Quantitation Report (Qedit)

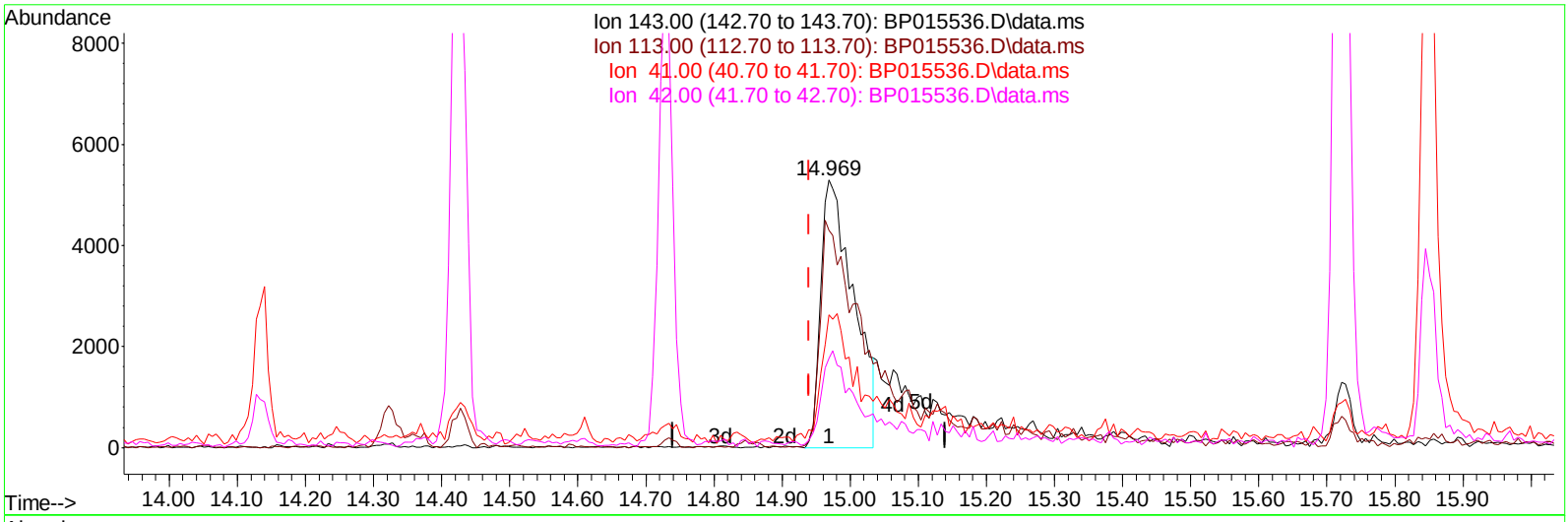
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061423\
Data File : BP015536.D
Acq On : 15 Jun 2023 01:37
Operator : MA/JU
Sample : 03104-10
Misc :
ALS Vial : 30 Sample Multiplier: 1

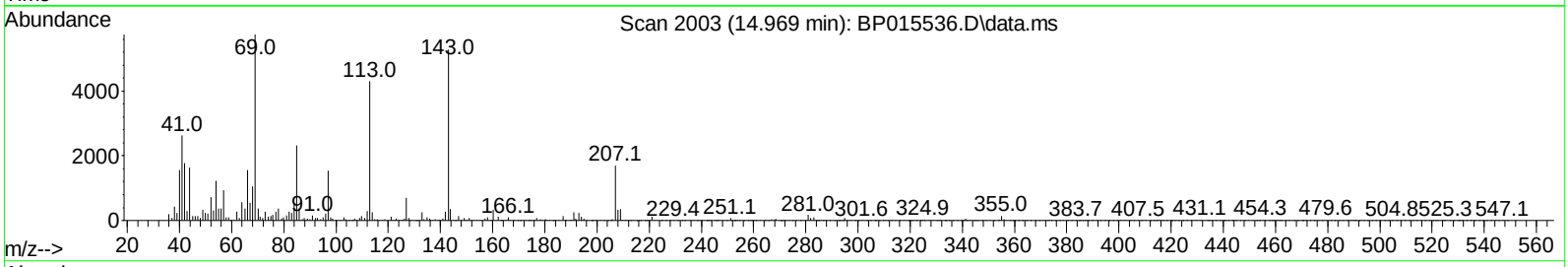
Quant Time: Jun 15 02:26:19 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Jun 09 23:33:38 2023
Response via : Initial Calibration

