

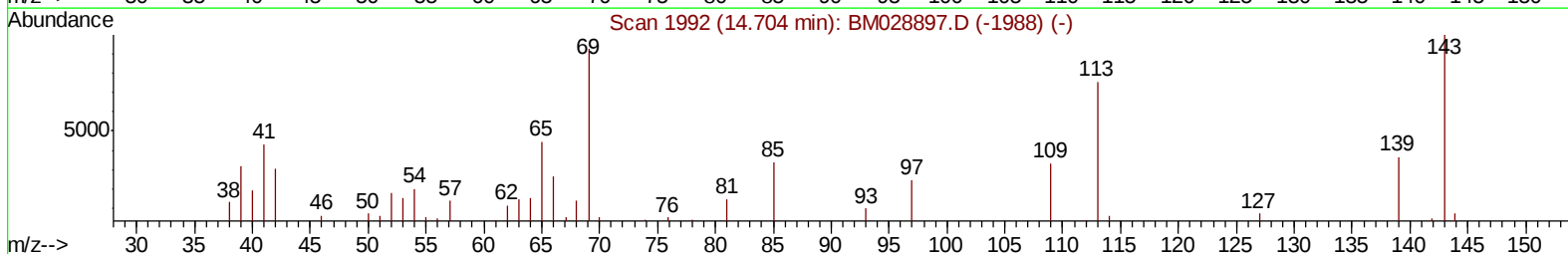
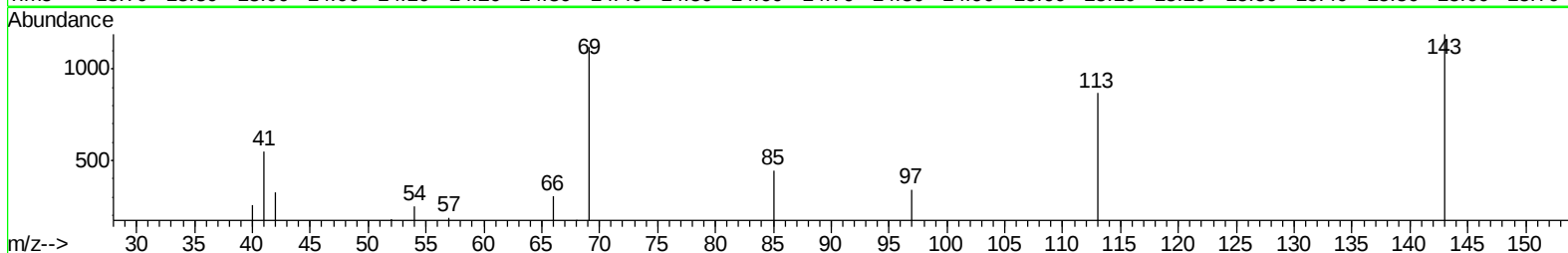
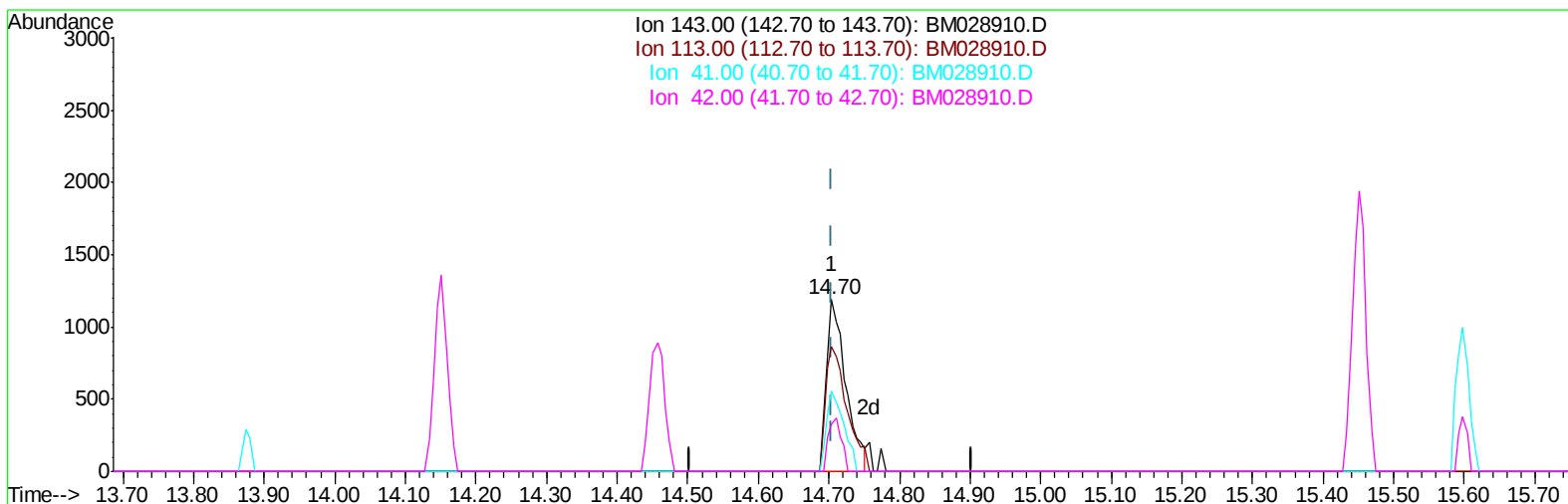
Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM022621\
 Data File : BM028910.D
 Acq On : 25 Feb 2021 23:18
 Operator : CG/JU
 Sample : M1482-08
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBG6M6

