

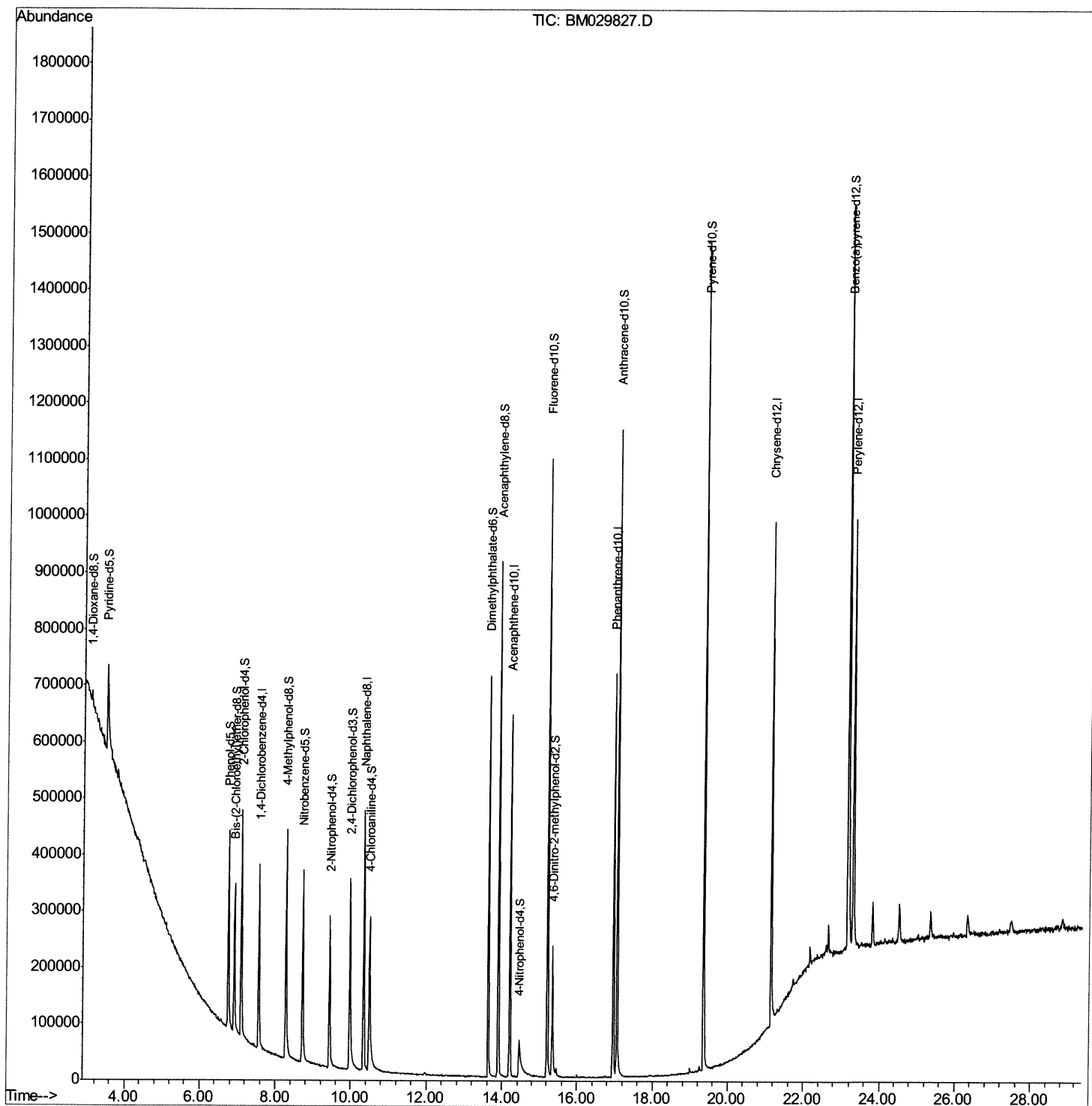
Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM050421\
Data File : BM029827.D
Acq On : 05 May 2021 15:18
Operator : CG/JU
Sample : PB135972BL
Misc :
ALS Vial : 46 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
SBLK972

Manual Integrations
APPROVED

mohammad
5/6/2021 4:48:51 PM

Quant Time: May 05 15:52:20 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)

