

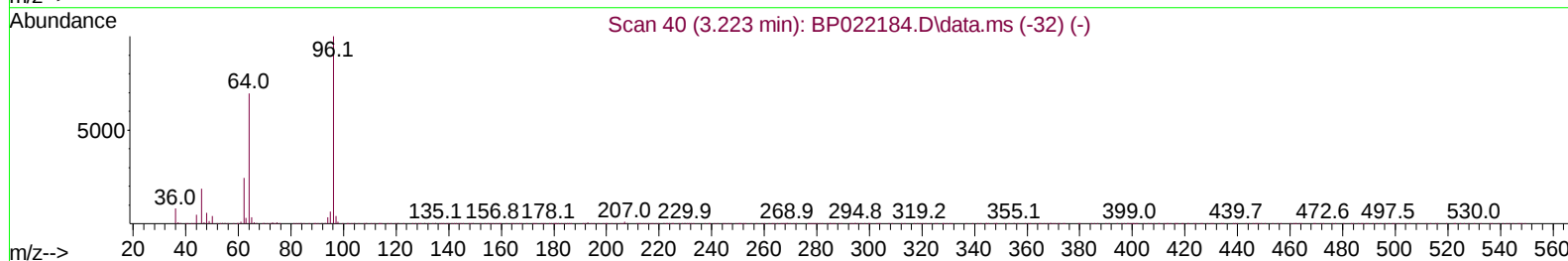
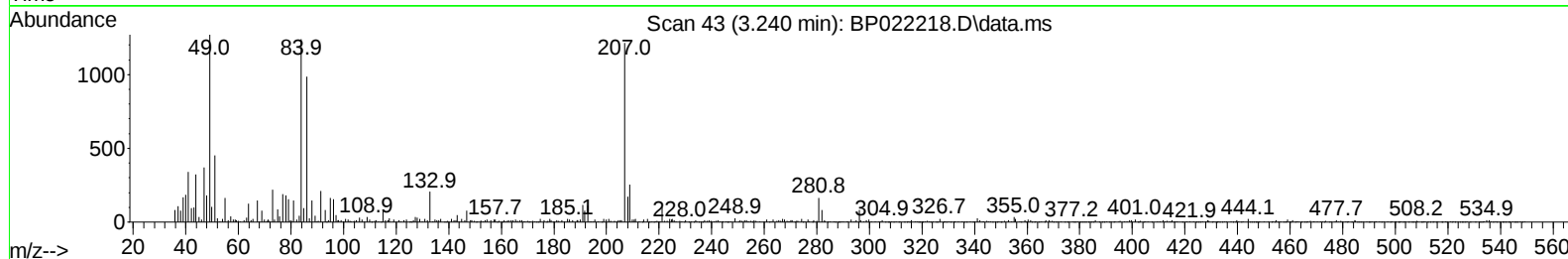
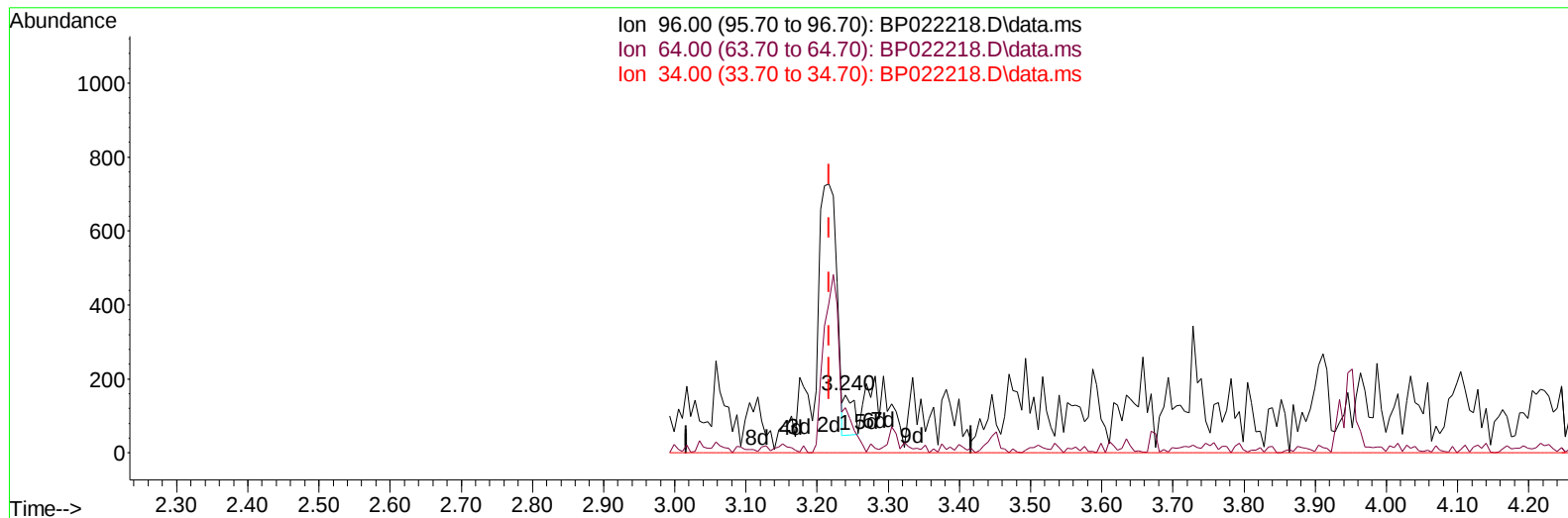
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022218.D
Acq On : 04 Oct 2024 07:24
Operator : RC/JU
Sample : P4017-20DL 10X
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDCD6DL