Manual Integrations
 APPROVED

mohammad
 2/26/2021 10:43:14 AM

Quant Time: Feb 26 00:12:57 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM022121MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Feb 25 23:08:07 2021
 Response via : Initial Calibration



TIC: BM028910.D

(52) 4-Nitrophenol-d4 (S)
 14.704min (+0.000) 4.90ng/ul
 response 2267

Ion	Exp%	Act%
143.00	100	100
113.00	75.60	73.00
41.00	43.10	46.43
42.00	30.60	27.50

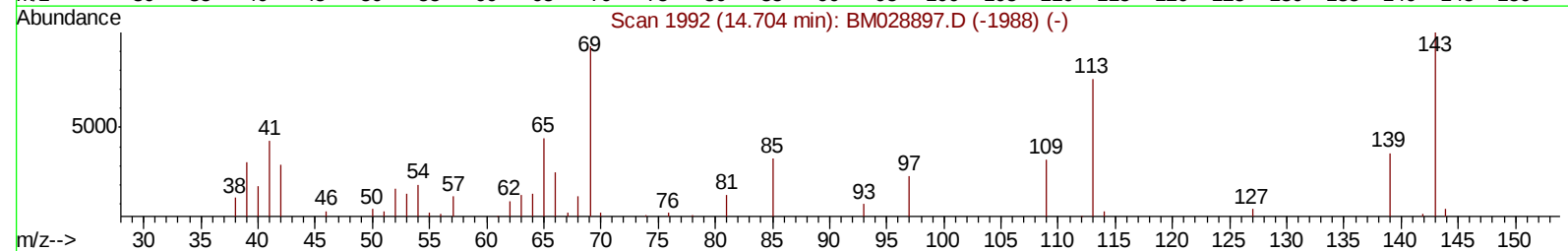
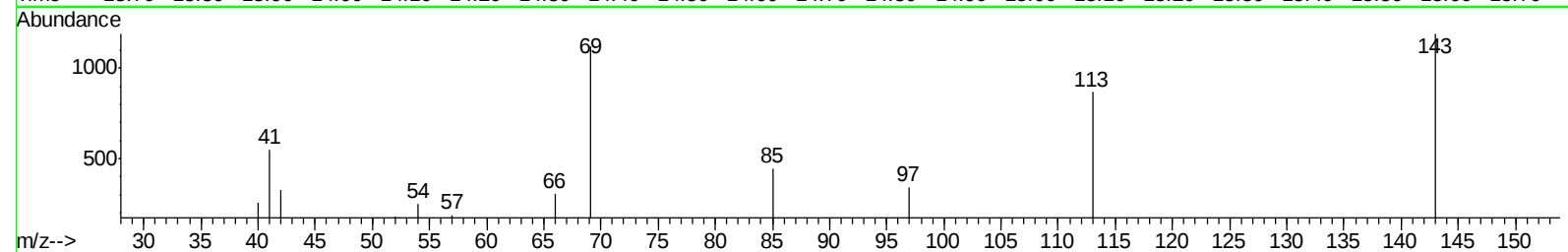
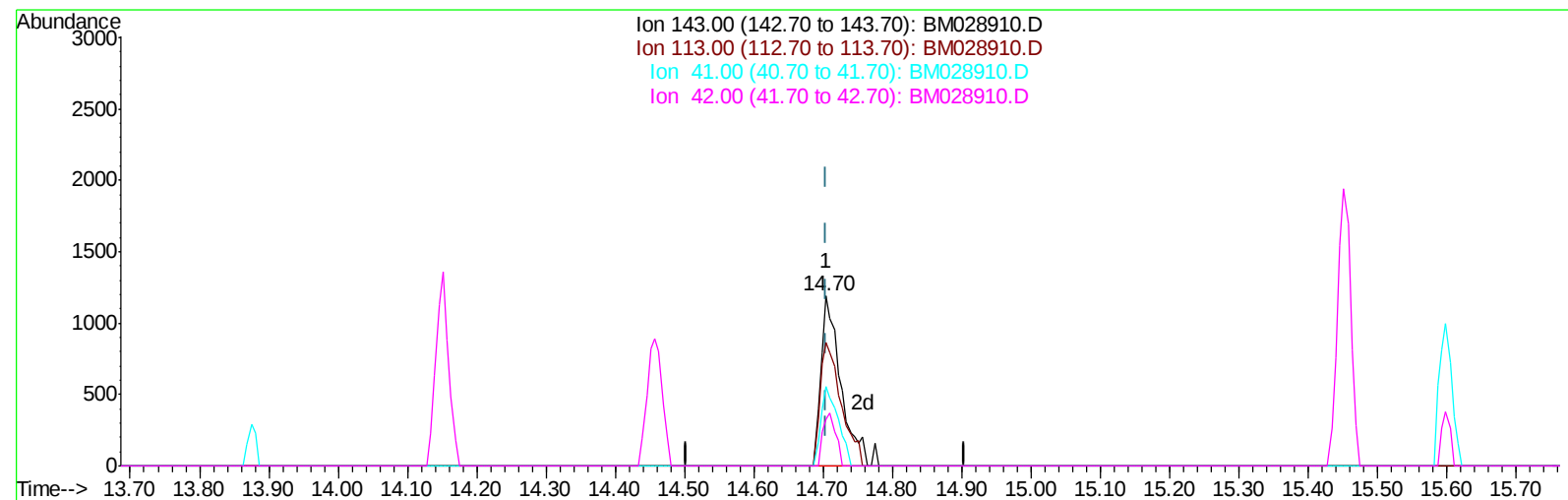
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mohammad
2/26/2021 10:43:14 AM



