

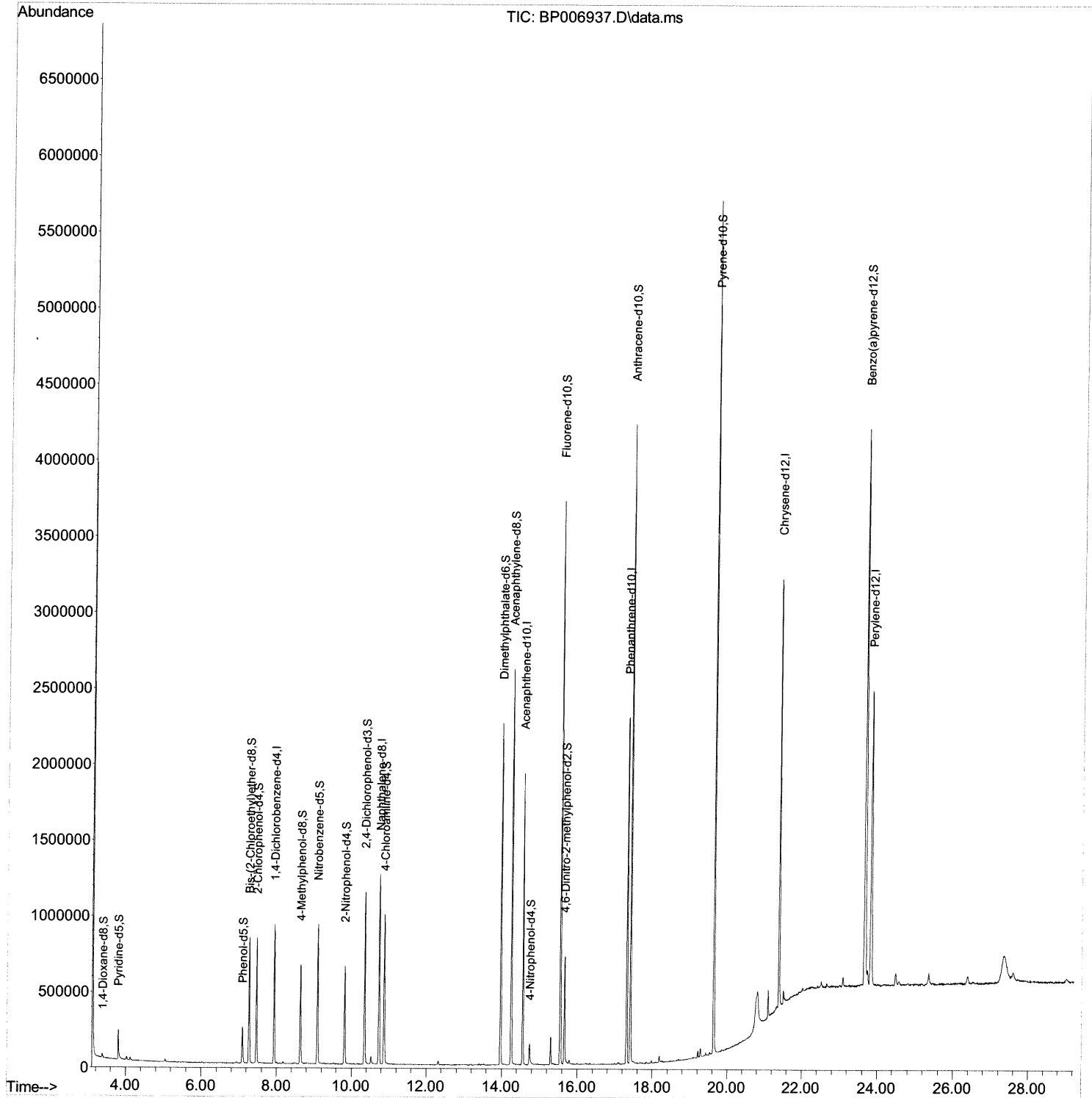
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090921\
Data File : BP006937.D
Acq On : 10 Sep 2021 12:53
Operator : CG/JU
Sample : M3651-10
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBKM4

Manual Integrations
APPROVED

mohammad
9/10/2021 3:26:31 PM

Quant Time: Sep 10 13:56:48 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)

