

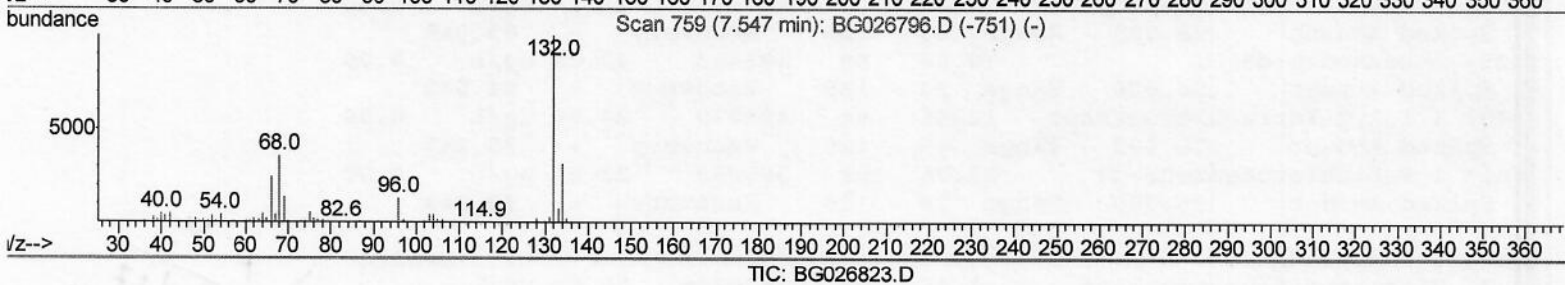
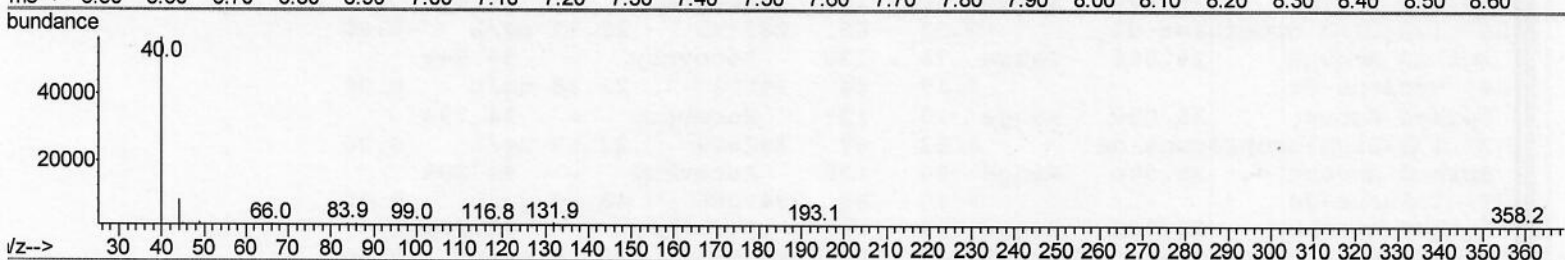
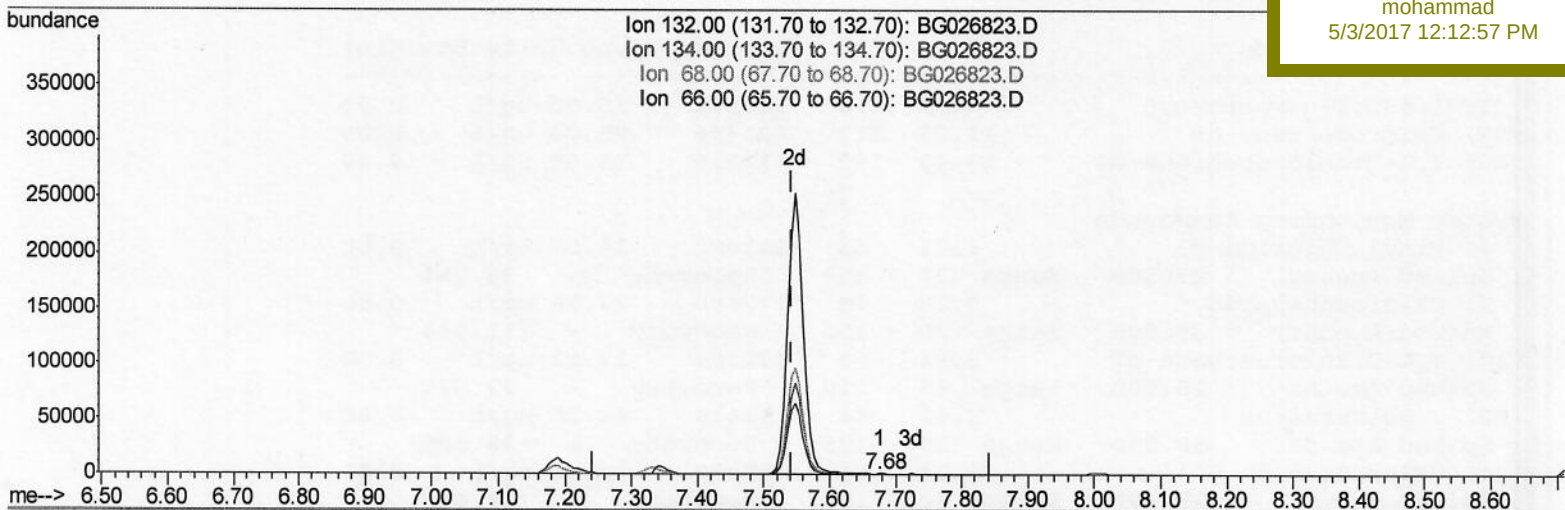
Data Path : Z:\HPCHEM1\BNA G\Data\BG050117\
 Data File : BG026823.D
 Acq On : 2 May 2017 1:46
 Operator : SJ/MA
 Sample : I2850-22
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Quant Time: May 15 15:39:18 2017
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM-EPA-BG042717.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed May 10 01:49:01 2017
 Response via : Initial Calibration

