

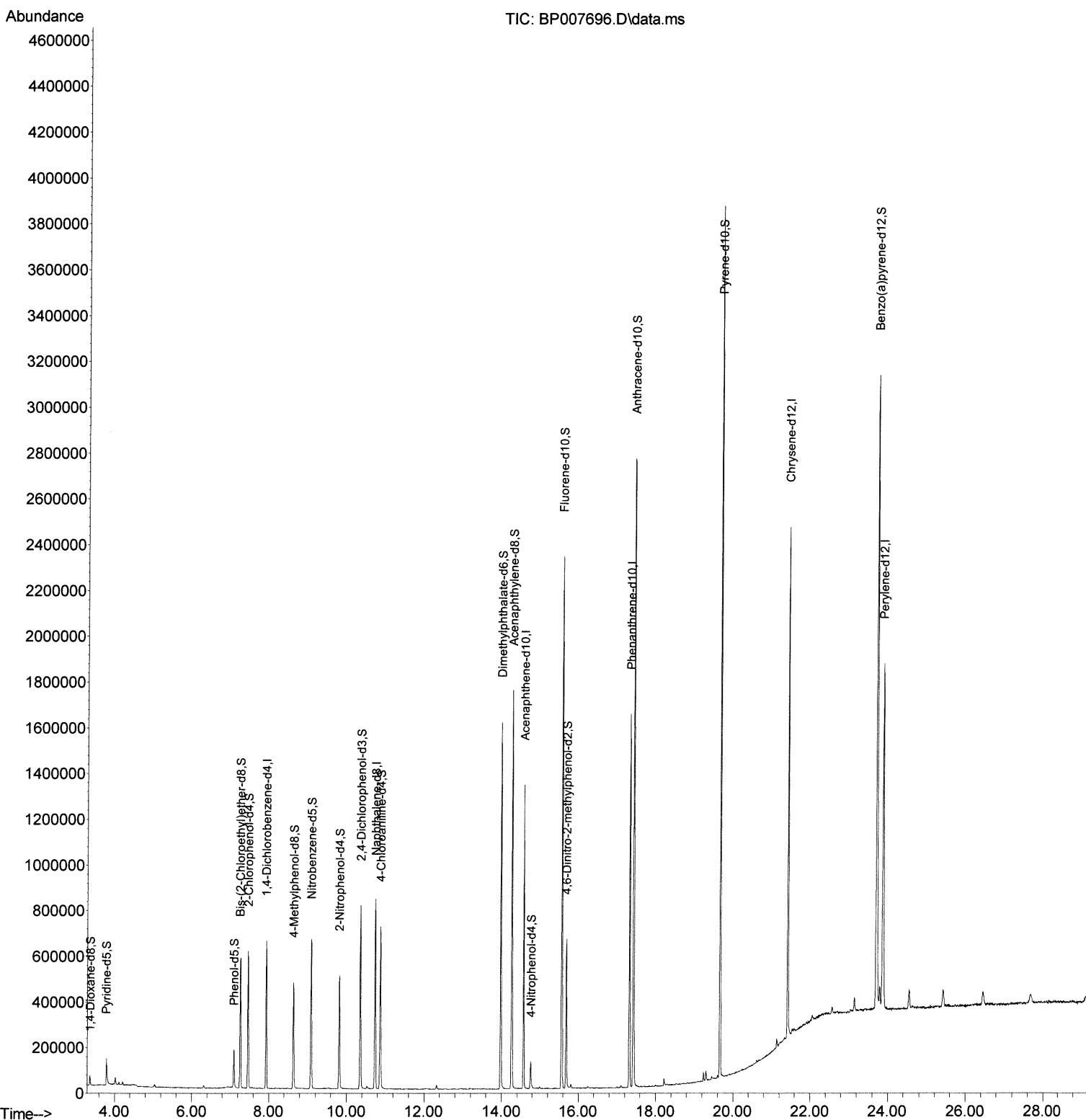
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102521\
Data File : BP007696.D
Acq On : 27 Oct 2021 01:16
Operator : CG/JU
Sample : M4303-03
Misc :
ALS Vial : 54 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
P005-TW001-02

Manual Integrations
APPROVED

mohammad
10/27/2021 11:50:58 AM

Quant Time: Oct 27 02:37:09 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



