

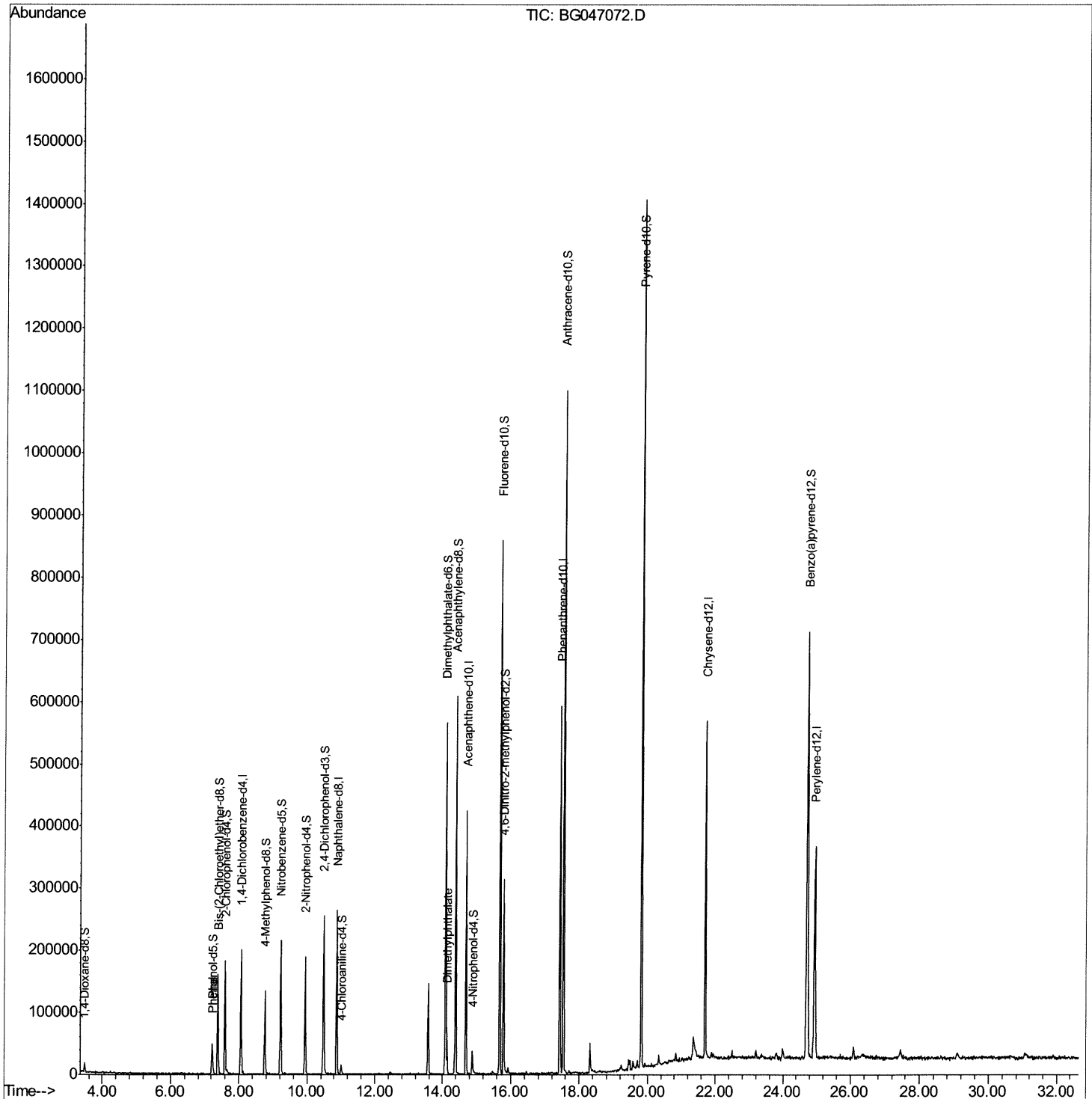
Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG102620\
Data File : BG047072.D
Acq On : 26 Oct 2020 10:57
Operator : CG/JU
Sample : L4499-13
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EW2C6

