

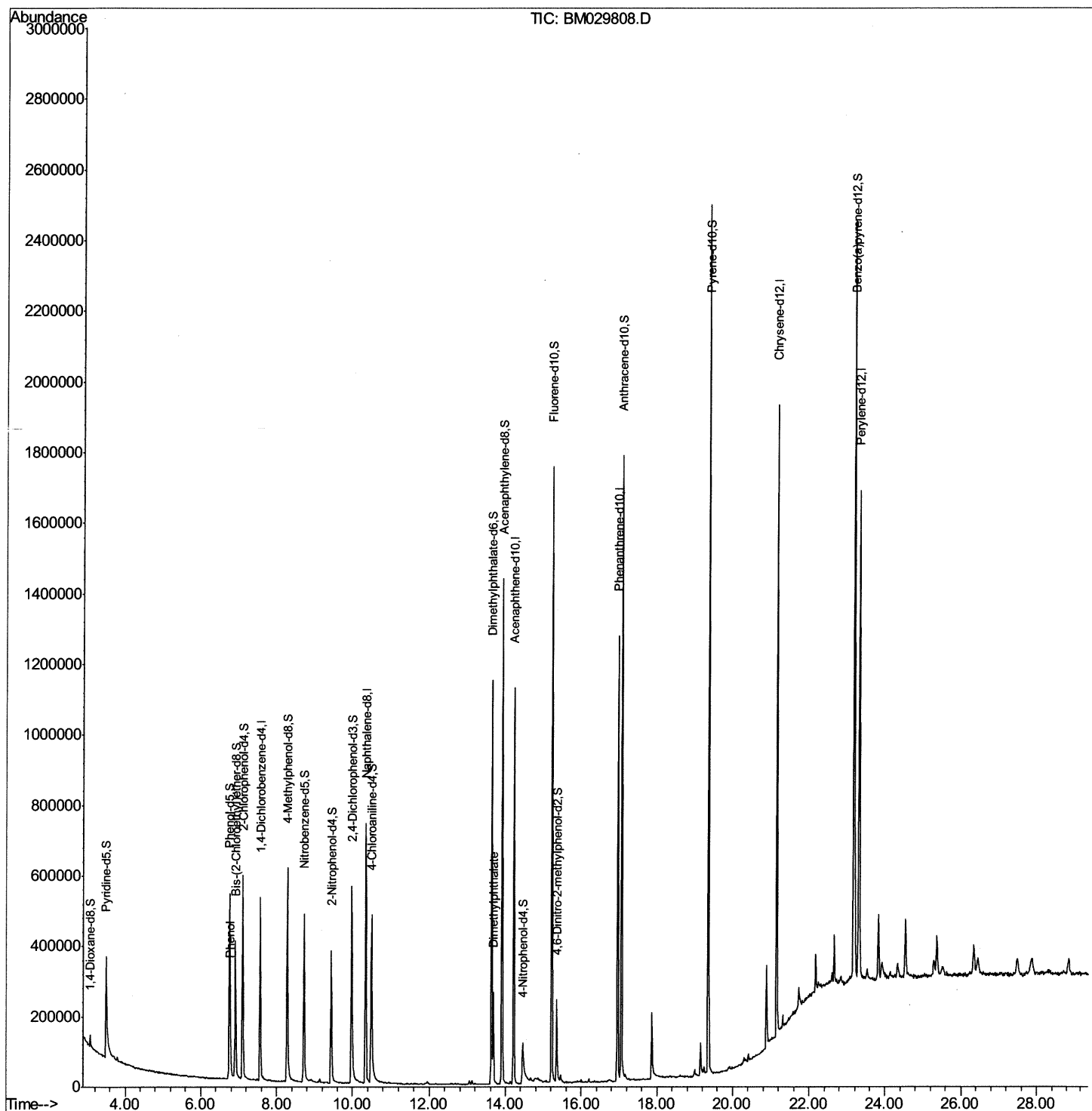
Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM050421\
Data File : BM029808.D
Acq On : 05 May 2021 03:04
Operator : CG/JU
Sample : M2121-01
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BG510

Manual Integrations
APPROVED

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

Quant Time: May 05 04:28:45 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM050421.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed May 05 03:36:11 2021
Response via : Initial Calibration



Quantitation Report (Qedit)

