Quantitation Report (Qedit)

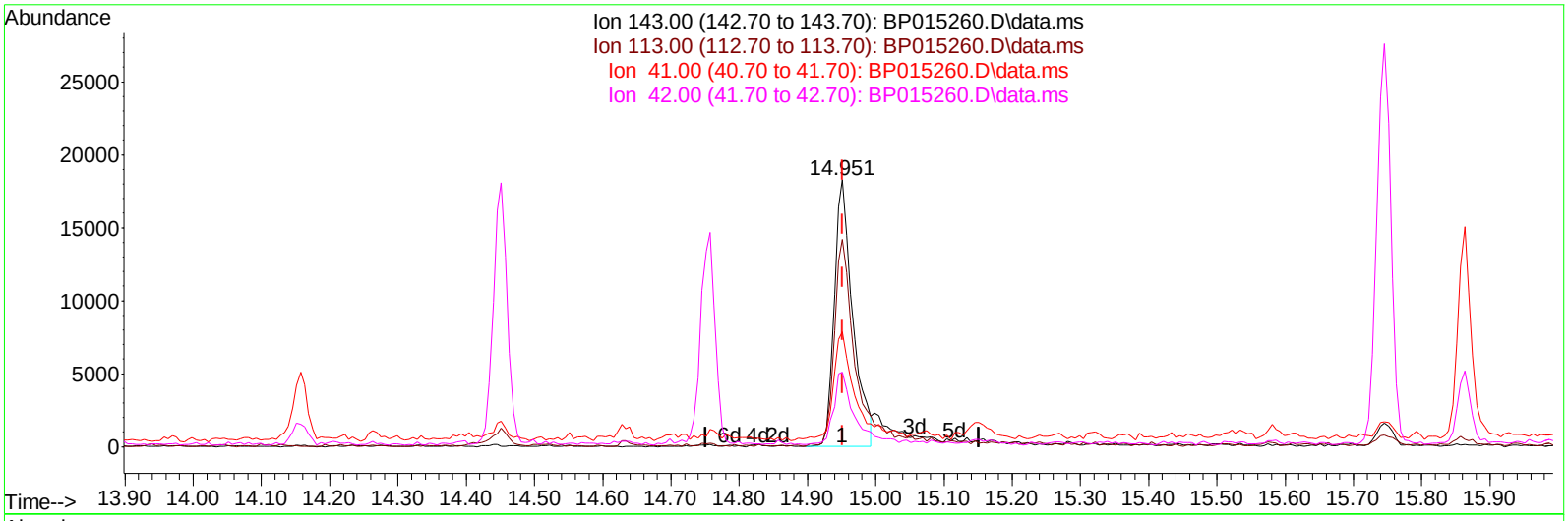
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052323\
Data File : BP015260.D
Acq On : 24 May 2023 02:35
Operator : CG/JU
Sample : 02861-15
Misc :
ALS Vial : 31 Sample Multiplier: 1

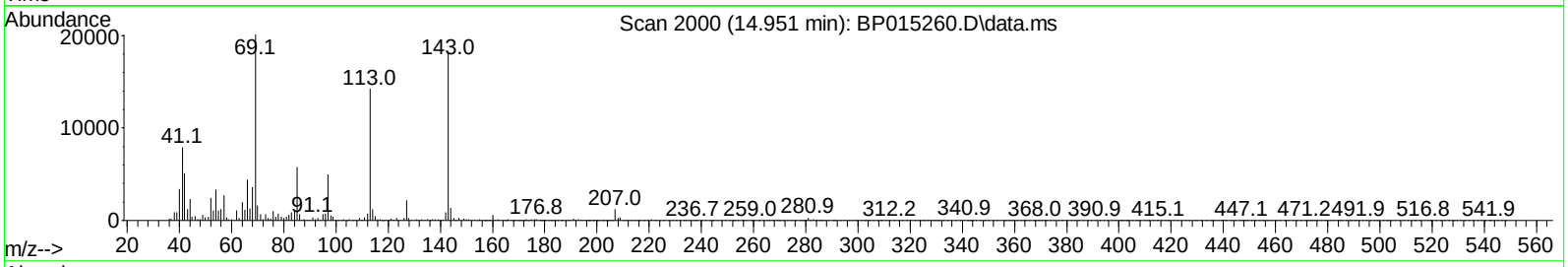
Quant Time: May 24 04:11:42 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051123.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue May 23 01:25:41 2023
Response via : Initial Calibration

