

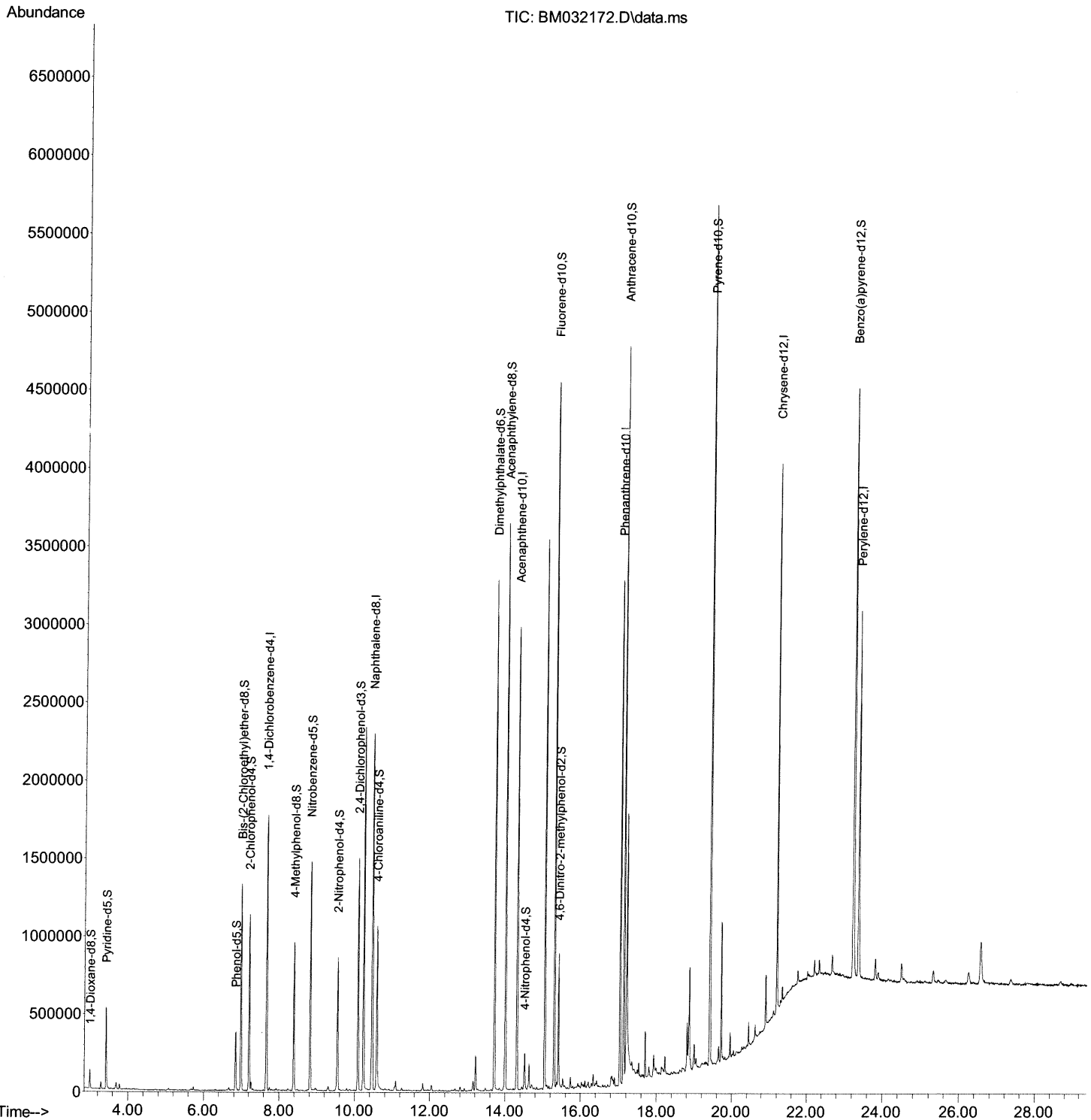
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092521\
Data File : BM032172.D
Acq On : 26 Sep 2021 02:40
Operator : CG/JU
Sample : M3739-12
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0B71

Manual Integrations
APPROVED

mohammad
9/27/2021 3:06:37 PM

Quant Time: Sep 26 03:18:01 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM092421.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat Sep 25 01:15:33 2021
Response via : Initial Calibration



Quantitation Report (Qedit)

