

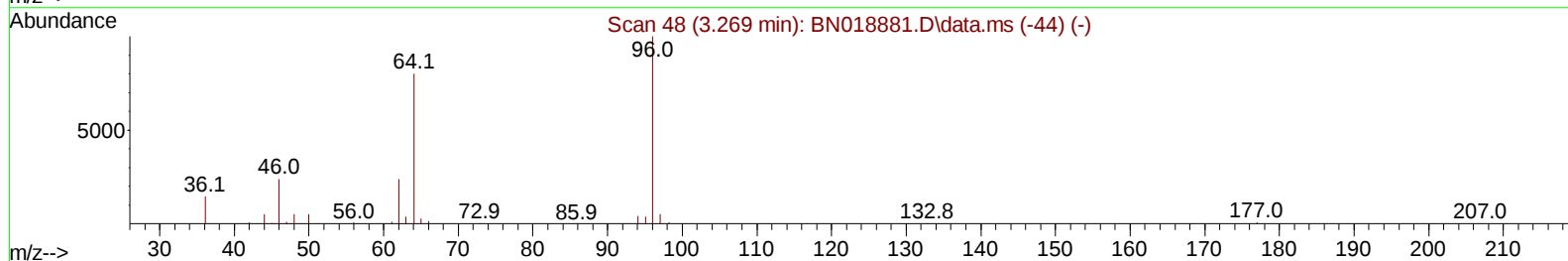
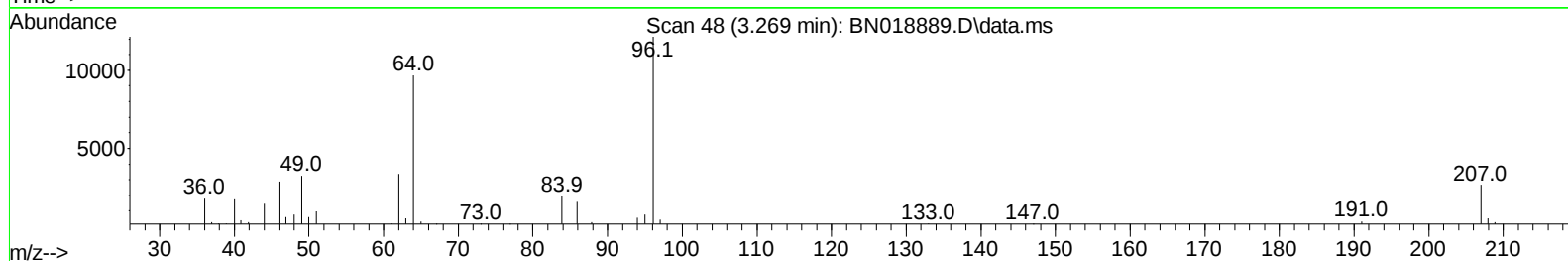
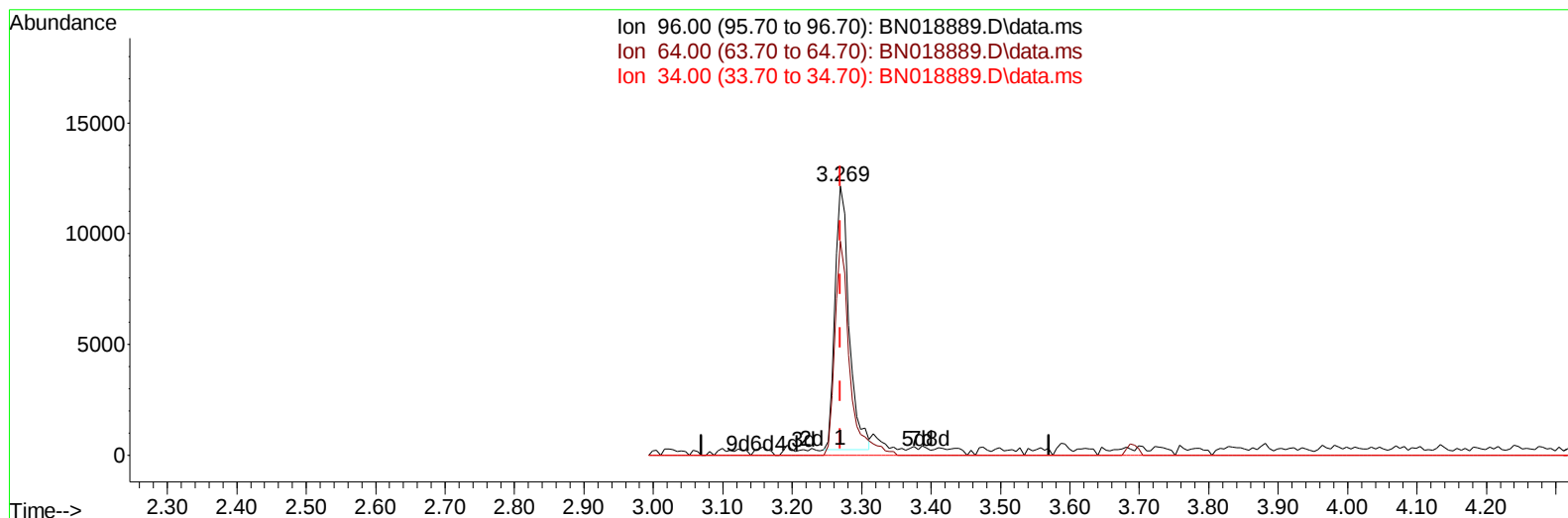
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN030922\
 Data File : BN018889.D
 Acq On : 09 Mar 2022 17:57
 Operator : CG/JU
 Sample : PB143227BL
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SBLK227

