

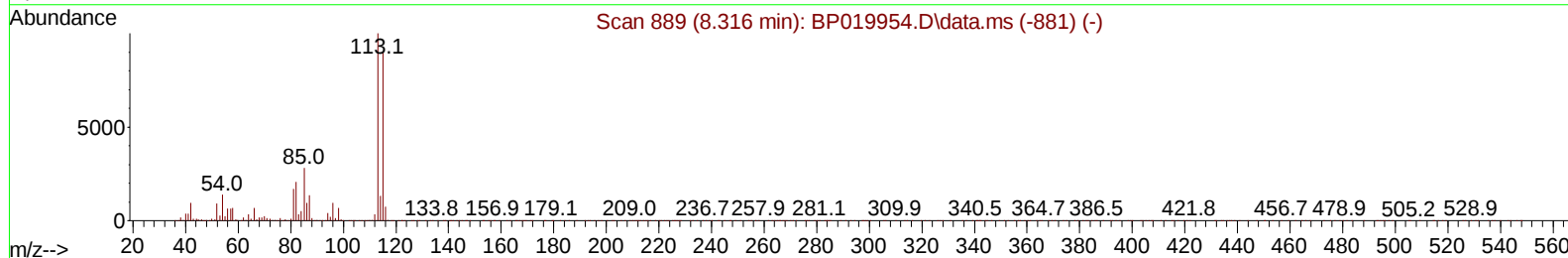
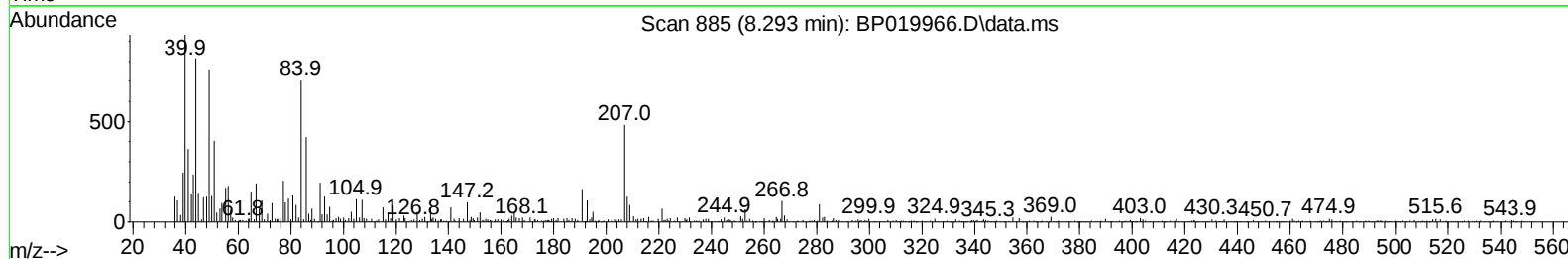
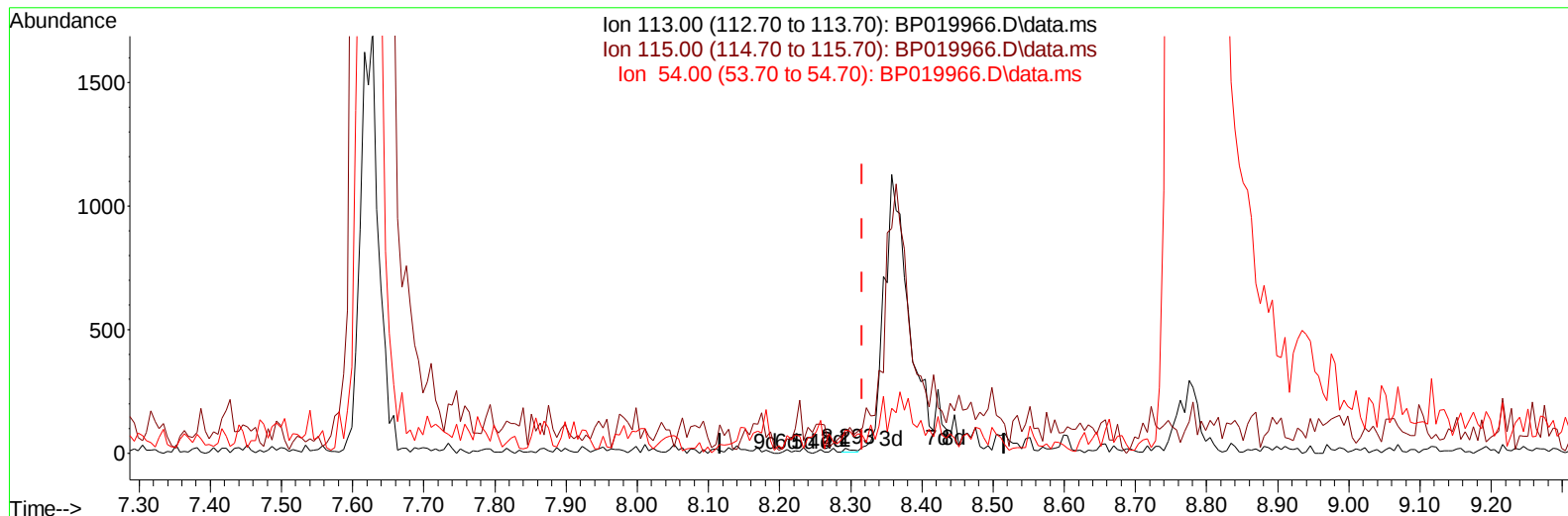
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032224\
Data File : BP019966.D
Acq On : 22 Mar 2024 22:06
Operator : MA/JU
Sample : P1813-13
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E0LY1

