

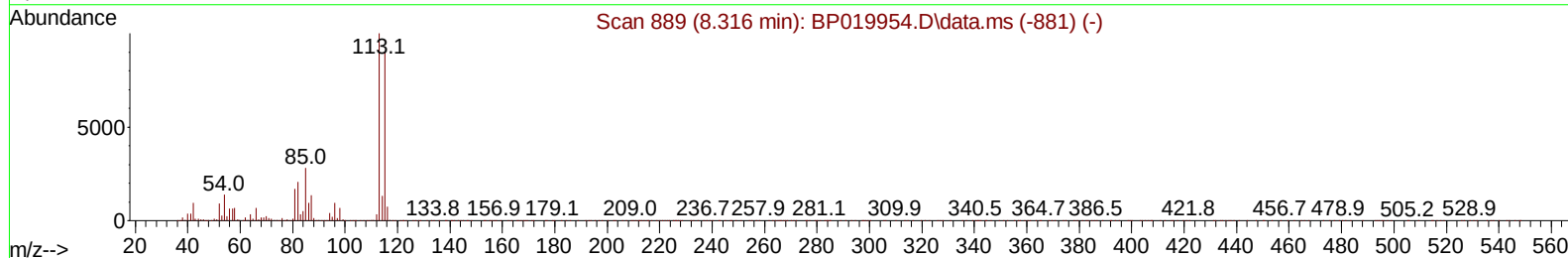
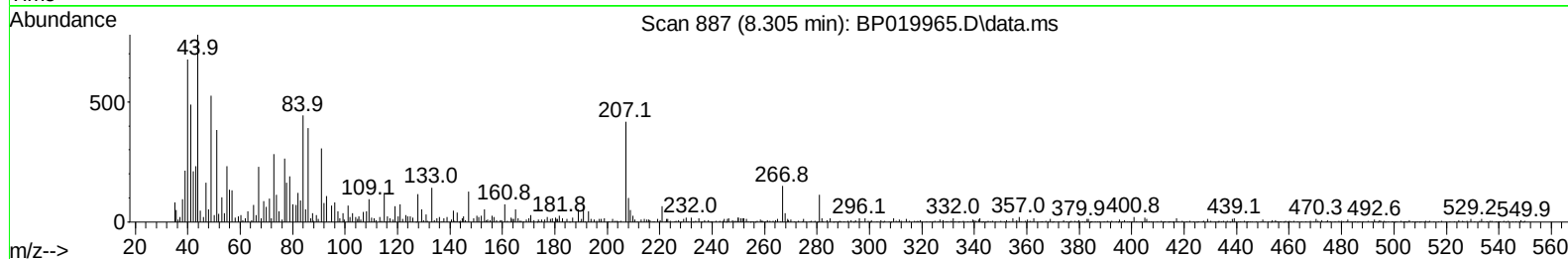
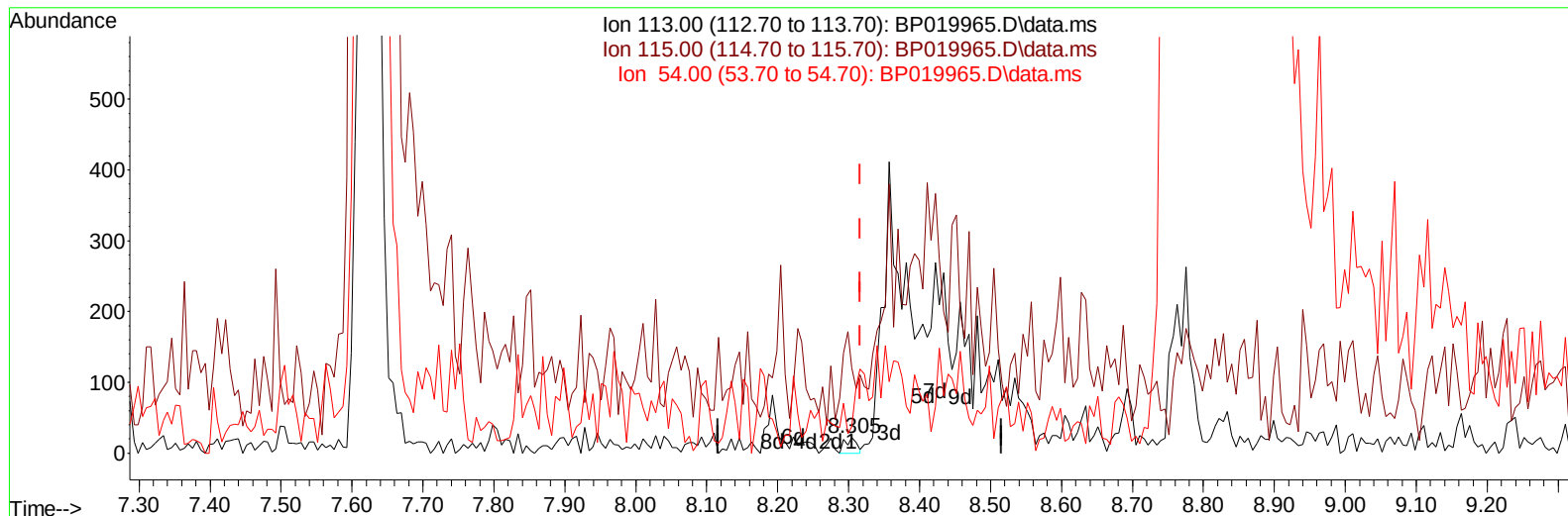
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032224\
Data File : BP019965.D
Acq On : 22 Mar 2024 21:24
Operator : MA/JU
Sample : P1813-12
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E0LY0