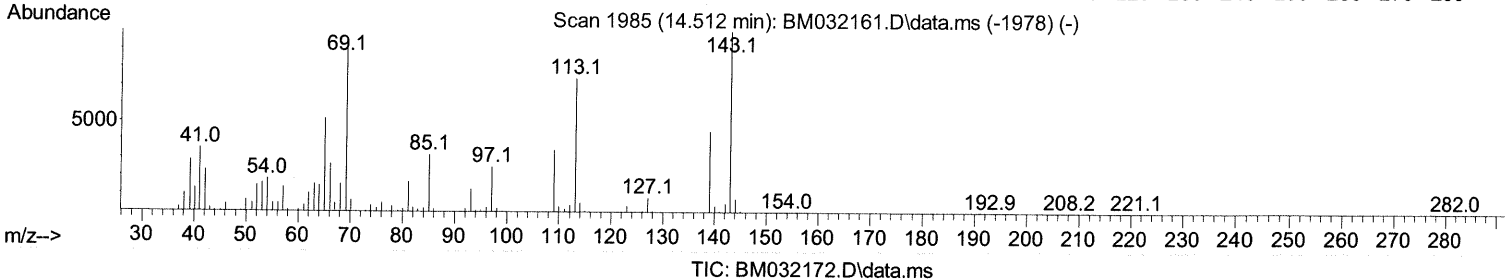
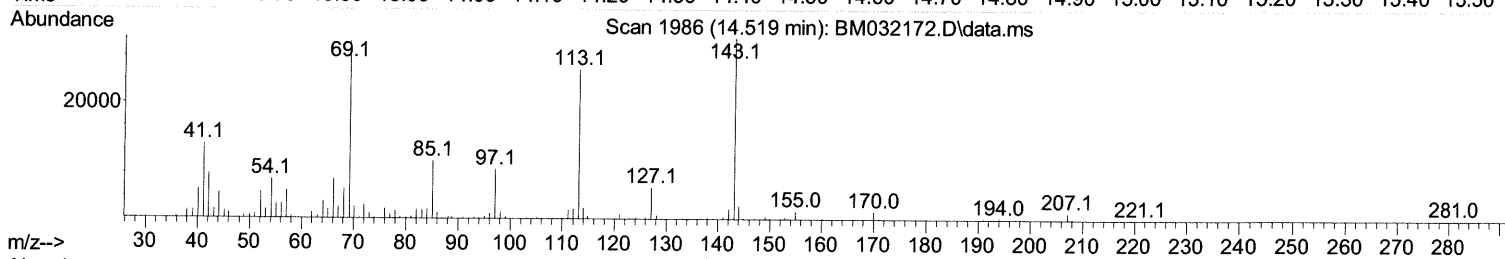
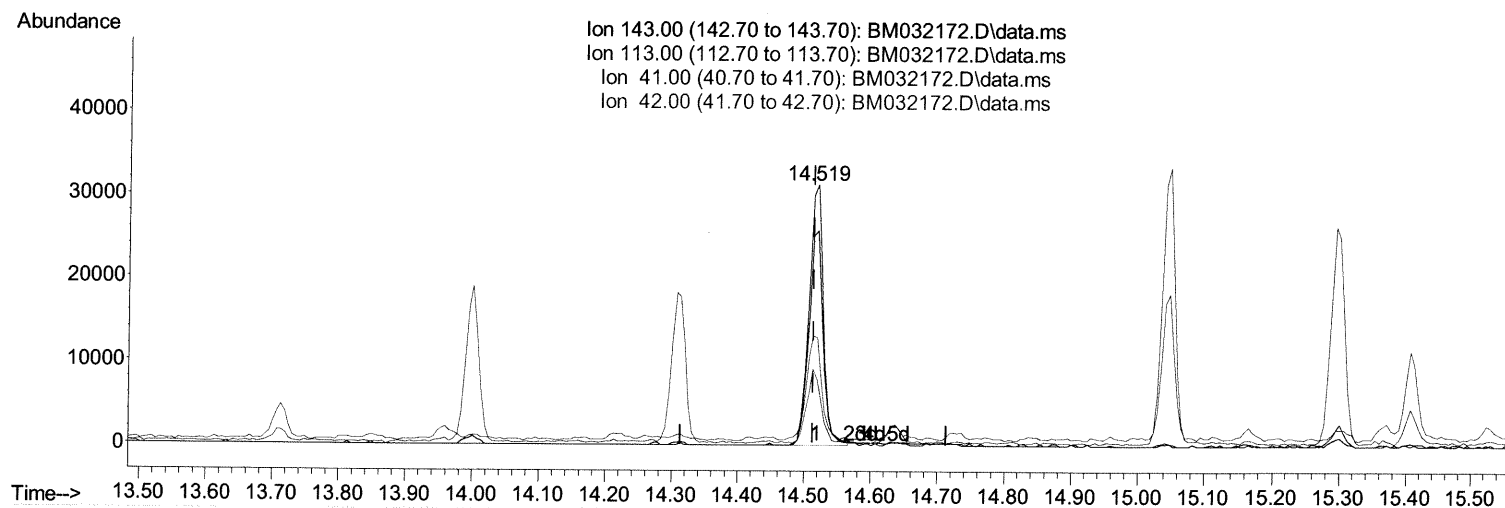
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(54) 4-Nitrophenol-d4 (S)

