

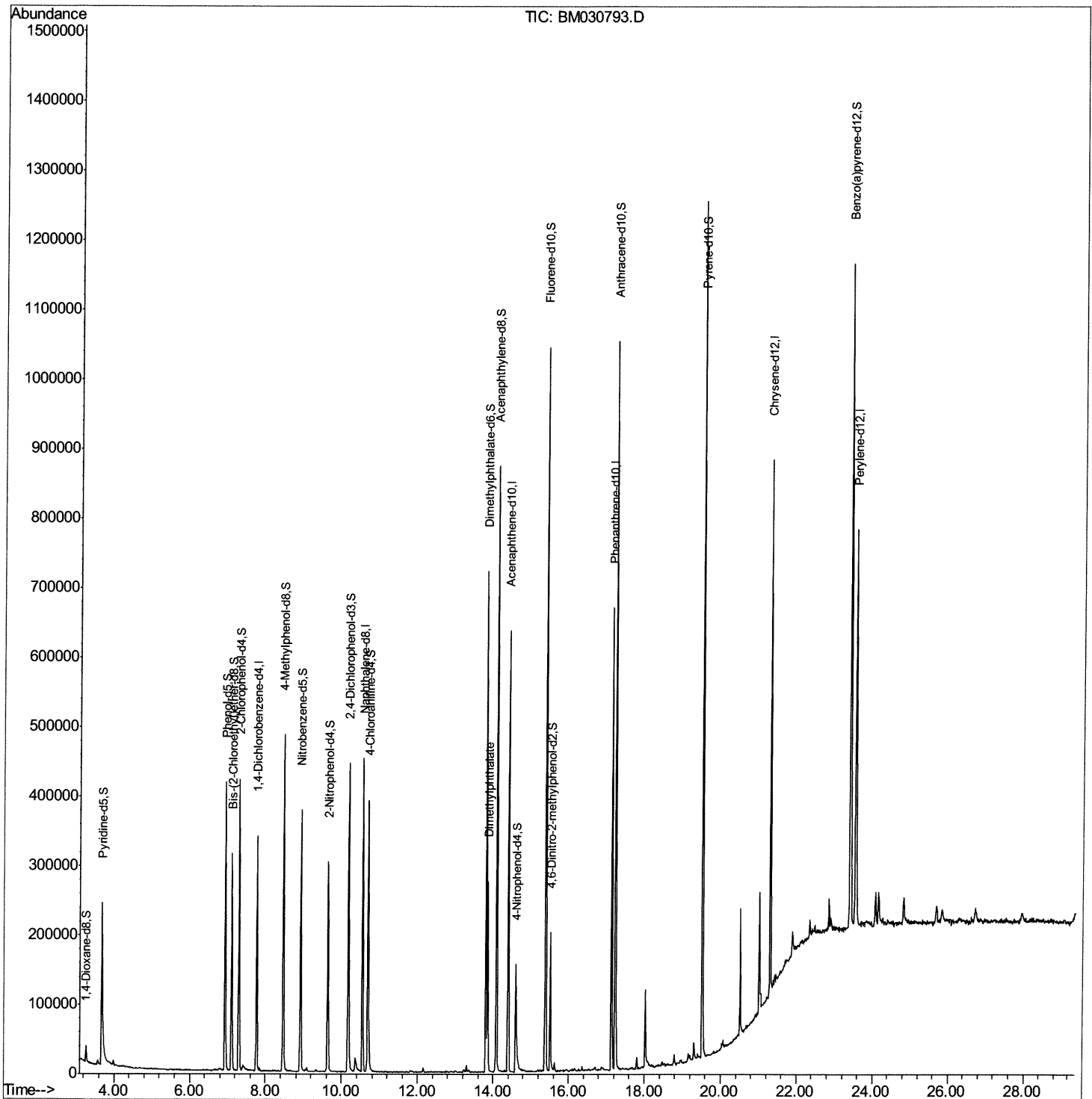
Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM063021\
Data File : BM030793.D
Acq On : 03 Jul 2021 02:35
Operator : CG/JU
Sample : M2869-14
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BGDY4

