

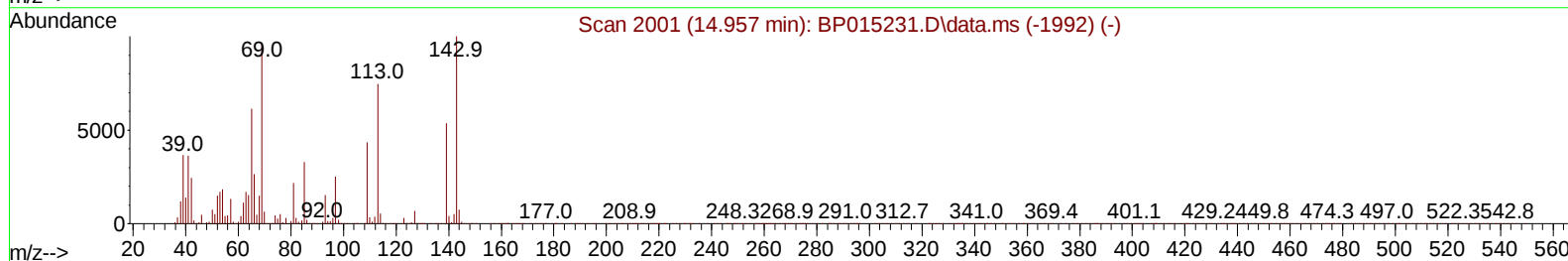
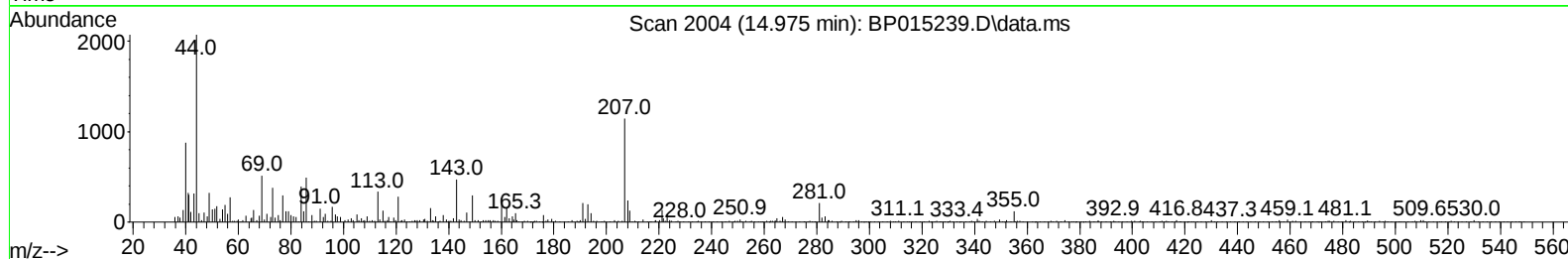
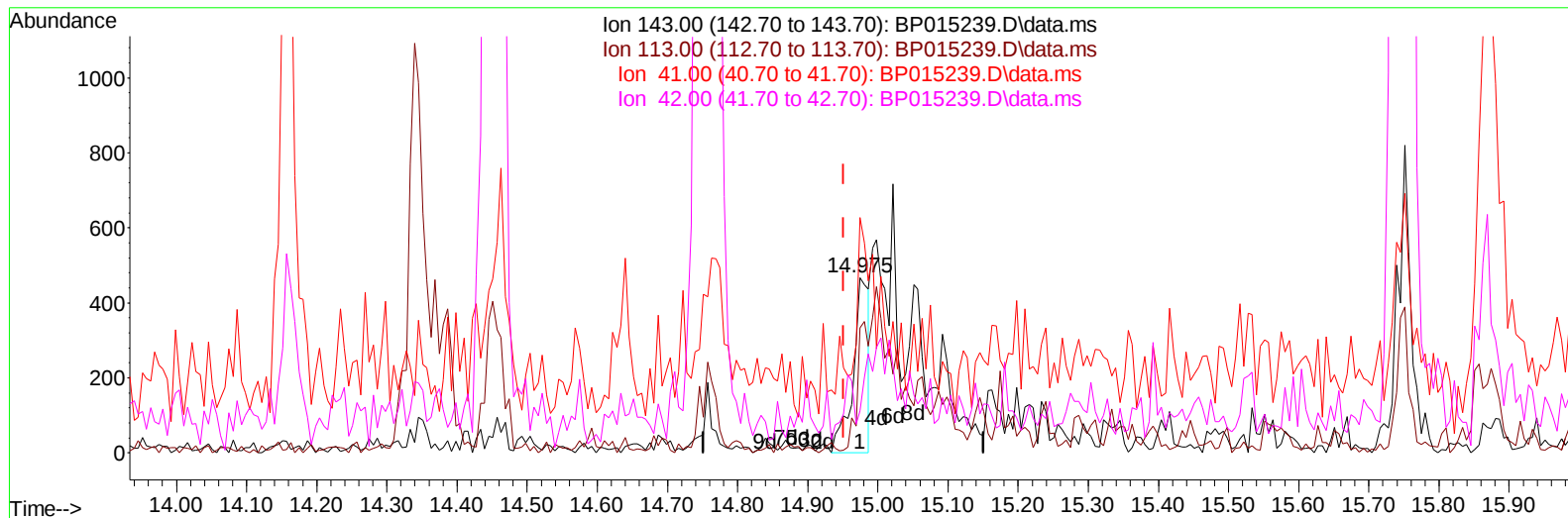
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052323\
Data File : BP015239.D
Acq On : 23 May 2023 14:18
Operator : CG/JU
Sample : 02861-06
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
H0FL1

