

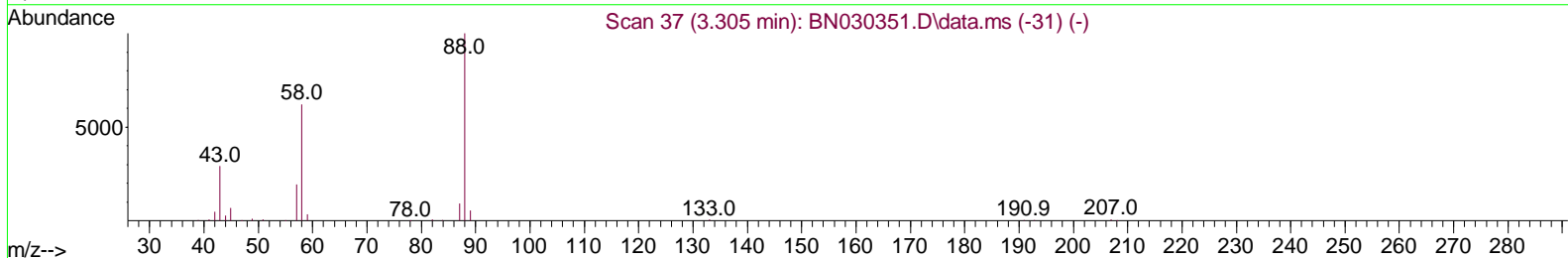
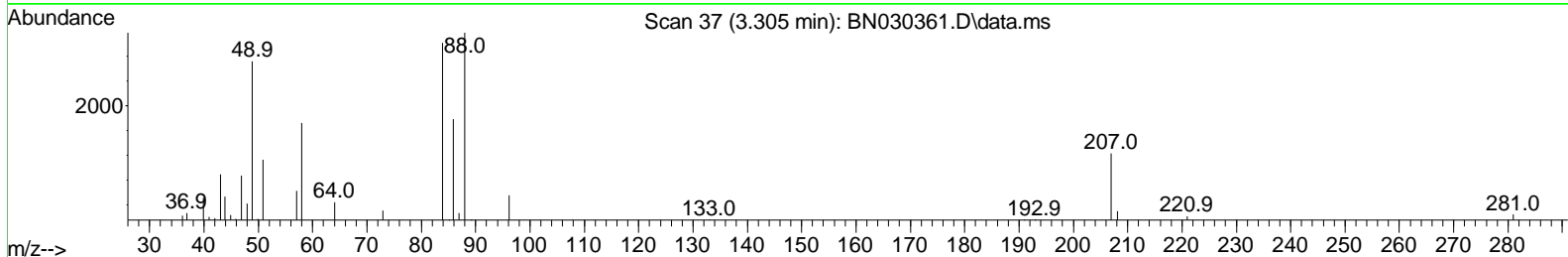
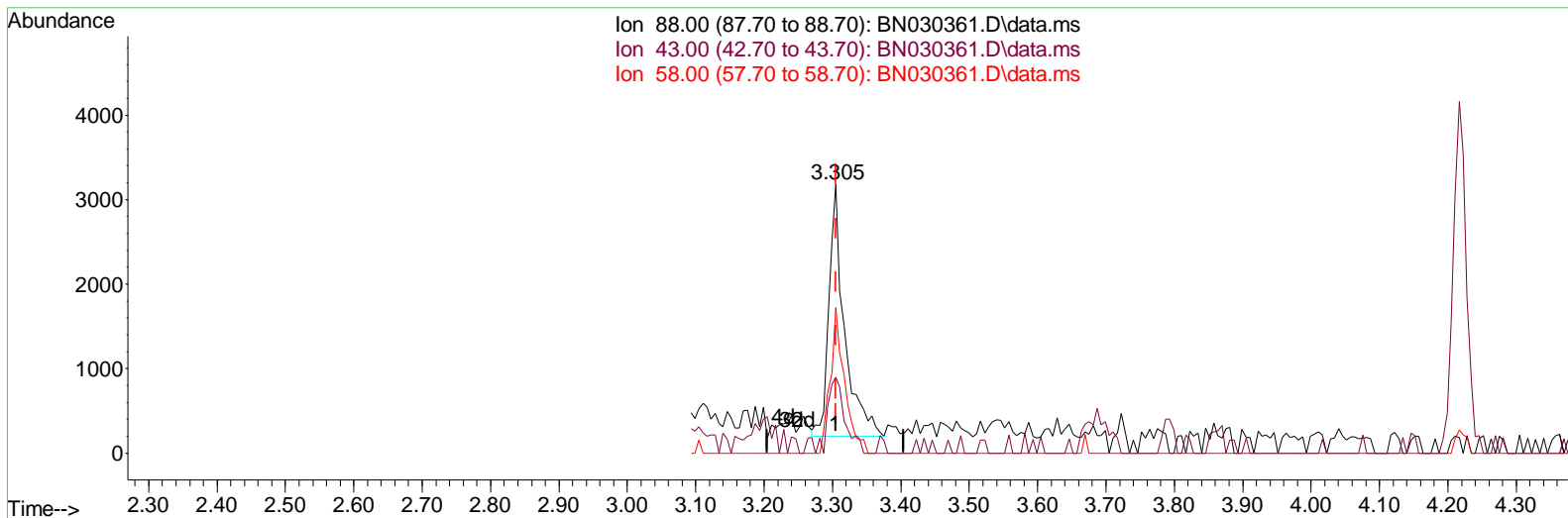
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN031124\
 Data File : BN030361.D
 Acq On : 11 Mar 2024 15:30
 Operator : MA/JU
 Sample : P1601-05
 Misc : SFAM-MDL-ME SOIL-01
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
ClientSampleId :
 MDL-MED-SOIL-QT1-2024-01

