

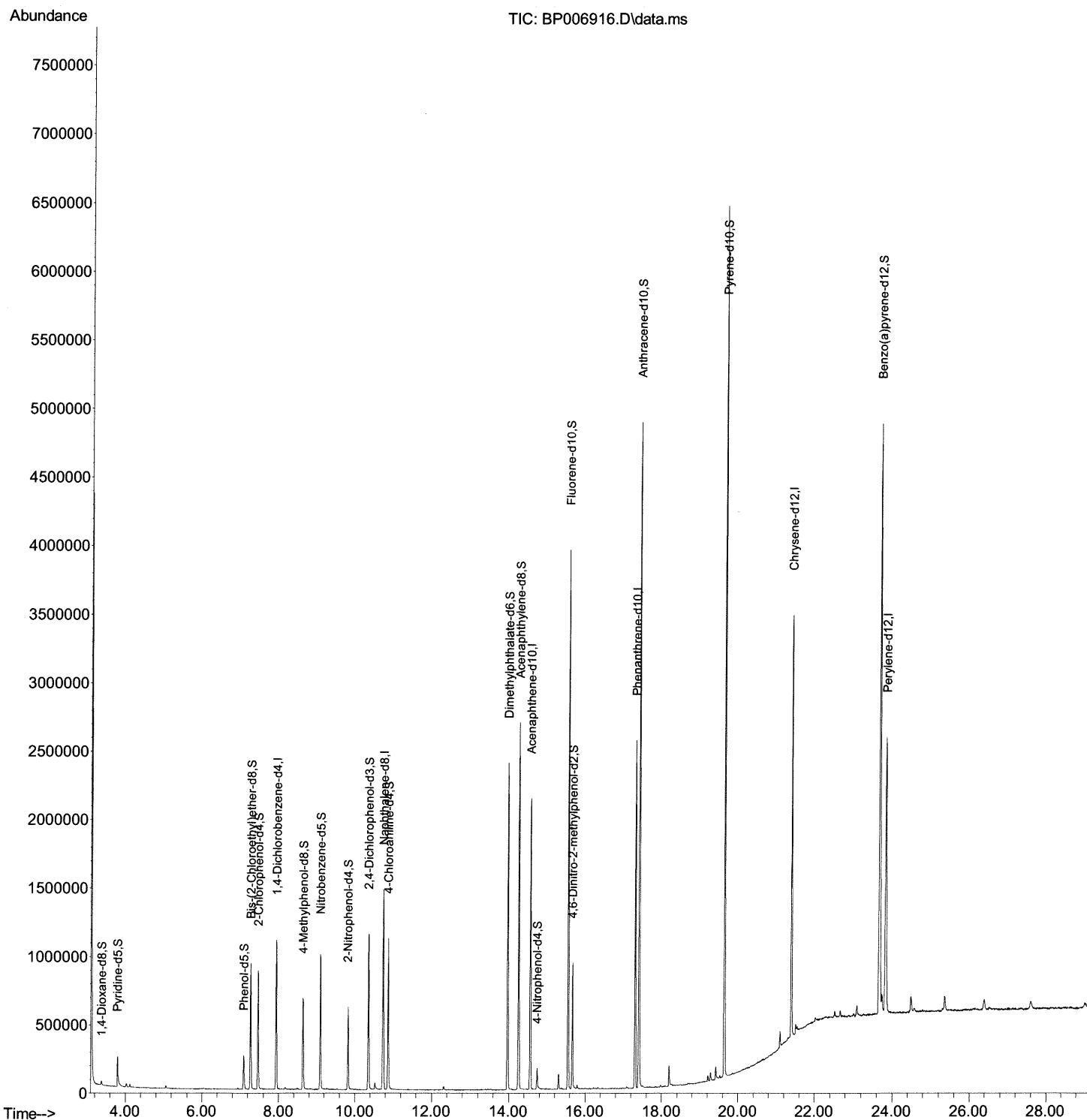
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090921\
Data File : BP006916.D
Acq On : 10 Sep 2021 00:23
Operator : CG/JU
Sample : M3651-13
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBKM7

Manual Integrations
APPROVED

mohammad
9/10/2021 3:26:09 PM

Quant Time: Sep 10 01:08:38 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP090821.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Sep 09 11:07:24 2021
Response via : Initial Calibration



Quantitation Report (Qedit)

