

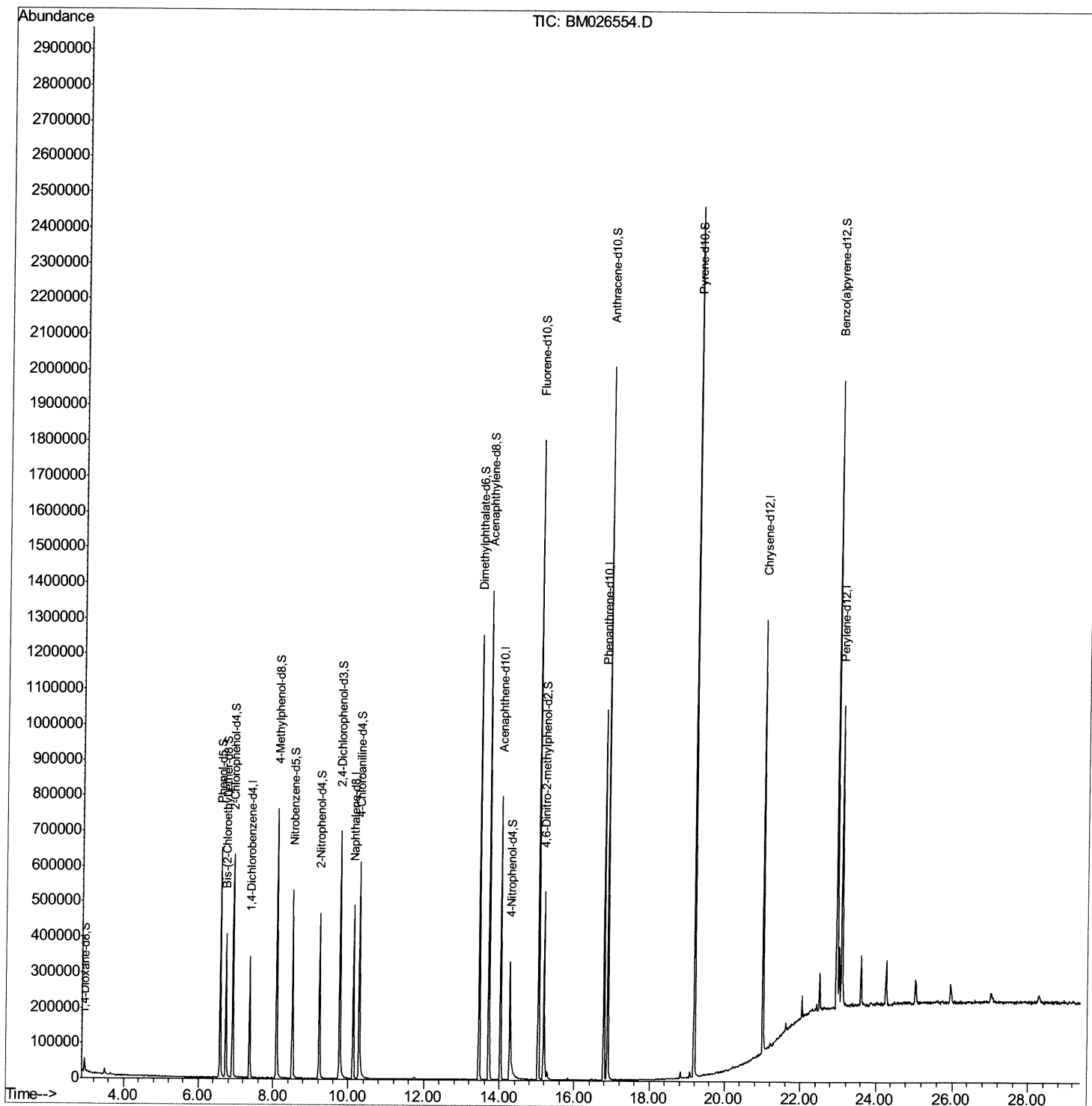
Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM062520\
Data File : BM026554.D
Acq On : 25 Jun 2020 10:56
Operator : JU/CG
Sample : PB129704BL
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
SBLK04

Manual Integrations
APPROVED

mohammad
6/26/2020 12:19:44 PM

Quant Time: Jun 25 12:25:42 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)

