

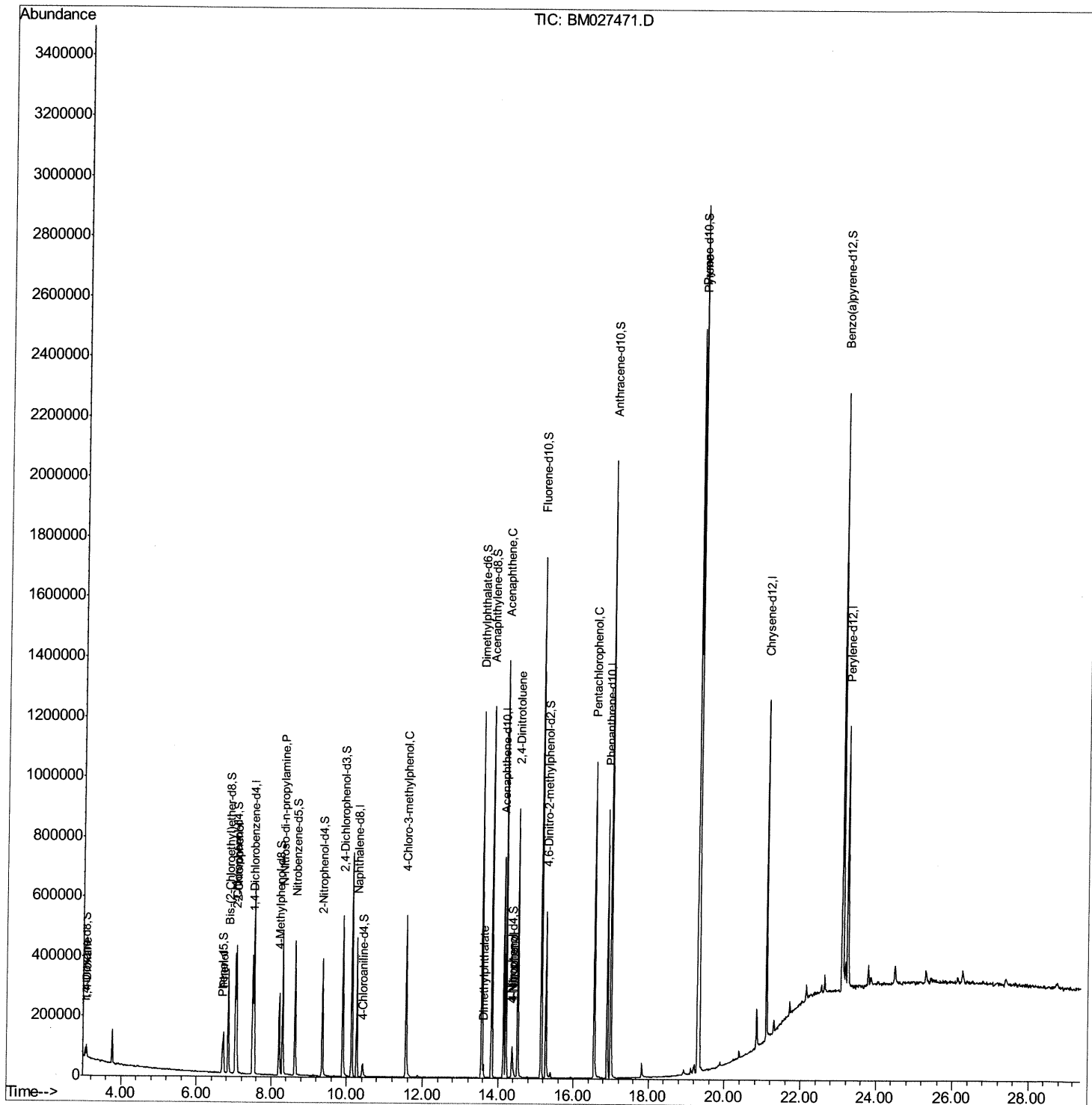
Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM091920\
Data File : BM027471.D
Acq On : 18 Sep 2020 20:50
Operator : JU/CG
Sample : L4001-05MS
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
ETYP2MS

