

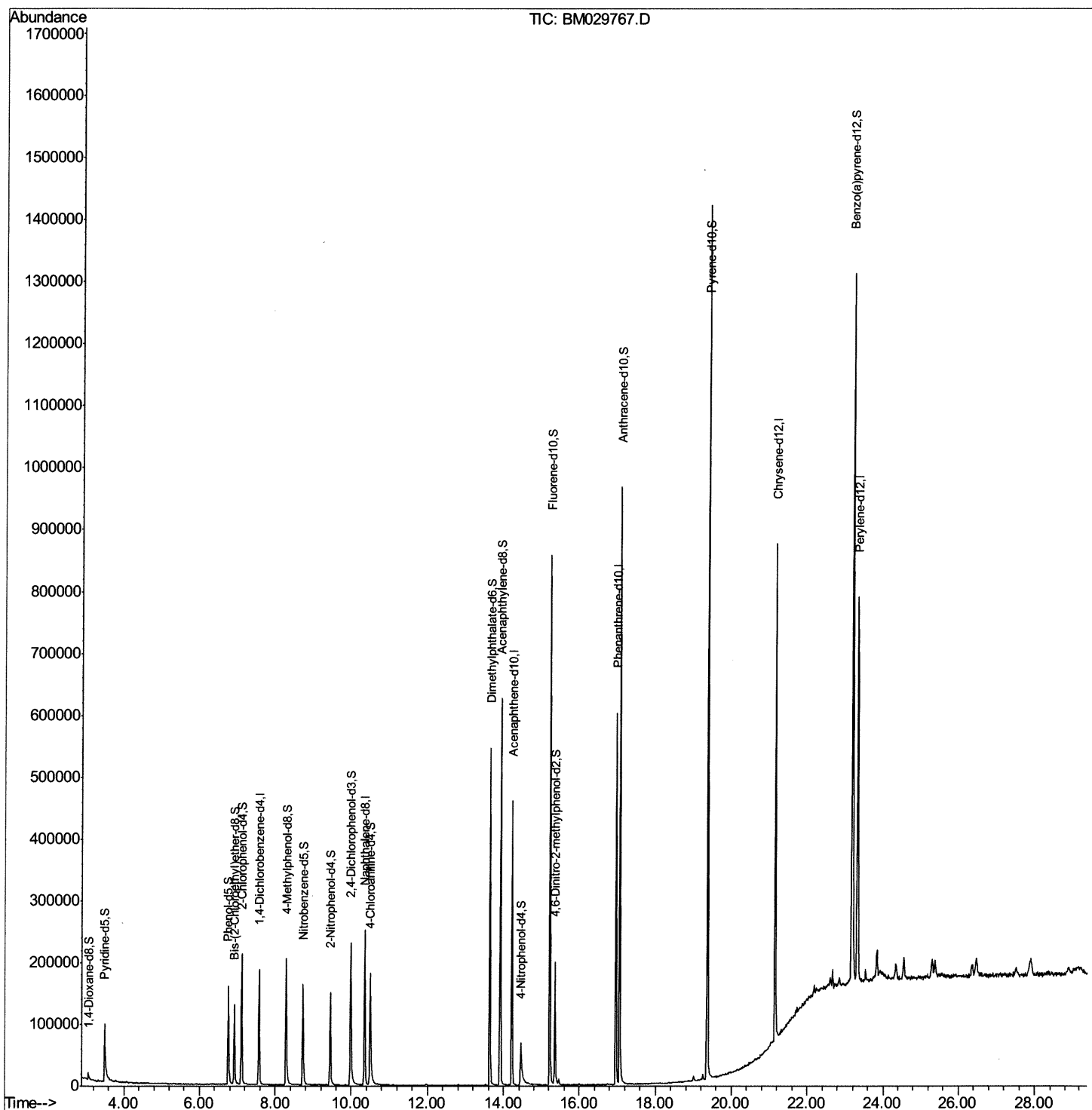
Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM050121\
Data File : BM029767.D
Acq On : 02 May 2021 04:29
Operator : CG/JU
Sample : PB135971BL
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
SBLK971

Manual Integrations
APPROVED

mohammad
5/3/2021 12:25:08 PM

Quant Time: May 02 05:33:45 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM050121.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat May 01 16:23:05 2021
Response via : Initial Calibration



