

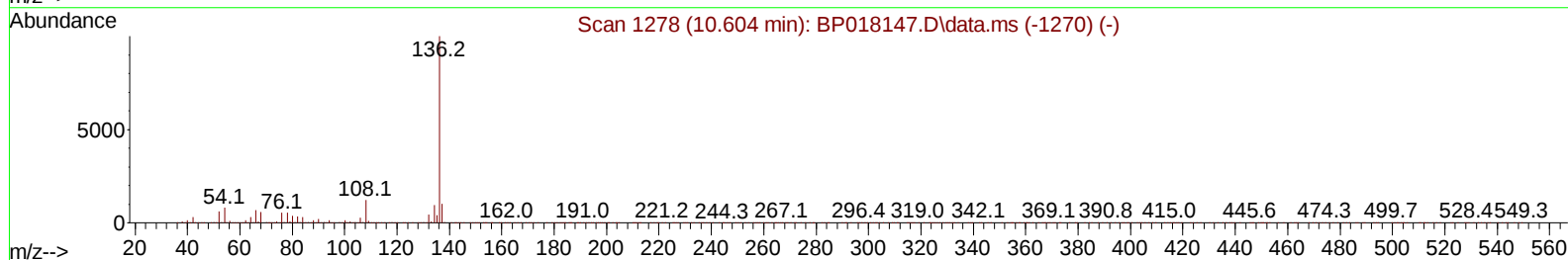
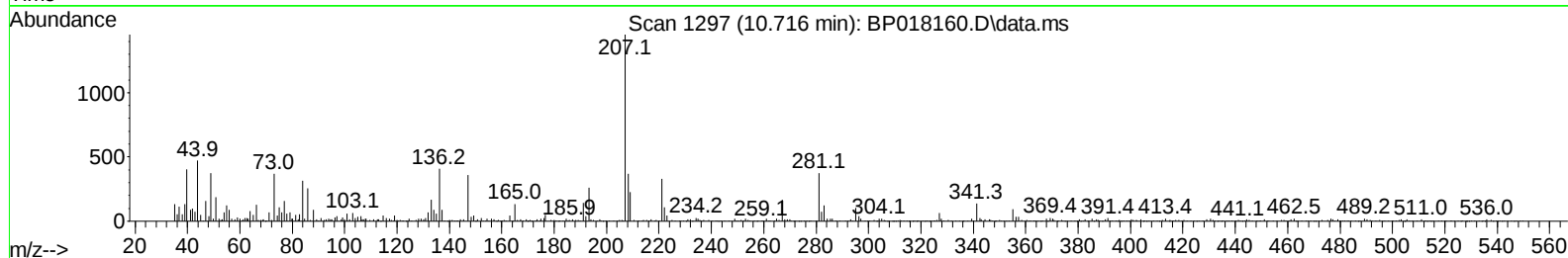
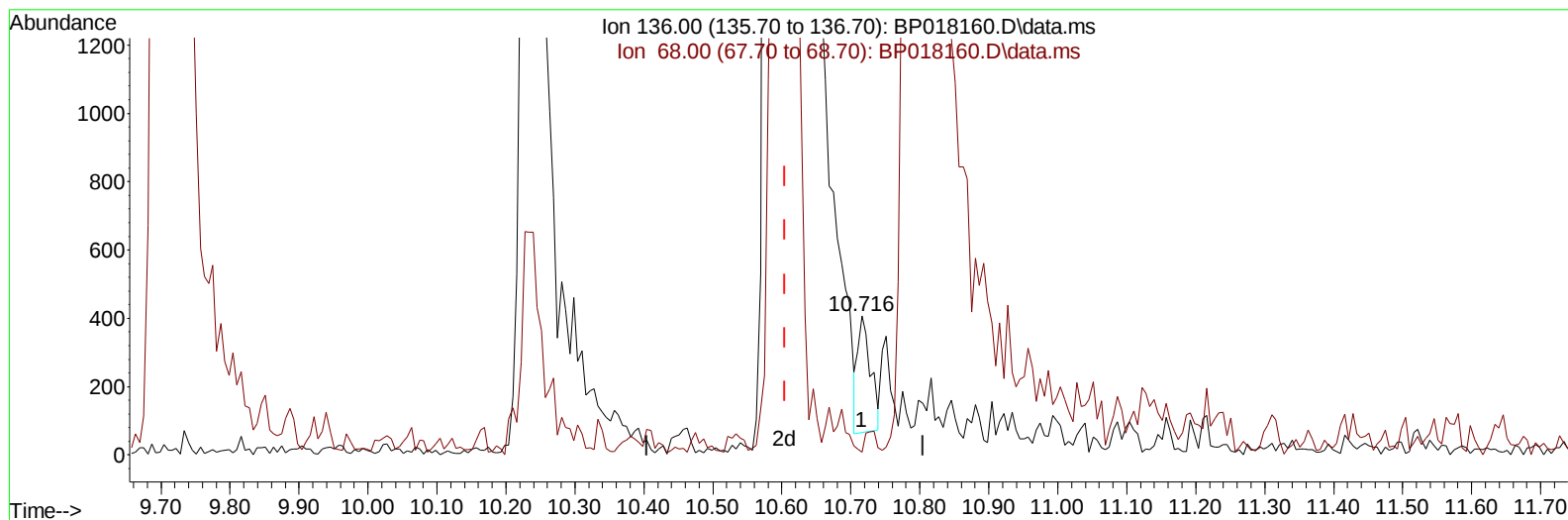
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111423\
 Data File : BP018160.D
 Acq On : 14 Nov 2023 17:51
 Operator : MA/JU
 Sample : 05337-06
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 GCDJ1