14.519min (+ 0.006) 4.50 ng/ul

response 46723

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.30	82.46
41.00	38.30	41.38
42.00	26.20	25.11

Quantitation Report (Qedit)

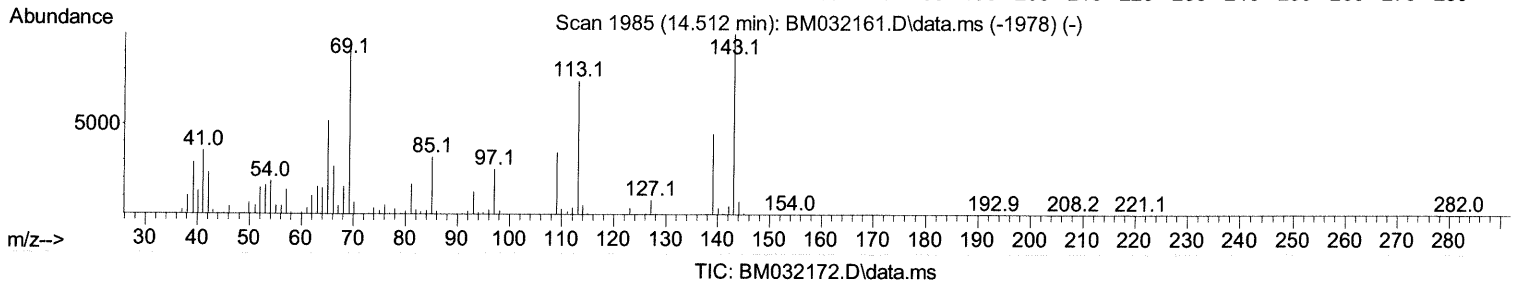
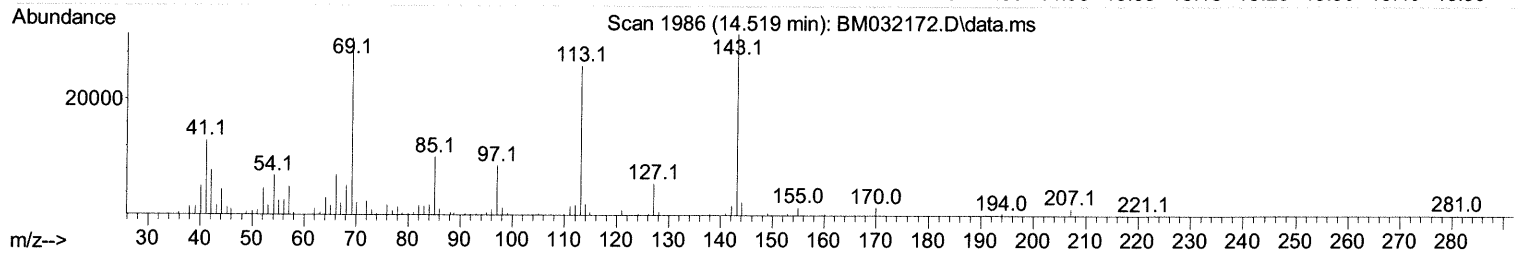
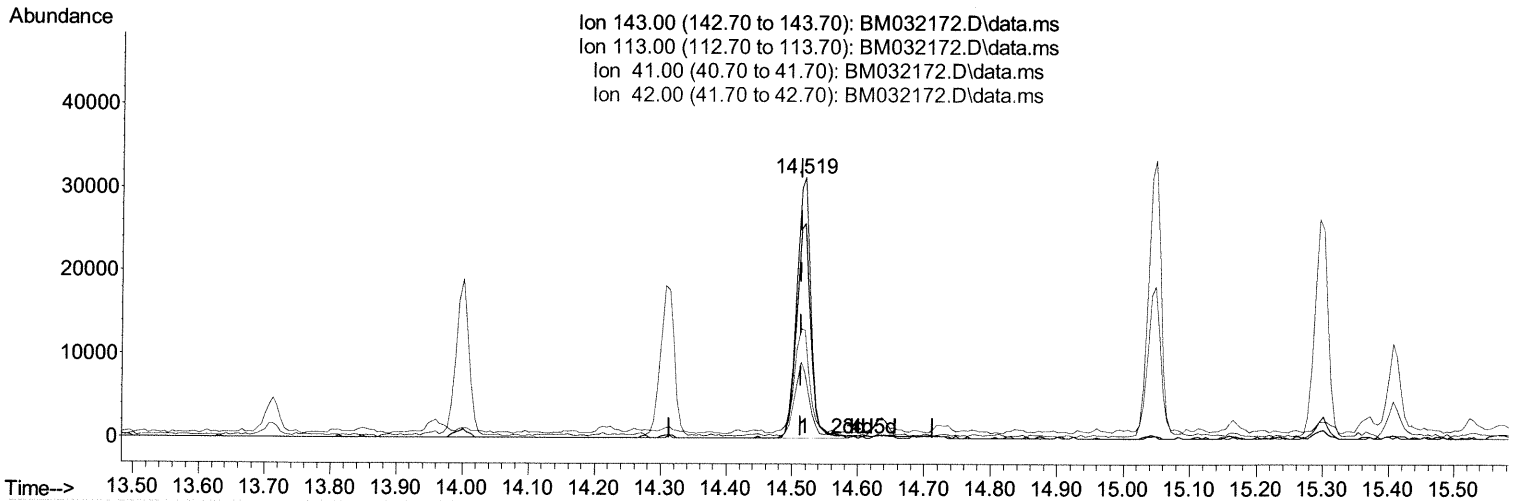
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Manual Integrations
 APPROVED