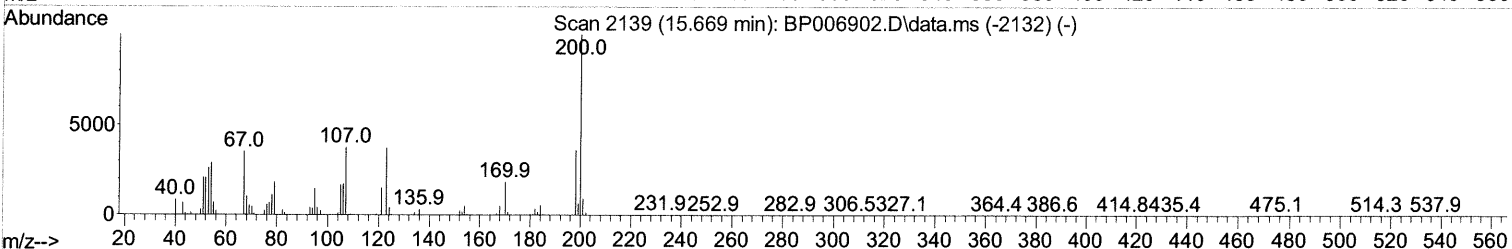
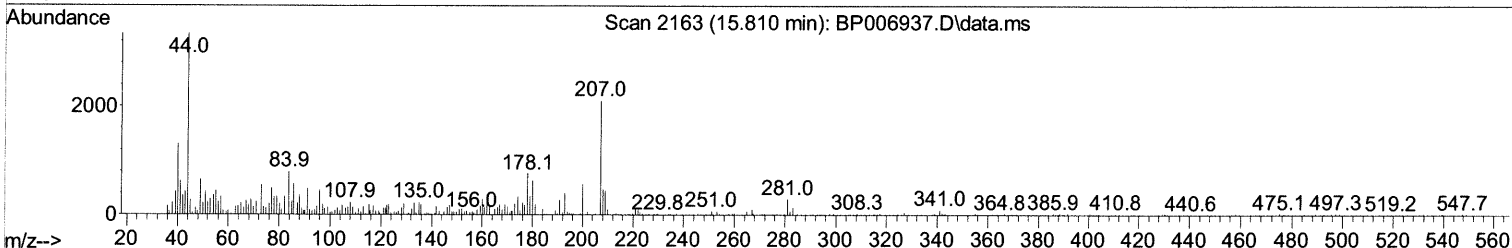
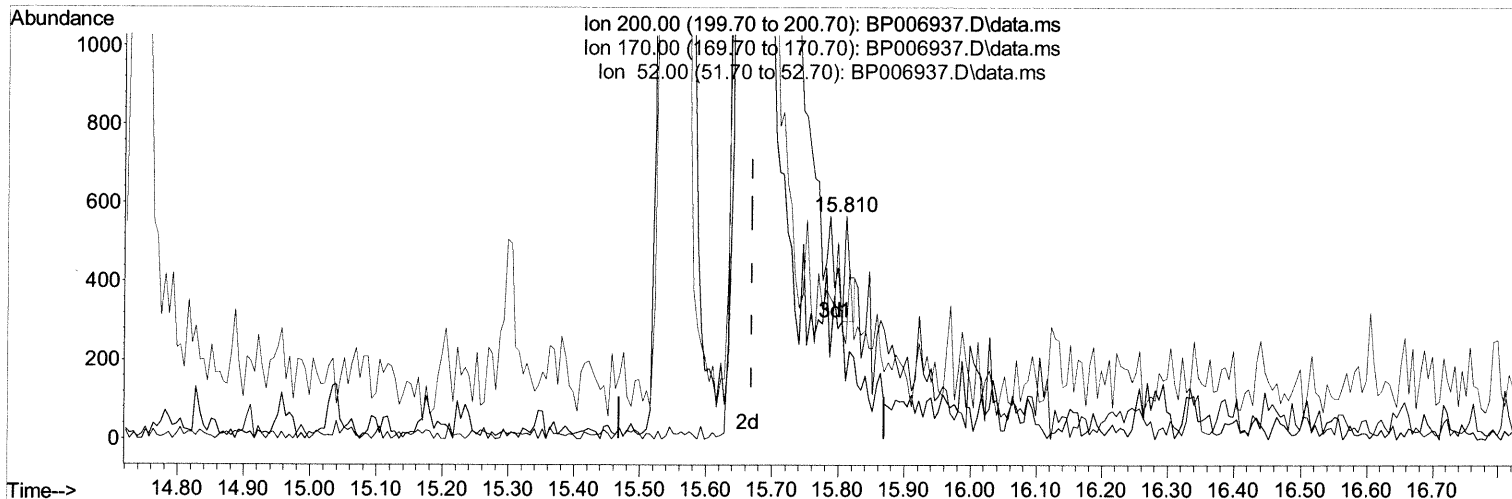
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Manual Integrations
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mohammad
 9/10/2021 3:26:31 PM

