

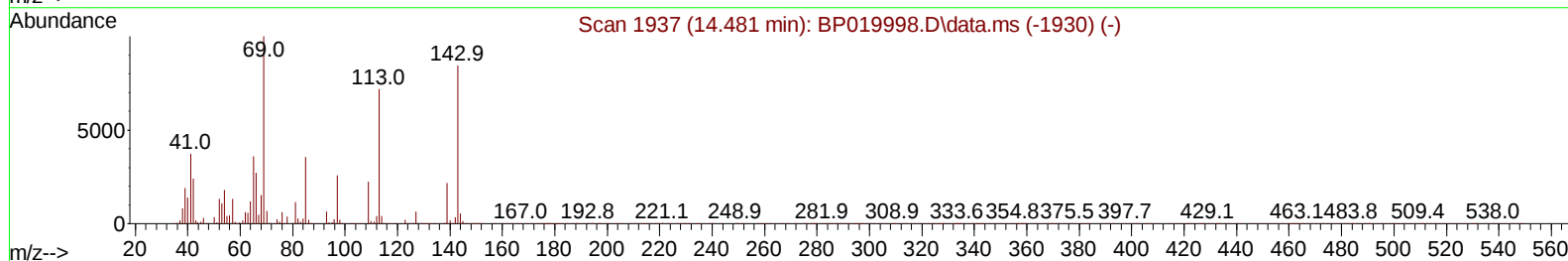
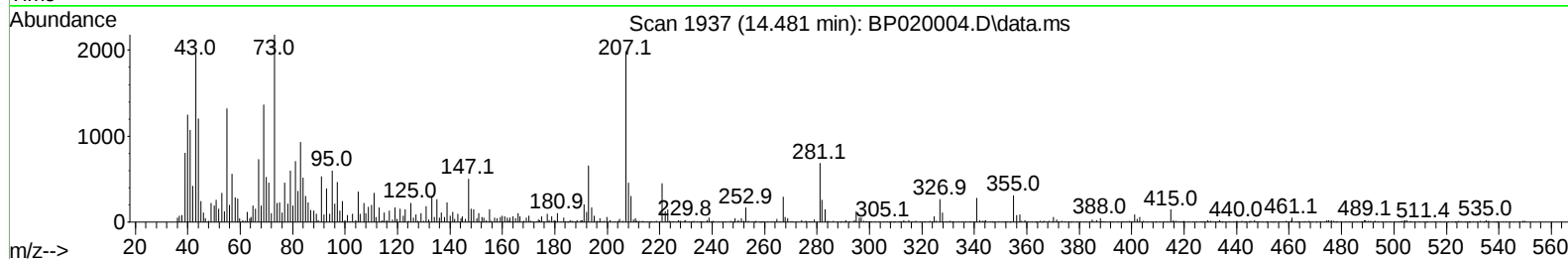
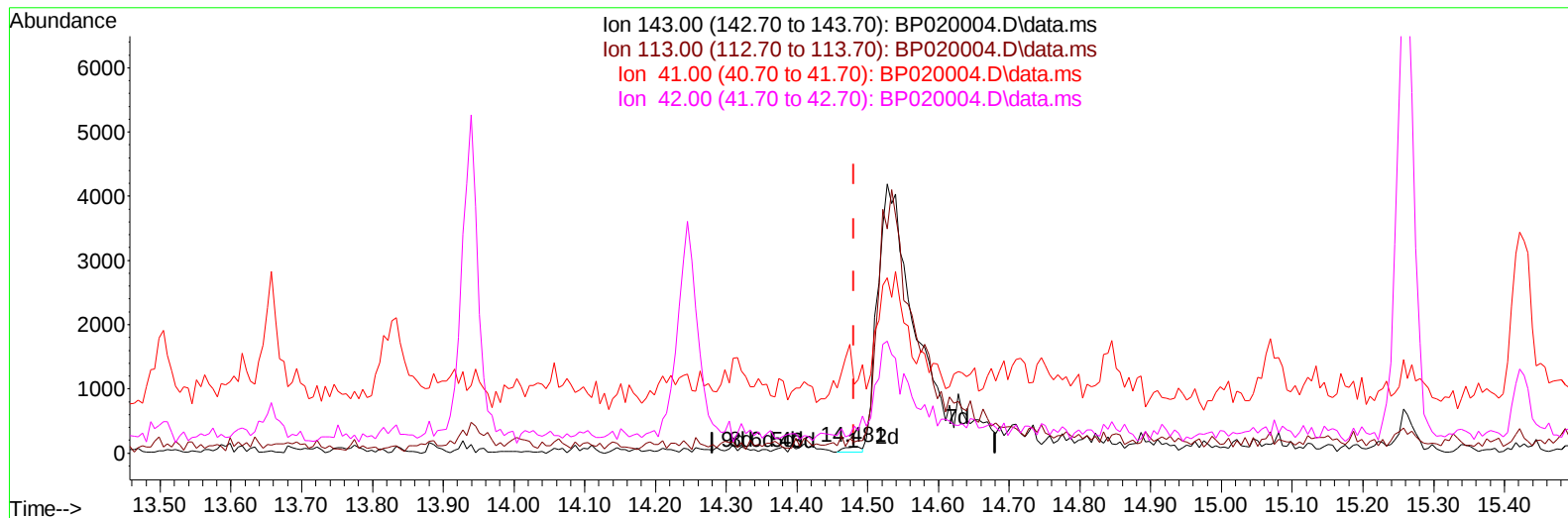
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032624\
Data File : BP020004.D
Acq On : 26 Mar 2024 19:52
Operator : MA/JU
Sample : P1848-04
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EOLY6

