

Quantitation Report (Qedit)

Data Path : \\74.0.250.170\SVOASRV\HPCHEM1\BNA G\DATA\BG110717\
Data File : BG030838.D
Acq On : 8 Nov 2017 10:07
Operator : SJ/JU
Sample : I6222-11
Misc :
ALS Vial : 29 Sample Multiplier: 1

Quant Time: Nov 08 11:04:08 2017
Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM-EPA-BG101417.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Nov 08 04:22:00 2017
Response via : Initial Calibration

Instrument :

BNA_G

ClientSampleId :

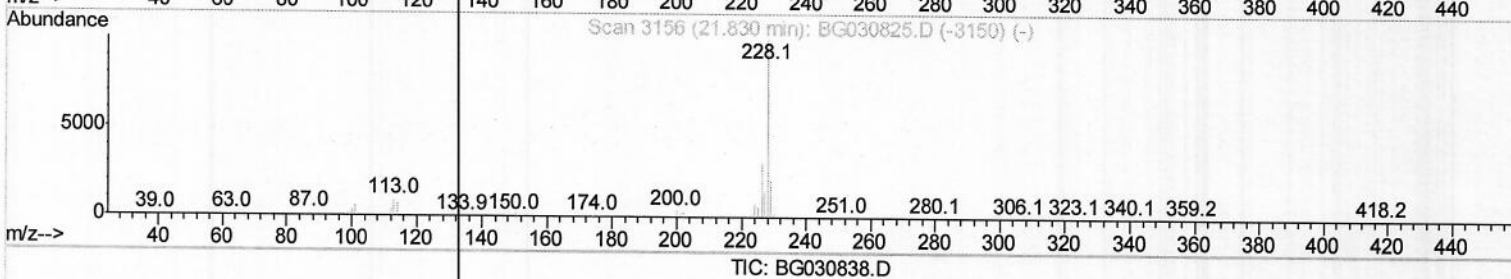
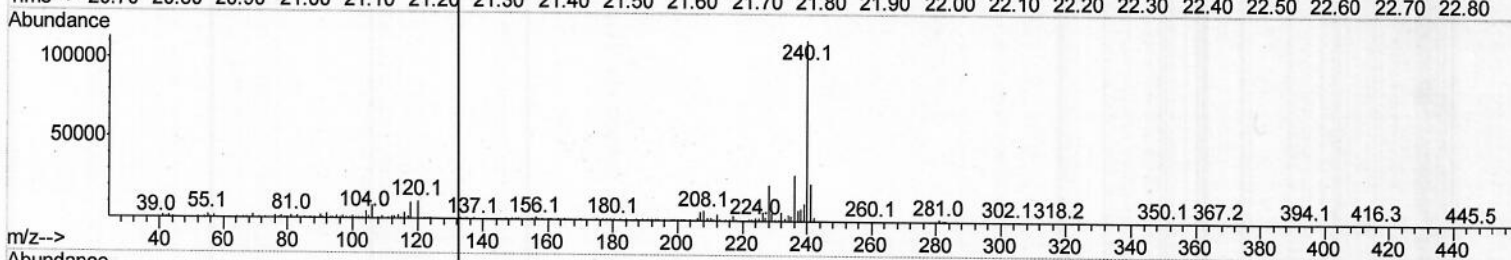
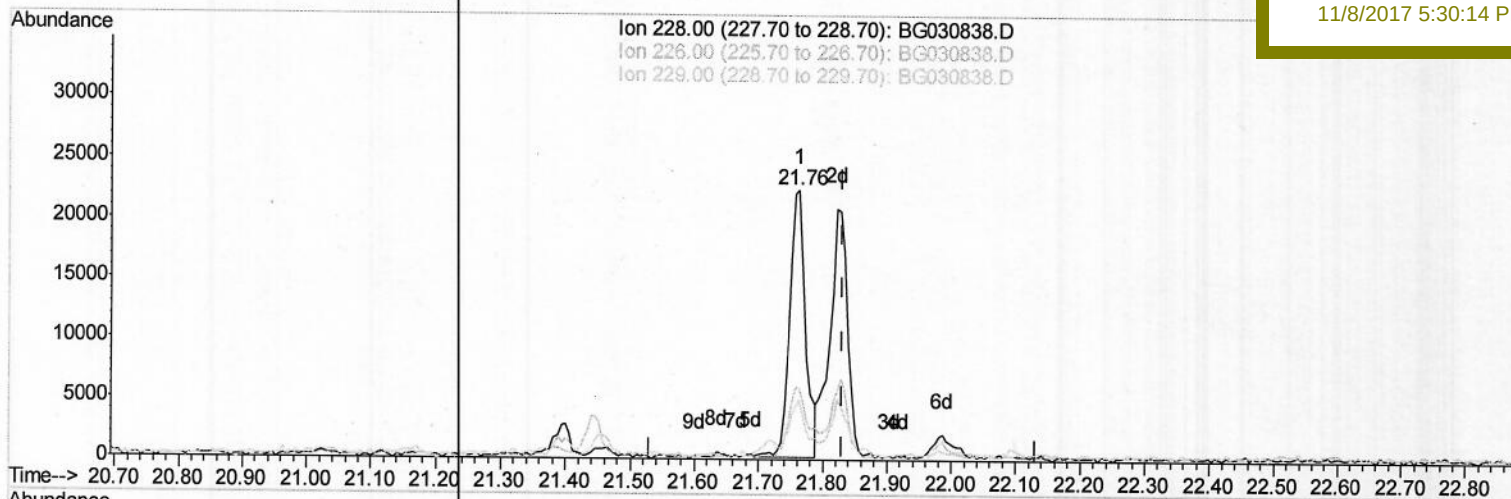
B0ET2