Manual Integrations APPROVED

Reviewed By : Yogesh Patel 03/12/2024
 Supervised By : mohammad ahmed 03/12/2024

Quant Time: Mar 11 16:00:23 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN030824.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Mar 08 13:57:34 2024
 Response via : Initial Calibration



TIC: BN030361.D\data.ms

(2) 1,4-Dioxane

3.305min (-0.000) 1.42 ng/uL

response	4699	
Ion	Exp%	Act%
88.00	100.00	100.00
43.00	30.00	28.21
58.00	61.20	54.19
0.00	0.00	0.00

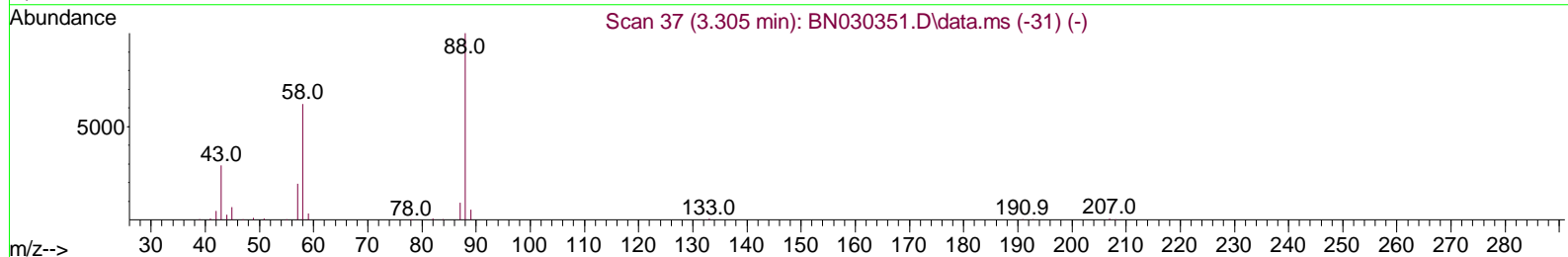
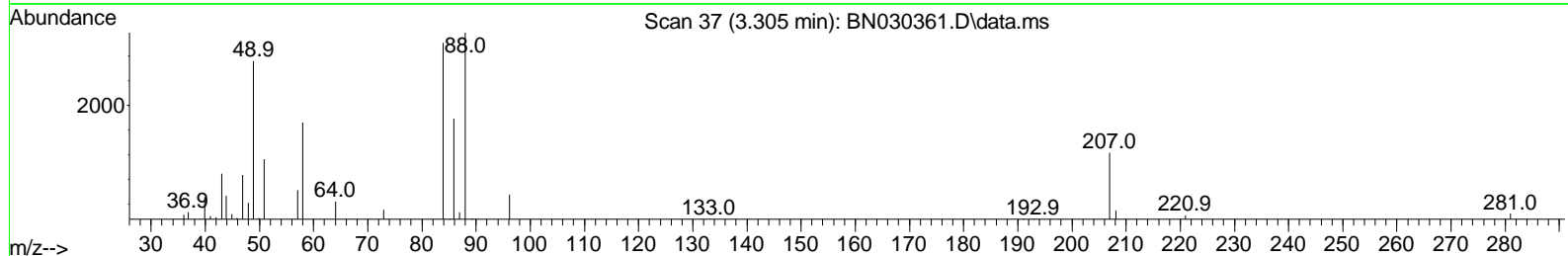
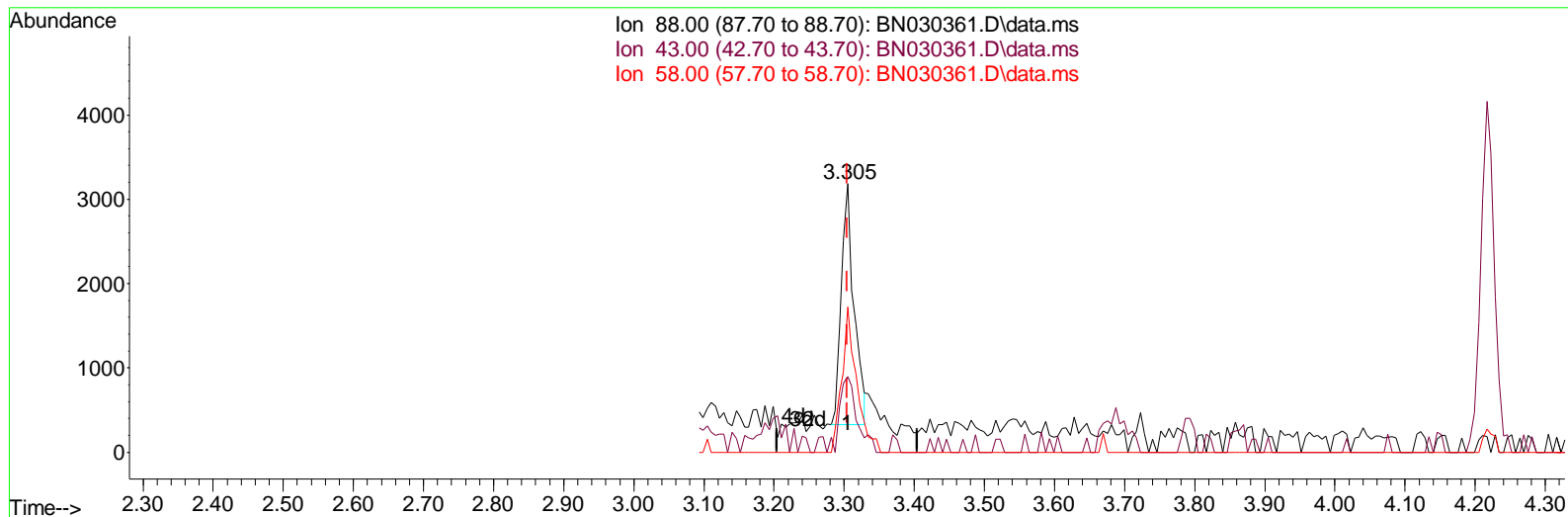
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN031124\
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 Sample : P1601-05
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 ALS Vial : 6 Sample Multiplier: 1

