

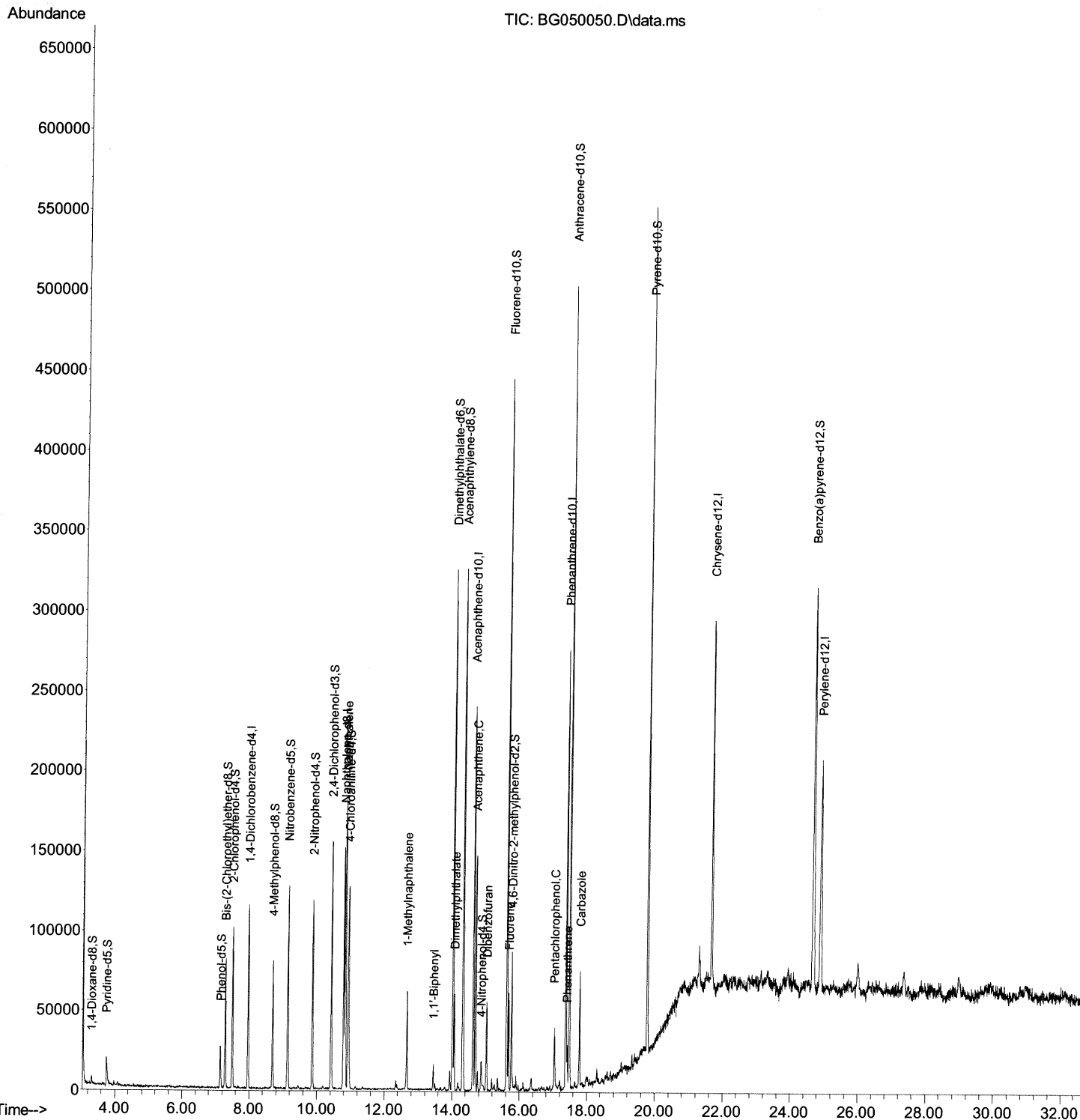
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG083021\
Data File : BG050050.D
Acq On : 31 Aug 2021 4:45
Operator : CG/JU
Sample : M3448-21
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBKG6

Manual Integrations
APPROVED

mohammad
8/31/2021 11:36:09 AM

Quant Time: Aug 31 05:58:18 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG081921.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Aug 30 14:59:55 2021
Response via : Initial Calibration



