

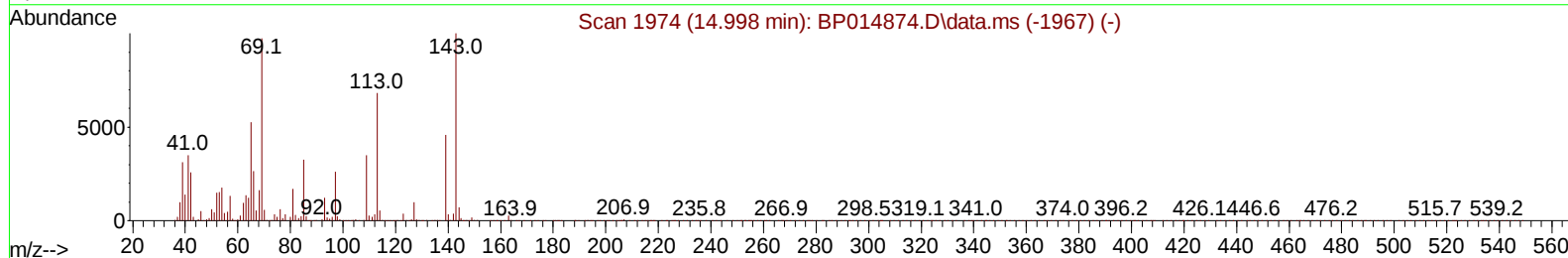
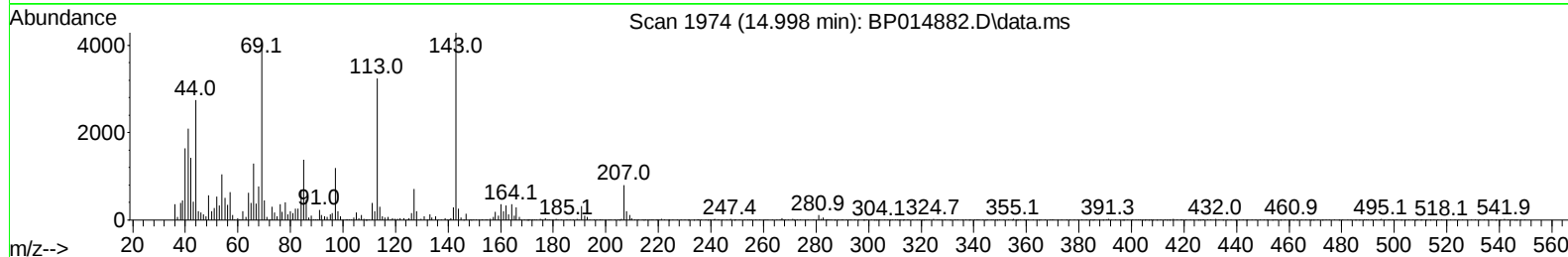
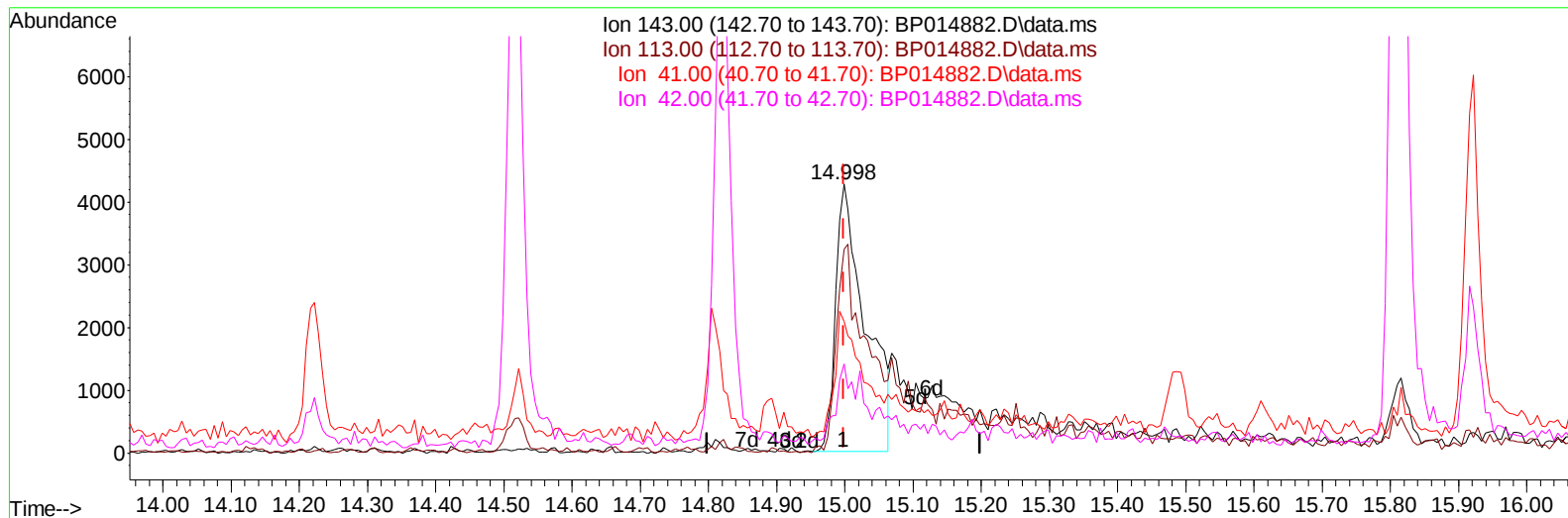
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042823\
Data File : BP014882.D
Acq On : 28 Apr 2023 19:22
Operator : CG/JU
Sample : 02488-16
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHE71

