

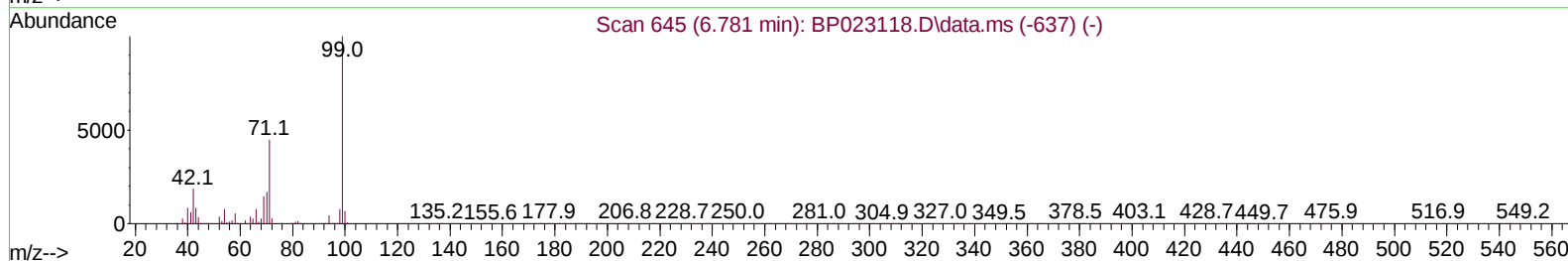
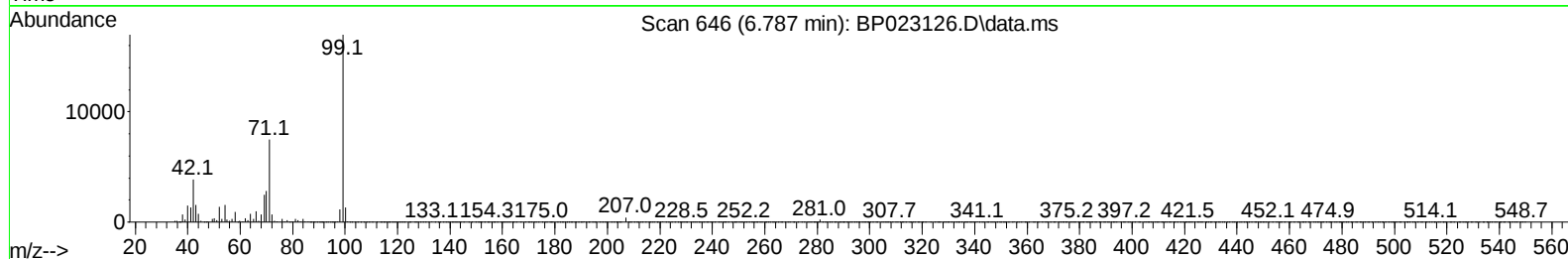
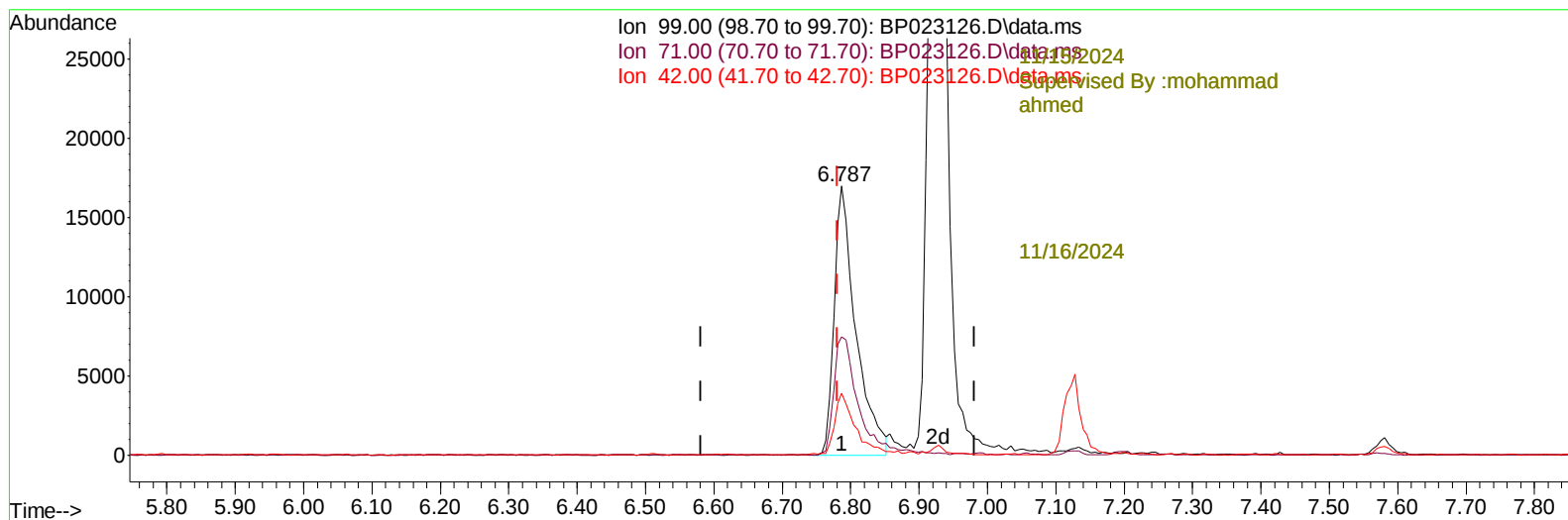
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111324\
Data File : BP023126.D
Acq On : 15 Nov 2024 05:57
Operator : RC/JU
Sample : P4740-05
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0CC2

