

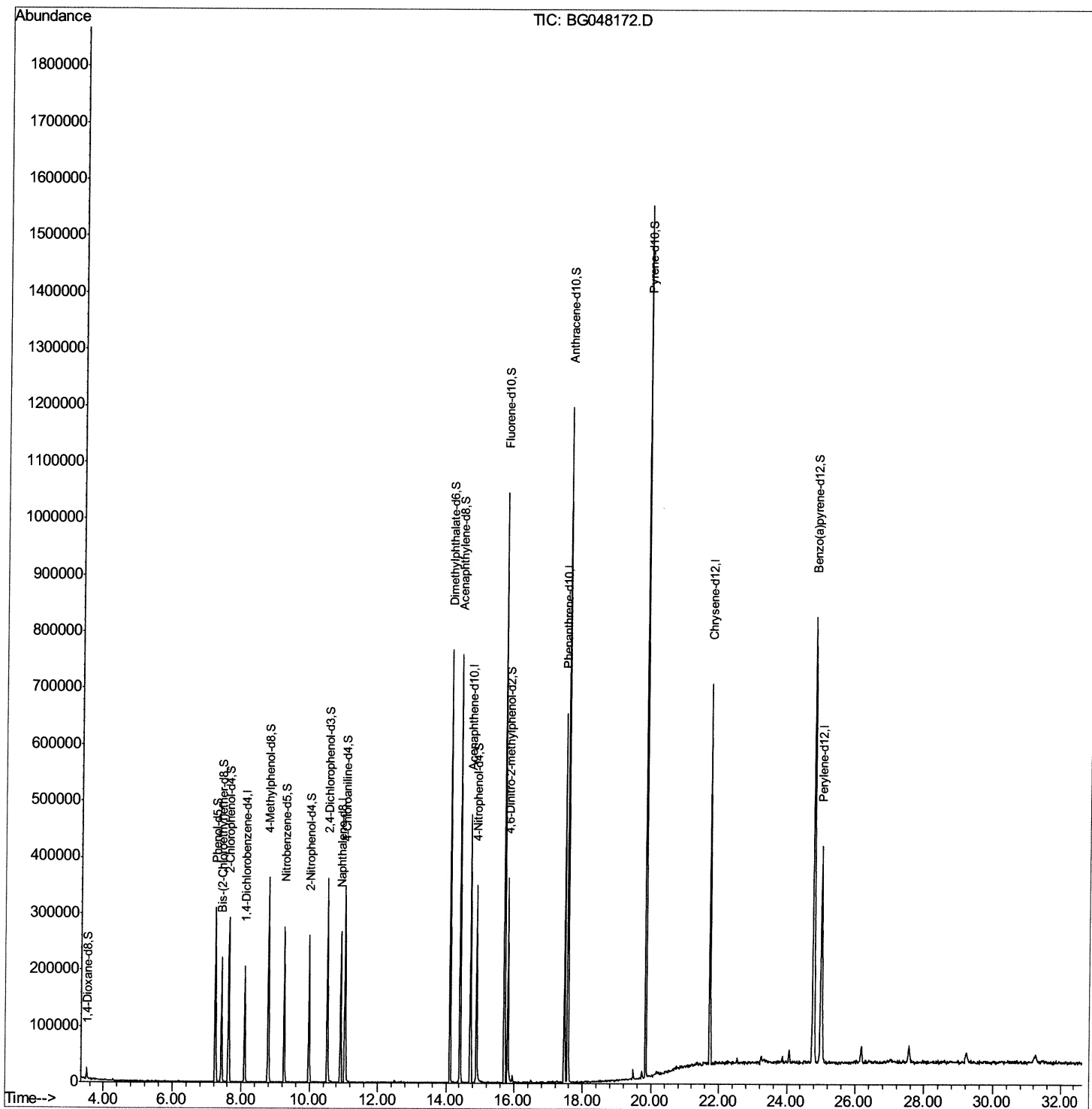
Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021621\
Data File : BG048172.D
Acq On : 16 Feb 2021 21:00
Operator : CG/JU
Sample : PB134674BL
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SBLK74

Manual Integrations
APPROVED

mohammad
2/17/2021 1:46:19 PM

Quant Time: Feb 17 01:08:15 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG021321MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Feb 16 23:25:09 2021
Response via : Initial Calibration



Quantitation Report (Qedit)