Manual Integrations
APPROVED

mohammad
10/26/2020 2:05:47 PM

Quant Time: Oct 26 11:36:02 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG102220MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Oct 26 10:36:50 2020
Response via : Initial Calibration



Quantitation Report (Qedit)

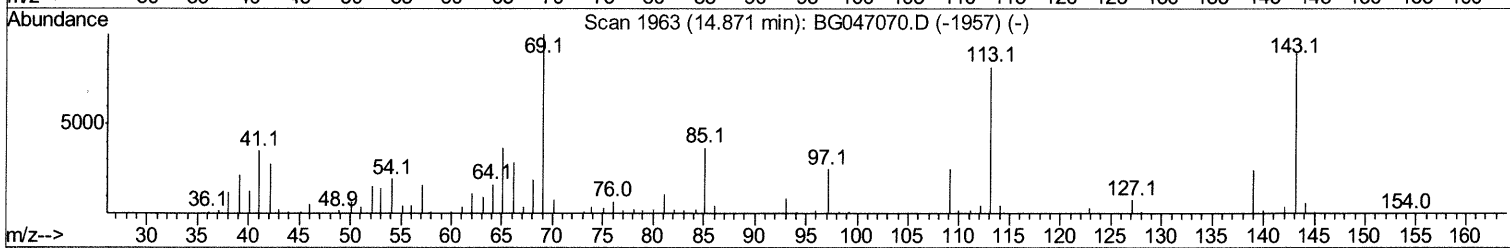
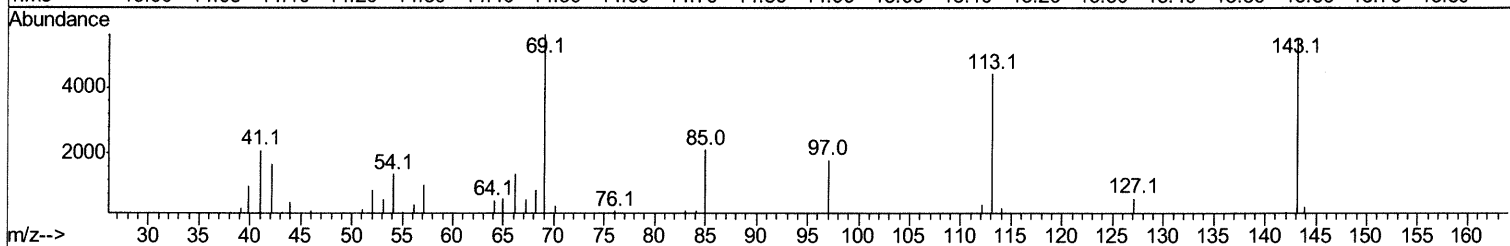
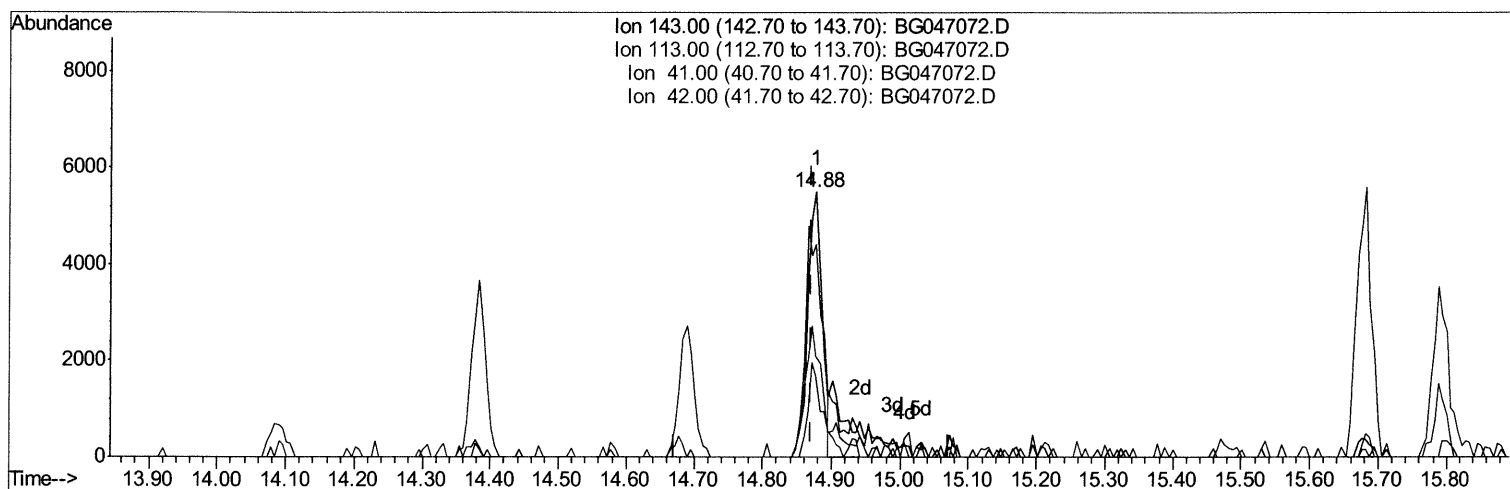
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TIC: BG047072.D