Instrument :
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Manual Integrations APPROVED

Quant Time: Mar 11 16:00:23 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN030824.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Mar 08 13:57:34 2024
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Reviewed By :Yogesh Patel 03/12/2024
 Supervised By :mohammad ahmed 03/12/2024



TIC: BN030361.D\data.ms

(2) 1,4-Dioxane

3.305min (-0.000) 1.10 ng/uL m

response	3634
Ion	Exp% Act%
88.00	100.00 100.00
43.00	30.00 28.21
58.00	61.20 54.19
0.00	0.00 0.00

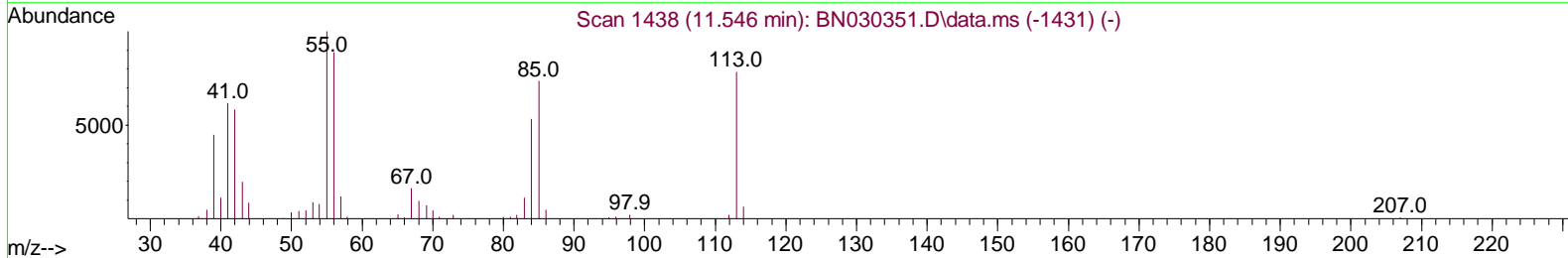
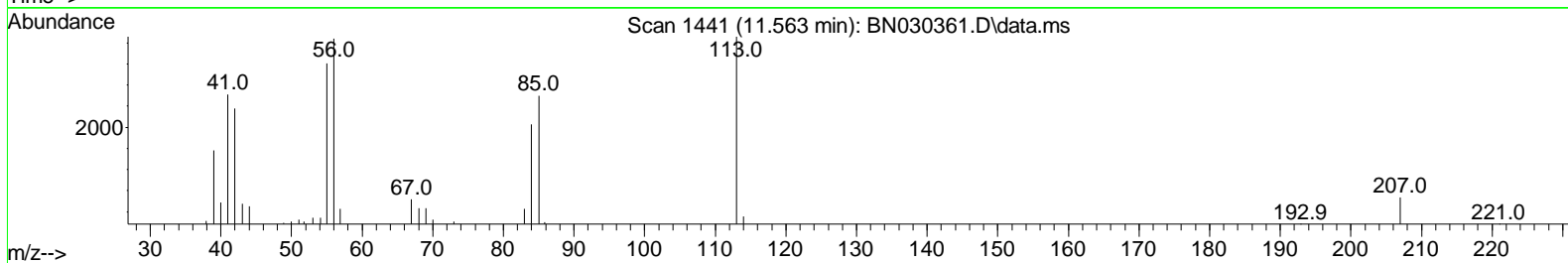
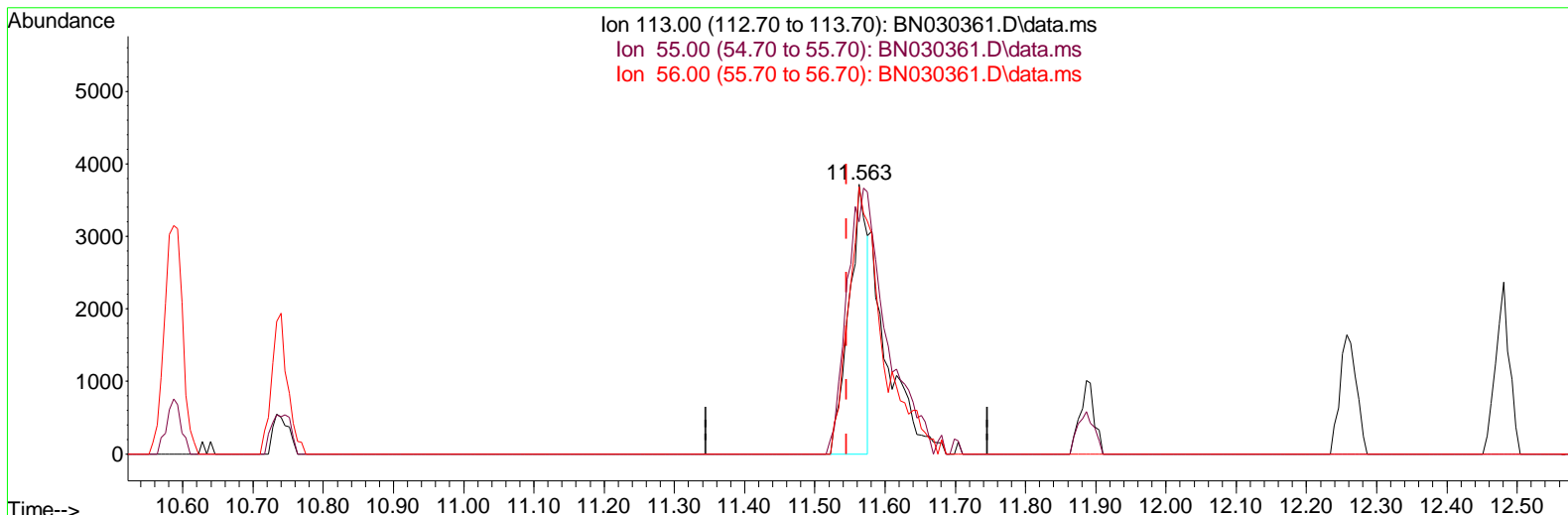
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN031124\
 Data File : BN030361.D
 Acq On : 11 Mar 2024 15:30
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TIC: BN030361.D\data.ms

