

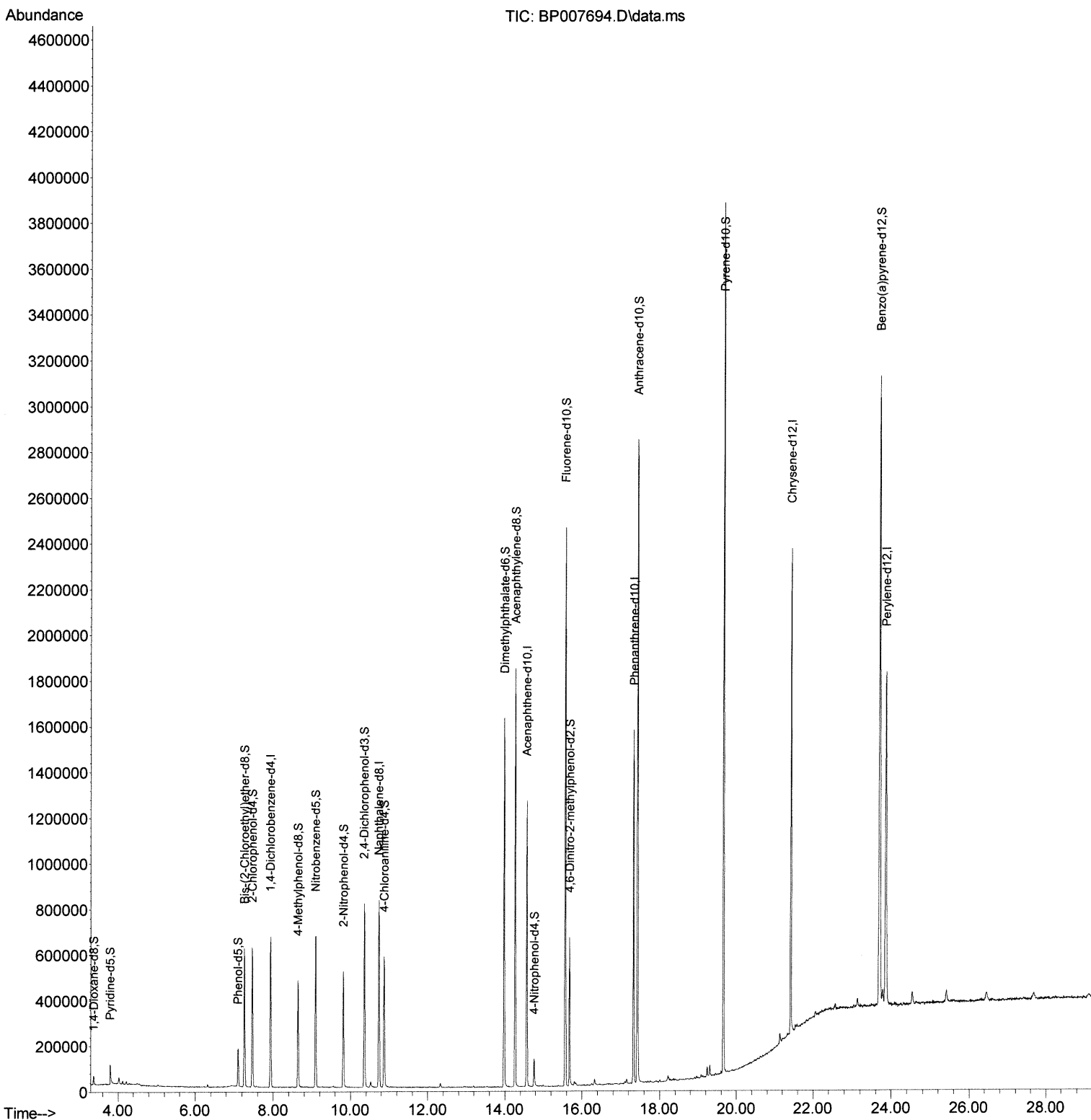
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102521\
Data File : BP007694.D
Acq On : 27 Oct 2021 00:08
Operator : CG/JU
Sample : M4303-01
Misc :
ALS Vial : 52 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
P003-TW001-02

Manual Integrations
APPROVED

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

Quant Time: Oct 27 00:57:46 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



