

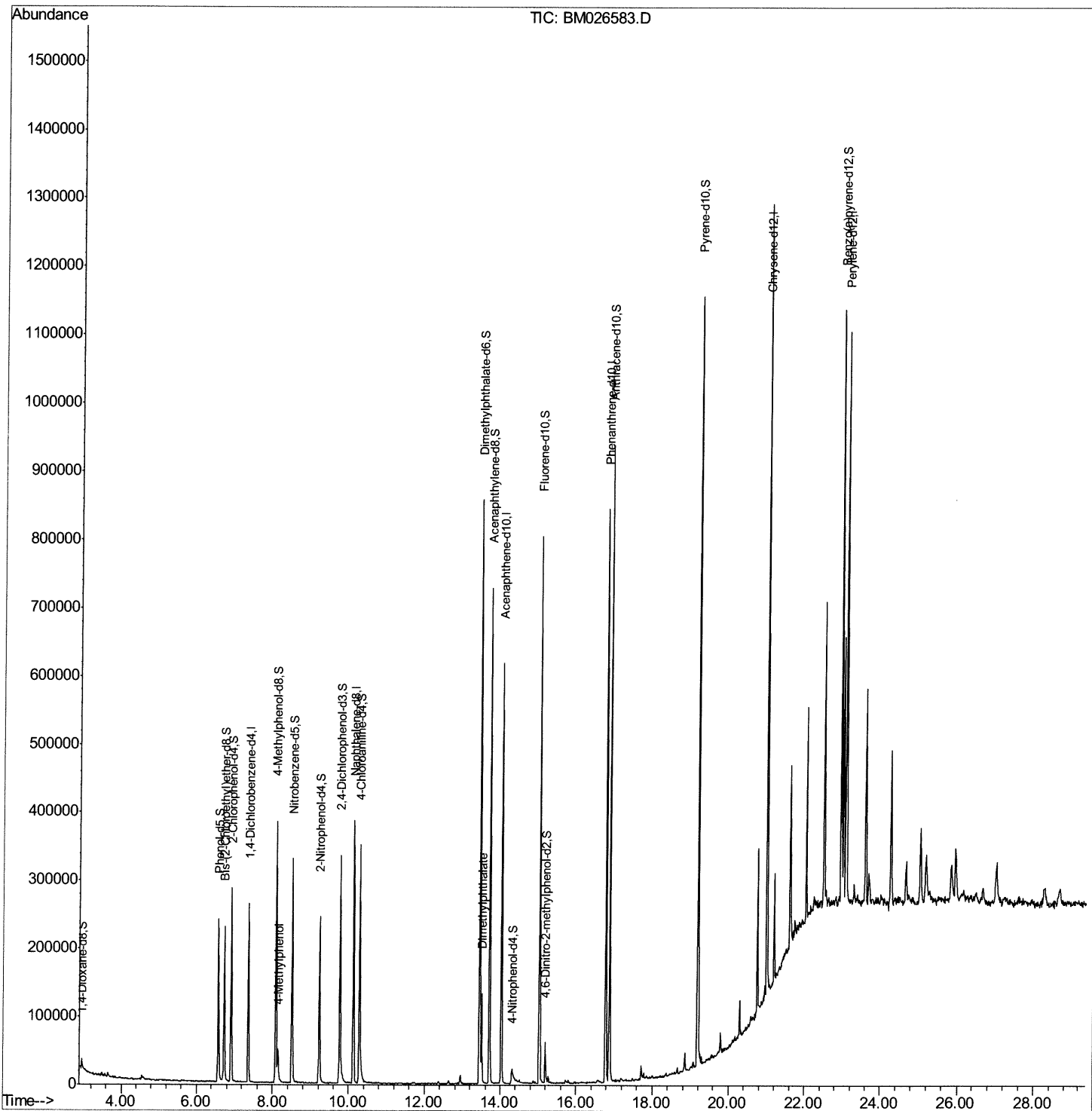
Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM062620\
Data File : BM026583.D
Acq On : 26 Jun 2020 11:36
Operator : JU/CG
Sample : L3062-17
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
ET150

Manual Integrations
APPROVED

mohammad
6/29/2020 3:17:23 PM

Quant Time: Jun 26 12:22:14 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061220MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Jun 26 11:31:11 2020
Response via : Initial Calibration