(52) 4-Nitrophenol-d4 (S)

14.877min (+0.006) 4.17ng/ul

response 8113

Ion	Exp%	Act%
143.00	100	100
113.00	89.40	80.27
41.00	40.00	38.13
42.00	29.30	30.63

Quantitation Report (Qedit)

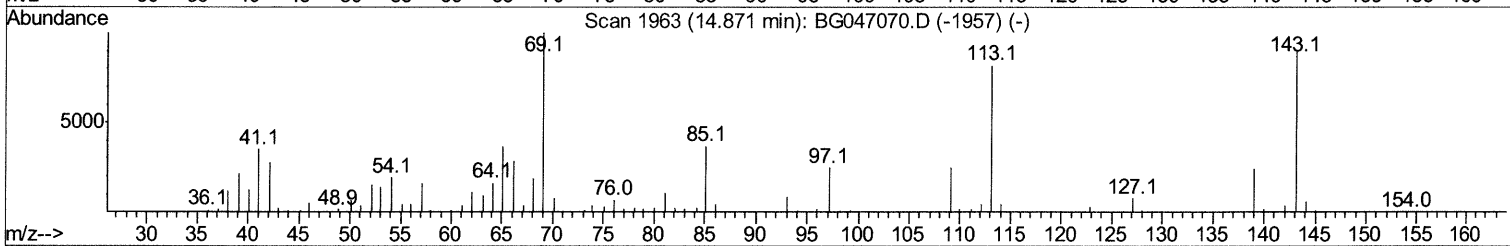
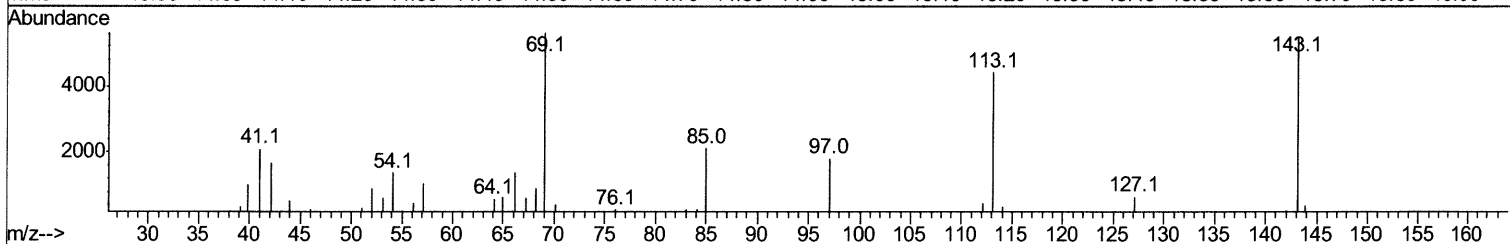
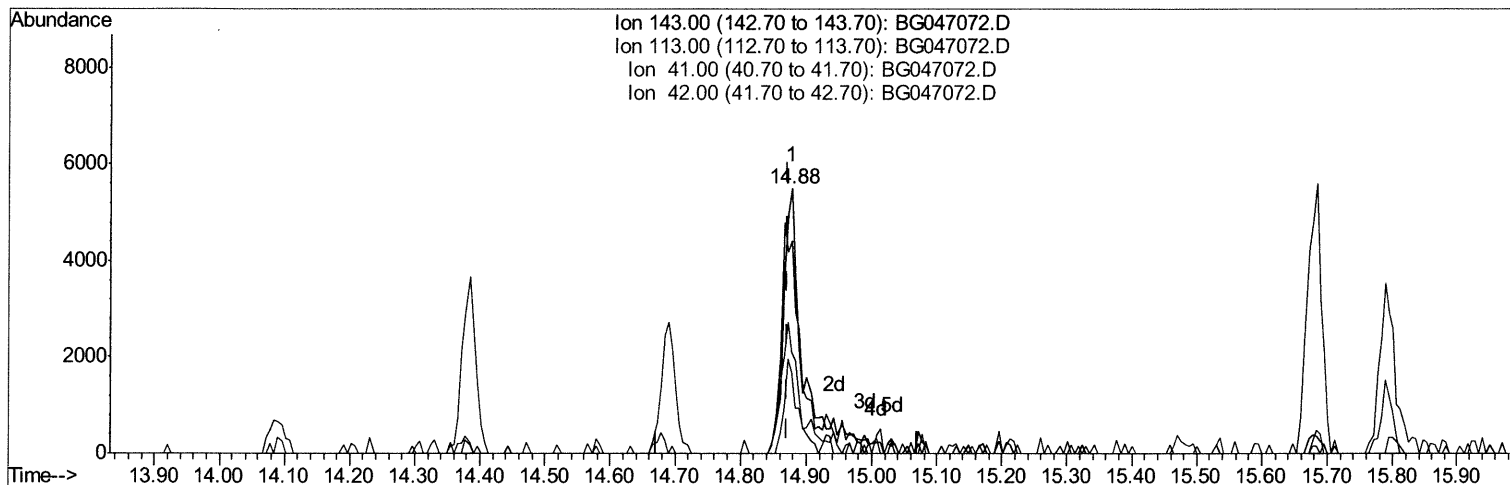
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TIC: BG047072.D

