

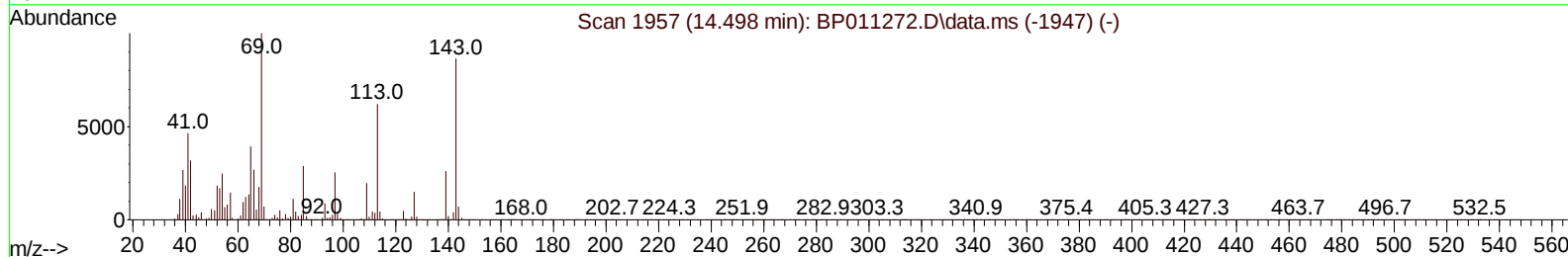
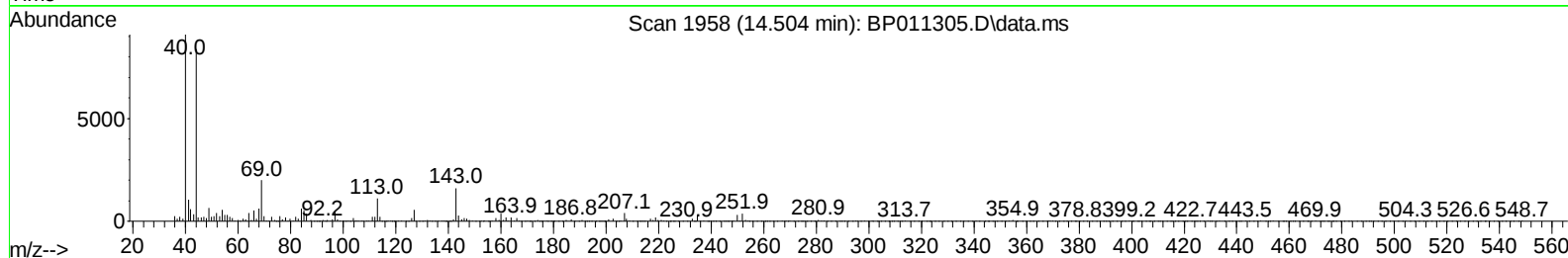
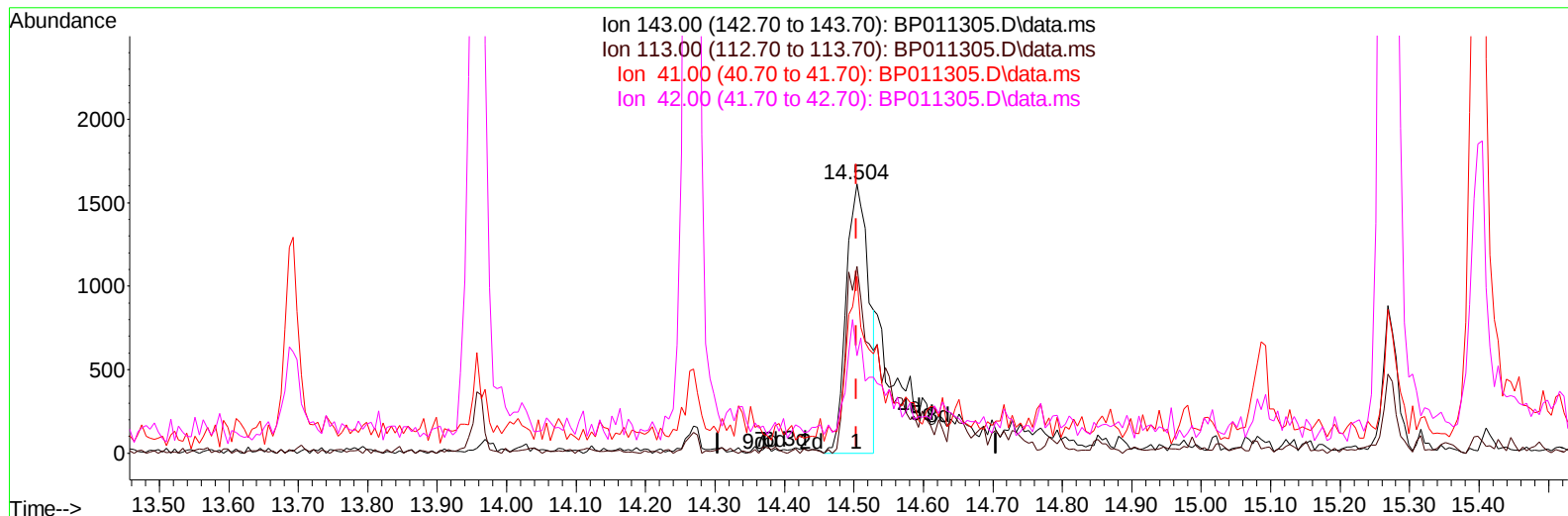
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080522\
Data File : BP011305.D
Acq On : 06 Aug 2022 07:49
Operator : CG/JU
Sample : N4026-21
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GBJJ6

Manual IntegrationsAPPROVED

Quant Time: Aug 08 01:10:46 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080222.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Aug 02 23:52:37 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 08/08/2022
Supervised By :Sohil Jodhani 08/10/2022



TIC: BP011305.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.504min (+ 0.000) 1.26 ng/u1

response 3627

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	69.40
41.00	54.40	65.61#
42.00	35.50	36.19

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.593	152	94650	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.387	136	384777	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.269	164	246271	20.000	ng/ul	# 0.00
64) Phenanthrene-d10	17.039	188	496261	20.000	ng/ul	0.00
79) Chrysene-d12	21.163	240	650655	20.000	ng/ul	-0.01
88) Perylene-d12	23.468	264	630669	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.087	96	2042	1.744	ng/uL	0.00
4) Pyridine-d5	3.511	84	9194	2.912	ng/ul	0.00
7) Phenol-d5	6.775	99	17197	2.180	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	6.940	67	137862	24.222	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.128	132	74354	11.565	ng/ul	-0.01
15) 4-Methylphenol-d8	8.316	113	37009	5.702	ng/ul	-0.01
21) Nitrobenzene-d5	8.758	128	112903	40.794	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.475	143	63948	24.650	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.016	165	82625	16.702	ng/ul	0.00
31) 4-Chloroaniline-d4	10.540	131	105045	13.725	ng/ul	0.00
46) Dimethylphthalate-d6	13.693	166	364712	21.623	ng/ul	0.00
49) Acenaphthylene-d8	13.957	160	595545	28.063	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.504	143	4448m	1.551	ng/ul	0.00
60) Fluorene-d10	15.269	176	539285	37.575	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.404	200	52168	22.639	ng/ul	-0.01
73) Anthracene-d10	17.134	188	1034328	47.424	ng/ul	-0.01
81) Pyrene-d10	19.398	212	1129293	34.328	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.321	264	1040990	33.124	ng/ul	-0.02

Target Compounds Qvalue

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

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