

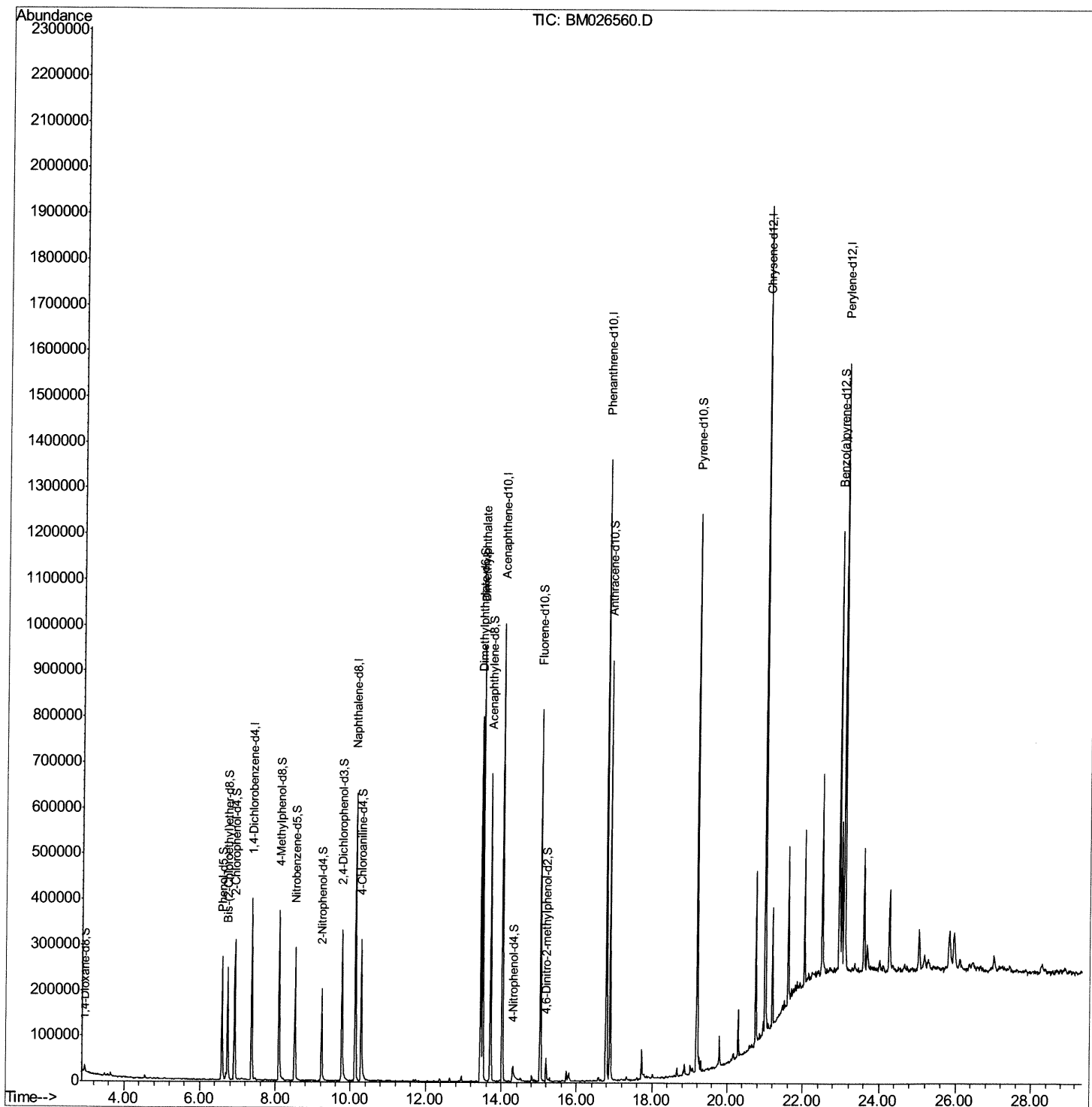
Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM062520\
Data File : BM026560.D
Acq On : 25 Jun 2020 16:00
Operator : JU/CG
Sample : L3062-08
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
ET141

Manual Integrations
APPROVED

mohammad
6/26/2020 12:19:56 PM

Quant Time: Jun 25 16:33:02 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061220MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Jun 25 10:53:13 2020
Response via : Initial Calibration



Quantitation Report (Qedit)