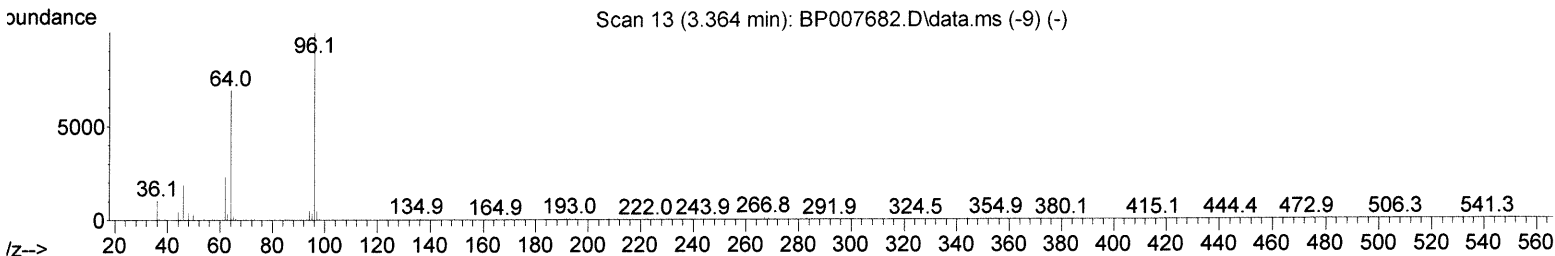
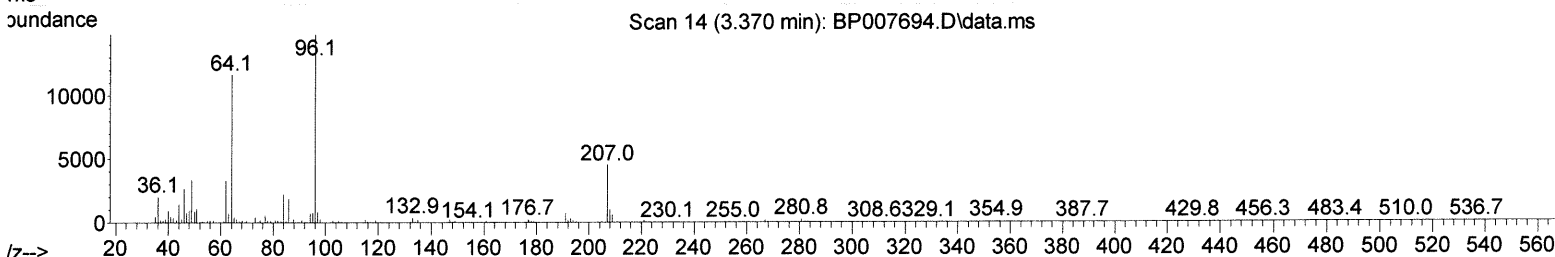
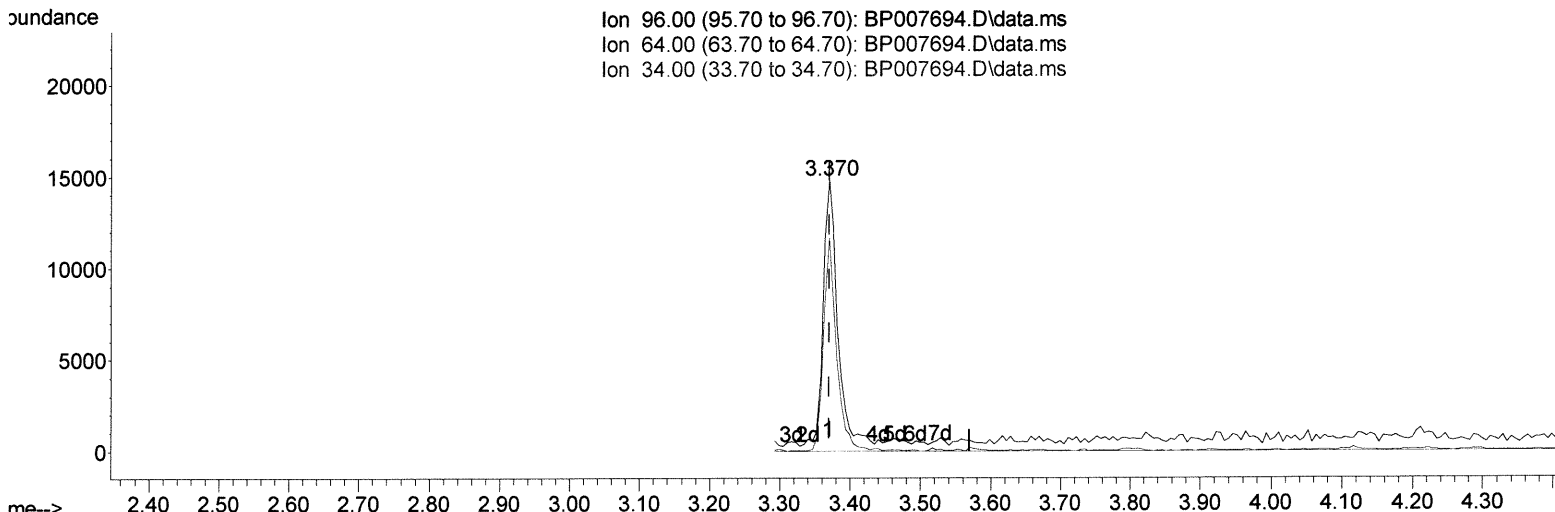
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TIC: BP007694.D\data.ms