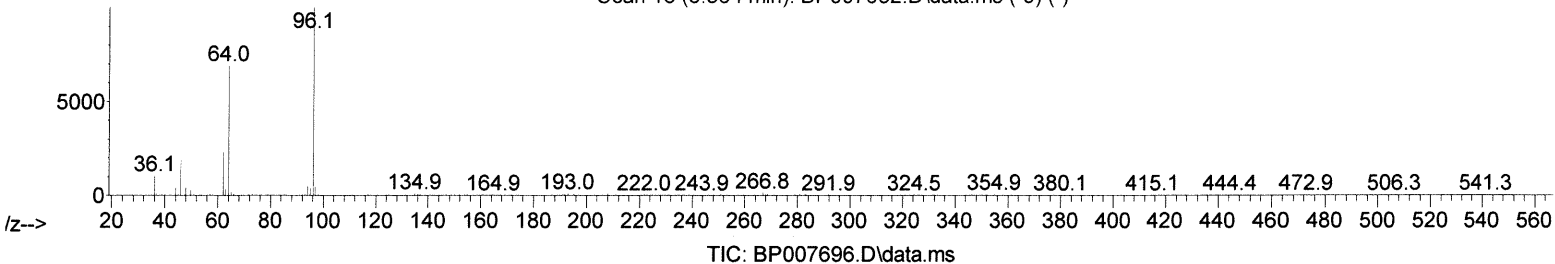
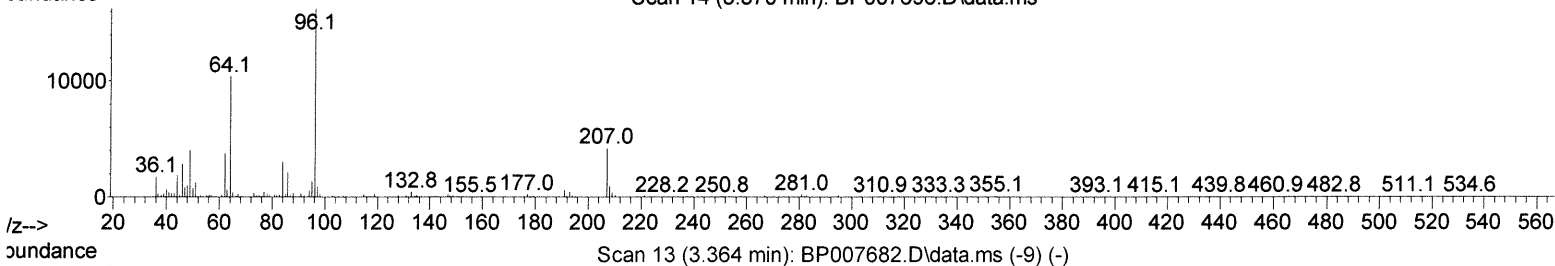
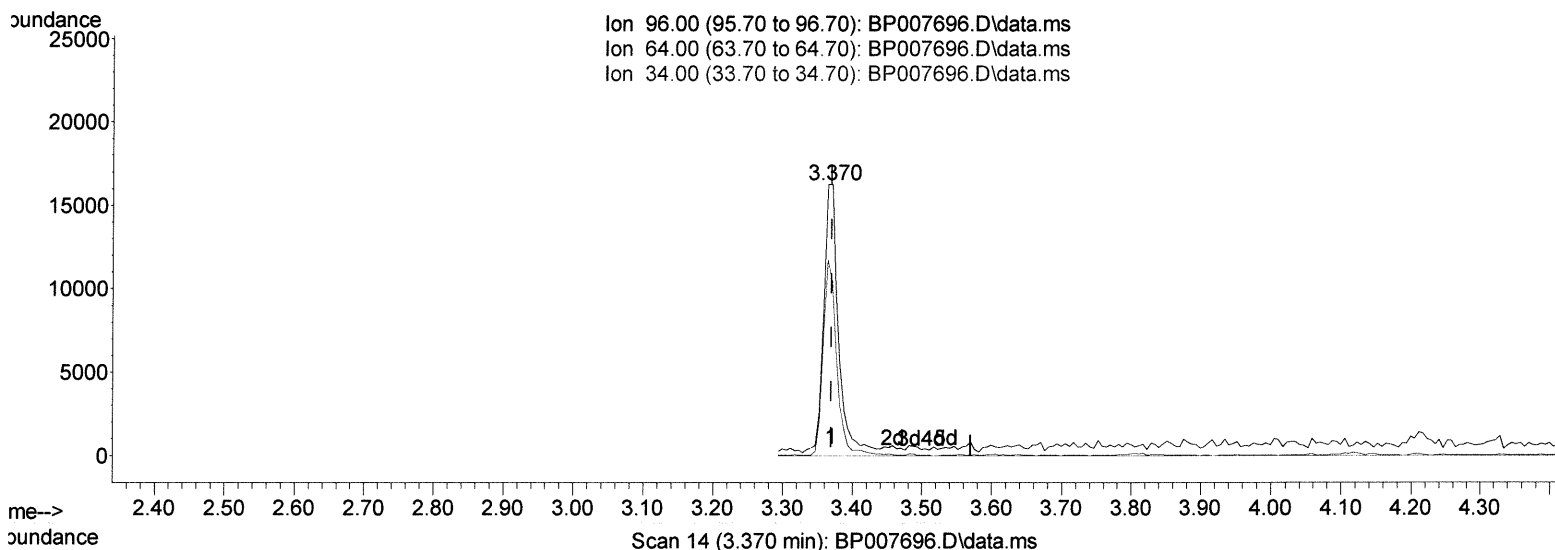
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(3) 1,4-Dioxane-d8 (S)

