

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM120419\
Quantitation Report (QT Reviewed)

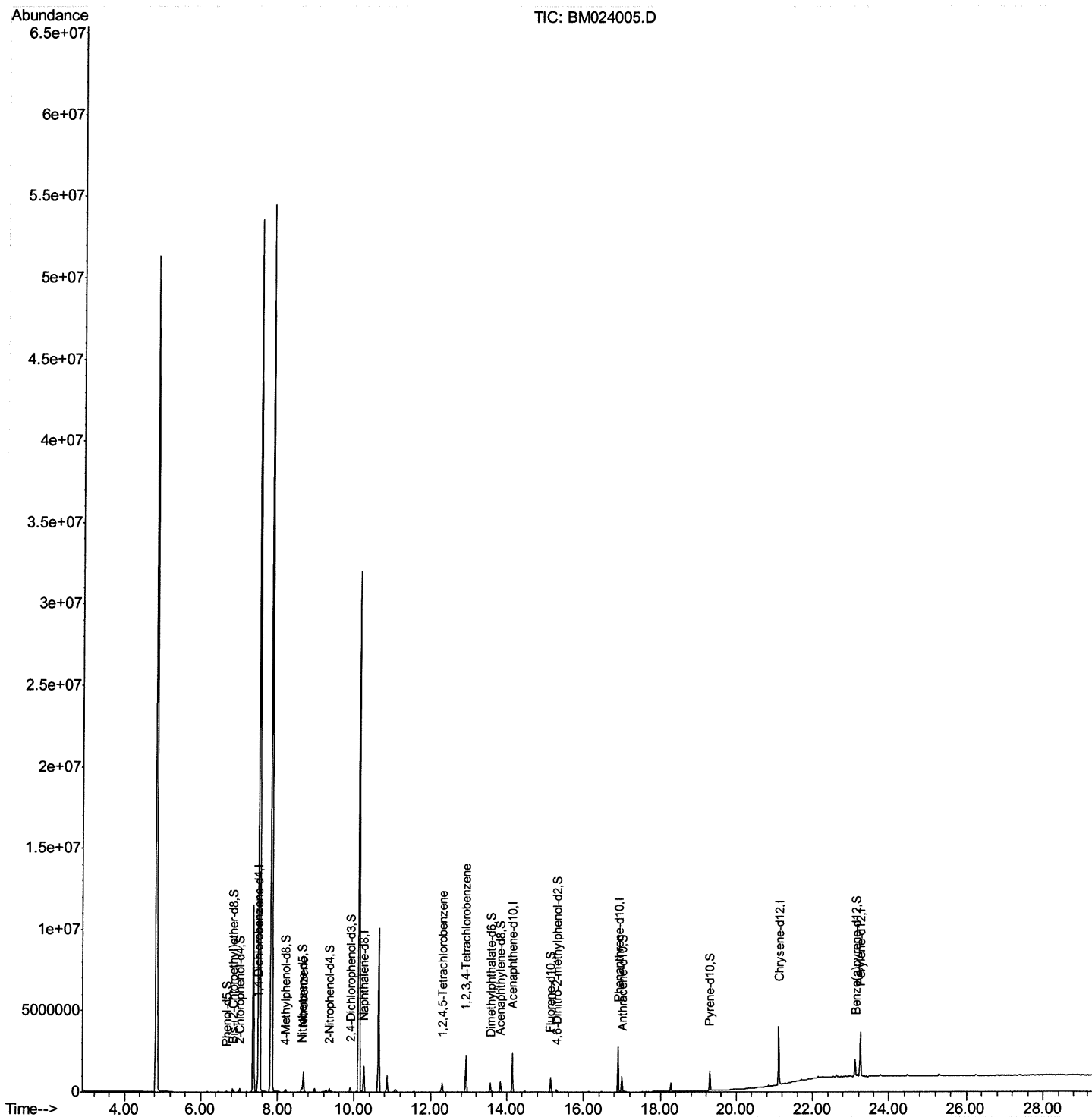
Data File : BM024005.D
Acq On : 04 Dec 2019 17:13
Operator : JU
Sample : K4649-06DL 5X
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
A5169DL

Manual Integrations
APPROVED

mohammad
12/5/2019 10:50:08 AM

Quant Time: Dec 05 03:02:42 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM112719MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Dec 05 02:46:26 2019
Response via : Initial Calibration



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM120419\
Quantitation Report (Qedit)

