

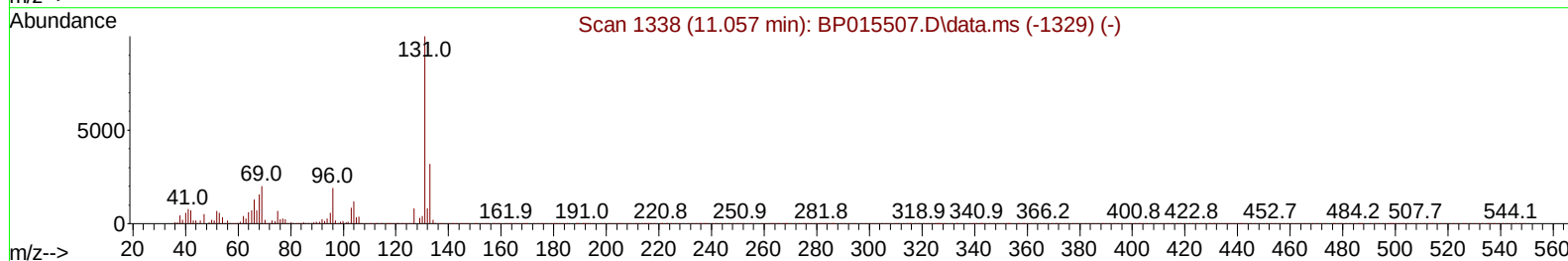
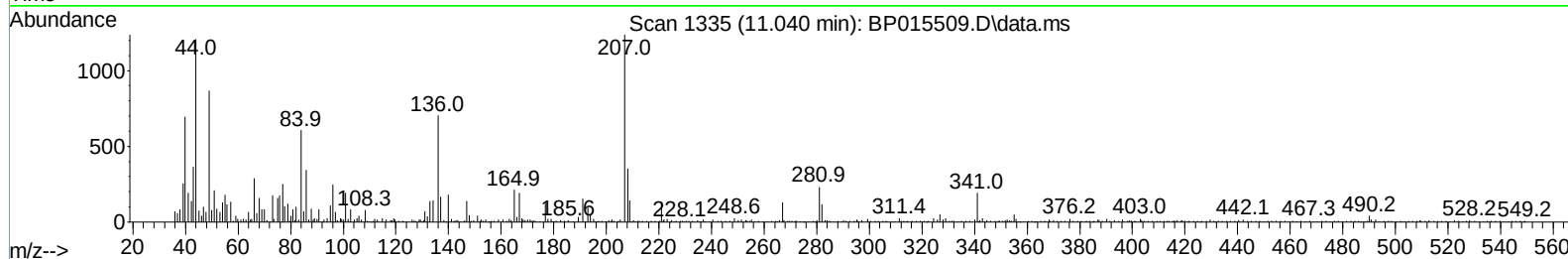
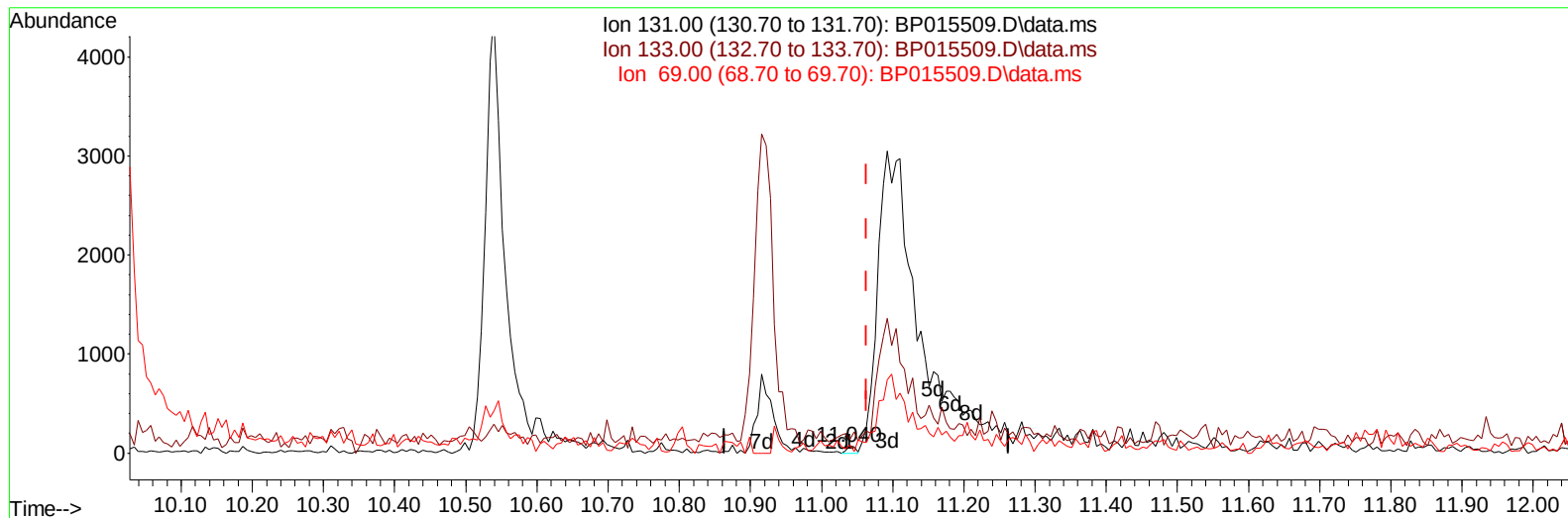
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061423\
Data File : BP015509.D
Acq On : 14 Jun 2023 10:25
Operator : MA/JU
Sample : 03098-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EZEP9

