

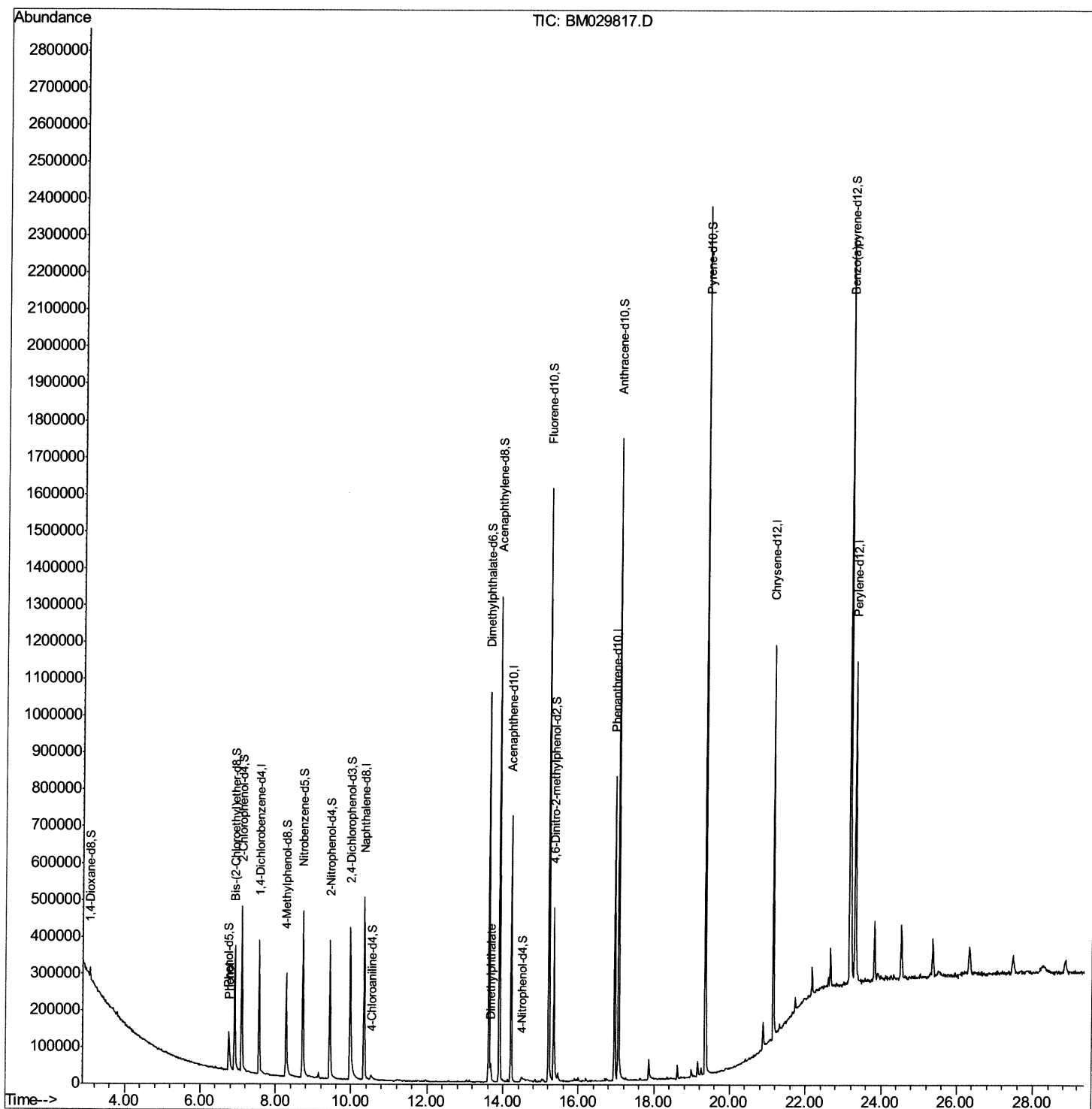
Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM050421\
Data File : BM029817.D
Acq On : 05 May 2021 09:06
Operator : CG/JU
Sample : M2127-09
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
E0AB2

Manual Integrations
APPROVED

mohammad
5/5/2021 11:27:30 AM

Quant Time: May 05 10:07:01 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)