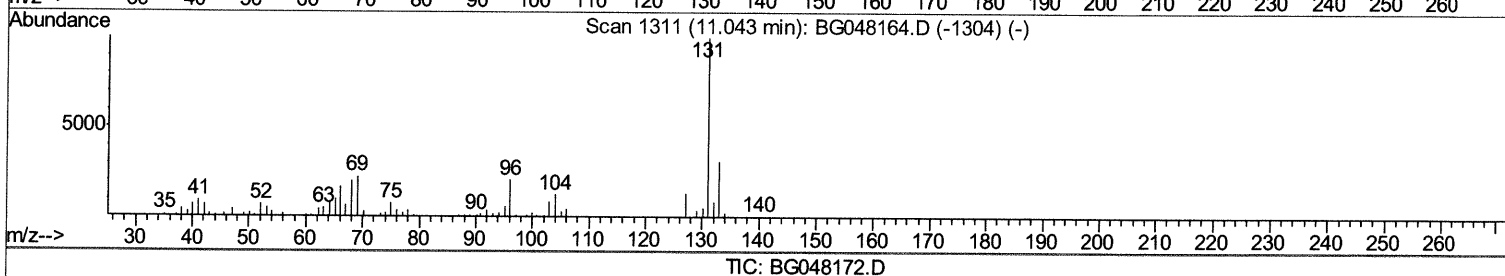
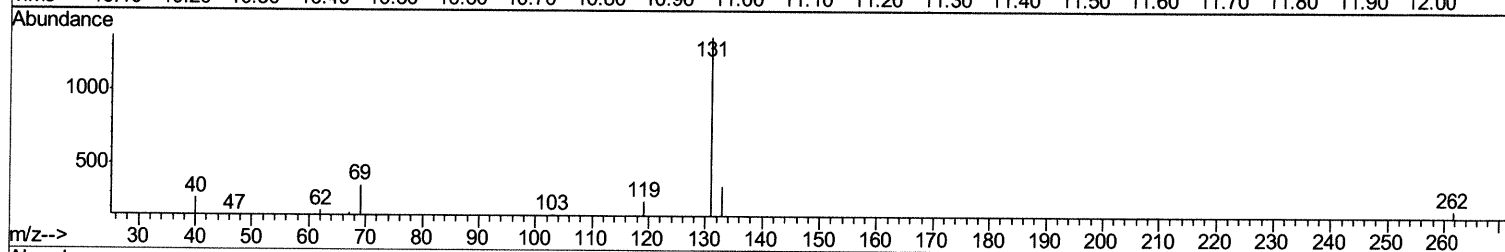
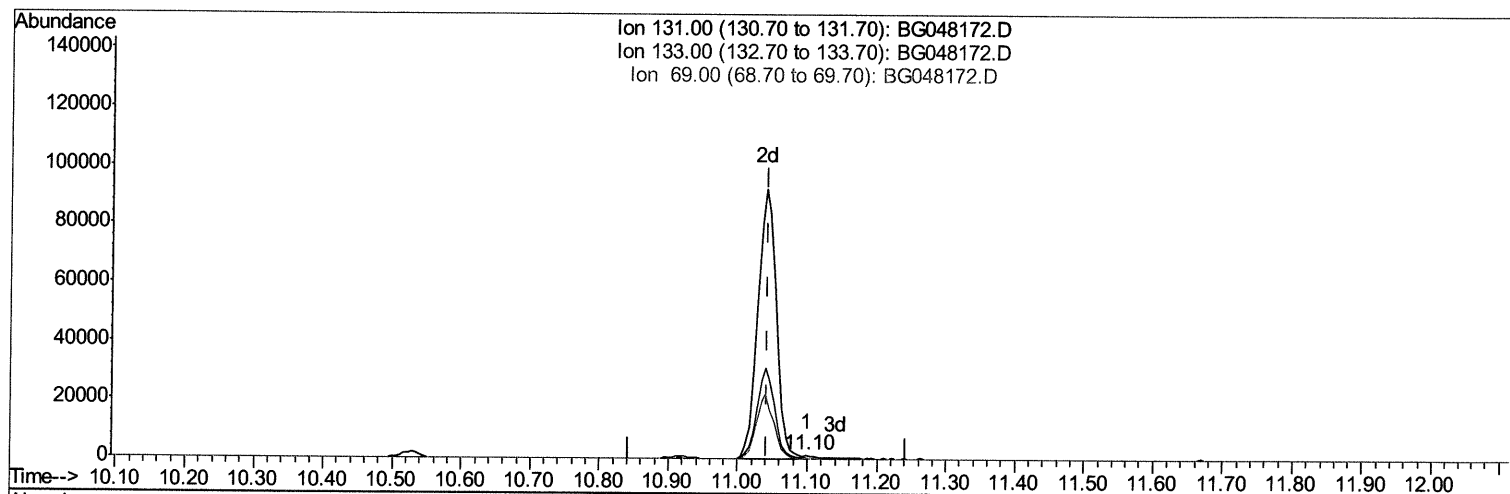
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(29) 4-Chloroaniline-d4 (S)

