Quantitation Report (Qedit)

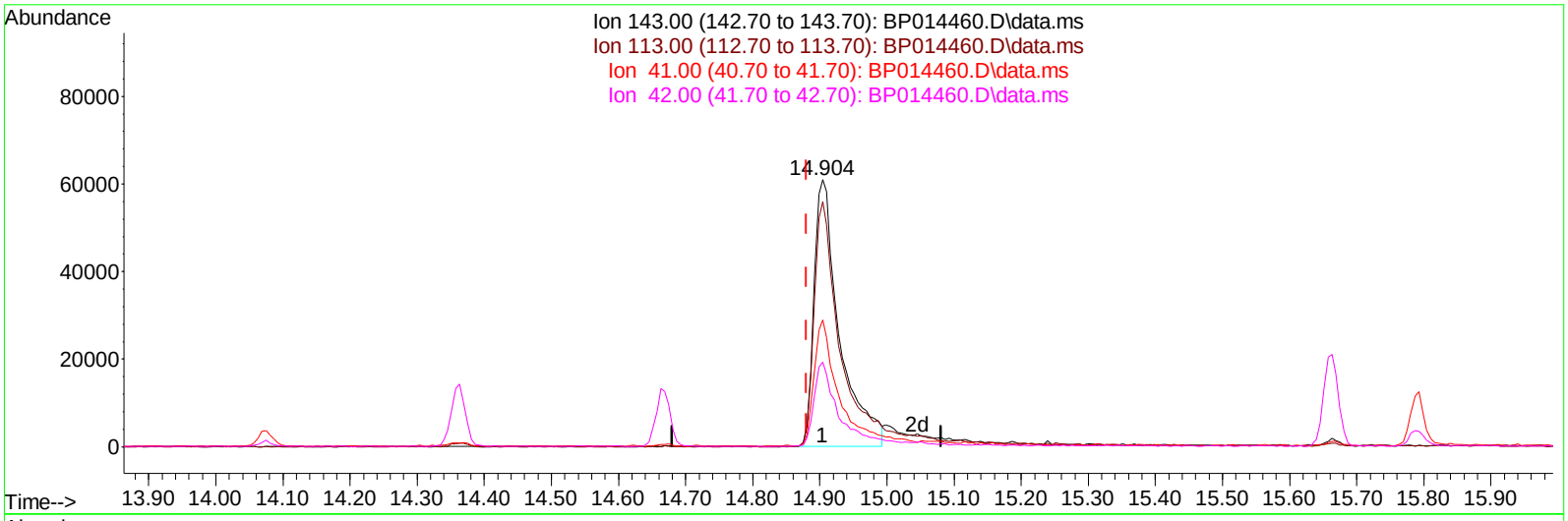
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP033023\
Data File : BP014460.D
Acq On : 01 Apr 2023 00:59
Operator : CG/JU
Sample : PB151752BL
Misc :
ALS Vial : 70 Sample Multiplier: 1

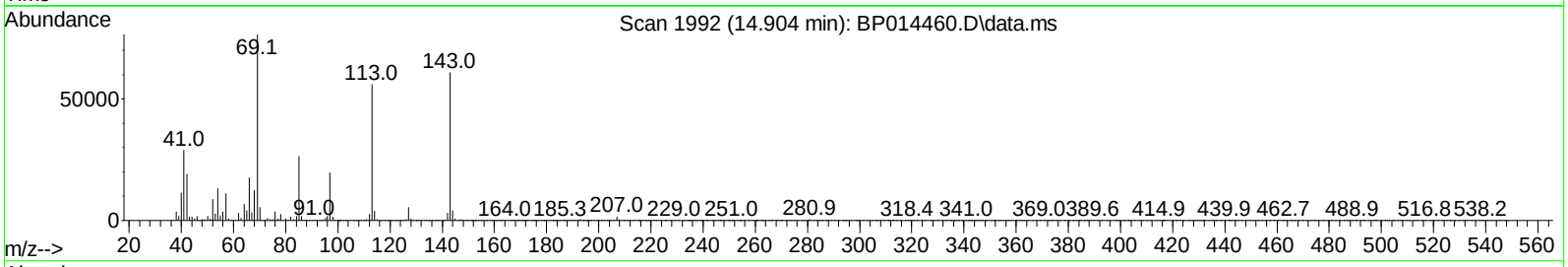
Quant Time: Apr 01 02:04:19 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Mar 31 04:41:29 2023
Response via : Initial Calibration