Manual IntegrationsAPPROVED

Quant Time: May 23 14:54:13 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051123.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue May 23 01:25:41 2023
Response via : Initial Calibration

Reviewed By :Christian Giraldo 05/24/2023
Supervised By :mohammad ahmed 05/24/2023



TIC: BP015239.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.975min (+ 0.023) 0.09 ng/u1

response 736

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	67.30	71.37
41.00	36.10	133.97#
42.00	24.50	24.15

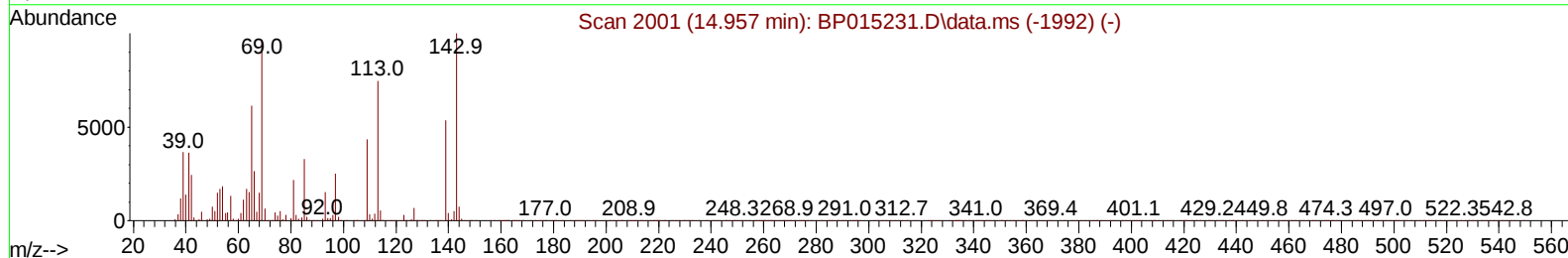
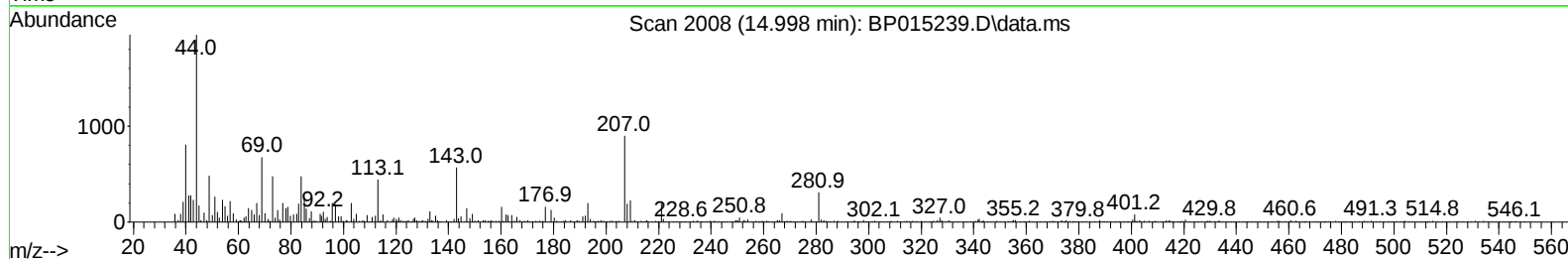
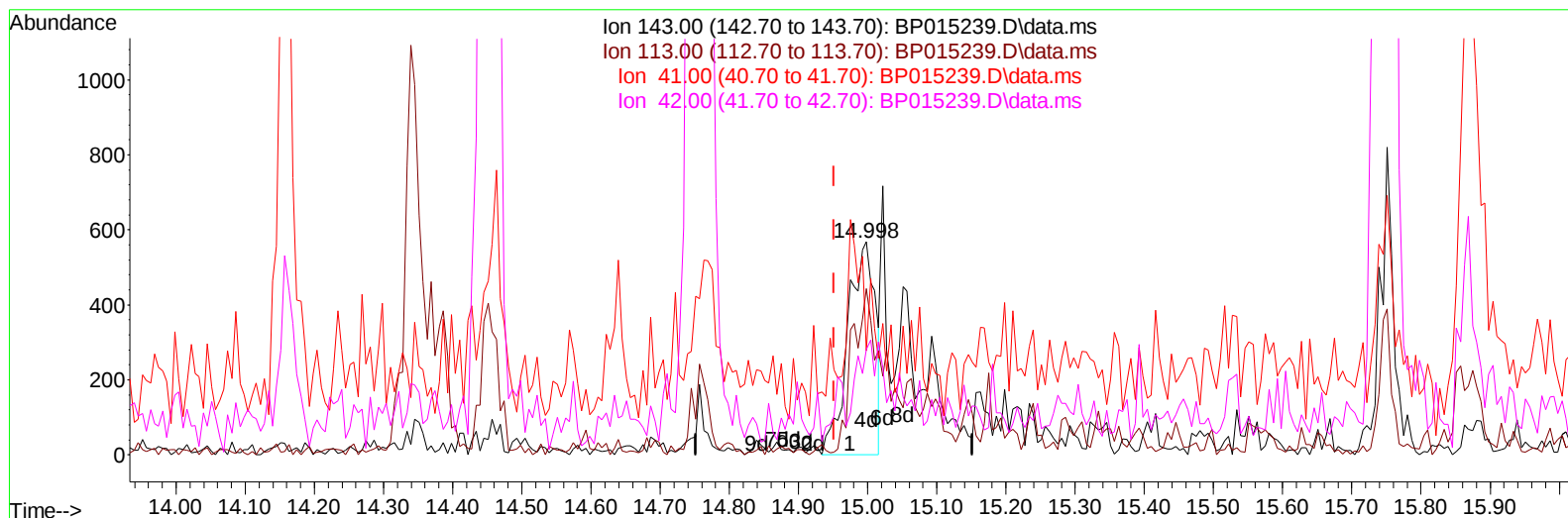
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 ALS Vial : 10 Sample Multiplier: 1

