

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061722\
 Data File : BM035508.D
 Acq On : 17 Jun 2022 23:48
 Operator : CG/JU
 Sample : N3226-16
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 EW4N6

Manual IntegrationsAPPROVED

Quant Time: Jun 18 04:04:21 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM061622.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Jun 17 01:58:49 2022
 Response via : Initial Calibration

Reviewed By :Jagrut

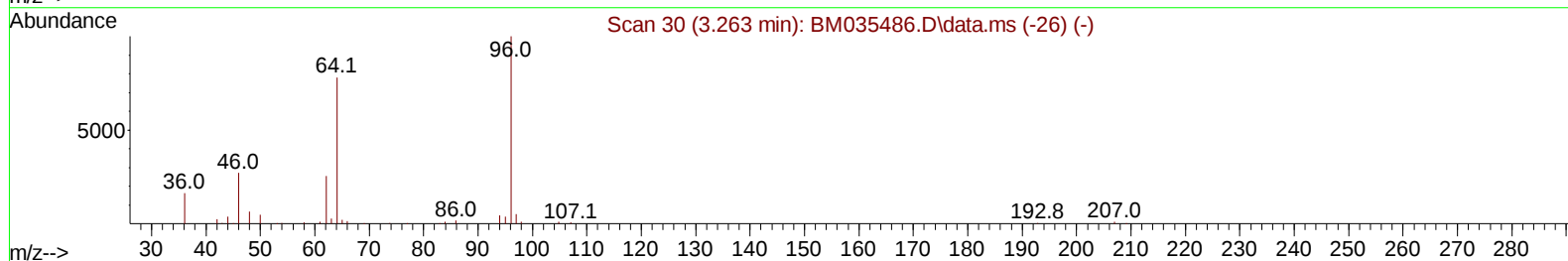
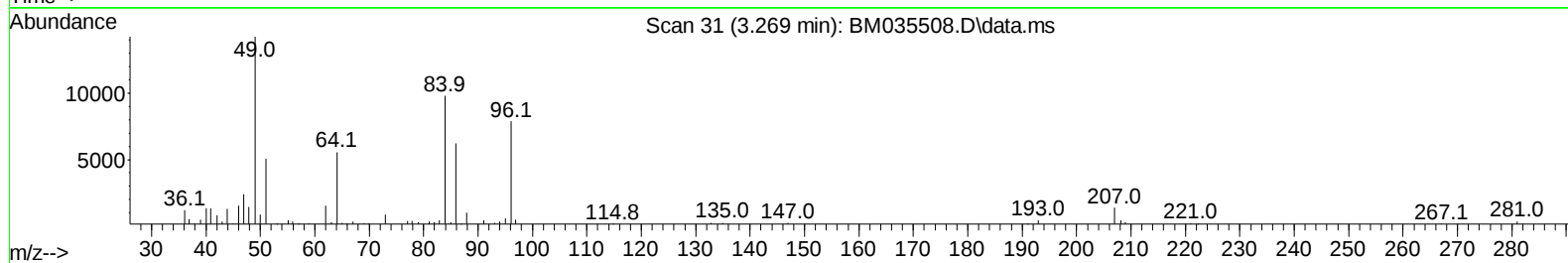
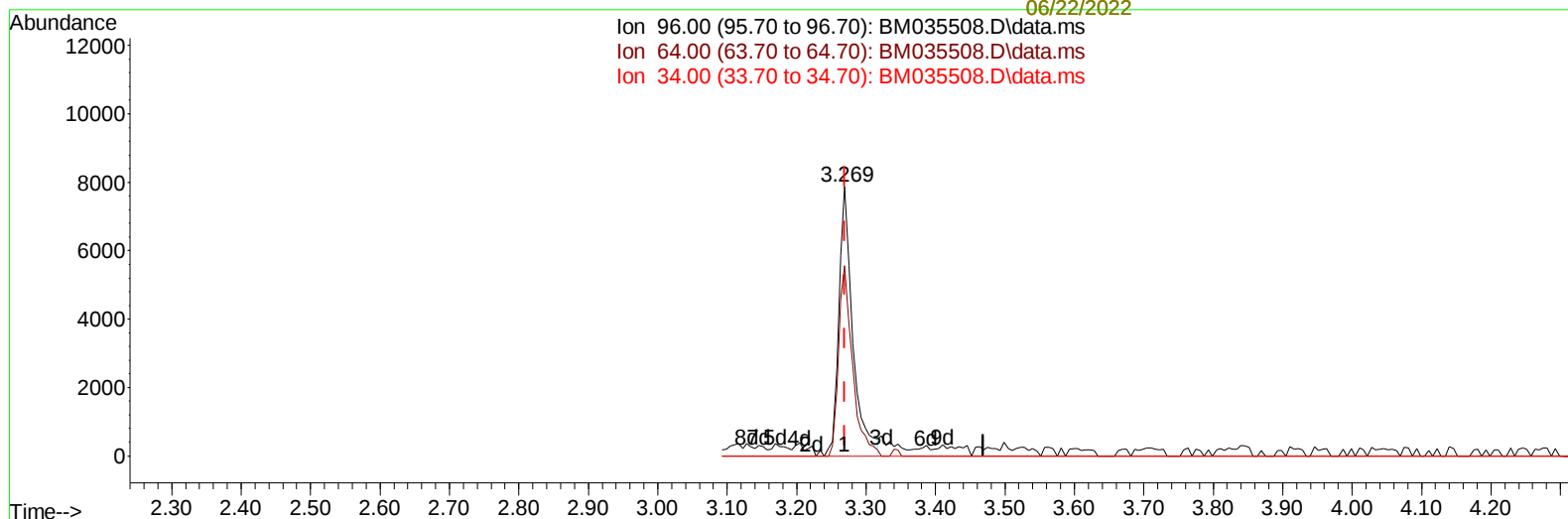
Upadhyay