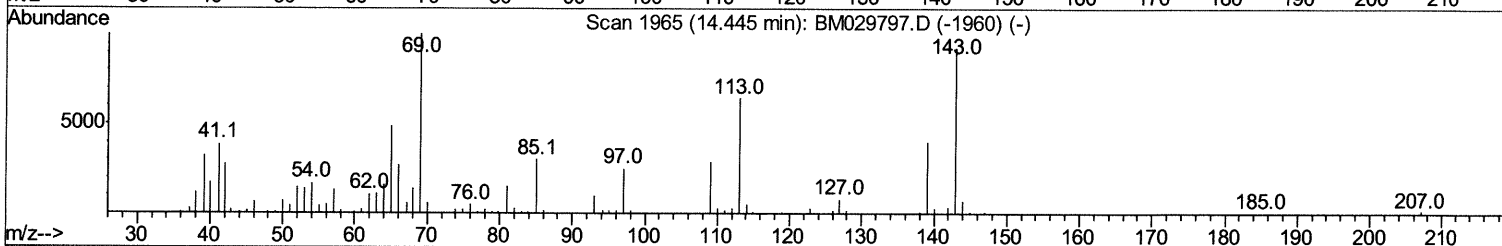
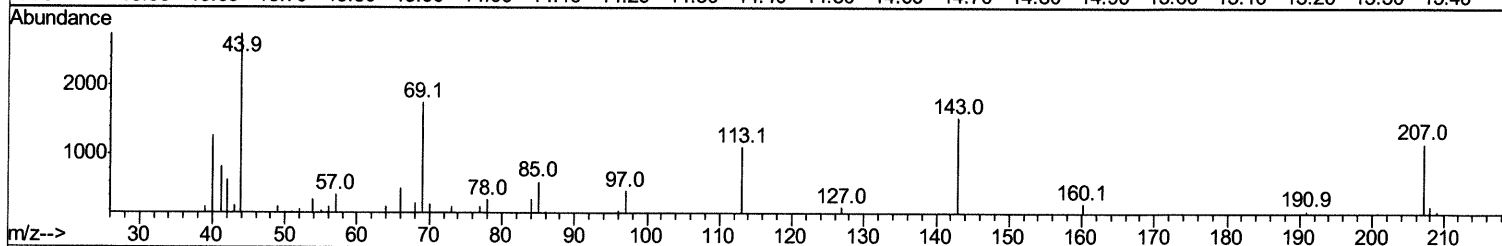
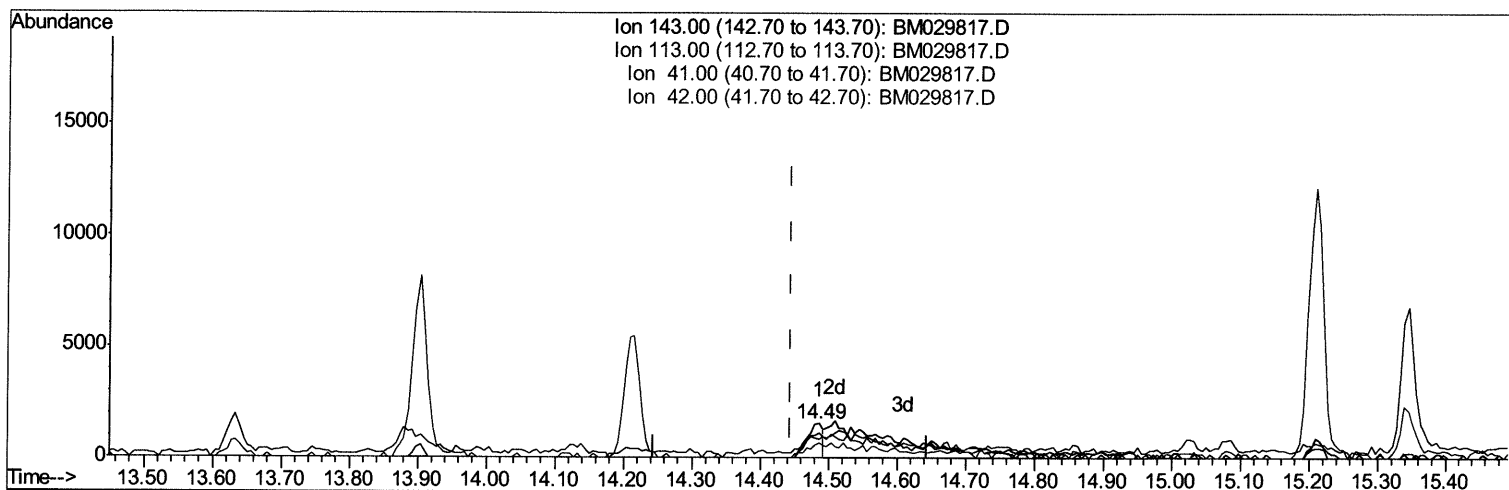
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ClientSampled :
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Manual Integrations
APPROVED

mohammad
 5/5/2021 11:27:30 AM

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 Response via : Initial Calibration