(34) Caprolactam

11.563min (+ 0.018) 2.94 ng/ul

response 6701

Ion	Exp%	Act%
113.00	100.00	100.00
55.00	127.60	86.30#
56.00	113.50	99.11
0.00	0.00	0.00

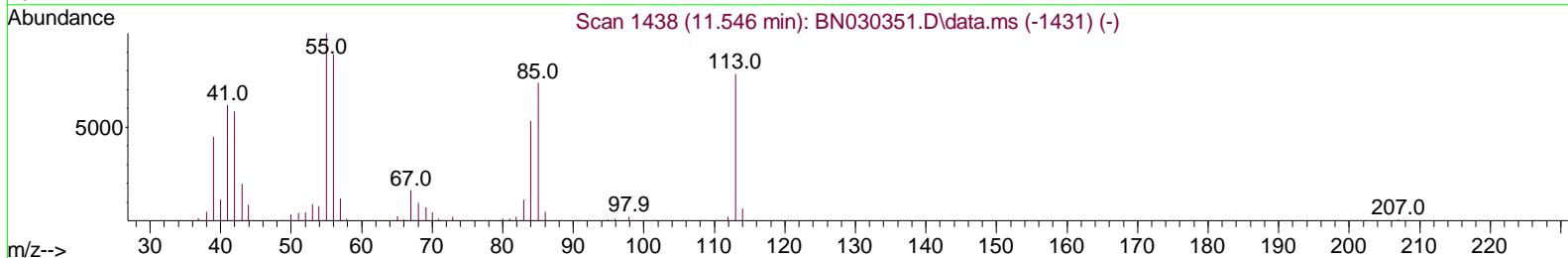
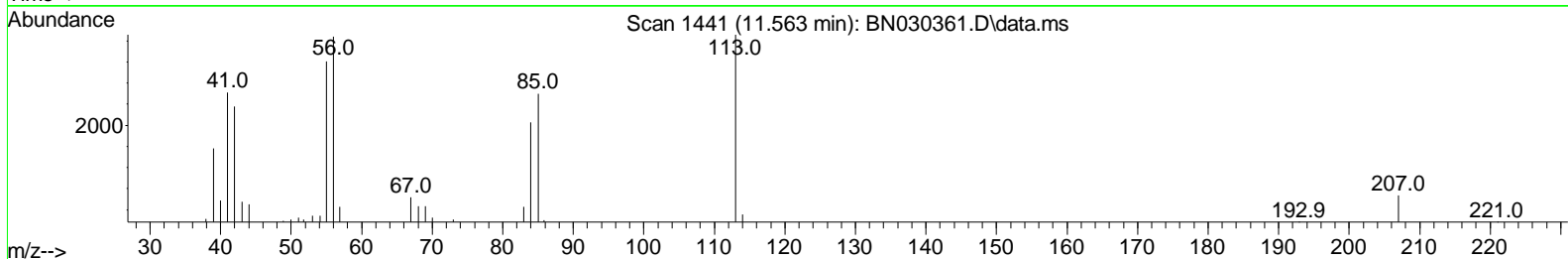
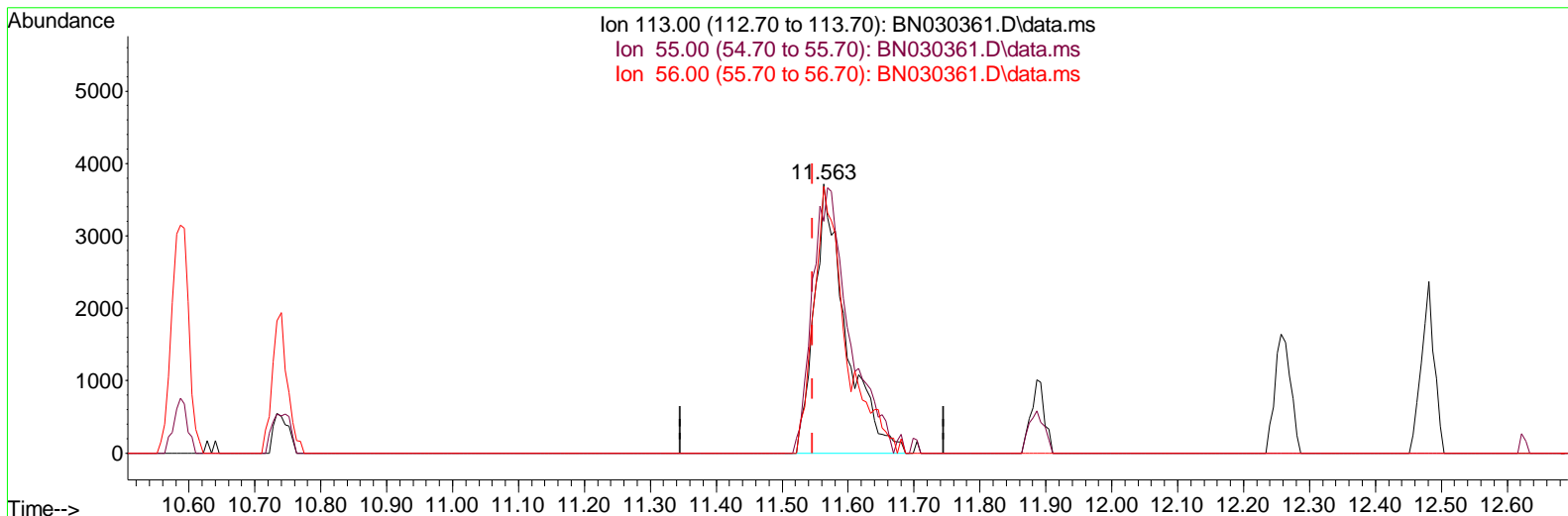
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN031124\
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Instrument :
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TIC: BN030361.D\data.ms

