

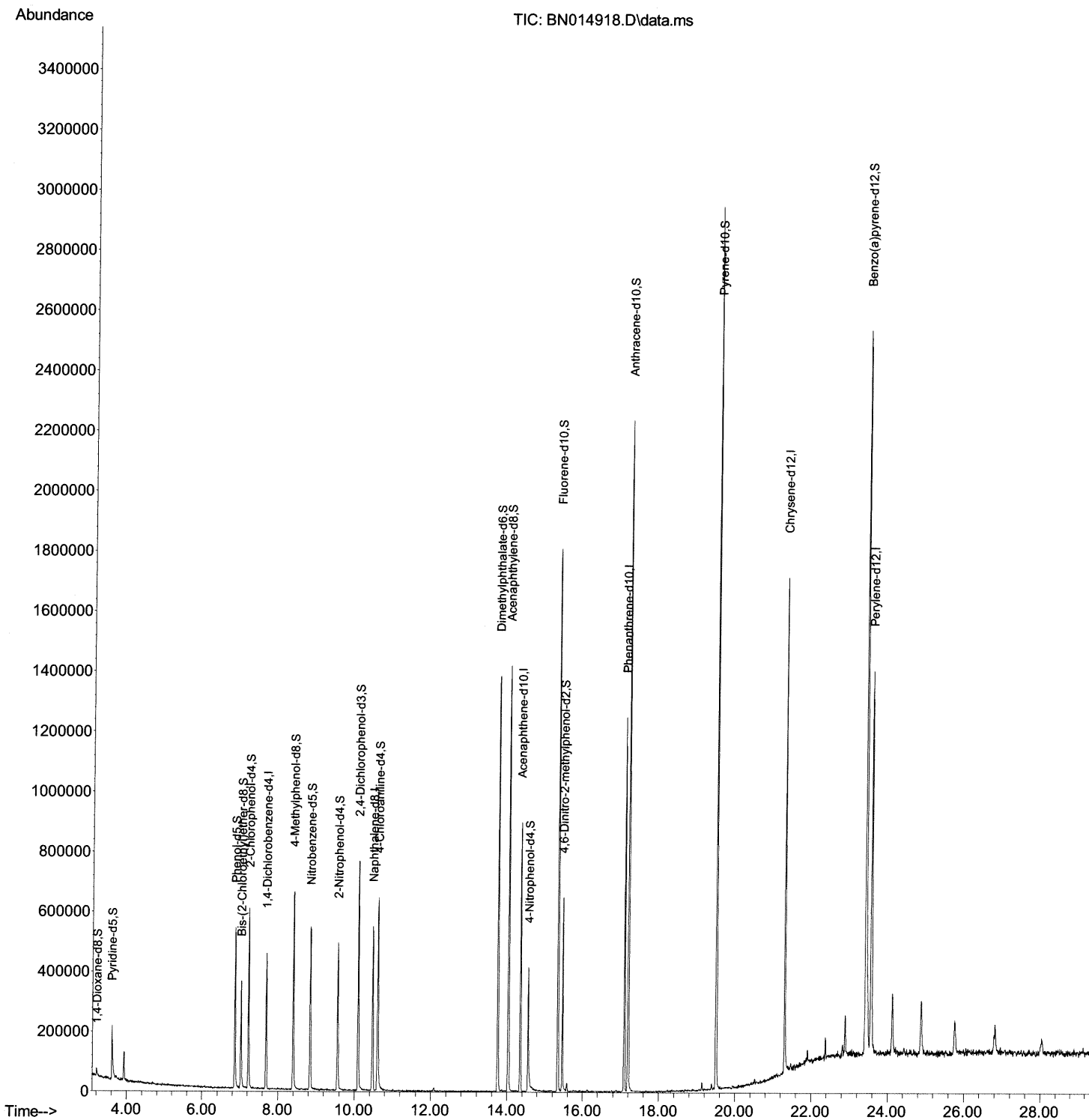
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061421\
Data File : BN014918.D
Acq On : 15 Jun 2021 08:44
Operator : CG/JU
Sample : PB137010BL
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
SBLK010