Manual IntegrationsAPPROVED

Quant Time: Mar 22 23:03:54 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032224.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Mar 22 22:40:53 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 03/26/2024
Supervised By :mohammad ahmed 03/26/2024



TIC: BP019966.D\data.ms

(15) 4-Methylphenol-d8 (S)

8.293min (-0.023) 0.00 ng/u1

response 13

Ion	Exp%	Act%
113.00	100.00	100.00
115.00	90.00	389.47#
54.00	13.80	500.00#
0.00	0.00	0.00

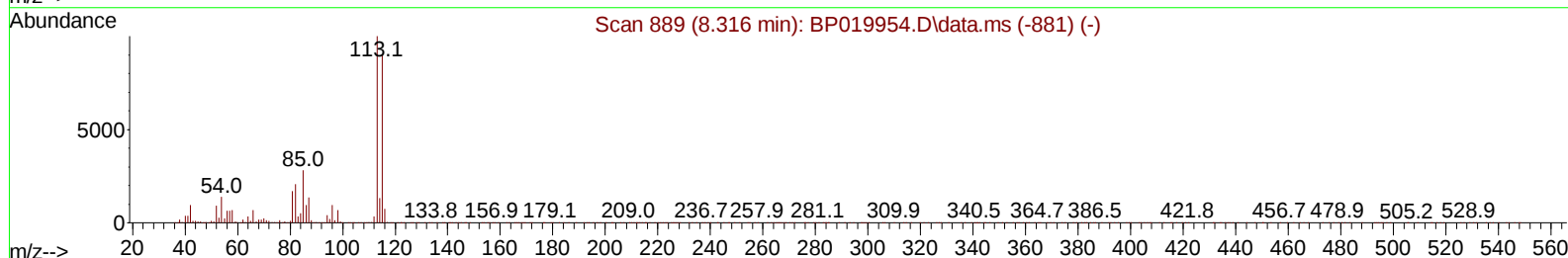
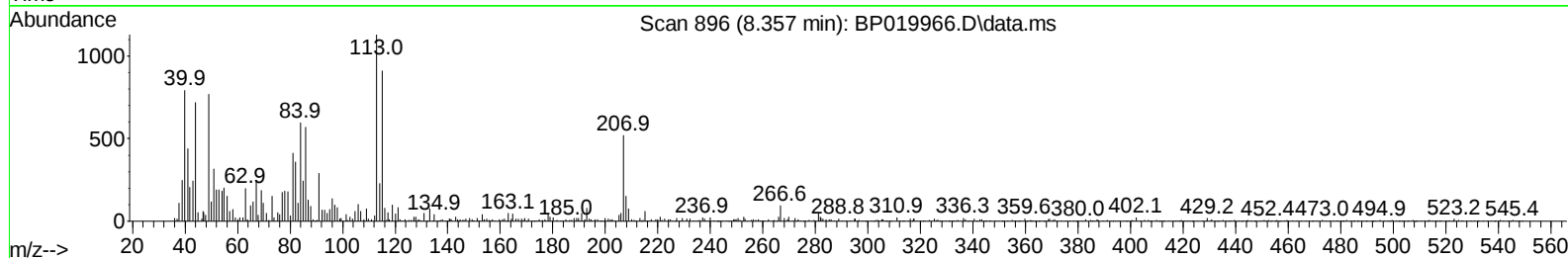
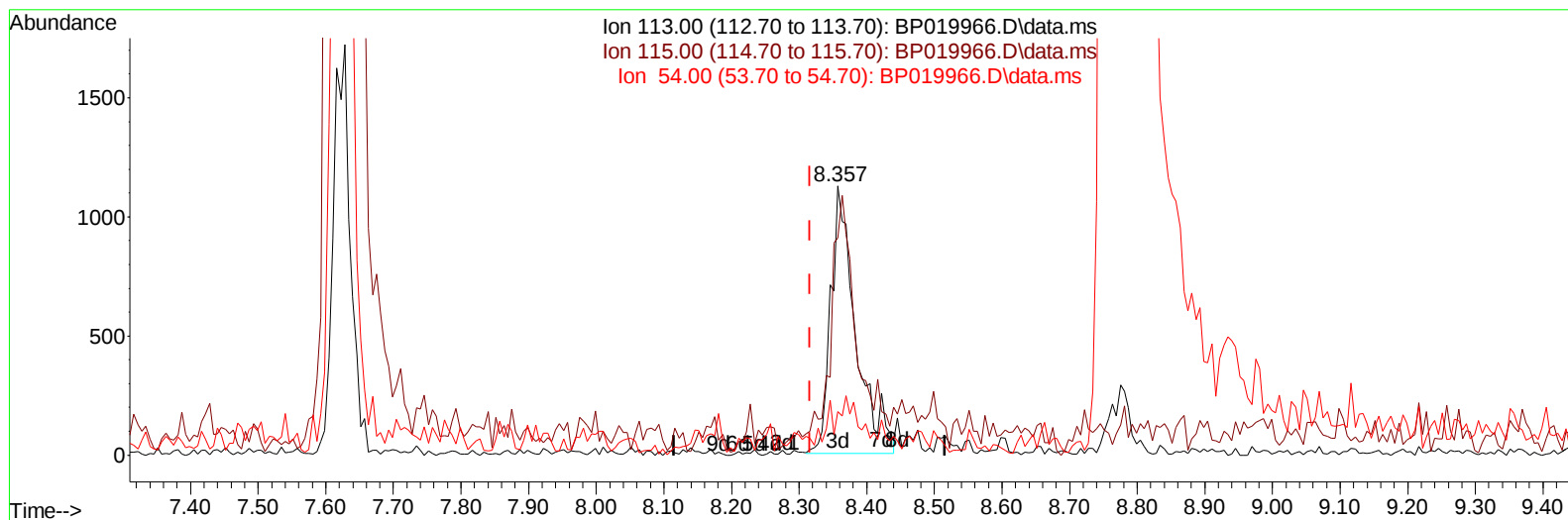
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TIC: BP019966.D\data.ms

