

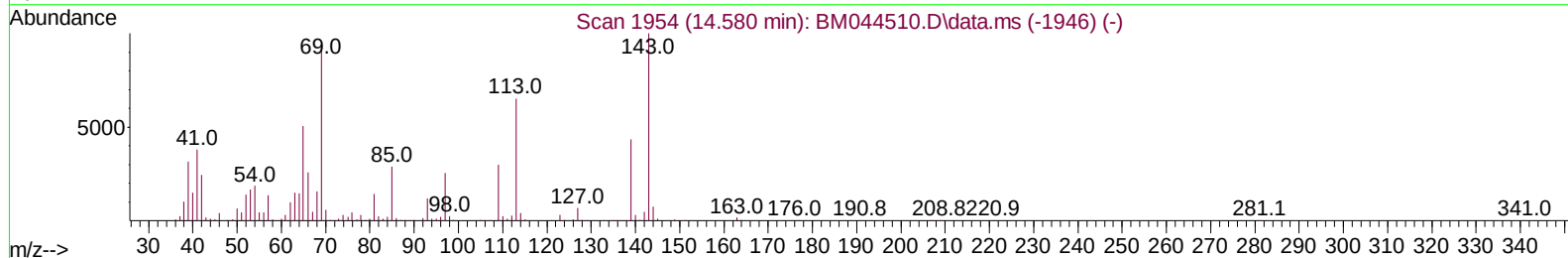
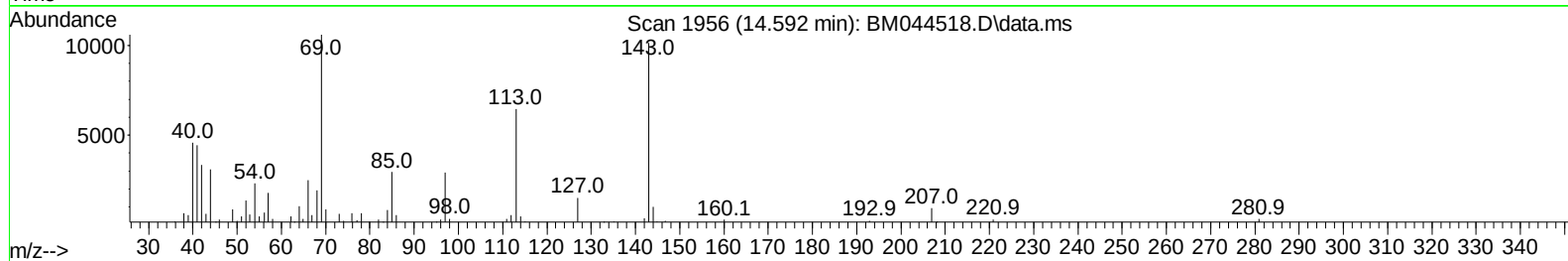
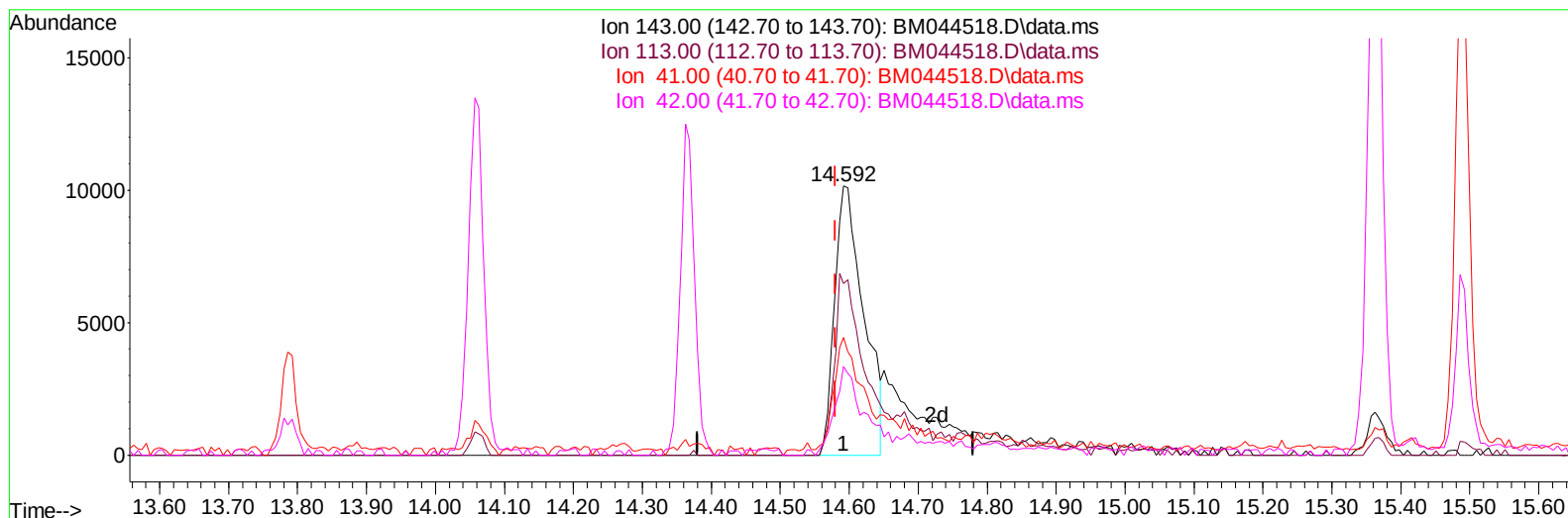
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030624\
Data File : BM044518.D
Acq On : 06 Mar 2024 16:33
Operator : MA/JU
Sample : P1642-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0R31