Quantitation Report (Qedit)

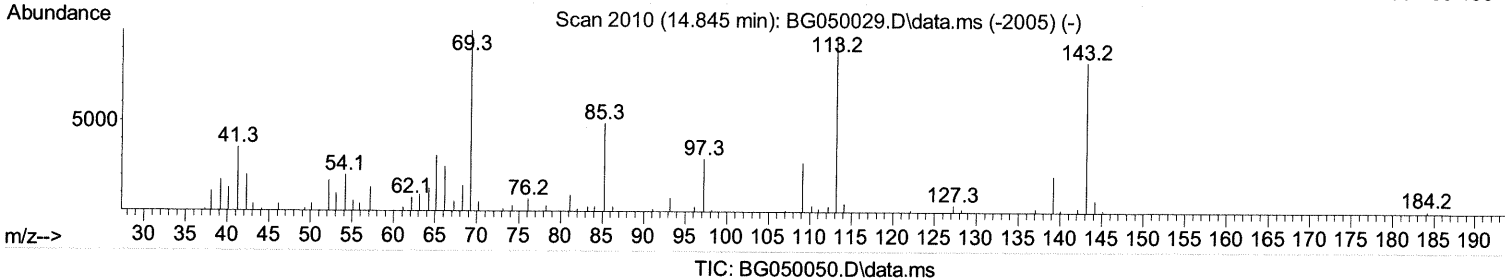
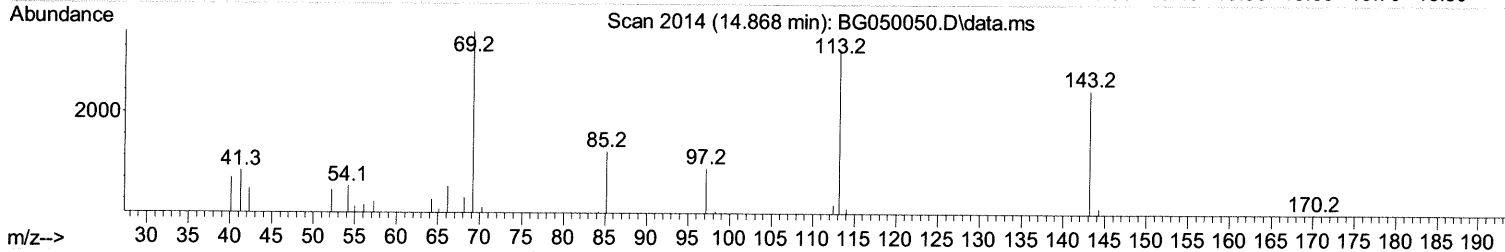
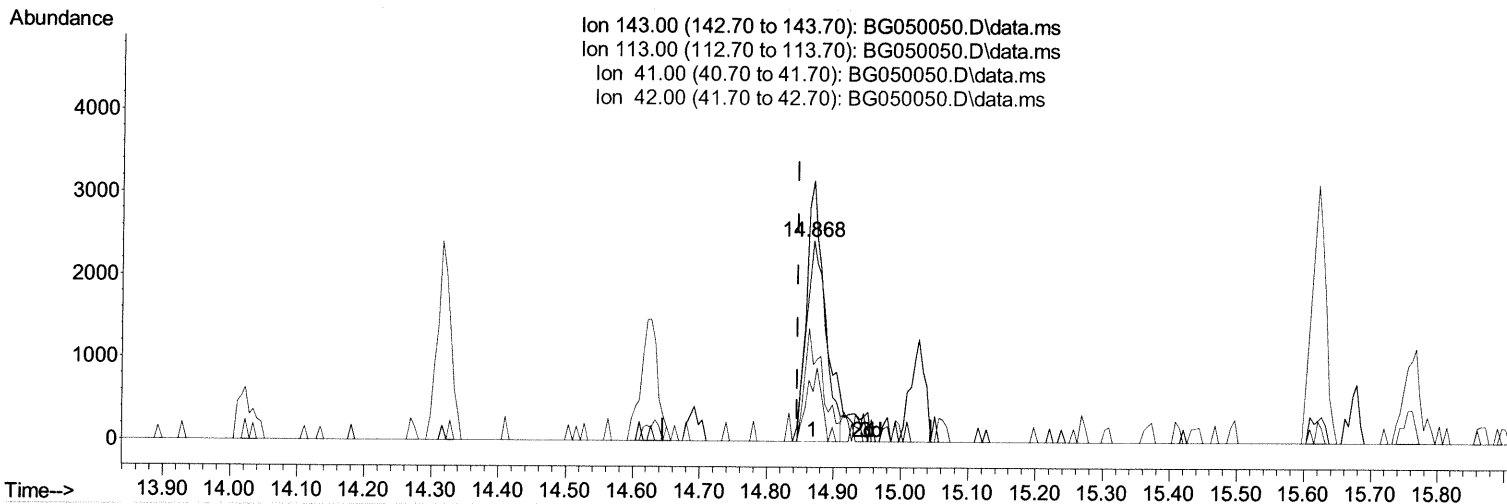
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(54) 4-Nitrophenol-d4 (S)

