

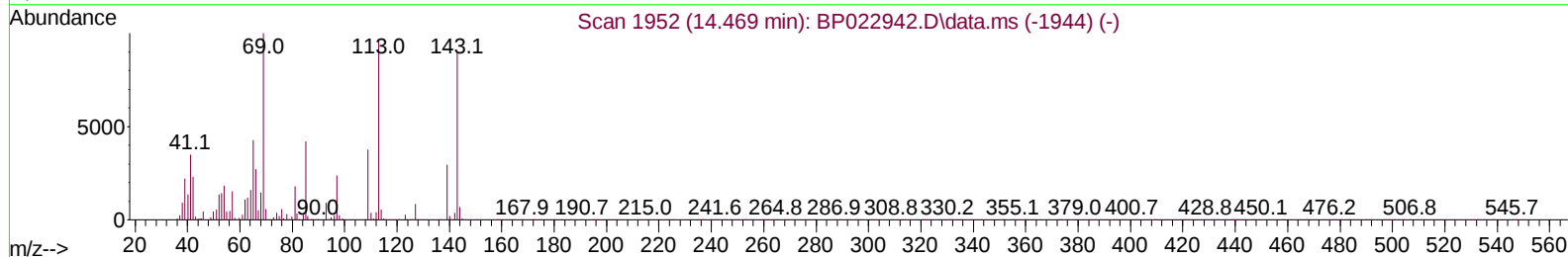
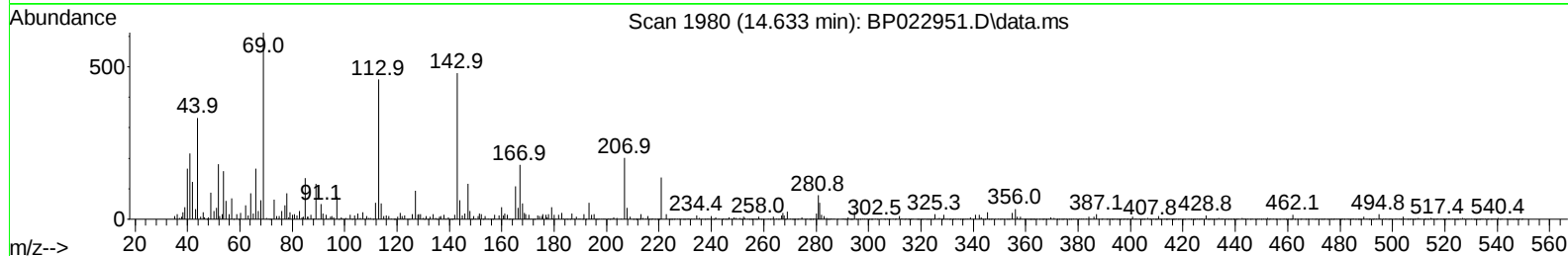
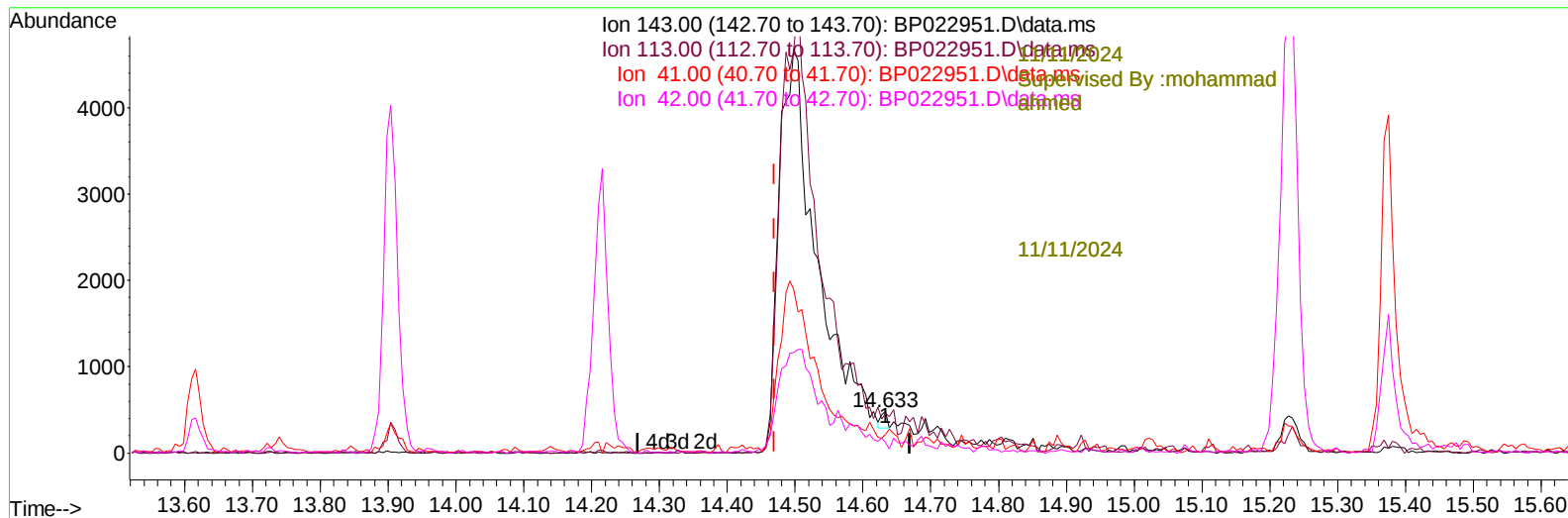
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110924\
Data File : BP022951.D
Acq On : 09 Nov 2024 04:26
Operator : RC/JU
Sample : P4744-09
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCP63

