

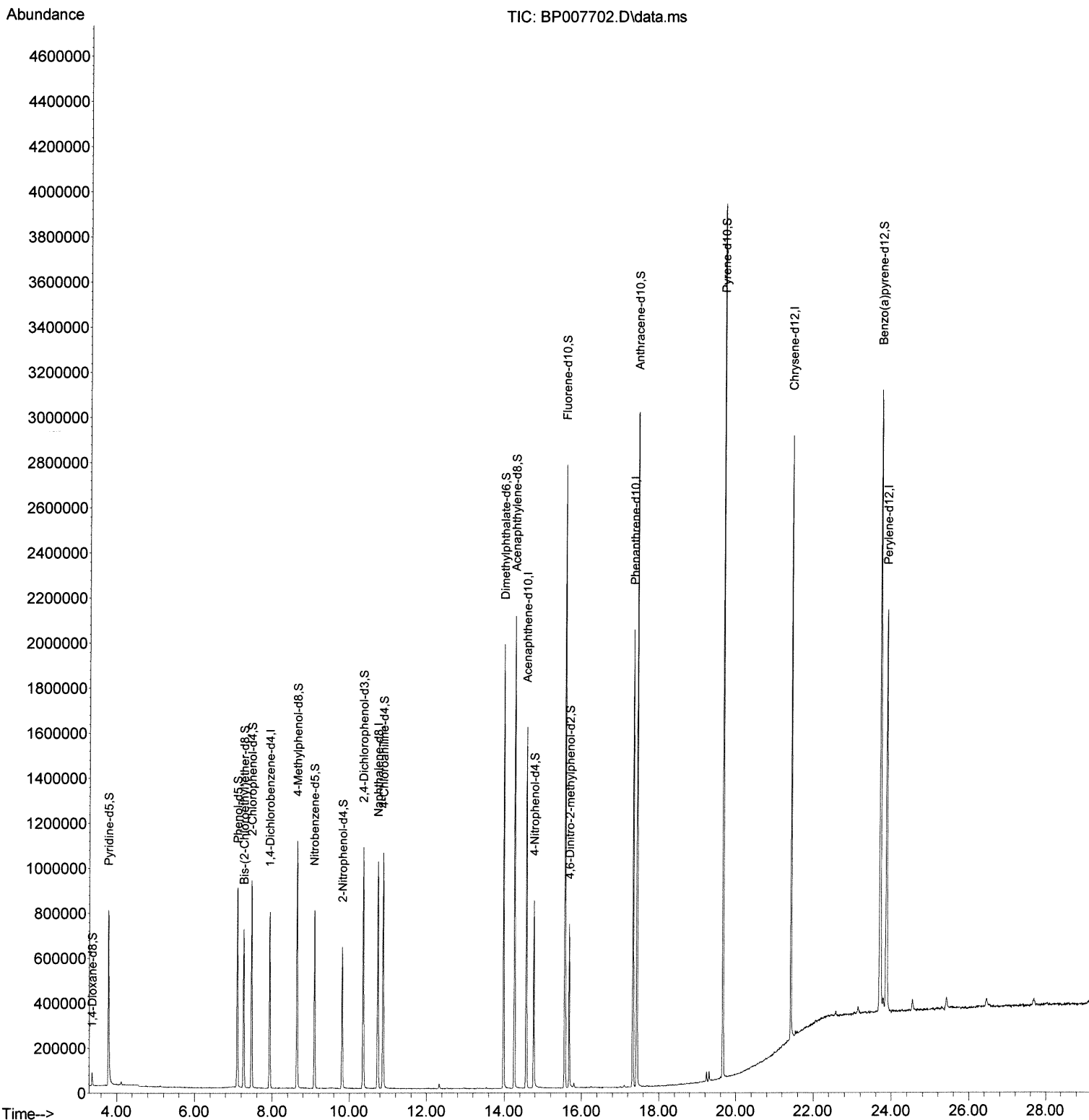
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102521\
Data File : BP007702.D
Acq On : 27 Oct 2021 05:13
Operator : CG/JU
Sample : PB140188BL
Misc :
ALS Vial : 60 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SBLK188

Manual Integrations
APPROVED

mohammad
10/27/2021 11:51:11 AM

Quant Time: Oct 27 05:55:19 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



