

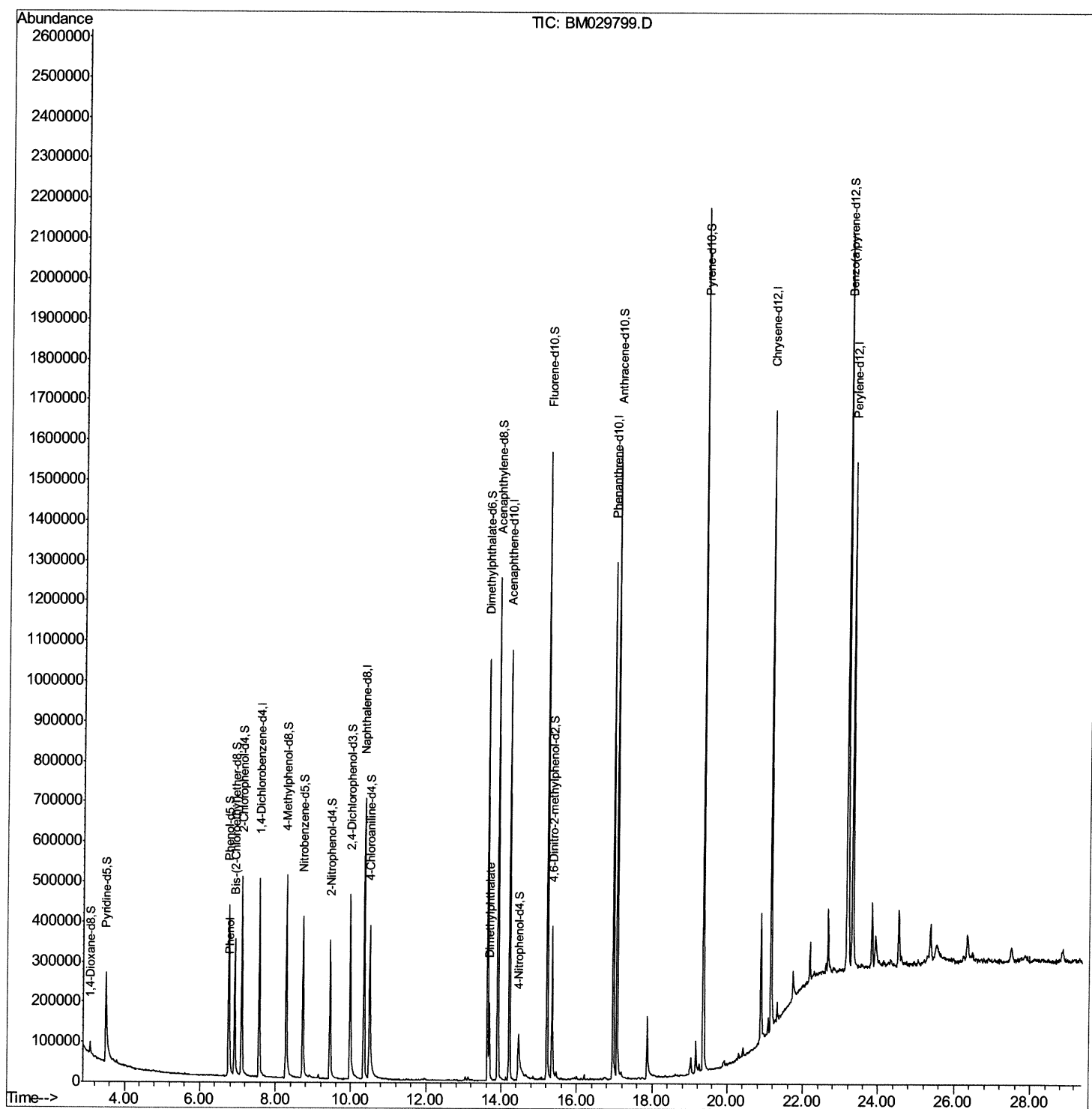
Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM050421\
Data File : BM029799.D
Acq On : 04 May 2021 21:38
Operator : CG/JU
Sample : M2121-07
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BG505

Manual Integrations
APPROVED

mohammad
5/5/2021 11:26:51 AM

Quant Time: May 05 02:39:59 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM050421.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue May 04 14:23:48 2021
Response via : Initial Calibration



Quantitation Report (Qedit)