(34) Caprolactam

11.563min (+ 0.018) 5.45 ng/ul m

response 12429

Ion	Exp%	Act%
113.00	100.00	100.00
55.00	127.60	86.30#
56.00	113.50	99.11
0.00	0.00	0.00

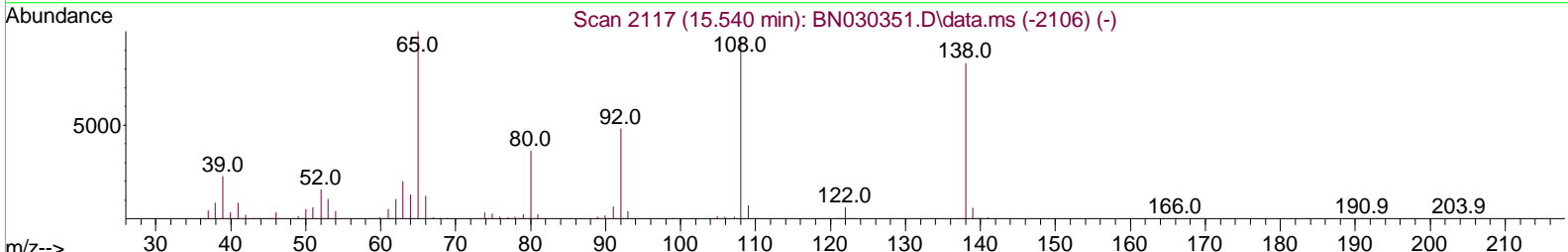
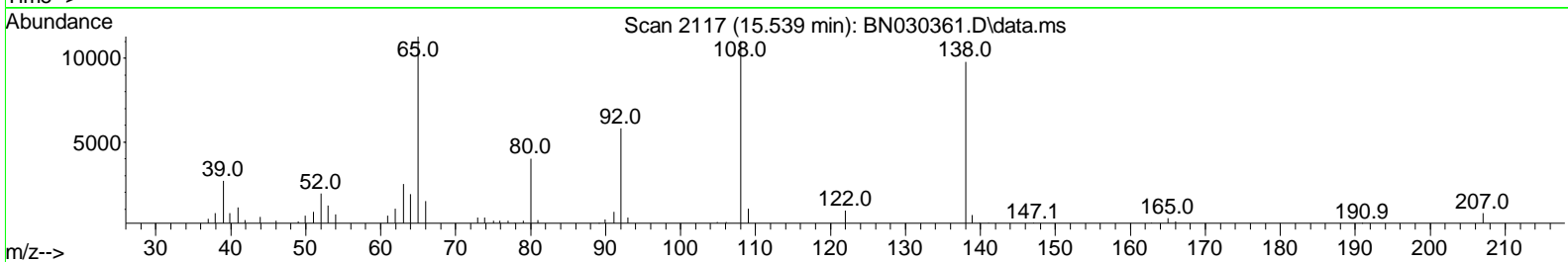
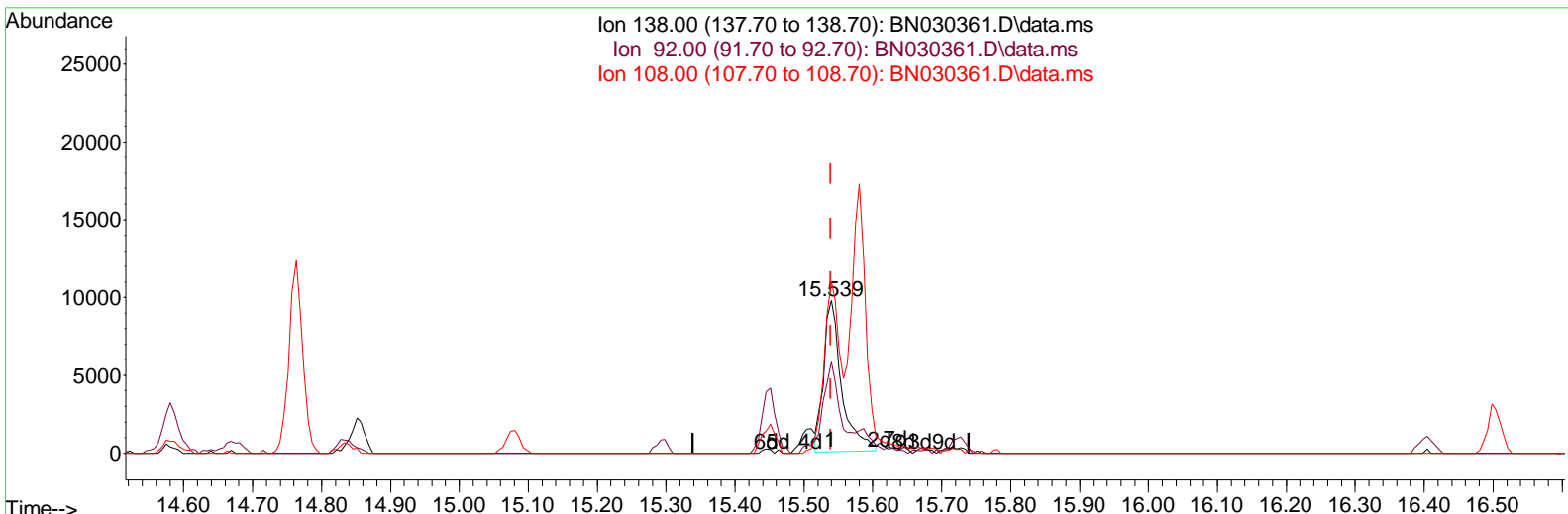
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN031124\
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Instrument :
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Reviewed By :Yogesh Patel 03/12/2024
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TIC: BN030361.D\data.ms