Manual IntegrationsAPPROVED

Quant Time: Jun 14 23:07:24 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Jun 09 23:33:38 2023
Response via : Initial Calibration

Reviewed By :Yogesh Patel 06/15/2023
Supervised By :mohammad ahmed 06/16/2023



TIC: BP015509.D\data.ms

(31) 4-Chloroaniline-d4 (S)

11.040min (-0.024) 0.00 ng/u1

response 57

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	32.20	192.96#
69.00	18.30	115.49#
0.00	0.00	0.00

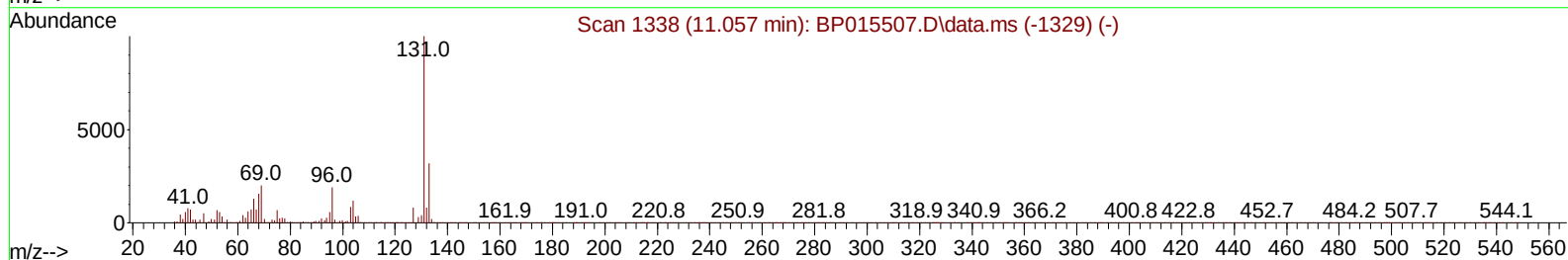
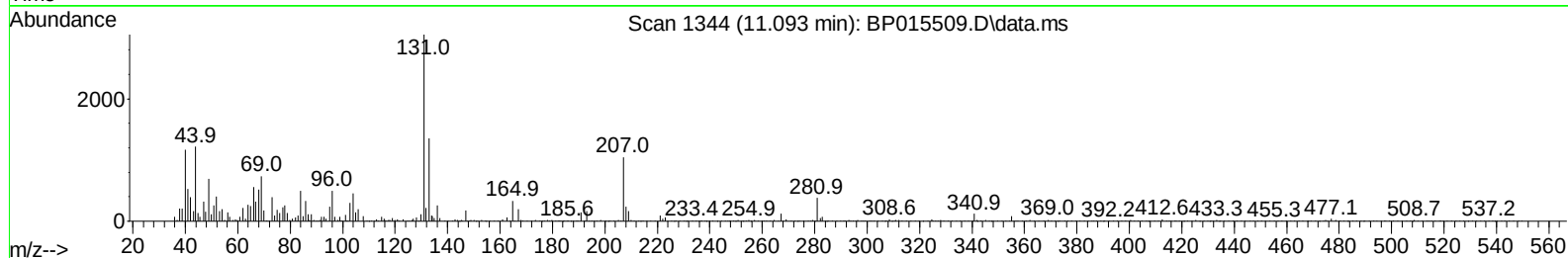
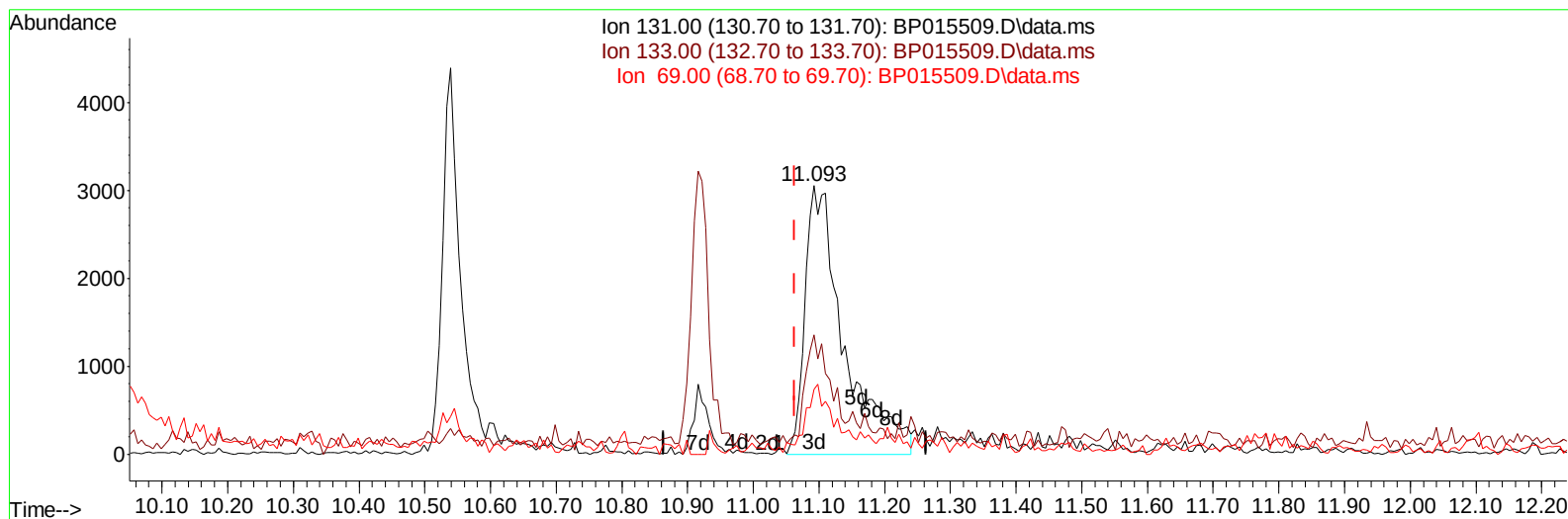
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TIC: BP015509.D\data.ms