Manual IntegrationsAPPROVED

Quant Time: Oct 04 08:29:56 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Oct 04 05:25:28 2024
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 10/04/2024
Supervised By :mohammad ahmed 10/05/2024



TIC: BP022218.D\data.ms

(3) 1,4-Dioxane-d8 (S)

3.240min (+ 0.024) 0.03 ng/uL

response 102

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	72.90	78.85
34.00	0.00	0.00
0.00	0.00	0.00

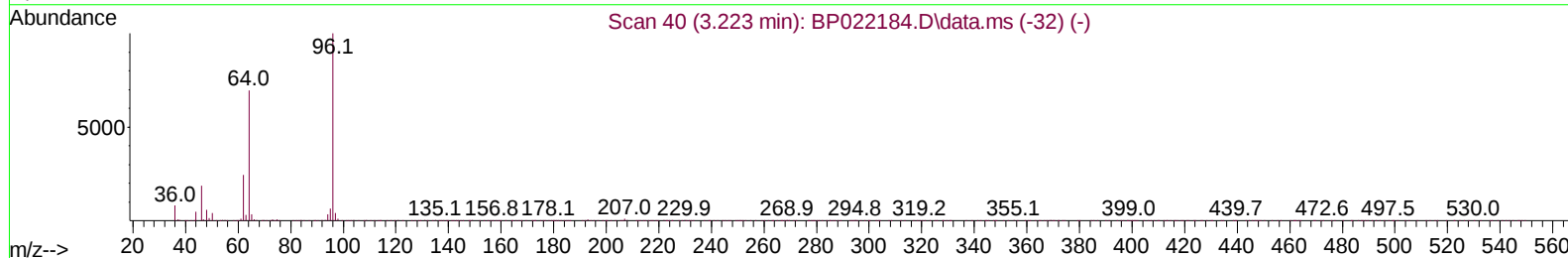
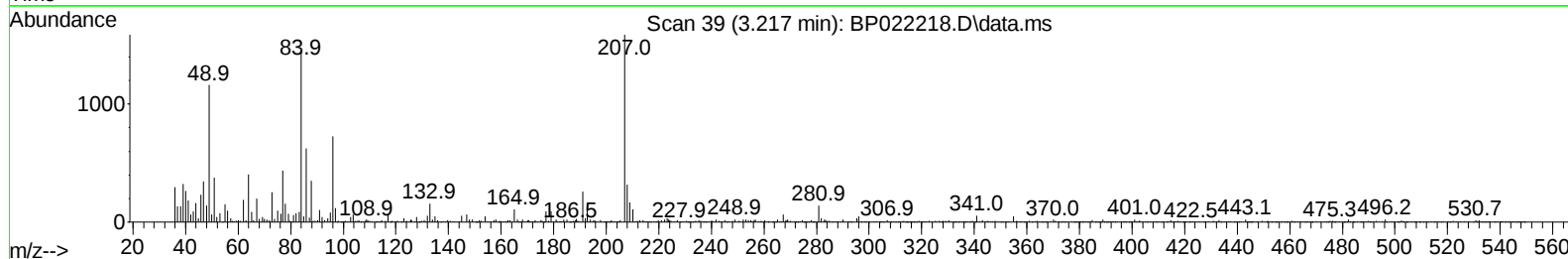
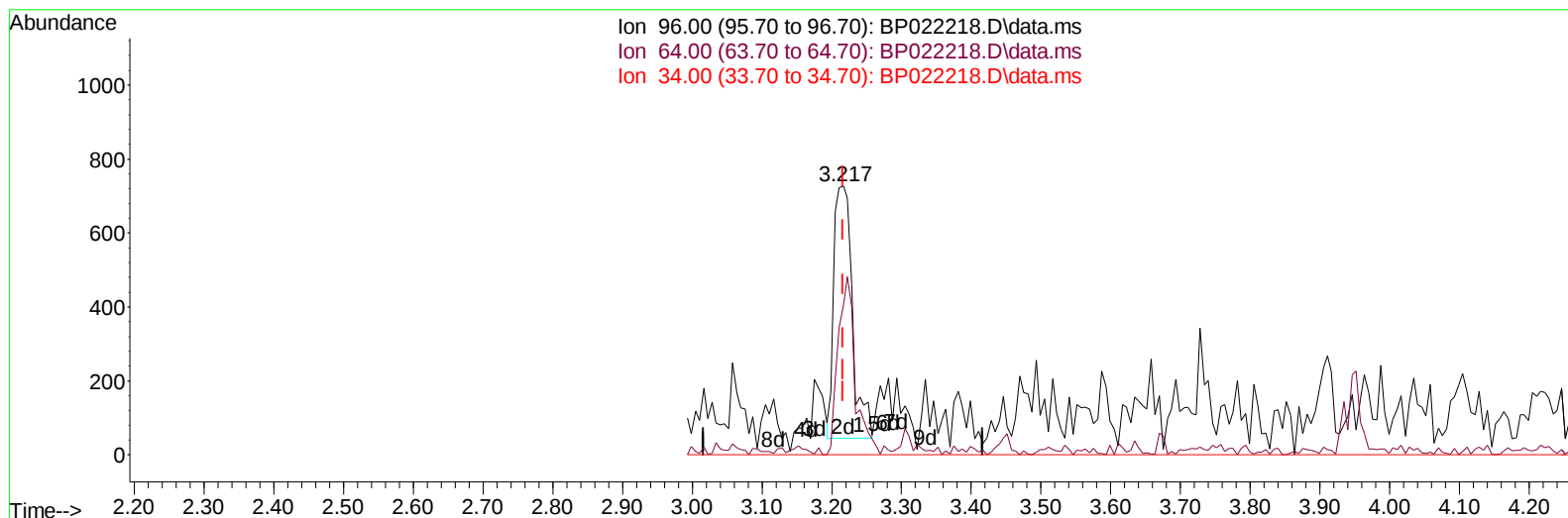
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TIC: BP022218.D\data.ms