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

3.370min (-0.000) 4.89 ng/uL

response 18995

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	69.20	78.54
34.00	0.00	0.00
0.00	0.00	0.00

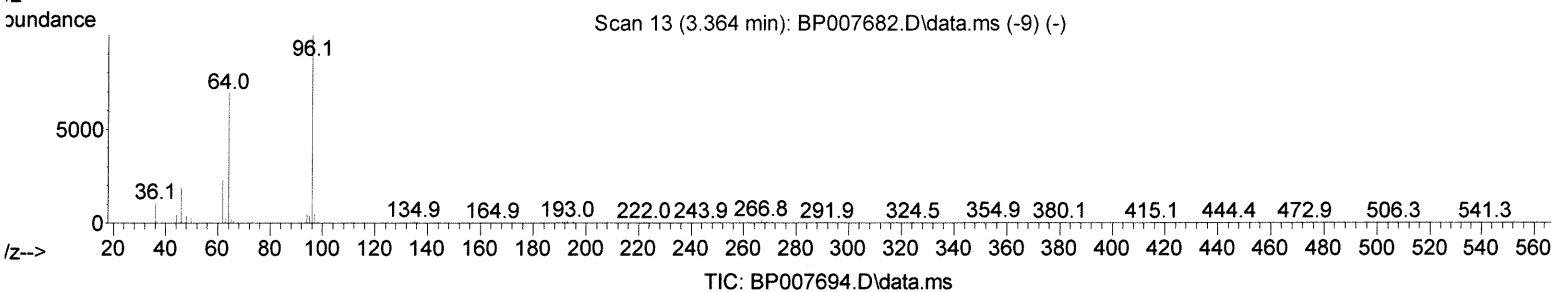
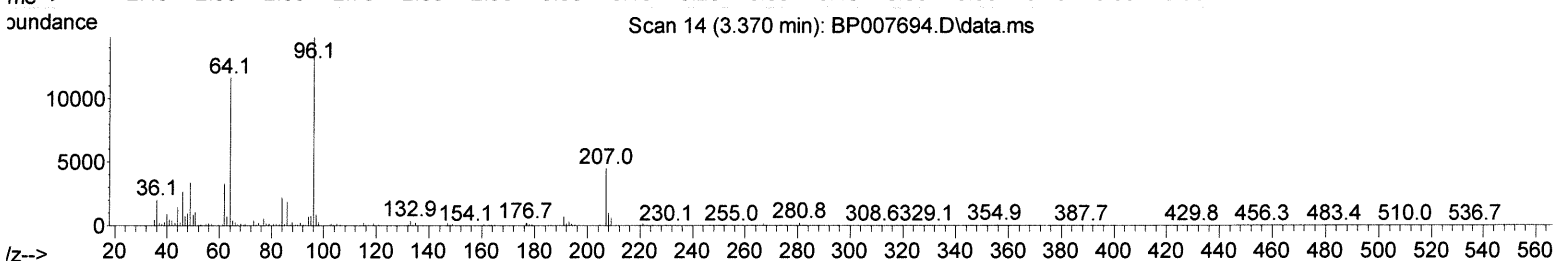
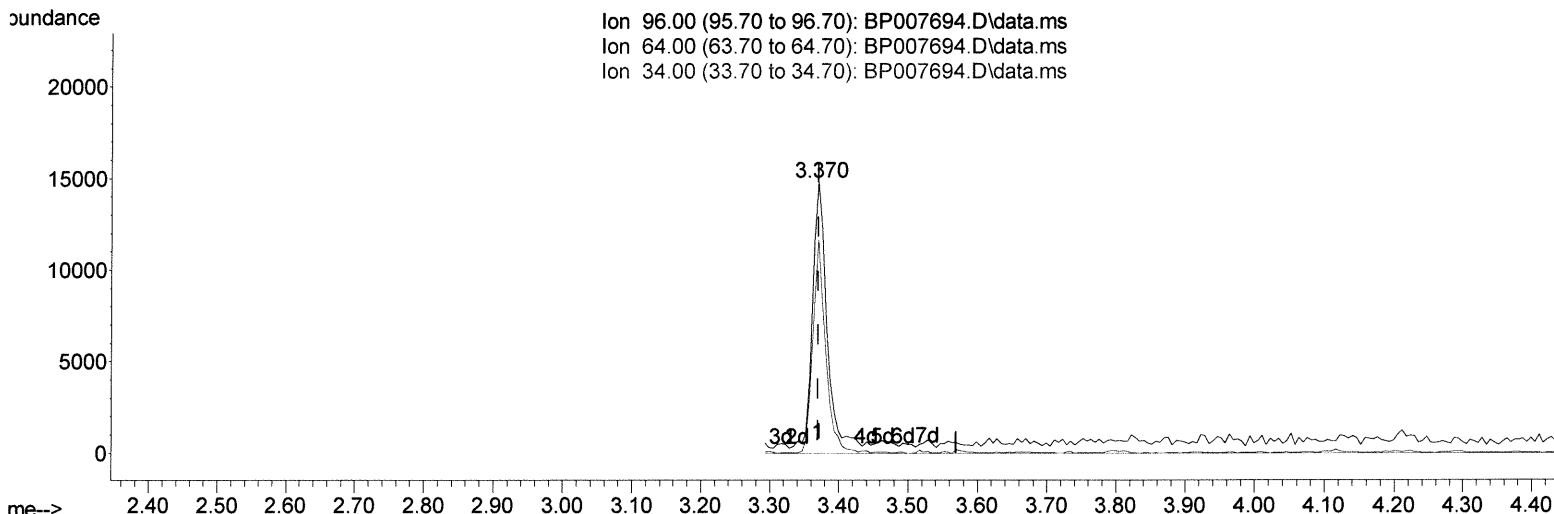
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(3) 1,4-Dioxane-d8 (S)

