Quantitation Report (Qedit)

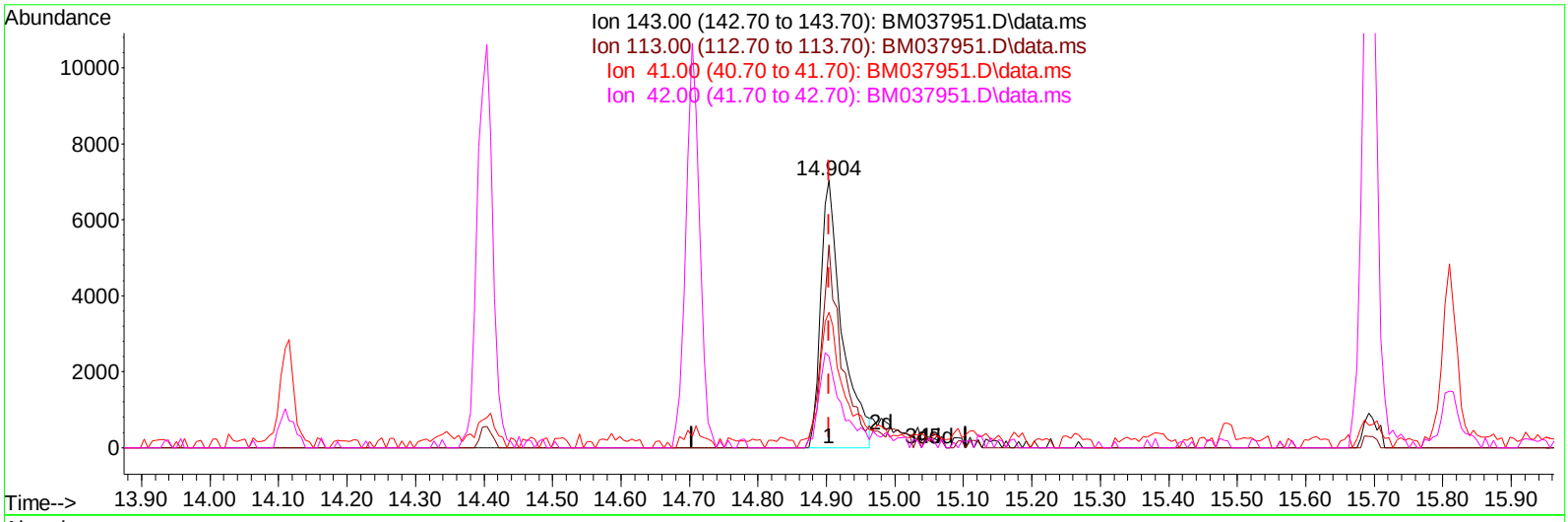
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121022\
Data File : BM037951.D
Acq On : 11 Dec 2022 01:46
Operator : CG/JU
Sample : N5927-01
Misc :
ALS Vial : 27 Sample Multiplier: 1

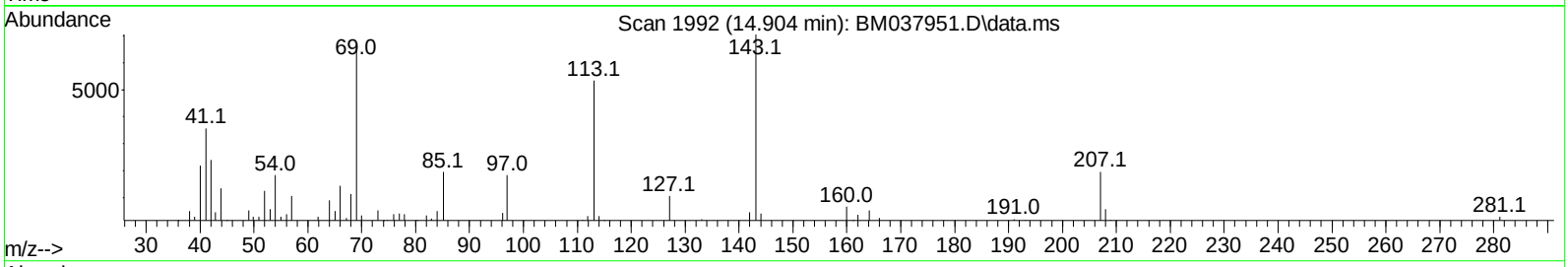
Quant Time: Dec 11 21:31:02 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM120722.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Dec 09 21:34:57 2022
Response via : Initial Calibration