06/18/2022

Supervised By :Yogesh

Patel

06/22/2022



TIC: BM035508.D\data.ms

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

3.269min (+ 0.000) 2.24 ng/uL

response 11008

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	75.40	70.52
34.00	0.00	0.00
0.00	0.00	0.00

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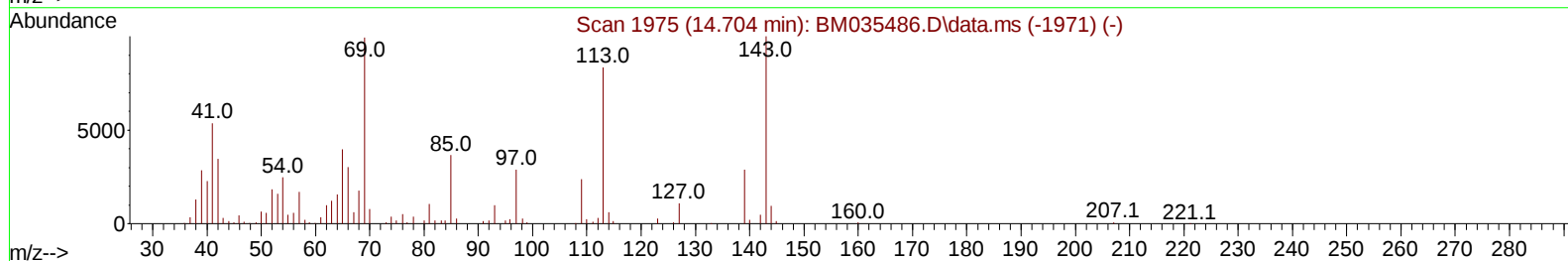
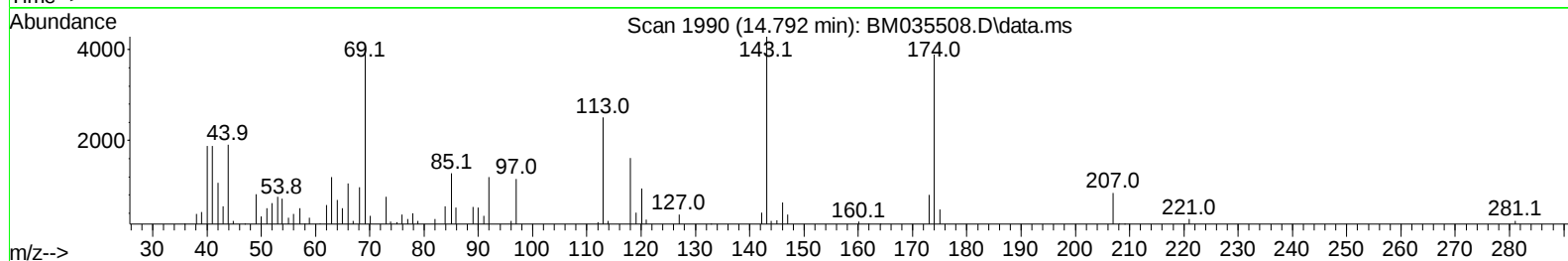
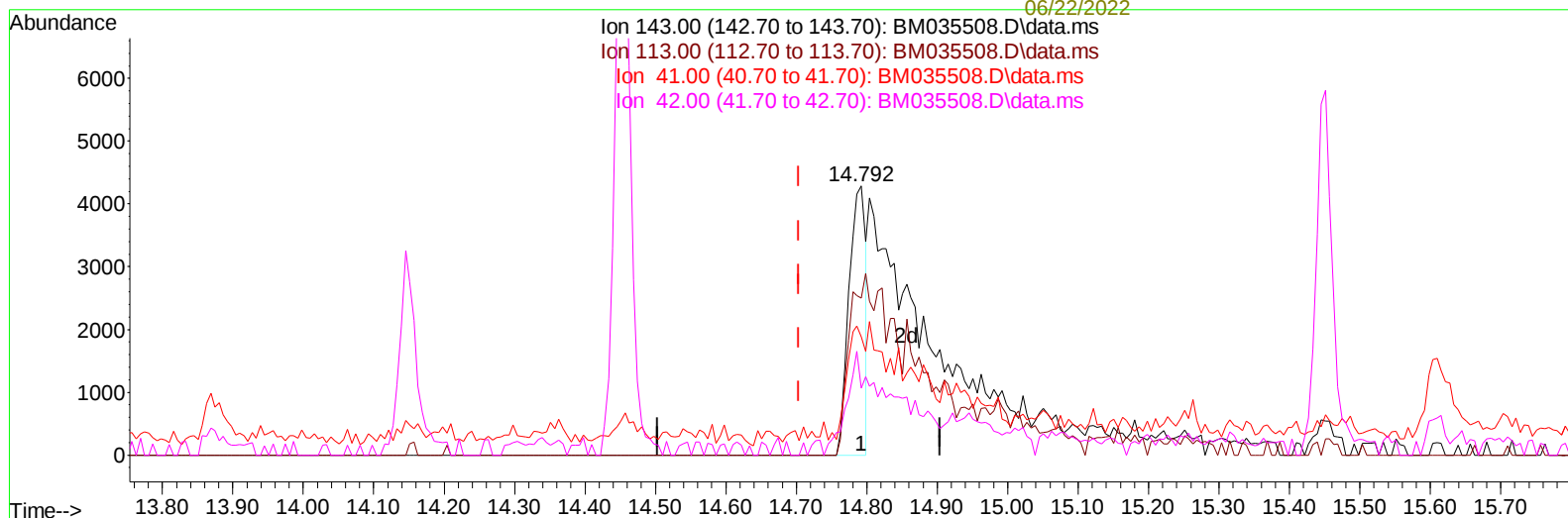
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(54) 4-Nitrophenol-d4 (S)