Quantitation Report (Qedit)

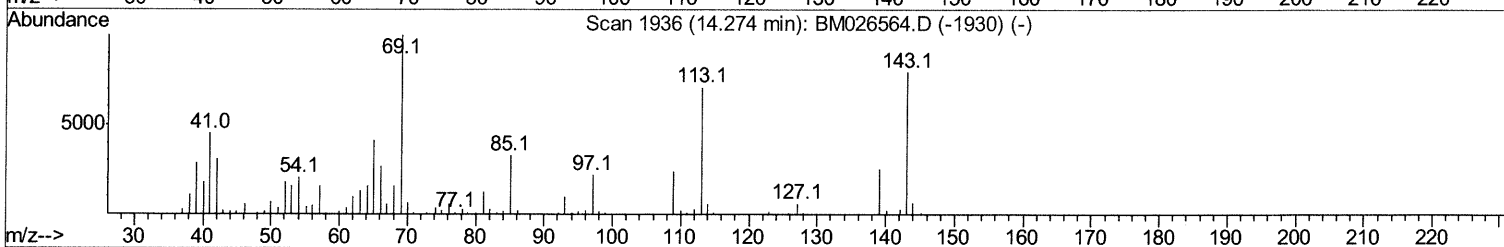
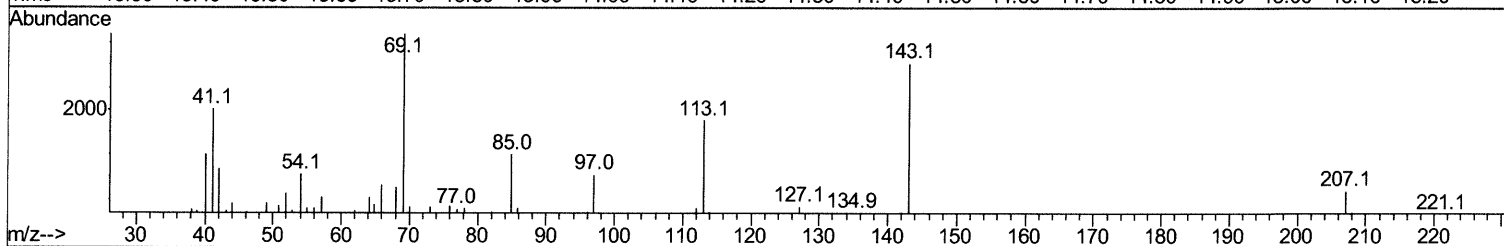
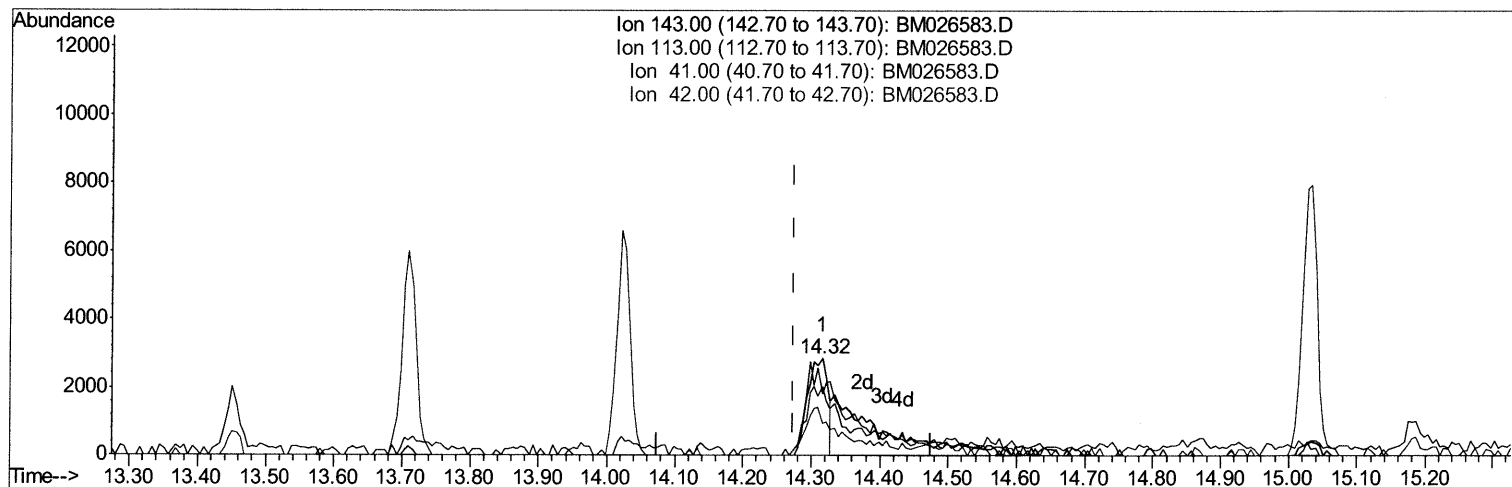
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TIC: BM026583.D