(31) 4-Chloroaniline-d4 (S)

11.093min (+ 0.029) 0.75 ng/ul m

response 12562

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	32.20	44.46#
69.00	18.30	24.15#
0.00	0.00	0.00

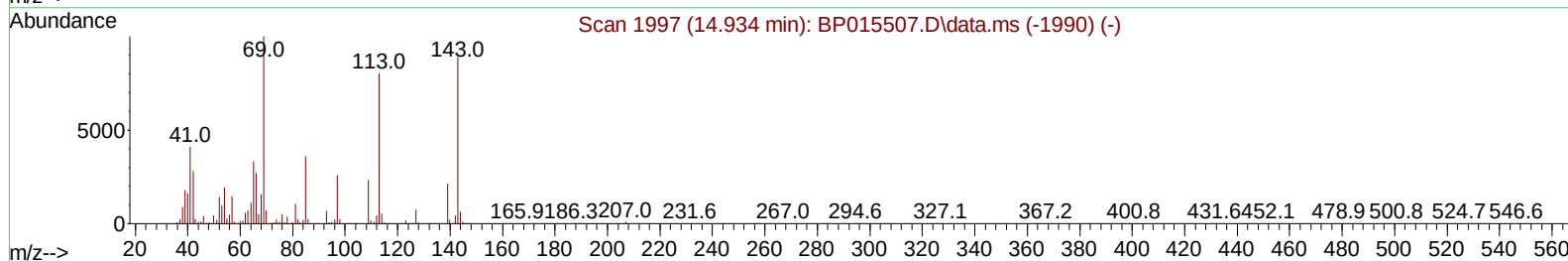
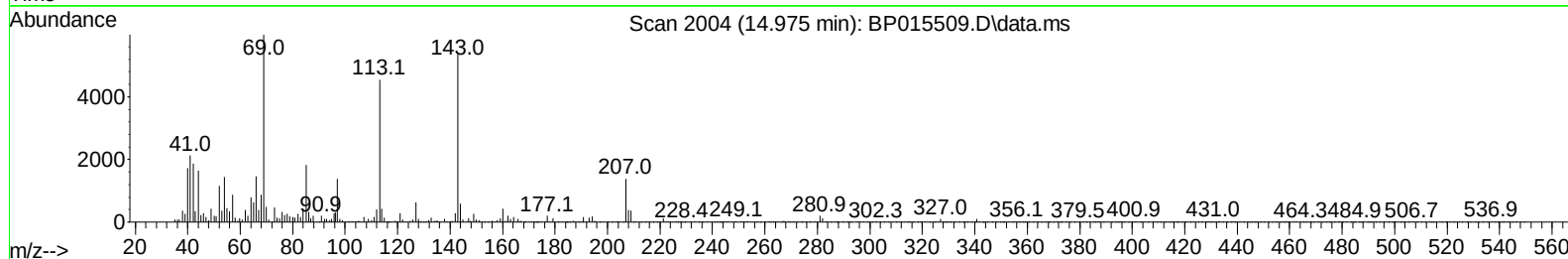
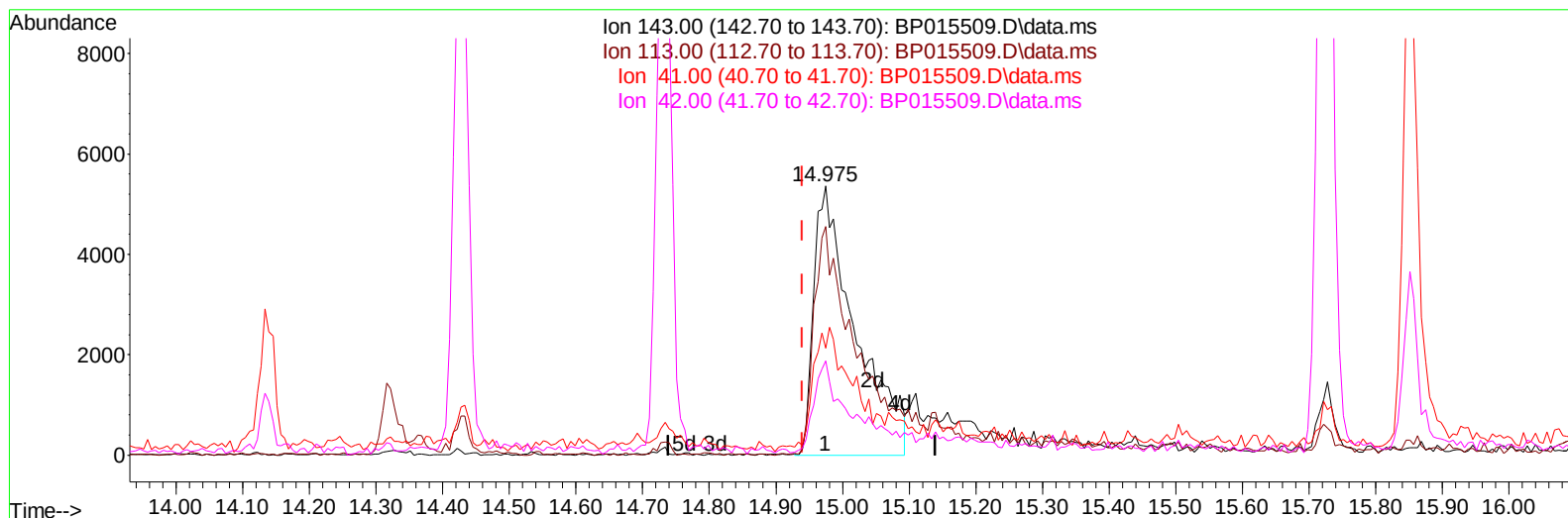
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Instrument :
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TIC: BP015509.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.975min (+ 0.035) 3.23 ng/ul m