3.370min (-0.000) 5.29 ng/uL m 10/29/21 JU

response	20537	
Ion	Exp%	Act%
96.00	100.00	100.00
64.00	69.20	78.54
34.00	0.00	0.00
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102521\
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 Operator : CG/JU
 Sample : M4303-01
 Misc :
 ALS Vial : 52 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
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Manual Integrations
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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	176318	20.000	ng/ul	0.00
20) Naphthalene-d8	10.734	136	690625	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.569	164	442302	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.328	188	956197	20.000	ng/ul	0.00
79) Chrysene-d12	21.416	240	1056690	20.000	ng/ul	0.00
88) Perylene-d12	23.863	264	1114513	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.370	96	20537m>	5.291	ng/uL>	0.00 10/24/21 JU
4) Pyridine-d5	3.793	84	49461	4.214	ng/ul	0.00
7) Phenol-d5	7.099	99	96852	7.232	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.263	67	259352	34.881	ng/ul	0.00
11) 2-Chlorophenol-d4	7.463	132	280623	26.942	ng/ul	-0.01
15) 4-Methylphenol-d8	8.640	113	172270	16.428	ng/ul	-0.01
21) Nitrobenzene-d5	9.093	128	161682	34.354	ng/ul	0.00
24) 2-Nitrophenol-d4	9.816	143	163305	31.769	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.357	165	298892	28.567	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.869	131	305785	22.103	ng/ul	-0.01
46) Dimethylphthalate-d6	13.981	166	1097705	36.639	ng/ul	-0.01
49) Acenaphthylene-d8	14.263	160	1246962	33.958	ng/ul	0.00
54) 4-Nitrophenol-d4	14.763	143	32967	6.003	ng/ul	-0.01
60) Fluorene-d10	15.563	176	948117	37.538	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.681	200	142418	27.845	ng/ul	0.00
73) Anthracene-d10	17.422	188	1645526	41.875	ng/ul	-0.01
81) Pyrene-d10	19.657	212	2054242	41.756	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.704	264	2144091	39.162	ng/ul	0.00

Target Compounds Qvalue

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