Quant Time: Sep 10 19:19:09 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



TIC: BP006937.D\data.ms

(65) 4,6-Dinitro-2-methylphenol-d2 (S)

15.810min (+ 0.141) 0.03 ng/u1

response 174

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	22.30	25.93
52.00	48.90	42.86
0.00	0.00	0.00

Quantitation Report (Qedit)

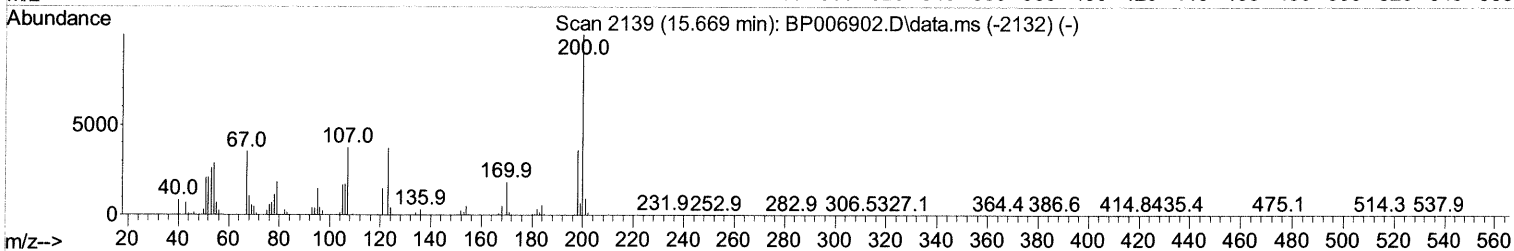
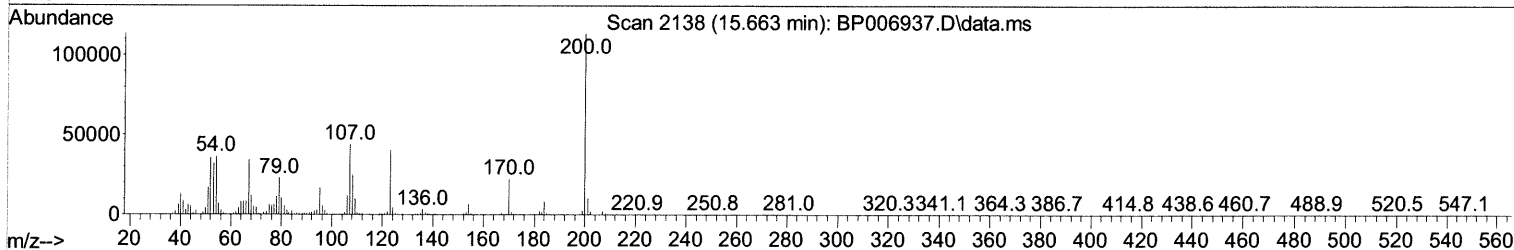
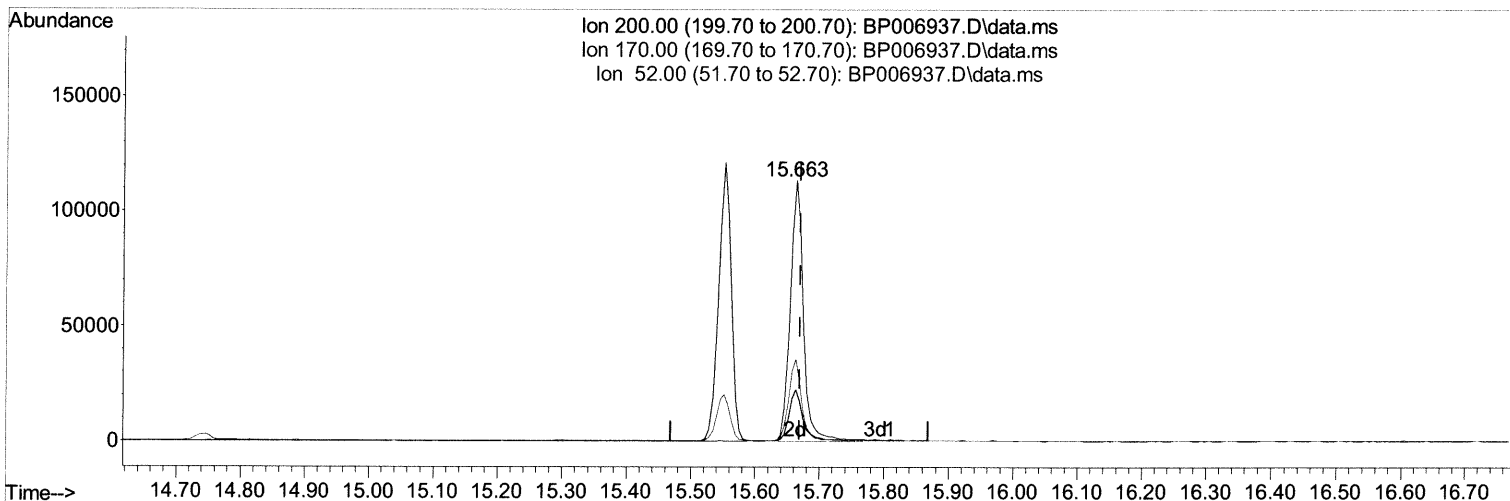
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Manual Integrations
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TIC: BP006937.D\data.ms