mohammad
 9/27/2021 3:06:37 PM

Quant Time: Sep 26 03:18:01 2021
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 QLast Update : Sat Sep 25 01:15:33 2021
 Response via : Initial Calibration



(54) 4-Nitrophenol-d4 (S)

14.519min (+ 0.006) 4.51 ng/ul m 09/29/21 JU

response 46912

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.30	82.46
41.00	38.30	41.38
42.00	26.20	25.11

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092521\
 Data File : BM032172.D
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 Operator : CG/JU
 Sample : M3739-12
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0B71

Manual Integrations
 APPROVED

mohammad
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.666	152	455374	20.000	ng/ul	0.00
20) Naphthalene-d8	10.460	136	1779935	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.313	164	983275	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.036	188	1774361	20.000	ng/ul	0.00
79) Chrysene-d12	21.195	240	1520587	20.000	ng/ul	0.00
88) Perylene-d12	23.359	264	1507643	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.990	96	56376	5.227	ng/ul	0.00
4) Pyridine-d5	3.419	84	260449	8.507	ng/ul	0.00
7) Phenol-d5	6.848	99	202864	5.612	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.990	67	548382	25.231	ng/ul	0.00
11) 2-Chlorophenol-d4	7.201	132	490577	17.481	ng/ul	0.00
15) 4-Methylphenol-d8	8.389	113	338521	11.977	ng/ul	0.00
21) Nitrobenzene-d5	8.819	128	325271	30.596	ng/ul	0.00
24) 2-Nitrophenol-d4	9.542	143	260148	23.142	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.083	165	518589	20.180	ng/ul	0.00
31) 4-Chloroaniline-d4	10.589	131	559641	16.113	ng/ul	0.00
46) Dimethylphthalate-d6	13.713	166	2030510	31.784	ng/ul	0.00
49) Acenaphthylene-d8	14.001	160	2361527	29.095	ng/ul	0.00
54) 4-Nitrophenol-d4	14.519	143	46912m	4.514	ng/ul	0.00
60) Fluorene-d10	15.301	176	1705694	31.290	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.413	200	164668	19.637	ng/ul	0.00
73) Anthracene-d10	17.136	188	2551543	34.464	ng/ul	0.00
81) Pyrene-d10	19.418	212	2722511	34.029	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.224	264	2350359	33.219	ng/ul	-0.01

09/29/21 JU

Target Compounds Qvalue

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