

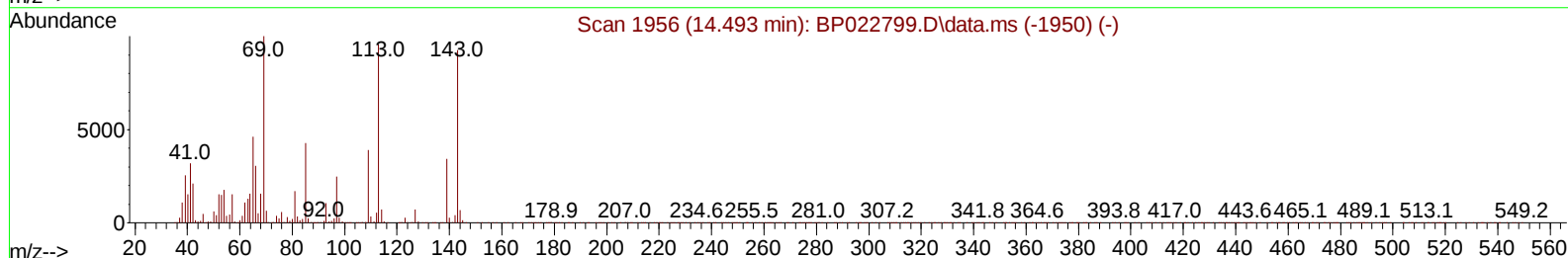
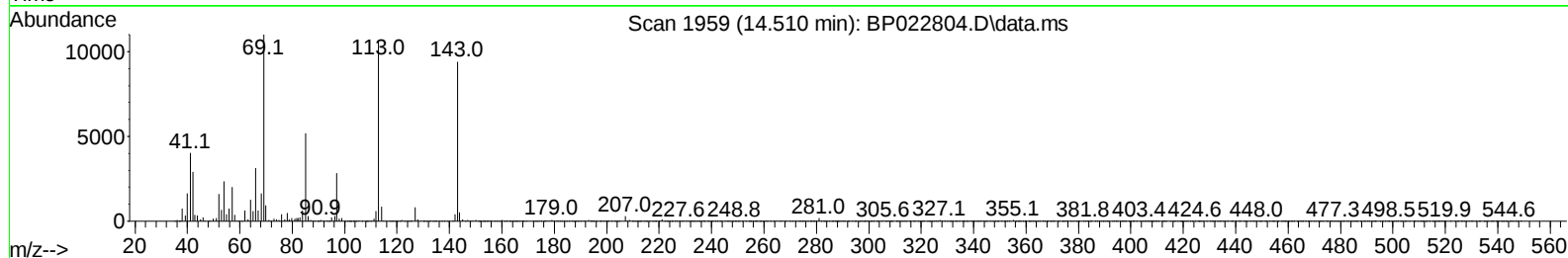
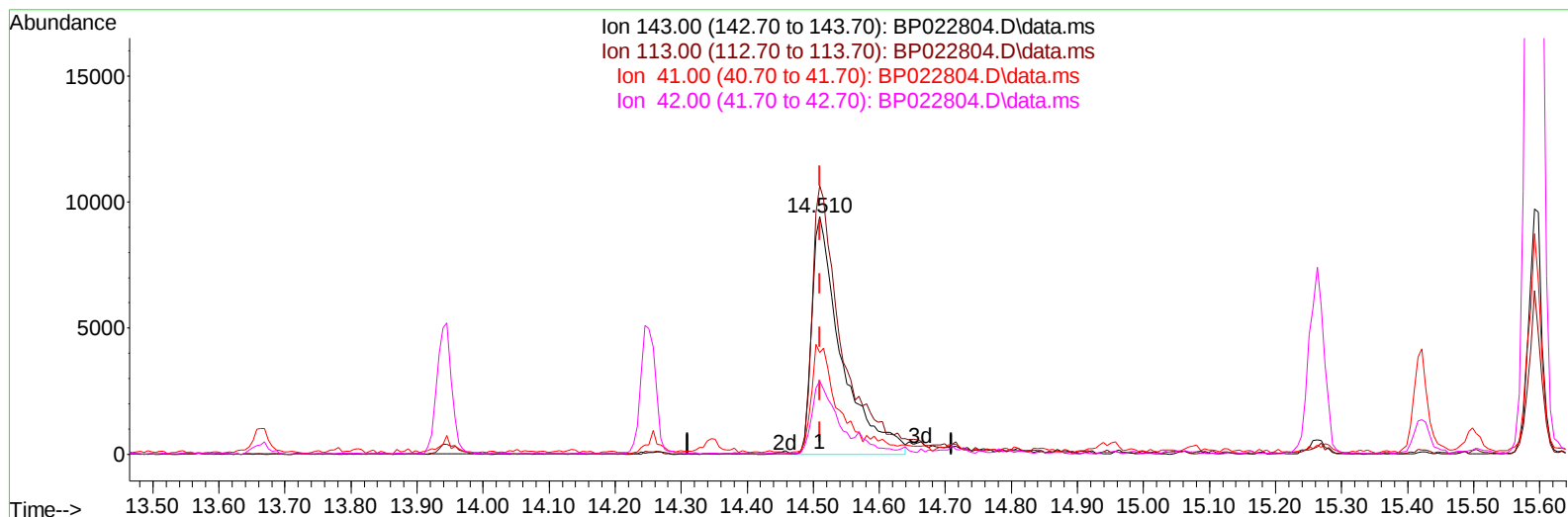
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP103024\
 Data File : BP022804.D
 Acq On : 01 Nov 2024 14:45
 Operator : RC/JU
 Sample : P4583-01
 Misc :
 ALS Vial : 40 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0PI5

Manual IntegrationsAPPROVED

Quant Time: Nov 01 16:16:42 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 30 10:56:39 2024
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 11/04/2024
 Supervised By :mohammad ahmed 11/05/2024



TIC: BP022804.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.510min (-0.000) 9.38 ng/ul m

response 29188

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	97.60	113.08
41.00	35.90	42.83
42.00	23.70	30.85#

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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 11/04/2024
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Quant Time: Nov 01 16:17:29 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.628	152	80588	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.387	136	302340	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.251	164	233495	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.069	188	578815	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.516	240	720845	20.000	ng/ul	# 0.02
88) Perylene-d12	24.768	264	842781	20.000	ng/ul	0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.175	96	7573	3.652	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	6.828	99	74725	11.015	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.969	67	108516	27.217	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.169	132	136009	28.256	ng/ul	-0.01
15) 4-Methylphenol-d8	8.340	113	121058	22.361	ng/ul	0.00
21) Nitrobenzene-d5	8.769	128	68798	29.594	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.487	143	83523	31.033	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.028	165	164030	28.675	ng/ul	0.00
31) 4-Chloroaniline-d4	10.534	131	29341	4.341	ng/ul	0.00
46) Dimethylphthalate-d6	13.663	166	547235	29.498	ng/ul	0.00
49) Acenaphthylene-d8	13.939	160	577850	29.326	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.510	143	29188m	9.378	ng/ul	0.00
60) Fluorene-d10	15.263	176	489807	30.440	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.422	200	92794	24.376	ng/ul	0.02
73) Anthracene-d10	17.169	188	916838	33.672	ng/ul	0.00
81) Pyrene-d10	19.551	212	1227220	32.194	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.551	264	1512407	33.602	ng/ul	0.03

Target Compounds Qvalue

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

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