response 23086

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	86.60	84.92
41.00	42.20	39.69
42.00	29.40	34.99

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 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 06/15/2023
 Supervised By :mohammad ahmed 06/16/2023

Quant Time: Jun 14 23:17:57 2023
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.093	152	152217	20.000	ng/ul	0.00
20) Naphthalene-d8	10.916	136	712730	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.734	164	505721	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.486	188	1178712	20.000	ng/ul	0.00
79) Chrysene-d12	21.568	240	1165638	20.000	ng/ul	0.00
88) Perylene-d12	24.115	264	1220986	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.405	96	16658	4.052	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	7.246	99	80248	5.723	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.410	67	226930	27.096	ng/ul	0.00
11) 2-Chlorophenol-d4	7.616	132	213356	20.985	ng/ul	0.00
15) 4-Methylphenol-d8	8.804	113	155640	13.523	ng/ul	0.00
21) Nitrobenzene-d5	9.269	128	143902	25.717	ng/ul	0.00
24) 2-Nitrophenol-d4	9.998	143	156077	24.425	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.540	165	254956	21.872	ng/ul	0.00
31) 4-Chloroaniline-d4	11.093	131	12562m	0.751	ng/ul	0.03
46) Dimethylphthalate-d6	14.139	166	1142097	28.639	ng/ul	0.00
49) Acenaphthylene-d8	14.428	160	1178758	27.032	ng/ul	0.00
54) 4-Nitrophenol-d4	14.975	143	23086m	3.233	ng/ul	0.04
60) Fluorene-d10	15.728	176	960602	28.565	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.851	200	183508	24.575	ng/ul	0.00
73) Anthracene-d10	17.586	188	1677709	31.283	ng/ul	0.00
81) Pyrene-d10	19.810	212	2120655	33.992	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.951	264	2007509	32.768	ng/ul	0.00

Target Compounds Qvalue

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

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