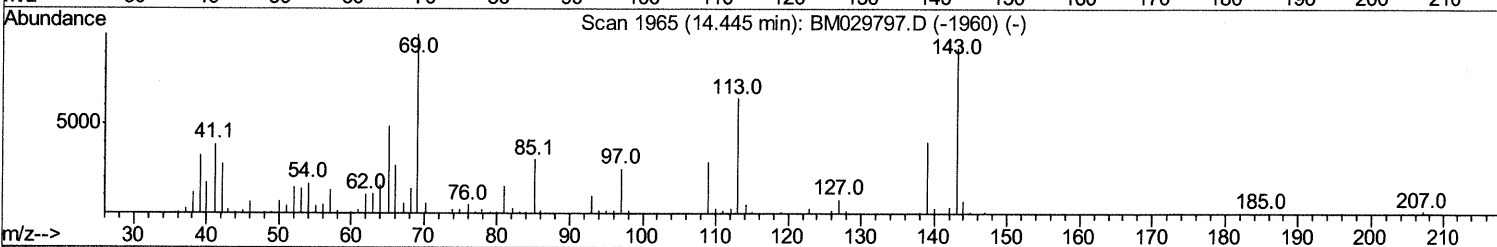
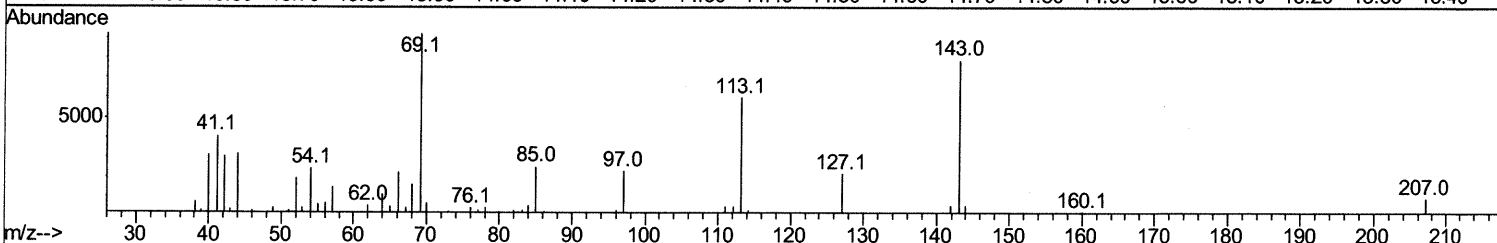
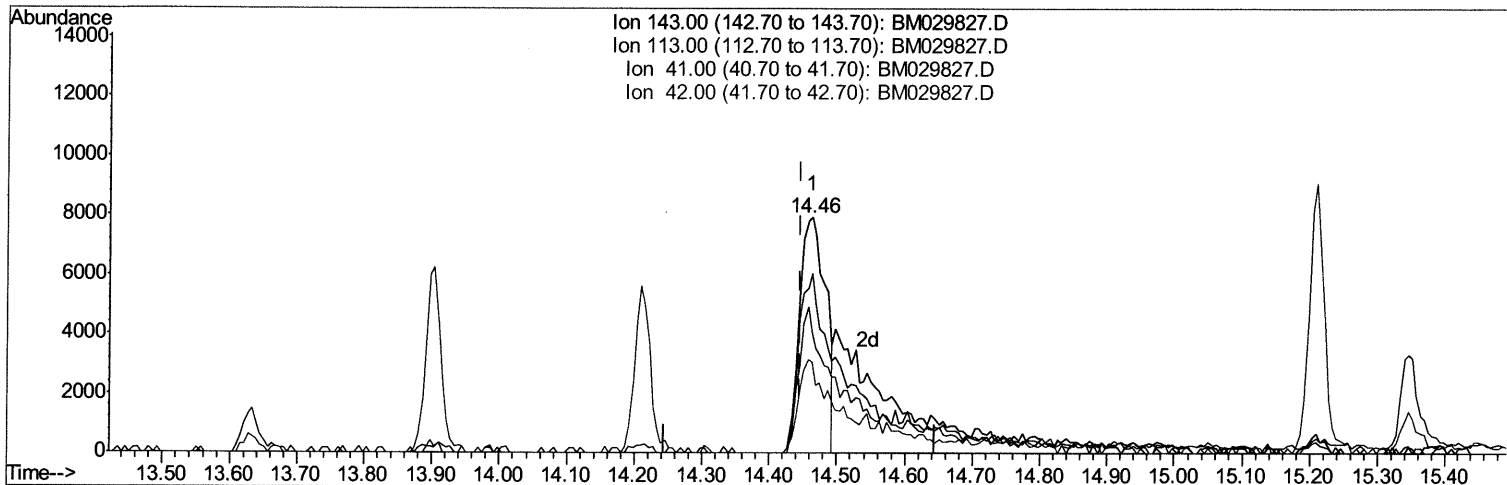
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TIC: BM029827.D

