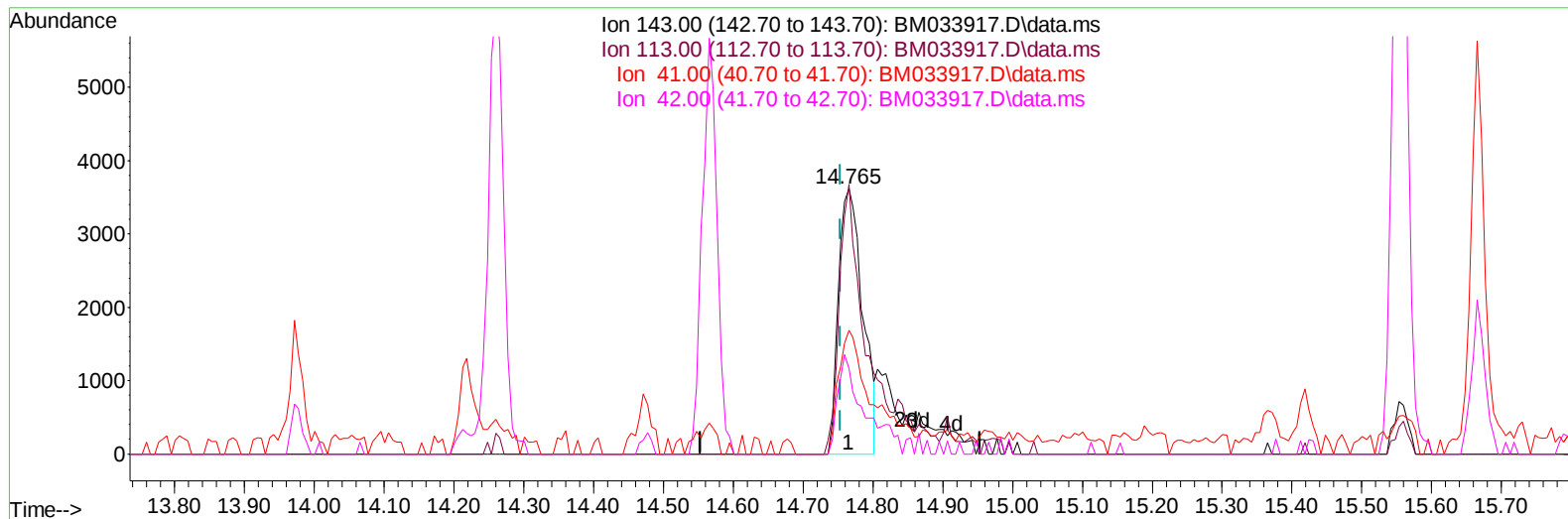
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM012822\
Data File : BM033917.D
Acq On : 28 Jan 2022 23:04
Operator : CG/JU
Sample : N1302-07
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0F24

Manual IntegrationsAPPROVED

Quant Time: Jan 29 00:17:21 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM010622.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Jan 28 23:26:29 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 01/31/2022
Supervised By :mohammad ahmed 02/01/2022



TIC: BM033917.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.765min (+ 0.012) 2.35 ng/u1

response 8647

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	105.00	101.49
41.00	57.20	46.56
42.00	39.50	32.23

