

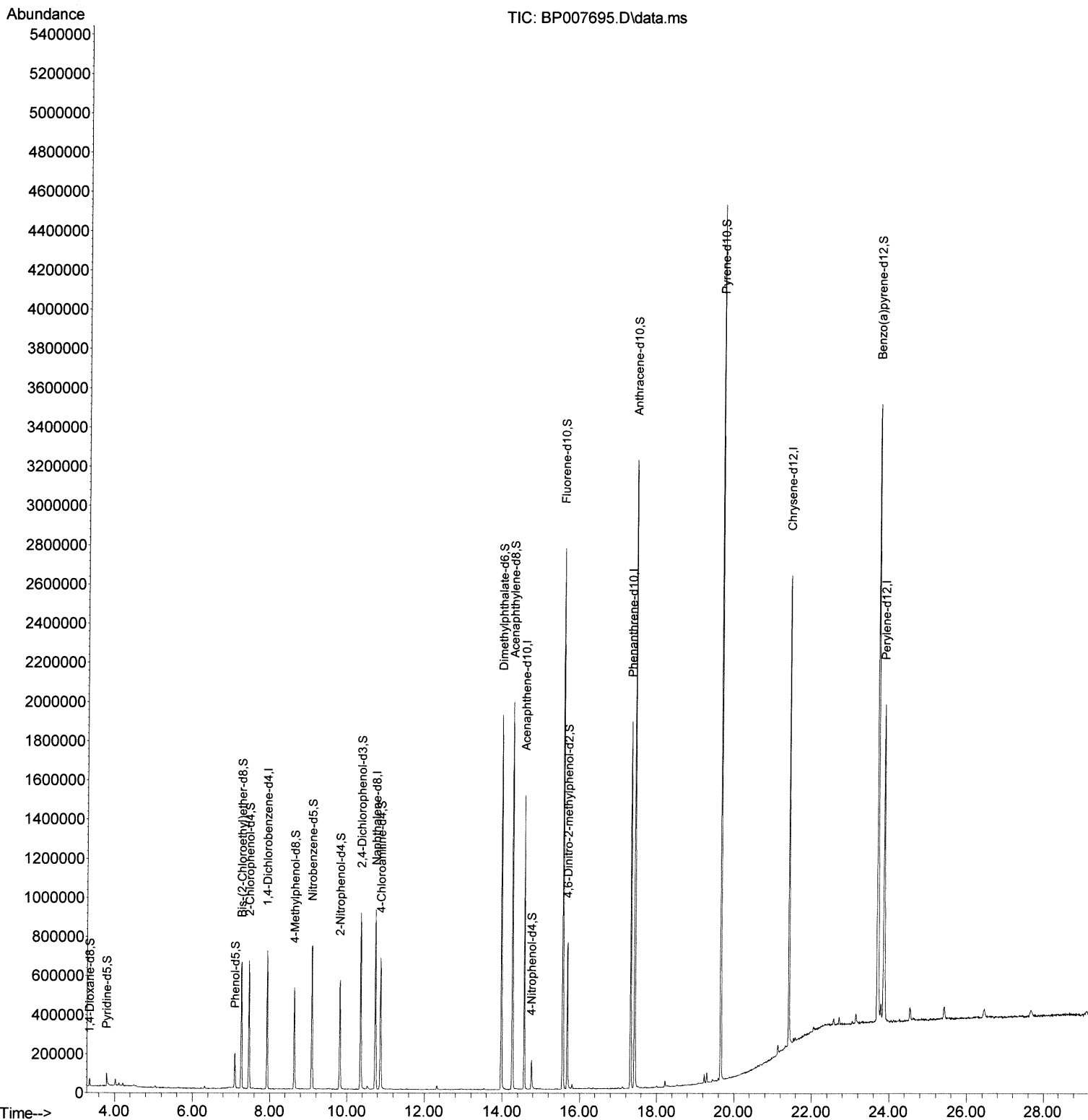
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Data File : BP007695.D
Acq On : 27 Oct 2021 00:42
Operator : CG/JU
Sample : M4303-02
Misc :
ALS Vial : 53 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
P004-TW001-02

Manual Integrations
APPROVED

mohammad
10/27/2021 11:50:53 AM

Quant Time: Oct 27 02:36:48 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP102521.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Oct 25 16:57:16 2021
Response via : Initial Calibration



