

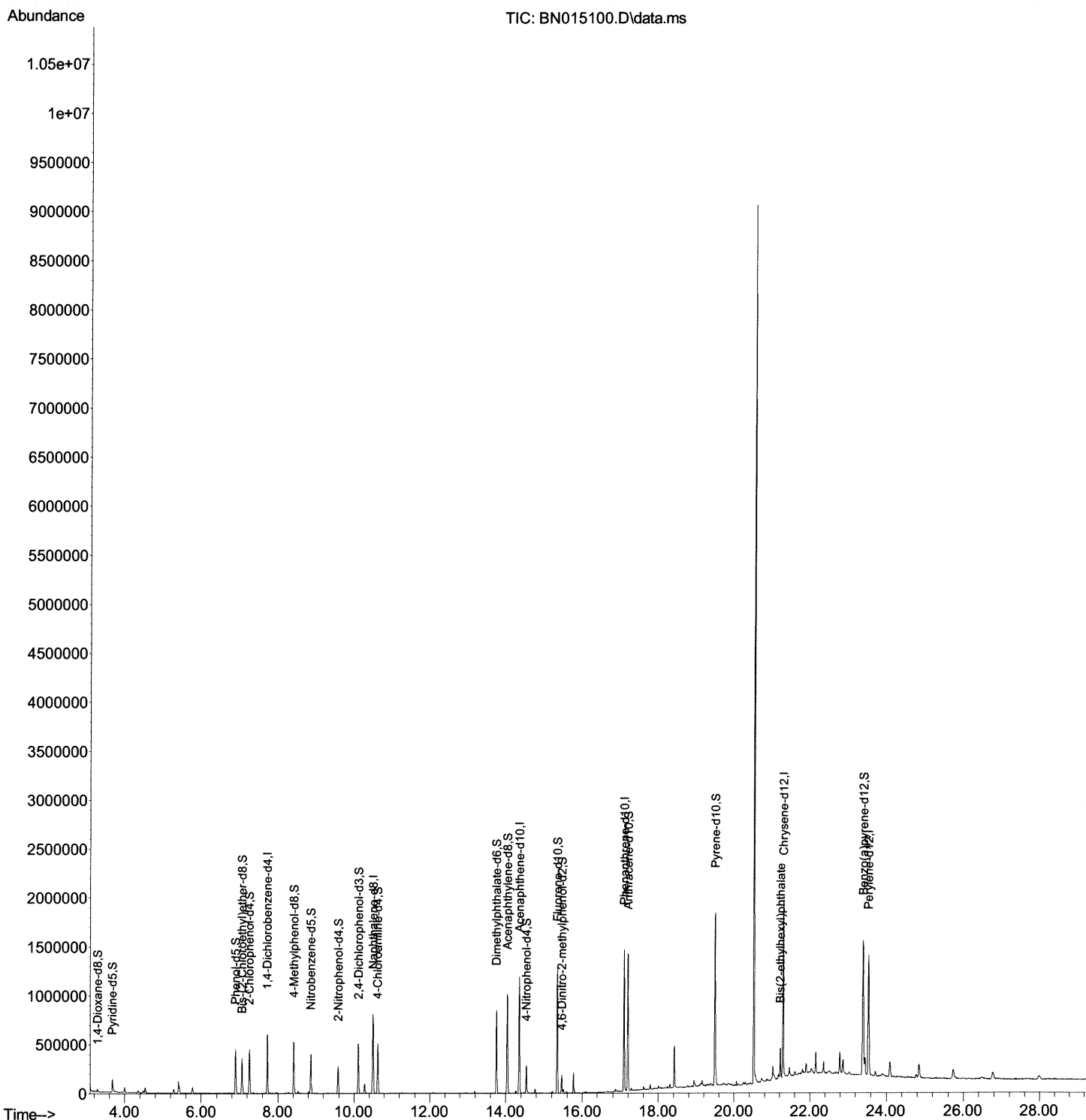
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062621\
Data File : BN015100.D
Acq On : 26 Jun 2021 05:04
Operator : CG/JU
Sample : M2704-12
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BGDD5

Manual Integrations
APPROVED

mohammad
6/28/2021 1:25:39 PM

Quant Time: Jun 26 06:11:57 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN061821MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Jun 24 14:43:13 2021
Response via : Initial Calibration