Manual IntegrationsAPPROVED

Quant Time: Nov 15 00:39:31 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Nov 13 15:50:18 2023
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 11/16/2023
 Supervised By :mohammad ahmed 11/16/2023



TIC: BP018160.D\data.ms

(20) Naphthalene-d8 (I)

10.716min (+ 0.111) 20.00 ng/ul

response 447

Ion	Exp%	Act%
136.00	100.00	100.00
68.00	4.60	5.15
0.00	0.00	0.00
0.00	0.00	0.00

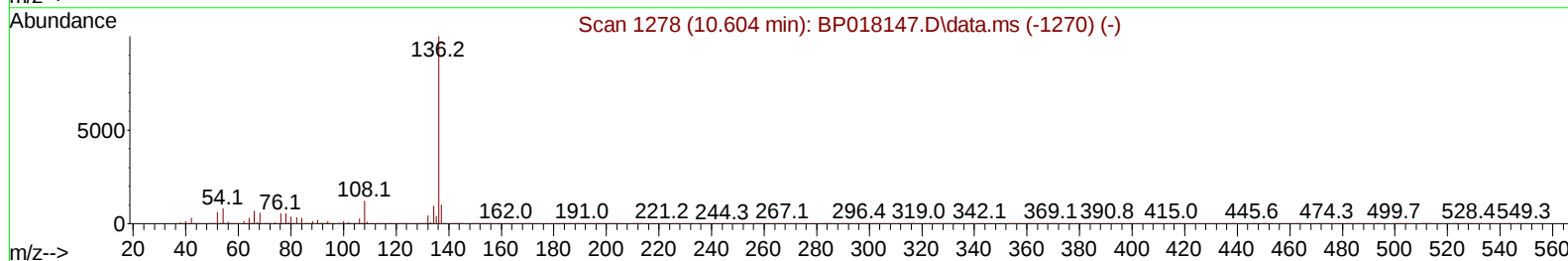
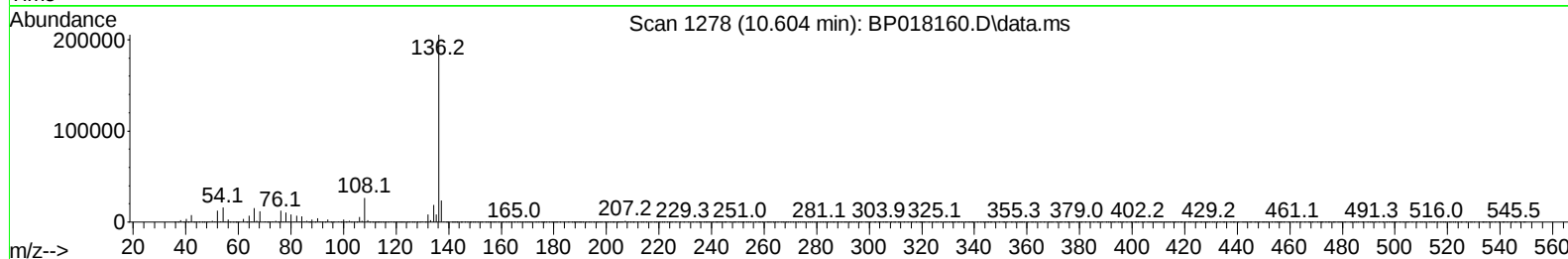
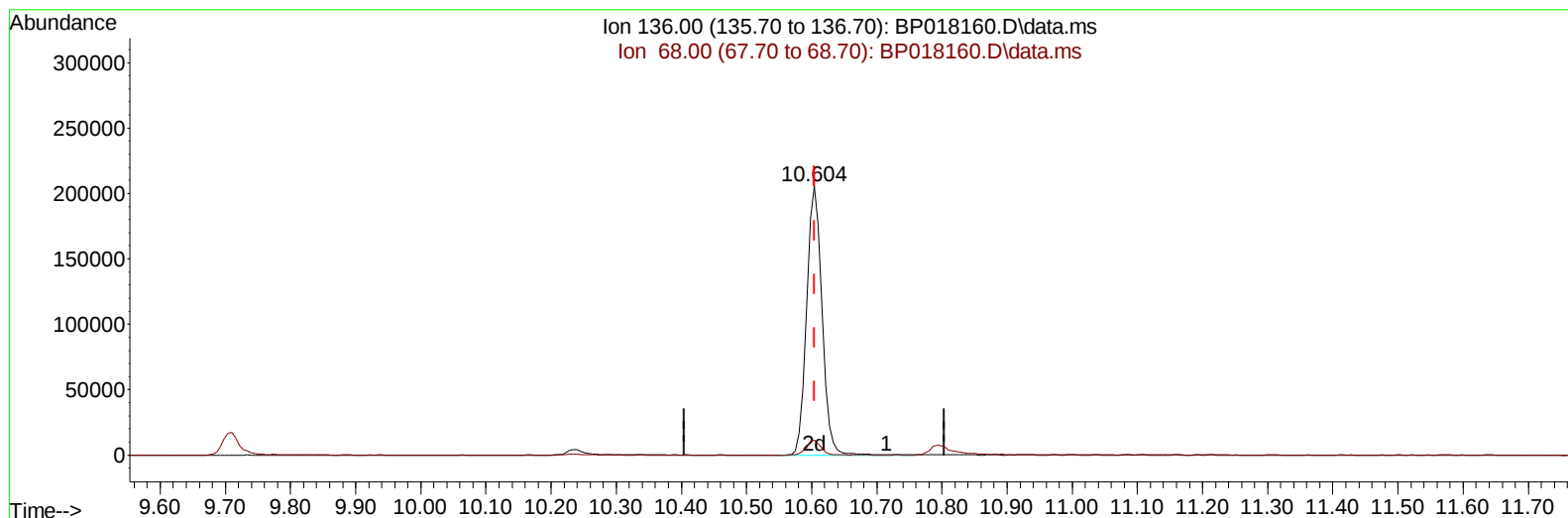
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(20) Naphthalene-d8 (I)

10.604min (-0.000) 20.00 ng/ul m

response 339886

Ion	Exp%	Act%
136.00	100.00	100.00
68.00	4.60	5.62#
0.00	0.00	0.00
0.00	0.00	0.00