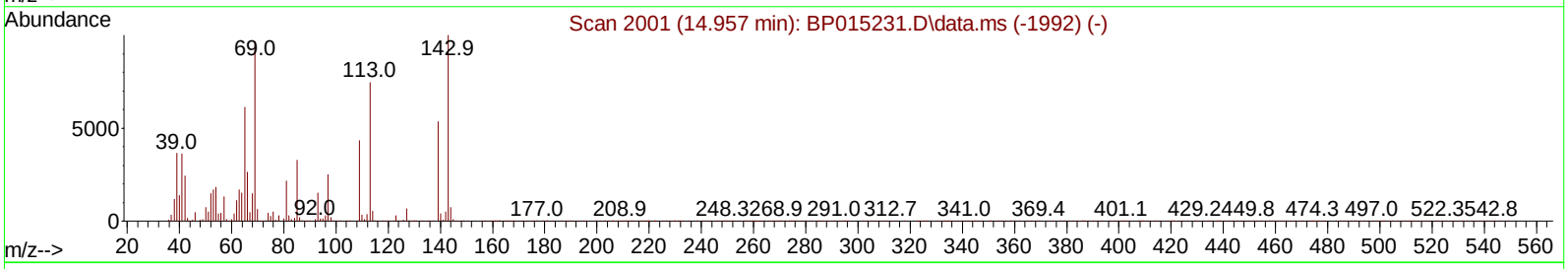
Instrument :
BNA_P
ClientSampleId :
H0FL8

Manual IntegrationsAPPROVED

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







TIC: BP015260.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.951min (0.000) 3.36 ng/ul

response 34333

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	67.30	77.72
41.00	36.10	43.10
42.00	24.50	27.90

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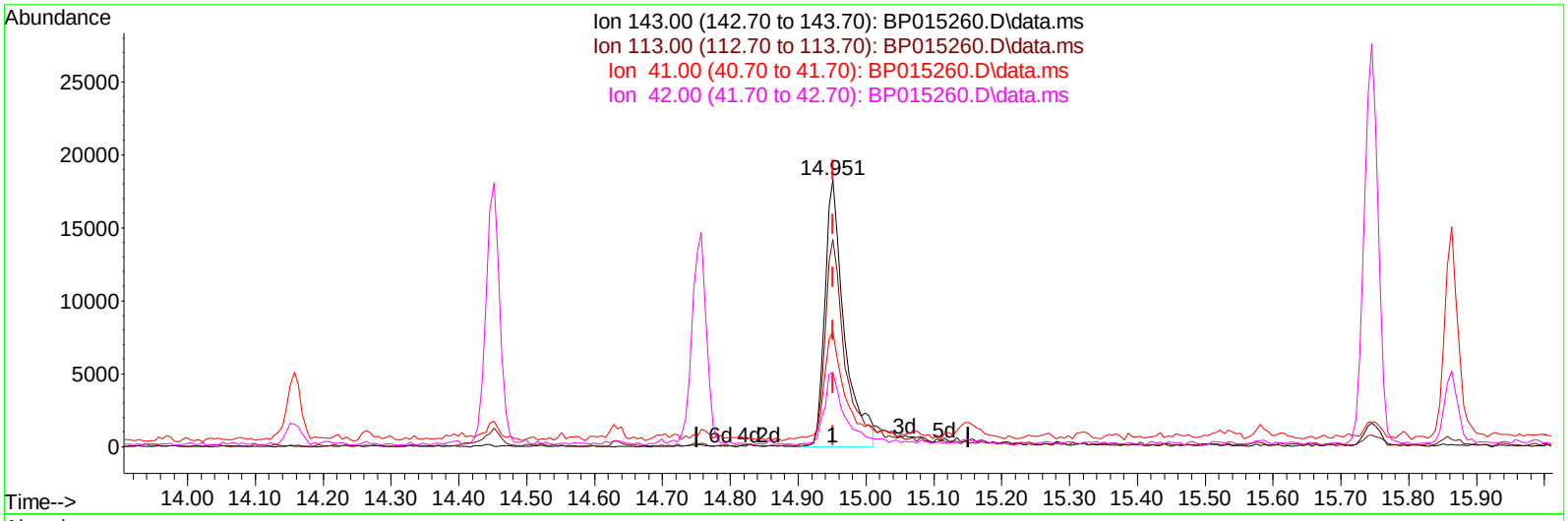

