

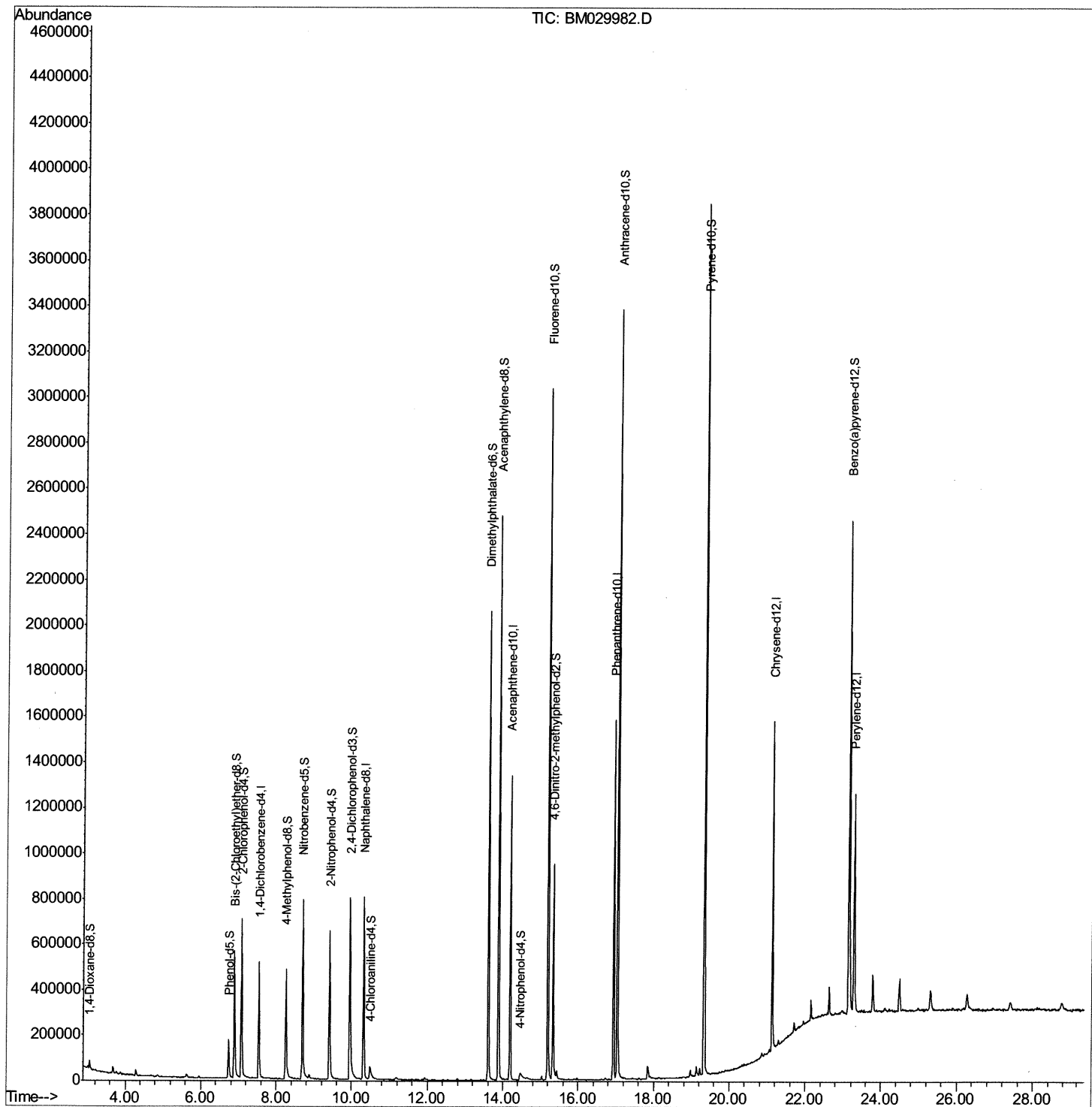
Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM051021\
Data File : BM029982.D
Acq On : 12 May 2021 15:33
Operator : CG/JU
Sample : M2197-15
Misc :
ALS Vial : 81 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
YARR8