TIC: BM028910.D

(52) 4-Nitrophenol-d4 (S)

14.704min (+0.000) 5.05ng/ul m

response 2339

Ion	Exp%	Act%
143.00	100	100
113.00	75.60	73.00
41.00	43.10	46.43
42.00	30.60	27.50

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.80	152	14650	20.00	ng/ul	0.00
18) Naphthalene-d8	10.60	136	60042	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.46	164	37091	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.20	188	71945	20.00	ng/ul	0.00
78) Chrysene-d12	21.36	240	56023	20.00	ng/ul	0.00
86) Perylene-d12	23.57	264	51693	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.15	96	1511	4.91	ng/uL	0.00
5) Phenol-d5	6.98	99	7716	6.92	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.13	67	20652	32.84	ng/ul	0.00
9) 2-Chlorophenol-d4	7.33	132	25928	27.05	ng/ul	0.00
13) 4-Methylphenol-d8	8.53	113	15001	16.02	ng/ul	0.00
19) Nitrobenzene-d5	8.97	128	15175	35.71	ng/ul	0.00
22) 2-Nitrophenol-d4	9.69	143	17415	36.11	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.23	165	27569	30.05	ng/ul	0.00
29) 4-Chloroaniline-d4	10.76	131	1978	1.82	ng/ul	0.00
44) Dimethylphthalate-d6	13.87	166	96481	36.08	ng/ul	0.00
47) Acenaphthylene-d8	14.15	160	120701	35.45	ng/ul	0.00
52) 4-Nitrophenol-d4	14.70	143	2339m	5.05	ng/ul	0.00
58) Fluorene-d10	15.45	176	84543	37.61	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.60	200	14609	33.20	ng/ul	0.00
71) Anthracene-d10	17.30	188	136842	41.76	ng/ul	0.00
79) Pyrene-d10	19.59	212	133550	47.83	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.43	264	111817	41.97	ng/ul	-0.01

Target Compounds	Ovalue

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

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