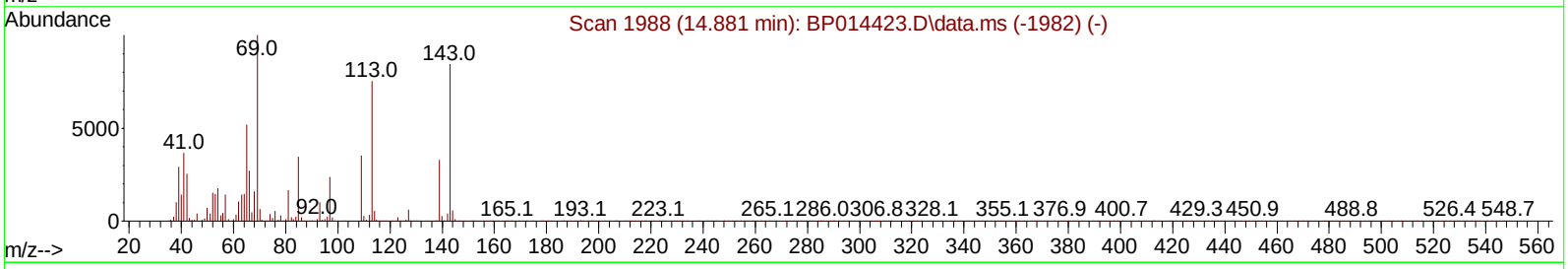
Instrument :
BNA_P
ClientSampleId :
SBLK752

Manual IntegrationsAPPROVED

Reviewed By :Christian Giraldo 04/03/2023
Supervised By :Jagrut Upadhyay 04/03/2023







TIC: BP014460.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.904min (+ 0.023) 23.34 ng/ul

response 160796

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	91.82
41.00	44.20	47.55
42.00	31.50	31.64

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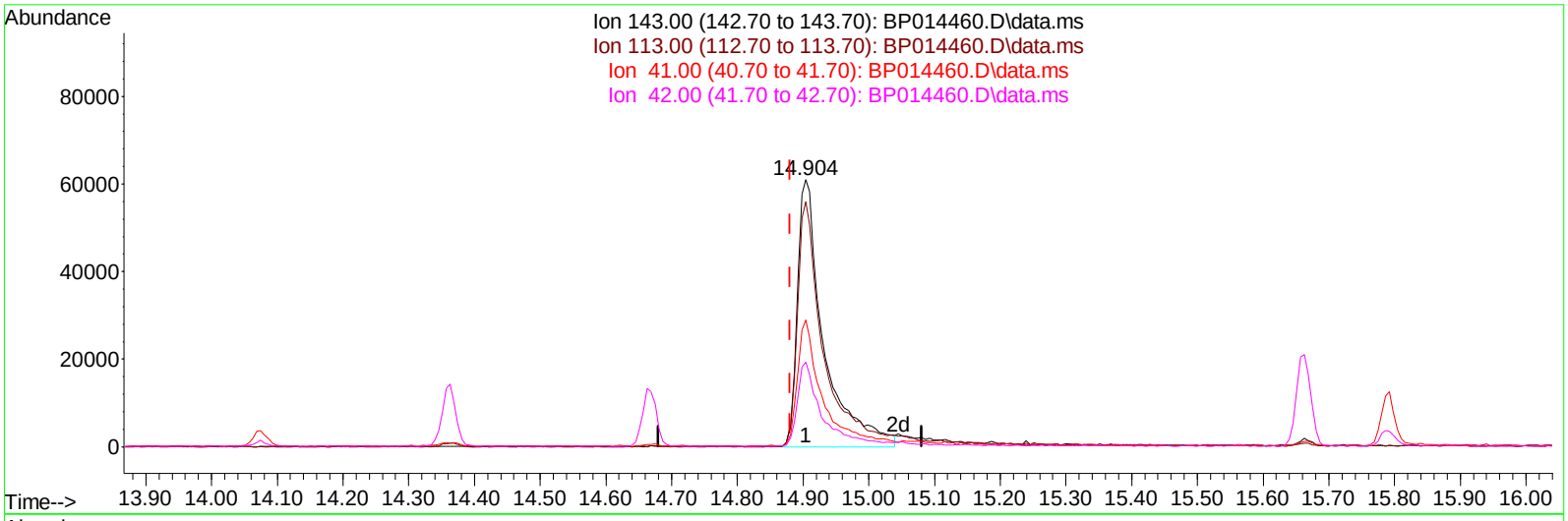