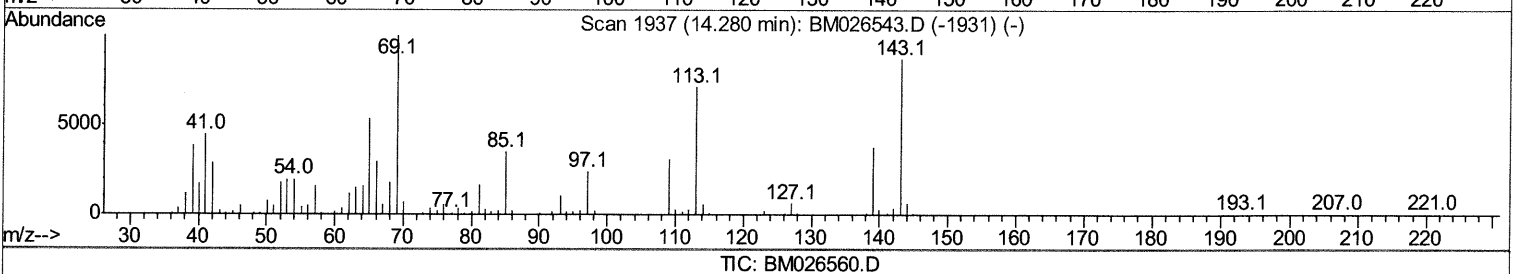
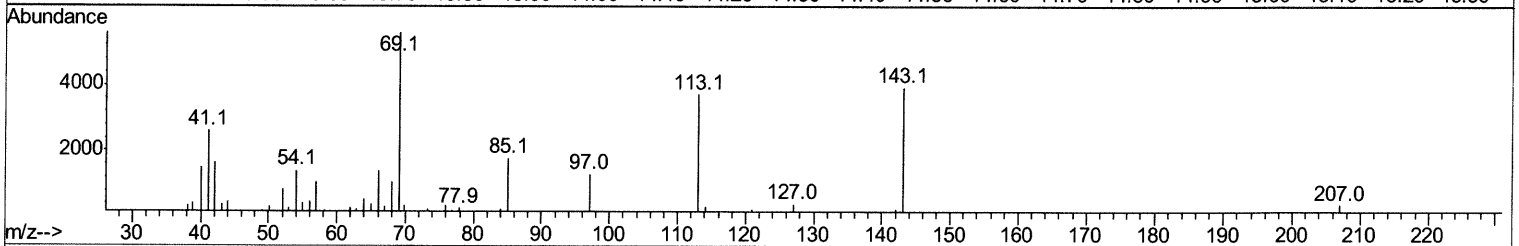
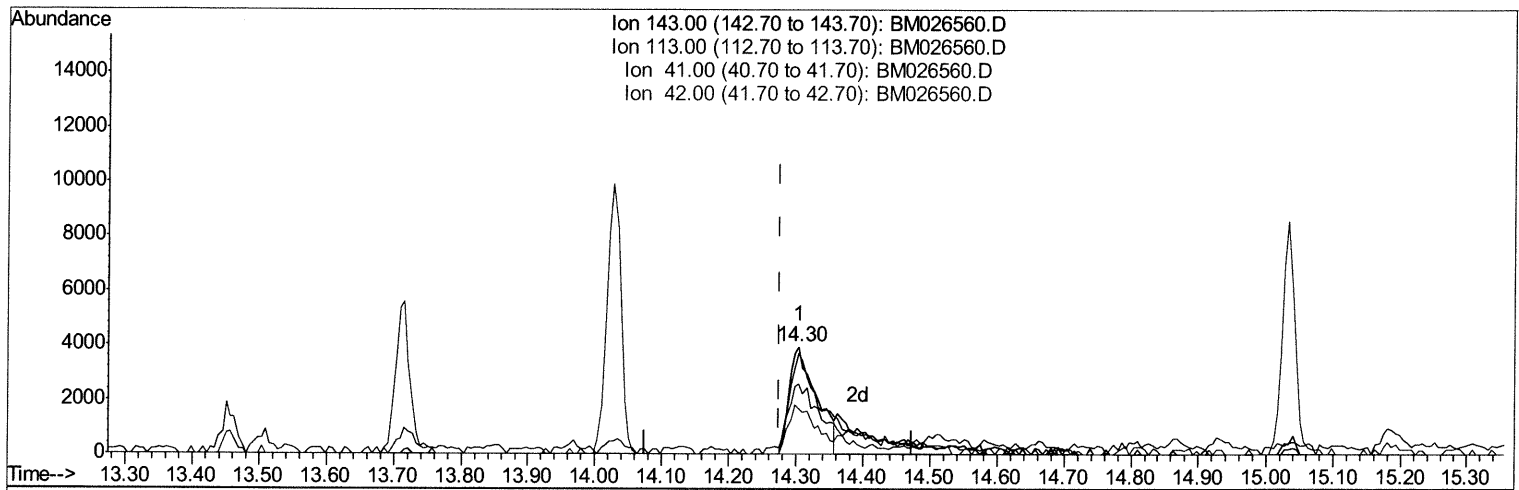
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Quant Time: Jun 25 16:32:04 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061220MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Jun 25 10:53:13 2020
Response via : Initial Calibration