Manual Integrations
APPROVED

Sohil
5/13/2021 10:30:42 AM

Quant Time: May 12 16:06:52 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM050721.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue May 11 15:21:38 2021
Response via : Initial Calibration



Quantitation Report (Qedit)

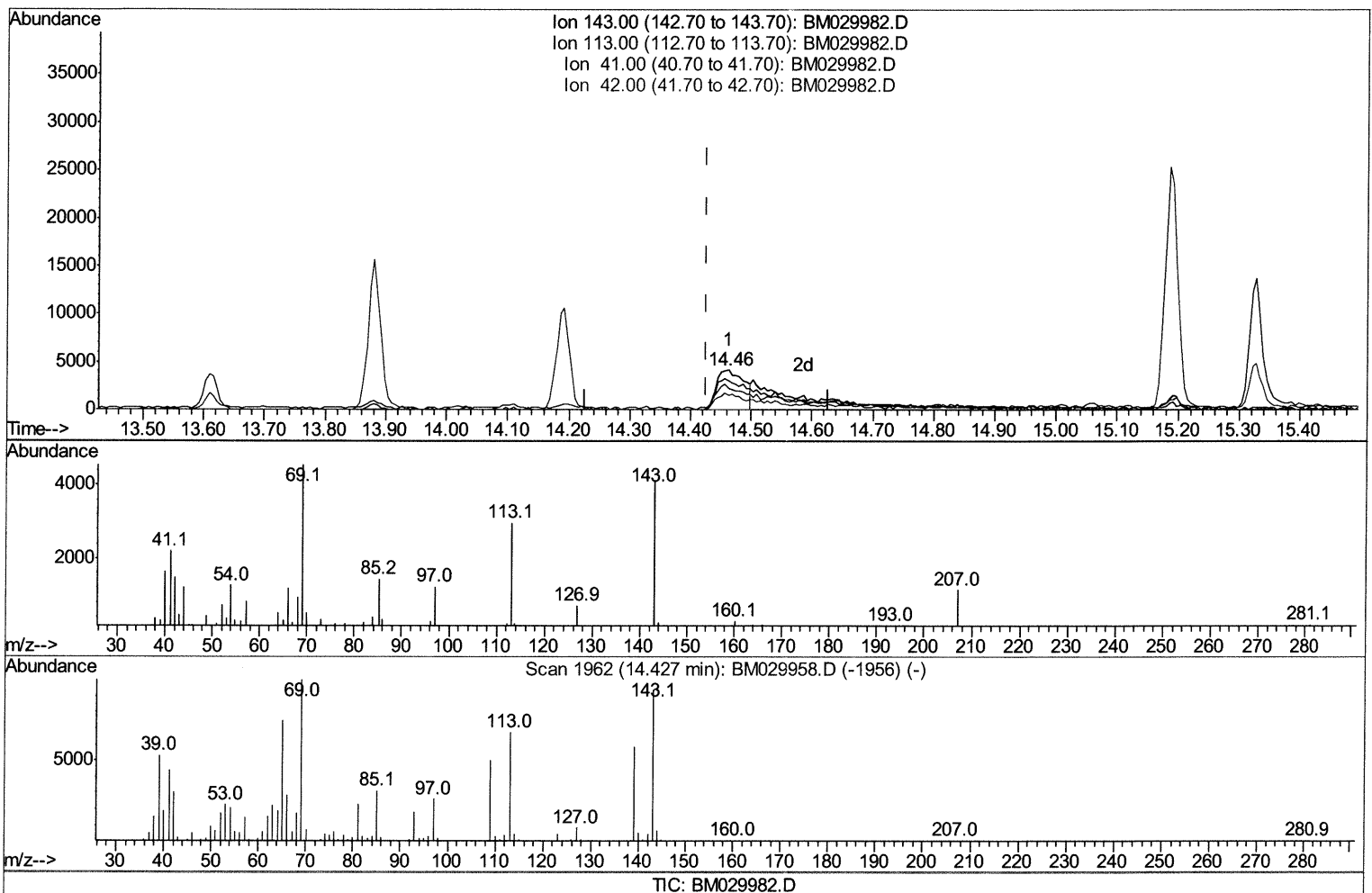
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ClientSampled :
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Manual Integrations
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Sohil
5/13/2021 10:30:42 AM

