

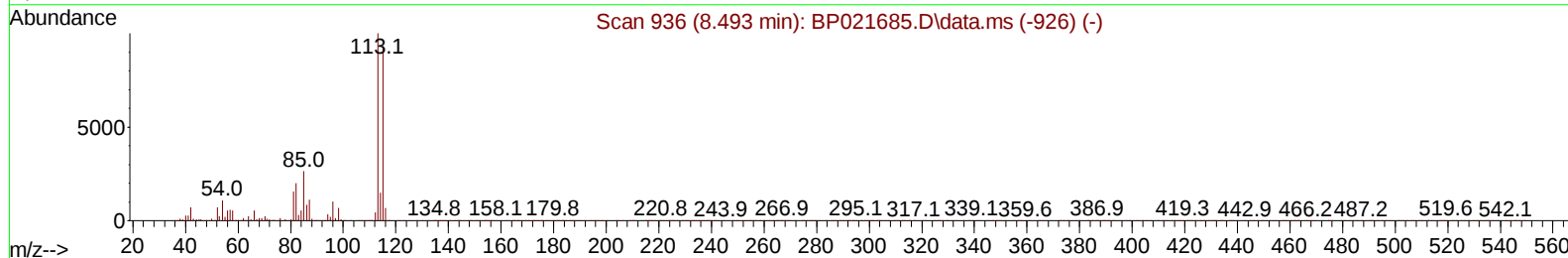
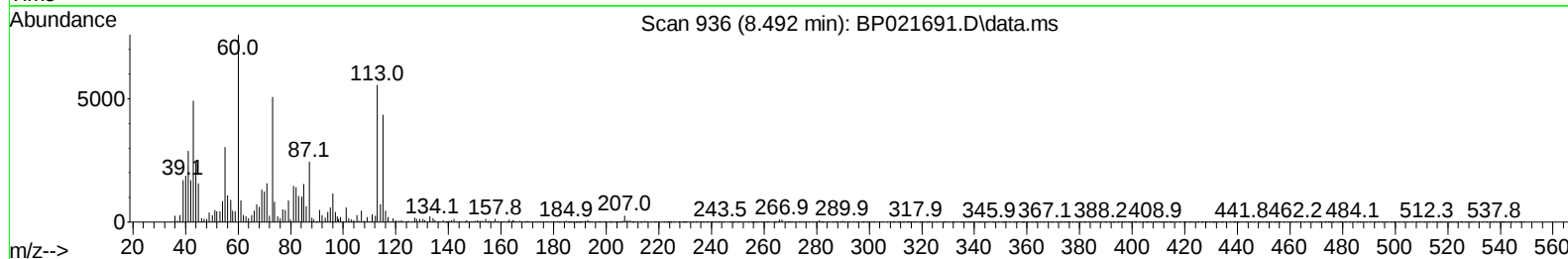
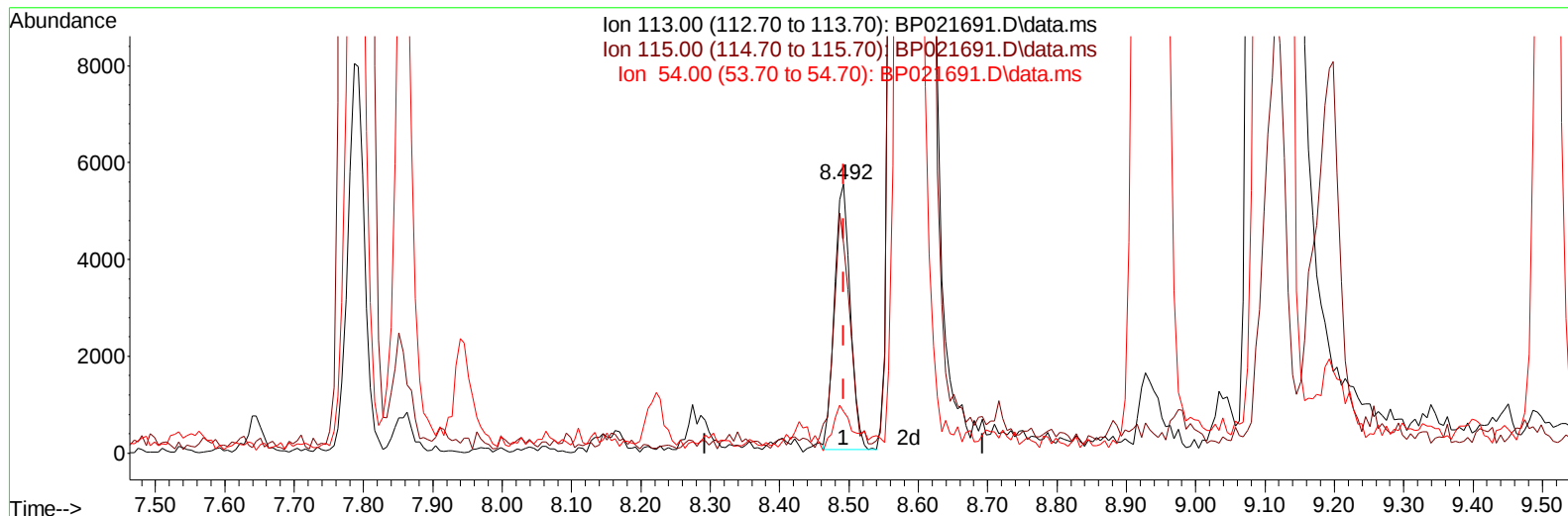
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082624\
Data File : BP021691.D
Acq On : 28 Aug 2024 05:57
Operator : RC/JU
Sample : P3697-04 10X
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCN88

