

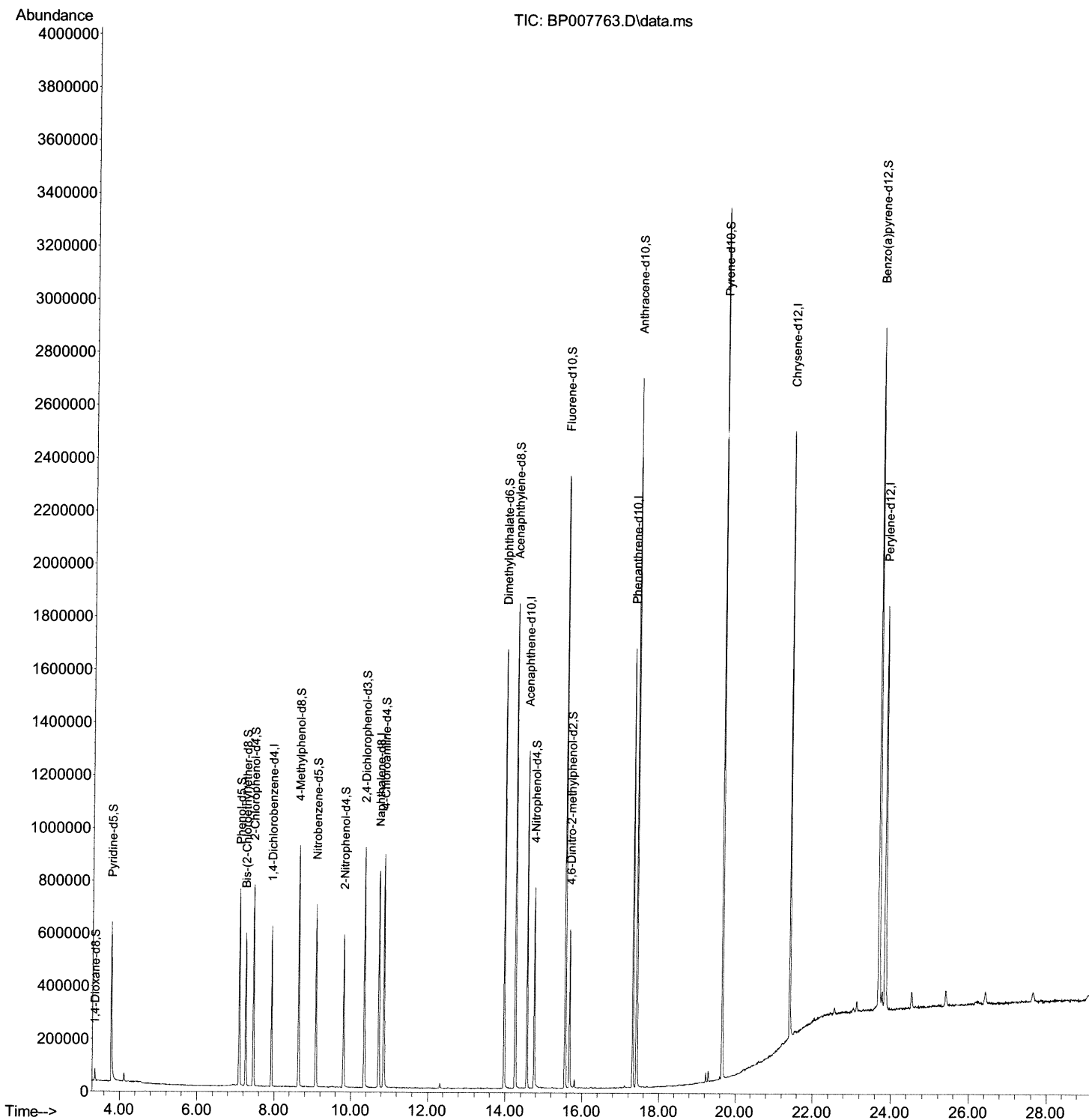
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102821\
Data File : BP007763.D
Acq On : 28 Oct 2021 21:17
Operator : CG/JU
Sample : PB140115BL
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SBLK115

Manual Integrations
APPROVED

mohammad
10/29/2021 2:25:32 PM

Quant Time: Oct 29 00:57:58 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



Quantitation Report (Qedit)