3.370min (-0.000) 5.58 ng/uL

response 21718

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

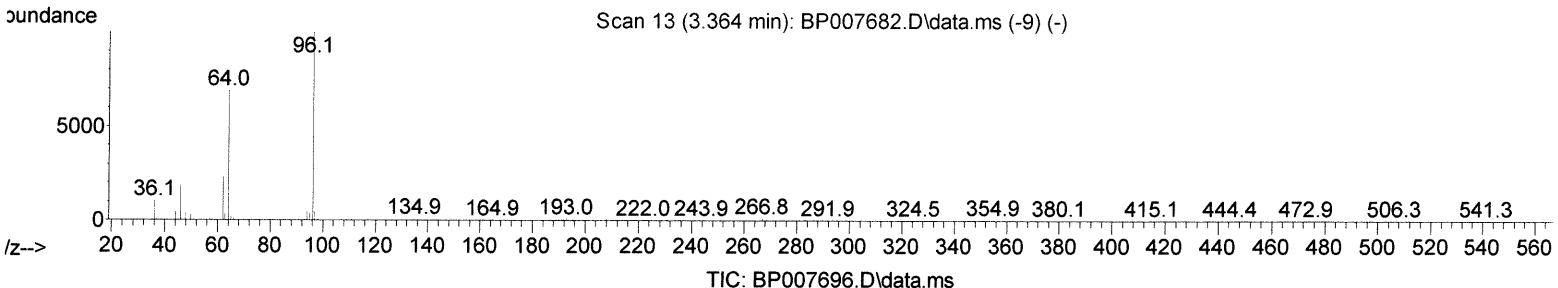
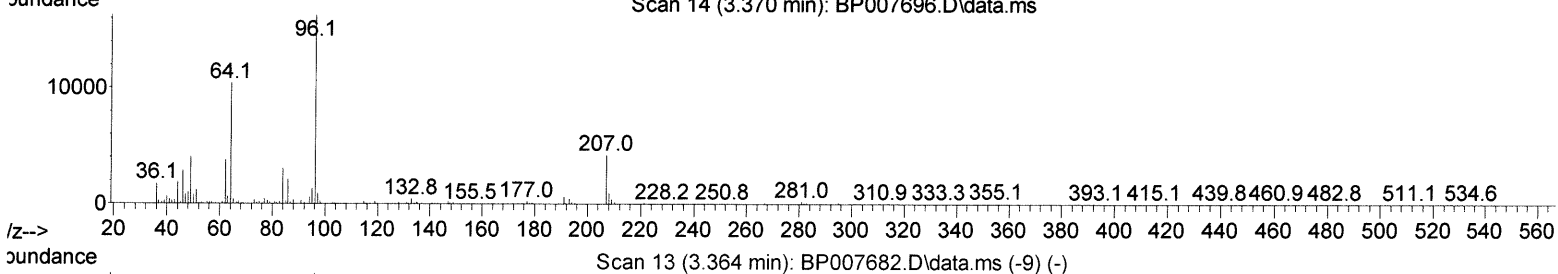
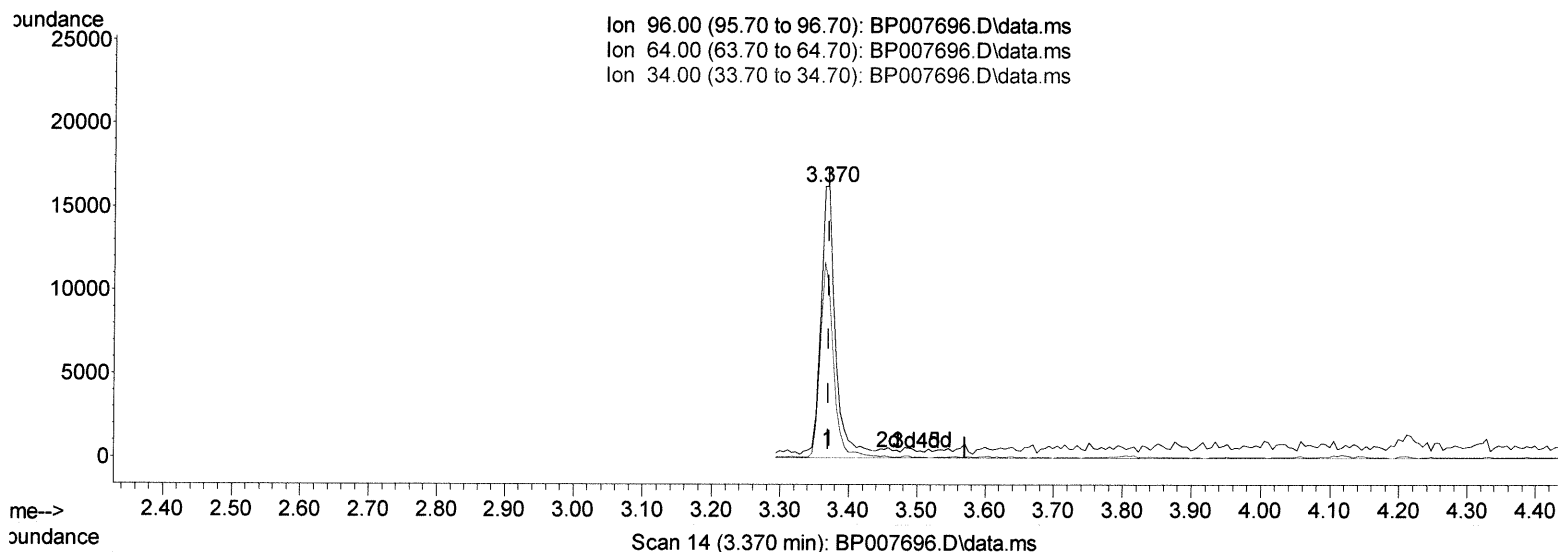
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Manual Integrations
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(3) 1,4-Dioxane-d8 (S)