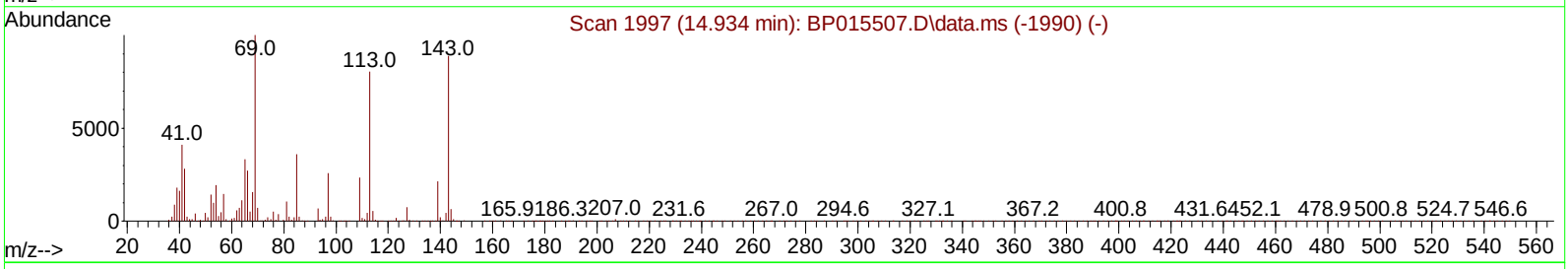
Instrument :
BNA_P
ClientSampleId :
C0NL6

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 06/15/2023
Supervised By :mohammad ahmed 06/16/2023







TIC: BP015536.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.969min (+ 0.029) 3.56 ng/ul

response 17745

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	86.60	81.33
41.00	42.20	49.67
42.00	29.40	33.45

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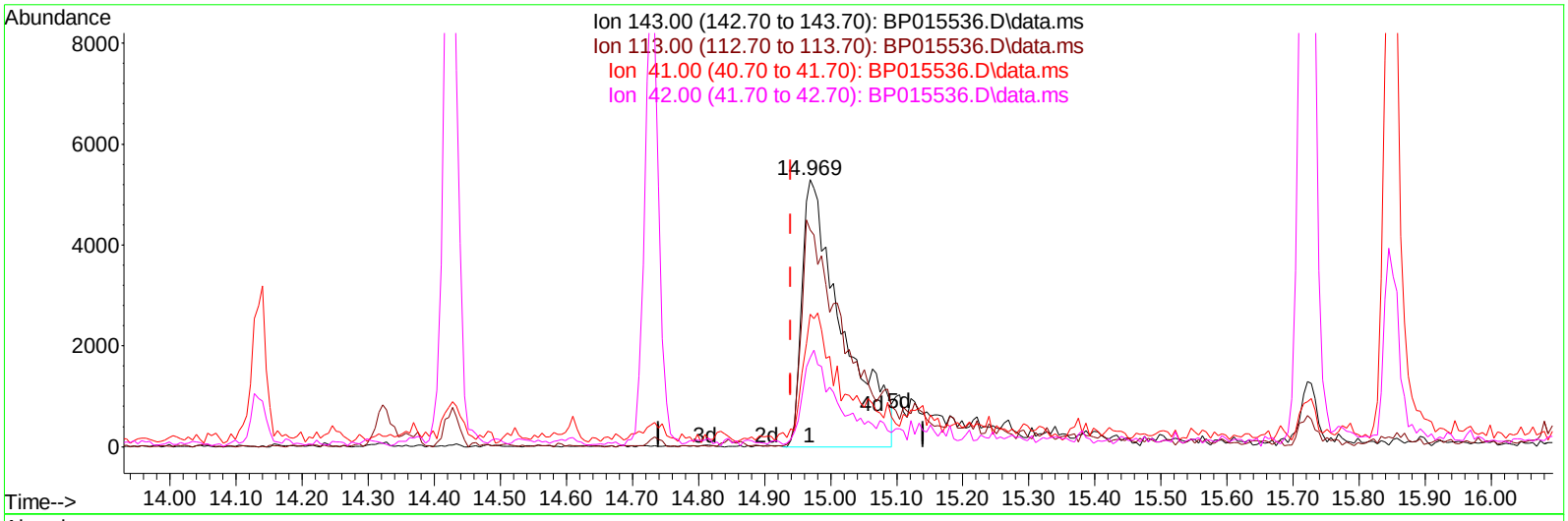

