

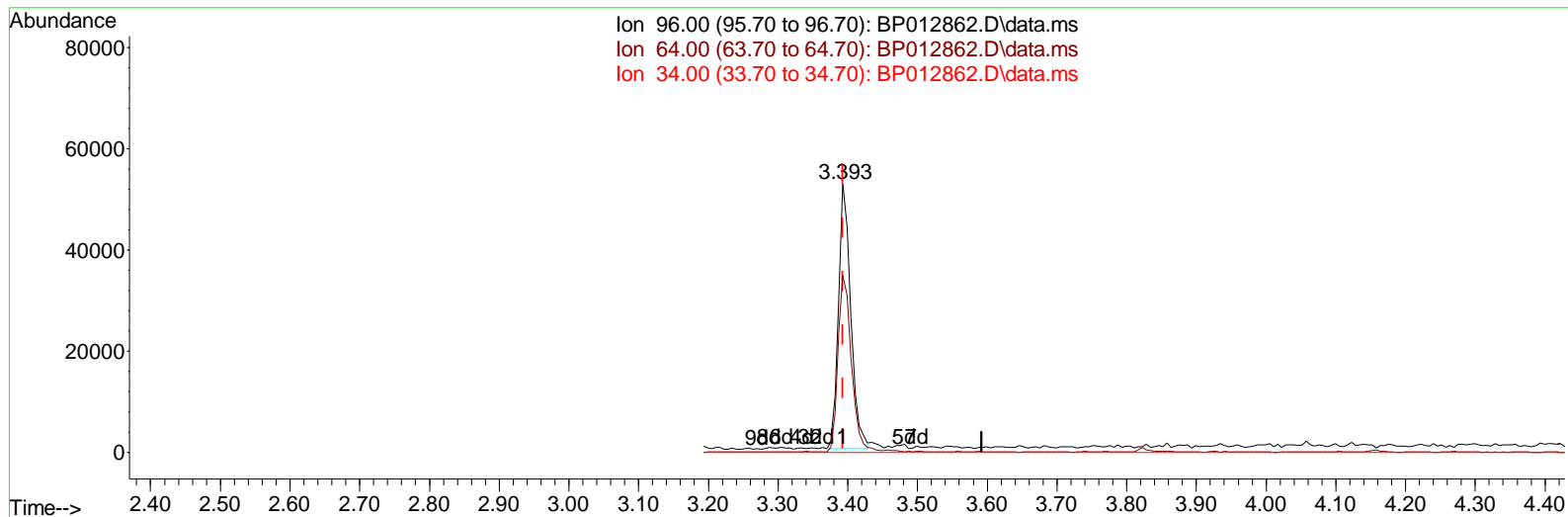
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112922\
 Data File : BP012862.D
 Acq On : 29 Nov 2022 20:54
 Operator : CG/JU
 Sample : N5725-05
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_P
ClientSampleId :
 BHB59

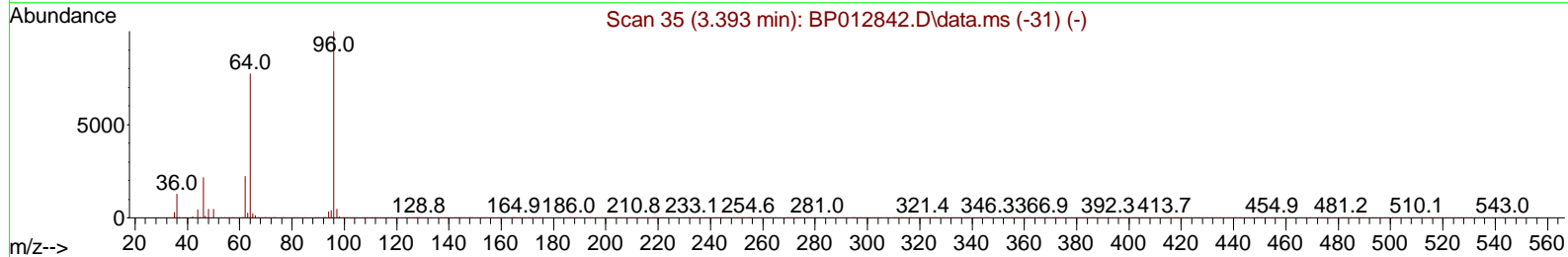
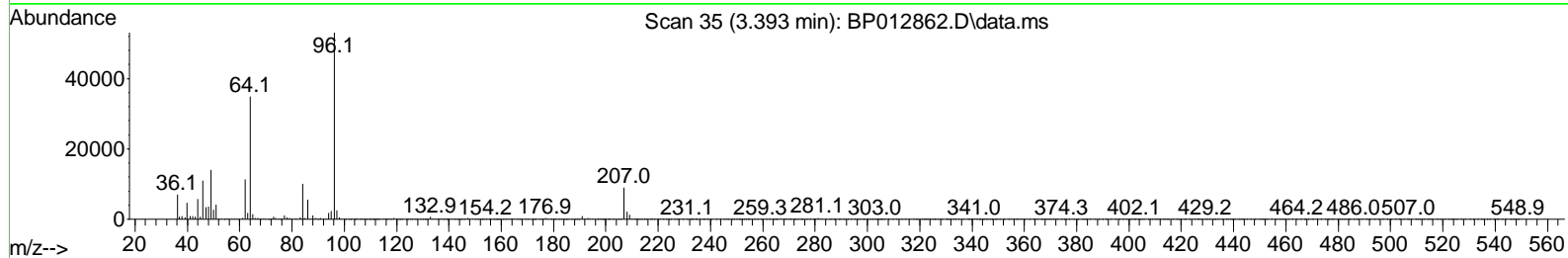
Manual Integrations APPROVED