TIC: BM029817.D

(54) 4-Nitrophenol-d4 (S)

14.486min (+0.041) 0.90ng/ul

response 2539

Ion	Exp%	Act%
143.00	100	100
113.00	67.10	72.50
41.00	45.40	54.60#
42.00	30.10	41.87#

Quantitation Report (Qedit)

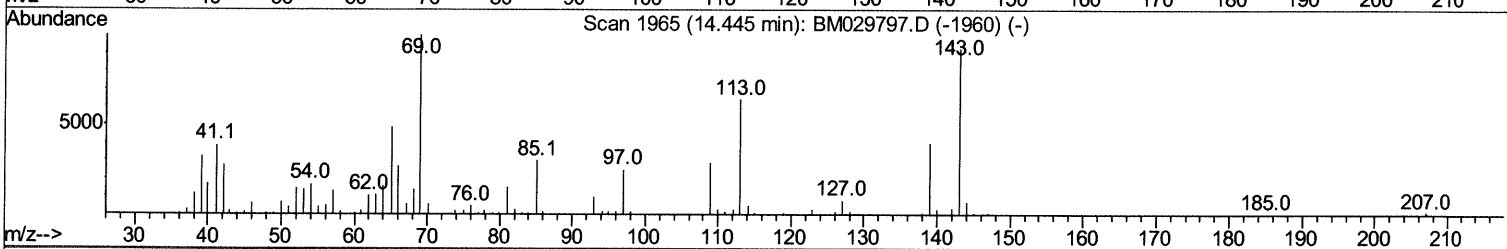
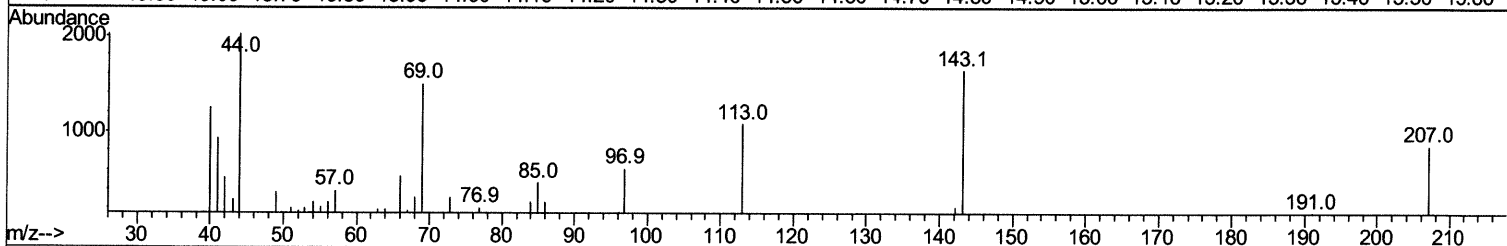
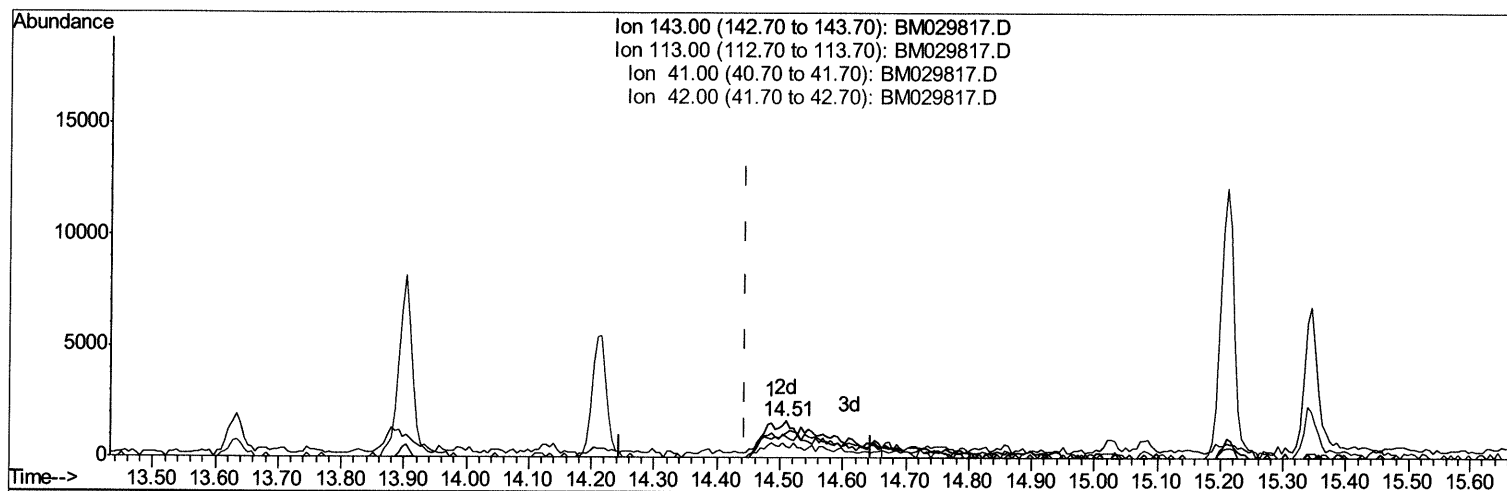
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Instrument :
BNA_M
ClientSampleId :
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Manual Integrations
APPROVED