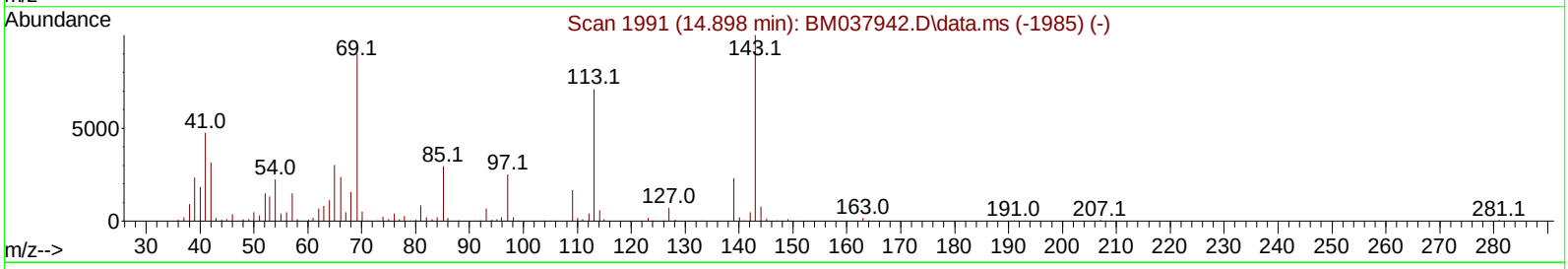
Instrument :
BNA_M
ClientSampleId :
EXZ85

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 12/12/2022
Supervised By :mohammad ahmed 12/12/2022







TIC: BM037951.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.904min (-0.000) 2.90 ng/ul

response 15198

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	60.30	75.75#
41.00	43.90	50.67
42.00	30.30	34.08

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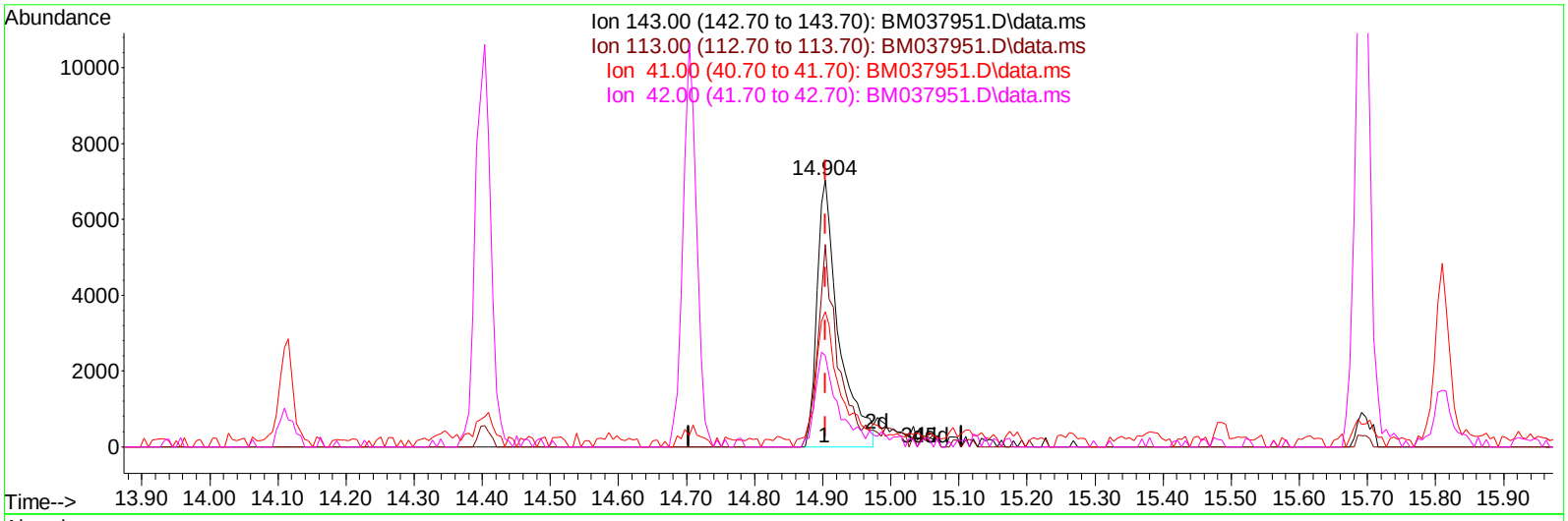