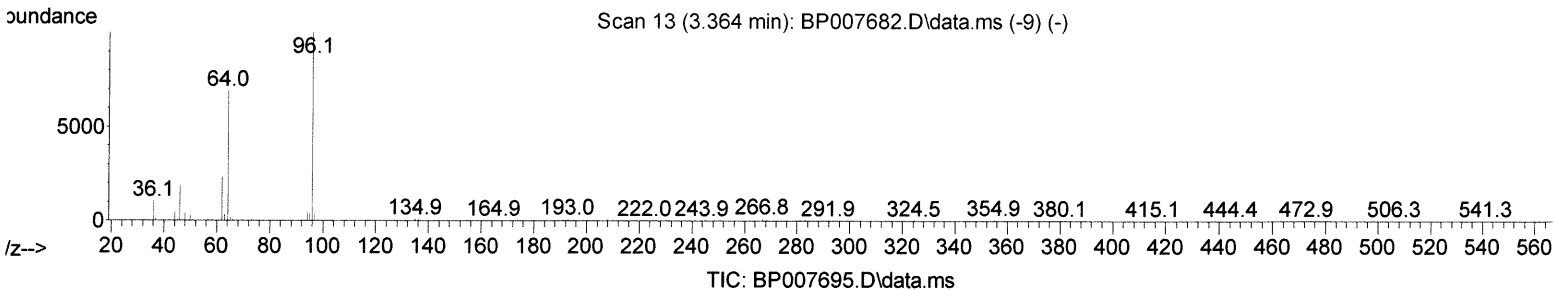
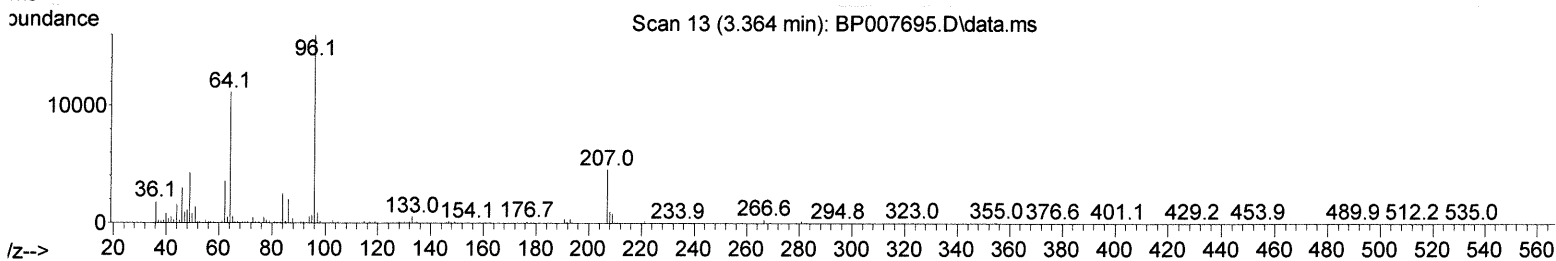
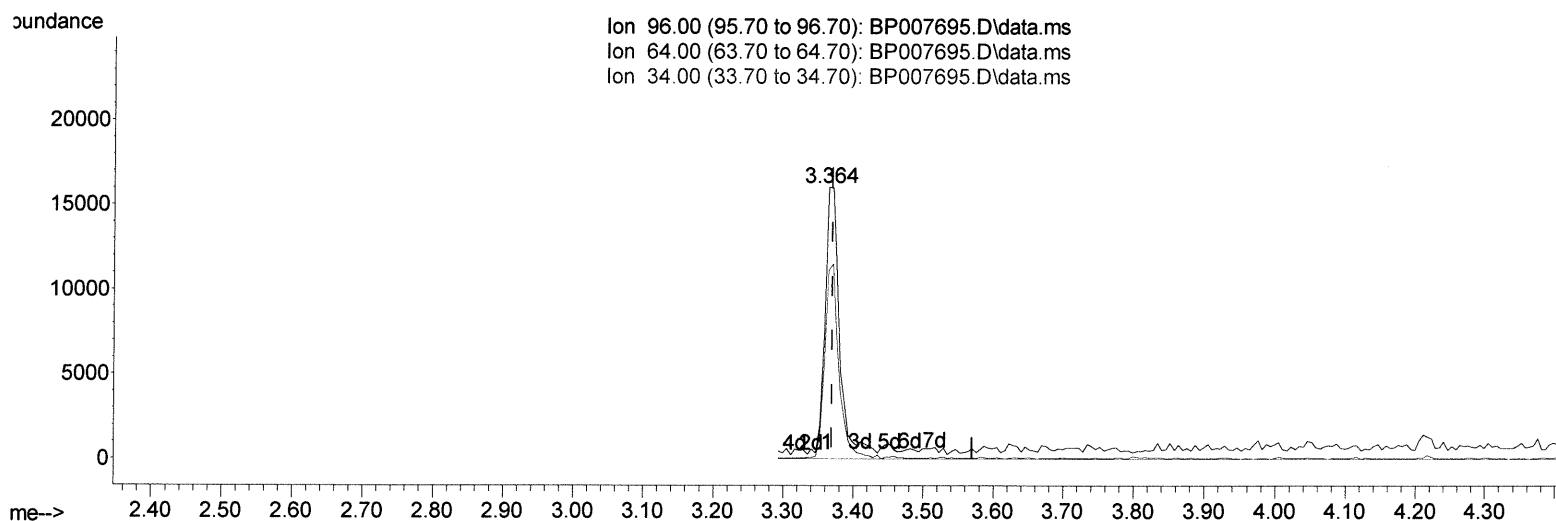
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(3) 1,4-Dioxane-d8 (S)