Manual Integrations

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Sohil

11/8/2017 5:30:14 PM



(82) Chrysene

21.763min (-0.067) 1.15ng/ul

response 38992

Ion	Exp%	Act%
228.00	100	100
226.00	30.20	26.19
229.00	19.00	22.25
0.00	0.00	0.00

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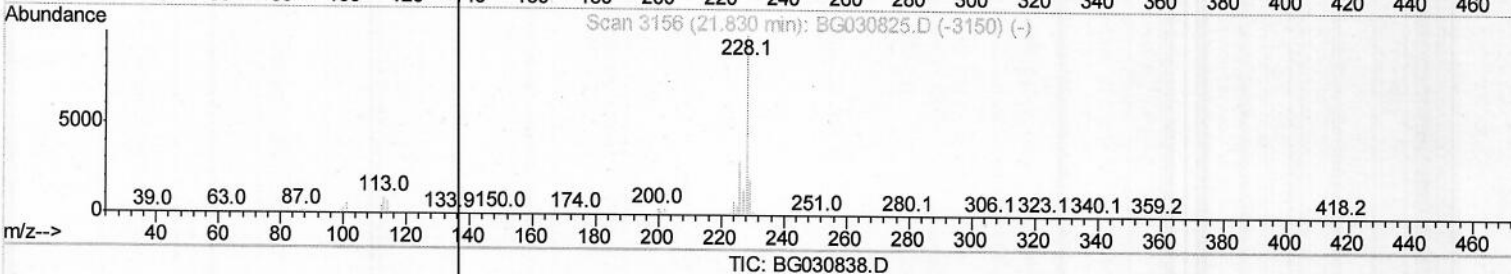
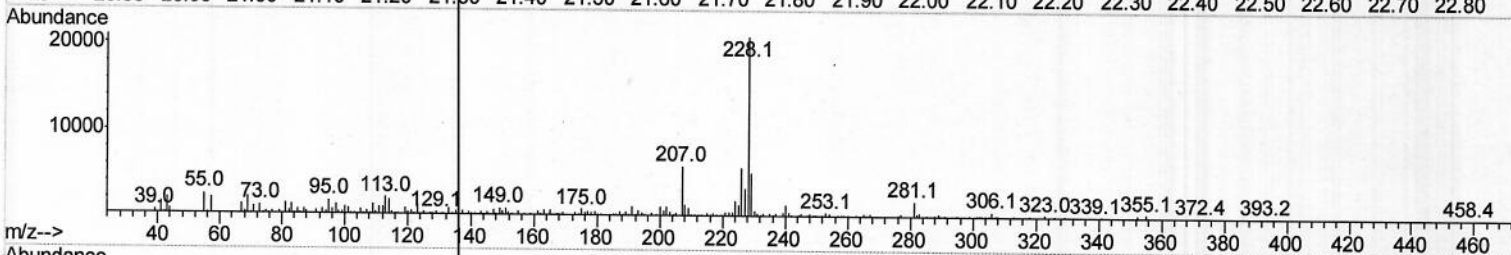
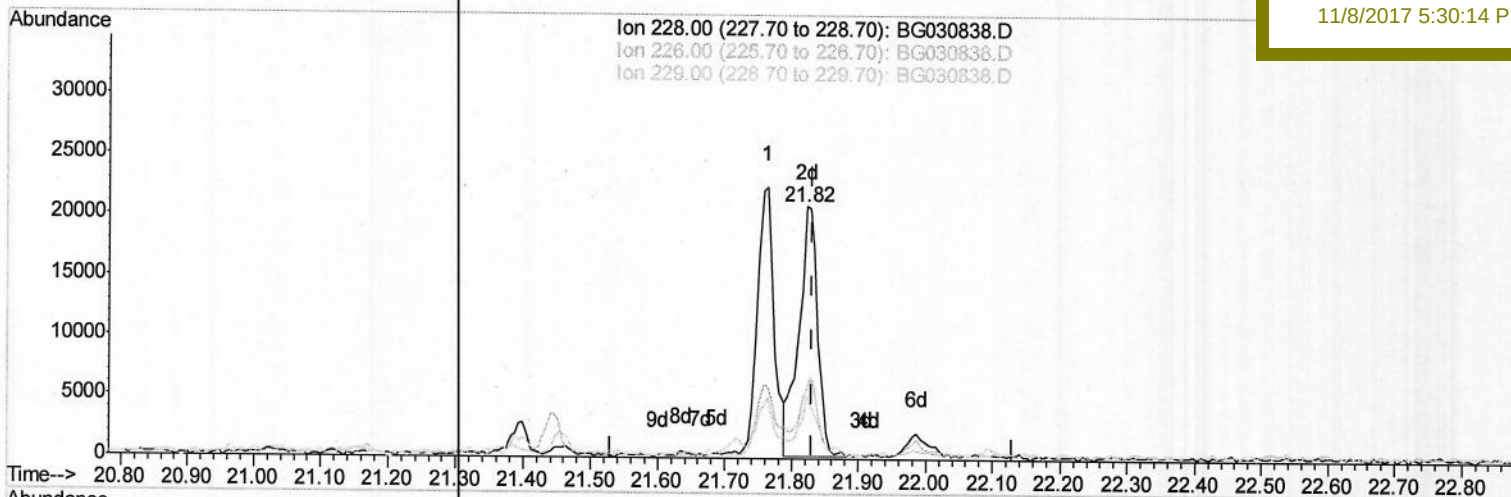
B0ET2