Quantitation Report (Qedit)

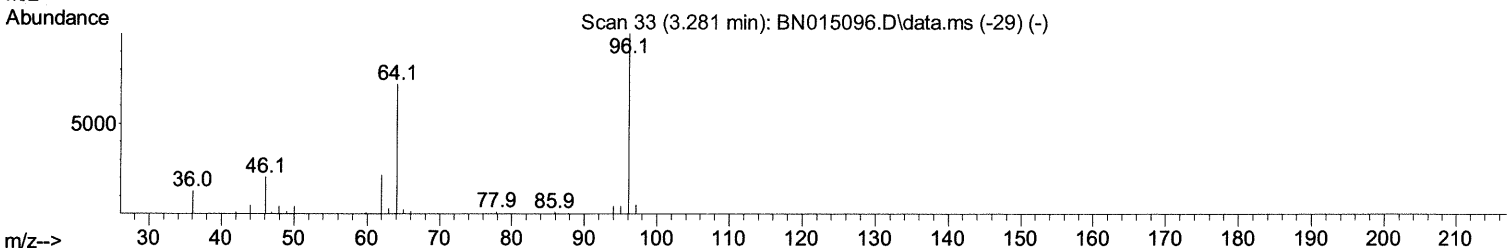
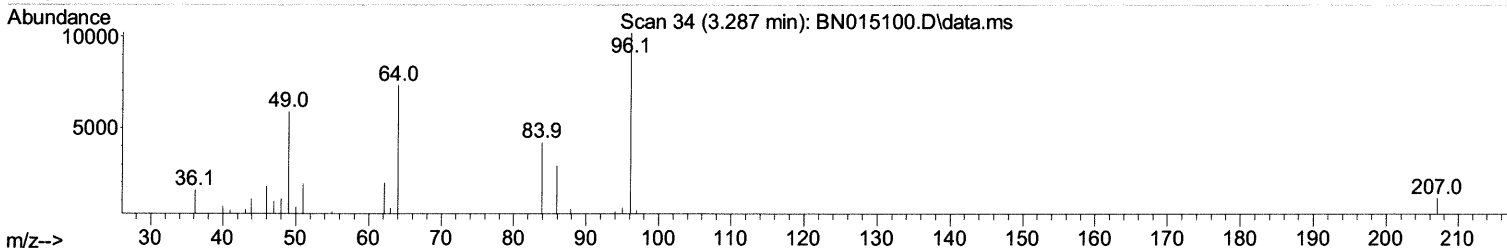
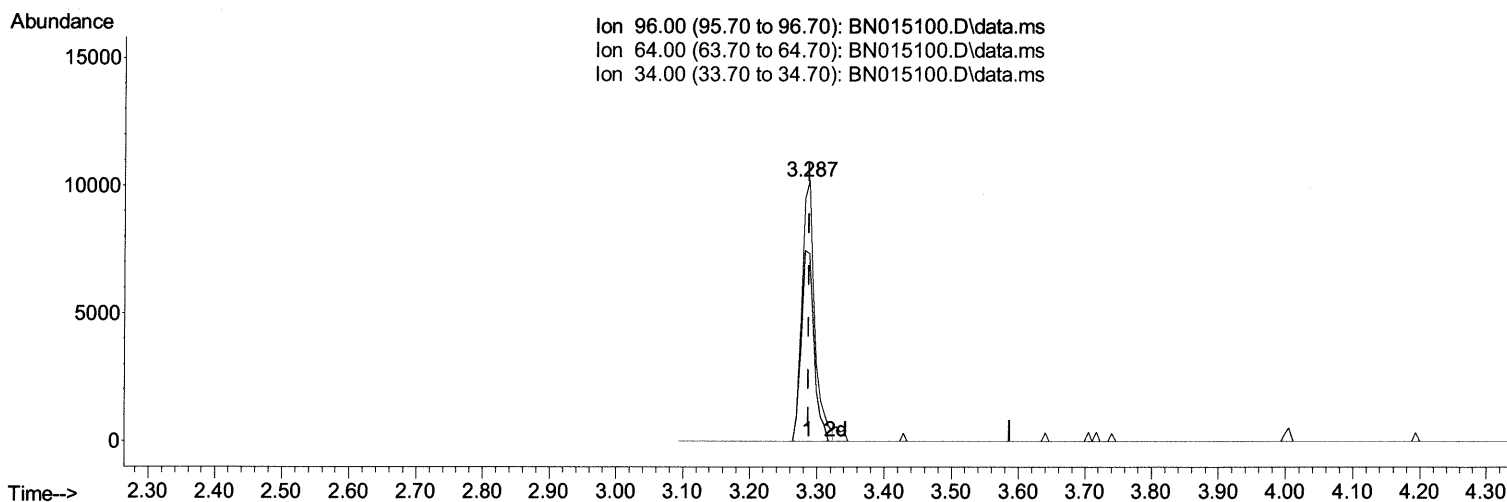
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TIC: BN015100.D\data.ms