Manual Integrations
APPROVED

mohammad
6/15/2021 2:42:15 PM

Quant Time: Jun 15 10:26:34 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN061221MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Jun 14 10:46:06 2021
Response via : Initial Calibration



Quantitation Report (Qedit)

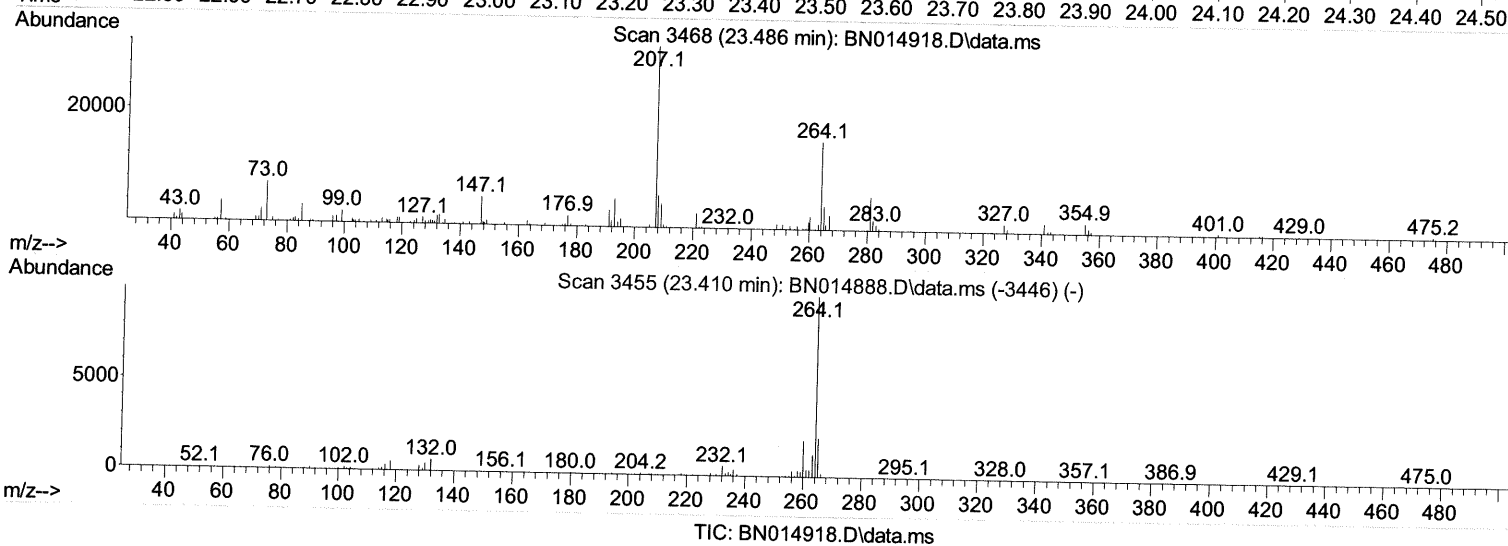
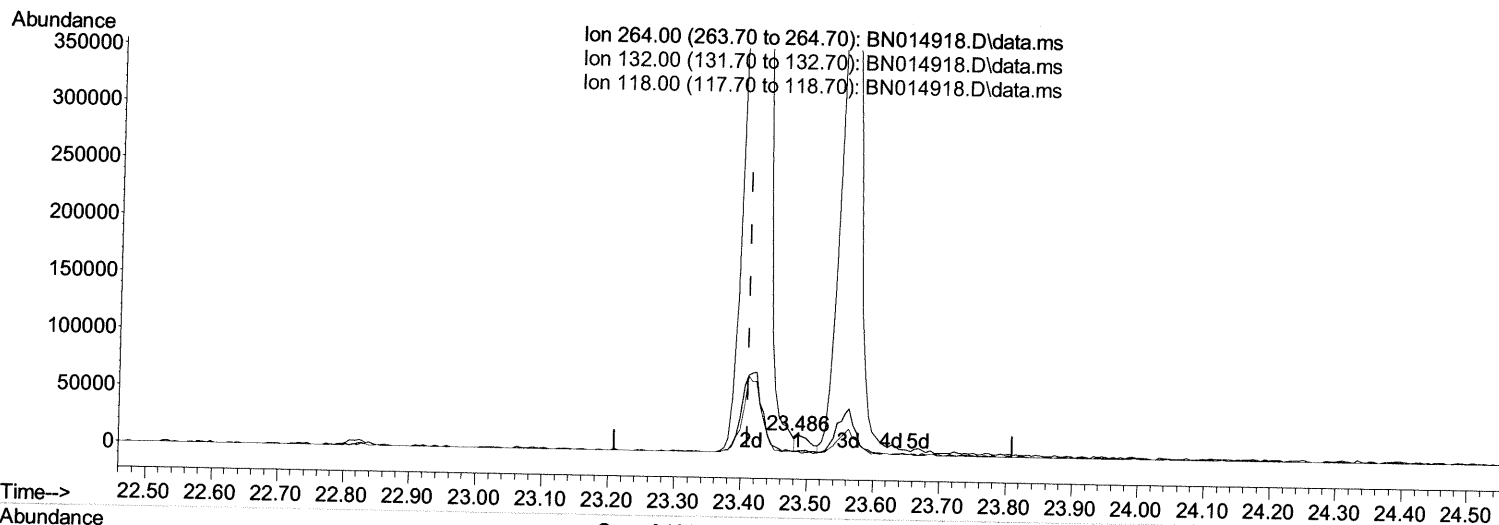
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(92) Benzo(a)pyrene-d12 (S)