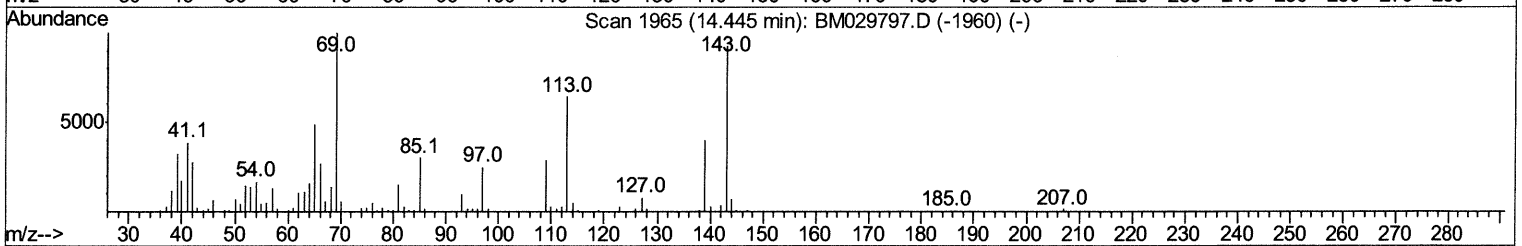
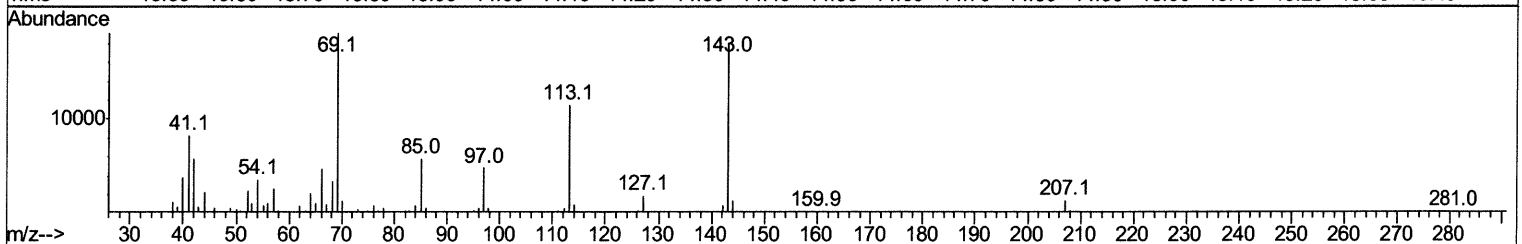
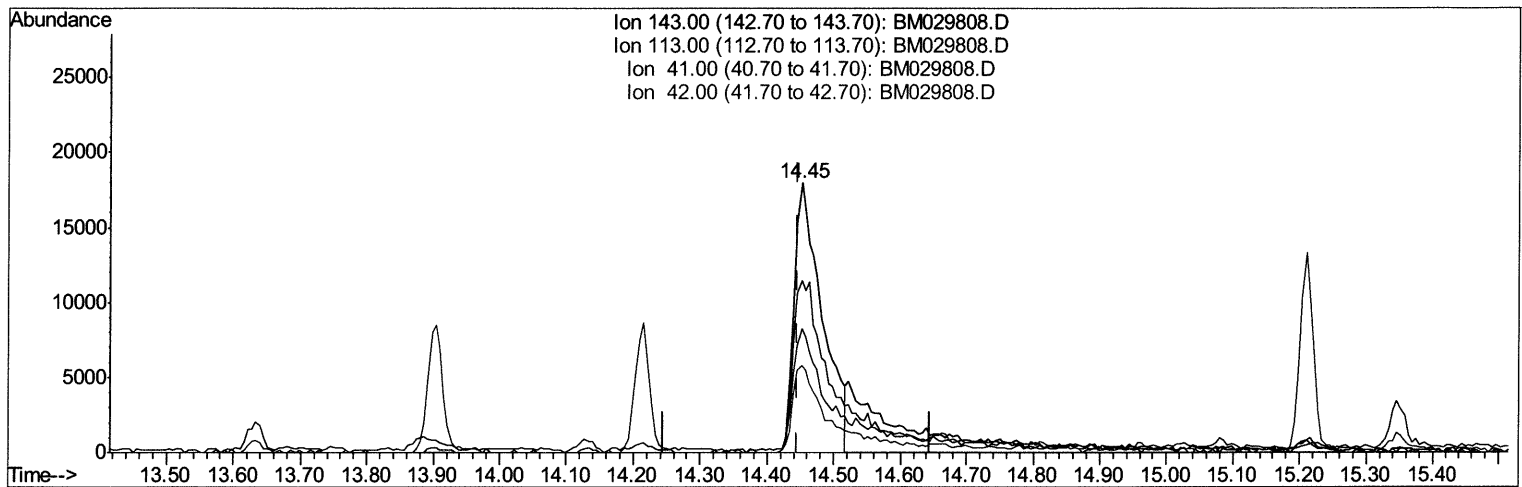
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TIC: BM029808.D

