

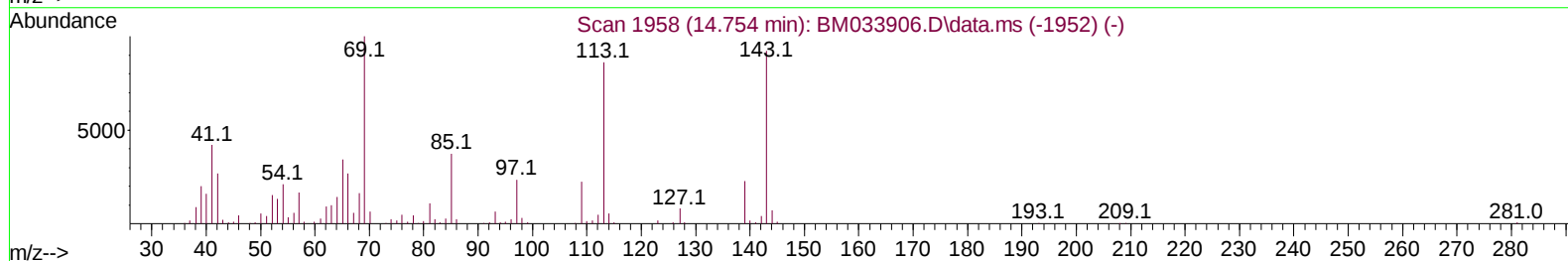
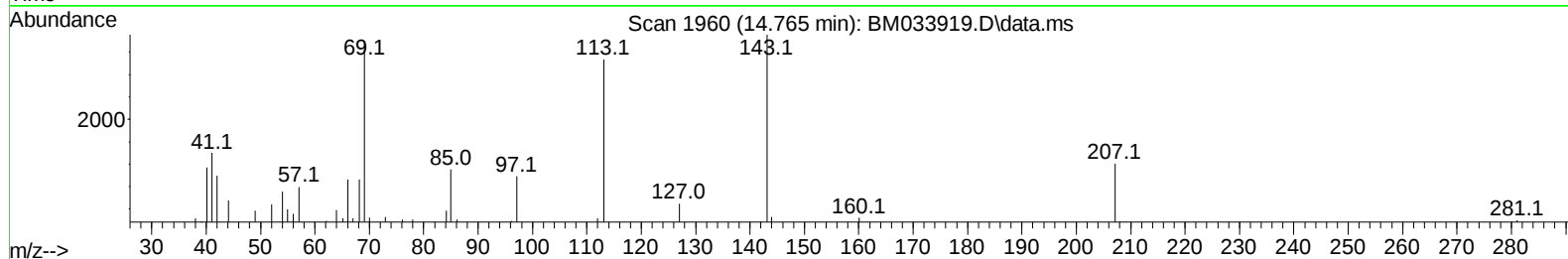
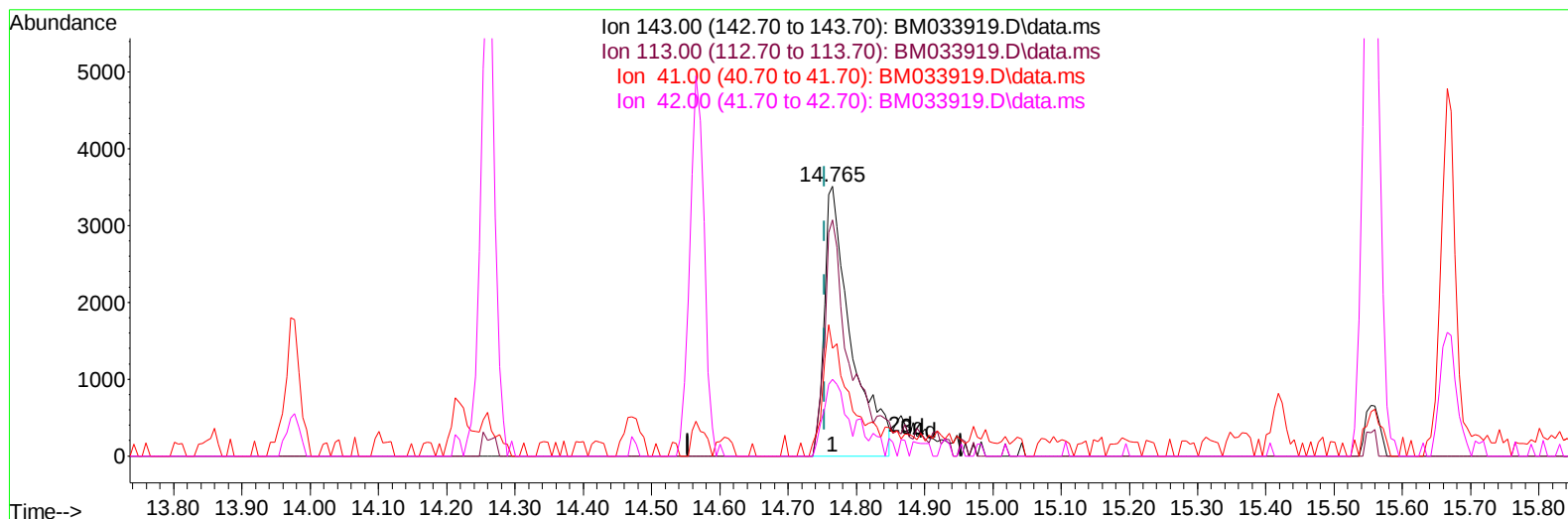
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM012822\
 Data File : BM033919.D
 Acq On : 29 Jan 2022 00:15
 Operator : CG/JU
 Sample : N1302-06
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0F23

Manual IntegrationsAPPROVED

Quant Time: Jan 29 00:53:16 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: BM033919.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.765min (+ 0.012) 2.88 ng/ul m

response 9487

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	105.00	87.52
41.00	57.20	39.93#
42.00	39.50	28.36#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.931	152	87382	20.000	ng/ul	0.00
20) Naphthalene-d8	10.742	136	353222	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.565	164	218128	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.301	188	457097	20.000	ng/ul	0.00
79) Chrysene-d12	21.465	240	460250	20.000	ng/ul	0.00
88) Perylene-d12	23.859	264	476831	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.290	96	9668	4.157	ng/uL	0.00
4) Pyridine-d5	3.719	84	59389	9.195	ng/ul	0.00
7) Phenol-d5	7.084	99	42311	5.391	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.254	67	152663	31.942	ng/ul	0.00
11) 2-Chlorophenol-d4	7.460	132	149836	24.971	ng/ul	0.00
15) 4-Methylphenol-d8	8.636	113	83955	12.923	ng/ul	0.00
21) Nitrobenzene-d5	9.089	128	94828	33.995	ng/ul	0.00
24) 2-Nitrophenol-d4	9.819	143	96716	32.715	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.360	165	157696	27.824	ng/ul	0.00
31) 4-Chloroaniline-d4	10.872	131	221941	26.344	ng/ul	0.00
46) Dimethylphthalate-d6	13.977	166	613205	35.681	ng/ul	0.00
49) Acenaphthylene-d8	14.260	160	701381	34.593	ng/ul	0.00
54) 4-Nitrophenol-d4	14.765	143	9487m	2.876	ng/ul	0.01
60) Fluorene-d10	15.554	176	518841	36.390	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.665	200	73193	26.919	ng/ul	0.00
73) Anthracene-d10	17.401	188	858244	38.694	ng/ul	0.00
81) Pyrene-d10	19.689	212	1027782	41.351	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.706	264	1025359	41.328	ng/ul	0.00

Target Compounds Qvalue

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

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