Manual IntegrationsAPPROVED

Quant Time: Mar 10 01:37:27 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN030822.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Mar 08 14:15:20 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 03/10/2022
 Supervised By :Yogesh Patel 03/11/2022



TIC: BN018889.D\data.ms

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

3.269min (-0.001) 5.47 ng/uL

response 16869

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	80.20	79.67
34.00	0.00	0.00
0.00	0.00	0.00

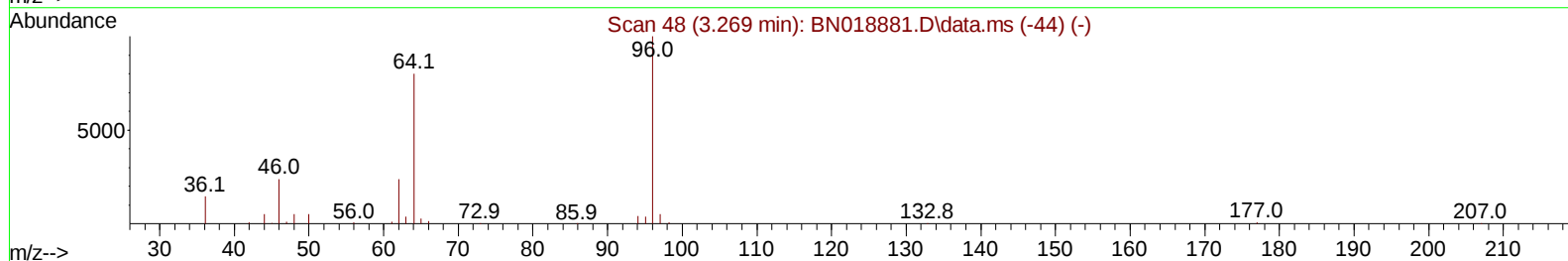
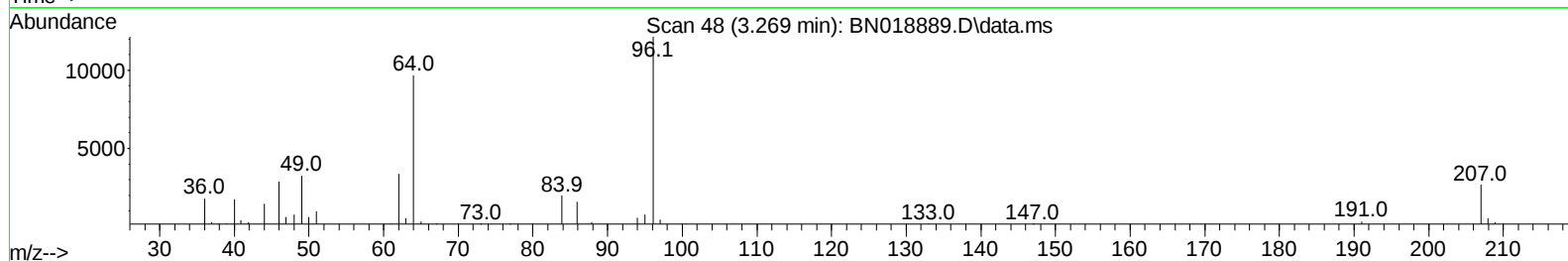
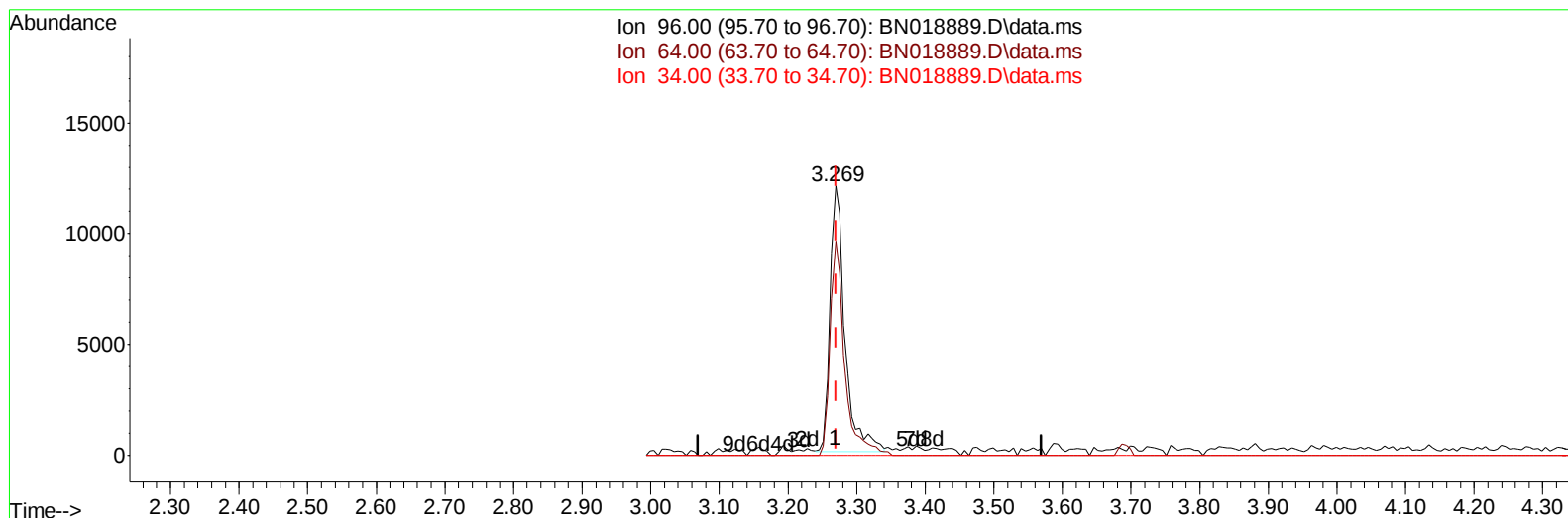
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TIC: BN018889.D\data.ms