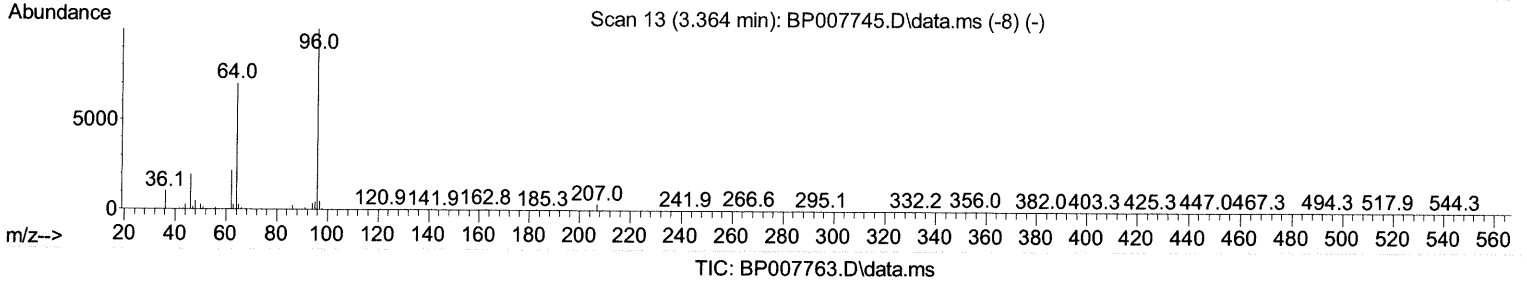
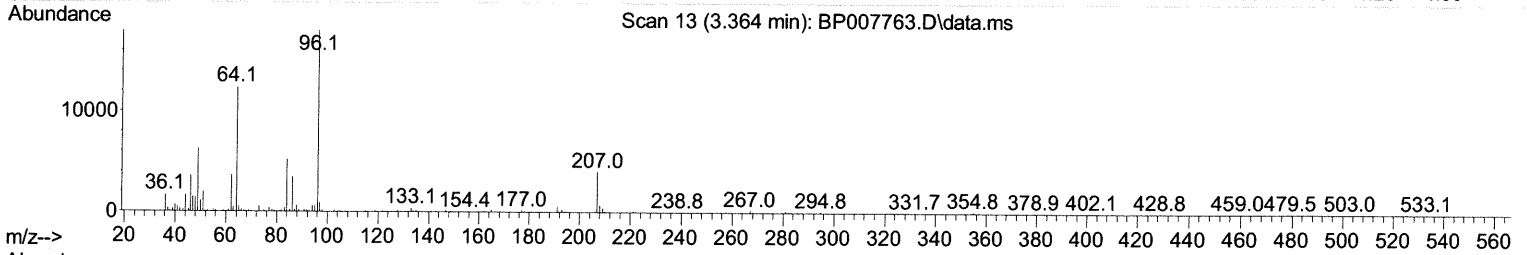
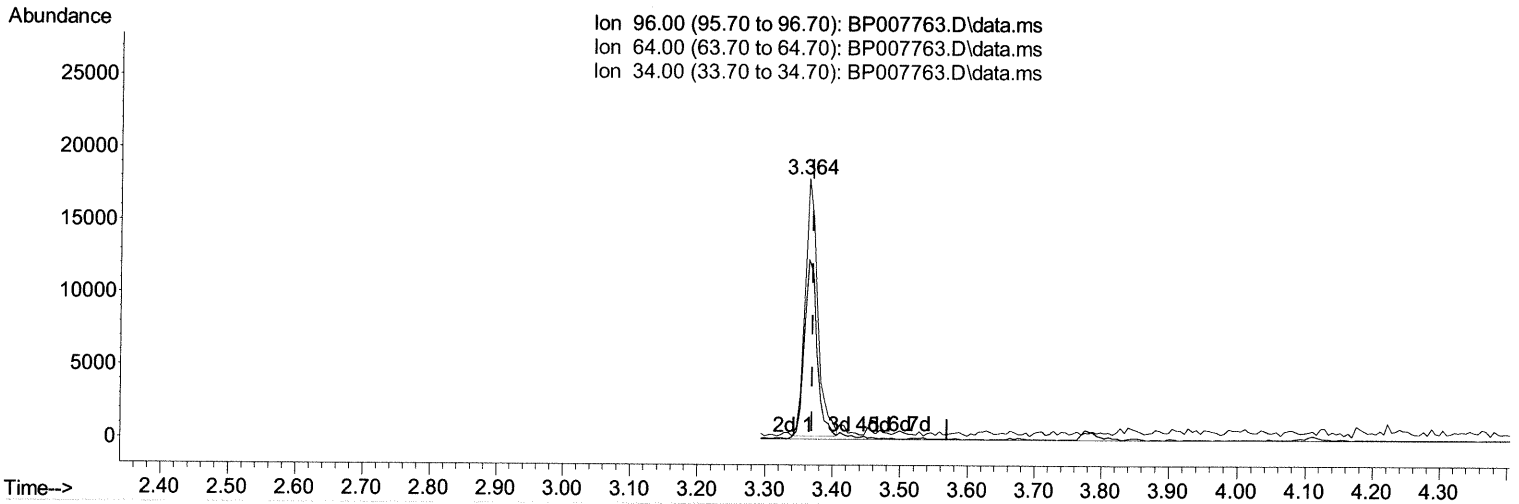
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TIC: BP007763.D\data.ms