Manual IntegrationsAPPROVED

Quant Time: Aug 28 13:56:52 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Aug 27 05:11:38 2024
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 08/28/2024
Supervised By :mohammad ahmed 08/29/2024



TIC: BP021691.D\data.ms

(15) 4-Methylphenol-d8 (S)

8.492min (-0.000) 0.27 ng/u1

response 8648

Ion	Exp%	Act%
113.00	100.00	100.00
115.00	91.10	78.20
54.00	10.60	15.38#
0.00	0.00	0.00

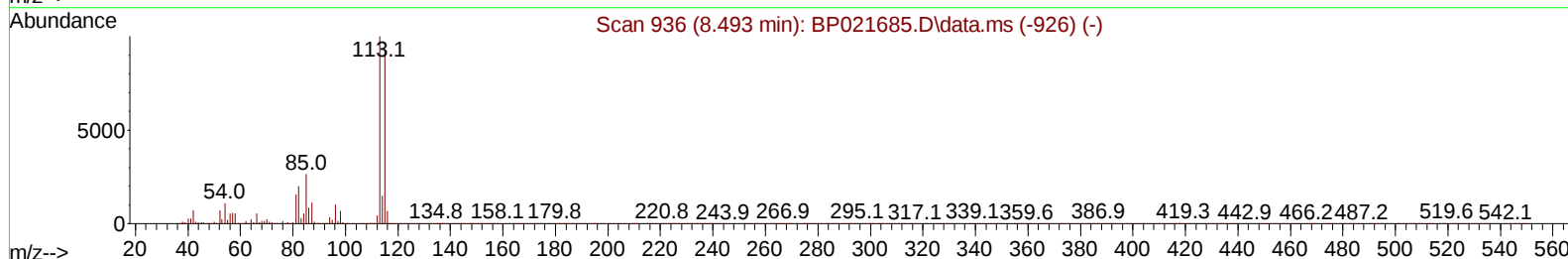
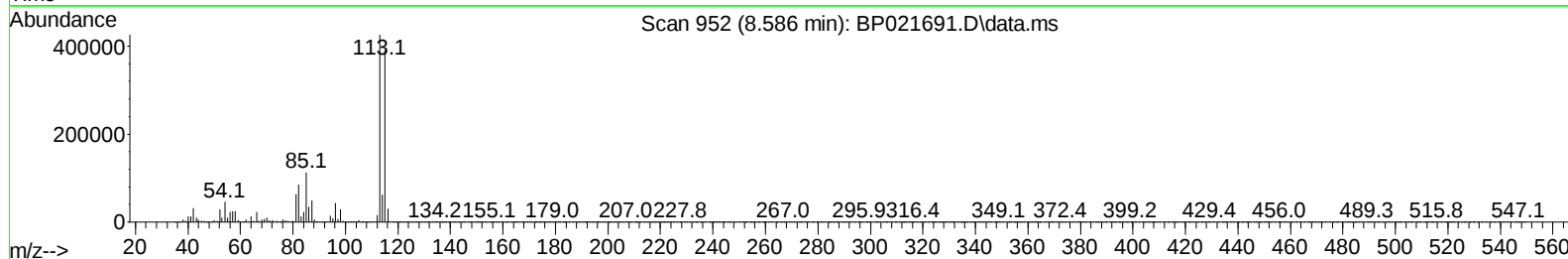
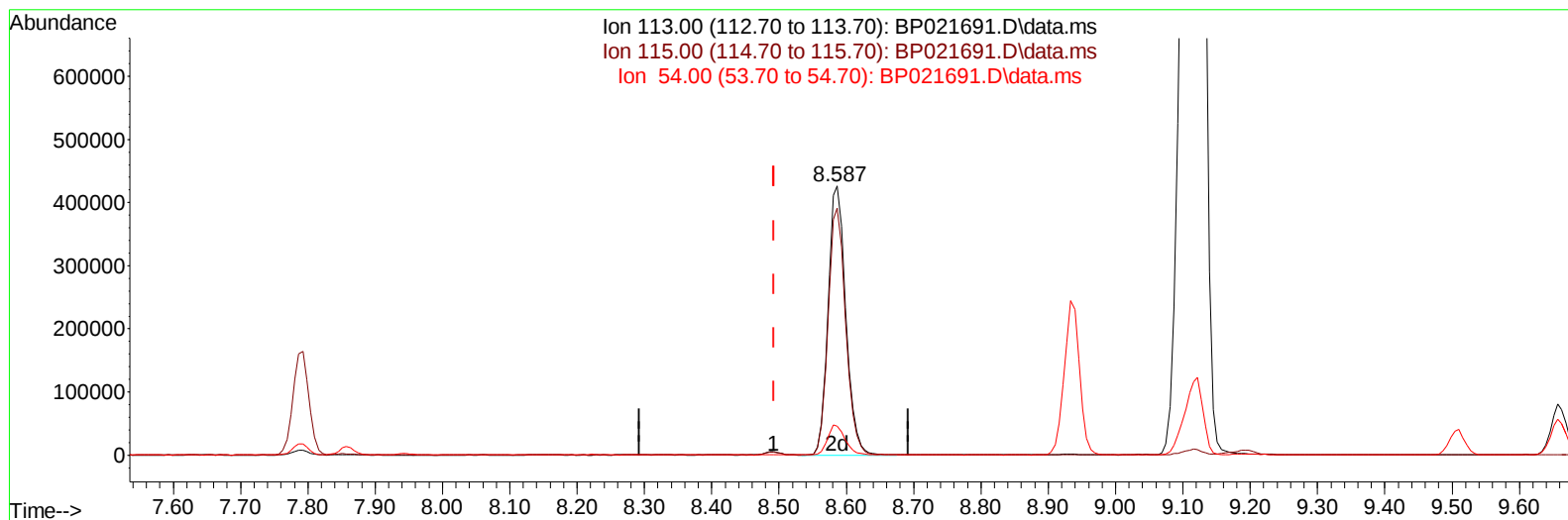
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TIC: BP021691.D\data.ms