Manual Integrations
APPROVED

mohammad
7/6/2021 9:25:24 AM

Quant Time: Jul 03 03:11:01 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM062221.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat Jul 03 02:38:59 2021
Response via : Initial Calibration



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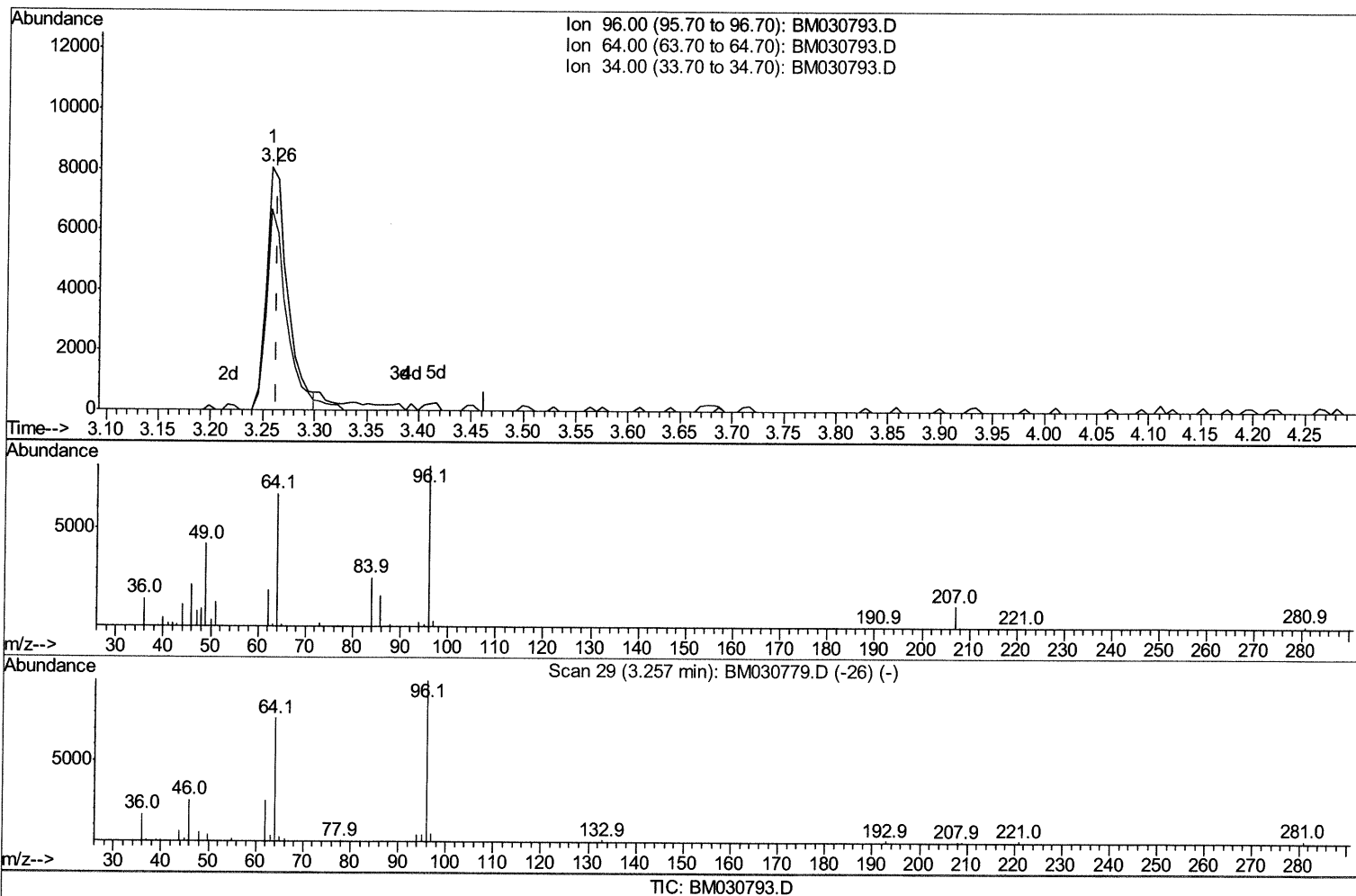
Quant Time: Jul 03 03:09:51 2021