Manual IntegrationsAPPROVED

Quant Time: Mar 26 23:28:38 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032224.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Mar 25 10:46:59 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 03/29/2024
Supervised By :mohammad ahmed 03/29/2024



TIC: BP020004.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.481min (-0.000) 0.07 ng/u1

response 151

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	81.50	170.71#
41.00	40.90	1083.84#
42.00	28.50	425.25#

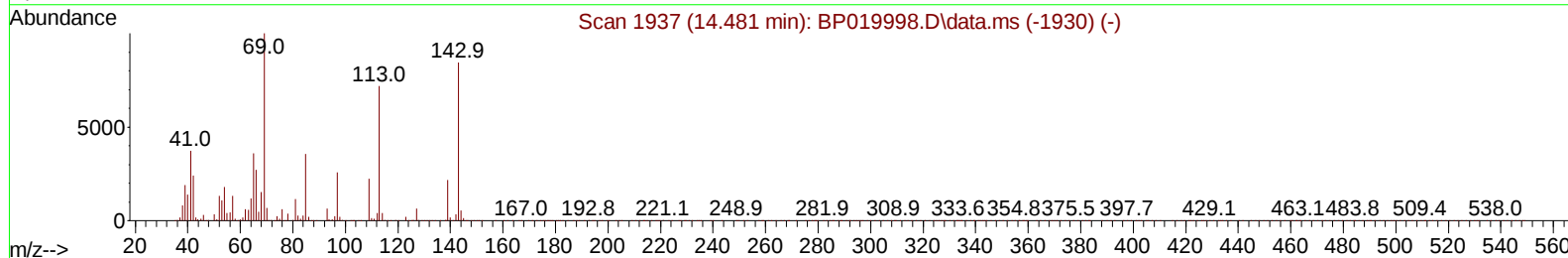
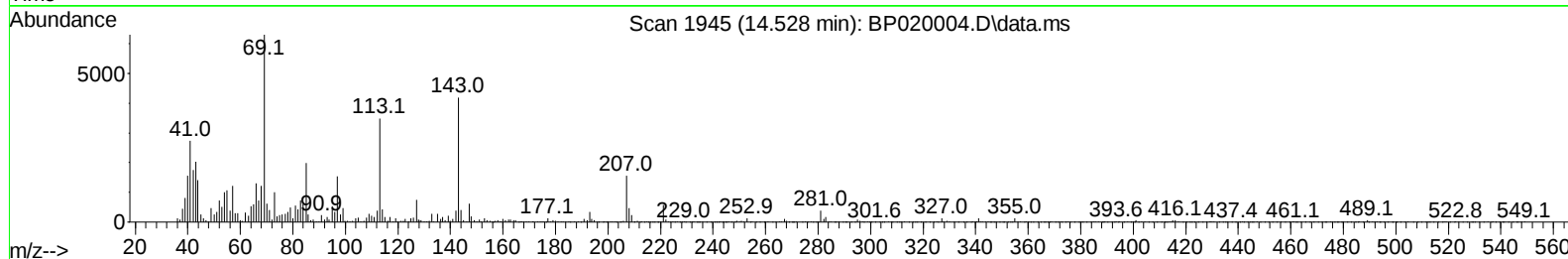
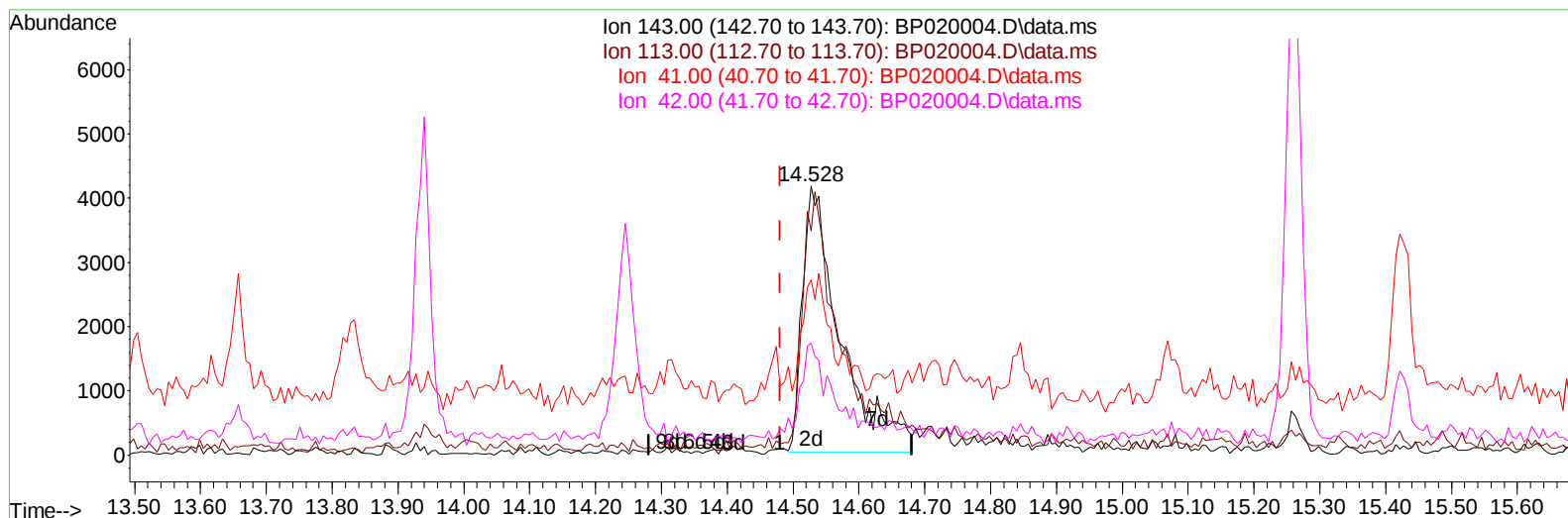
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Instrument :
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TIC: BP020004.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.528min (+ 0.047) 7.60 ng/ul m