(3) 1,4-Dioxane-d8 (S)

3.287min (-0.000) 3.22 ng/uL

response 13563

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	75.20	71.95
34.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

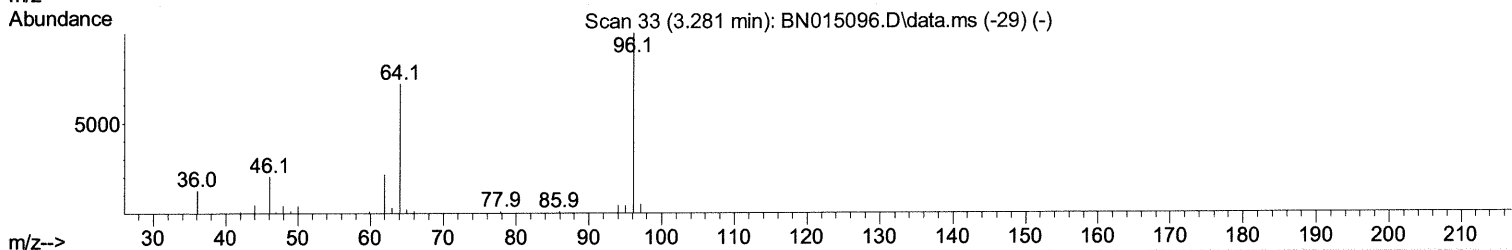
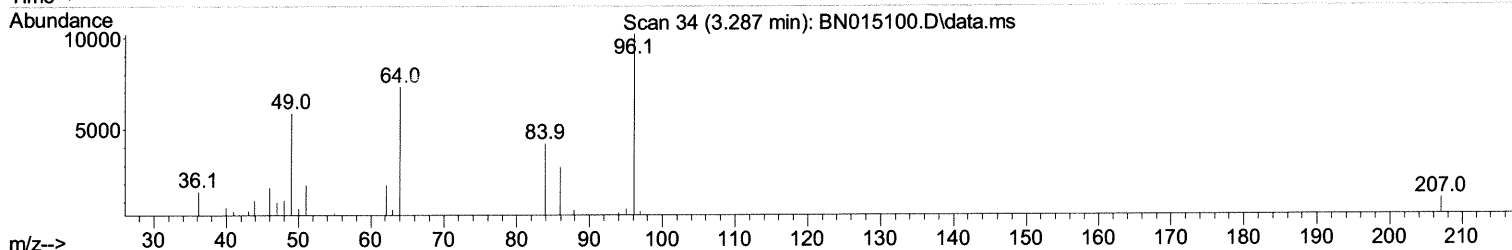
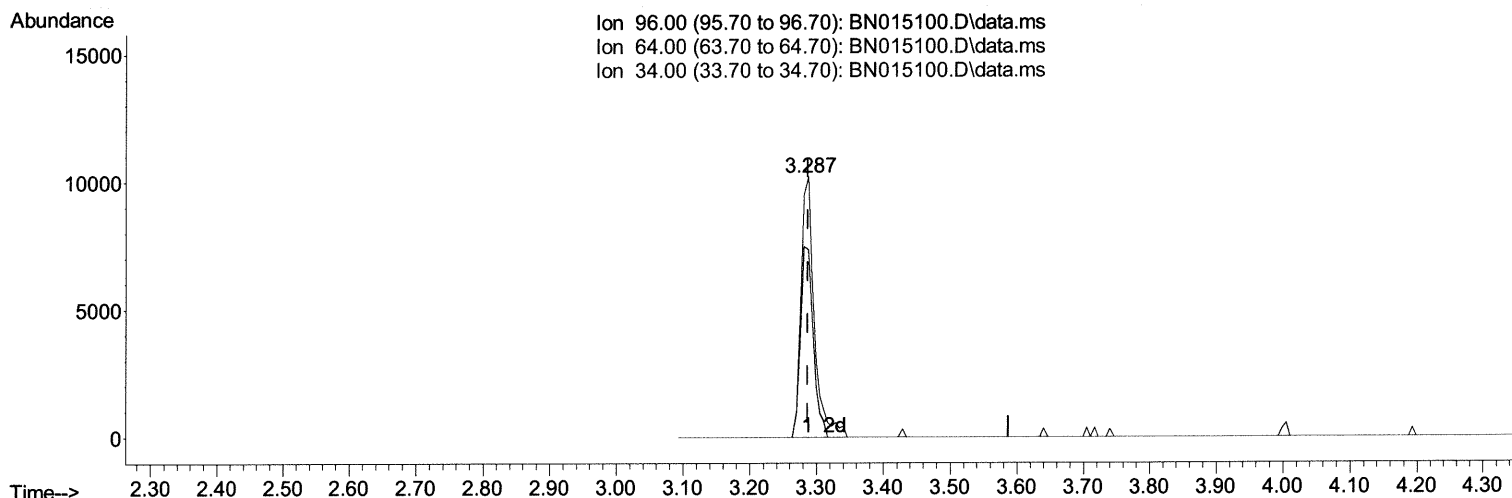
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Misc :
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Manual Integrations
APPROVED

