

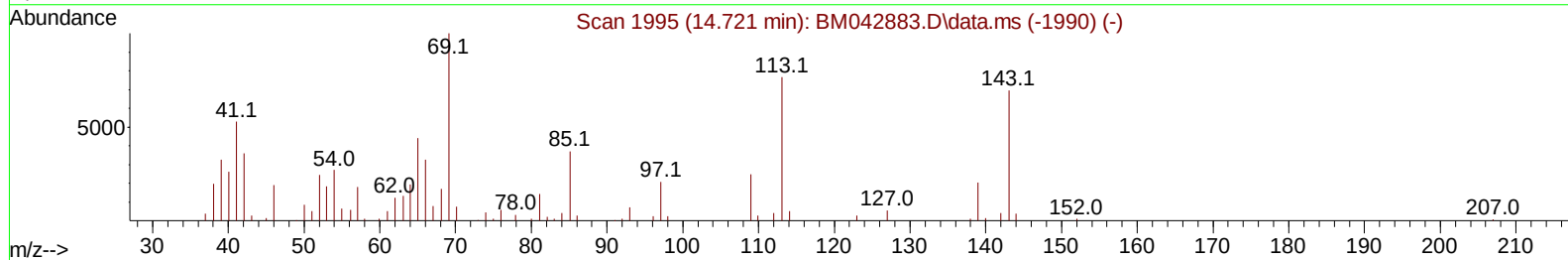
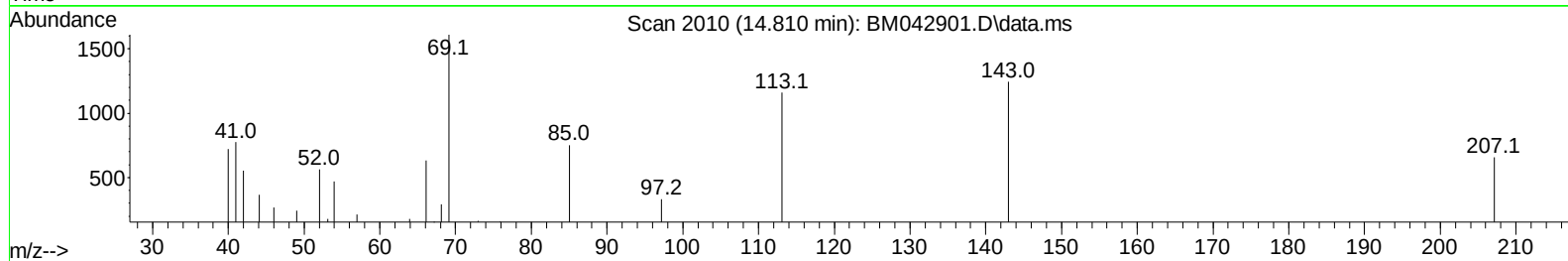
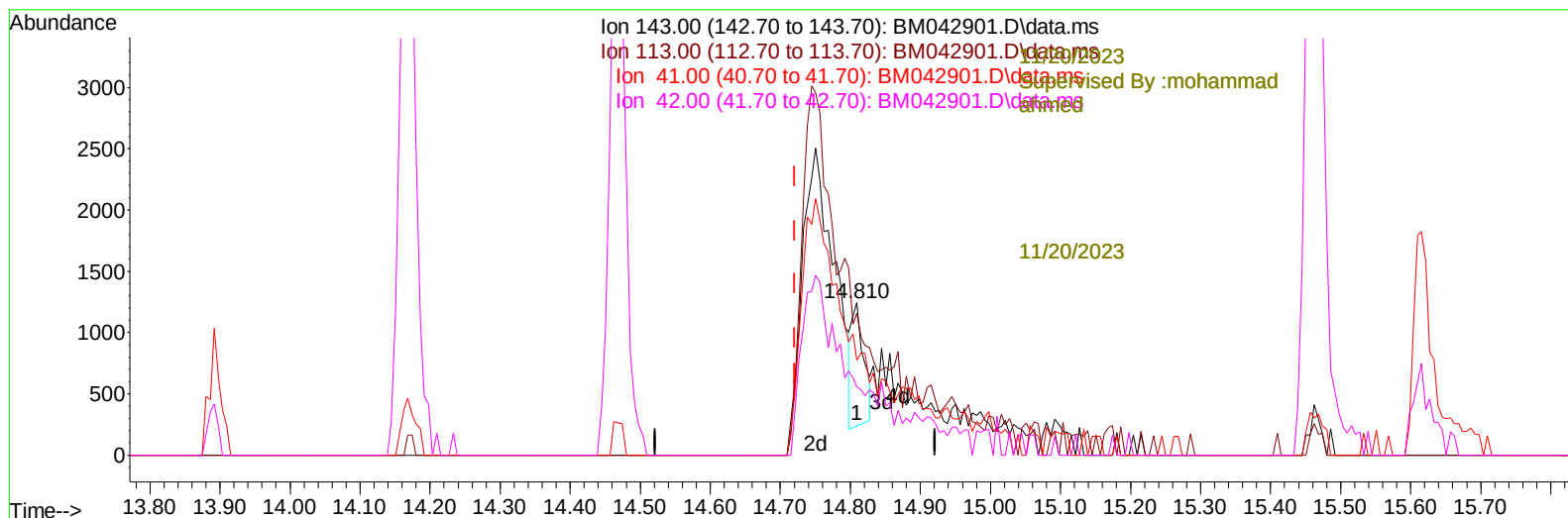
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111723\
Data File : BM042901.D
Acq On : 19 Nov 2023 12:00
Operator : MA/JU
Sample : PB157215BL
Misc :
ALS Vial : 58 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
SBLK215