3.370min (-0.000) 6.08 ng/uL m 1029130

response 23669

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

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Manual Integrations
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mohammad
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 Last Update : Mon Oct 25 16:57:16 2021
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.934	152	176788	20.000	ng/ul	0.00
20) Naphthalene-d8	10.740	136	706262	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.569	164	458933	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.328	188	999103	20.000	ng/ul	0.00
79) Chrysene-d12	21.416	240	1074787	20.000	ng/ul	0.00
88) Perylene-d12	23.863	264	1119903	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.370	96	23669m	6.082	ng/uL	0.00 10/29/21 JU
4) Pyridine-d5	3.787	84	63115	5.363	ng/ul	0.00
7) Phenol-d5	7.093	99	95720	7.128	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	7.258	67	252122	33.818	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.464	132	269390	25.794	ng/ul	-0.01
15) 4-Methylphenol-d8	8.640	113	172782	16.433	ng/ul	-0.01
21) Nitrobenzene-d5	9.093	128	155932	32.398	ng/ul	0.00
24) 2-Nitrophenol-d4	9.816	143	161574	30.737	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.358	165	303091	28.326	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.869	131	379010	26.789	ng/ul	-0.01
46) Dimethylphthalate-d6	13.981	166	1068861	34.383	ng/ul	-0.01
49) Acenaphthylene-d8	14.263	160	1236240	32.446	ng/ul	0.00
54) 4-Nitrophenol-d4	14.763	143	31942	5.605	ng/ul	-0.01
60) Fluorene-d10	15.563	176	916495	34.971	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.681	200	141724	26.519	ng/ul	0.00
73) Anthracene-d10	17.422	188	1641064	39.968	ng/ul	-0.01
81) Pyrene-d10	19.657	212	2012123	40.211	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.704	264	2085914	37.916	ng/ul	0.00

Target Compounds Qvalue

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