(52) 4-Nitrophenol-d4 (S)

14.877min (+0.006) 5.89ng/ul m 10/26/20 JU

response 11455

Ion	Exp%	Act%
143.00	100	100
113.00	89.40	80.27
41.00	40.00	38.13
42.00	29.30	30.63

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG102620\
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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	8.06	152	51257	20.00	ng/ul	0.00
18) Naphthalene-d8	10.87	136	210851	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.69	164	150166	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.43	188	363870	20.00	ng/ul	0.00
78) Chrysene-d12	21.70	240	364440	20.00	ng/ul	0.00
86) Perylene-d12	24.94	264	377675	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.47	96	7062	5.56	ng/uL	0.00
5) Phenol-d5	7.22	99	28859	6.49	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.38	67	73018	28.54	ng/ul	0.00
9) 2-Chlorophenol-d4	7.59	132	79344	22.94	ng/ul	0.00
13) 4-Methylphenol-d8	8.76	113	51936	14.63	ng/ul	0.00
19) Nitrobenzene-d5	9.22	128	48175	27.26	ng/ul	0.00
22) 2-Nitrophenol-d4	9.95	143	57452	27.83	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.48	165	101442	25.18	ng/ul	0.00
29) 4-Chloroaniline-d4	11.01	131	8421	2.45	ng/ul	0.00
44) Dimethylphthalate-d6	14.09	166	380520	30.82	ng/ul	0.00
47) Acenaphthylene-d8	14.38	160	455652	31.19	ng/ul	0.00
52) 4-Nitrophenol-d4	14.88	143	11455m >	5.89	ng/ul >	0.00
58) Fluorene-d10	15.68	176	365938	31.92	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.79	200	80075	31.91	ng/ul	0.00
71) Anthracene-d10	17.53	188	657524	37.28	ng/ul	0.00
79) Pyrene-d10	19.81	212	791668	38.08	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.71	264	771943	37.18	ng/ul	0.00

10/26/2020

Target Compounds

Target Compounds	R.T.	QIon	Response	Conc	Units	Ovalue
6) Phenol	7.23	94	5031	1.159	ng/ul#	80
45) Dimethylphthalate	14.13	163	47592	3.841	ng/ul#	94

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