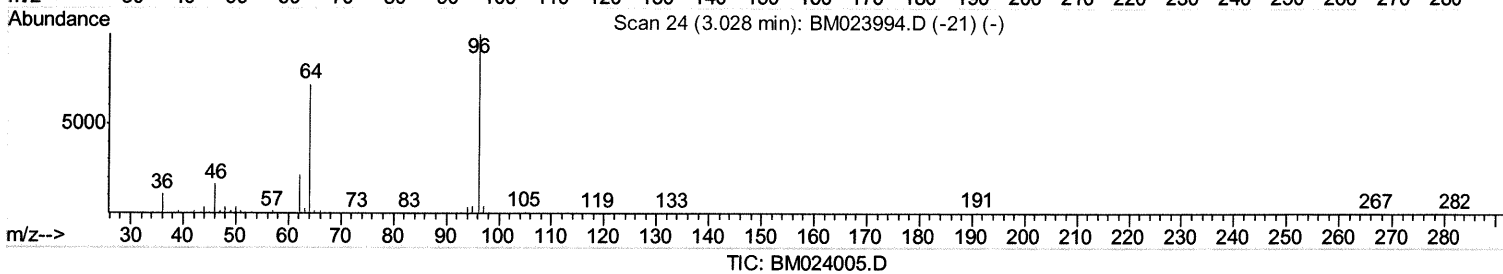
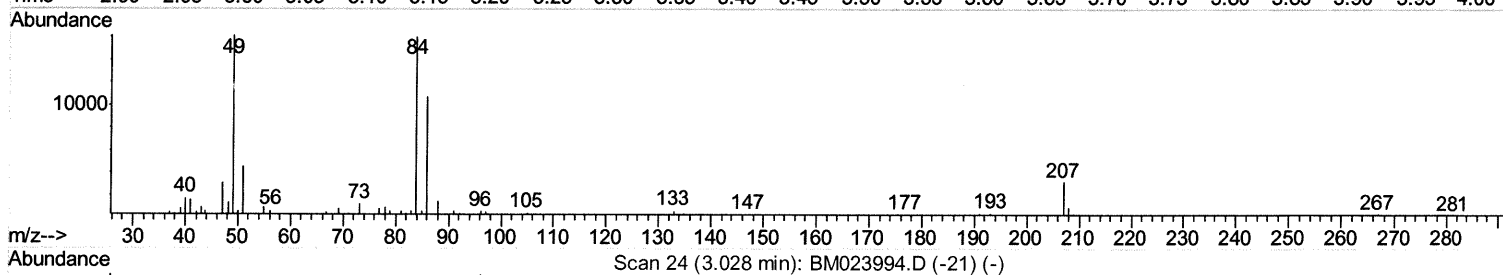
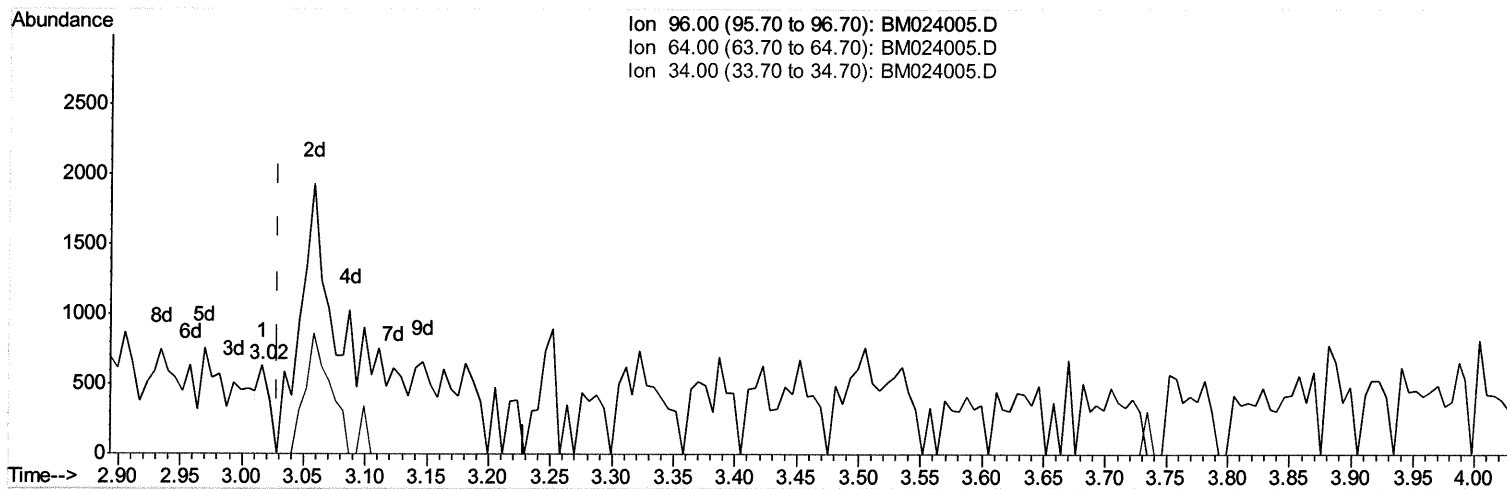
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(3) 1,4-Dioxane-d8 (S)