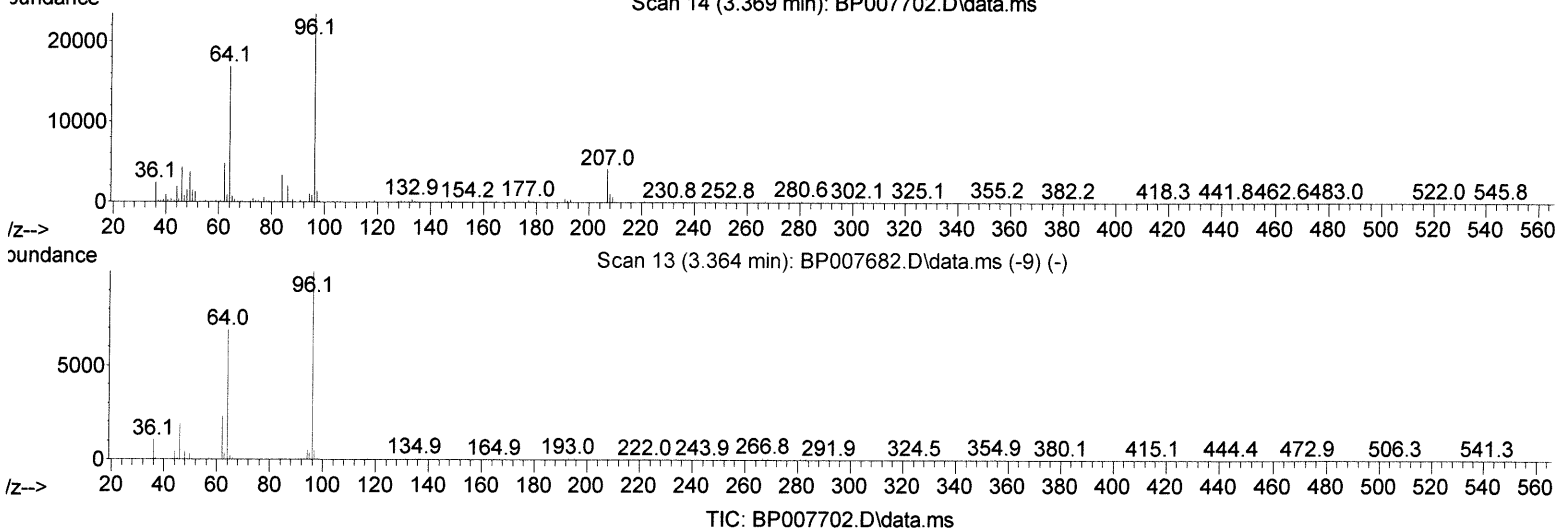
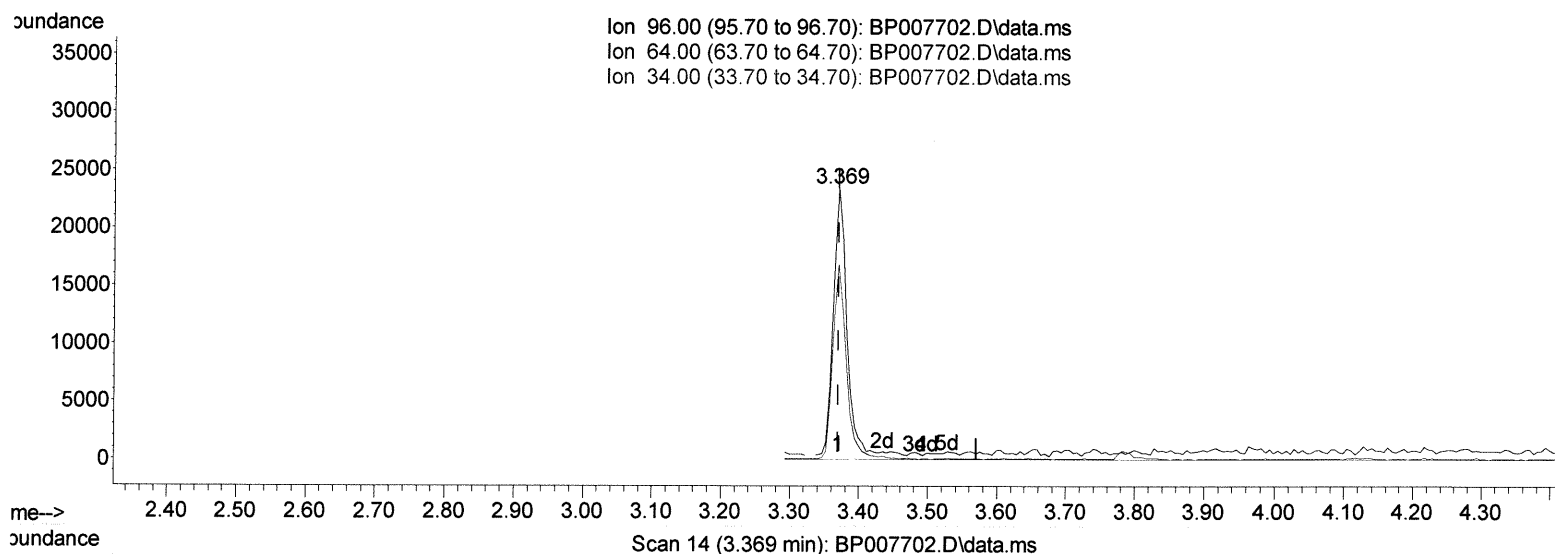
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(3) 1,4-Dioxane-d8 (S)