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Manual IntegrationsAPPROVED

Quant Time: May 23 23:03:08 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051123.MA.M
 Quant Title : SVOA CALIBRATION
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 Supervised By :mohammad ahmed 05/24/2023



TIC: BP015239.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.998min (+ 0.047) 0.19 ng/ul m

response 1566

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	67.30	78.17
41.00	36.10	48.94#
42.00	24.50	49.47#

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052323\
 Data File : BP015239.D
 Acq On : 23 May 2023 14:18
 Operator : CG/JU
 Sample : 02861-06
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 H0FL1

Manual IntegrationsAPPROVED

Reviewed By :Christian Giraldo 05/24/2023
 Supervised By :mohammad ahmed 05/24/2023

Quant Time: May 24 04:23:54 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051123.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue May 23 01:25:41 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.122	152	195913	20.000	ng/ul	0.00
20) Naphthalene-d8	10.951	136	856032	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.757	164	550883	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.510	188	1117222	20.000	ng/ul	0.00
79) Chrysene-d12	21.586	240	802129	20.000	ng/ul	0.00
88) Perylene-d12	24.139	264	884479	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.428	96	18940	3.756	ng/uL	0.00
4) Pyridine-d5	3.887	84	45315	3.170	ng/ul	0.02
7) Phenol-d5	7.269	99	33124	1.821	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.440	67	132962	12.656	ng/ul	0.00
11) 2-Chlorophenol-d4	7.646	132	122562	9.164	ng/ul	0.00
15) 4-Methylphenol-d8	8.834	113	74227	5.121	ng/ul	0.00
21) Nitrobenzene-d5	9.299	128	87112	12.950	ng/ul	0.00
24) 2-Nitrophenol-d4	10.022	143	85504	11.424	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.563	165	148676	11.008	ng/ul	0.00
31) 4-Chloroaniline-d4	11.087	131	150189	7.516	ng/ul	0.00
46) Dimethylphthalate-d6	14.163	166	610990	14.677	ng/ul	0.00
49) Acenaphthylene-d8	14.451	160	656489	13.789	ng/ul	0.00
54) 4-Nitrophenol-d4	14.998	143	1566m	0.194	ng/ul	0.05
60) Fluorene-d10	15.751	176	507117	14.437	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.875	200	19292	2.740	ng/ul	0.00
73) Anthracene-d10	17.610	188	764410	14.942	ng/ul	0.00
81) Pyrene-d10	19.827	212	827832	17.794	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.974	264	706059	16.081	ng/ul	0.00

Target Compounds Qvalue

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

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Manual IntegrationsAPPROVED

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