Quant Time: May 12 16:04:55 2021
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Response via : Initial Calibration



(54) 4-Nitrophenol-d4 (S)

14.463min (+0.035) 2.42ng/ul

response 13123

Ion	Exp%	Act%
143.00	100	100
113.00	71.10	72.12
41.00	44.70	53.38
42.00	28.30	36.04#

Quantitation Report (Qedit)

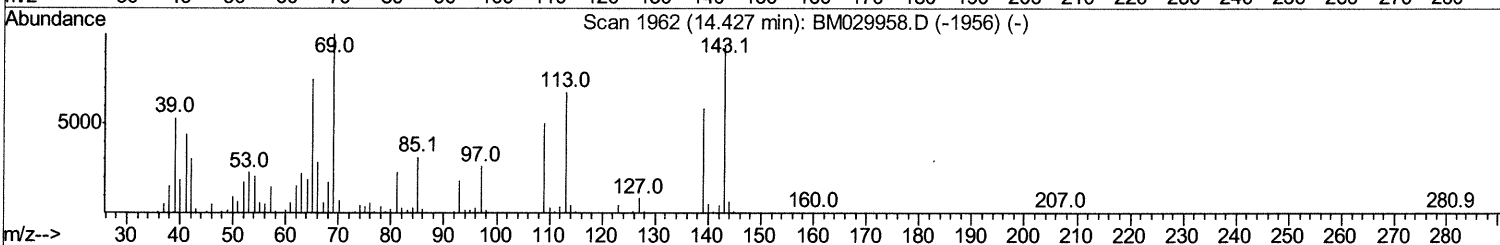
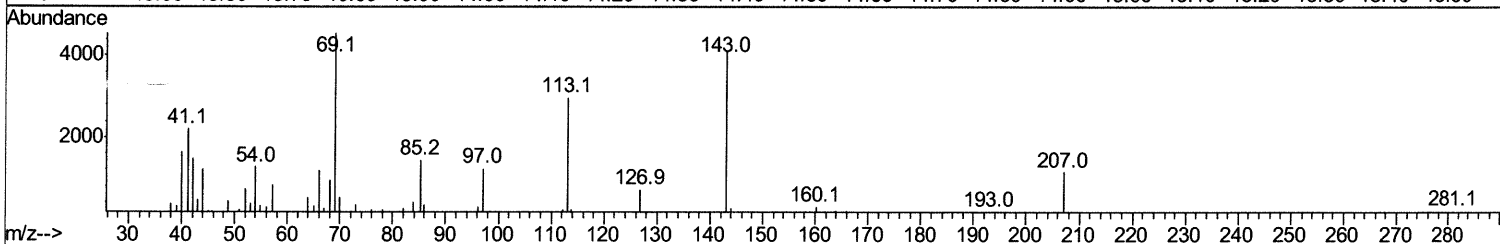
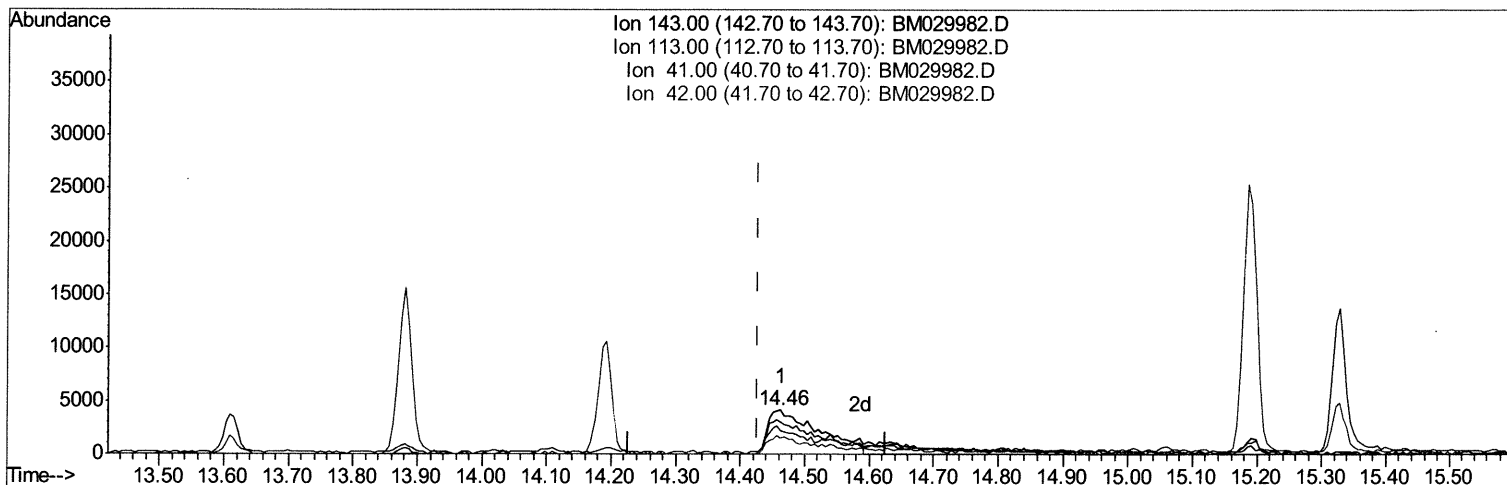
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Instrument :
 BNA_M
 ClientSampleId :
 YARR8