Quantitation Report (Qedit)

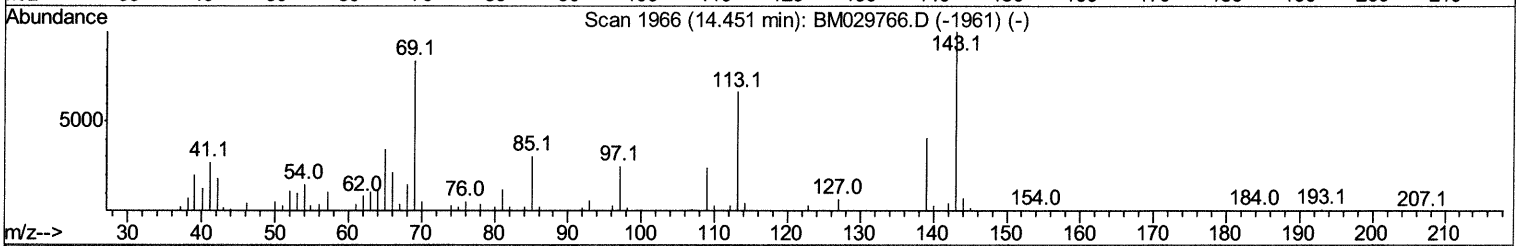
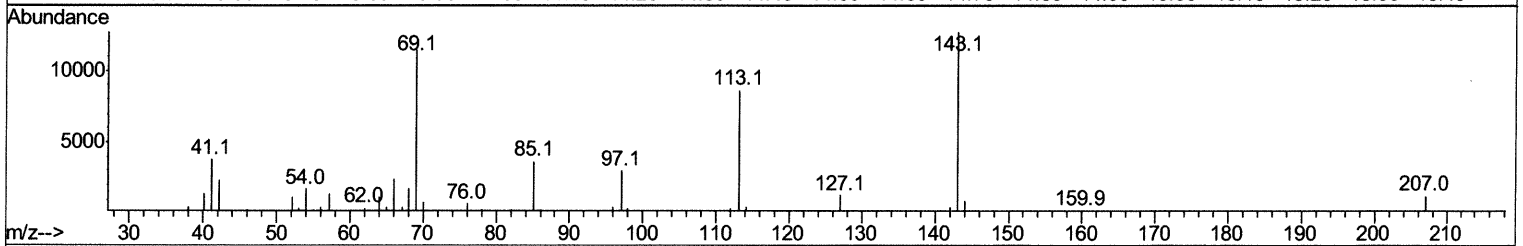
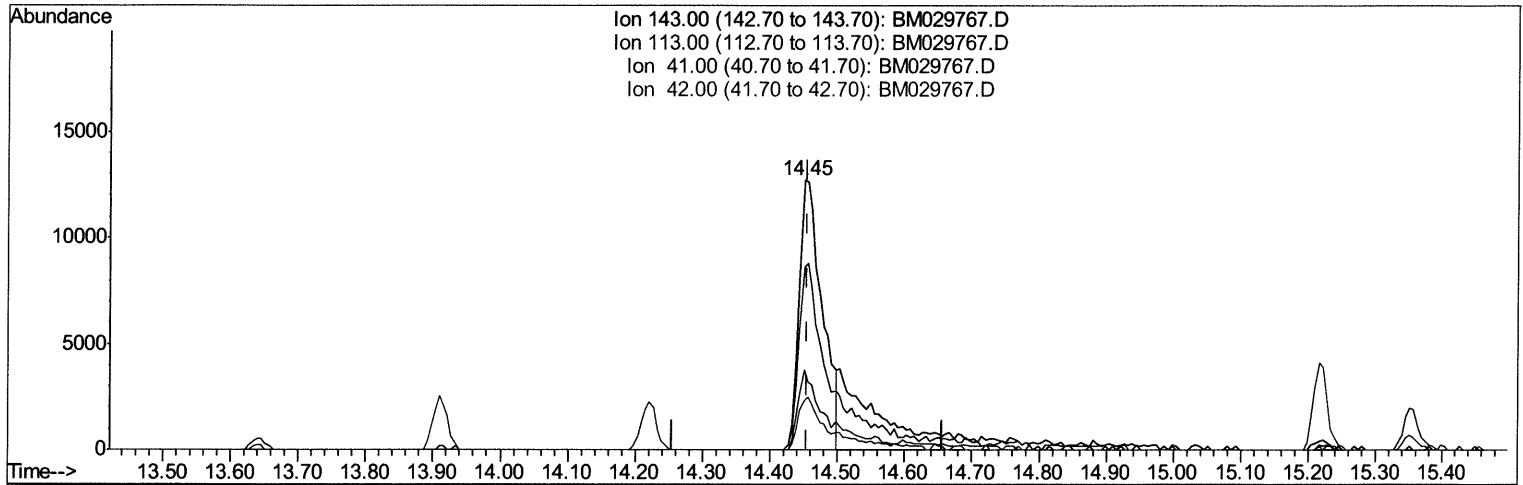
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TIC: BM029767.D