(52) 4-Nitrophenol-d4 (S)
14.304min (+0.029) 2.62ng/ul
response 11199

Ion	Exp%	Act%
143.00	100	100
113.00	81.70	94.54
41.00	49.50	66.02#
42.00	34.20	42.21#

Quantitation Report (Qedit)

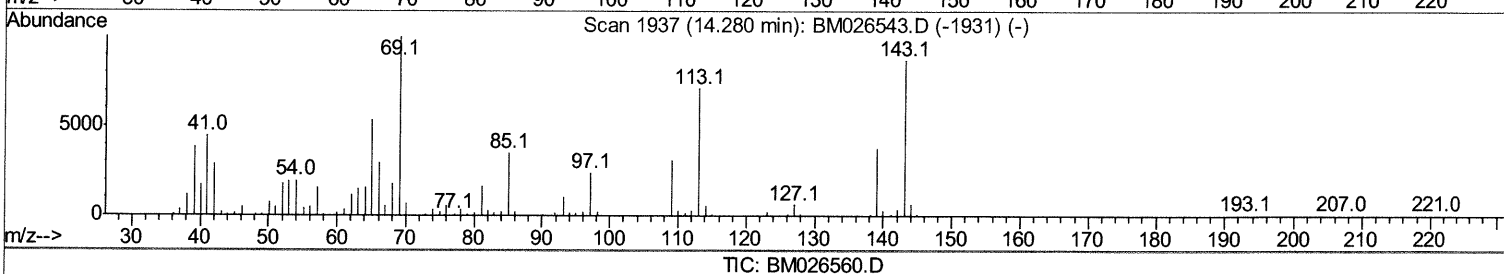
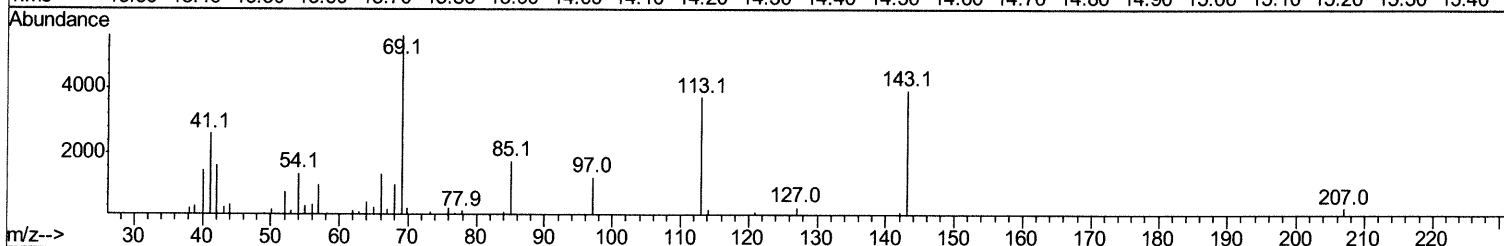
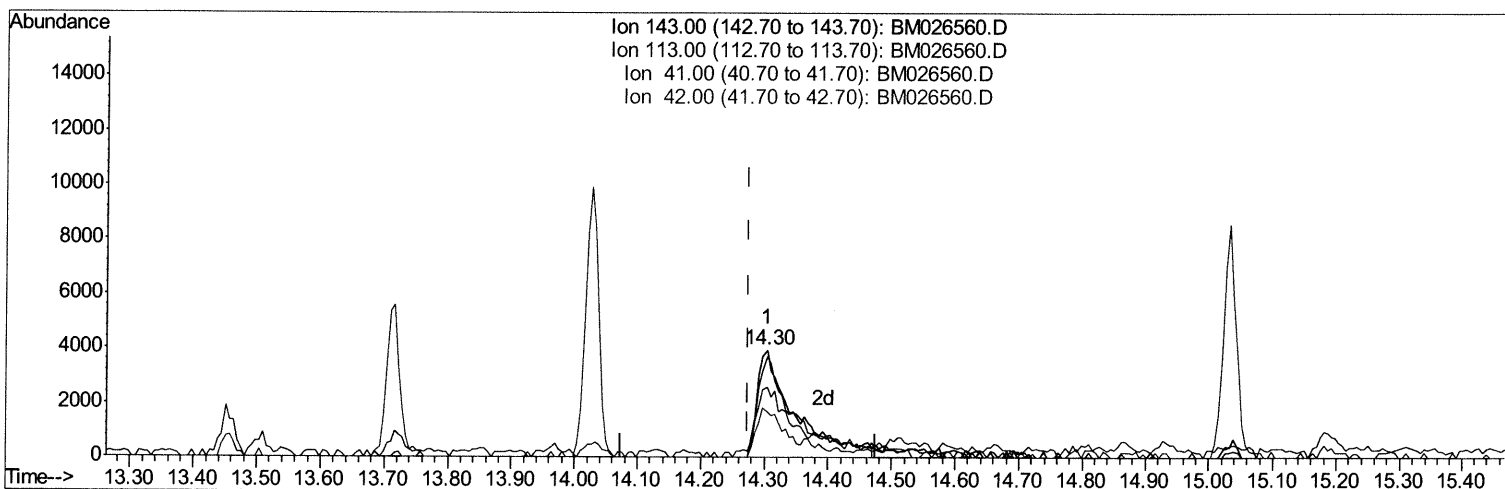
Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM062520\
Data File : BM026560.D
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Manual Integrations
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Quant Time: Jun 25 16:32:04 2020
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Quant Title : SVOA CALIBRATION
QLast Update : Thu Jun 25 10:53:13 2020
Response via : Initial Calibration