14.792min (+ 0.088) 1.36 ng/ul

response 6942

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	61.40	58.49
41.00	41.00	43.96
42.00	29.30	24.94

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.810	152	196415	20.000	ng/ul	0.00
20) Naphthalene-d8	10.610	136	822703	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.457	164	489542	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.198	188	999083	20.000	ng/ul	0.00
79) Chrysene-d12	21.386	240	1036700	20.000	ng/ul	0.00
88) Perylene-d12	23.721	264	1027609	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.269	96	11572m	2.352	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	7.010	99	137341	9.124	ng/ul	0.01
9) Bis-(2-Chloroethyl)eth...	7.151	67	100492	10.735	ng/ul	0.00
11) 2-Chlorophenol-d4	7.357	132	142516	10.931	ng/ul	0.00
15) 4-Methylphenol-d8	8.551	113	102260	8.115	ng/ul	0.01
21) Nitrobenzene-d5	8.981	128	63853	10.427	ng/ul	0.00
24) 2-Nitrophenol-d4	9.710	143	67690	10.251	ng/ul	0.01
28) 2,4-Dichlorophenol-d3	10.275	165	100329	8.948	ng/ul	0.02
31) 4-Chloroaniline-d4	10.781	131	141519	8.022	ng/ul	0.02
46) Dimethylphthalate-d6	13.874	166	400900	12.039	ng/ul	0.00
49) Acenaphthylene-d8	14.151	160	466739	11.311	ng/ul	0.00
54) 4-Nitrophenol-d4	14.792	143	28574m	5.586	ng/ul	0.09
60) Fluorene-d10	15.451	176	360498	12.503	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.610	200	29280	4.916	ng/ul	0.02
73) Anthracene-d10	17.304	188	561513	12.504	ng/ul	0.00
81) Pyrene-d10	19.598	212	691154	13.561	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.574	264	723277	14.593	ng/ul	0.00

Target Compounds Qvalue

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

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