(54) 4-Nitrophenol-d4 (S)

14.451min (-0.006) 16.89ng/ul

response 30043

Ion	Exp%	Act%
143.00	100	100
113.00	71.30	67.67
41.00	32.40	29.74
42.00	22.10	18.08

Quantitation Report (Qedit)

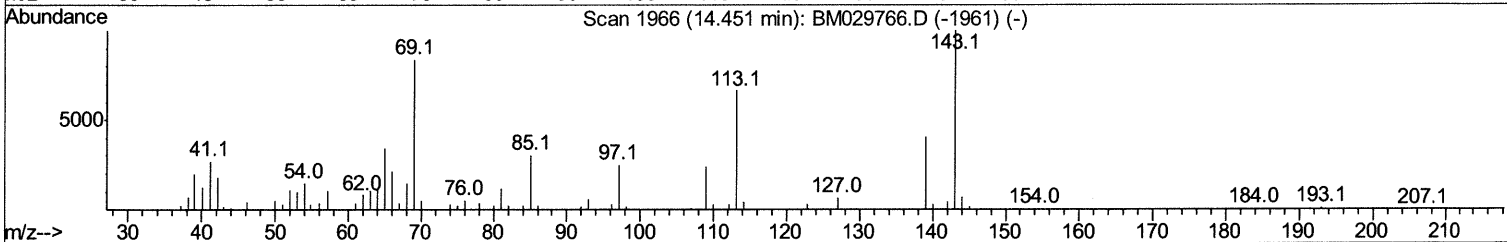
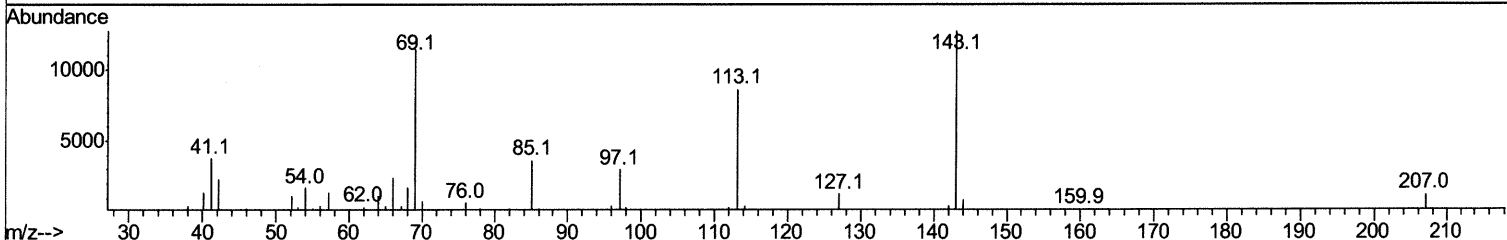
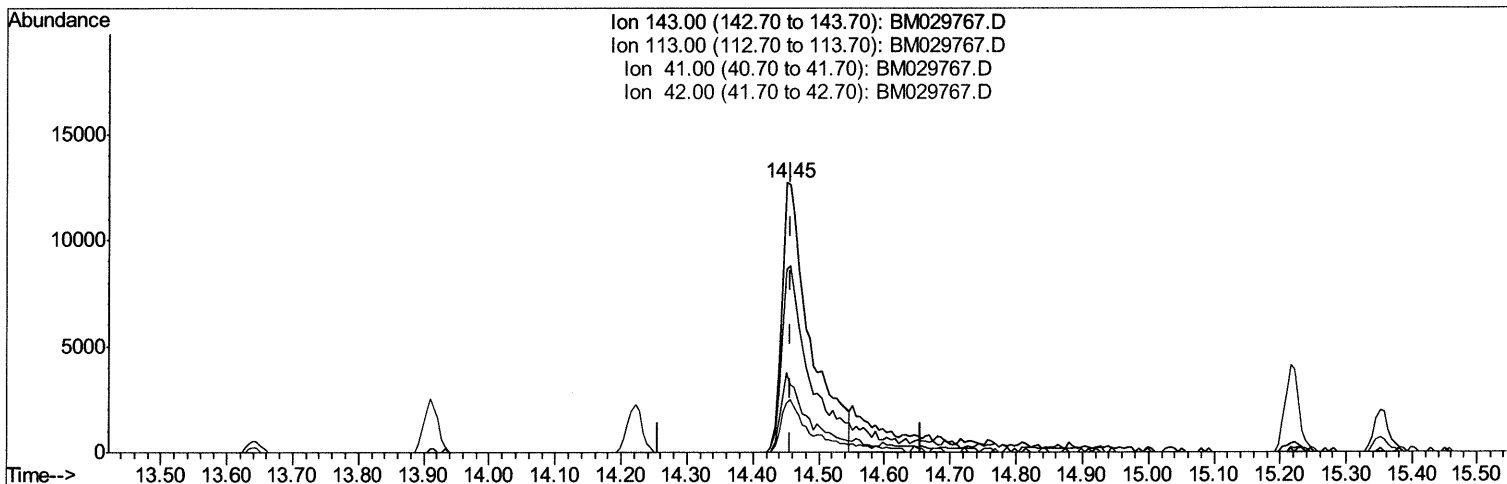
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(54) 4-Nitrophenol-d4 (S)

