

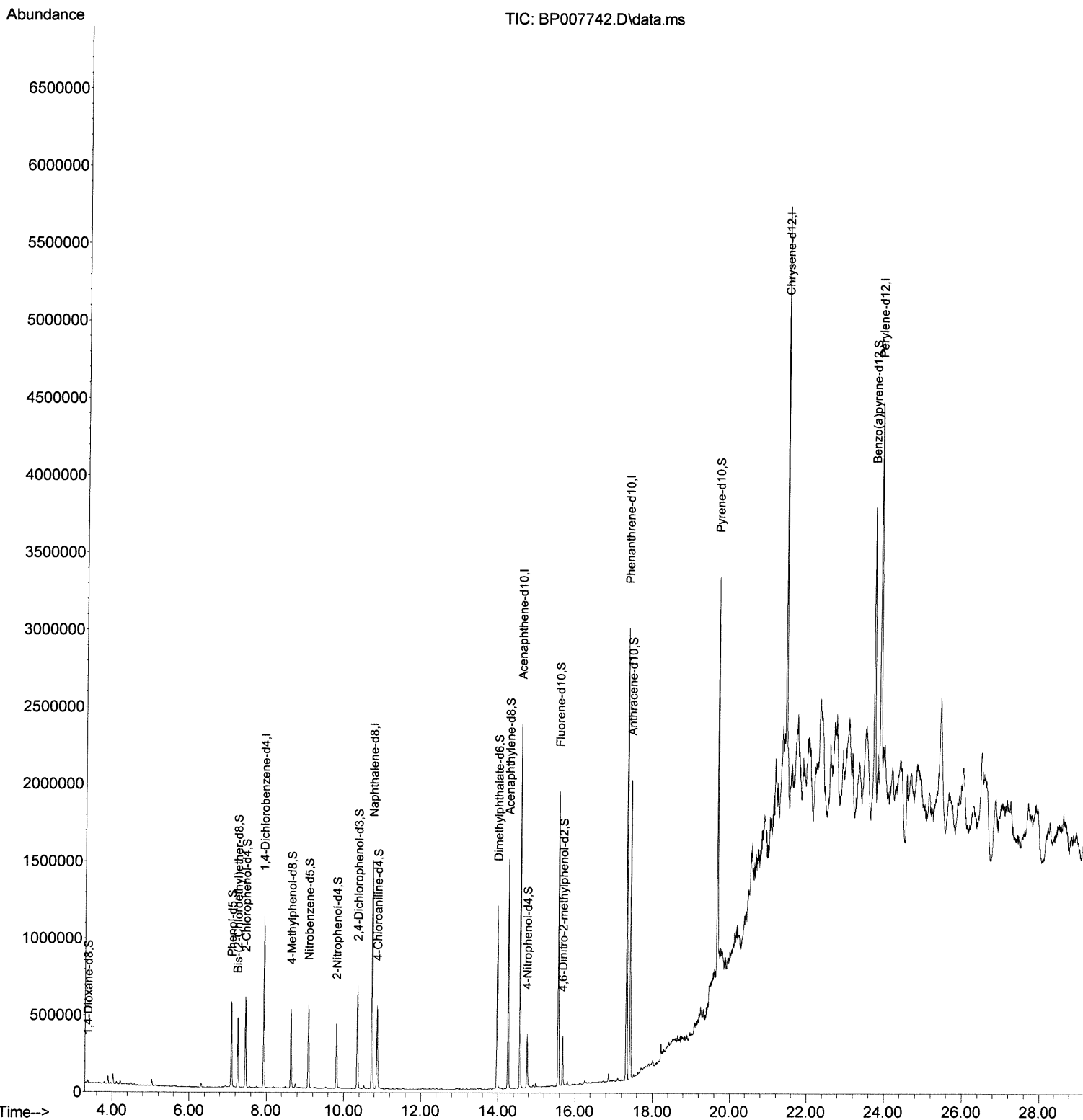
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102521\
Data File : BP007742.D
Acq On : 28 Oct 2021 04:35
Operator : CG/JU
Sample : M4215-16
Misc :
ALS Vial : 99 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
JDGF3

Manual Integrations
APPROVED

mohammad
10/28/2021 11:12:32 AM

Quant Time: Oct 28 06:13:40 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP102521.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Oct 27 15:37:43 2021
Response via : Initial Calibration