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

3.258min (-0.006) 5.24ng/uL

response 11434

Ion	Exp%	Act%
96.00	100	100
64.00	80.00	82.83
34.00	0.00	0.00
0.00	0.00	0.00

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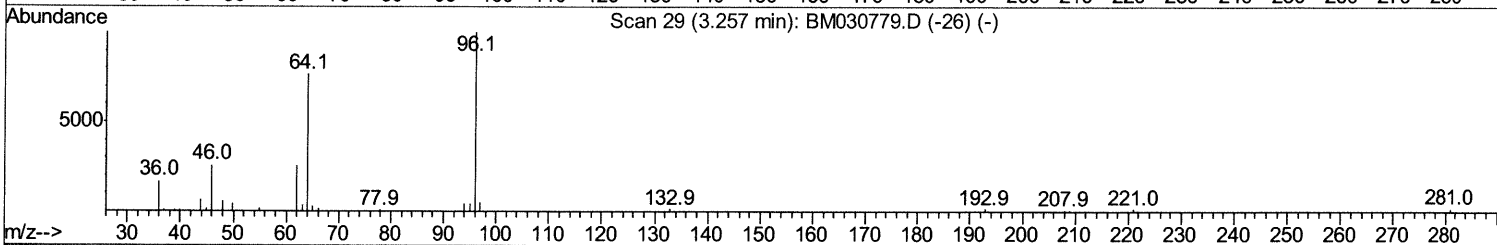
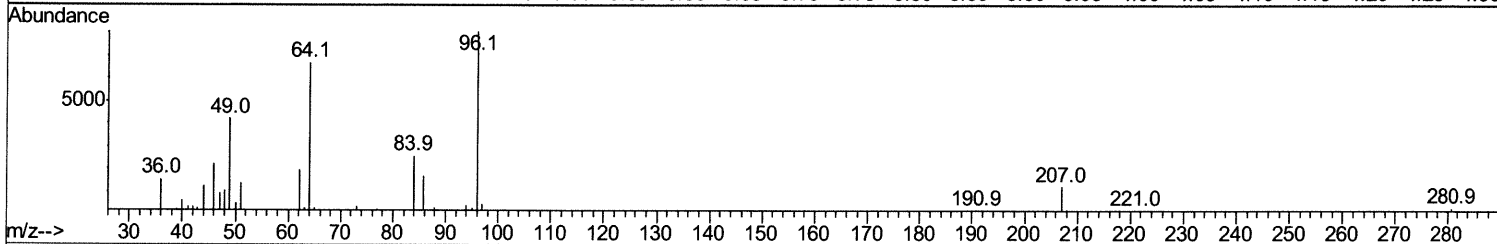
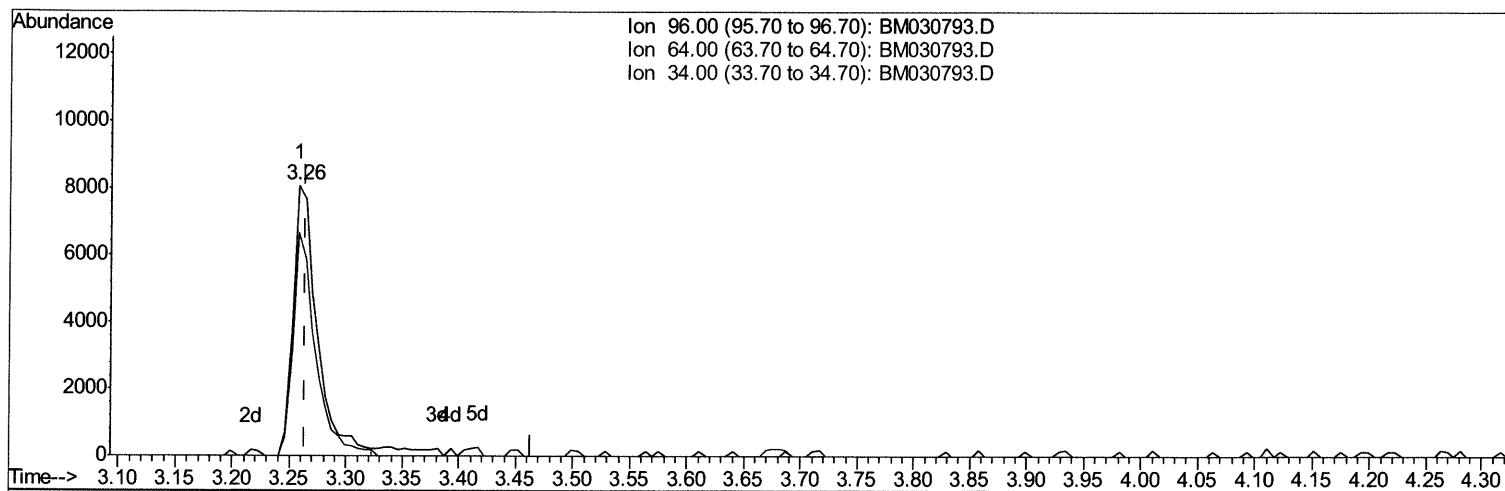
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TIC: BM030793.D