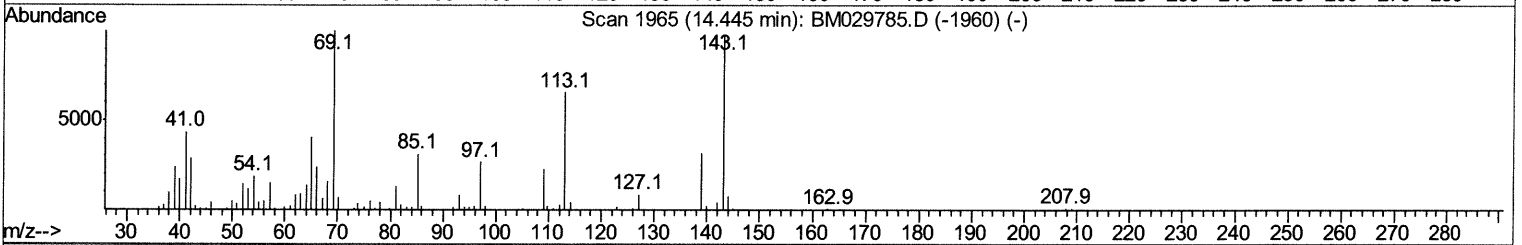
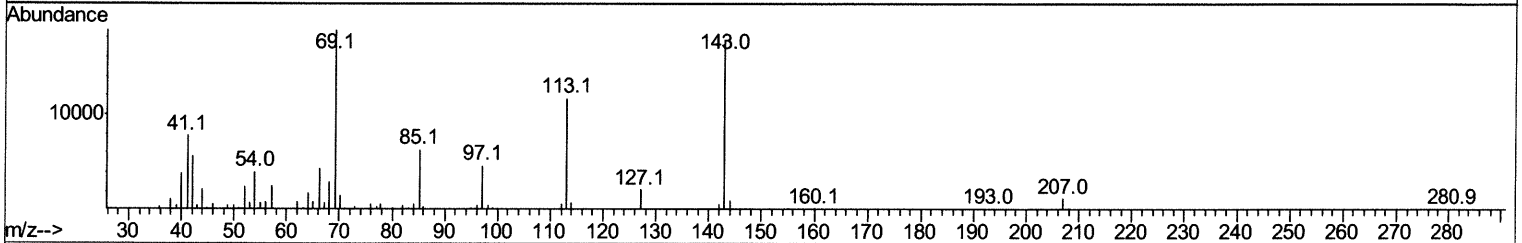
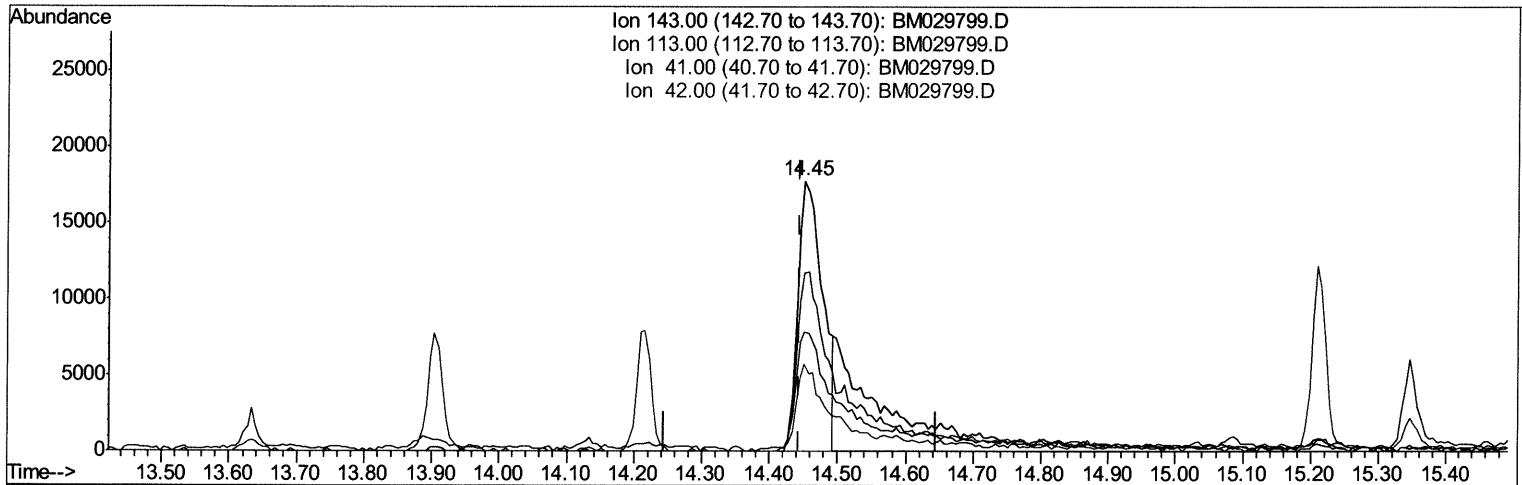
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Quant Time: Mav 05 02:39:02 2021
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QLast Update : Tue May 04 14:23:48 2021
Response via : Initial Calibration



