

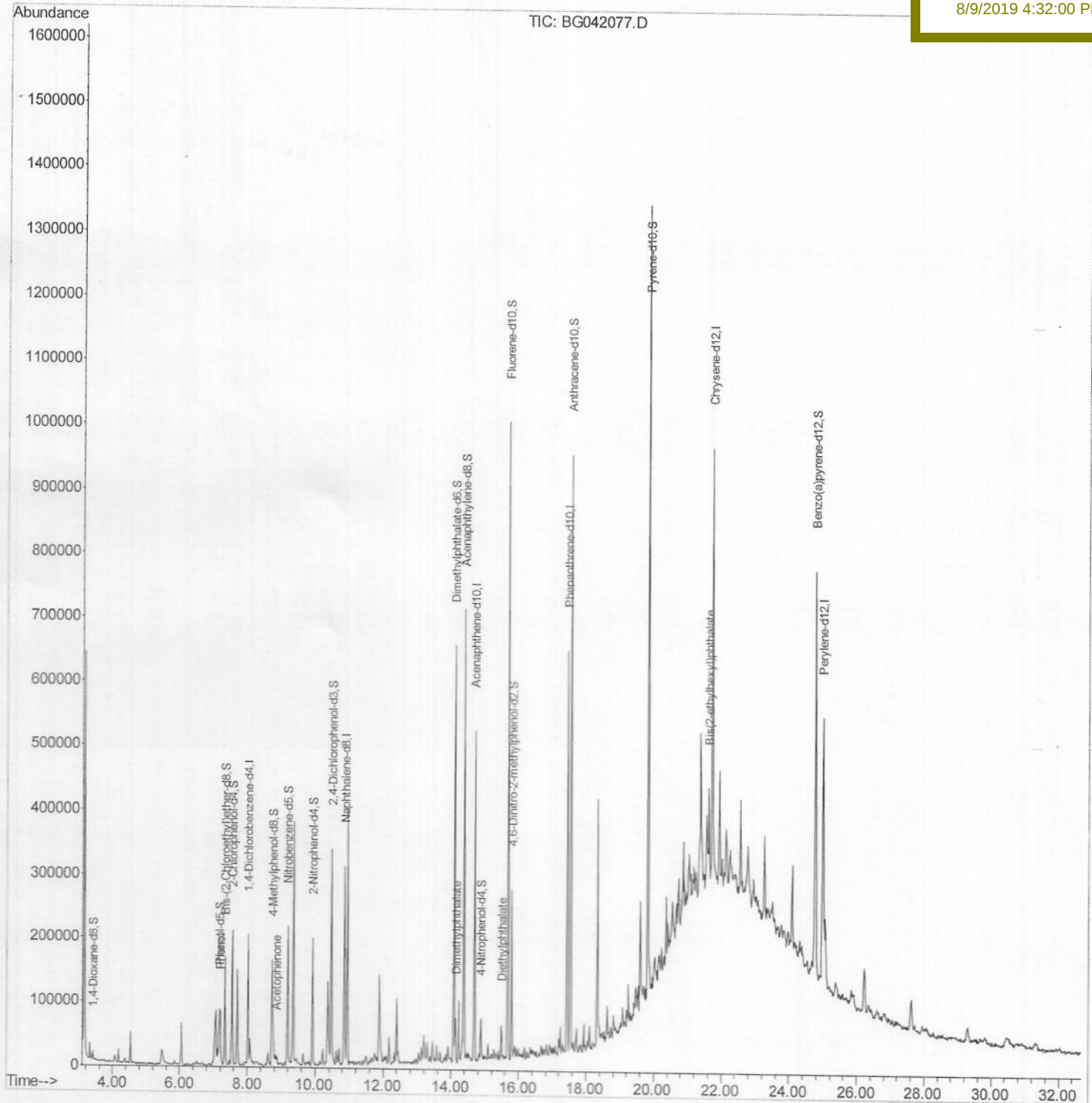
Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG080819\
Data File : BG042077.D
Acq On : 8 Aug 2019 21:21
Operator : HP/JU
Sample : K4117-13
Misc :
ALS Vial : 7 Sample Multiplier: 1

Quant Time: Aug 09 09:14:11 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG080319MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Aug 08 15:44:01 2019
Response via : Initial Calibration

Instrument :
BNA_G
ClientSampleId :
JLLH4

Manual Integrations
APPROVED

mohammad
8/9/2019 4:32:00 PM



Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG080819\

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Quant Title : SVOA CALIBRATION