Manual IntegrationsAPPROVED

Quant Time: Nov 15 07:42:12 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110924.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Nov 13 04:49:01 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 11/15/2024
Supervised By :mohammad ahmed 11/18/2024



TIC: BP023126.D\data.ms

(7) Phenol-d5 (S)

6.787min (+ 0.006) 7.95 ng/u1

response 37230

Ion	Exp%	Act%
99.00	100.00	100.00
71.00	44.70	43.96
42.00	21.10	22.89
0.00	0.00	0.00

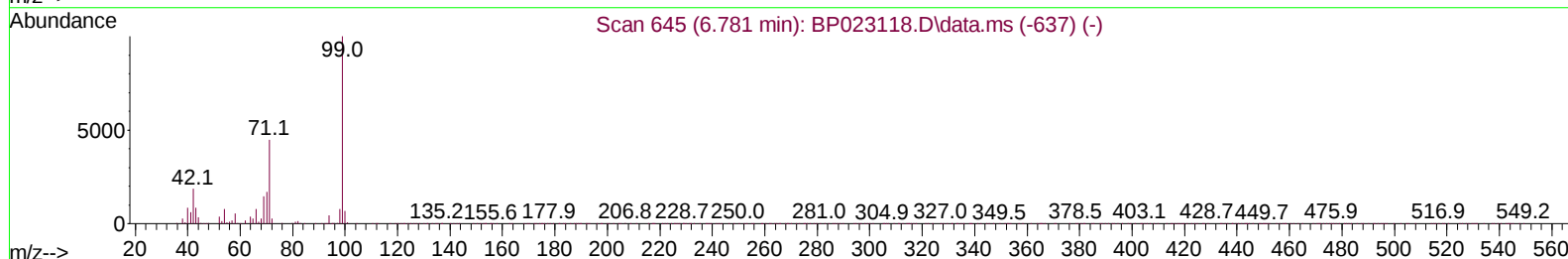
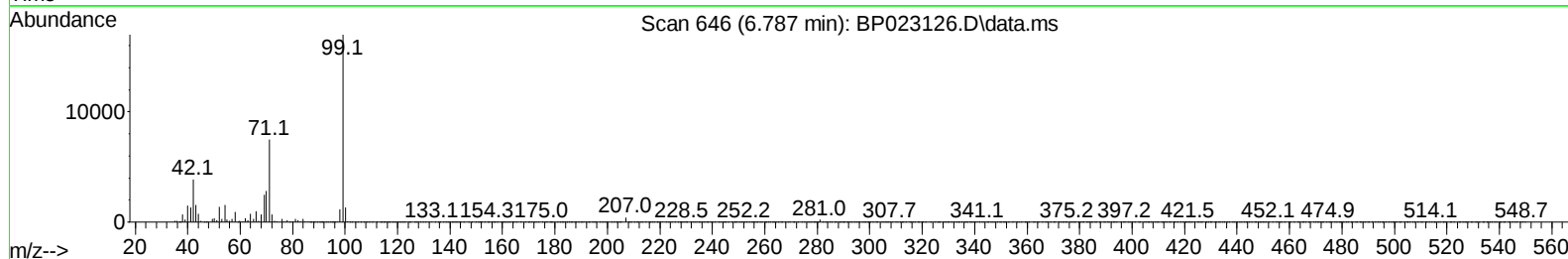
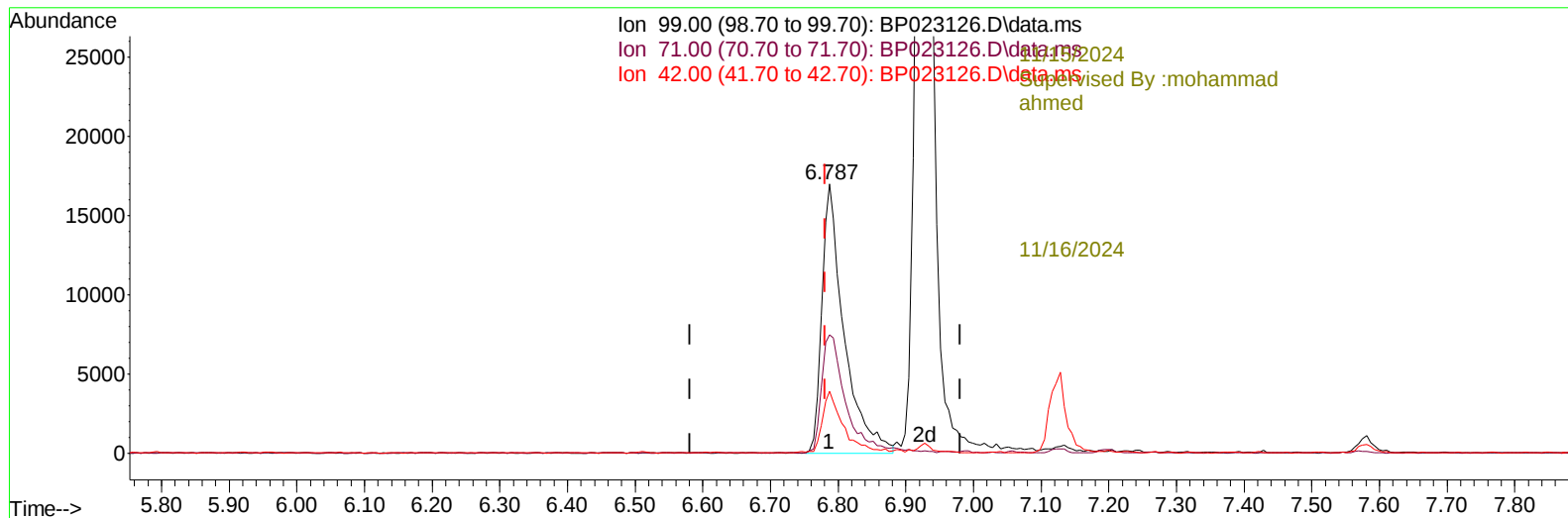
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TIC: BP023126.D\data.ms

(7) Phenol-d5 (S)

6.787min (+ 0.006) 8.25 ng/u1 m

response 38621

Ion	Exp%	Act%
99.00	100.00	100.00
71.00	44.70	43.96
42.00	21.10	22.89
0.00	0.00	0.00