Manual Integrations
APPROVED

mohammad
9/21/2020 10:47:56 AM

Quant Time: Sep 19 01:01:20 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM090420MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat Sep 19 00:24:00 2020
Response via : Initial Calibration



Quantitation Report (Qedit)

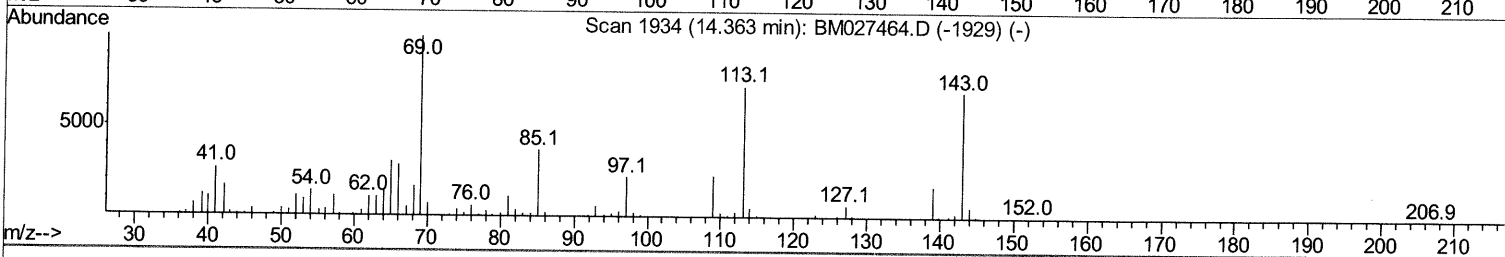
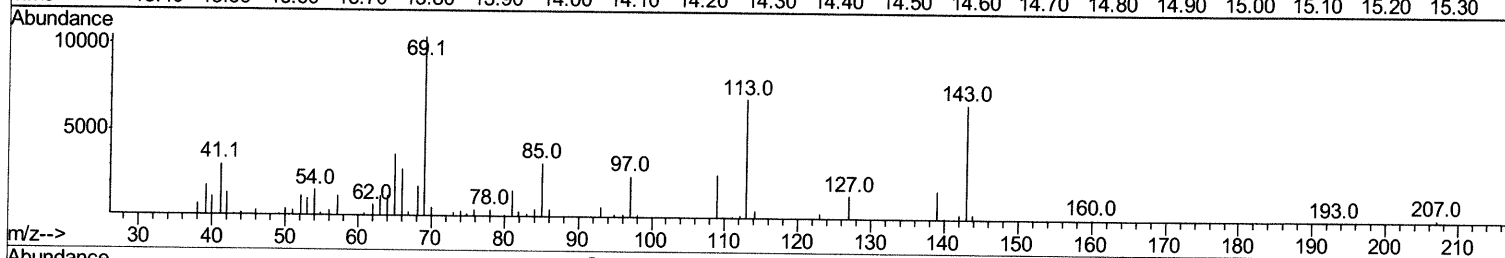
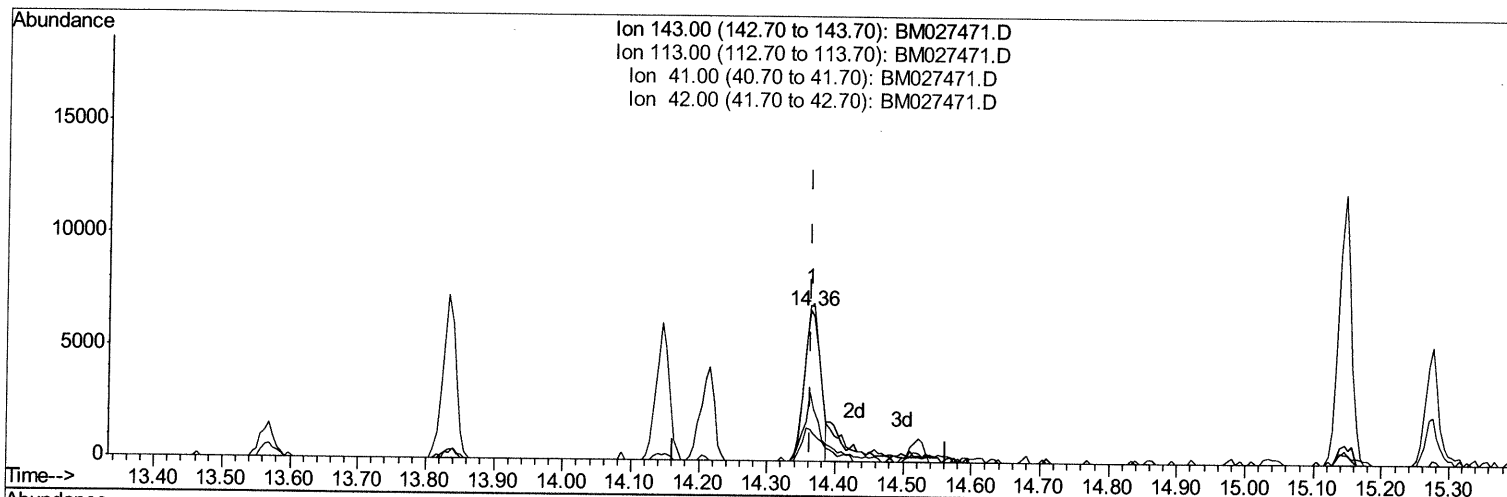
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TIC: BM027471.D