(63) 4-Nitroaniline

15.539min (-0.000) 4.09 ng/ul

response 17667

Ion	Exp%	Act%
138.00	100.00	100.00
92.00	57.80	59.60
108.00	116.50	112.78
0.00	0.00	0.00

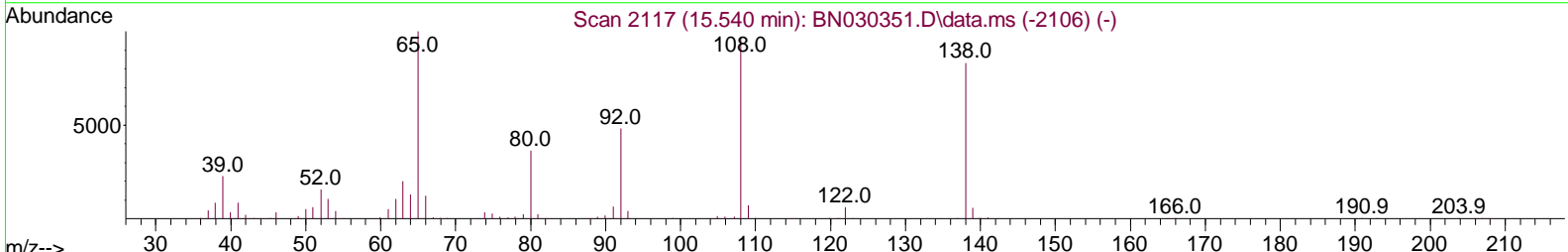
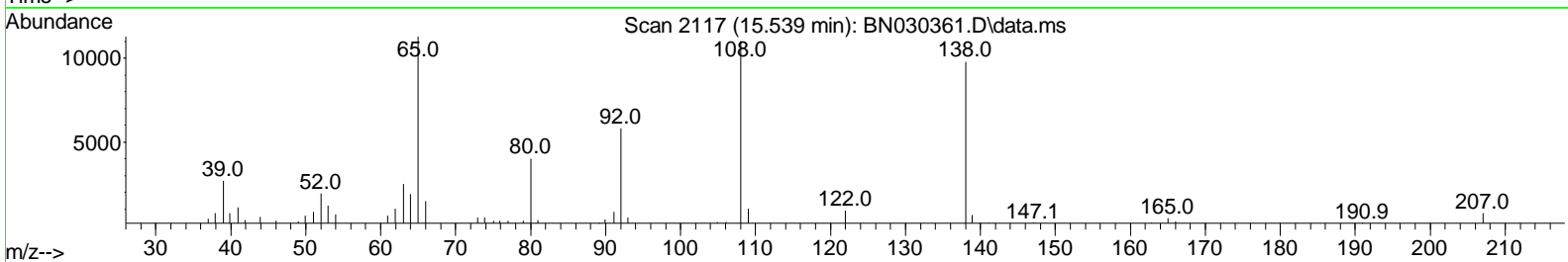
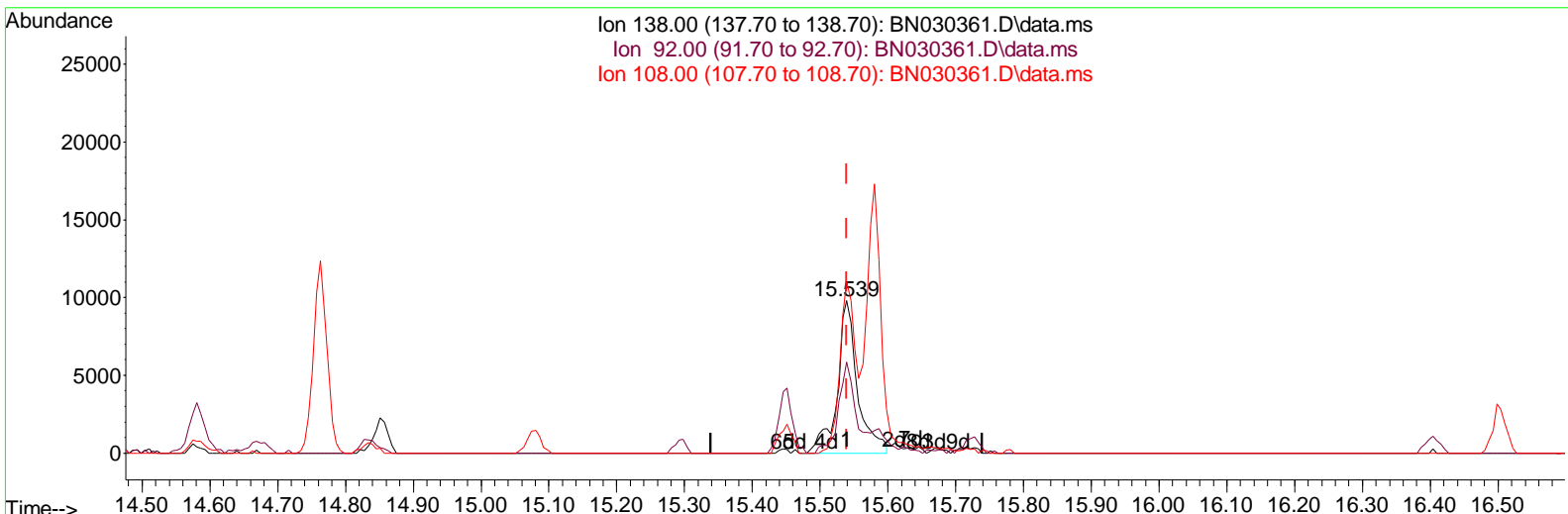
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN031124\
 Data File : BN030361.D
 Acq On : 11 Mar 2024 15:30
 Operator : MA/JU
 Sample : P1601-05
 Misc : SFAM-MDL-ME SOIL-01
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
ClientSampleId :
 MDL-MED-SOIL-QT1-2024-01