(15) 4-Methylphenol-d8 (S)

8.586min (+ 0.094) 23.98 ng/ul m

response 779233

Ion	Exp%	Act%
113.00	100.00	100.00
115.00	91.10	91.72
54.00	10.60	10.75
0.00	0.00	0.00

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 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 GCN88

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 08/28/2024
 Supervised By :mohammad ahmed 08/29/2024

Quant Time: Aug 28 13:58:15 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Aug 27 05:11:38 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.792	152	394057	20.000	ng/ul	0.00
20) Naphthalene-d8	10.581	136	1683181	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.439	164	1081520	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.251	188	2215495	20.000	ng/ul	0.00
79) Chrysene-d12	21.709	240	2207678	20.000	ng/ul	0.02
88) Perylene-d12	25.103	264	2446167	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.275	96	40486	3.411	ng/uL	0.00
4) Pyridine-d5	3.693	84	204298	5.951	ng/ul	0.00
7) Phenol-d5	7.034	99	501309	11.921	ng/ul	0.08
9) Bis-(2-Chloroethyl)eth...	7.122	67	569892	24.901	ng/ul	0.00
11) 2-Chlorophenol-d4	7.334	132	756901	27.878	ng/ul	0.00
15) 4-Methylphenol-d8	8.586	113	779233m	23.979	ng/ul	0.09
21) Nitrobenzene-d5	8.934	128	354011	26.142	ng/ul	0.00
24) 2-Nitrophenol-d4	9.657	143	380100	27.652	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.263	165	792734	29.265	ng/ul	0.06
31) 4-Chloroaniline-d4	10.728	131	357922	9.082	ng/ul	0.02
46) Dimethylphthalate-d6	13.839	166	2245962	27.140	ng/ul	0.00
49) Acenaphthylene-d8	14.127	160	2512204	27.039	ng/ul	0.00
54) 4-Nitrophenol-d4	14.674	143	201110	12.774	ng/ul	0.05
60) Fluorene-d10	15.445	176	1930004	27.740	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.563	200	324223	26.608	ng/ul	0.00
73) Anthracene-d10	17.351	188	2967919	28.227	ng/ul	0.00
81) Pyrene-d10	19.727	212	3640552	28.847	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.886	264	3745373	28.714	ng/ul	0.01
Target Compounds						
					Qvalue	
71) Pentachlorophenol	16.915	266	12429366	668.870	ng/ul	99

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

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