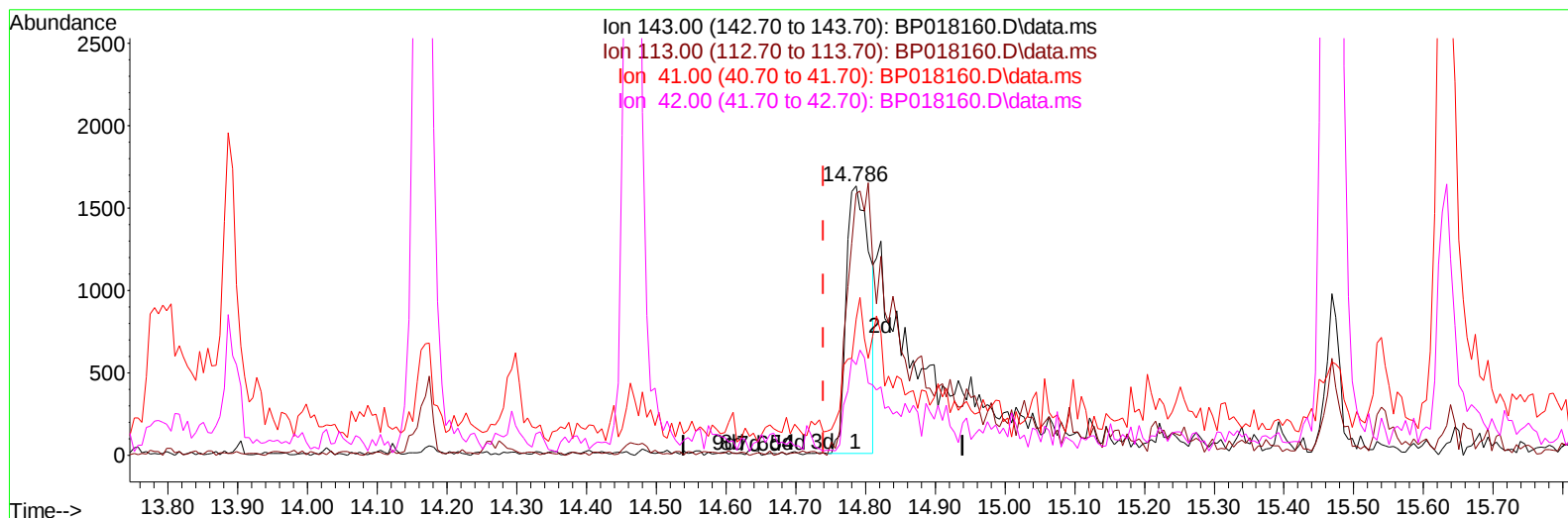
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Sample : 05337-06
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDJ1

Manual IntegrationsAPPROVED

Quant Time: Nov 15 01:25:37 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Nov 13 15:50:18 2023
Response via : Initial Calibration

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Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111423\
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 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 11/16/2023
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Quant Time: Nov 15 01:26:24 2023
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.787	152	84124	20.000	ng/ul	0.00
20) Naphthalene-d8	10.604	136	339886m	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.469	164	214000	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.251	188	538464	20.000	ng/ul	0.00
79) Chrysene-d12	21.380	240	510284	20.000	ng/ul	#-0.01
88) Perylene-d12	23.850	264	543339	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.163	96	6885	2.887	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	6.963	99	39564	4.871	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.140	67	145426	30.470	ng/ul	0.00
11) 2-Chlorophenol-d4	7.310	132	130606	20.307	ng/ul	0.00
15) 4-Methylphenol-d8	8.516	113	82345	12.373	ng/ul	0.00
21) Nitrobenzene-d5	8.992	128	83680	27.327	ng/ul	0.00
24) 2-Nitrophenol-d4	9.704	143	91460	26.364	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.234	165	143951	23.281	ng/ul	0.00
31) 4-Chloroaniline-d4	10.792	131	114328	13.483	ng/ul	0.00
46) Dimethylphthalate-d6	13.892	166	624047	31.605	ng/ul	0.00
49) Acenaphthylene-d8	14.169	160	620662	28.860	ng/ul	0.00
54) 4-Nitrophenol-d4	14.786	143	7679m	2.603	ng/ul	0.05
60) Fluorene-d10	15.469	176	530236	32.526	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.633	200	78500	20.166	ng/ul	-0.03
73) Anthracene-d10	17.351	188	997870	34.265	ng/ul	0.00
81) Pyrene-d10	19.604	212	1225520	35.955	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.686	264	1137970	36.019	ng/ul	-0.02

Target Compounds Qvalue

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

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