Reviewed By : Jagrut Upadhyay 11/30/2022
 Supervised By : mohammad ahmed 11/30/2022

Quant Time: Nov 29 22:59:24 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Nov 29 22:53:34 2022
 Response via : Initial Calibration



Ion 96.00 (95.70 to 96.70): BP012862.D\data.ms
 Ion 64.00 (63.70 to 64.70): BP012862.D\data.ms
 Ion 34.00 (33.70 to 34.70): BP012862.D\data.ms



TIC: BP012862.D\data.ms

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

3.393min (+ 0.000) 4.40 ng/uL

response	64545	
Ion	Exp%	Act%
96.00	100.00	100.00
64.00	72.90	65.92
34.00	0.00	0.00
0.00	0.00	0.00

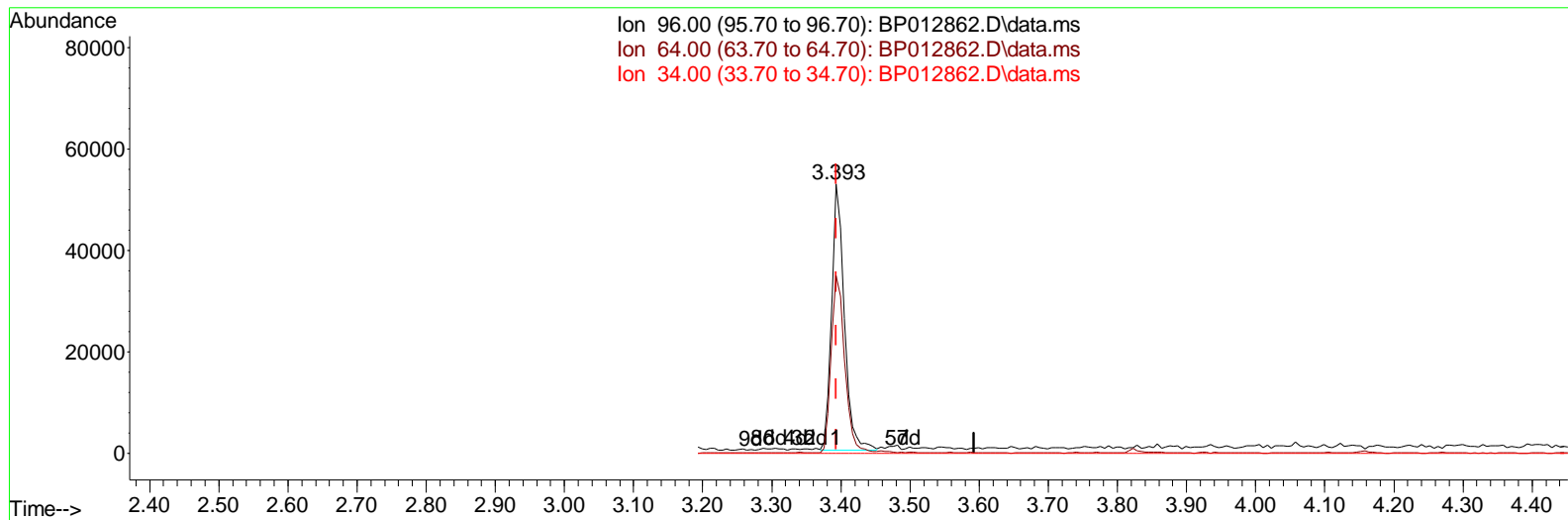
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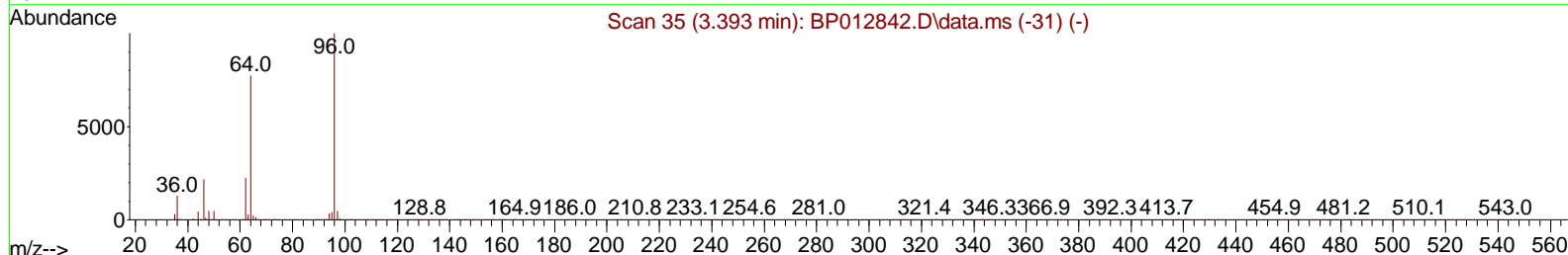
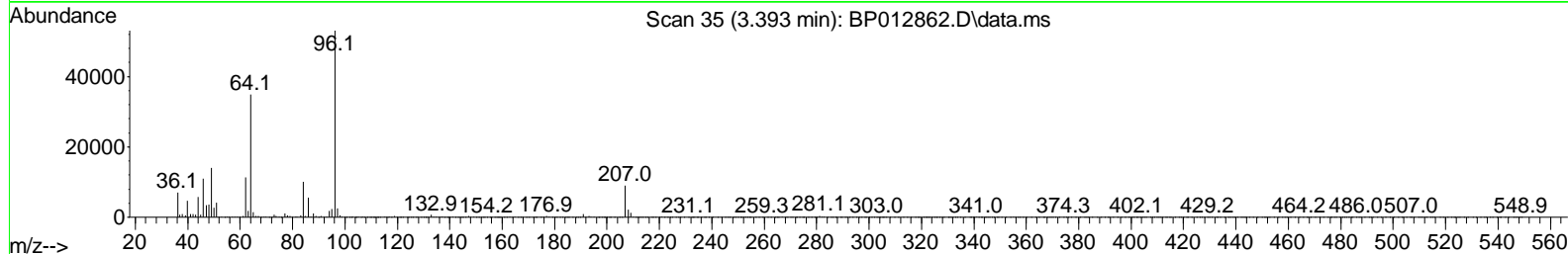
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Ion 96.00 (95.70 to 96.70): BP012862.D\data.ms
 Ion 64.00 (63.70 to 64.70): BP012862.D\data.ms
 Ion 34.00 (33.70 to 34.70): BP012862.D\data.ms