(15) 4-Methylphenol-d8 (S)

8.357min (+ 0.041) 0.45 ng/u1 m

response 2867

Ion	Exp%	Act%
113.00	100.00	100.00
115.00	90.00	80.69
54.00	13.80	16.12
0.00	0.00	0.00

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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 03/26/2024
 Supervised By :mohammad ahmed 03/26/2024

Quant Time: Mar 22 23:28:42 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032224.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Mar 22 22:40:53 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.622	152	98712	20.000	ng/ul	0.00
20) Naphthalene-d8	10.369	136	383544	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.263	164	247698	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.075	188	515964	20.000	ng/ul	-0.02
79) Chrysene-d12	21.551	240	504104	20.000	ng/ul	0.00
88) Perylene-d12	24.833	264	590263	20.000	ng/ul	-0.03
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.246	96	12008	5.330	ng/uL	0.00
4) Pyridine-d5	3.646	84	112674	16.563	ng/ul	0.00
7) Phenol-d5	6.852	99	233	0.028	ng/ul	0.04
9) Bis-(2-Chloroethyl)eth...	6.975	67	179449	33.084	ng/ul	0.00
11) 2-Chlorophenol-d4	7.146	132	25	0.004	ng/ul	-0.01
15) 4-Methylphenol-d8	8.357	113	2867m	0.445	ng/ul	0.04
21) Nitrobenzene-d5	8.775	128	100144	34.065	ng/ul	0.00
24) 2-Nitrophenol-d4	9.463	143	59	0.018	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.004	165	95	0.016	ng/ul	0.00
31) 4-Chloroaniline-d4	10.534	131	240119	30.663	ng/ul	0.01
46) Dimethylphthalate-d6	13.663	166	635599	36.599	ng/ul	0.00
49) Acenaphthylene-d8	13.945	160	709645	34.980	ng/ul	0.00
54) 4-Nitrophenol-d4	14.504	143	49	0.018	ng/ul	0.00
60) Fluorene-d10	15.275	176	560977	37.920	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.422	200	12	0.004	ng/ul	-0.01
73) Anthracene-d10	17.186	188	964887	42.246	ng/ul	0.00
81) Pyrene-d10	19.575	212	1164639	39.377	ng/ul	-0.01
92) Benzo(a)pyrene-d12	24.610	264	1213046	42.777	ng/ul	-0.02

Target Compounds Qvalue

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

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