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

3.364min (-0.006) 6.39 ng/uL

response 22993

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	69.20	68.98
34.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

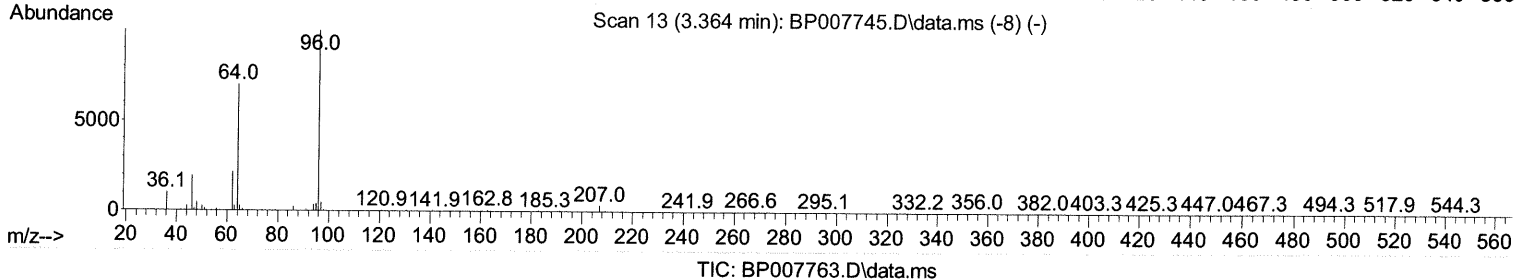
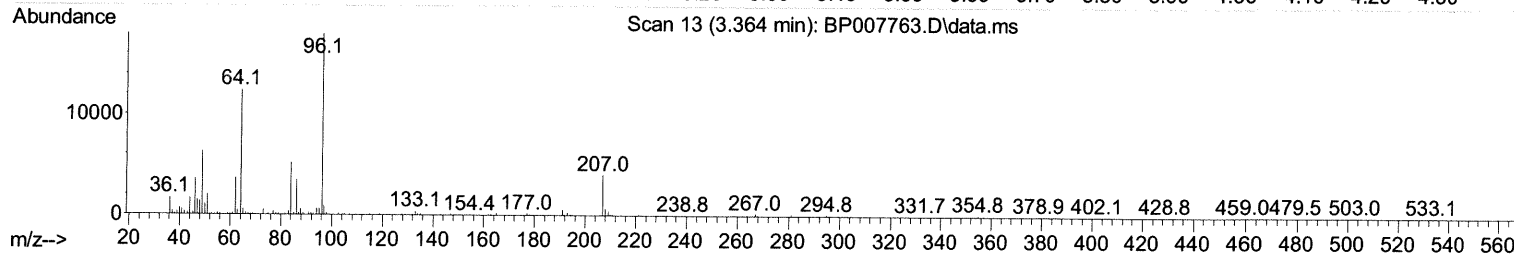
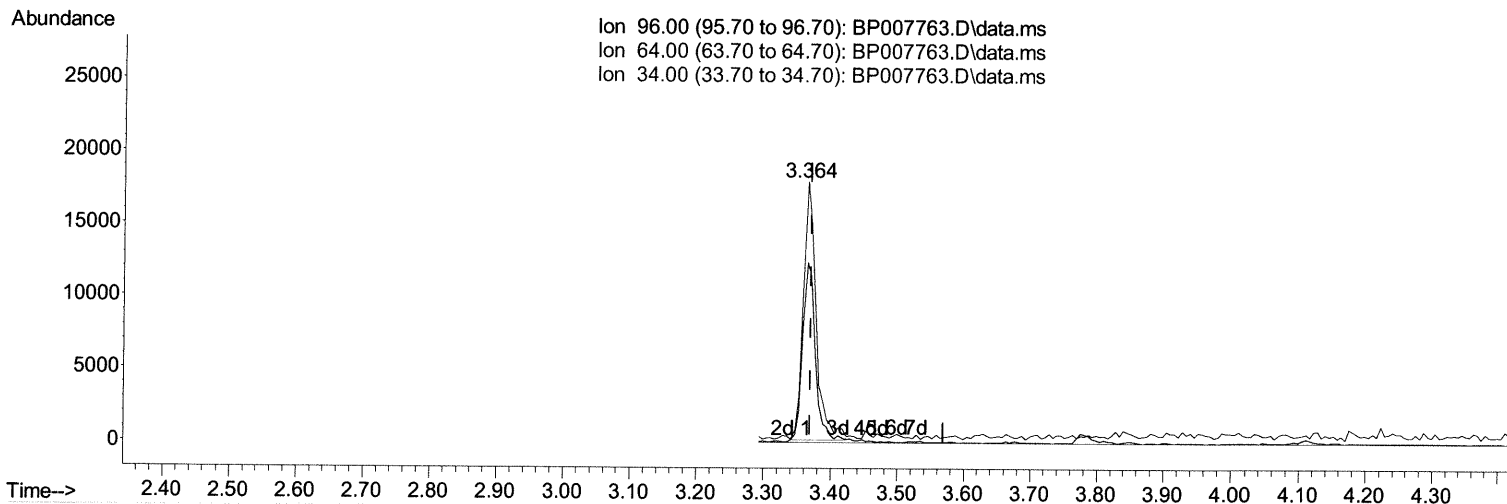
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Manual Integrations
APPROVED