14.868min (+ 0.023) 5.32 ng/ul

response 5084

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	121.00	130.05
41.00	45.30	38.25
42.00	25.60	24.57

Quantitation Report (Qedit)

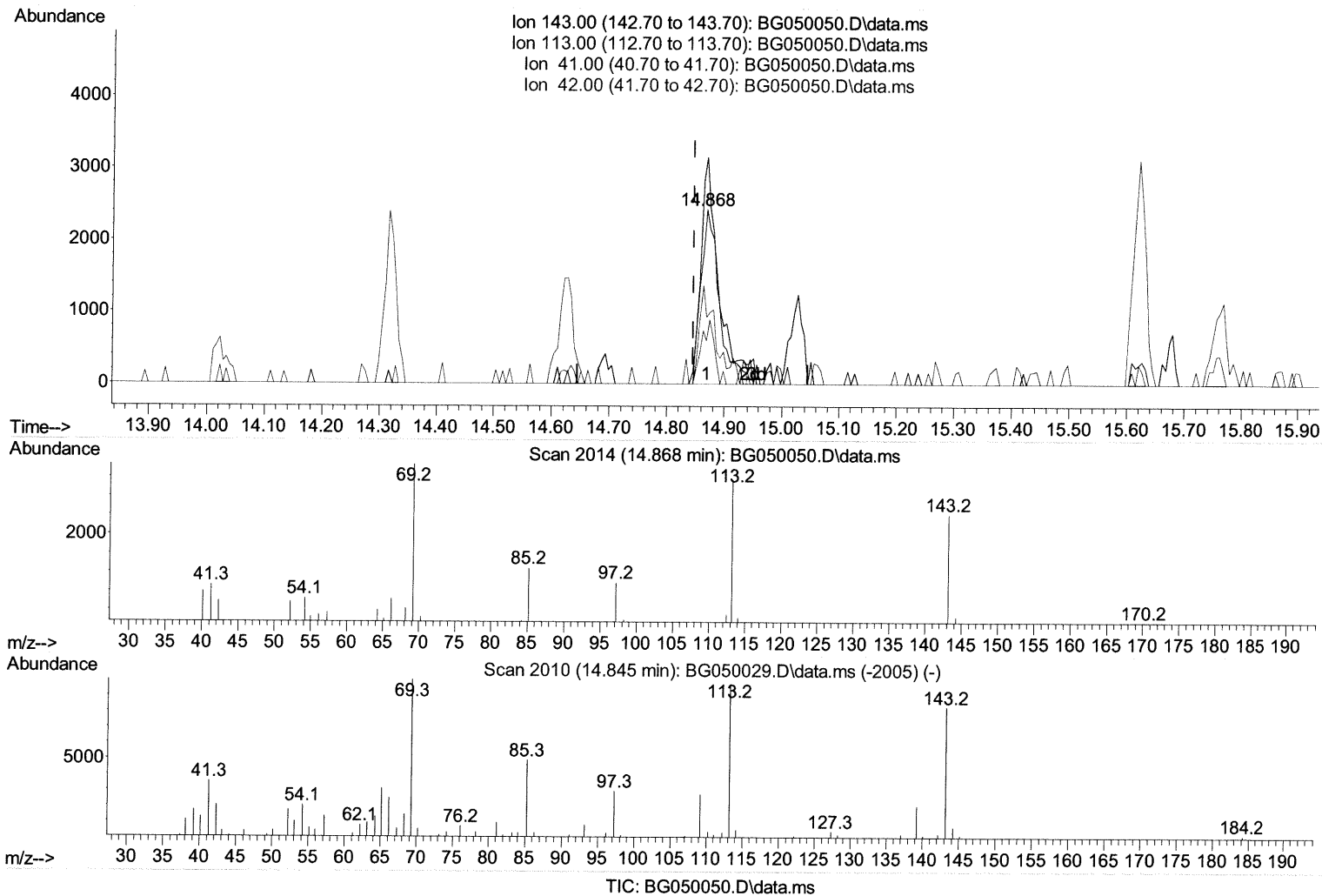
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Data File : BG050050.D
Acq On : 31 Aug 2021 4:45
Operator : CG/JU
Sample : M3448-21
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBKG6