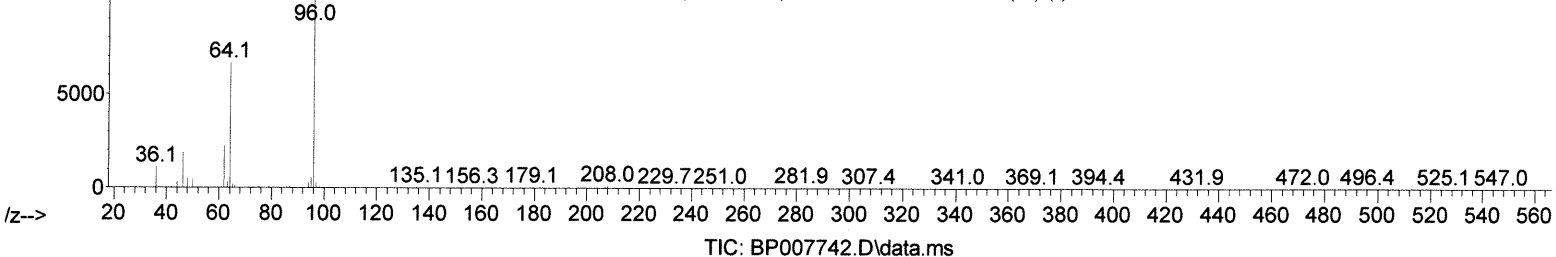
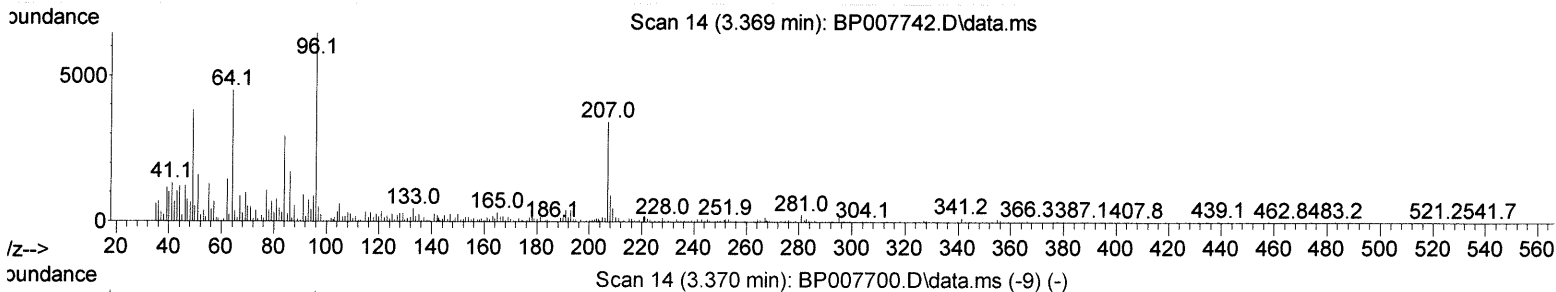
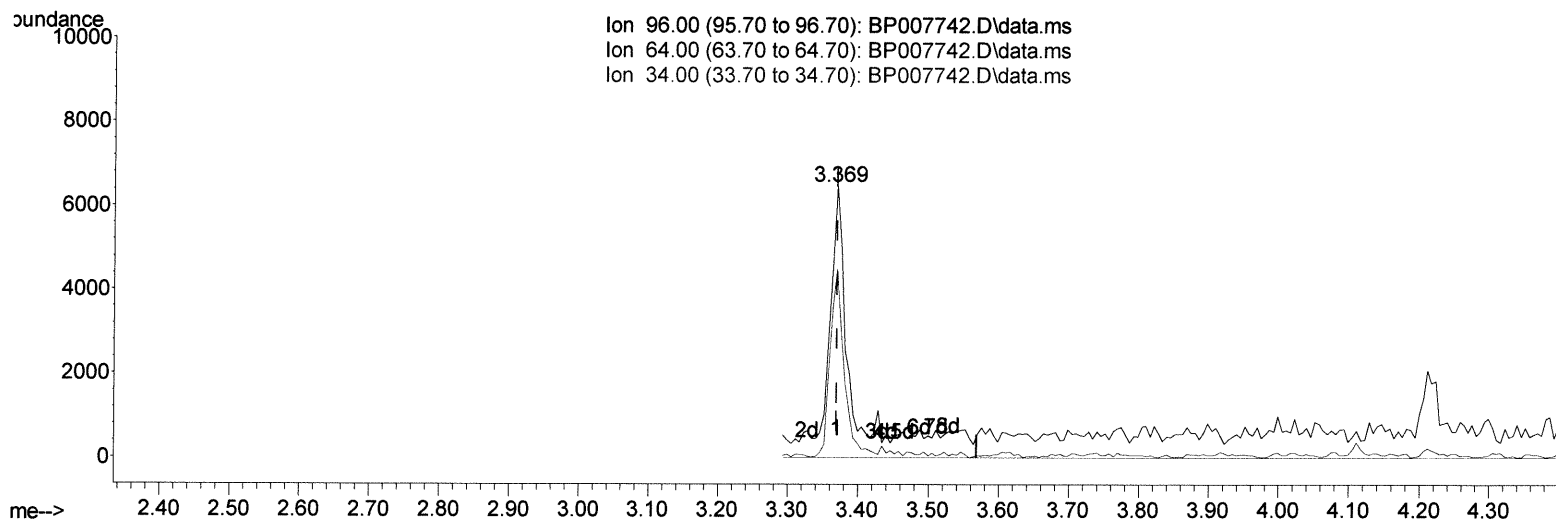
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TIC: BP007742.D\data.ms