(3) 1,4-Dioxane-d8 (S)

3.217min (-0.000) 0.40 ng/uL m

response 1252

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	72.90	55.85#
34.00	0.00	0.00
0.00	0.00	0.00

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Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 10/04/2024
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Quant Time: Oct 04 08:31:55 2024
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 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Oct 04 05:25:28 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.693	152	121434	20.000	ng/ul	0.00
20) Naphthalene-d8	10.451	136	458453	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.310	164	377074	20.000	ng/ul	0.01
64) Phenanthrene-d10	17.110	188	880976	20.000	ng/ul	0.02
79) Chrysene-d12	21.527	240	999439	20.000	ng/ul	0.00
88) Perylene-d12	24.804	264	1176726	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.217	96	1252m	0.404	ng/uL	0.00
4) Pyridine-d5	3.646	84	7307	0.825	ng/ul	0.02
7) Phenol-d5	6.875	99	20491	2.063	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.028	67	15142	2.476	ng/ul	0.00
11) 2-Chlorophenol-d4	7.234	132	15750	2.191	ng/ul	0.00
15) 4-Methylphenol-d8	8.399	113	13686	1.739	ng/ul	0.00
21) Nitrobenzene-d5	8.834	128	8937	2.605	ng/ul	0.00
24) 2-Nitrophenol-d4	9.546	143	7721	1.985	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.087	165	16437	1.945	ng/ul	0.00
31) 4-Chloroaniline-d4	10.569	131	26980	2.676	ng/ul	-0.02
46) Dimethylphthalate-d6	13.716	166	65366	2.243	ng/ul	0.01
49) Acenaphthylene-d8	13.998	160	68461	2.218	ng/ul	0.01
54) 4-Nitrophenol-d4	14.528	143	6534	1.311	ng/ul	0.01
60) Fluorene-d10	15.316	176	59452	2.306	ng/ul	0.01
65) 4,6-Dinitro-2-methylph...	15.445	200	9445	1.727	ng/ul	0.01
73) Anthracene-d10	17.204	188	95699	2.327	ng/ul	0.00
81) Pyrene-d10	19.580	212	125190	2.418	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.574	264	129037	2.085	ng/ul	0.00
Target Compounds						
					Qvalue	
30) Naphthalene	10.504	128	246903	10.235	ng/ul	98
36) 2-Methylnaphthalene	12.110	142	705879	39.376	ng/ul	100
37) 1-Methylnaphthalene	12.328	142	372954	20.628	ng/ul	100
58) 2,3,4,6-Tetrachlorophenol	14.945	232	832368	95.547	ng/ul	98
71) Pentachlorophenol	16.775	266	6383366	744.784	ng/ul	99
72) Phenanthrene	17.151	178	102725	2.224	ng/ul	97

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

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