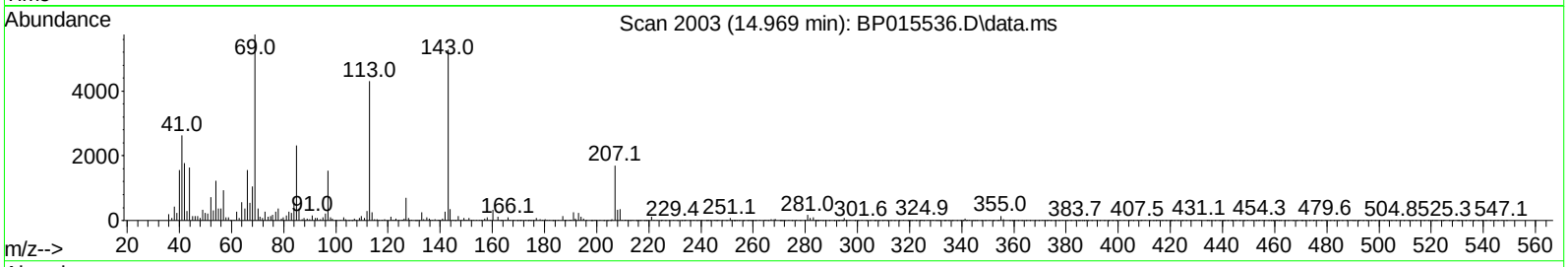

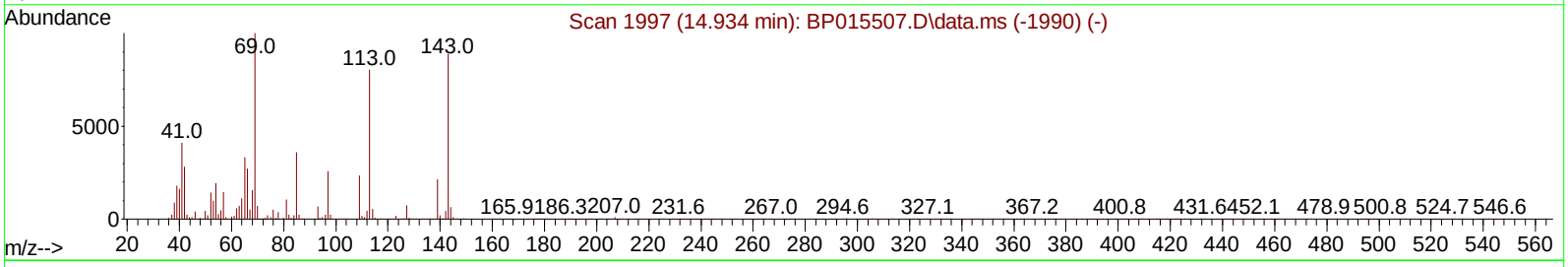

TIC: BP015536.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.969min (+ 0.029) 4.47 ng/ul m

response 22259

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	86.60	81.33
41.00	42.20	49.67
42.00	29.40	33.45

Quantitation Report (QT Reviewed)

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.087	152	105827	20.000	ng/ul	0.00
20) Naphthalene-d8	10.910	136	503094	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.734	164	352810	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.486	188	827951	20.000	ng/ul	0.00
79) Chrysene-d12	21.568	240	839094	20.000	ng/ul	0.00
88) Perylene-d12	24.115	264	906091	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.405	96	15586	5.454	ng/uL	0.00
4) Pyridine-d5	3.864	84	41007	5.524	ng/ul	0.02
7) Phenol-d5	7.246	99	70326	7.213	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.410	67	221476	38.037	ng/ul	0.00
11) 2-Chlorophenol-d4	7.610	132	207489	29.354	ng/ul	-0.01
15) 4-Methylphenol-d8	8.799	113	146824	18.349	ng/ul	-0.01
21) Nitrobenzene-d5	9.263	128	145647	36.875	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.993	143	162504	36.027	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.534	165	266169	32.349	ng/ul	0.00
31) 4-Chloroaniline-d4	11.057	131	244608	20.711	ng/ul	0.00
46) Dimethylphthalate-d6	14.134	166	1131876	40.684	ng/ul	0.00
49) Acenaphthylene-d8	14.428	160	1208905	39.739	ng/ul	0.00
54) 4-Nitrophenol-d4	14.969	143	22259m	4.468	ng/ul	0.03
60) Fluorene-d10	15.722	176	967625	41.244	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.851	200	198929	37.926	ng/ul	0.00
73) Anthracene-d10	17.586	188	1743491	46.283	ng/ul	0.00
81) Pyrene-d10	19.810	212	2194336	48.861	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.951	264	2166398	47.651	ng/ul	0.00

Target Compounds Qvalue

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

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