Instrument :
 BNA_G
 ClientSampleId :
 JHSZ2

Manual Integrations
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(9) 2-Chlorophenol-d4 (S)

7.676min (+0.133) 0.08ng/ul

response 981

Ion	Exp%	Act%
132.00	100	100
134.00	38.90	33.20
68.00	51.70	43.18
66.00	47.60	52.36

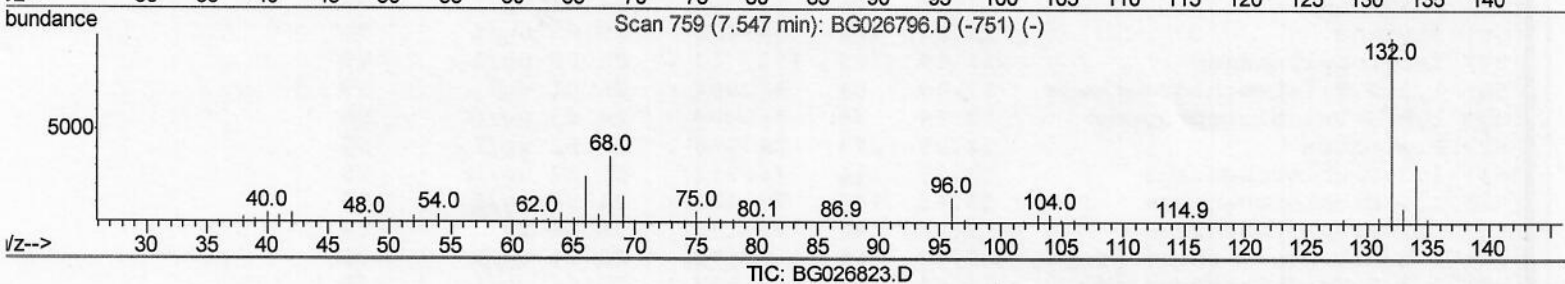
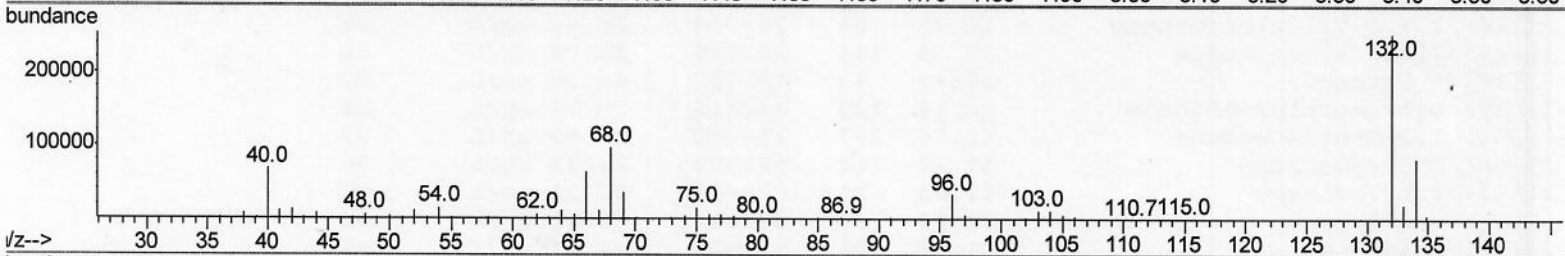
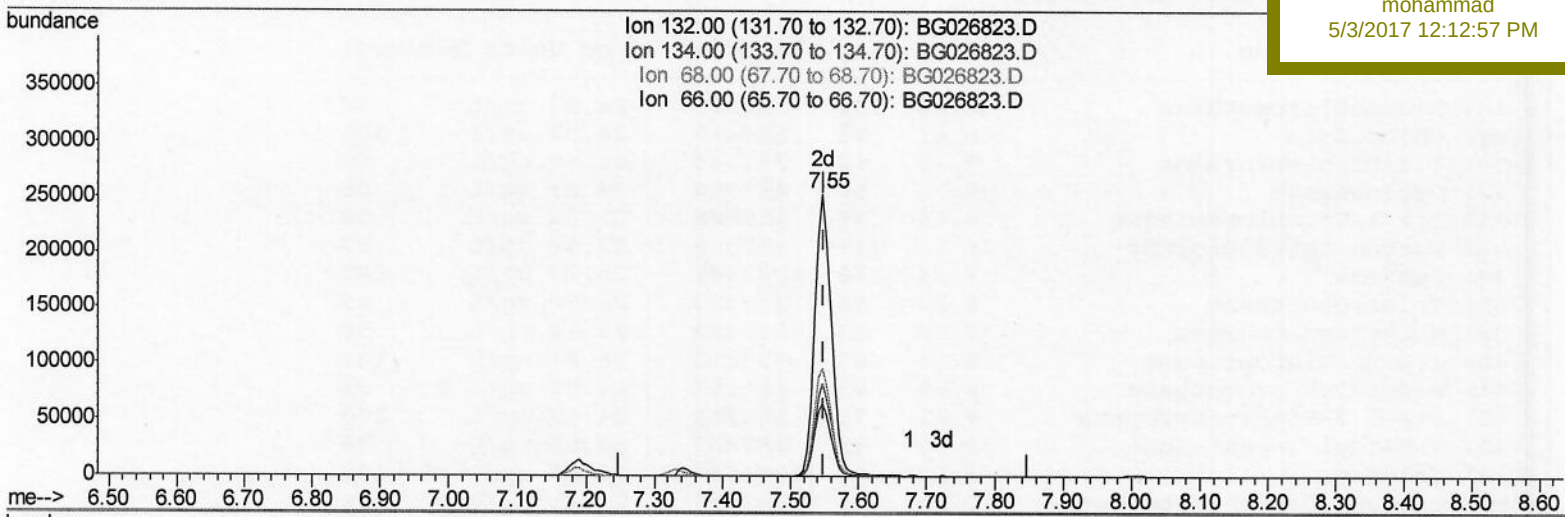
Data Path : Z:\HPCHEM1\BNA G\Data\BG050117\
 Data File : BG026823.D
 Acq On : 2 May 2017 1:46
 Operator : SJ/MA
 Sample : I2850-22
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Quant Time: May 02 03:33:23 2017
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM-EPA-BG042717.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue May 02 03:22:01 2017
 Response via : Initial Calibration