(54) 4-Nitrophenol-d4 (S)

14.463min (+0.018) 8.77ng/ul

response 21303

Ion	Exp%	Act%
143.00	100	100
113.00	67.10	76.05
41.00	45.40	51.34
42.00	30.10	38.56#

Quantitation Report (Qedit)

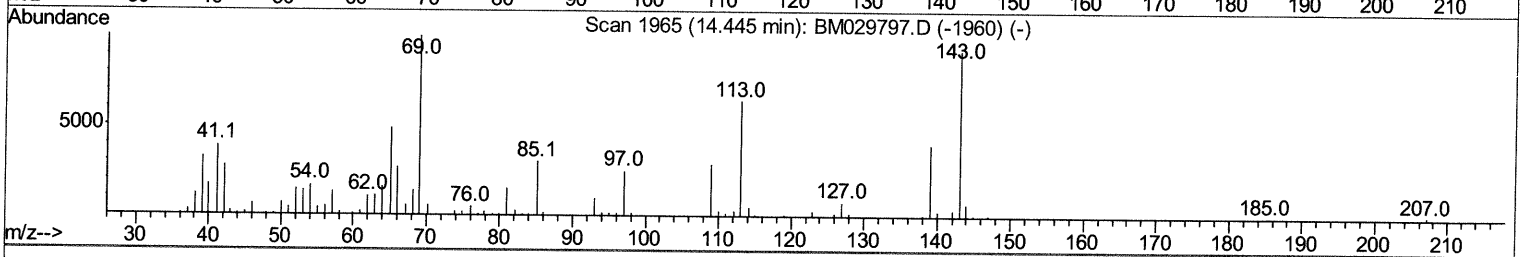
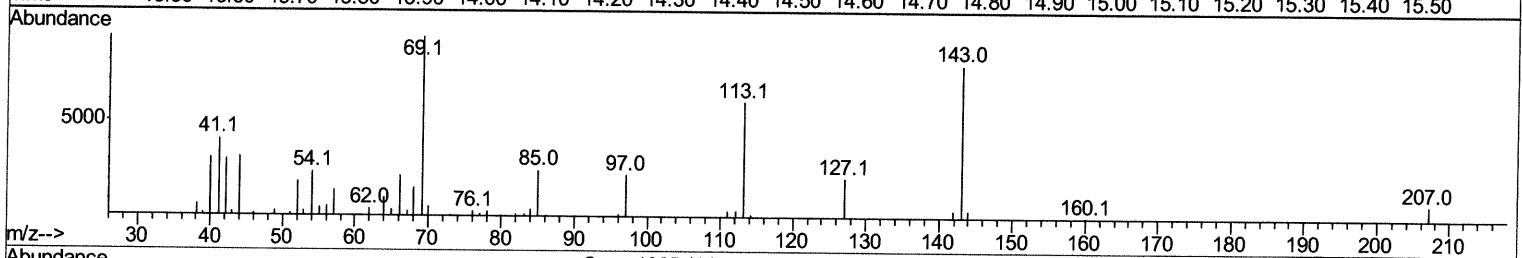
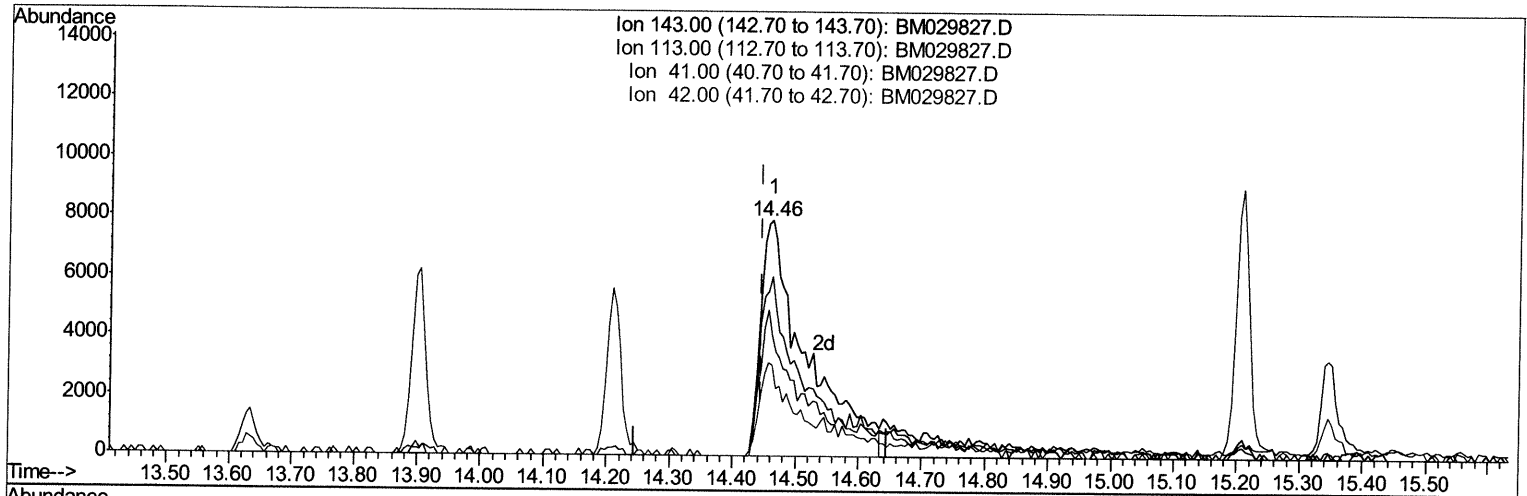
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Manual Integrations
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Response via : Initial Calibration