TIC: BP012862.D\data.ms

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

3.393min (+ 0.000) 4.54 ng/uL m

response	66574	
Ion	Exp%	Act%
96.00	100.00	100.00
64.00	72.90	65.92
34.00	0.00	0.00
0.00	0.00	0.00

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Compound	R. T.	QI on	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Di chl orobenzene-d4	8.011	152	628874	20.000	ng/ul	0.00
20) Naphthal ene-d8	10.828	136	2874853	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.646	164	1961837	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.398	188	4178879	20.000	ng/ul	0.00
79) Chrysene-d12	21.486	240	3382610	20.000	ng/ul	0.00
88) Peryl ene-d12	23.992	264	2793147	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Di oxane-d8	3.393	96	66574m	4.536	ng/uL	0.00
4) Pyri di ne-d5	3.823	84	332961	7.581	ng/ul	0.00
7) Phenol -d5	7.158	99	382726	7.406	ng/ul	0.00
9) Bi s-(2-Chl oroethyl)eth. . .	7.334	67	953902	29.957	ng/ul	0.00
11) 2-Chl orophenol -d4	7.540	132	902448	22.918	ng/ul	0.00
15) 4-Methyl phenol -d8	8.711	113	657413	15.790	ng/ul	0.00
21) Ni trobenzene-d5	9.175	128	592927	33.697	ng/ul	0.00
24) 2-Ni trophenol -d4	9.905	143	599743	32.100	ng/ul	0.00
28) 2,4-Di chl orophenol -d3	10.440	165	1115523	25.589	ng/ul	0.00
31) 4-Chl oroani li ne-d4	10.958	131	894651	14.987	ng/ul	0.00
46) Di methyl phthal ate-d6	14.051	166	4380797	32.441	ng/ul	0.00
49) Acenaphthyl ene-d8	14.340	160	4753702	30.171	ng/ul	0.00
54) 4-Ni trophenol -d4	14.828	143	120564	5.008	ng/ul	0.00
60) Fl uorene-d10	15.640	176	3834307	32.422	ng/ul	0.00
65) 4,6-Di ni tro-2-methyl ph. . .	15.751	200	539506	27.012	ng/ul	0.00
73) Anthracene-d10	17.498	188	6256545	34.012	ng/ul	0.00
81) Pyrene-d10	19.728	212	6933379	35.276	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.833	264	4726779	34.452	ng/ul	0.00

Target Compounds Qvalue

(#) = qual ifier out of range (m) = manual i ntegrati on (+) = signal s summed

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