3.369min (-0.001) 6.60 ng/uL

response 30761

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	69.20	71.97
34.00	0.00	0.00
0.00	0.00	0.00

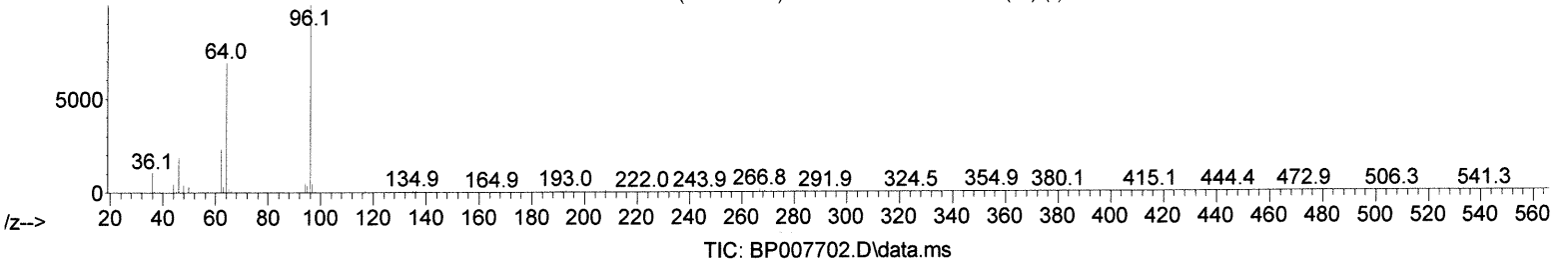
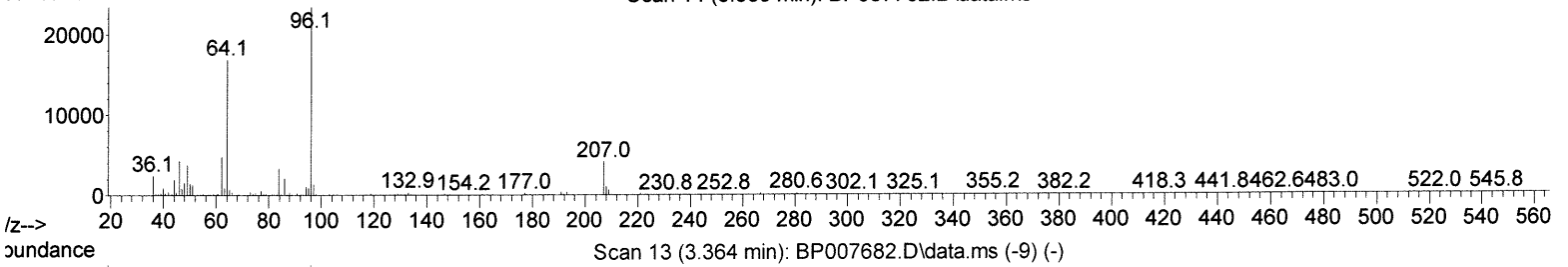
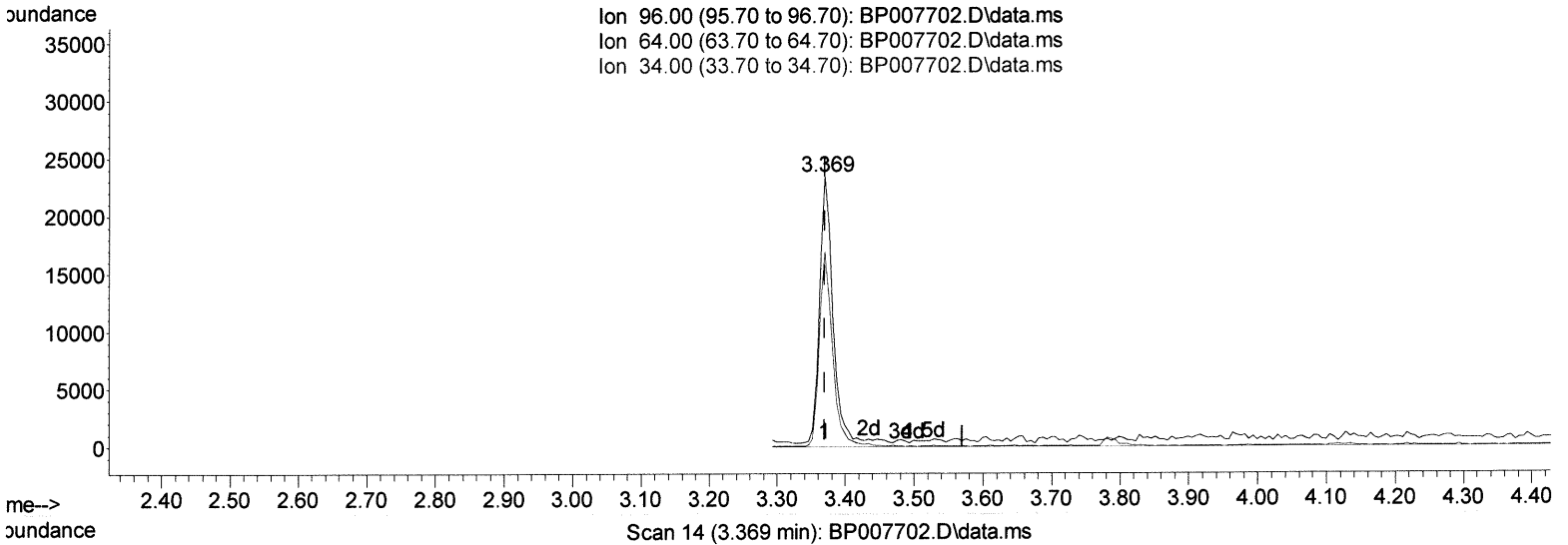
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Manual Integrations
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 Response via : Initial Calibration