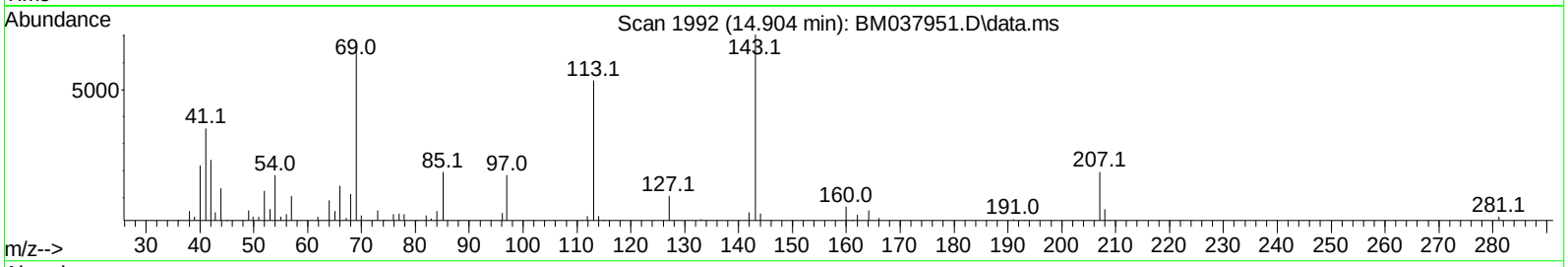

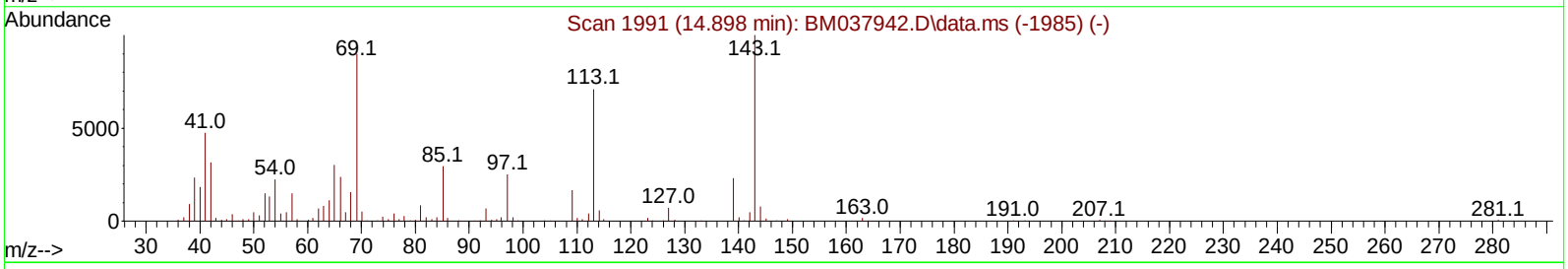

TIC: BM037951.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.904min (-0.000) 2.99 ng/ul m

response 15667

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	60.30	75.75#
41.00	43.90	50.67
42.00	30.30	34.08

Quantitation Report (QT Reviewed)

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.075	152	207136	20.000	ng/ul	0.00
20) Naphthalene-d8	10.898	136	804187	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.704	164	430742	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.445	188	788796	20.000	ng/ul	0.00
79) Chrysene-d12	21.615	240	712221	20.000	ng/ul	0.00
88) Perylene-d12	24.097	264	785720	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.422	96	21662	4.748	ng/uL	0.00
4) Pyridine-d5	3.863	84	105850	7.822	ng/ul	0.00
7) Phenol-d5	7.228	99	101102	6.007	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	7.398	67	303289	29.591	ng/ul	0.00
11) 2-Chlorophenol-d4	7.604	132	315137	22.823	ng/ul	0.00
15) 4-Methylphenol-d8	8.786	113	177257	13.491	ng/ul	0.00
21) Nitrobenzene-d5	9.251	128	207301	32.558	ng/ul	0.00
24) 2-Nitrophenol-d4	9.975	143	187221	28.058	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.510	165	320583	25.295	ng/ul	-0.01
31) 4-Chloroaniline-d4	11.033	131	418546	23.924	ng/ul	0.00
46) Dimethylphthalate-d6	14.115	166	969361	31.566	ng/ul	0.00
49) Acenaphthylene-d8	14.404	160	1182507	31.507	ng/ul	0.00
54) 4-Nitrophenol-d4	14.904	143	15667m	2.994	ng/ul	0.00
60) Fluorene-d10	15.692	176	850291	31.078	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.809	200	63228	14.993	ng/ul	0.00
73) Anthracene-d10	17.545	188	1298334	35.267	ng/ul	0.00
81) Pyrene-d10	19.827	212	1462149	34.176	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.932	264	1431324	36.471	ng/ul	-0.01

Target Compounds Qvalue

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

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