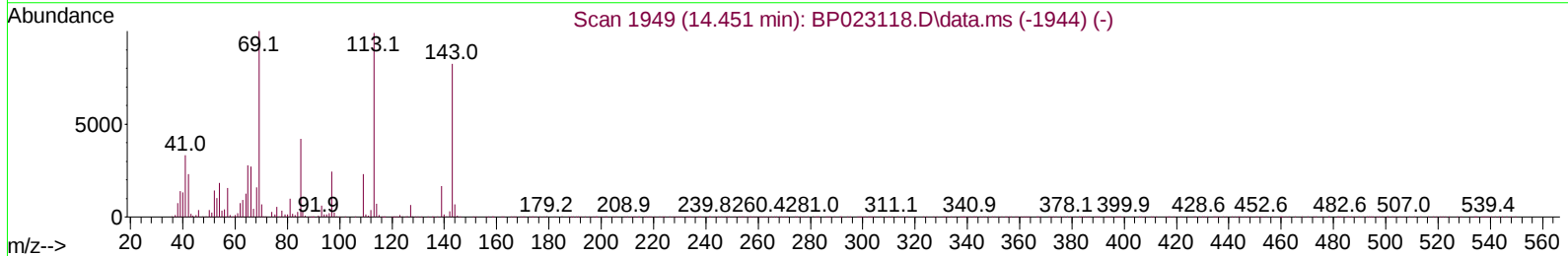
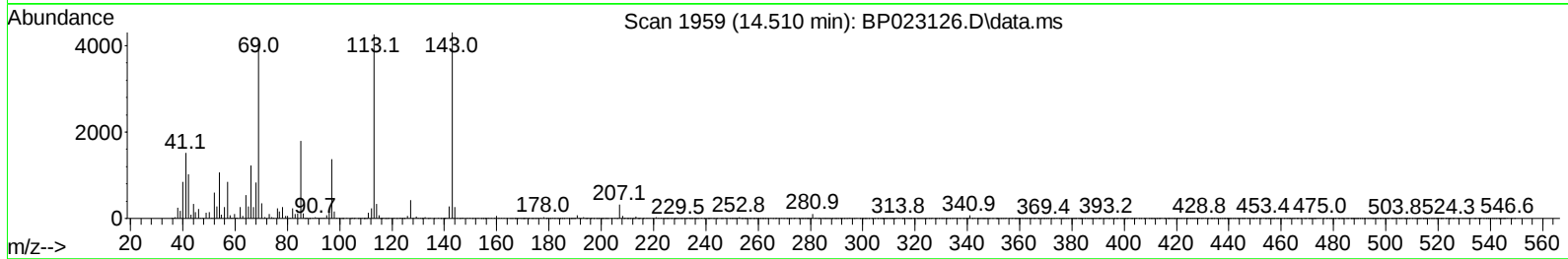
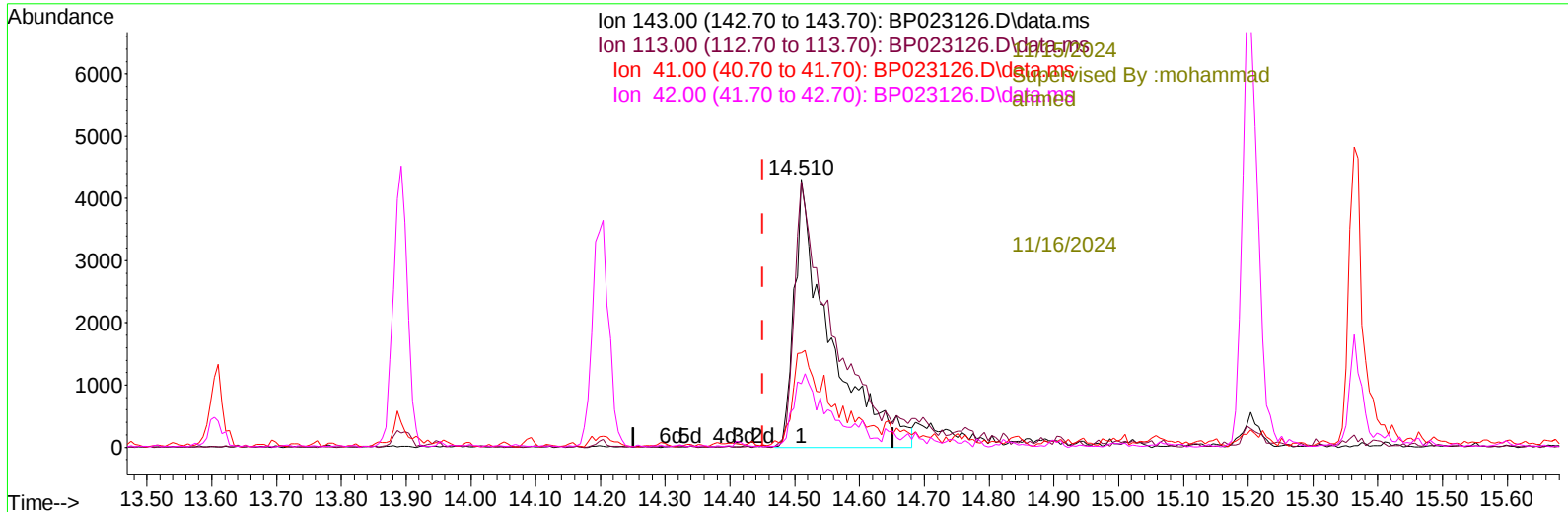
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ALS Vial : 50 Sample Multiplier: 1