3.364min (-0.006) 5.19 ng/uL

response 21612

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	69.20	69.42
34.00	0.00	0.00
0.00	0.00	0.00

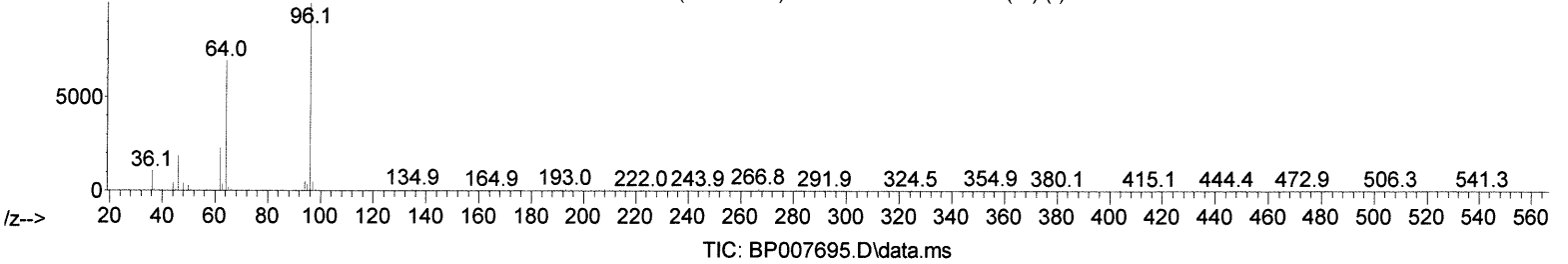
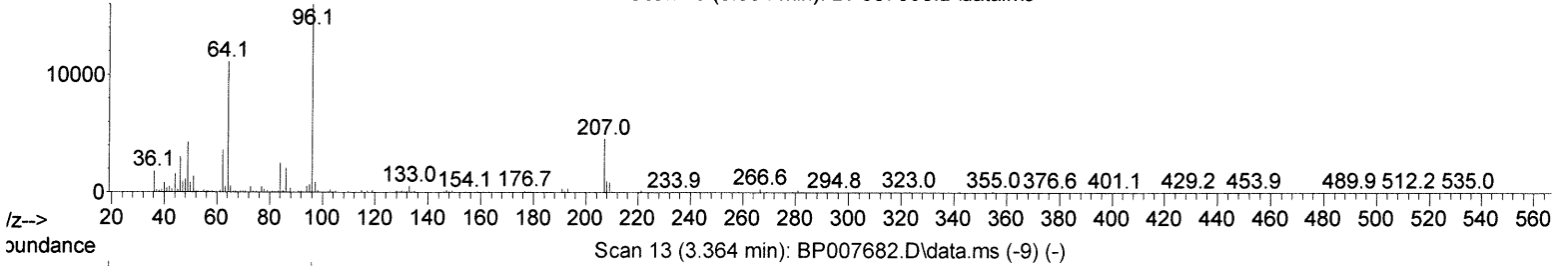
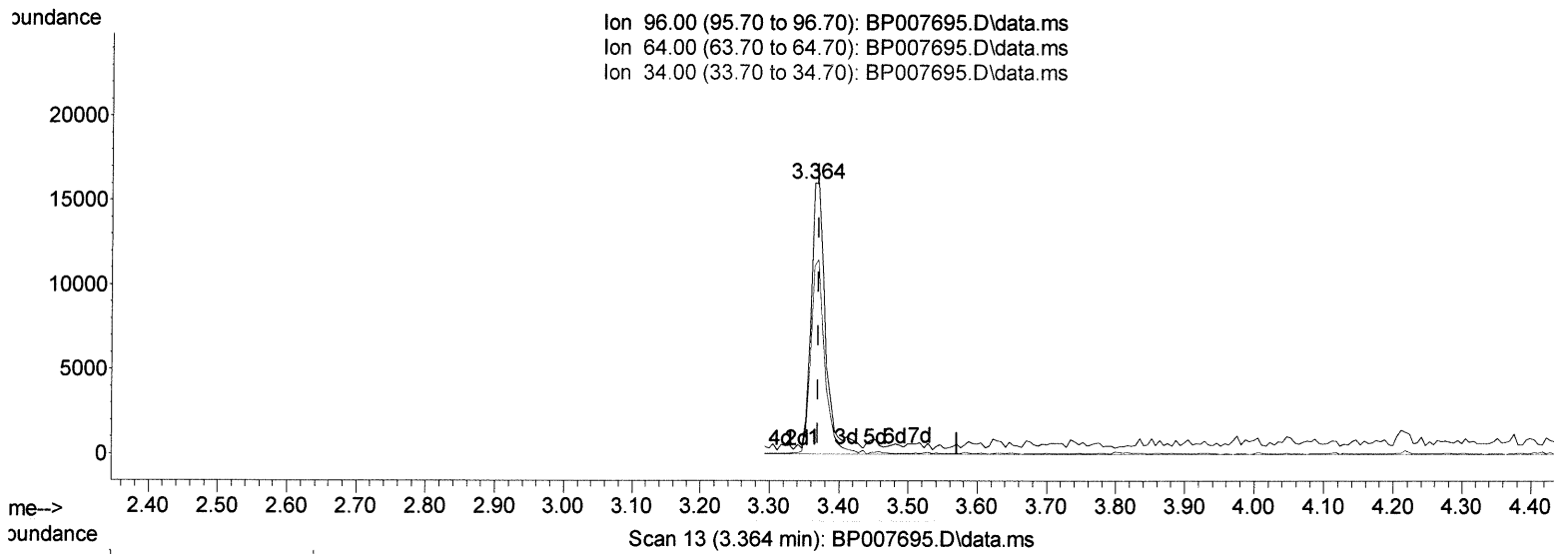
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Manual Integrations
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(3) 1,4-Dioxane-d8 (S)