Manual Integrations APPROVED

Reviewed By : Yogesh Patel 03/12/2024
 Supervised By : mohammad ahmed 03/12/2024

Quant Time: Mar 11 16:00:23 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN030824.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Mar 08 13:57:34 2024
 Response via : Initial Calibration



TIC: BN030361.D\data.ms

(63) 4-Nitroaniline

15.539min (-0.000) 4.70 ng/ul m

response 20297

Ion	Exp%	Act%
138.00	100.00	100.00
92.00	57.80	59.60
108.00	116.50	112.78
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BNO31124\
 Data File : BNO30361.D
 Acq On : 11 Mar 2024 15:30
 Operator : MA/JU
 Sample : P1601-05
 Misc : SFAM-MDL-ME SOIL-01
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
ClientSampleId :
 MDL-MED-SOIL-QT1-2024-01

Manual Integrations APPROVED

Reviewed By :Yogesh Patel 03/12/2024
 Supervised By :mohammad ahmed 03/12/2024

Quant Time: Mar 11 16:03:05 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BNO30824.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Mar 08 13:57:34 2024
 Response via : Initial Calibration

Compound	R.T.	QI on	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.793	152	116038	20.000	ng/ul	0.00
20) Naphthalene-d8	10.587	136	488730	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.451	164	300907	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.210	188	668268	20.000	ng/ul	0.00
79) Chrysene-d12	21.433	240	598265	20.000	ng/ul	0.00
88) Perylene-d12	23.786	264	566724	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.269	96	15073	4.384	ng/uL	0.00
4) Pyridine-d5	3.675	84	185874	22.646	ng/ul	0.00
7) Phenol-d5	6.963	99	209174	20.317	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.140	67	117490	20.098	ng/ul	0.00
11) 2-Chlorophenol-d4	7.322	132	180553	21.893	ng/ul	0.00
15) 4-Methylphenol-d8	8.505	113	170845	21.054	ng/ul	0.00
21) Nitrobenzene-d5	8.963	128	88916	21.804	ng/ul	0.00
24) 2-Nitrophenol-d4	9.675	143	94309	22.058	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.210	165	164656	21.594	ng/ul	0.00
31) 4-Chloroaniline-d4	10.740	131	240125	21.218	ng/ul	0.00
46) Dimethylphthalate-d6	13.875	166	549228	22.638	ng/ul	0.00
49) Acenaphthylene-d8	14.140	160	606316	22.060	ng/ul	0.00
54) 4-Nitrophenol-d4	14.657	143	90879	21.127	ng/ul	0.00
60) Fluorene-d10	15.451	176	451966	22.709	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.581	200	83102	19.909	ng/ul	0.00
73) Anthracene-d10	17.310	188	651113	20.065	ng/ul	0.00
81) Pyrene-d10	19.622	212	810805	21.341	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.633	264	682869	23.005	ng/ul	0.00
Target Compounds						
2) 1,4-Dioxane	3.305	88	3634m	1.101	ng/uL	
5) Pyridine	3.699	79	39560	4.737	ng/ul #	80
6) Benzaldehyde	6.946	77	29658	6.181	ng/ul	95
8) Phenol	6.987	94	47338	4.621	ng/ul	99
10) Bis(2-Chloroethyl)ether	7.234	93	37528	4.584	ng/ul	99
12) 2-Chlorophenol	7.352	128	40501	4.773	ng/ul	97
13) 2-Methylphenol	8.240	108	34549	4.498	ng/ul	100
14) 2,2'-oxybis(1-Chloropr...	8.328	45	33647	4.520	ng/ul	95
16) Acetophenone	8.628	105	65167	4.997	ng/ul	98
17) N-Nitrosodimethylpropyl a...	8.610	70	32055	5.137	ng/ul	96
18) 4-Methylphenol	8.563	108	38908	4.794	ng/ul	99
19) Hexachloroethane	8.857	117	18555	4.707	ng/ul	97
22) Nitrobenzene	8.999	77	50082	4.759	ng/ul	98
23) Isophorone	9.528	82	85506	4.700	ng/ul	99
25) 2-Nitrophenol	9.710	139	22926	4.978	ng/ul	94
26) 2,4-Dimethylphenol	9.769	107	33114	3.483	ng/ul	100
27) Bis(2-Chloroethoxy)met...	10.022	93	49973	4.683	ng/ul	97
29) 2,4-Dichlorophenol	10.234	162	38146	4.981	ng/ul	91
30) Naphthalene	10.640	128	140889	5.050	ng/ul	98
32) 4-Chloroaniline	10.763	127	54318	4.935	ng/ul	98
33) Hexachlorobutadiene	10.910	225	26065	4.879	ng/ul	99
34) Caprolactam	11.563	113	12429m	5.450	ng/ul	
35) 4-Chloro-3-methylphenol	11.887	107	41994	4.846	ng/ul	89