Instrument :
 BNA_G
 ClientSampled :
 JHSZ2

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TIC: BG026823.D

(9) 2-Chlorophenol-d4 (S)

7.547min (-0.002) 35.06ng/ul m > S J 5/5/17

response 427590

Ion	Exp%	Act%
132.00	100	100
134.00	38.90	32.23
68.00	51.70	37.78#
66.00	47.60	24.81#

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 Operator : SJ/MA
 Sample : I2850-22
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Quant Time: May 02 03:33:23 2017
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM-EPA-BG042717.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue May 02 03:22:01 2017
 Response via : Initial Calibration

Instrument :
 BNA_G
 ClientSampleId :
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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
1) 1,4-Dichlorobenzene-d4	8.01	152	161005	20.00	ng/ul	0.00
18) Naphthalene-d8	10.81	136	830508	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.62	164	480350	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.36	188	914538	20.00	ng/ul	0.00
75) Chrysene-d12	21.60	240	831720	20.00	ng/ul	0.00
83) Perylene-d12	24.77	264	872937	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev (Min)
3) 1,4-Dioxane-d8	3.43	96	23177	6.48	ng/uL	0.00
5) Phenol-d5	7.19	99	542379	37.75	ng/ul	0.00
7) Bis-(2-Chloroethyl) ether-d	7.33	67	320263	37.90	ng/ul	0.00
9) 2-Chlorophenol-d4	7.55	132	427590m	35.06	ng/ul	0.00
13) 4-Methylphenol-d8	8.72	113	460642	38.63	ng/ul	0.00
19) Nitrobenzene-d5	9.17	128	238929	34.42	ng/ul	0.00
22) 2-Nitrophenol-d4	9.89	143	254701	33.73	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.44	165	417887	34.83	ng/ul	0.00
29) 4-Chloroaniline-d4	10.94	131	614509	38.27	ng/ul	0.00
43) Dimethylphthalate-d6	14.02	166	1301819	33.67	ng/ul	0.00
46) Acenaphthylene-d8	14.32	160	1723893	35.50	ng/ul	0.00
51) 4-Nitrophenol-d4	14.84	143	187239	26.87	ng/ul	0.00
57) Fluorene-d10	15.61	176	1147326	34.00	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.73	200	144375	26.00	ng/ul	0.00
70) Anthracene-d10	17.45	188	1559440	35.14	ng/ul	0.00
76) Pyrene-d10	19.74	212	1518195	34.58	ng/ul	0.00
87) Benzo(a)pyrene-d12	24.55	264	1500820	36.36	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Ovalue
6) Phenol	7.22	94	25509	1.75	ng/ul#	79
44) Dimethylphthalate	14.07	163	217273	5.74	ng/ul	97

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

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