Manual IntegrationsAPPROVED

Quant Time: Mar 06 17:19:00 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM030524.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Mar 05 23:05:11 2024
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 03/07/2024
Supervised By :mohammad ahmed 03/09/2024



TIC: BM044518.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.592min (+ 0.012) 3.12 ng/ul

response 29693

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	63.70	63.70
41.00	38.40	43.63
42.00	26.60	33.01#

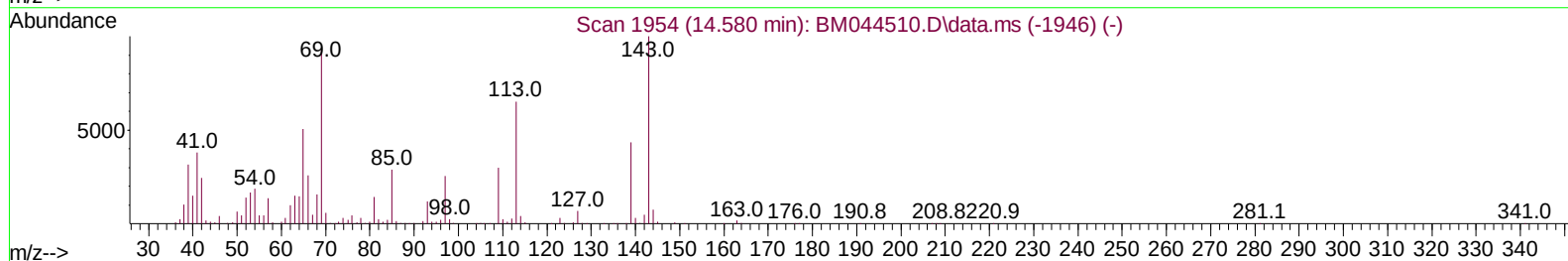
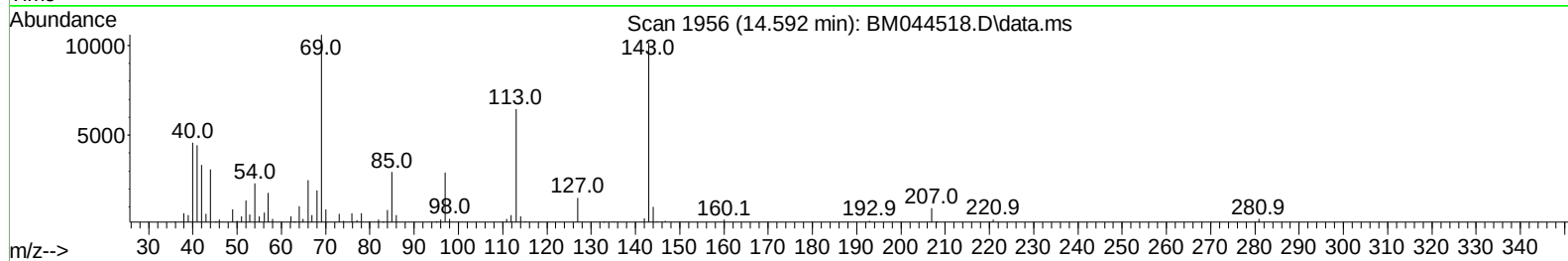
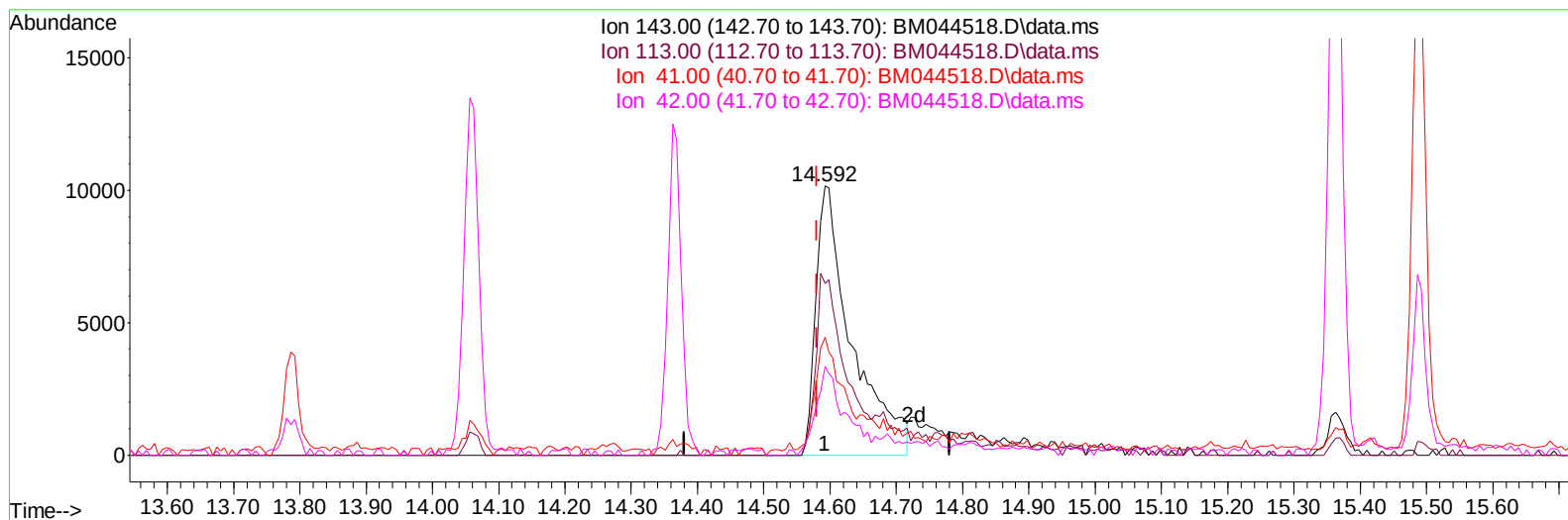
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(54) 4-Nitrophenol-d4 (S)