Manual IntegrationsAPPROVED

Quant Time: Mar 22 23:03:43 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: BP019965.D\data.ms

(15) 4-Methylphenol-d8 (S)

8.305min (-0.011) 0.00 ng/u1

response 25

Ion	Exp%	Act%
113.00	100.00	100.00
115.00	90.00	563.64#
54.00	13.80	168.18#
0.00	0.00	0.00

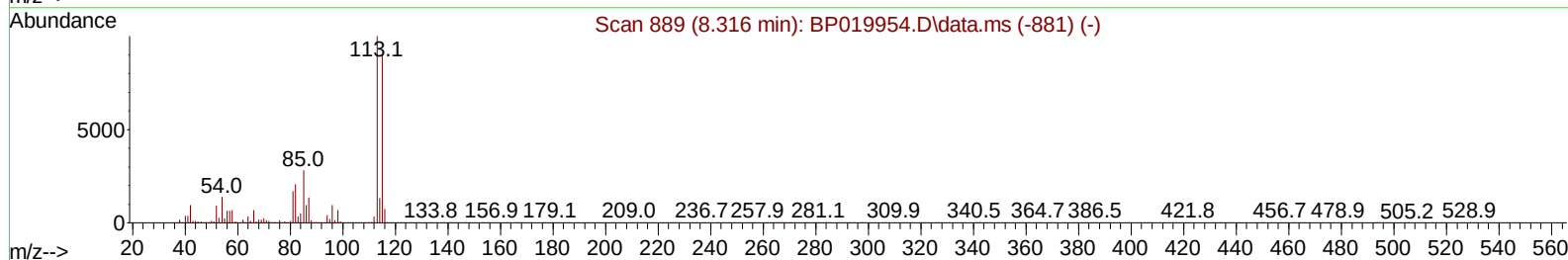
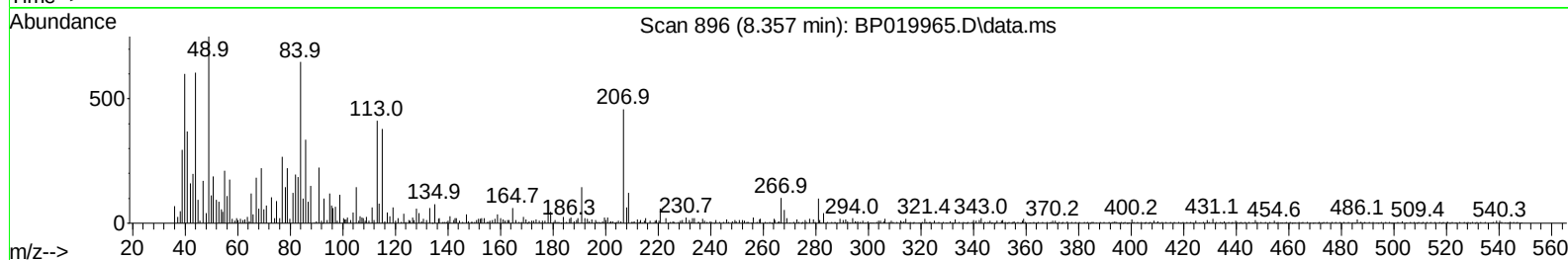
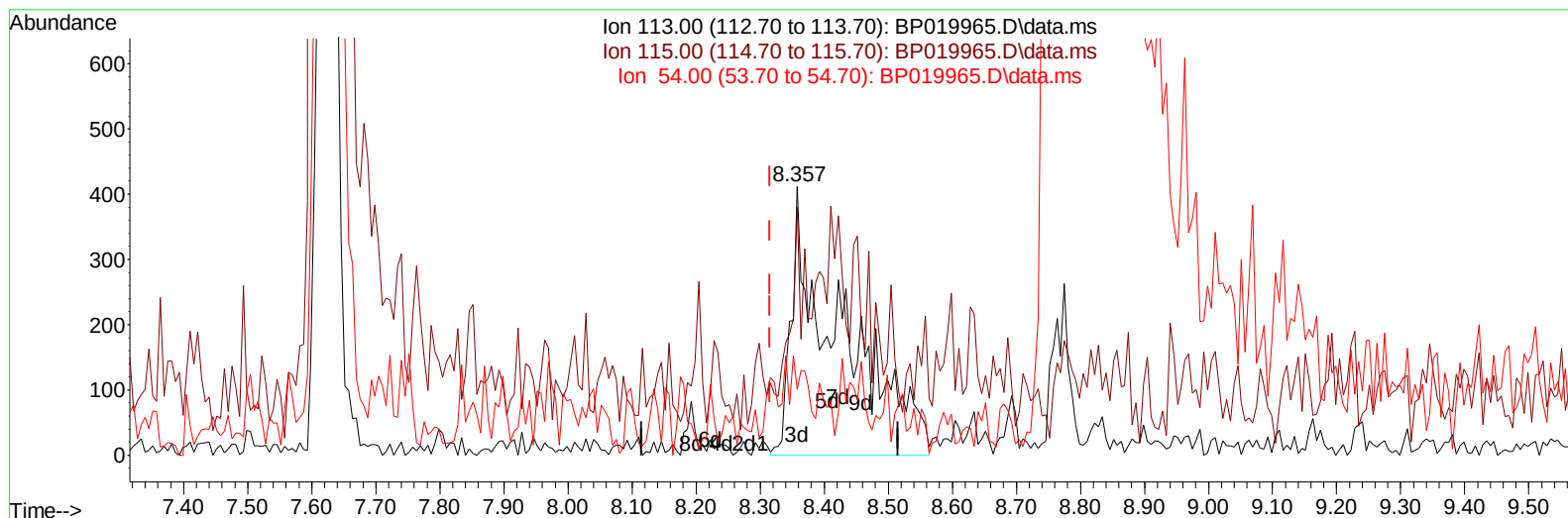
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TIC: BP019965.D\data.ms