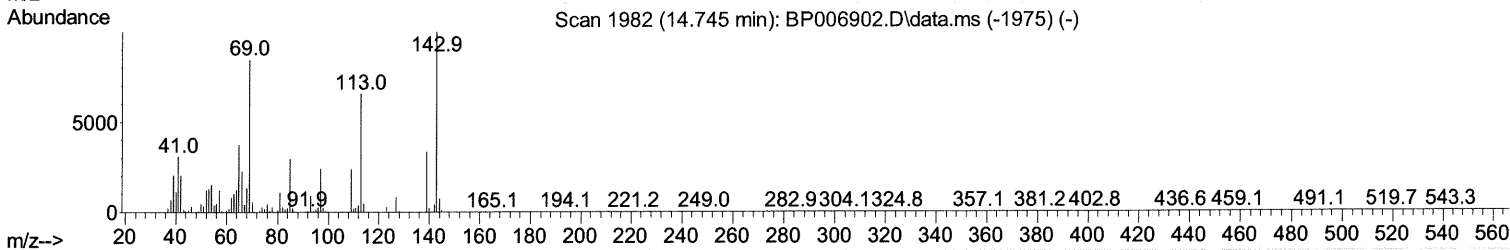
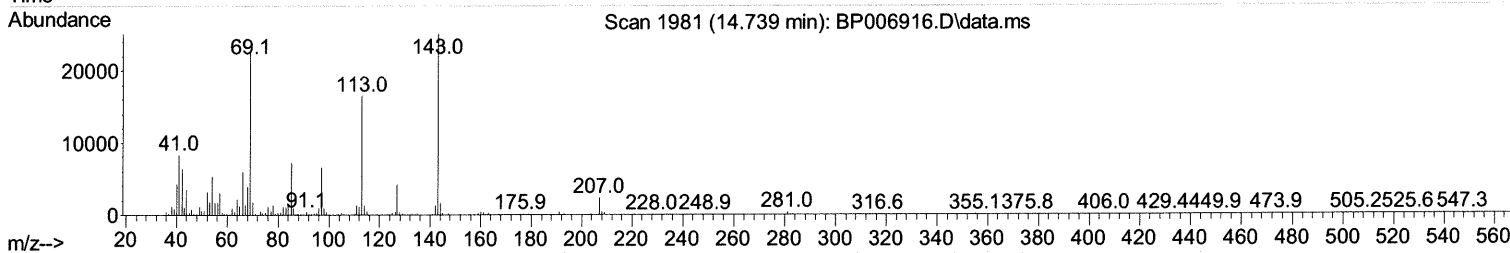
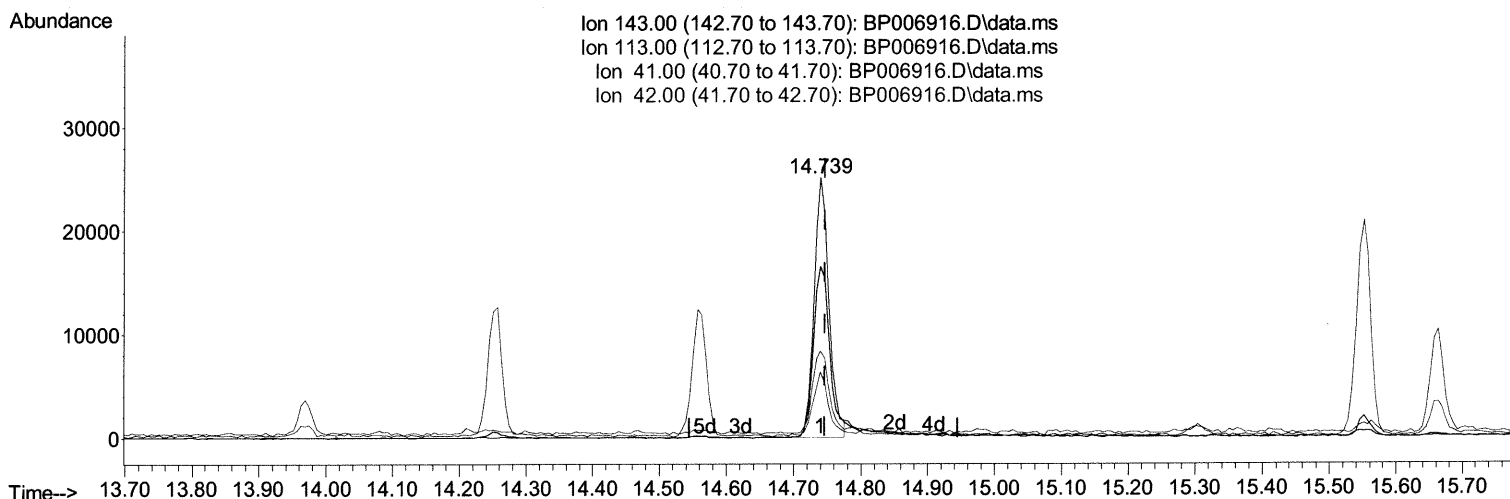
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TIC: BP006916.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.739min (-0.006) 4.57 ng/ul