14.592min (+ 0.012) 4.00 ng/ul m

response 37979

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	63.70	63.70
41.00	38.40	43.63
42.00	26.60	33.01#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.734	152	215950	20.000	ng/ul	0.00
20) Naphthalene-d8	10.510	136	910930	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.369	164	584139	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.115	188	1233703	20.000	ng/ul	0.00
79) Chrysene-d12	21.303	240	1111072	20.000	ng/ul	0.00
88) Perylene-d12	23.556	264	1251504	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.287	96	27216	4.275	ng/uL	0.00
4) Pyridine-d5	3.693	84	83324	4.446	ng/ul	0.01
7) Phenol-d5	6.904	99	121496	5.472	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.081	67	414262	30.207	ng/ul	0.00
11) 2-Chlorophenol-d4	7.263	132	369055	22.620	ng/ul	0.00
15) 4-Methylphenol-d8	8.434	113	231201	13.260	ng/ul	0.00
21) Nitrobenzene-d5	8.892	128	240927	31.409	ng/ul	0.00
24) 2-Nitrophenol-d4	9.604	143	247789	29.966	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.133	165	406570	28.889	ng/ul	0.00
31) 4-Chloroaniline-d4	10.663	131	183021	7.830	ng/ul	0.00
46) Dimethylphthalate-d6	13.786	166	1584443	34.578	ng/ul	0.00
49) Acenaphthylene-d8	14.063	160	1735045	33.109	ng/ul	0.00
54) 4-Nitrophenol-d4	14.592	143	37979m	3.997	ng/ul	0.01
60) Fluorene-d10	15.363	176	1396358	36.263	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.492	200	280087	34.975	ng/ul	0.00
73) Anthracene-d10	17.215	188	2547838	43.206	ng/ul	0.00
81) Pyrene-d10	19.515	212	2942678	43.182	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.415	264	2816627	42.818	ng/ul	0.00

Target Compounds Qvalue

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

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