Manual Integrations
APPROVED

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Quant Title : SVOA CALIBRATION
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Response via : Initial Calibration



(54) 4-Nitrophenol-d4 (S)

14.868min (+ 0.023) 5.75 ng/ul m 09/01/21JU

response 5497

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	121.00	130.05
41.00	45.30	38.25
42.00	25.60	24.57

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Manual Integrations
 APPROVED

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Compound	R.T.	QIon	Response	Conc Units	Dev(Min)
Internal Standards					
1) 1,4-Dichlorobenzene-d4	7.966	152	31737	20.000 ng/ul	0.01
20) Naphthalene-d8	10.786	136	127138	20.000 ng/ul	0.00
38) Acenaphthene-d10	14.627	164	86378	20.000 ng/ul	0.00
64) Phenanthrene-d10	17.382	188	169560	20.000 ng/ul	0.00
79) Chrysene-d12	21.653	240	141597	20.000 ng/ul	0.00
88) Perylene-d12	24.890	264	150234	20.000 ng/ul	0.02
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.313	96	2820	4.582 ng/uL	0.00
4) Pyridine-d5	3.760	84	13475	7.269 ng/ul	0.02
7) Phenol-d5	7.138	99	16296	6.047 ng/ul	0.01
9) Bis-(2-Chloroethyl)eth...	7.278	67	39429	26.335 ng/ul	0.00
11) 2-Chlorophenol-d4	7.496	132	44748	22.191 ng/ul	0.00
15) 4-Methylphenol-d8	8.688	113	30677	14.454 ng/ul	0.00
21) Nitrobenzene-d5	9.135	128	31251	31.705 ng/ul	0.00
24) 2-Nitrophenol-d4	9.863	143	32486	27.679 ng/ul	0.01
28) 2,4-Dichlorophenol-d3	10.415	165	57587	27.509 ng/ul	0.01
31) 4-Chloroaniline-d4	10.926	131	69262	24.308 ng/ul	0.00
46) Dimethylphthalate-d6	14.022	166	206374	31.219 ng/ul	0.00
49) Acenaphthylene-d8	14.322	160	247373	31.915 ng/ul	0.00
54) 4-Nitrophenol-d4	14.868	143	5497m	5.747 ng/ul	0.02
60) Fluorene-d10	15.620	176	179763	32.073 ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.761	200	18993	17.411 ng/ul	0.00
73) Anthracene-d10	17.482	188	283627	37.250 ng/ul	0.00
81) Pyrene-d10	19.773	212	290315	37.941 ng/ul	0.00
92) Benzo(a)pyrene-d12	24.667	264	272398	34.176 ng/ul	0.02
Target Compounds					
30) Naphthalene	10.838	128	152788	24.043 ng/ul	98
37) 1-Methylnaphthalene	12.665	142	28259	5.927 ng/ul#	86
43) 1,1'-Biphenyl	13.452	154	9352	1.550 ng/ul#	96
47) Dimethylphthalate	14.069	163	36905	5.641 ng/ul#	98
52) Acenaphthene	14.692	153	63595	12.375 ng/ul	97
56) Dibenzofuran	15.027	168	35697	5.016 ng/ul	93
61) Fluorene	15.679	166	27316	4.550 ng/ul	93
71) Pentachlorophenol	17.042	266	8749	9.539 ng/ul	87
72) Phenanthrene	17.418	178	16962	2.007 ng/ul#	93
77) Carbazole	17.788	167	46327	6.078 ng/ul	96

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