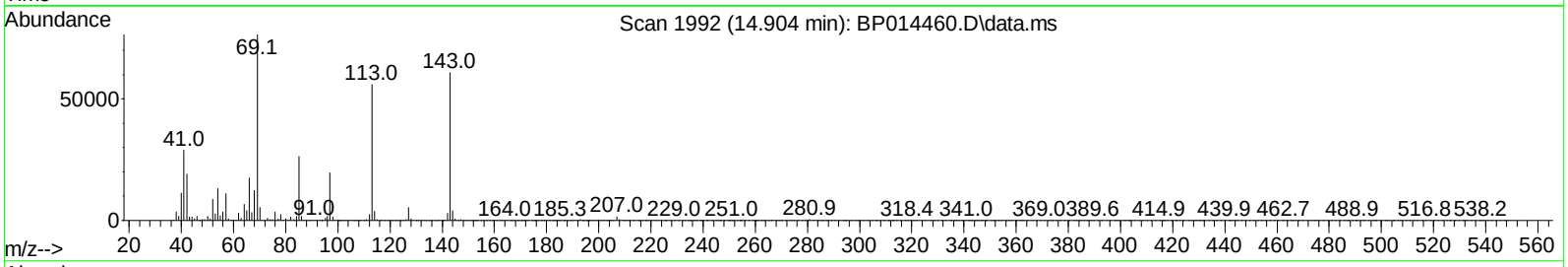

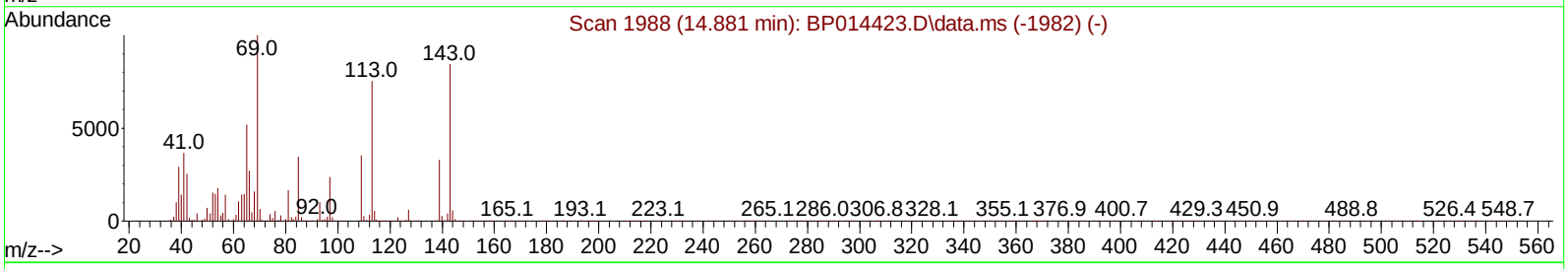

TIC: BP014460.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.904min (+ 0.023) 24.87 ng/ul m

response 171311

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	91.82
41.00	44.20	47.55
42.00	31.50	31.64

Quantitation Report (QT Reviewed)

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.010	152	176257	20.000	ng/ul	0.00
20) Naphthalene-d8	10.834	136	805489	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.669	164	562436	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.428	188	1249826	20.000	ng/ul	0.00
79) Chrysene-d12	21.516	240	941149	20.000	ng/ul	0.00
88) Perylene-d12	24.039	264	936324	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.387	96	25553	5.622	ng/uL	0.00
4) Pyridine-d5	3.828	84	110191	9.460	ng/ul	0.01
7) Phenol-d5	7.187	99	497748	30.621	ng/ul	0.01
9) Bis-(2-Chloroethyl)eth...	7.340	67	315540	30.255	ng/ul	0.00
11) 2-Chlorophenol-d4	7.546	132	368820	31.392	ng/ul	0.00
15) 4-Methylphenol-d8	8.740	113	413562	31.541	ng/ul	0.01
21) Nitrobenzene-d5	9.193	128	190072	29.230	ng/ul	0.00
24) 2-Nitrophenol-d4	9.916	143	227792	30.593	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.469	165	385338	28.416	ng/ul	0.01
31) 4-Chloroaniline-d4	10.999	131	53074	2.971	ng/ul	0.02
46) Dimethylphthalate-d6	14.075	166	1414380	31.015	ng/ul	0.00
49) Acenaphthylene-d8	14.363	160	1494451	29.534	ng/ul	0.00
54) 4-Nitrophenol-d4	14.904	143	171311m	24.866	ng/ul	0.02
60) Fluorene-d10	15.663	176	1161574	30.544	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.792	200	206073	22.408	ng/ul	0.00
73) Anthracene-d10	17.528	188	1933613	31.968	ng/ul	0.00
81) Pyrene-d10	19.757	212	2131791	33.722	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.874	264	1618248	32.530	ng/ul	0.00
Target Compounds						Qvalue

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

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