Manual IntegrationsAPPROVED

Quant Time: Nov 09 05:16:34 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110924.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Nov 08 23:49:48 2024
Response via : Initial Calibration

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



TIC: BP022951.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.633min (+ 0.165) 0.08 ng/u1

response 121

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	95.62
41.00	39.20	44.89
42.00	25.90	25.68

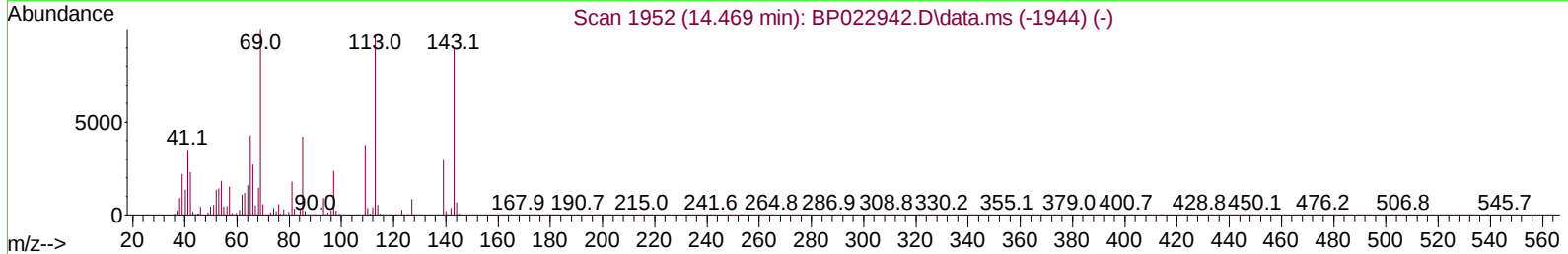
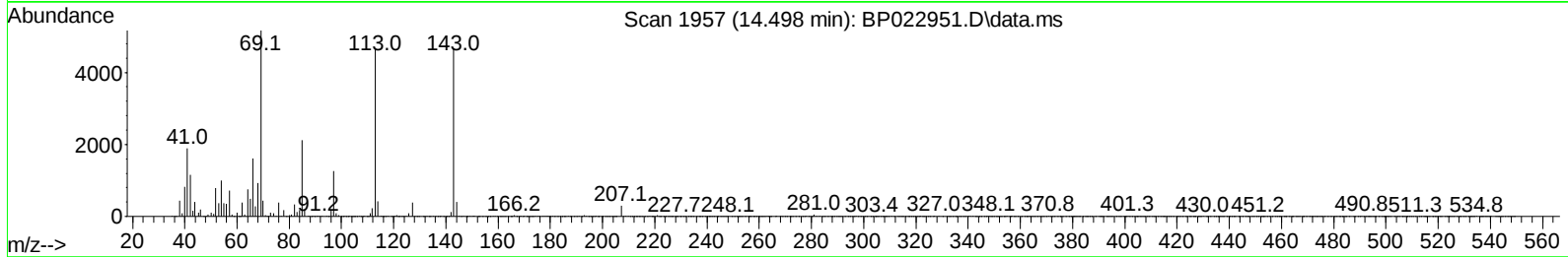
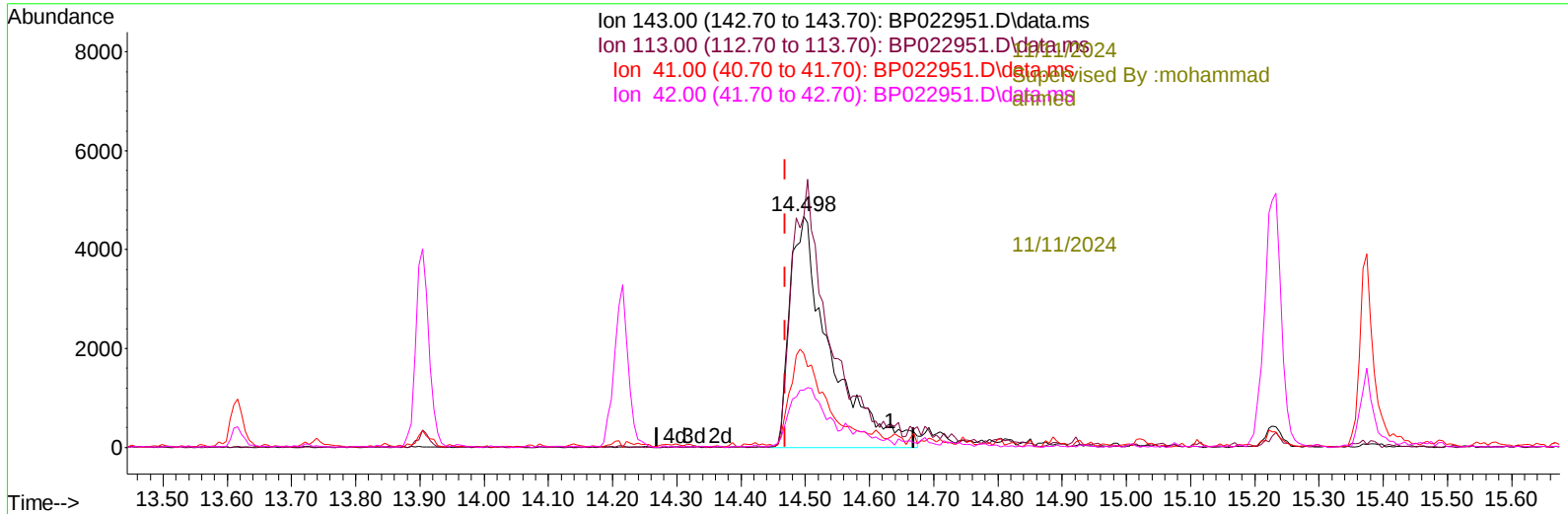
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TIC: BP022951.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.498min (+ 0.029) 13.45 ng/ul m