response 16573

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	81.50	83.26
41.00	40.90	65.19#
42.00	28.50	41.71#

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 Sample : P1848-04
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 E0LY6

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 03/29/2024
 Supervised By :mohammad ahmed 03/29/2024

Quant Time: Mar 26 23:40:08 2024
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 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Mar 25 10:46:59 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.616	152	76276	20.000	ng/ul	0.00
20) Naphthalene-d8	10.369	136	302633	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.245	164	197953	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.069	188	438490	20.000	ng/ul	0.00
79) Chrysene-d12	21.533	240	446683	20.000	ng/ul	-0.01
88) Perylene-d12	24.815	264	488406	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.240	96	10593	6.085	ng/uL	0.00
4) Pyridine-d5	3.652	84	48236	9.176	ng/ul	0.01
7) Phenol-d5	6.804	99	60389	9.387	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.975	67	157021	37.464	ng/ul	0.00
11) 2-Chlorophenol-d4	7.157	132	135406	29.309	ng/ul	0.00
15) 4-Methylphenol-d8	8.310	113	102124	20.519	ng/ul	0.00
21) Nitrobenzene-d5	8.769	128	84202	36.300	ng/ul	0.01
24) 2-Nitrophenol-d4	9.475	143	88508	34.103	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.004	165	160509	33.864	ng/ul	0.00
31) 4-Chloroaniline-d4	10.516	131	292726	47.375	ng/ul	0.00
46) Dimethylphthalate-d6	13.657	166	574027	41.359	ng/ul	0.00
49) Acenaphthylene-d8	13.939	160	616615	38.033	ng/ul	0.00
54) 4-Nitrophenol-d4	14.528	143	16573m	7.596	ng/ul	0.05
60) Fluorene-d10	15.263	176	485906	41.099	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.428	200	51023	19.439	ng/ul	0.00
73) Anthracene-d10	17.163	188	780200	40.195	ng/ul	0.00
81) Pyrene-d10	19.563	212	1005868	38.381	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.580	264	984421	41.955	ng/ul	-0.03

Target Compounds Qvalue

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

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