(15) 4-Methylphenol-d8 (S)

8.357min (+ 0.041) 0.34 ng/ul m

response 2162

Ion	Exp%	Act%
113.00	100.00	100.00
115.00	90.00	92.23
54.00	13.80	13.35
0.00	0.00	0.00

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Quant Time: Mar 22 23:26:41 2024
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 QLast Update : Fri Mar 22 22:40:53 2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.622	152	98144	20.000	ng/ul	0.00
20) Naphthalene-d8	10.381	136	408329	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.257	164	301580	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.075	188	676551	20.000	ng/ul	-0.02
79) Chrysene-d12	21.539	240	636110	20.000	ng/ul	-0.01
88) Perylene-d12	24.839	264	679100	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.240	96	11078	4.946	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	6.852	99	227	0.027	ng/ul	0.04
9) Bis-(2-Chloroethyl)eth...	6.981	67	168680	31.279	ng/ul	0.00
11) 2-Chlorophenol-d4	7.169	132	80	0.013	ng/ul	0.01
15) 4-Methylphenol-d8	8.357	113	2162m	0.338	ng/ul	0.04
21) Nitrobenzene-d5	8.769	128	92801	29.651	ng/ul	0.00
24) 2-Nitrophenol-d4	9.463	143	16	0.005	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	9.887	165	18	0.003	ng/ul	-0.12
31) 4-Chloroaniline-d4	10.546	131	52	0.006	ng/ul	0.02
46) Dimethylphthalate-d6	13.669	166	729953	34.522	ng/ul	0.00
49) Acenaphthylene-d8	13.945	160	612452	24.796	ng/ul	0.00
54) 4-Nitrophenol-d4	14.498	143	47	0.014	ng/ul	0.00
60) Fluorene-d10	15.281	176	602049	33.425	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.404	200	25	0.006	ng/ul	-0.03
73) Anthracene-d10	17.192	188	970958	32.421	ng/ul	0.00
81) Pyrene-d10	19.575	212	1211083	32.450	ng/ul	-0.01
92) Benzo(a)pyrene-d12	24.610	264	1014714	31.102	ng/ul	-0.02

Target Compounds Qvalue

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

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