11.099min (+0.056) 0.26ng/ul

response 973

Ion	Exp%	Act%
131.00	100	100
133.00	24.60	25.88
69.00	23.70	25.88
0.00	0.00	0.00

Quantitation Report (Qedit)

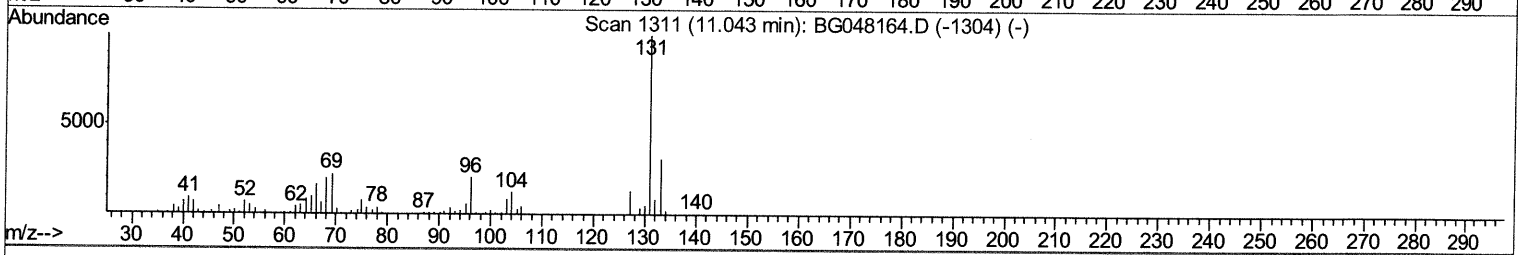
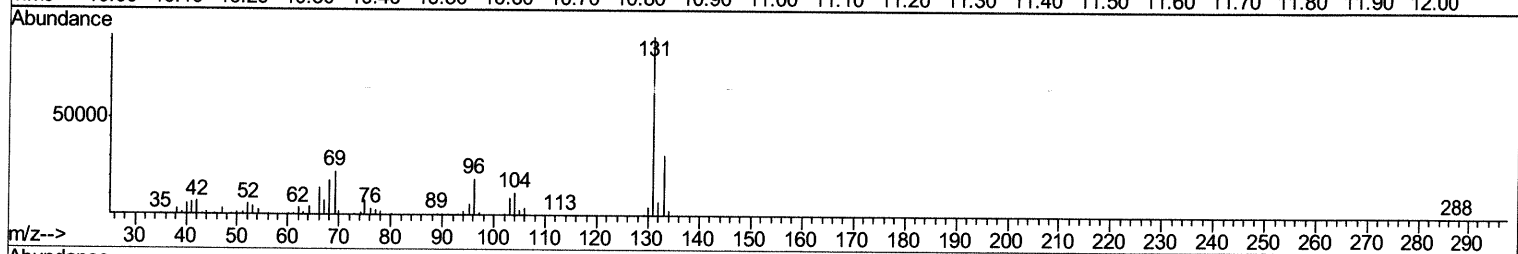
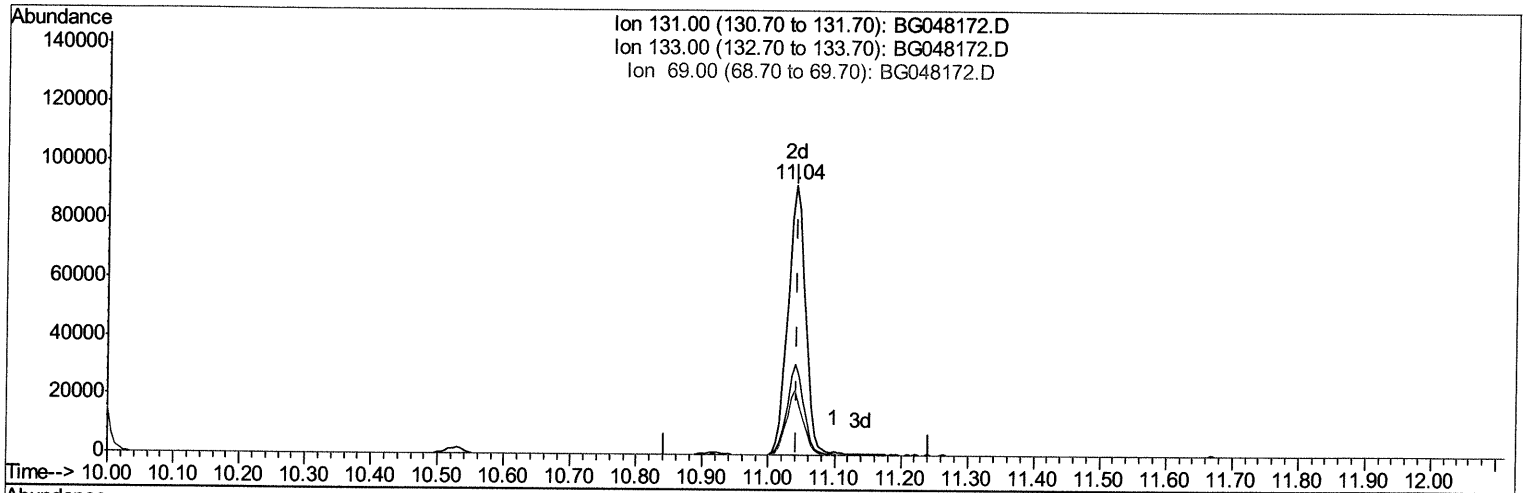
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Manual Integrations
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Response via : Initial Calibration