Manual IntegrationsAPPROVED

Quant Time: Nov 19 21:48:02 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM111323.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat Nov 18 22:06:09 2023
Response via : Initial Calibration

Reviewed By :Yogesh Patel 11/20/2023
Supervised By :mohammad ahmed 11/20/2023



TIC: BM042901.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.810min (+ 0.088) 1.52 ng/ul

response 1196

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	110.80	93.32
41.00	73.50	62.40
42.00	48.00	44.69

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Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)
-----							-11/20/2023
Internal Standards							Supervised By :mohammad
1) 1,4-Dichlorobenzene-d4	7.810	152		31036	20.000	ng/ul	0.00
20) Naphthalene-d8	10.622	136		120606	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.468	164		73581	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.221	188		159218	20.000	ng/ul	0.00
79) Chrysene-d12	21.409	240		157127	20.000	ng/ul	0.00
88) Perylene-d12	23.762	264		205569	20.000	ng/ul	0.00
System Monitoring Compounds							11/20/2023
3) 1,4-Dioxane-d8	3.169	96		5309	5.367	ng/uL	0.00
4) Pyridine-d5	3.604	84		62638	25.459	ng/ul	0.00
7) Phenol-d5	6.981	99		78005	25.825	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.163	67		61566	29.006	ng/ul	0.00
11) 2-Chlorophenol-d4	7.334	132		58830	25.241	ng/ul	0.00
15) 4-Methylphenol-d8	8.533	113		56601	23.799	ng/ul	0.00
21) Nitrobenzene-d5	9.016	128		25436	24.055	ng/ul	0.00
24) 2-Nitrophenol-d4	9.727	143		28463	23.010	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.251	165		48248	20.977	ng/ul	0.00
31) 4-Chloroaniline-d4	10.804	131		57768	19.781	ng/ul	0.00
46) Dimethylphthalate-d6	13.892	166		182218	26.166	ng/ul	0.00
49) Acenaphthylene-d8	14.168	160		195768	26.274	ng/ul	0.00
54) 4-Nitrophenol-d4	14.751	143		13042m	16.576	ng/ul	0.03
60) Fluorene-d10	15.462	176		158750	27.529	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.615	200		18204	15.094	ng/ul	0.00
73) Anthracene-d10	17.321	188		237093	26.571	ng/ul	0.00
81) Pyrene-d10	19.621	212		311875	30.311	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.609	264		330863	26.908	ng/ul	0.00

Target Compounds Qvalue

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

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