response 38447

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	85.80	65.96#
41.00	40.70	33.39
42.00	26.80	25.39

Quantitation Report (Qedit)

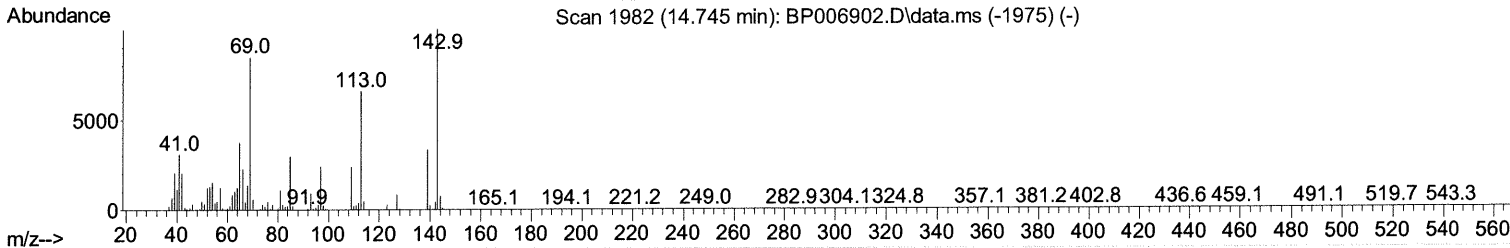
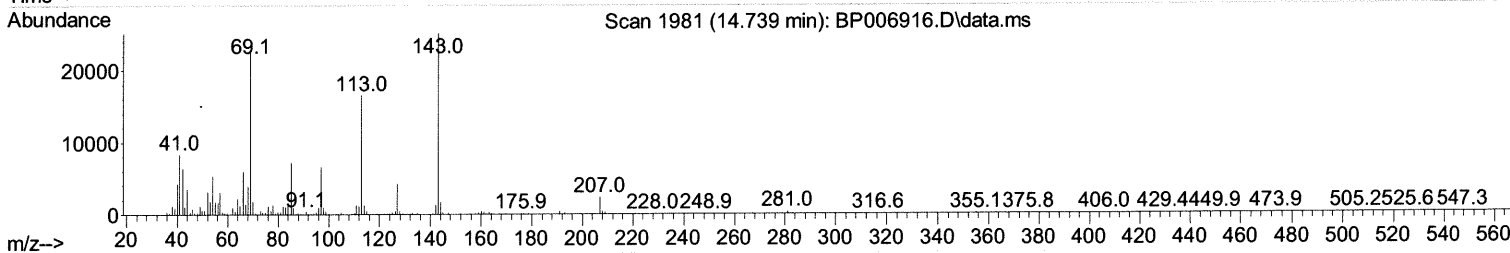
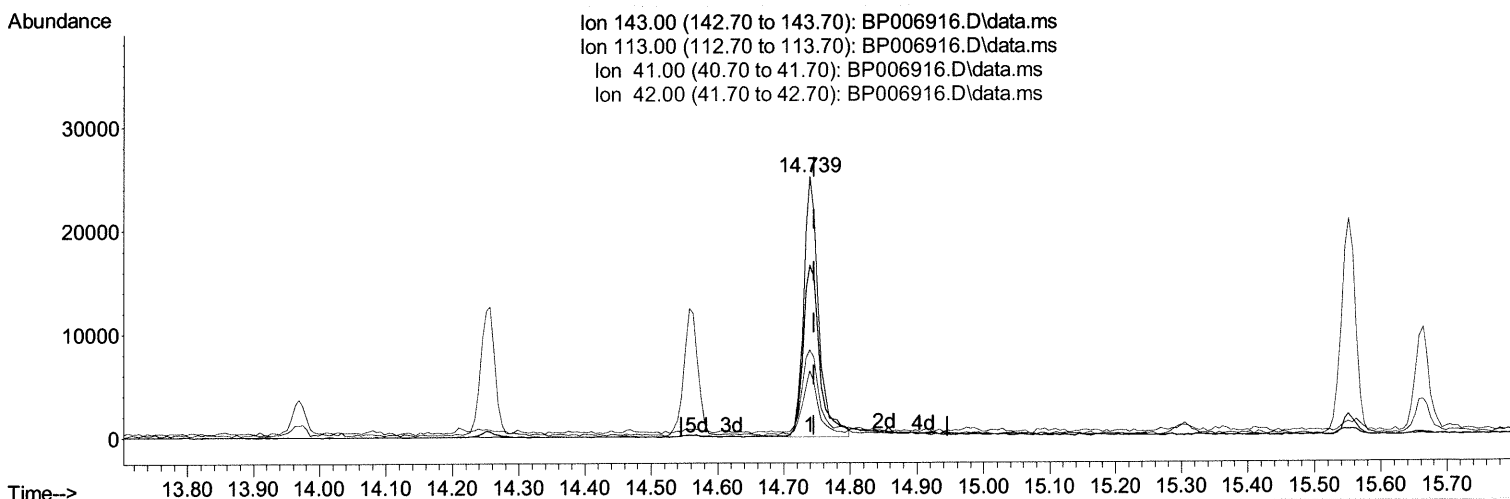
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Manual Integrations
 APPROVED