(54) 4-Nitrophenol-d4 (S)

14.451min (+0.006) 12.48ng/ul

response 52374

Ion	Exp%	Act%
143.00	100	100
113.00	67.10	63.81
41.00	45.40	46.05
42.00	30.10	32.08

Quantitation Report (Qedit)

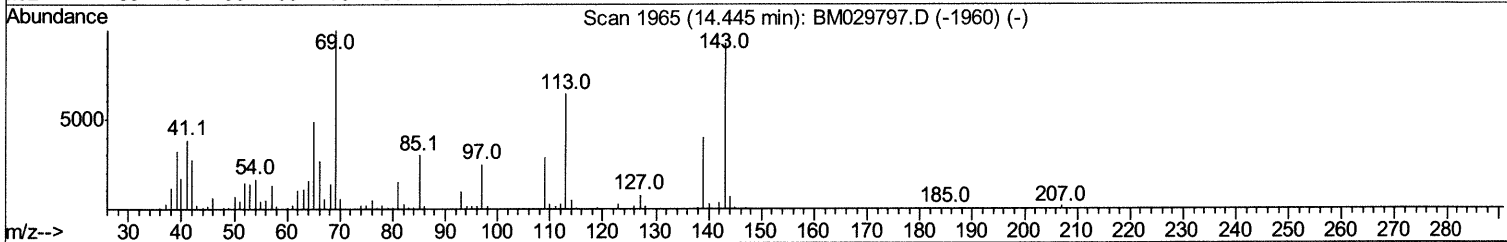
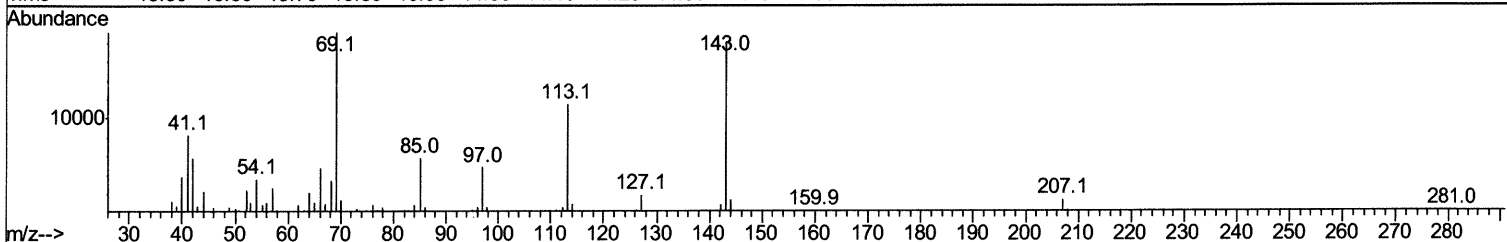
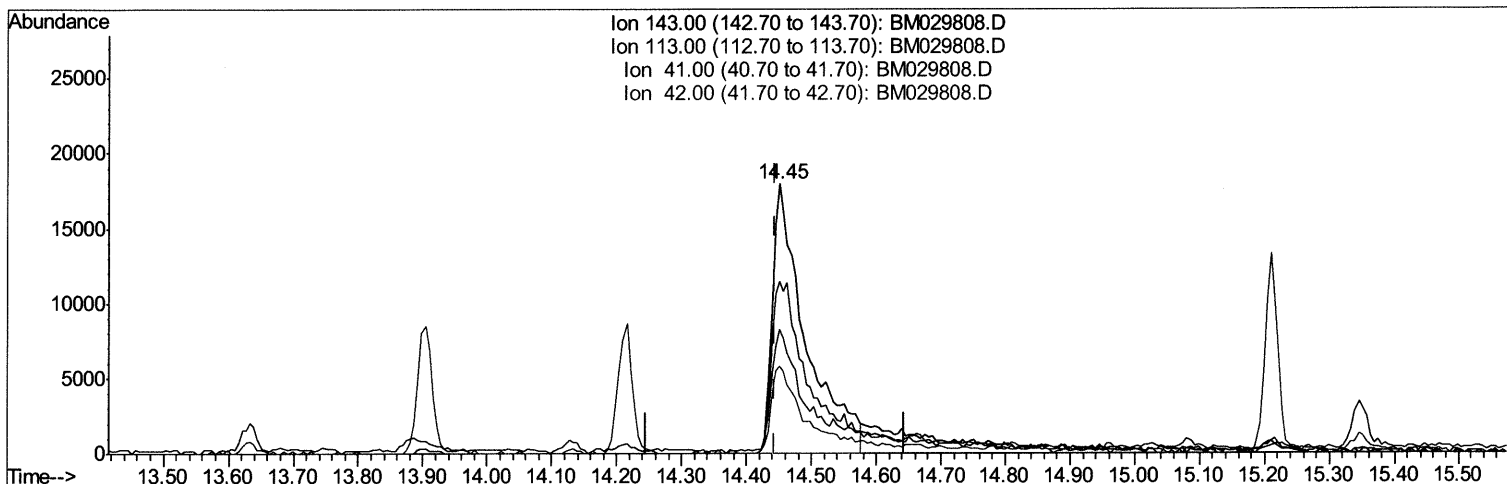
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TIC: BM029808.D