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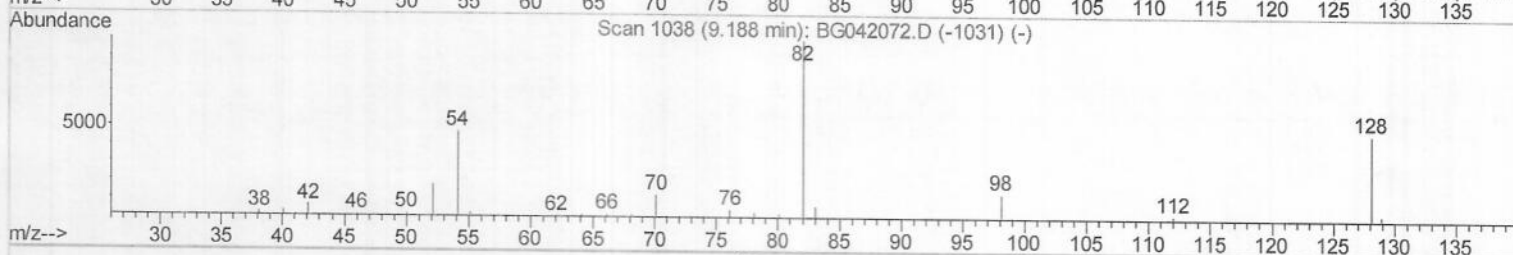
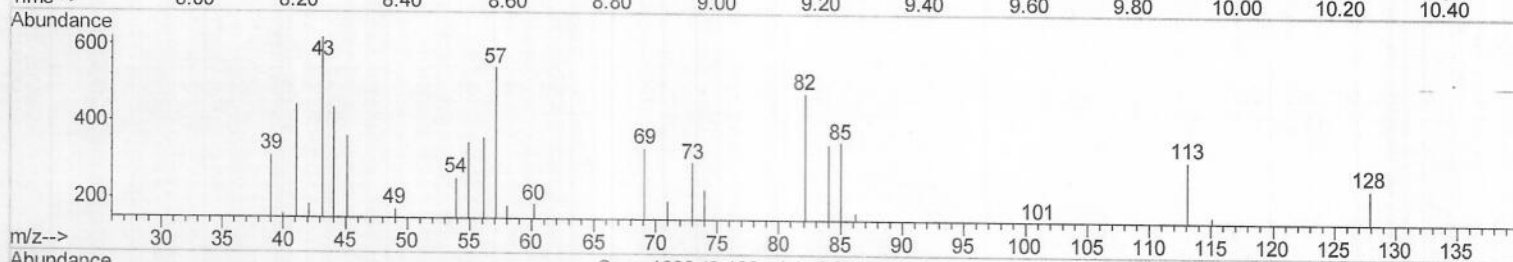
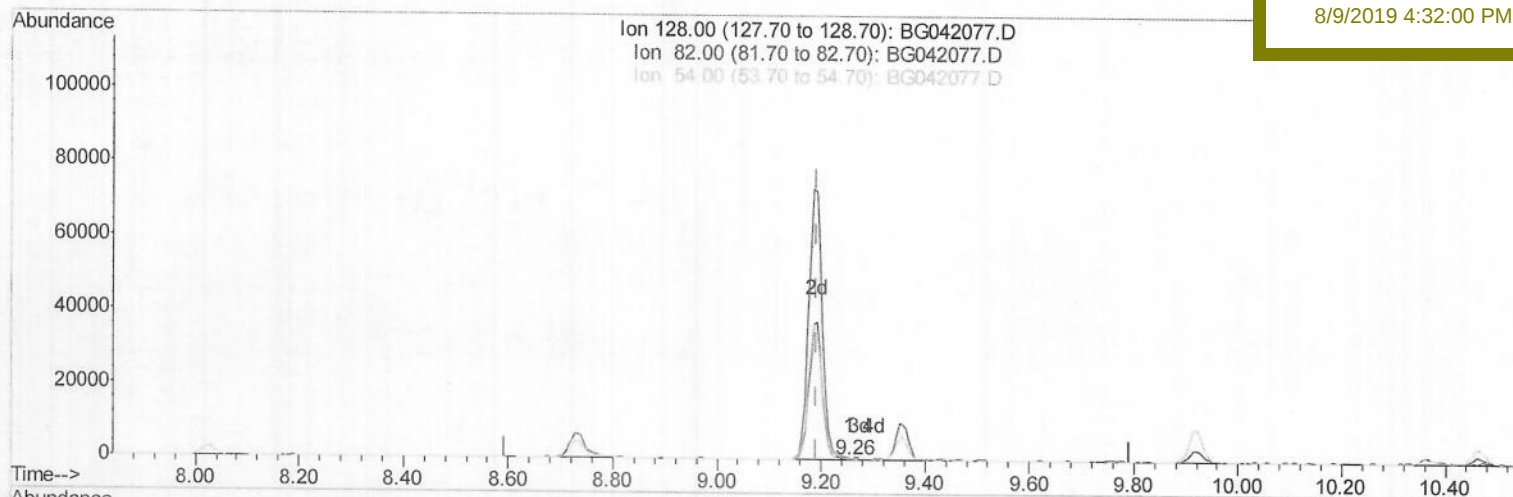
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Manual Integrations
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TIC: BG042077.D