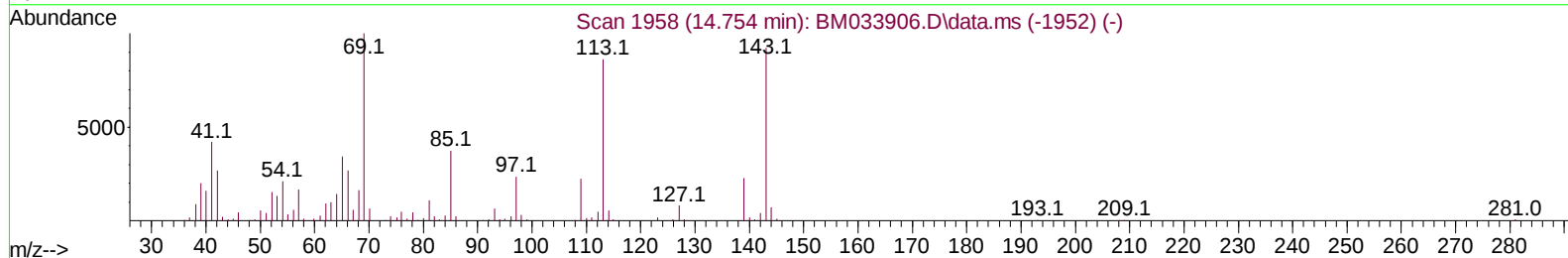
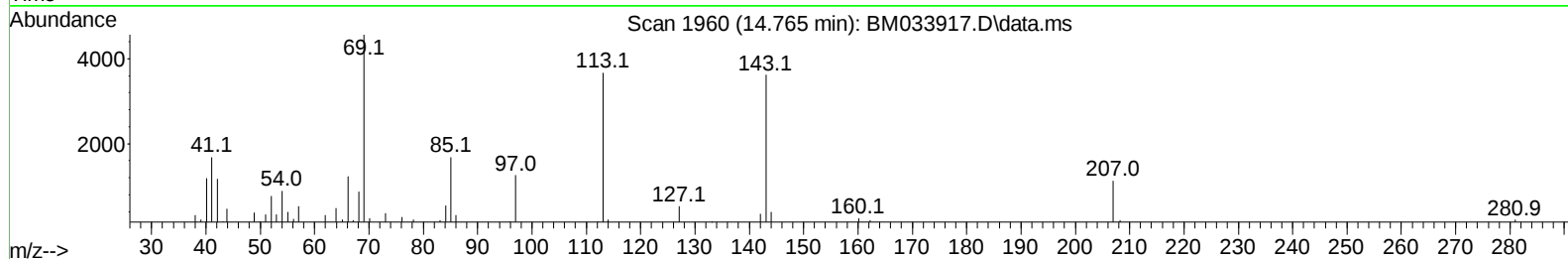
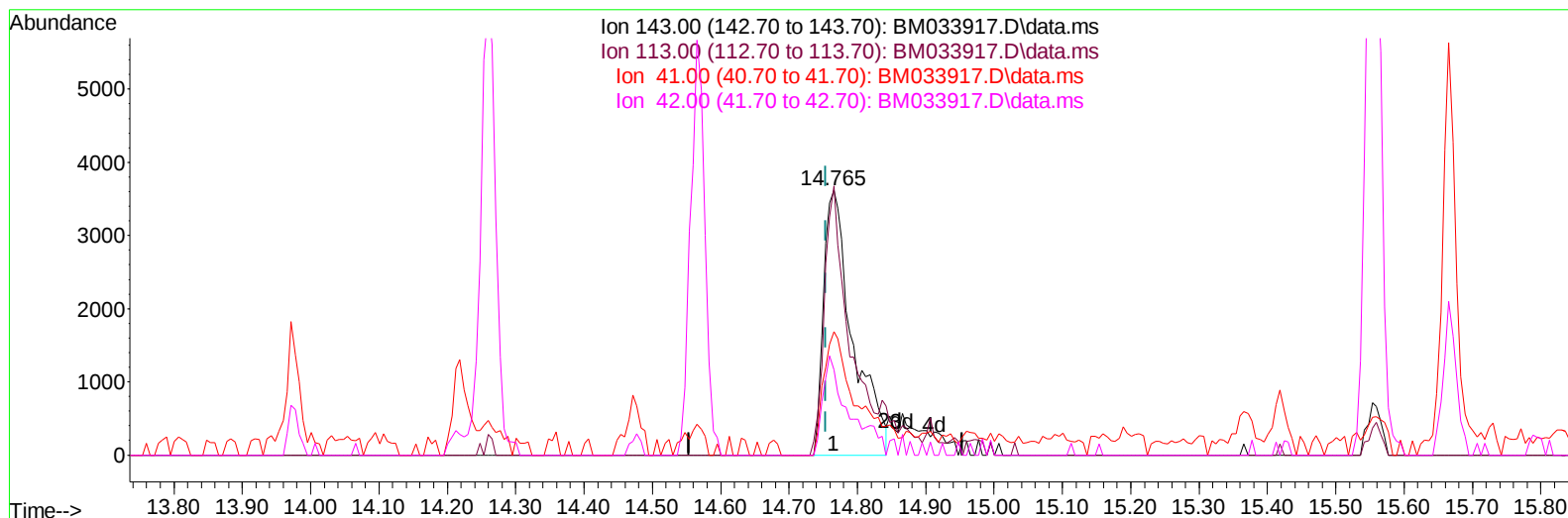
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TIC: BM033917.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.765min (+ 0.012) 2.90 ng/ul m

response 10683

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	105.00	101.49
41.00	57.20	46.56
42.00	39.50	32.23

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.931	152	94920	20.000	ng/ul	0.00
20) Naphthalene-d8	10.742	136	388234	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.565	164	243308	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.301	188	506901	20.000	ng/ul	0.00
79) Chrysene-d12	21.465	240	500456	20.000	ng/ul	0.00
88) Perylene-d12	23.865	264	518678	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.290	96	9253	3.662	ng/uL	0.00
4) Pyridine-d5	3.719	84	58547	8.345	ng/ul	0.00
7) Phenol-d5	7.084	99	44794	5.254	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.254	67	146483	28.215	ng/ul	0.00
11) 2-Chlorophenol-d4	7.454	132	145965	22.394	ng/ul	0.00
15) 4-Methylphenol-d8	8.636	113	85574	12.126	ng/ul	0.00
21) Nitrobenzene-d5	9.089	128	91056	29.699	ng/ul	0.00
24) 2-Nitrophenol-d4	9.819	143	94572	29.104	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.360	165	153817	24.692	ng/ul	0.00
31) 4-Chloroaniline-d4	10.872	131	230703	24.914	ng/ul	0.00
46) Dimethylphthalate-d6	13.977	166	607001	31.665	ng/ul	0.00
49) Acenaphthylene-d8	14.260	160	685074	30.292	ng/ul	0.00
54) 4-Nitrophenol-d4	14.765	143	10683m	2.904	ng/ul	0.01
60) Fluorene-d10	15.554	176	518323	32.591	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.665	200	80537	26.710	ng/ul	0.00
73) Anthracene-d10	17.401	188	896813	36.461	ng/ul	0.00
81) Pyrene-d10	19.689	212	1051968	38.924	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.706	264	1013736	37.563	ng/ul	0.00

Target Compounds Qvalue

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

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