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TIC: BP023126.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.510min (+ 0.059) 7.82 ng/ul m

response 16443

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	99.09
41.00	39.20	35.24
42.00	25.90	23.72

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 Misc :
 ALS Vial : 50 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0CC2

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 11/15/2024
 Supervised By :mohammad ahmed 11/18/2024

Quant Time: Nov 15 07:43:39 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110924.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Nov 13 04:49:01 2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
-----11/15/2024						
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.575	152	63311	20.000	ng/ul	0.00
20) Naphthalene-d8	10.328	136	238502	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.198	164	194147	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.004	188	500913	20.000	ng/ul	-0.01
79) Chrysene-d12	21.439	240	600109	20.000	ng/ul	0.00
88) Perylene-d12	24.657	264	702657	20.000	ng/ul	0.00
-----11/16/2024						
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.134	96	6443	4.892	ng/uL	0.00
4) Pyridine-d5	3.558	84	21338	5.453	ng/ul	0.01
7) Phenol-d5	6.787	99	38621m	8.251	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.928	67	91086	33.122	ng/ul	0.00
11) 2-Chlorophenol-d4	7.128	132	98428	28.712	ng/ul	0.00
15) 4-Methylphenol-d8	8.293	113	72792	18.898	ng/ul	0.00
21) Nitrobenzene-d5	8.722	128	57515	34.639	ng/ul	0.00
24) 2-Nitrophenol-d4	9.434	143	67123	33.909	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.975	165	135278	32.633	ng/ul	0.00
31) 4-Chloroaniline-d4	10.487	131	86083	17.406	ng/ul	0.00
46) Dimethylphthalate-d6	13.610	166	545355	38.974	ng/ul	0.00
49) Acenaphthylene-d8	13.892	160	503715	34.500	ng/ul	0.00
54) 4-Nitrophenol-d4	14.510	143	16443m	7.819	ng/ul	0.06
60) Fluorene-d10	15.204	176	462899	37.732	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.363	200	104290	34.755	ng/ul	0.00
73) Anthracene-d10	17.116	188	875264	41.268	ng/ul	0.00
81) Pyrene-d10	19.486	212	1191868	43.119	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.445	264	1394105	41.652	ng/ul	0.00

Target Compounds Qvalue

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

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