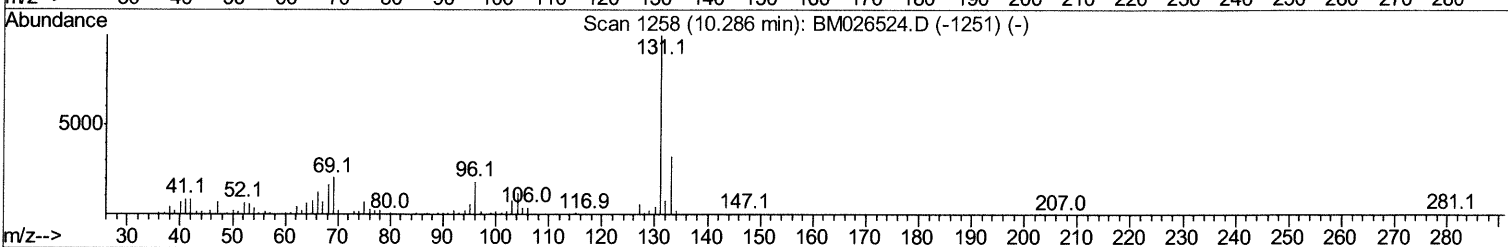
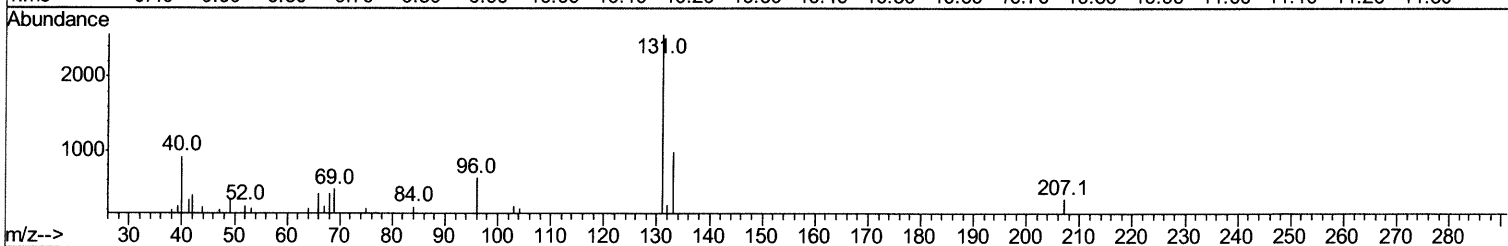
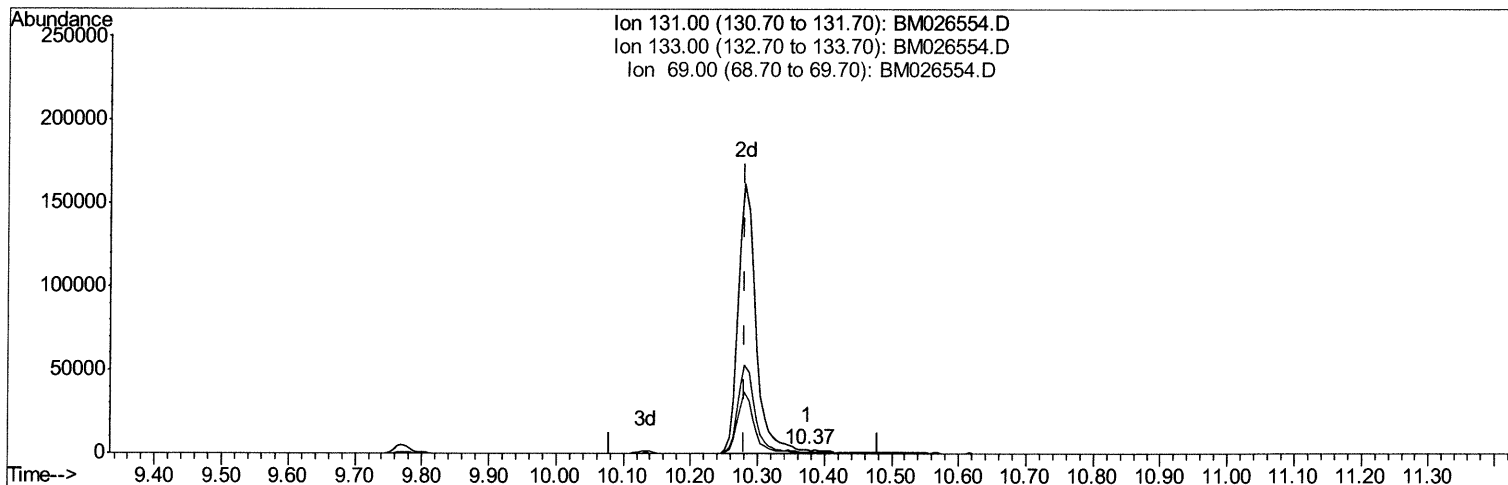
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Quant Time: Jun 25 12:01:51 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061220MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jun 25 10:53:13 2020
 Response via : Initial Calibration