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Response via : Initial Calibration



TIC: BM029817.D

(54) 4-Nitrophenol-d4 (S)

14.510min (+0.065) 4.37ng/ul m 05/06/21 JU

response 12362

Ion	Exp%	Act%
143.00	100	100
113.00	67.10	65.77
41.00	45.40	57.06#
42.00	30.10	31.91

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM050421\
 Data File : BM029817.D
 Acq On : 05 May 2021 09:06
 Operator : CG/JU
 Sample : M2127-09
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 E0AB2

Manual Integrations
 APPROVED

mohammad
 5/5/2021 11:27:30 AM

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.57	152	106465	20.00	ng/ul	0.00
20) Naphthalene-d8	10.34	136	407854	20.00	ng/ul	0.00
38) Acenaphthene-d10	14.22	164	246024	20.00	ng/ul	0.00
64) Phenanthrene-d10	16.96	188	491209	20.00	ng/ul	0.00
79) Chrysene-d12	21.14	240	518667	20.00	ng/ul	0.00
88) Perylene-d12	23.31	264	614379	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.08	96	15933	5.53	ng/uL	0.00
4) Pyridine-d5	0.00	84	0d	0.00	ng/ul	
7) Phenol-d5	6.76	99	73479	7.95	ng/ul	0.00
9) Bis-(2-Chloroethyl) ether-d	6.91	67	185484	32.48	ng/ul	0.00
11) 2-Chlorophenol-d4	7.10	132	211273	28.69	ng/ul	0.00
15) 4-Methylphenol-d8	8.28	113	127638	17.30	ng/ul	0.00
21) Nitrobenzene-d5	8.72	128	113028	37.30	ng/ul	0.00
24) 2-Nitrophenol-d4	9.44	143	127033	38.39	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.97	165	194424	31.42	ng/ul	0.00
31) 4-Chloroaniline-d4	10.53	131	10209	1.13	ng/ul	0.04
46) Dimethylphthalate-d6	13.63	166	707367	38.49	ng/ul	0.00
49) Acenaphthylene-d8	13.90	160	913645	38.02	ng/ul	0.00
54) 4-Nitrophenol-d4	14.51	143	12362m>	4.37	ng/ul>	0.06
60) Fluorene-d10	15.21	176	621461	39.75	ng/ul	0.00
65) 4,6-Dinitro-2-methylphenol	15.34	200	97569	29.59	ng/ul	0.00
73) Anthracene-d10	17.06	188	991032	40.79	ng/ul	0.00
81) Pyrene-d10	19.36	212	1162637	43.69	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.17	264	1408664	40.83	ng/ul	0.00

05/06/21 JU

Target Compounds

					Ovalue
8) Phenol	6.78	94	25966	2.760	ng/ul 97
47) Dimethylphthalate	13.68	163	25716	1.403	ng/ul 98

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