

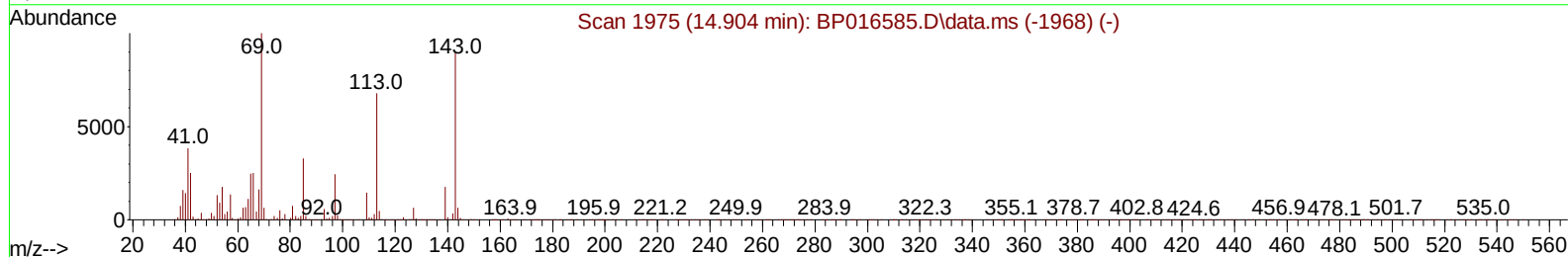
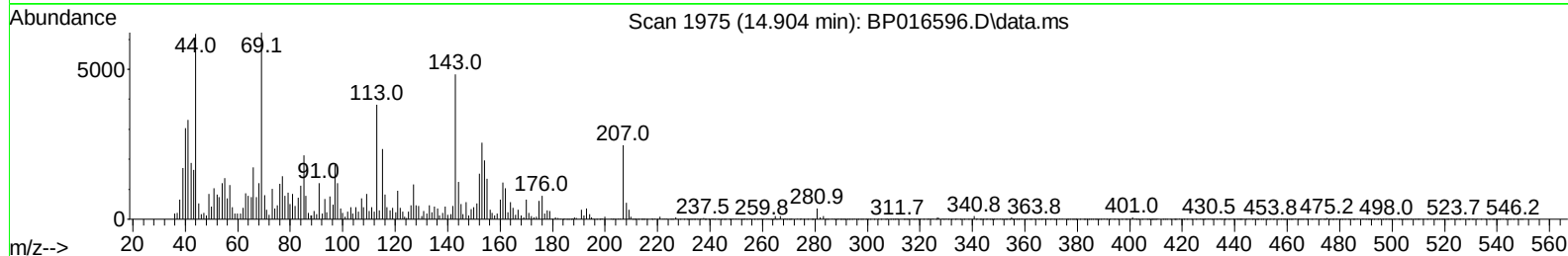
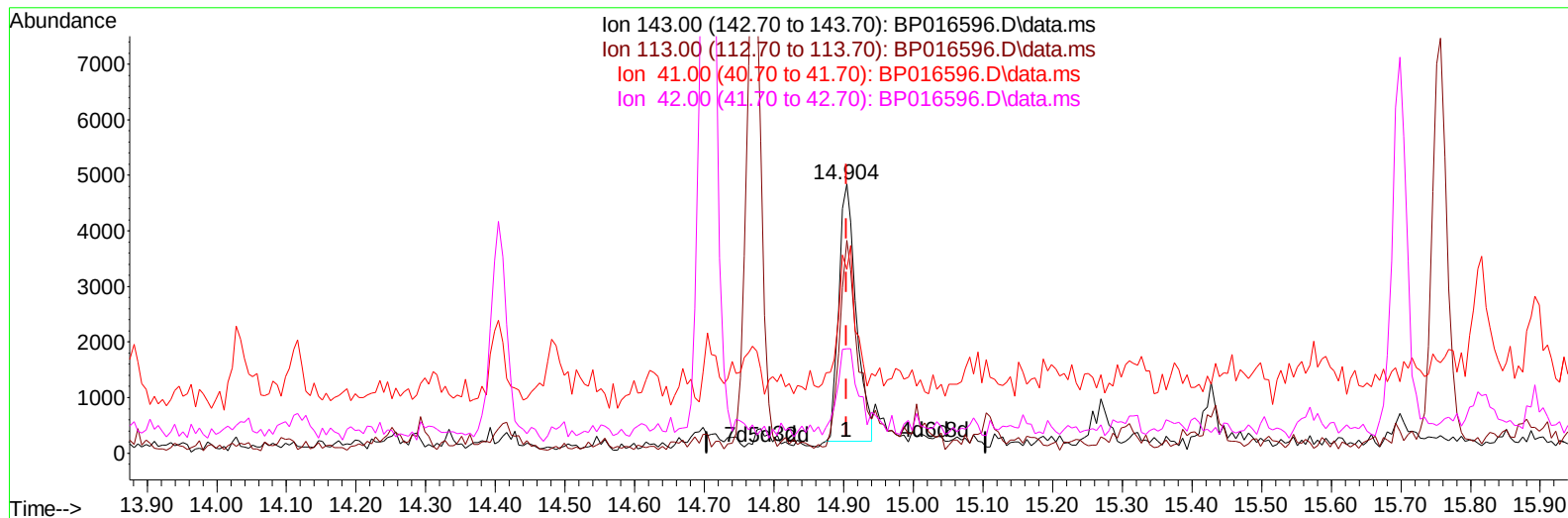
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081523\
Data File : BP016596.D
Acq On : 15 Aug 2023 16:45
Operator : MA/JU
Sample : 03962-01DL 5X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AF4DL