3.017min (-0.012) 0.05ng/uL

response 361

Ion	Exp%	Act%
96.00	100	100
64.00	71.10	0.00#
34.00	0.00	0.00
0.00	0.00	0.00

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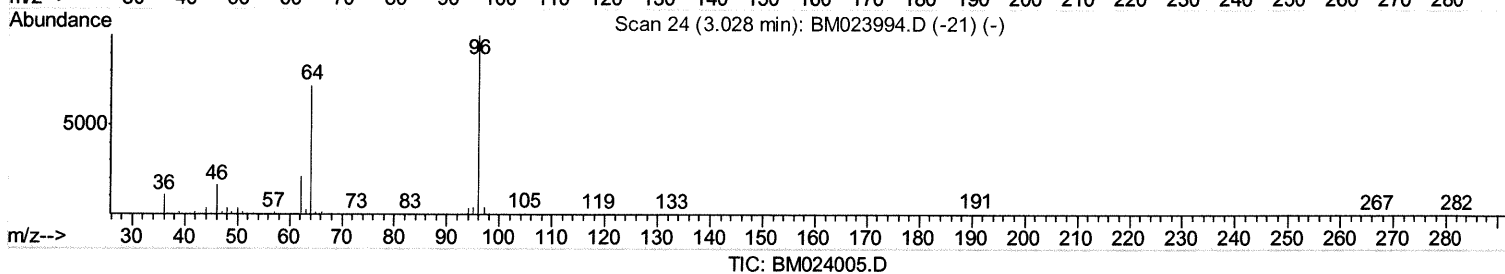
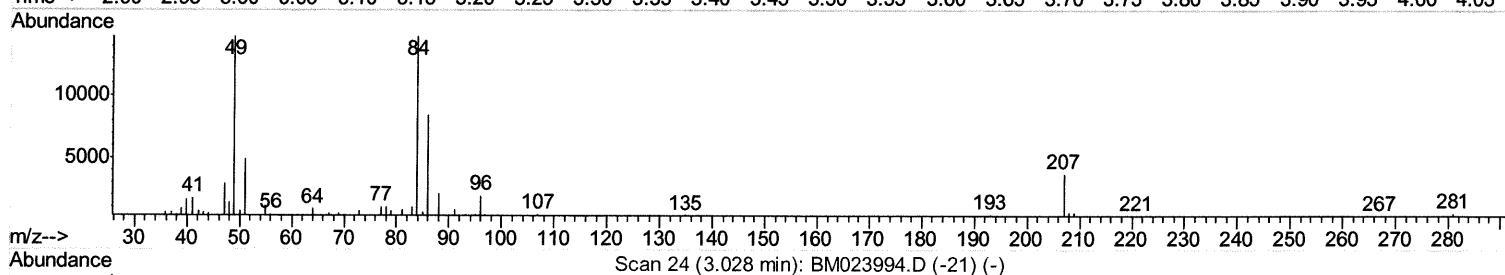
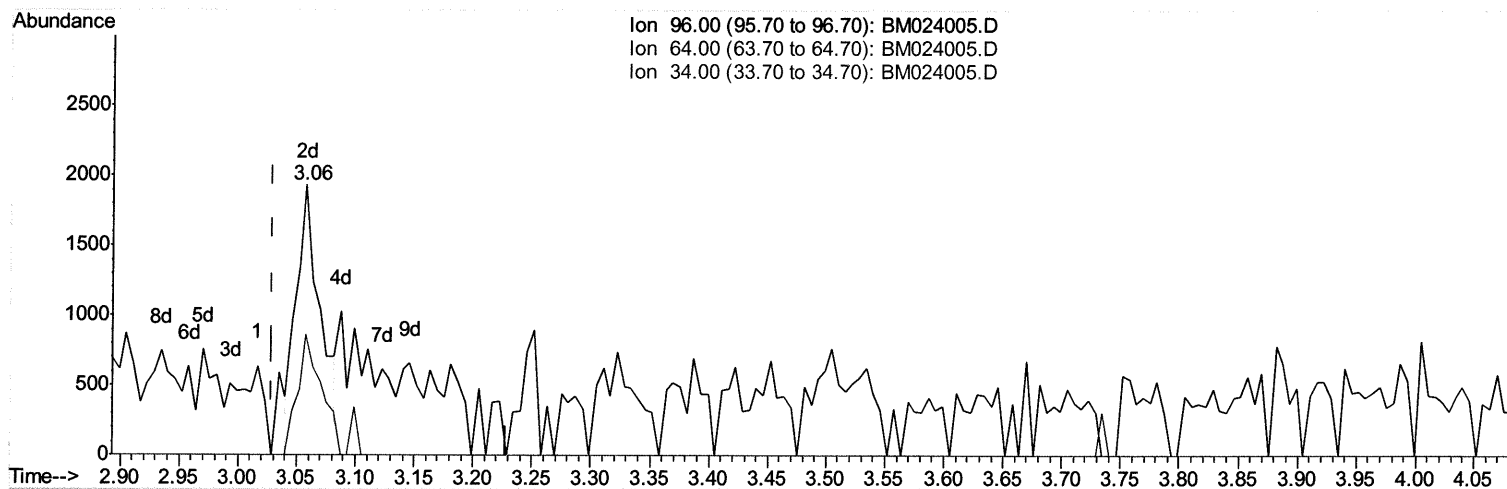
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Manual Integrations
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(3) 1,4-Dioxane-d8 (S)