(52) 4-Nitrophenol-d4 (S)

14.363min (-0.000) 3.55ng/ul

response 11274

Ion	Exp%	Act%
143.00	100	100
113.00	105.60	103.27
41.00	38.00	45.34
42.00	24.80	20.96

Quantitation Report (Qedit)

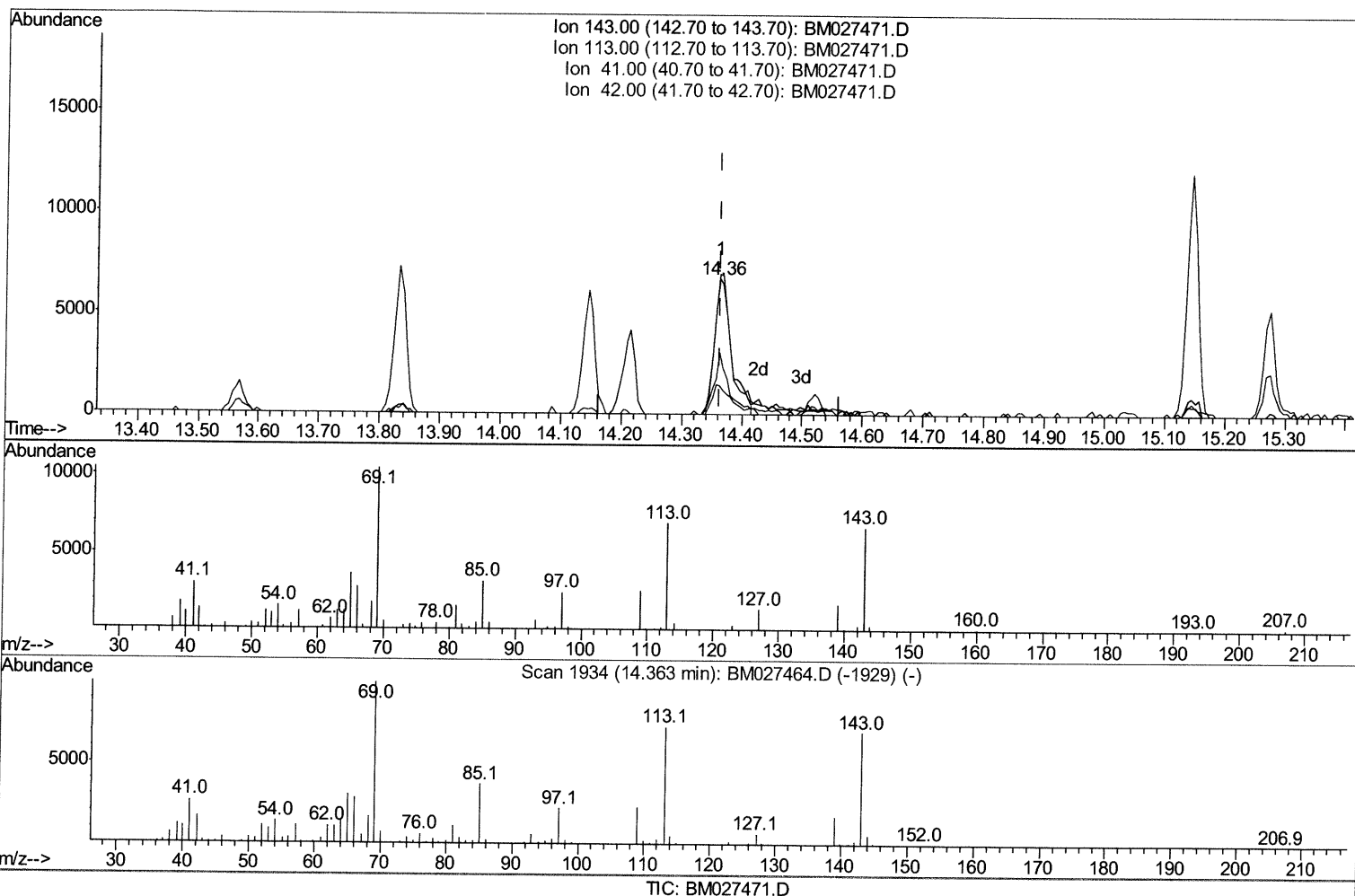
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(52) 4-Nitrophenol-d4 (S)