(52) 4-Nitrophenol-d4 (S)

14.316min (+0.041) 2.23ng/ul

response 5919

Ion	Exp%	Act%
143.00	100	100
113.00	81.70	63.96#
41.00	49.50	70.95#
42.00	34.20	33.36

Quantitation Report (Qedit)

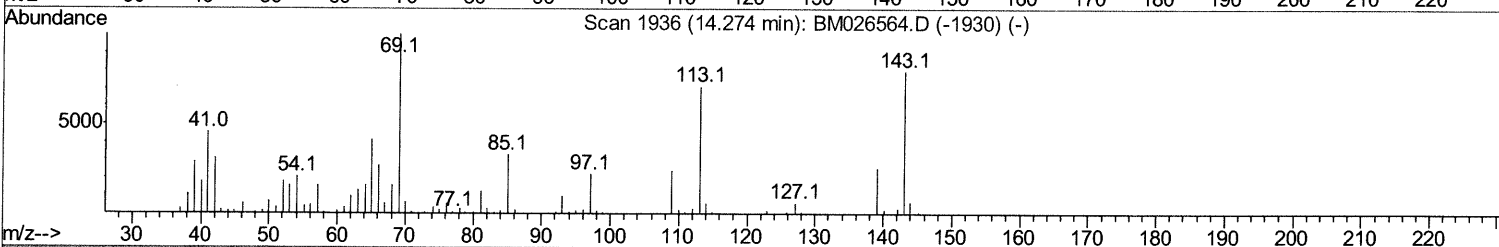
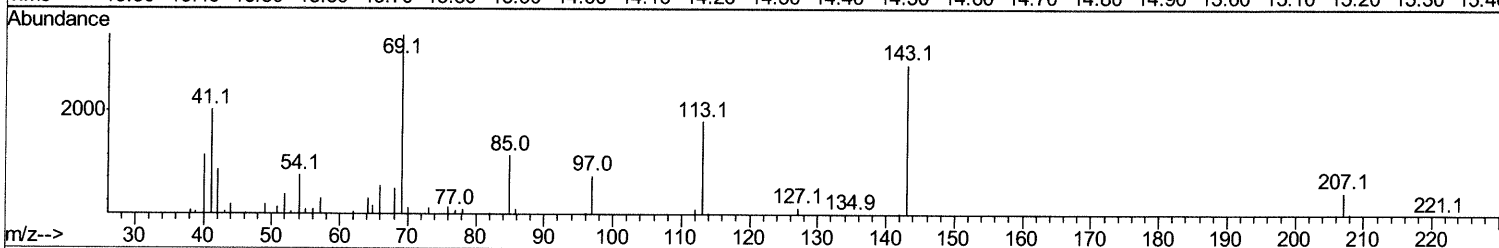
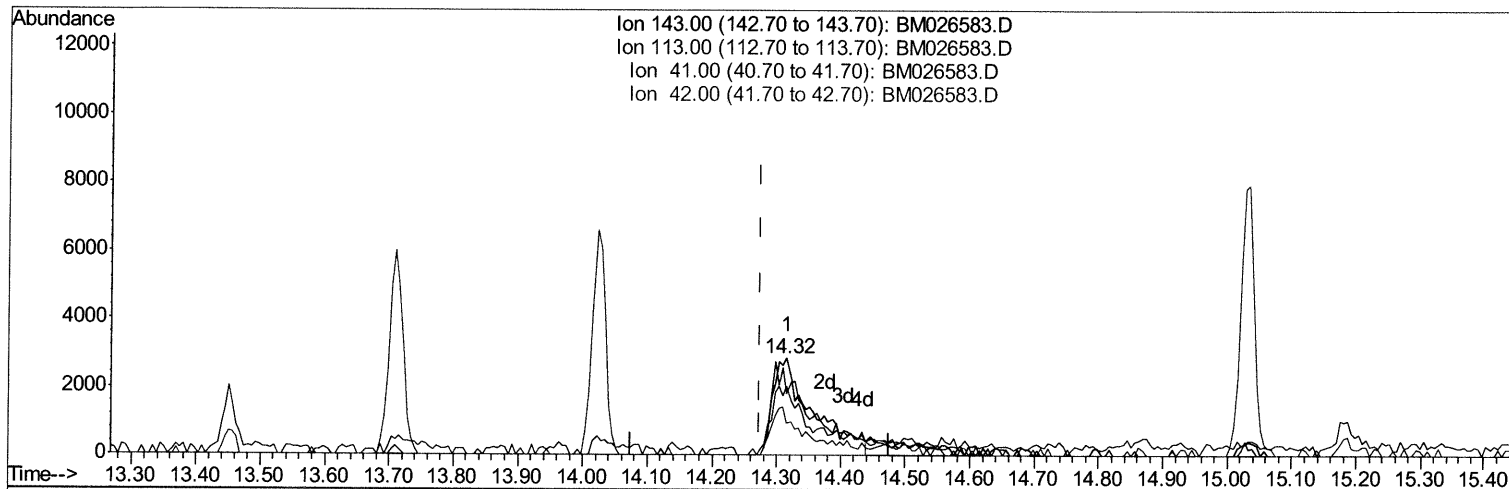
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Manual Integrations
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 Response via : Initial Calibration