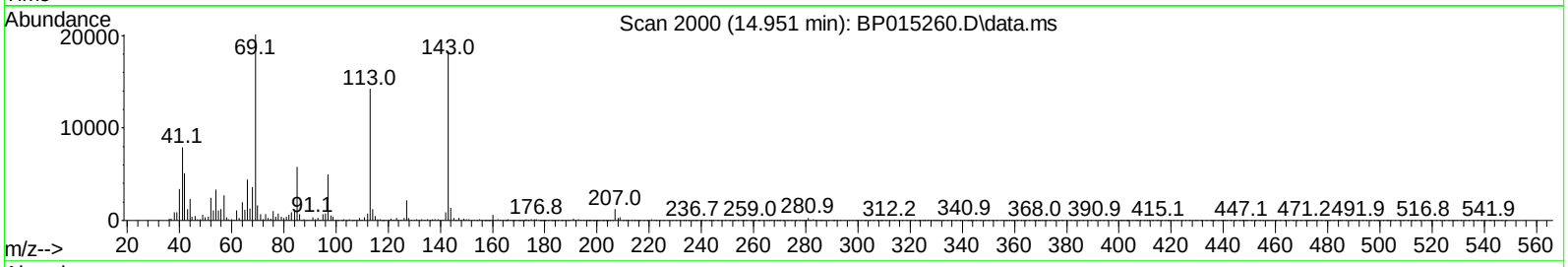

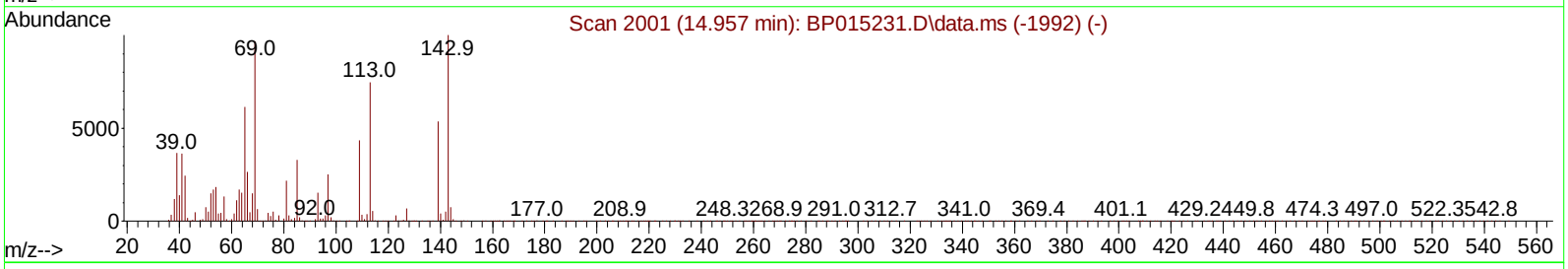

TIC: BP015260.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.951min (0.000) 3.57 ng/ul m

response 36536

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	67.30	77.72
41.00	36.10	43.10
42.00	24.50	27.90

Quantitation Report (QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052323\
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Acq On : 24 May 2023 02:35
Operator : CG/JU
Sample : 02861-15
Misc :
ALS Vial : 31 Sample Multiplier: 1

Quant Time: May 24 04:15:57 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
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.116	152	251309	20.000	ng/ul	0.00
20) Naphthalene-d8	10.946	136	1099405	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.757	164	696950	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.510	188	1460431	20.000	ng/ul	0.00
79) Chrysene-d12	21.580	240	1035999	20.000	ng/ul	0.00
88) Perylene-d12	24.139	264	1117367	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.428	96	27934	4.319	ng/uL	0.00
4) Pyridine-d5	3.870	84	163097	8.896	ng/ul	0.00
7) Phenol-d5	7.269	99	128816	5.521	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.434	67	406686	30.178	ng/ul	0.00
11) 2-Chlorophenol-d4	7.640	132	393651	22.945	ng/ul	0.00
15) 4-Methylphenol-d8	8.828	113	262457	14.116	ng/ul	0.00
21) Nitrobenzene-d5	9.293	128	274946	31.826	ng/ul	0.00
24) 2-Nitrophenol-d4	10.022	143	295858	30.778	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.557	165	508507	29.316	ng/ul	0.00
31) 4-Chloroaniline-d4	11.081	131	463817	18.073	ng/ul	0.00
46) Dimethylphthalate-d6	14.157	166	1921079	36.477	ng/ul	0.00
49) Acenaphthylene-d8	14.451	160	2059639	34.194	ng/ul	0.00
54) 4-Nitrophenol-d4	14.951	143	36536m	3.573	ng/ul	0.00
60) Fluorene-d10	15.745	176	1568776	35.301	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.863	200	243497	26.456	ng/ul	0.00
73) Anthracene-d10	17.604	188	2331795	34.868	ng/ul	0.00
81) Pyrene-d10	19.828	212	2533459	42.164	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.969	264	2102908	37.913	ng/ul	0.00
Target Compounds						Qvalue

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

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