TIC: BM026554.D

(29) 4-Chloroaniline-d4 (S)

10.375min (+0.094) 0.52ng/ul

response 3319

Ion	Exp%	Act%
131.00	100	100
133.00	32.50	38.08
69.00	18.60	18.55
0.00	0.00	0.00

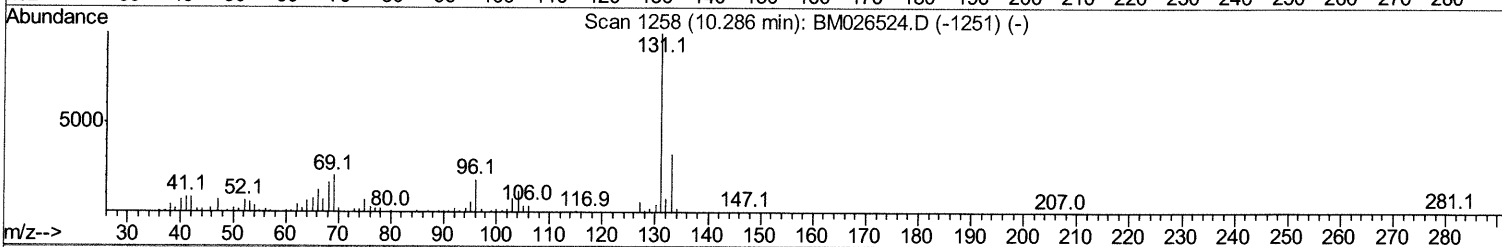
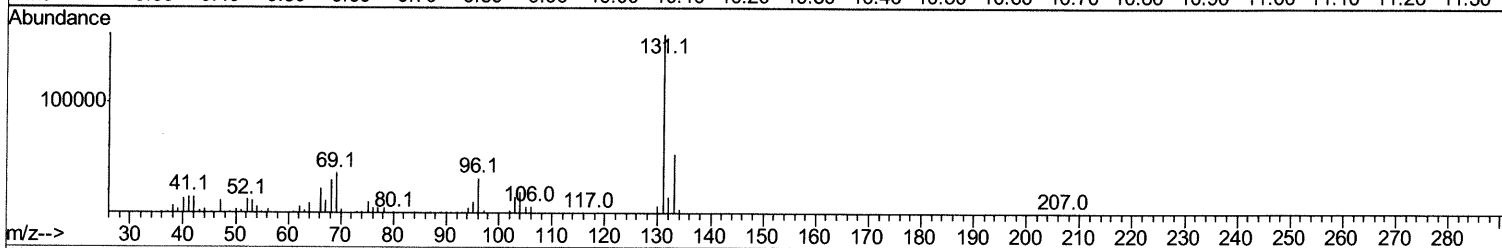
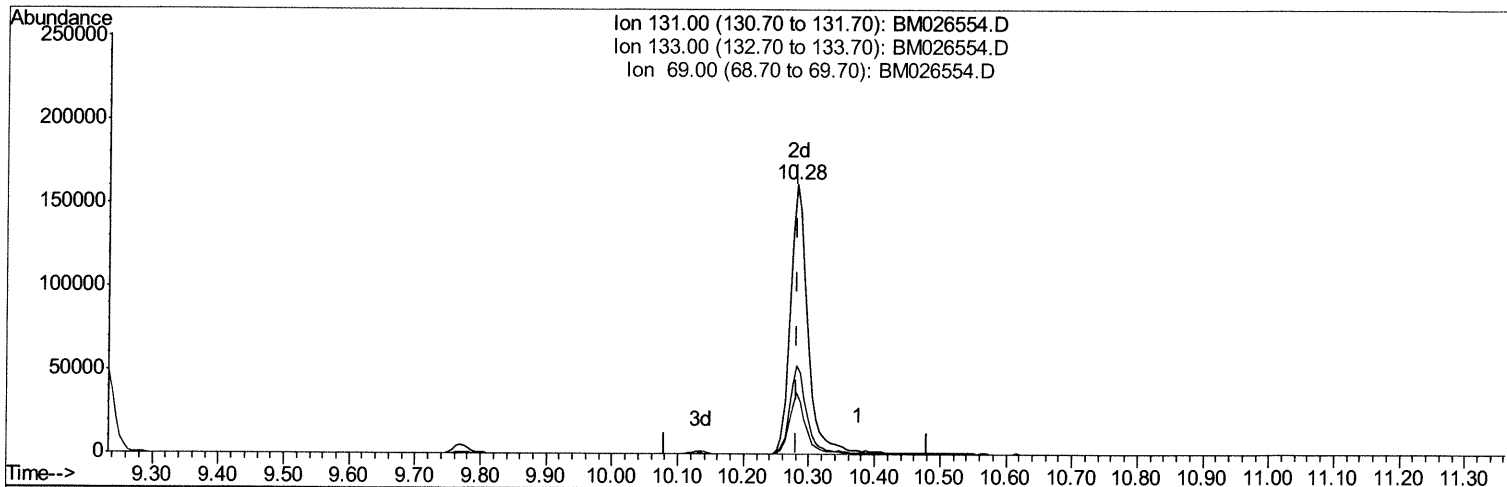
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Manual Integrations
APPROVED