Manual Integrations
 APPROVED

Sohil
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Quant Time: May 12 16:04:55 2021
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 Quant Title : SVOA CALIBRATION
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 Response via : Initial Calibration



TIC: BM029982.D

(54) 4-Nitrophenol-d4 (S)

14.463min (+0.035) 4.24ng/ul m 05/13/21

response 23005

Ion	Exp%	Act%
143.00	100	100
113.00	71.10	72.12
41.00	44.70	53.38
42.00	28.30	36.04#

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM051021\
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 Sample : M2197-15
 Misc :
 ALS Vial : 81 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 YARR8

Manual Integrations
 APPROVED

Sohil
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 Quant Title : SVOA CALIBRATION
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 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.55	152	140831	20.00	ng/ul	0.00
20) Naphthalene-d8	10.32	136	636392	20.00	ng/ul	-0.01
38) Acenaphthene-d10	14.19	164	430145	20.00	ng/ul	-0.01
64) Phenanthrene-d10	16.94	188	873684	20.00	ng/ul	0.00
79) Chrysene-d12	21.13	240	713986	20.00	ng/ul	0.00
88) Perylene-d12	23.28	264	614198	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.06	96	20869	5.92	ng/uL	0.00
4) Pyridine-d5	0.00	84	0d	0.00	ng/ul	
7) Phenol-d5	6.73	99	112209	8.57	ng/ul	0.00
9) Bis-(2-Chloroethyl)ether-d	6.89	67	299716	36.90	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.08	132	320913	31.75	ng/ul	0.00
15) 4-Methylphenol-d8	8.26	113	197983	18.28	ng/ul	-0.01
21) Nitrobenzene-d5	8.70	128	181295	41.34	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.42	143	207360	41.89	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.95	165	340493	34.31	ng/ul	0.00
31) 4-Chloroaniline-d4	10.49	131	45332	3.00	ng/ul	0.01
46) Dimethylphthalate-d6	13.61	166	1338613	40.15	ng/ul	-0.01
49) Acenaphthylene-d8	13.88	160	1627110	38.51	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.46	143	23005m>	4.24	ng/ul>	0.0405/13/21JU
60) Fluorene-d10	15.19	176	1154724	40.80	ng/ul	0.00
65) 4,6-Dinitro-2-methylphenol	15.33	200	189698	33.43	ng/ul	0.00
73) Anthracene-d10	17.04	188	1796774	40.84	ng/ul	0.00
81) Pyrene-d10	19.33	212	1932086	52.75	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.14	264	1419078	40.76	ng/ul	0.00

Target Compounds Ovalue

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