(19) Nitrobenzene-d5 (S)

9.256min (+0.063) 0.06ng/ul

response 137

Ion	Exp%	Act%
128.00	100	100
82.00	223.40	202.53
54.00	125.40	107.17
0.00	0.00	0.00

Quantitation Report (Qedit)

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BNA_G

ClientSampleId :

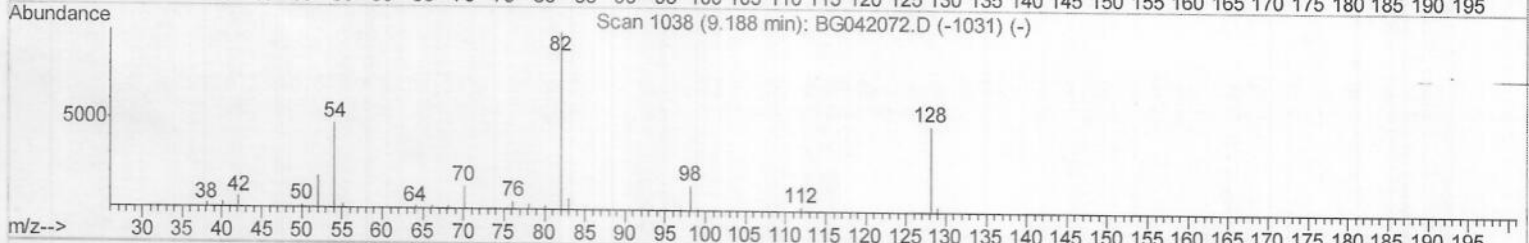
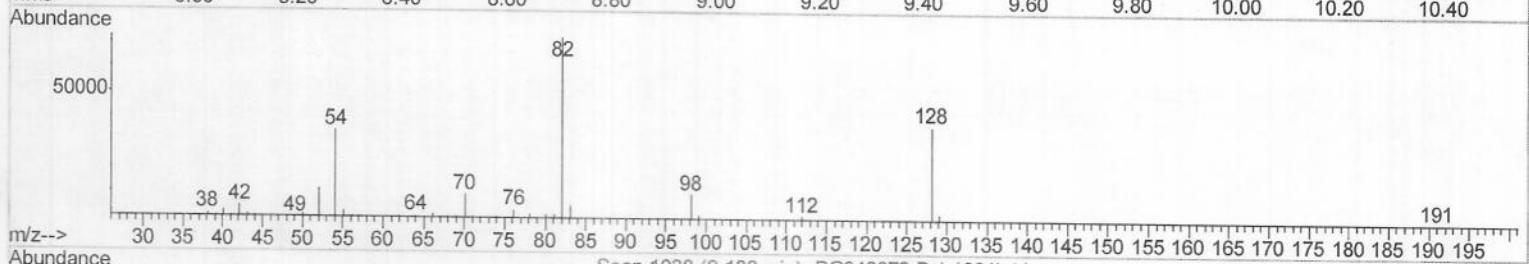
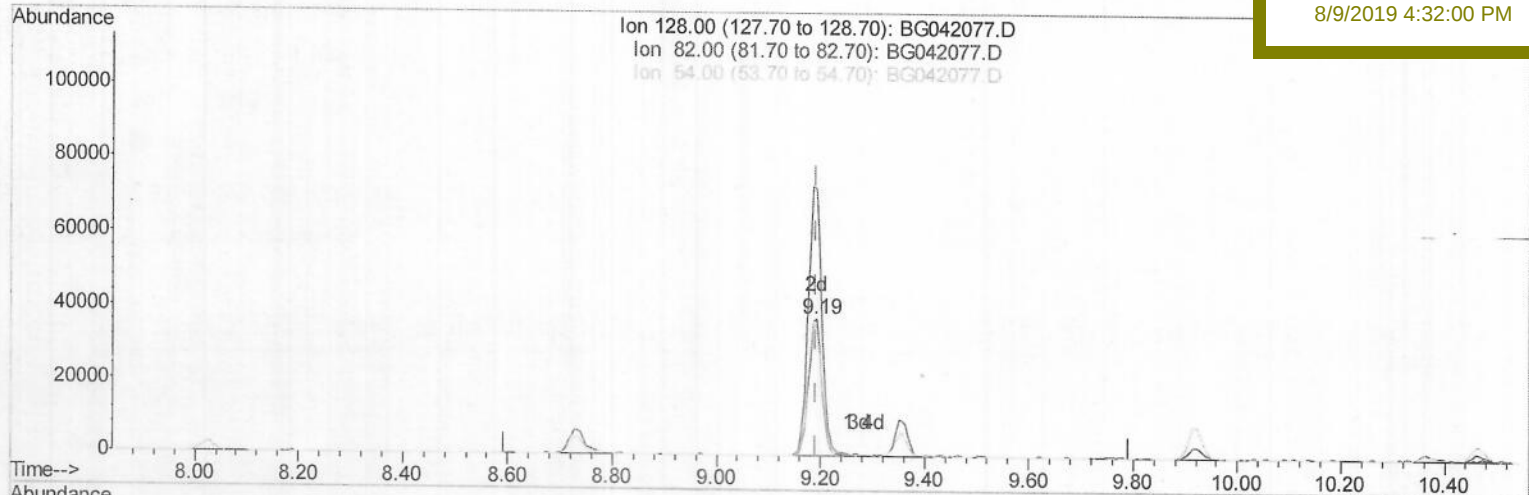
JLLH4