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10/29/2021 2:25:32 PM

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Response via : Initial Calibration



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

3.364min (-0.006) 6.60 ng/uL m 10/30/21JU

response 23756

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	69.20	68.98
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	7.928	152	163552	20.000	ng/ul	0.00
20) Naphthalene-d8	10.734	136	677822	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.563	164	455689	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.322	188	1002119	20.000	ng/ul	0.00
79) Chrysene-d12	21.410	240	1132859	20.000	ng/ul	-0.01
88) Perylene-d12	23.851	264	1207038	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.364	96	23756m	> 6.598	ng/uL	> 0.00
4) Pyridine-d5	3.781	84	347630	31.928	ng/ul	0.00
7) Phenol-d5	7.093	99	423563	34.096	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.258	67	241202	34.972	ng/ul	0.00
11) 2-Chlorophenol-d4	7.458	132	350452	36.272	ng/ul	0.00
15) 4-Methylphenol-d8	8.634	113	338767	34.826	ng/ul	-0.01
21) Nitrobenzene-d5	9.087	128	168091	36.390	ng/ul	0.00
24) 2-Nitrophenol-d4	9.810	143	184757	36.621	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.352	165	342807	33.383	ng/ul	0.00
31) 4-Chloroaniline-d4	10.863	131	466652	34.368	ng/ul	0.00
46) Dimethylphthalate-d6	13.975	166	1097819	35.566	ng/ul	0.00
49) Acenaphthylene-d8	14.257	160	1278106	33.783	ng/ul	0.00
54) 4-Nitrophenol-d4	14.757	143	181090	32.005	ng/ul	0.00
60) Fluorene-d10	15.557	176	926977	35.623	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.675	200	138311	25.803	ng/ul	0.00
73) Anthracene-d10	17.422	188	1535504	37.285	ng/ul	0.00
81) Pyrene-d10	19.657	212	1866723	35.393	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.692	264	2055513	34.666	ng/ul	-0.02

Target Compounds Qvalue

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