mohammad
6/26/2020 12:19:44 PM

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Quant Title : SVOA CALIBRATION
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Response via : Initial Calibration



TIC: BM026554.D

(29) 4-Chloroaniline-d4 (S)

10.281min (-0.000) 46.51ng/ul m 06/29/2020

response 298078

Ion	Exp%	Act%
131.00	100	100
133.00	32.50	32.82
69.00	18.60	22.53#
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM062520\
 Data File : BM026554.D
 Acq On : 25 Jun 2020 10:56
 Operator : JU/CG
 Sample : PB129704BL
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK04

Manual Integrations
 APPROVED

mohammad
 6/26/2020 12:19:44 PM

Quant Time: Jun 25 12:25:42 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	80410	20.00	ng/ul	0.00
18) Naphthalene-d8	10.13	136	341506	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.03	164	232703	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.79	188	543035	20.00	ng/ul	0.00
78) Chrysene-d12	21.00	240	578228	20.00	ng/ul	0.00
86) Perylene-d12	23.09	264	543977	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	2.96	96	18672	8.14	ng/uL	0.00
5) Phenol-d5	6.58	99	306424	41.37	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.73	67	182373	38.92	ng/ul	0.00
9) 2-Chlorophenol-d4	6.92	132	227022	37.60	ng/ul	0.00
13) 4-Methylphenol-d8	8.10	113	246036	42.20	ng/ul	0.00
19) Nitrobenzene-d5	8.52	128	102150	33.67	ng/ul	0.00
22) 2-Nitrophenol-d4	9.23	143	115463	34.58	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.77	165	231372	37.64	ng/ul	0.00
29) 4-Chloroaniline-d4	10.28	131	298078m	46.51	ng/ul	0.00
44) Dimethylphthalate-d6	13.46	166	723810	36.99	ng/ul	0.00
47) Acenaphthylene-d8	13.72	160	826067	35.20	ng/ul	0.00
52) 4-Nitrophenol-d4	14.27	143	87784	27.05	ng/ul	0.00
58) Fluorene-d10	15.03	176	617410	36.77	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.18	200	94451	26.58	ng/ul	0.00
71) Anthracene-d10	16.89	188	996429	36.05	ng/ul	0.00
79) Pyrene-d10	19.19	212	1176793	37.11	ng/ul	0.00
90) Benzo(a)pyrene-d12	22.96	264	1115744	37.09	ng/ul	0.00

06/29/20 JU

Target Compounds Ovalue

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