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Quant Title : SVOA CALIBRATION
QLast Update : Thu Jun 24 14:43:13 2021
Response via : Initial Calibration



TIC: BN015100.D\data.ms

(3) 1,4-Dioxane-d8 (S)

3.287min (-0.000) 3.33 ng/uL m 06/24/21 J4

response 14038

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	75.20	71.95
34.00	0.00	0.00
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062621\
 Data File : BN015100.D
 Acq On : 26 Jun 2021 05:04
 Operator : CG/JU
 Sample : M2704-12
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampled :
 BGDD5

Manual Integrations
 APPROVED

mohammad
 6/28/2021 1:25:39 PM

Quant Time: Jun 26 06:11:57 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN061821MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jun 24 14:43:13 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.728	152	163901	20.000	ng/ul	0.00
20) Naphthalene-d8	10.498	136	670366	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.351	164	407275	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.098	188	818272	20.000	ng/ul	0.00
79) Chrysene-d12	21.292	240	810678	20.000	ng/ul	0.00
88) Perylene-d12	23.533	264	820237	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.287	96	14038m	3.333	ng/ul	0.00
4) Pyridine-d5	3.681	84	73369	6.461	ng/ul	0.00
7) Phenol-d5	6.893	99	258025	18.138	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.063	67	152909	18.303	ng/ul	0.00
11) 2-Chlorophenol-d4	7.257	132	206522	19.146	ng/ul	0.00
15) 4-Methylphenol-d8	8.416	113	191957	17.484	ng/ul	0.00
21) Nitrobenzene-d5	8.869	128	95935	20.921	ng/ul	0.00
24) 2-Nitrophenol-d4	9.587	143	90431	22.139	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.116	165	181725	18.849	ng/ul	0.00
31) 4-Chloroaniline-d4	10.628	131	260427	17.430	ng/ul	0.00
46) Dimethylphthalate-d6	13.757	166	567384	20.288	ng/ul	0.00
49) Acenaphthylene-d8	14.039	160	723893	19.939	ng/ul	0.00
54) 4-Nitrophenol-d4	14.539	143	65689	13.466	ng/ul	0.00
60) Fluorene-d10	15.345	176	503402	20.690	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.457	200	35415	10.479	ng/ul	0.00
73) Anthracene-d10	17.192	188	788633	21.057	ng/ul	0.00
81) Pyrene-d10	19.498	212	924691	21.762	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.392	264	937936	23.163	ng/ul	-0.01
Target Compounds						
86) Bis(2-ethylhexyl)phtha...	21.221	149	104352	3.419	ng/ul	99

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