Manual IntegrationsAPPROVED

Quant Time: Aug 15 18:18:34 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP081423.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Aug 14 18:26:05 2023
Response via : Initial Calibration

Reviewed By :Yogesh Patel 08/16/2023
Supervised By :Sohil Jodhani 08/16/2023



TIC: BP016596.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.904min (+ 0.000) 0.91 ng/ul

response 8038

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	77.80	79.03
41.00	46.00	68.21#
42.00	30.30	38.90#

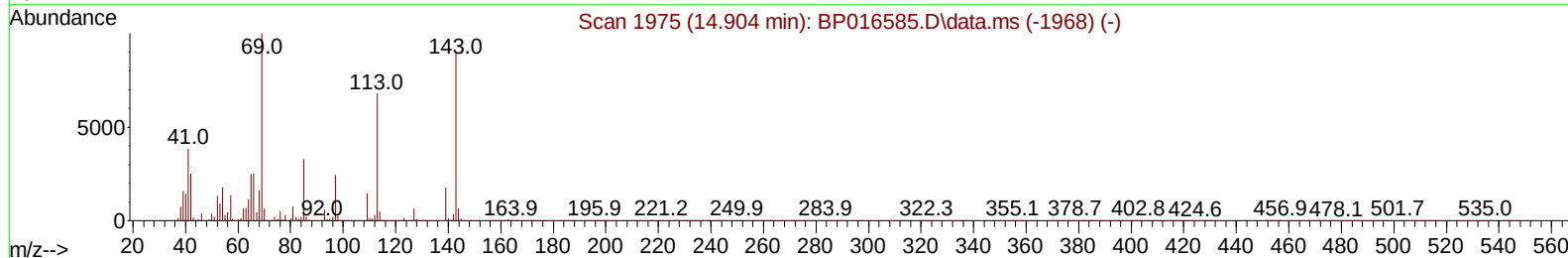
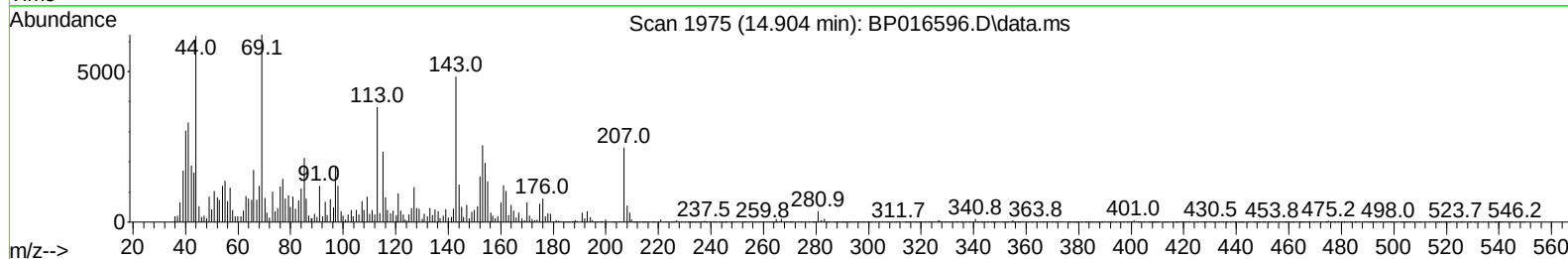
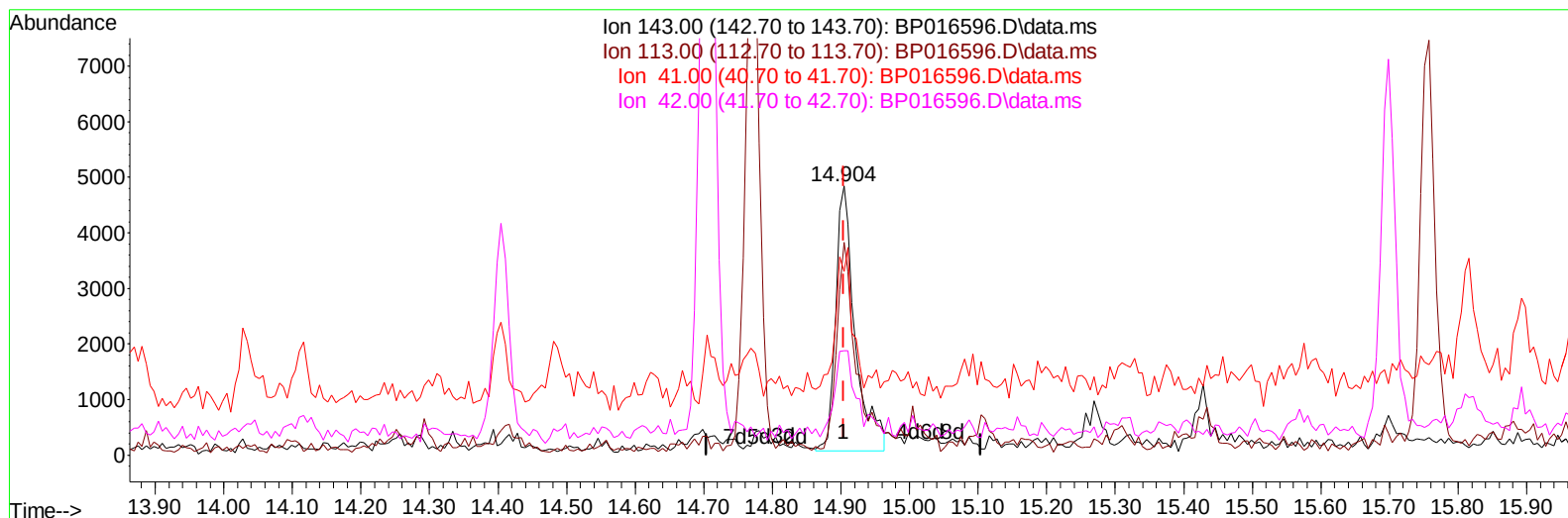
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TIC: BP016596.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.904min (+ 0.000) 1.07 ng/ul m