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

3.258min (-0.006) 5.47ng/uL m

response 11932

Ion	Exp%	Act%
96.00	100	100
64.00	80.00	82.83
34.00	0.00	0.00
0.00	0.00	0.00

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
1) 1,4-Dichlorobenzene-d4	7.77	152	88394	20.00	ng/ul	0.00
20) Naphthalene-d8	10.55	136	357994	20.00	ng/ul	0.00
38) Acenaphthene-d10	14.40	164	210389	20.00	ng/ul	0.00
64) Phenanthrene-d10	17.14	188	389359	20.00	ng/ul	0.00
79) Chrysene-d12	21.30	240	352660	20.00	ng/ul	0.00
88) Perylene-d12	23.55	264	386741	20.00	ng/ul	-0.01

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.26	96	11932m	5.47	ng/uL	0.00
4) Pyridine-d5	3.68	84	147213	25.13	ng/ul	0.00
7) Phenol-d5	6.93	99	217542	28.54	ng/ul	0.00
9) Bis-(2-Chloroethyl)ether-d	7.10	67	142129	29.66	ng/ul	0.00
11) 2-Chlorophenol-d4	7.30	132	179819	30.56	ng/ul	0.00
15) 4-Methylphenol-d8	8.47	113	176977	29.83	ng/ul	0.00
21) Nitrobenzene-d5	8.92	128	81433	30.48	ng/ul	0.00
24) 2-Nitrophenol-d4	9.64	143	88635	29.86	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.18	165	169094	29.73	ng/ul	0.00
31) 4-Chloroaniline-d4	10.69	131	217792	27.46	ng/ul	0.00
46) Dimethylphthalate-d6	13.81	166	476746	32.82	ng/ul	0.00
49) Acenaphthylene-d8	14.09	160	613116	32.51	ng/ul	0.00
54) 4-Nitrophenol-d4	14.60	143	51893	18.86	ng/ul	0.00
60) Fluorene-d10	15.39	176	415435	32.24	ng/ul	0.00
65) 4,6-Dinitro-2-methylphenol	15.52	200	43144	16.94	ng/ul	0.00
73) Anthracene-d10	17.23	188	579622	31.90	ng/ul	0.00
81) Pyrene-d10	19.52	212	612026	33.12	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.41	264	670228	33.33	ng/ul	0.00

Target Compounds				Ovalue	
47) Dimethylphthalate	13.86	163	173387	12.182	ng/ul 99

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