3.058min (+0.029) 0.38ng/uL m JU 12/10/19

response 2793

Ion	Exp%	Act%
96.00	100	100
64.00	71.10	44.67#
34.00	0.00	0.00
0.00	0.00	0.00

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.49	152	326397	20.00	ng/ul	0.00
18) Naphthalene-d8	10.25	136	1291682	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.14	164	770734	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.89	188	1621990	20.00	ng/ul	0.00
78) Chrysene-d12	21.10	240	1623622	20.00	ng/ul	0.00
86) Perylene-d12	23.25	264	1930889	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.06	96	2793m	0.38	ng/uL	> 0.03	JU 12/10/19
5) Phenol-d5	6.68	99	33144	1.16	ng/ul	0.01	
7) Bis-(2-Chloroethyl)ether-d	6.83	67	101939	6.13	ng/ul	0.00	
9) 2-Chlorophenol-d4	7.02	132	106583	4.82	ng/ul	0.00	
13) 4-Methylphenol-d8	8.20	113	60611	2.69	ng/ul	0.00	
19) Nitrobenzene-d5	8.63	128	63267	6.31	ng/ul	0.00	
22) 2-Nitrophenol-d4	9.35	143	68791	6.04	ng/ul	0.00	
26) 2,4-Dichlorophenol-d3	9.89	165	118993	5.51	ng/ul	0.00	
29) 4-Chloroaniline-d4	10.41	131	7955	0.34	ng/ul	0.01	
44) Dimethylphthalate-d6	13.56	166	403968	6.75	ng/ul	0.00	
47) Acenaphthylene-d8	13.83	160	456981	6.51	ng/ul	0.00	
52) 4-Nitrophenol-d4	14.42	143	5020	0.50	ng/ul	0.04	
58) Fluorene-d10	15.14	176	354330	7.03	ng/ul	0.00	
63) 4,6-Dinitro-2-methylphenol	15.29	200	32619	3.20	ng/ul	0.00	
71) Anthracene-d10	16.99	188	545875	7.05	ng/ul	0.00	
79) Pyrene-d10	19.30	212	627164	7.48	ng/ul	0.00	
90) Benzo(a)pyrene-d12	23.12	264	732989	7.18	ng/ul	0.00	

Target Compounds

					Qvalue
20) Nitrobenzene	8.68	77	754574	30.983	ng/ul 98
37) 1,2,4,5-Tetrachlorobenzene	12.31	216	218471	9.088	ng/ul 99
73) 1,2,3,4-Tetrachlorobenzene	12.93	216	709649	29.090	ng/uL 100

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