response 9427

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	77.80	79.03
41.00	46.00	68.21#
42.00	30.30	38.90#

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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 08/16/2023
 Supervised By :Sohil Jodhani 08/16/2023

Quant Time: Aug 16 02:35:33 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP081423.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Aug 14 18:26:05 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.058	152	205290	20.000	ng/ul	0.00
20) Naphthalene-d8	10.887	136	956713	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.704	164	600338	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.475	188	1397678	20.000	ng/ul	0.00
79) Chrysene-d12	21.569	240	1306916	20.000	ng/ul	0.00
88) Perylene-d12	24.157	264	1318904	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.393	96	2745	0.508	ng/uL	0.00
4) Pyridine-d5	3.840	84	44989	2.737	ng/ul	0.00
7) Phenol-d5	7.193	99	30355	1.604	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.393	67	92744	7.440	ng/ul	0.00
11) 2-Chlorophenol-d4	7.575	132	76727	5.644	ng/ul	0.00
15) 4-Methylphenol-d8	8.758	113	55407	3.682	ng/ul	0.00
21) Nitrobenzene-d5	9.252	128	50093	7.441	ng/ul	0.00
24) 2-Nitrophenol-d4	9.969	143	41484	6.488	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.493	165	85847	6.823	ng/ul	0.00
31) 4-Chloroaniline-d4	11.046	131	131940	6.112	ng/ul	0.00
46) Dimethylphthalate-d6	14.116	166	367378	8.461	ng/ul	0.00
49) Acenaphthylene-d8	14.404	160	404458	8.252	ng/ul	0.00
54) 4-Nitrophenol-d4	14.904	143	9427m	1.066	ng/ul	0.00
60) Fluorene-d10	15.698	176	333129	8.999	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.816	200	25137	3.780	ng/ul	0.00
73) Anthracene-d10	17.569	188	541719	8.960	ng/ul	0.00
81) Pyrene-d10	19.804	212	640248	9.349	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.980	264	555269	8.785	ng/ul	0.00
Target Compounds						
					Qvalue	
52) Acenaphthene	14.769	153	1719771	44.510	ng/ul	97
61) Fluorene	15.757	166	485133	11.223	ng/ul	99
74) Anthracene	17.610	178	90482	1.223	ng/ul	99
77) Carbazole	17.887	167	82946	1.217	ng/ul	97
80) Fluoranthene	19.475	202	353469	4.325	ng/ul	100
82) Pyrene	19.834	202	188226	2.181	ng/ul	97

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

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