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

3.369min (-0.001) 6.71 ng/uL m 10/29/21 JU

response 31289

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	69.20	71.97
34.00	0.00	0.00
0.00	0.00	0.00

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Manual Integrations
 APPROVED

mohammad
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.934	152	211805	20.000	ng/ul	0.00
20) Naphthalene-d8	10.734	136	850777	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.569	164	560318	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.328	188	1214741	20.000	ng/ul	0.00
79) Chrysene-d12	21.416	240	1290762	20.000	ng/ul	0.00
88) Perylene-d12	23.863	264	1325395	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.369	96	31289m	> 6.711	ng/uL	> 0.00
4) Pyridine-d5	3.781	84	461330	32.718	ng/ul	0.00
7) Phenol-d5	7.099	99	521059	32.389	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.263	67	307130	34.386	ng/ul	0.00
11) 2-Chlorophenol-d4	7.463	132	421909	33.719	ng/ul	-0.01
15) 4-Methylphenol-d8	8.640	113	409585	32.514	ng/ul	-0.01
21) Nitrobenzene-d5	9.087	128	192730	33.242	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.816	143	207338	32.743	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.357	165	395084	30.652	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.869	131	554339	32.526	ng/ul	-0.01
46) Dimethylphthalate-d6	13.981	166	1278968	33.697	ng/ul	-0.01
49) Acenaphthylene-d8	14.263	160	1486673	31.958	ng/ul	0.00
54) 4-Nitrophenol-d4	14.763	143	204354	29.373	ng/ul	-0.01
60) Fluorene-d10	15.563	176	1079682	33.743	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.681	200	160170	24.650	ng/ul	0.00
73) Anthracene-d10	17.422	188	1739455	34.844	ng/ul	-0.01
81) Pyrene-d10	19.663	212	2125729	35.374	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.704	264	2142917	32.913	ng/ul	0.00

Target Compounds Qvalue

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