TIC: BM029799.D

(54) 4-Nitrophenol-d4 (S)

14.451min (+0.006) 11.25ng/ul

response 44923

Ion	Exp%	Act%
143.00	100	100
113.00	67.10	65.76
41.00	45.40	44.03
42.00	30.10	32.09

Quantitation Report (Qedit)

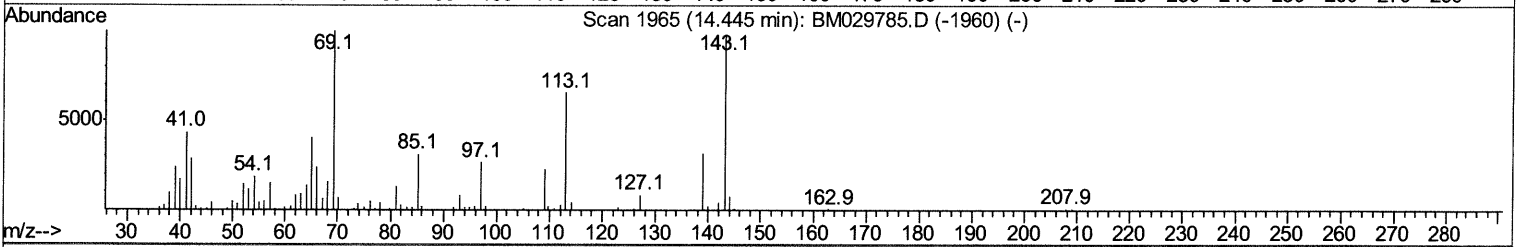
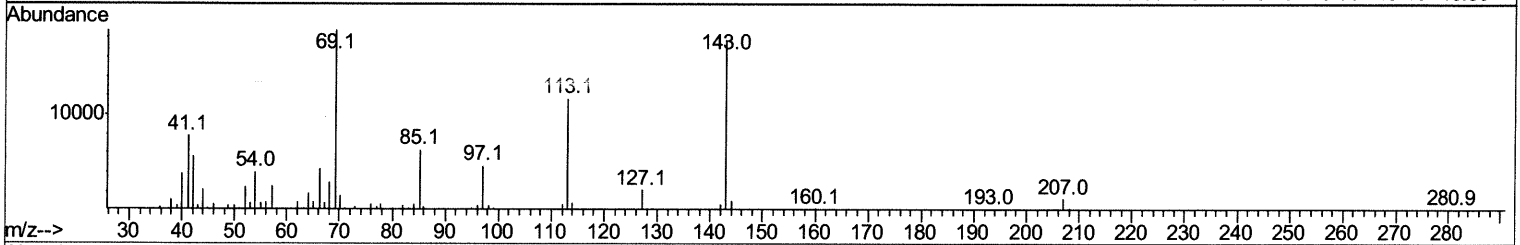
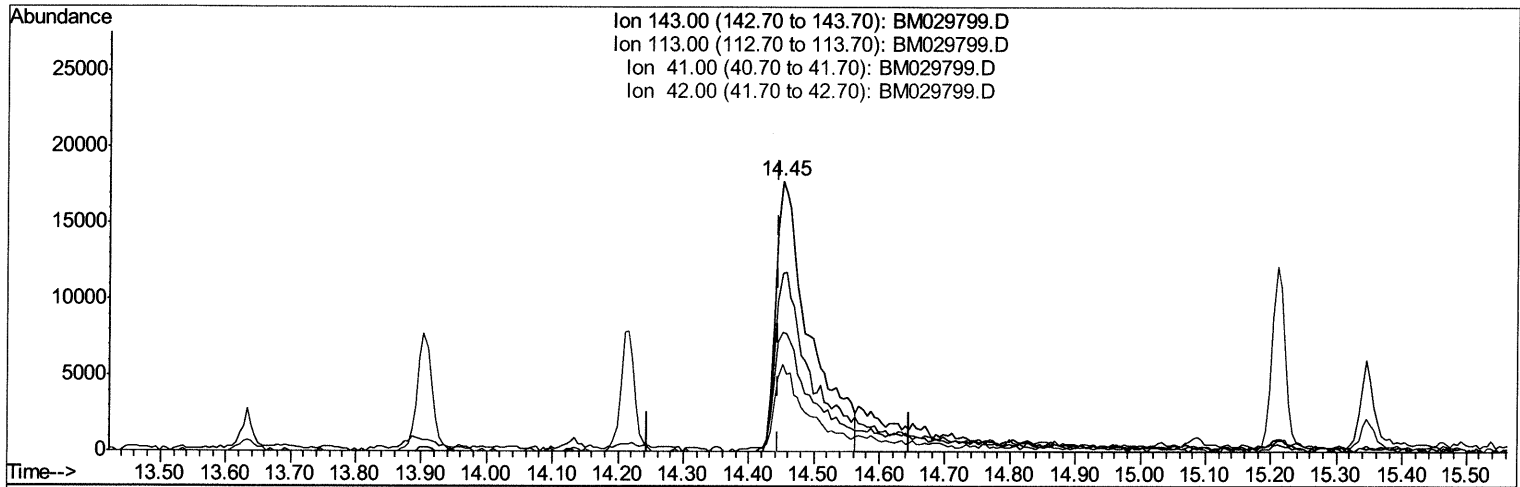
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TIC: BM029799.D