TIC: BM026583.D

(52) 4-Nitrophenol-d4 (S)

14.316min (+0.041) 4.62ng/ul m 06/30/20 JU

response 12252

Ion	Exp%	Act%
143.00	100	100
113.00	81.70	63.96#
41.00	49.50	70.95#
42.00	34.20	33.36

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM062620\
 Data File : BM026583.D
 Acq On : 26 Jun 2020 11:36
 Operator : JU/CG
 Sample : L3062-17
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 ET150

Manual Integrations
 APPROVED

mohammad
 6/29/2020 3:17:23 PM

Quant Time: Jun 26 12:22:14 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061220MA.M
 Quant Title : SVOA CALIBRATION
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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.37	152	60543	20.00	ng/ul	0.00
18) Naphthalene-d8	10.12	136	277203	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.03	164	190165	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.78	188	435291	20.00	ng/ul	0.00
78) Chrysene-d12	21.00	240	511458	20.00	ng/ul	0.00
86) Perylene-d12	23.08	264	504683	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	2.96	96	6671	3.86	ng/uL	0.00
5) Phenol-d5	6.57	99	126356	22.66	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.73	67	109660	31.08	ng/ul	0.00
9) 2-Chlorophenol-d4	6.91	132	108135	23.78	ng/ul	0.00
13) 4-Methylphenol-d8	8.09	113	130762	29.79	ng/ul	0.00
19) Nitrobenzene-d5	8.52	128	63126	25.63	ng/ul	0.00
22) 2-Nitrophenol-d4	9.23	143	61298	22.62	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.77	165	117279	23.51	ng/ul	0.00
29) 4-Chloroaniline-d4	10.28	131	185775	35.71	ng/ul	0.00
44) Dimethylphthalate-d6	13.45	166	511693	32.00	ng/ul	0.00
47) Acenaphthylene-d8	13.71	160	444043	23.15	ng/ul	0.00
52) 4-Nitrophenol-d4	14.32	143	12252m>	4.62	ng/ul>	0.04
58) Fluorene-d10	15.03	176	302443	22.04	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.19	200	13120	4.61	ng/ul	0.00
71) Anthracene-d10	16.88	188	478242	21.58	ng/ul	0.00
79) Pyrene-d10	19.19	212	548814	19.57	ng/ul	0.00
90) Benzo(a)pyrene-d12	22.95	264	525811	18.84	ng/ul	0.00

06/30/20 JU

Target Compounds

					Ovalue
16) 4-Methylphenol	8.15	108	15205	3.163	ng/ul 82
45) Dimethylphthalate	13.50	163	77424	4.663	ng/ul 99

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