(52) 4-Nitrophenol-d4 (S)

14.304min (+0.029) 3.73ng/ul m

06/24/20 JU

response 15965

Ion	Exp%	Act%
143.00	100	100
113.00	81.70	94.54
41.00	49.50	66.02#
42.00	34.20	42.21#

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM062520\
 Data File : BM026560.D
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 Sample : L3062-08
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 ET141

Manual Integrations
 APPROVED

mohammad
 6/26/2020 12:19:56 PM

Quant Time: Jun 25 16:33:02 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061220MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jun 25 10:53:13 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.37	152	94043	20.00	ng/ul	0.00
18) Naphthalene-d8	10.13	136	438855	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.03	164	306860	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.79	188	695105	20.00	ng/ul	0.00
78) Chrysene-d12	21.00	240	808946	20.00	ng/ul	0.00
86) Perylene-d12	23.09	264	846921	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	2.96	96	8094	3.02	ng/uL	0.00
5) Phenol-d5	6.58	99	137329	15.85	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.73	67	113012	20.62	ng/ul	0.00
9) 2-Chlorophenol-d4	6.92	132	116554	16.50	ng/ul	0.00
13) 4-Methylphenol-d8	8.10	113	125726	18.44	ng/ul	0.00
19) Nitrobenzene-d5	8.52	128	56635	14.52	ng/ul	0.00
22) 2-Nitrophenol-d4	9.23	143	49775	11.60	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.77	165	117037	14.82	ng/ul	0.00
29) 4-Chloroaniline-d4	10.29	131	163172	19.81	ng/ul	0.00
44) Dimethylphthalate-d6	13.46	166	455603	17.66	ng/ul	0.00
47) Acenaphthylene-d8	13.72	160	409592	13.23	ng/ul	0.00
52) 4-Nitrophenol-d4	14.30	143	15965m>	3.73	ng/ul>	0.03
58) Fluorene-d10	15.03	176	290371	13.11	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.18	200	10099	2.22	ng/ul	0.00
71) Anthracene-d10	16.89	188	485668	13.73	ng/ul	0.00
79) Pyrene-d10	19.19	212	577630	13.02	ng/ul	0.00
90) Benzo(a)pyrene-d12	22.96	264	623737	13.32	ng/ul	0.00

06/29/20 JU

Target Compounds					Ovalue
45) Dimethylphthalate	13.50	163	523853	19.551	ng/ul 99

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