(54) 4-Nitrophenol-d4 (S)

14.451min (+0.006) 15.15ng/ul m 05/06/21 JU

response 63611

Ion	Exp%	Act%
143.00	100	100
113.00	67.10	63.81
41.00	45.40	46.05
42.00	30.10	32.08

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM050421\
 Data File : BM029808.D
 Acq On : 05 May 2021 03:04
 Operator : CG/JU
 Sample : M2121-01
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BG510

Manual Integrations
 APPROVED

mohammad
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 Quant Title : SVOA CALIBRATION
 QLast Update : Wed May 05 03:36:11 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.57	152	148465	20.00	ng/ul	0.00
20) Naphthalene-d8	10.34	136	612823	20.00	ng/ul	0.00
38) Acenaphthene-d10	14.22	164	365405	20.00	ng/ul	0.00
64) Phenanthrene-d10	16.96	188	765762	20.00	ng/ul	0.00
79) Chrysene-d12	21.14	240	855575	20.00	ng/ul	0.00
88) Perylene-d12	23.31	264	935338	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.08	96	19174	4.78	ng/uL	0.00
4) Pyridine-d5	3.49	84	210444	22.89	ng/ul	0.00
7) Phenol-d5	6.75	99	317815	24.65	ng/ul	0.00
9) Bis-(2-Chloroethyl) ether-d	6.91	67	202612	25.44	ng/ul	0.00
11) 2-Chlorophenol-d4	7.10	132	273703	26.65	ng/ul	0.00
15) 4-Methylphenol-d8	8.28	113	254763	24.76	ng/ul	0.00
21) Nitrobenzene-d5	8.72	128	122547	26.91	ng/ul	0.00
24) 2-Nitrophenol-d4	9.44	143	129975	26.14	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.97	165	232482	25.01	ng/ul	0.00
31) 4-Chloroaniline-d4	10.49	131	339112	25.03	ng/ul	0.00
46) Dimethylphthalate-d6	13.63	166	764384	28.01	ng/ul	0.00
49) Acenaphthylene-d8	13.90	160	989082	27.71	ng/ul	0.00
54) 4-Nitrophenol-d4	14.45	143	63611m >	15.15	ng/ul >	0.0005/06/21JU
60) Fluorene-d10	15.21	176	662437	28.53	ng/ul	0.00
65) 4,6-Dinitro-2-methylphenol	15.35	200	53675	10.44	ng/ul	0.00
73) Anthracene-d10	17.06	188	1033039	27.28	ng/ul	0.00
81) Pyrene-d10	19.36	212	1195607	27.24	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.17	264	1460650	27.81	ng/ul	0.00

Target Compounds

					Ovalue
8) Phenol	6.78	94	80984	6.174	ng/ul 98
47) Dimethylphthalate	13.68	163	162005	5.949	ng/ul 100

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