14.451min (-0.006) 21.08ng/ul m 05/04/21 JU

response 37500

Ion	Exp%	Act%
143.00	100	100
113.00	71.30	67.67
41.00	32.40	29.74
42.00	22.10	18.08

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Manual Integrations
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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.57	152	59932	20.00	ng/ul	0.00
20) Naphthalene-d8	10.35	136	218786	20.00	ng/ul	0.00
38) Acenaphthene-d10	14.22	164	169734	20.00	ng/ul	0.00
64) Phenanthrene-d10	16.97	188	391298	20.00	ng/ul	0.00
79) Chrysene-d12	21.15	240	452789	20.00	ng/ul	0.00
88) Perylene-d12	23.31	264	486920	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.08	96	6804	5.55	ng/uL	0.00
4) Pyridine-d5	3.49	84	72713	27.39	ng/ul	0.00
7) Phenol-d5	6.76	99	111575	27.65	ng/ul	0.00
9) Bis-(2-Chloroethyl)ether-d	6.92	67	59943	28.36	ng/ul	0.00
11) 2-Chlorophenol-d4	7.11	132	107102	29.34	ng/ul	0.00
15) 4-Methylphenol-d8	8.29	113	94721	27.40	ng/ul	0.00
21) Nitrobenzene-d5	8.73	128	47669	29.92	ng/ul	0.00
24) 2-Nitrophenol-d4	9.45	143	60899	30.19	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.99	165	116456	27.63	ng/ul	0.00
31) 4-Chloroaniline-d4	10.50	131	130708	27.50	ng/ul	0.00
46) Dimethylphthalate-d6	13.64	166	433183	32.22	ng/ul	0.00
49) Acenaphthylene-d8	13.92	160	510073	30.77	ng/ul	0.00
54) 4-Nitrophenol-d4	14.45	143	37500m>	21.08	ng/ul>	0.00
60) Fluorene-d10	15.22	176	379781	32.65	ng/ul	0.00
65) 4,6-Dinitro-2-methylphenol	15.35	200	68108	23.94	ng/ul	0.00
73) Anthracene-d10	17.07	188	629468	31.91	ng/ul	0.00
81) Pyrene-d10	19.36	212	781496	33.80	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.18	264	892062	31.93	ng/ul	0.00

05/04/21 JU

Target Compounds Ovalue

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