Manual IntegrationsAPPROVED

Quant Time: Apr 29 03:18:07 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042823.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat Apr 29 00:40:33 2023
Response via : Initial Calibration

Reviewed By :Christian Giraldo 05/01/2023
Supervised By :Jagrut Upadhyay 05/01/2023



TIC: BP014882.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.998min (-0.000) 2.46 ng/u1

response 12883

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	68.20	75.73
41.00	36.10	48.89#
42.00	26.80	33.12#

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Quant Time: Apr 29 03:28:08 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042823.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Apr 29 00:40:33 2023
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.175	152	190688	20.000	ng/ul	0.00
20) Naphthalene-d8	11.010	136	796582	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.822	164	491929	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.575	188	1037026	20.000	ng/ul	0.00
79) Chrysene-d12	21.657	240	943151	20.000	ng/ul	0.00
88) Perylene-d12	24.262	264	1083157	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.469	96	25079	4.740	ng/uL	0.00
4) Pyridine-d5	3.922	84	45389	3.569	ng/ul	0.02
7) Phenol-d5	7.316	99	97627	6.541	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.493	67	308881	32.998	ng/ul	0.00
11) 2-Chlorophenol-d4	7.699	132	296461	26.235	ng/ul	0.00
15) 4-Methylphenol-d8	8.881	113	174528	15.347	ng/ul	0.00
21) Nitrobenzene-d5	9.351	128	174537	30.516	ng/ul	0.00
24) 2-Nitrophenol-d4	10.081	143	161980	30.851	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.622	165	294146	26.085	ng/ul	0.00
31) 4-Chloroaniline-d4	11.145	131	277884	17.045	ng/ul	0.00
46) Dimethylphthalate-d6	14.222	166	1238535	35.131	ng/ul	0.00
49) Acenaphthylene-d8	14.516	160	1445457	34.324	ng/ul	0.00
54) 4-Nitrophenol-d4	14.998	143	15539m	2.962	ng/ul	0.00
60) Fluorene-d10	15.816	176	1164891	37.906	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.922	200	106293	24.232	ng/ul	0.00
73) Anthracene-d10	17.675	188	1988122	41.978	ng/ul	0.00
81) Pyrene-d10	19.892	212	2340748	44.685	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.098	264	2347127	43.872	ng/ul	0.00

Target Compounds Qvalue

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

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