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BNO31124\
 Data File : BNO30361.D
 Acq On : 11 Mar 2024 15:30
 Operator : MA/JU
 Sample : P1601-05
 Mi sc : SFAM-MDL-ME SOI L-01
 ALS Vi al : 6 Sample Multi plier: 1

Instrument :
 BNA_N
ClientSampleId :
 MDL-MED-SOIL-QT1-2024-01

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 03/12/2024
 Supervised By :mohammad ahmed 03/12/2024

Quant Time: Mar 11 16:03:05 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BNO30824.MA.M
 Quant Title : SVOA CALI BRATI ON
 QLast Update : Fri Mar 08 13:57:34 2024
 Response via : Ini tial Cal i brati on

Compound	R. T.	QI on	Response	Conc	Units	Dev(Mi n)
36) 2-Methyl naphthal ene	12.257	142	93895	5.168	ng/ul	98
37) 1-Methyl naphthal ene	12.481	142	97660	5.437	ng/ul	98
39) 1, 2, 4, 5-Tetrachl oroben. . .	12.622	216	45789	4.819	ng/ul	98
40) Hexachl orocycl opentadi ene	12.598	237	31774	4.464	ng/ul #	93
41) 2, 4, 6-Tri chl orophenol	12.875	196	28159	4.671	ng/ul	95
42) 2, 4, 5-Tri chl orophenol	12.945	196	30954	4.699	ng/ul	97
43) 1, 1' -Bi phenyl	13.281	154	125057	4.924	ng/ul	97
44) 2-Chl oronaphthal ene	13.322	162	97413	4.939	ng/ul	95
45) 2-Ni troani li ne	13.540	65	24094	4.237	ng/ul	97
47) Di methyl phthal ate	13.922	163	130321	5.285	ng/ul	97
48) 2, 6-Di ni trotol uene	14.045	165	22887	4.806	ng/ul	92
50) Acenaphthyl ene	14.169	152	159808	4.997	ng/ul	97
51) 3-Ni troani li ne	14.375	138	19500	4.179	ng/ul	92
52) Acenaphthene	14.516	153	106377	5.021	ng/ul	99
53) 2, 4-Di ni trophenol	14.581	184	15838	5.919	ng/ul	96
55) 4-Ni trophenol	14.675	109	25941	5.319	ng/ul	92
56) Di benzofuran	14.851	168	144336	5.108	ng/ul	99
57) 2, 4-Di ni trotol uene	14.834	165	34122	5.021	ng/ul	97
58) 2, 3, 4, 6-Tetrachl orophenol	15.081	232	25379	4.787	ng/ul	92
59) Di ethyl phthal ate	15.292	149	135661	5.148	ng/ul	98
61) Fl uorene	15.504	166	124089	5.319	ng/ul	91
62) 4-Chl orophenyl -phenyl e. . .	15.510	204	57802	5.175	ng/ul	94
63) 4-Ni troani li ne	15.539	138	20297m	4.698	ng/ul	
66) 4, 6-Di ni tro-2-methyl ph. . .	15.592	198	24261	5.387	ng/ul #	94
67) N-Ni trosodi phenyl ami ne	15.722	169	98162	4.656	ng/ul	99
68) 4-Bromophenyl -phenyl ether	16.404	248	33848	4.727	ng/ul	93
69) Hexachl orobenzene	16.504	284	39365	5.020	ng/ul	98
70) Atrazi ne	16.692	200	42366	5.838	ng/ul	95
71) Pentachl orophenol	16.857	266	28035	5.284	ng/ul	94
72) Phenanthrene	17.251	178	196921	5.053	ng/ul	99
74) Anthracene	17.345	178	182759	4.610	ng/ul	99
75) 1, 2, 3, 4-Tetrachl oroben. . .	13.240	216	44794	4.490	ng/uL	96
76) Pentachl orobenzene	14.763	250	46288	4.975	ng/uL#	89
77) Carbazol e	17.628	167	163478	4.643	ng/ul	98
78) Di -n-butyl phthal ate	18.204	149	226977	4.639	ng/ul	100
80) Fl uoranthene	19.286	202	217031	4.662	ng/ul #	96
82) Pyrene	19.651	202	231509	4.838	ng/ul	99
83) Butyl benzyl phthal ate	20.580	149	99897	4.358	ng/ul	99
84) 3, 3' -Di chl orobenzi di ne	21.363	252	59821	4.601	ng/ul	99
85) Benzo(a)anthracene	21.416	228	222836	4.964	ng/ul	98
86) Bi s(2-ethyl hexyl)phtha. . .	21.369	149	159687	4.602	ng/ul	98
87) Chrysene	21.469	228	210816	5.063	ng/ul	97
89) Di -n-octyl phthal ate	22.292	149	264572	4.987	ng/ul	100
90) Benzo(b)fl uoranthene	23.068	252	198531	5.313	ng/ul	96
91) Benzo(k)fl uoranthene	23.121	252	191036	4.950	ng/ul	95
93) Benzo(a)pyrene	23.680	252	176667	4.996	ng/ul	97
94) I ndeno(1, 2, 3-cd)pyrene	26.204	276	186384	4.432	ng/ul	93
95) Di benzo(a, h)anthracene	26.233	278	151575	4.421	ng/ul	96
96) Benzo(g, h, i)peryl ene	26.945	276	148556	4.361	ng/ul	99

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

Instrument :

BNA_N

ClientSampleId :

MDL-MED-SOIL-QT1-2024-01

Manual IntegrationsAPPROVED

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Supervised By :mohammad ahmed 03/12/2024

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