TIC: BG048172.D

(29) 4-Chloroaniline-d4 (S)

11.040min (-0.003) 46.44ng/ul m 02/18/21 JU

response 171599

Ion	Exp%	Act%
131.00	100	100
133.00	24.60	33.62#
69.00	23.70	24.19
0.00	0.00	0.00

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021621\
 Data File : BG048172.D
 Acq On : 16 Feb 2021 21:00
 Operator : CG/JU
 Sample : PB134674BL
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK74

Manual Integrations
 APPROVED

mohammad
 2/17/2021 1:46:19 PM

Quant Time: Feb 17 01:08:15 2021
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 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Feb 16 23:25:09 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.10	152	54559	20.00	ng/ul	0.00
18) Naphthalene-d8	10.91	136	211327	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.72	164	162195	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.47	188	404950	20.00	ng/ul	0.00
78) Chrysene-d12	21.74	240	437770	20.00	ng/ul	0.00
86) Perylene-d12	25.00	264	429196	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.51	96	10265	7.83	ng/uL	0.00
5) Phenol-d5	7.24	99	164074	35.69	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.42	67	104833	38.29	ng/ul	0.00
9) 2-Chlorophenol-d4	7.63	132	120727	37.41	ng/ul	0.00
13) 4-Methylphenol-d8	8.80	113	130808	36.27	ng/ul	0.00
19) Nitrobenzene-d5	9.26	128	62503	39.60	ng/ul	0.00
22) 2-Nitrophenol-d4	9.99	143	73713	38.61	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.52	165	132024	35.30	ng/ul	0.00
29) 4-Chloroaniline-d4	11.04	131	171599m>	46.44	ng/ul>	0.00
44) Dimethylphthalate-d6	14.12	166	485892	40.44	ng/ul	0.00
47) Acenaphthylene-d8	14.42	160	551169	37.53	ng/ul	0.00
52) 4-Nitrophenol-d4	14.90	143	78810	38.61	ng/ul	0.00
58) Fluorene-d10	15.72	176	423380	39.51	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.82	200	87379	35.19	ng/ul	0.00
71) Anthracene-d10	17.57	188	705017	40.61	ng/ul	0.00
79) Pyrene-d10	19.85	212	872692	39.92	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.78	264	909312	41.53	ng/ul	0.00

02/18/21JU

Target Compounds Ovalue

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