(65) 4,6-Dinitro-2-methylphenol-d2 (S)

15.663min (-0.006) 26.40 ng/ul m 09/14/21 JU

response 163654

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	22.30	19.78
52.00	48.90	31.42#
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090921\
 Data File : BP006937.D
 Acq On : 10 Sep 2021 12:53
 Operator : CG/JU
 Sample : M3651-10
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBKM4

Manual Integrations
 APPROVED

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.934	152	259291	20.000	ng/ul	0.00
20) Naphthalene-d8	10.740	136	1059291	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.557	164	668066	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.310	188	1401136	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.386	240	1413760	20.000	ng/ul	-0.01
88) Perylene-d12	23.821	264	1474126	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.370	96	15191	2.331	ng/ul	0.00
4) Pyridine-d5	3.793	84	125374	7.384	ng/ul	0.01
7) Phenol-d5	7.093	99	144262	7.570	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.263	67	368087	33.631	ng/ul	0.00
11) 2-Chlorophenol-d4	7.463	132	401279	26.083	ng/ul	0.00
15) 4-Methylphenol-d8	8.634	113	259012	17.307	ng/ul	0.00
21) Nitrobenzene-d5	9.093	128	231105	36.019	ng/ul	0.00
24) 2-Nitrophenol-d4	9.816	143	217655	34.456	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.352	165	422413	28.360	ng/ul	0.00
31) 4-Chloroaniline-d4	10.869	131	551951	26.796	ng/ul	0.00
46) Dimethylphthalate-d6	13.969	166	1540687	35.607	ng/ul	0.00
49) Acenaphthylene-d8	14.251	160	1838524	33.549	ng/ul	0.00
54) 4-Nitrophenol-d4	14.745	143	37947	5.048	ng/ul	0.00
60) Fluorene-d10	15.551	176	1400279	36.668	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.663	200	163654m	26.403	ng/ul	> 0.00 09/14/21 JU
73) Anthracene-d10	17.404	188	2407272	41.499	ng/ul	-0.01
81) Pyrene-d10	19.639	212	2841547	42.917	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.668	264	2839411	39.927	ng/ul	-0.01

Target Compounds Qvalue

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