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11/8/2017 5:30:14 PM



(82) Chrysene

21.822min (-0.008) 1.22ng/ul m

> 5511/9/17

response 41160

Ion	Exp%	Act%
228.00	100	100
226.00	30.20	27.38
229.00	19.00	24.32#
0.00	0.00	0.00

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Quant Time: Nov 08 11:07:48 2017
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Instrument :
 BNA_G
 Client Sampled :
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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.15	152	82930	20.00	ng/ul	0.00
18) Naphthalene-d8	10.96	136	388511	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.77	164	250640	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.51	188	579152	20.00	ng/ul	0.00
75) Chrysene-d12	21.78	240	593949	20.00	ng/ul	0.00
83) Perylene-d12	25.09	264	611866	20.00	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.53	96	5087	2.49	ng/uL	0.00
5) Phenol-d5	7.32	99	119473	15.01	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.46	67	73707	14.80	ng/ul	0.00
9) 2-Chlorophenol-d4	7.69	132	100769	17.95	ng/ul	0.00
13) 4-Methylphenol-d8	8.86	113	85495	13.30	ng/ul	0.00
19) Nitrobenzene-d5	9.32	128	51309	18.40	ng/ul	0.00
22) 2-Nitrophenol-d4	10.04	143	63128	22.09	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.59	165	113447	17.42	ng/ul	0.00
29) 4-Chloroaniline-d4	11.10	131	108742	14.37	ng/ul	0.00
43) Dimethylphthalate-d6	14.17	166	372956	17.41	ng/ul	0.00
46) Acenaphthylene-d8	14.47	160	459242	17.62	ng/ul	0.00
51) 4-Nitrophenol-d4	14.98	143	52231	13.20	ng/ul	0.00
57) Fluorene-d10	15.75	176	346013	19.72	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.88	200	43780	15.49	ng/ul	0.00
70) Anthracene-d10	17.61	188	522949	19.09	ng/ul	0.00
76) Pyrene-d10	19.88	212	595338	19.36	ng/ul	0.00
87) Benzo(a)pyrene-d12	24.87	264	567815	19.59	ng/ul	0.00
Target Compounds						
6) Phenol	7.34	94	16128	1.958	ng/ul#	86
44) Dimethylphthalate	14.21	163	163843	7.839	ng/ul	98
69) Phenanthrene	17.55	178	31417	1.003	ng/ul	96
74) Fluoranthene	19.55	202	60475	1.728	ng/ul	99
77) Pvrene	19.91	202	56026	1.407	ng/ul	96
80) Benzo(a)anthracene	21.76	228	39379	1.058	ng/ul	97
82) Chrysene	21.82	228	41160m	1.217	ng/ul	93
85) Benzo(b)fluoranthene	24.04	252	64860	1.766	ng/ul#	93
88) Benzo(a)pvrene	24.93	252	46442	1.299	ng/ul#	93

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

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