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

3.369min (+ 0.000) 1.17 ng/uL

response 7979

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	69.20	69.55
34.00	0.00	0.00
0.00	0.00	0.00

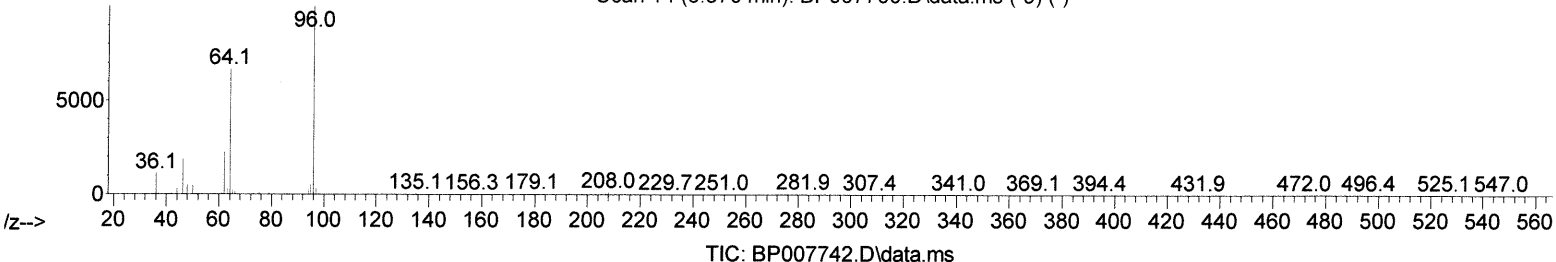
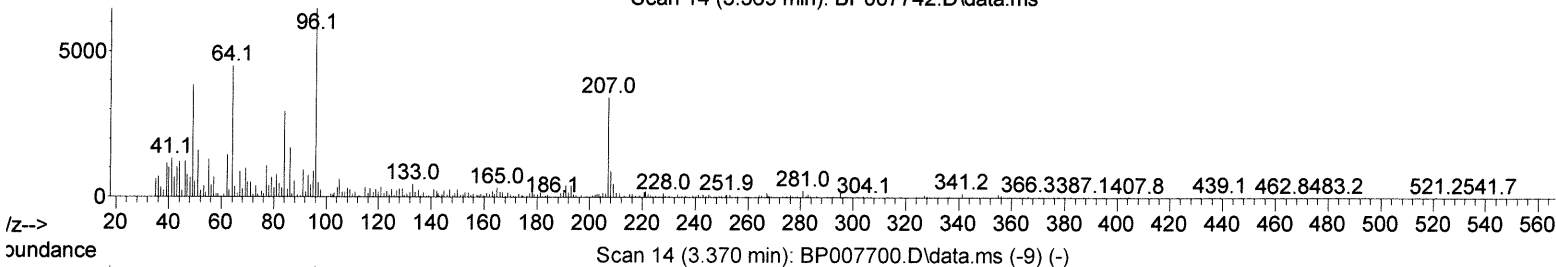
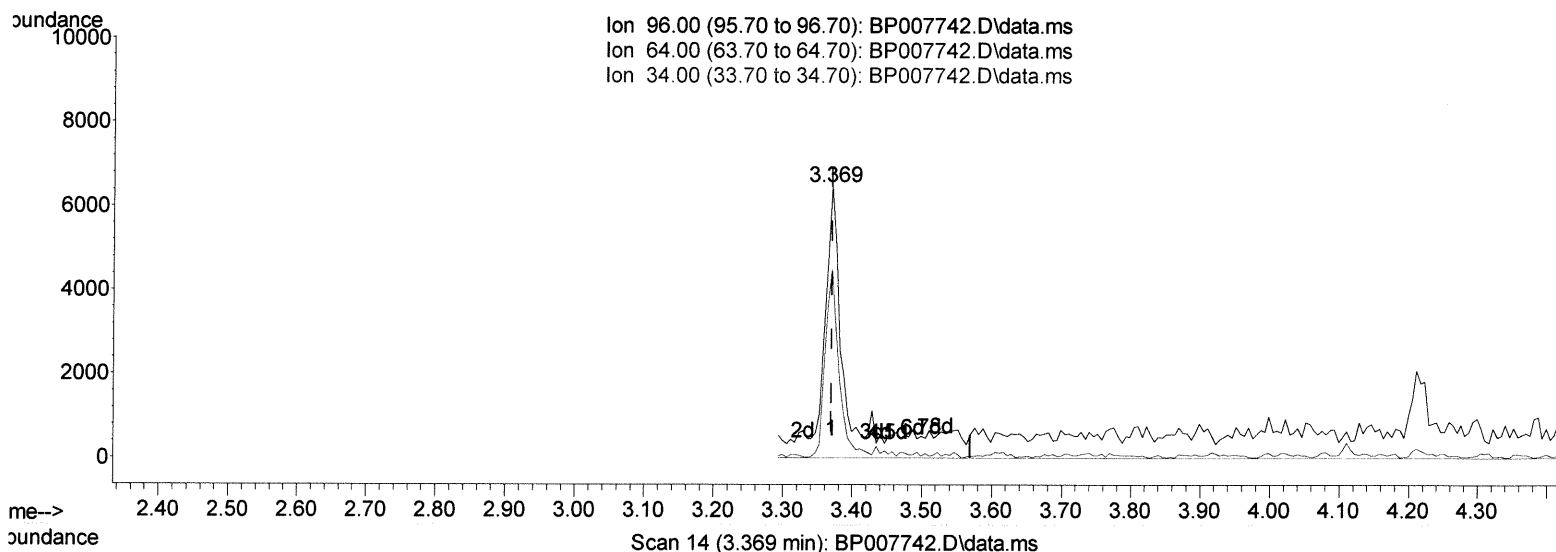
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Manual Integrations
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(3) 1,4-Dioxane-d8 (S)