(54) 4-Nitrophenol-d4 (S)

14.451min (+0.006) 15.95ng/ul m 05/06/21 JU

response 63704

Ion	Exp%	Act%
143.00	100	100
113.00	67.10	65.76
41.00	45.40	44.03
42.00	30.10	32.09

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM050421\
 Data File : BM029799.D
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 Operator : CG/JU
 Sample : M2121-07
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BG505

Manual Integrations
 APPROVED

mohammad
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Quant Time: May 05 02:39:59 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM050421.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue May 04 14:23:48 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.57	152	142826	20.00	ng/ul	0.00
20) Naphthalene-d8	10.35	136	588489	20.00	ng/ul	0.00
38) Acenaphthene-d10	14.22	164	347544	20.00	ng/ul	0.00
64) Phenanthrene-d10	16.96	188	715260	20.00	ng/ul	0.00
79) Chrysene-d12	21.15	240	774354	20.00	ng/ul	0.00
88) Perylene-d12	23.31	264	857034	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.08	96	15820	4.10	ng/uL	0.00
4) Pividine-d5	3.50	84	171164	19.35	ng/ul	0.00
7) Phenol-d5	6.76	99	277523	22.38	ng/ul	0.00
9) Bis-(2-Chloroethyl)ether-d	6.92	67	176197	23.00	ng/ul	0.00
11) 2-Chlorophenol-d4	7.10	132	234065	23.69	ng/ul	0.00
15) 4-Methylphenol-d8	8.28	113	225903	22.82	ng/ul	0.00
21) Nitrobenzene-d5	8.72	128	105421	24.11	ng/ul	0.00
24) 2-Nitrophenol-d4	9.44	143	119983	25.13	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.97	165	197482	22.12	ng/ul	0.00
31) 4-Chloroaniline-d4	10.49	131	285348	21.93	ng/ul	0.00
46) Dimethylphthalate-d6	13.63	166	699225	26.94	ng/ul	0.00
49) Acenaphthylene-d8	13.91	160	895404	26.38	ng/ul	0.00
54) 4-Nitrophenol-d4	14.45	143	63704m >	15.95	ng/ul >	0.00
60) Fluorene-d10	15.22	176	610780	27.65	ng/ul	0.00
65) 4,6-Dinitro-2-methylphenol	15.35	200	85815	17.87	ng/ul	0.00
73) Anthracene-d10	17.06	188	922868	26.09	ng/ul	0.00
81) Pyrene-d10	19.36	212	1069872	26.93	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.17	264	1273215	26.46	ng/ul	0.00

05/06/21 JU

Target Compounds

				Ovalue	
8) Phenol	6.78	94	61473	4.871	ng/ul 99
47) Dimethylphthalate	13.68	163	124946	4.824	ng/ul 100

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