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 9/10/2021 3:26:09 PM

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TIC: BP006916.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.739min (-0.006) 4.77 ng/ul m 09/13/21 JU

response 40076

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	85.80	65.96#
41.00	40.70	33.39
42.00	26.80	25.39

Quantitation Report (QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090921\
 Data File : BP006916.D
 Acq On : 10 Sep 2021 00:23
 Operator : CG/JU
 Sample : M3651-13
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBKM7

Manual Integrations
 APPROVED

mohammad
 9/10/2021 3:26:09 PM

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Compound	R.T.	QIon	Response	Conc Units	Dev(Min)
Internal Standards					
1) 1,4-Dichlorobenzene-d4	7.939	152	310049	20.000 ng/ul	0.00
20) Naphthalene-d8	10.739	136	1236997	20.000 ng/ul	0.00
38) Acenaphthene-d10	14.563	164	747277	20.000 ng/ul	0.00
64) Phenanthrene-d10	17.310	188	1514245	20.000 ng/ul	# 0.00
79) Chrysene-d12	21.392	240	1477523	20.000 ng/ul	# 0.00
88) Perylene-d12	23.827	264	1533831	20.000 ng/ul	-0.01
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.369	96	16832	2.160 ng/uL	0.00
4) Pyridine-d5	3.793	84	146611	7.221 ng/ul	0.01
7) Phenol-d5	7.092	99	147724	6.483 ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.269	67	405941	31.017 ng/ul	0.00
11) 2-Chlorophenol-d4	7.463	132	409743	22.273 ng/ul	0.00
15) 4-Methylphenol-d8	8.634	113	258885	14.467 ng/ul	0.00
21) Nitrobenzene-d5	9.098	128	243380	32.483 ng/ul	0.00
24) 2-Nitrophenol-d4	9.816	143	206861	28.043 ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.351	165	424508	24.406 ng/ul	0.00
31) 4-Chloroaniline-d4	10.869	131	611156	25.408 ng/ul	0.00
46) Dimethylphthalate-d6	13.969	166	1591336	32.879 ng/ul	0.00
49) Acenaphthylene-d8	14.257	160	1924362	31.393 ng/ul	0.00
54) 4-Nitrophenol-d4	14.739	143	40076m>	4.766 ng/ul>	0.00 09/13/21 JU
60) Fluorene-d10	15.551	176	1519304	35.568 ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.663	200	189586	28.302 ng/ul	0.00
73) Anthracene-d10	17.410	188	2743547	43.764 ng/ul	0.00
81) Pyrene-d10	19.639	212	3151403	45.543 ng/ul	0.00
92) Benzo(a)pyrene-d12	23.668	264	3102272	41.925 ng/ul	-0.01

Target Compounds Qvalue

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