23.486min (+ 0.076) 0.44 ng/ul

response 21097

Ion	Exp%	Act%
264.00	100.00	100.00
132.00	9.00	10.76
118.00	7.50	8.15
0.00	0.00	0.00

```
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```

mohammad
6/15/2021 2:42:15 PM

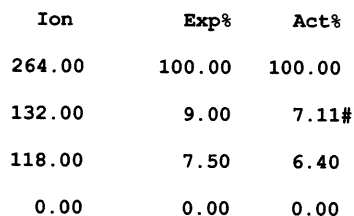
Abundance

Ion 264.00 (263.70 to 264.70): BN014918.D\data.ms
Ion 132.00 (131.70 to 132.70): BN014918.D\data.ms
Ion 118.00 (117.70 to 118.70): BN014918.D\data.ms

23416

2d 1 3d 4d5d

Time-->



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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.687	152	112109	20.000	ng/ul	0.00
20) Naphthalene-d8	10.475	136	394634	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.346	164	280334	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.098	188	663087	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.304	240	839398	20.000	ng/ul	0.00
88) Perylene-d12	23.563	264	965474	20.000	ng/ul	0.01

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.205	96	15079	7.152	ng/uL	0.00
4) Pyridine-d5	3.611	84	128460	26.195	ng/ul	0.00
7) Phenol-d5	6.864	99	286757	34.148	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.028	67	140721	35.372	ng/ul	0.00
11) 2-Chlorophenol-d4	7.228	132	228403	35.473	ng/ul	0.00
15) 4-Methylphenol-d8	8.399	113	219736	30.706	ng/ul	0.00
21) Nitrobenzene-d5	8.846	128	115270	38.926	ng/ul	0.00
24) 2-Nitrophenol-d4	9.563	143	122252	36.403	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.099	165	237060	35.060	ng/ul	0.00
31) 4-Chloroaniline-d4	10.616	131	282184	35.238	ng/ul	0.00
46) Dimethylphthalate-d6	13.757	166	864607	38.796	ng/ul	0.00
49) Acenaphthylene-d8	14.034	160	951514	39.062	ng/ul	0.00
54) 4-Nitrophenol-d4	14.557	143	86356	27.585	ng/ul	0.00
60) Fluorene-d10	15.345	176	703723	38.709	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.463	200	130044	32.810	ng/ul	0.00
73) Anthracene-d10	17.198	188	1169497	38.972	ng/ul	0.00
81) Pyrene-d10	19.504	212	1525099	39.392	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.416	264	1976061m	40.954	ng/ul	> 0.00 @ 6/16/21 JU

Target Compounds

Qvalue

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