3.369min (+ 0.000) 1.19 ng/uL m 10/29/2021

response 8102

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	69.20	69.55
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	310403	20.000	ng/ul	0.00
20) Naphthalene-d8	10.740	136	1282925	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.569	164	839452	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.328	188	1835688	20.000	ng/ul	0.00
79) Chrysene-d12	21.416	240	1740931	20.000	ng/ul	0.00
88) Perylene-d12	23.863	264	1928842	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.369	96	8102m	1.186	ng/uL	> 0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	7.093	99	329252	13.965	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.258	67	190174	14.528	ng/ul	0.00
11) 2-Chlorophenol-d4	7.463	132	282689	15.416	ng/ul	0.00
15) 4-Methylphenol-d8	8.640	113	197339	10.689	ng/ul	0.00
21) Nitrobenzene-d5	9.087	128	136756	15.642	ng/ul	0.00
24) 2-Nitrophenol-d4	9.810	143	147358	15.432	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.357	165	263804	13.573	ng/ul	0.00
31) 4-Chloroaniline-d4	10.869	131	297478	11.575	ng/ul	0.00
46) Dimethylphthalate-d6	13.981	166	835973	14.702	ng/ul	0.00
49) Acenaphthylene-d8	14.263	160	1060007	15.210	ng/ul	0.00
54) 4-Nitrophenol-d4	14.757	143	90495	8.682	ng/ul	0.00
60) Fluorene-d10	15.563	176	769203	16.046	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.681	200	78260	7.970	ng/ul	0.00
73) Anthracene-d10	17.422	188	1194469	15.833	ng/ul	0.00
81) Pyrene-d10	19.657	212	1405433	17.340	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.704	264	1431506	15.108	ng/ul	0.00

Target Compounds Qvalue

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