response 20363

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	99.83
41.00	39.20	40.58
42.00	25.90	24.88

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 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
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Manual IntegrationsAPPROVED

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

Quant Time: Nov 09 05:17:36 2024
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 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Nov 08 23:49:48 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
-----11/11/2024						
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.593	152	47613	20.000	ng/ul	0.00
20) Naphthalene-d8	10.345	136	178788	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.216	164	139812	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.016	188	356679	20.000	ng/ul	-0.02
79) Chrysene-d12	21.451	240	445756	20.000	ng/ul	0.00
88) Perylene-d12	24.680	264	558656	20.000	ng/ul	0.0011/11/2024
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.146	96	7189	7.258	ng/uL	0.00
4) Pyridine-d5	3.575	84	14577	4.953	ng/ul	0.02
7) Phenol-d5	6.798	99	44941	12.767	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.940	67	79369	38.377	ng/ul	0.00
11) 2-Chlorophenol-d4	7.140	132	95816	37.166	ng/ul	0.00
15) 4-Methylphenol-d8	8.304	113	79647	27.495	ng/ul	0.00
21) Nitrobenzene-d5	8.734	128	48245	38.761	ng/ul	0.00
24) 2-Nitrophenol-d4	9.445	143	58983	39.749	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.992	165	112332	36.149	ng/ul	0.00
31) 4-Chloroaniline-d4	10.498	131	58062	15.662	ng/ul	0.00
46) Dimethylphthalate-d6	13.616	166	434910	43.160	ng/ul	0.00
49) Acenaphthylene-d8	13.904	160	411253	39.114	ng/ul	0.00
54) 4-Nitrophenol-d4	14.498	143	20363m	13.447	ng/ul	0.03
60) Fluorene-d10	15.227	176	364492	41.257	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.375	200	81214	38.009	ng/ul	0.00
73) Anthracene-d10	17.127	188	650857	43.097	ng/ul	0.00
81) Pyrene-d10	19.492	212	882399	42.977	ng/ul	-0.02
92) Benzo(a)pyrene-d12	24.468	264	1156628	43.464	ng/ul	0.00

Target Compounds Qvalue

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

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