14.363min (-0.000) 4.21ng/ul m 09/21/20 JU

response 13356

Ion	Exp%	Act%
143.00	100	100
113.00	105.60	103.27
41.00	38.00	45.34
42.00	24.80	20.96

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Manual Integrations
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 QLast Update : Sat Sep 19 00:24:00 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.50	152	95501	20.00	ng/ul	0.00
18) Naphthalene-d8	10.26	136	335889	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.15	164	226393	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.90	188	512907	20.00	ng/ul	0.00
78) Chrysene-d12	21.10	240	541420	20.00	ng/ul	0.00
86) Perylene-d12	23.24	264	622665	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.06	96	14367	6.69	ng/uL	0.00
5) Phenol-d5	6.69	99	49310	6.26	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.85	67	153429	29.35	ng/ul	0.00
9) 2-Chlorophenol-d4	7.04	132	146539	25.42	ng/ul	0.00
13) 4-Methylphenol-d8	8.22	113	88859	13.84	ng/ul	0.00
19) Nitrobenzene-d5	8.64	128	87369	34.55	ng/ul	0.00
22) 2-Nitrophenol-d4	9.36	143	97277	35.08	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.90	165	179734	30.11	ng/ul	0.00
29) 4-Chloroaniline-d4	10.41	131	25457	3.61	ng/ul	0.00
44) Dimethylphthalate-d6	13.57	166	710452	39.44	ng/ul	0.00
47) Acenaphthylene-d8	13.83	160	790930	38.87	ng/ul	0.00
52) 4-Nitrophenol-d4	14.36	143	13356m>	4.21	ng/ul>	0.00
58) Fluorene-d10	15.14	176	610246	37.82	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.27	200	110304	33.60	ng/ul	0.00
71) Anthracene-d10	16.99	188	1052999	43.41	ng/ul	0.00
79) Pyrene-d10	19.30	212	1229885	50.99	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.10	264	1389171	41.59	ng/ul	0.00

Target Compounds

					Ovalue	
2) 1,4-Dioxane	3.09	88	18981	7.389	ng/uL#	92
6) Phenol	6.72	94	67905	8.117	ng/ul	98
10) 2-Chlorophenol	7.07	128	154938	25.772	ng/ul	98
15) N-Nitroso-di-n-propylamine	8.30	70	201272	32.645	ng/ul	99
33) 4-Chloro-3-methylphenol	11.56	107	176679	27.916	ng/ul	97
45) Dimethylphthalate	13.62	163	24295	1.280	ng/ul	97
50) Acenaphthene	14.21	153	525885	35.646	ng/ul	99
53) 4-Nitrophenol	14.38	109	19653	5.944	ng/ul	87
55) 2,4-Dinitrotoluene	14.52	165	209663	37.575	ng/ul	97
69) Pentachlorophenol	16.56	266	167265	37.918	ng/ul	98
80) Pyrene	19.33	202	1567431	47.856	ng/ul	98

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