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

3.269min (-0.001) 5.86 ng/uL m

response 18088

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	80.20	79.67
34.00	0.00	0.00
0.00	0.00	0.00

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.904	152	123412	20.000	ng/ul	0.00
20) Naphthalene-d8	10.710	136	576831	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.539	164	410030	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.275	188	919723	20.000	ng/ul	0.00
79) Chrysene-d12	21.451	240	1031835	20.000	ng/ul	0.00
88) Perylene-d12	23.845	264	1110844	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.269	96	18088m	5.862	ng/uL	0.00
4) Pyridine-d5	3.687	84	265667	28.138	ng/ul	0.00
7) Phenol-d5	7.057	99	332886	27.628	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.222	67	218530	30.112	ng/ul	0.00
11) 2-Chlorophenol-d4	7.434	132	258642	29.954	ng/ul	0.00
15) 4-Methylphenol-d8	8.604	113	278466	28.423	ng/ul	0.00
21) Nitrobenzene-d5	9.063	128	131180	29.606	ng/ul	0.00
24) 2-Nitrophenol-d4	9.787	143	149663	30.743	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.328	165	266585	28.629	ng/ul	0.00
31) 4-Chloroaniline-d4	10.840	131	420398	29.982	ng/ul	0.00
46) Dimethylphthalate-d6	13.939	166	981827	31.626	ng/ul	0.00
49) Acenaphthylene-d8	14.234	160	1197227	30.982	ng/ul	0.00
54) 4-Nitrophenol-d4	14.716	143	157297	25.019	ng/ul	0.00
60) Fluorene-d10	15.528	176	846484	31.369	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.639	200	138240	24.253	ng/ul	0.00
73) Anthracene-d10	17.375	188	1390167	31.601	ng/ul	0.00
81) Pyrene-d10	19.669	212	1700336	30.451	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.692	264	1922506	33.207	ng/ul	0.00

Target Compounds Qvalue

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

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