3.364min (-0.006) 5.33 ng/uL m 10/29/21 JU

response 22207

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	69.20	69.42
34.00	0.00	0.00
0.00	0.00	0.00

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Manual Integrations
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.934	152	189186	20.000	ng/uL	0.00
20) Naphthalene-d8	10.740	136	775782	20.000	ng/uL	0.00
38) Acenaphthene-d10	14.569	164	513413	20.000	ng/uL	-0.01
64) Phenanthrene-d10	17.322	188	1105657	20.000	ng/uL	-0.01
79) Chrysene-d12	21.416	240	1195440	20.000	ng/uL	0.00
88) Perylene-d12	23.857	264	1249907	20.000	ng/uL	-0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.364	96	22207m	5.332	ng/uL	> 0.00
4) Pyridine-d5	3.793	84	36417	2.892	ng/uL	0.00
7) Phenol-d5	7.099	99	104069	7.242	ng/uL	0.00
9) Bis-(2-Chloroethyl)eth...	7.263	67	277852	34.827	ng/uL	0.00
11) 2-Chlorophenol-d4	7.463	132	297562	26.625	ng/uL	-0.01
15) 4-Methylphenol-d8	8.640	113	192131	17.075	ng/uL	-0.01
21) Nitrobenzene-d5	9.093	128	182975	34.610	ng/uL	0.00
24) 2-Nitrophenol-d4	9.816	143	182279	31.568	ng/uL	-0.01
28) 2,4-Dichlorophenol-d3	10.357	165	338283	28.782	ng/uL	-0.01
31) 4-Chloroaniline-d4	10.869	131	362438	23.322	ng/uL	-0.01
46) Dimethylphthalate-d6	13.981	166	1264840	36.370	ng/uL	-0.01
49) Acenaphthylene-d8	14.263	160	1426992	33.478	ng/uL	0.00
54) 4-Nitrophenol-d4	14.757	143	38224	5.996	ng/uL	-0.02
60) Fluorene-d10	15.563	176	1084585	36.993	ng/uL	-0.01
65) 4,6-Dinitro-2-methylph...	15.681	200	168953	28.567	ng/uL	0.00
73) Anthracene-d10	17.422	188	1903562	41.894	ng/uL	-0.01
81) Pyrene-d10	19.657	212	2331510	41.891	ng/uL	0.00
92) Benzo(a)pyrene-d12	23.704	264	2394176	38.992	ng/uL	0.00

Target Compounds Qvalue

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