Manual Integrations

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TIC: BG042077.D

(19) Nitrobenzene-d5 (S)

9.192min (-0.001) 31.77ng/ul m) *JU*
08/12/19

response 67473

Ion	Exp%	Act%
128.00	100	100
82.00	223.40	194.09
54.00	125.40	92.72#
0.00	0.00	0.00

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 Operator : HP/JU
 Sample : K4117-13
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Quant Time: Aug 09 09:14:11 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG080319MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Aug 08 15:44:01 2019
 Response via : Initial Calibration

Instrument :
 BNA_G
 Client Sampled :
 JLLH4

Manual Integrations
 APPROVED

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 8/9/2019 4:32:00 PM

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.03	152	67091	20.00	ng/ul	0.00
18) Naphthalene-d8	10.85	136	282124	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.69	164	198216	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.43	188	420901	20.00	ng/ul	0.00
77) Chrysene-d12	21.72	240	407155	20.00	ng/ul	0.00
85) Perylene-d12	24.98	264	474998	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.41	96	7076	4.61	ng/uL	0.00
5) Phenol-d5	7.18	99	47087	8.94	ng/ul	0.00
7) Bis-(2-Chloroethyl) ether-d	7.34	67	76307	25.91	ng/ul	0.00
9) 2-Chlorophenol-d4	7.55	132	117869	26.18	ng/ul	0.00
13) 4-Methylphenol-d8	8.73	113	74140	16.77	ng/ul	0.00
19) Nitrobenzene-d5	9.19	128	67473m)	31.77	ng/ul	0.00
22) 2-Nitrophenol-d4	9.92	143	82640	33.51	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.47	165	153710	30.88	ng/ul	0.00
29) 4-Chloroaniline-d4	0.00	131	0d	0.00	ng/ul	0.00
43) Dimethylphthalate-d6	14.09	166	498026	32.40	ng/ul	0.00
46) Acenaphthylene-d8	14.38	160	546456	31.12	ng/ul	0.00
51) 4-Nitrophenol-d4	14.87	143	24010	9.30	ng/ul	0.00
57) Fluorene-d10	15.68	176	431203	31.53	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.79	200	79639	29.49	ng/ul	0.00
70) Anthracene-d10	17.53	188	619593	30.68	ng/ul	0.00
78) Pyrene-d10	19.82	212	702768	30.93	ng/ul	0.00
89) Benzo(a)pyrene-d12	24.76	264	731125	29.46	ng/ul	0.00

Target Compounds

					Qvalue
6) Phenol	7.20	94	43566	8.016	ng/ul 97
14) Acetophenone	8.84	105	7302	1.074	ng/ul 93
44) Dimethylphthalate	14.13	163	42634	2.793	ng/ul 99
56) Diethylphthalate	15.50	149	21367	1.407	ng/ul 95
83) Bis(2-ethylhexyl)phthalate	21.61	149	46812	3.183	ng/ul 99

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