TIC: BM029827.D

(54) 4-Nitrophenol-d4 (S)

14.463min (+0.018) 16.39ng/ul m 05/07/21 JU

response 39796

Ion	Exp%	Act%
143.00	100	100
113.00	67.10	76.05
41.00	45.40	51.34
42.00	30.10	38.56#

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM050421\
 Data File : BM029827.D
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 Operator : CG/JU
 Sample : PB135972BL
 Misc :
 ALS Vial : 46 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
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Manual Integrations
 APPROVED

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.57	152	93224	20.00	ng/ul	0.00
20) Naphthalene-d8	10.34	136	375119	20.00	ng/ul	0.00
38) Acenaphthene-d10	14.21	164	211322	20.00	ng/ul	0.00
64) Phenanthrene-d10	16.96	188	417005	20.00	ng/ul	0.00
79) Chrysene-d12	21.14	240	437064	20.00	ng/ul	0.00
88) Perylene-d12	23.30	264	514008	20.00	ng/ul	-0.01

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.08	96	11813	4.69	ng/uL	0.00
4) Pyridine-d5	3.50	84	131004	22.69	ng/ul	0.00
7) Phenol-d5	6.75	99	224195	27.70	ng/ul	0.00
9) Bis-(2-Chloroethyl) ether-d	6.91	67	140139	28.03	ng/ul	0.00
11) 2-Chlorophenol-d4	7.10	132	190398	29.53	ng/ul	0.00
15) 4-Methylphenol-d8	8.28	113	171739	26.58	ng/ul	0.00
21) Nitrobenzene-d5	8.72	128	81482	29.23	ng/ul	0.00
24) 2-Nitrophenol-d4	9.44	143	86381	28.38	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.97	165	137714	24.20	ng/ul	0.00
31) 4-Chloroaniline-d4	10.49	131	194411	23.44	ng/ul	0.00
46) Dimethylphthalate-d6	13.63	166	483999	30.66	ng/ul	0.00
49) Acenaphthylene-d8	13.90	160	616452	29.86	ng/ul	0.00
54) 4-Nitrophenol-d4	14.46	143	39796m>	16.39	ng/ul>	0.02 05/07/21 JU
60) Fluorene-d10	15.21	176	409903	30.52	ng/ul	0.00
65) 4,6-Dinitro-2-methylphenol	15.34	200	48435	17.30	ng/ul	0.00
73) Anthracene-d10	17.06	188	640568	31.06	ng/ul	0.00
81) Pyrene-d10	19.36	212	735081	32.78	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.17	264	902731	31.28	ng/ul	0.00

Target Compounds Ovalue

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