

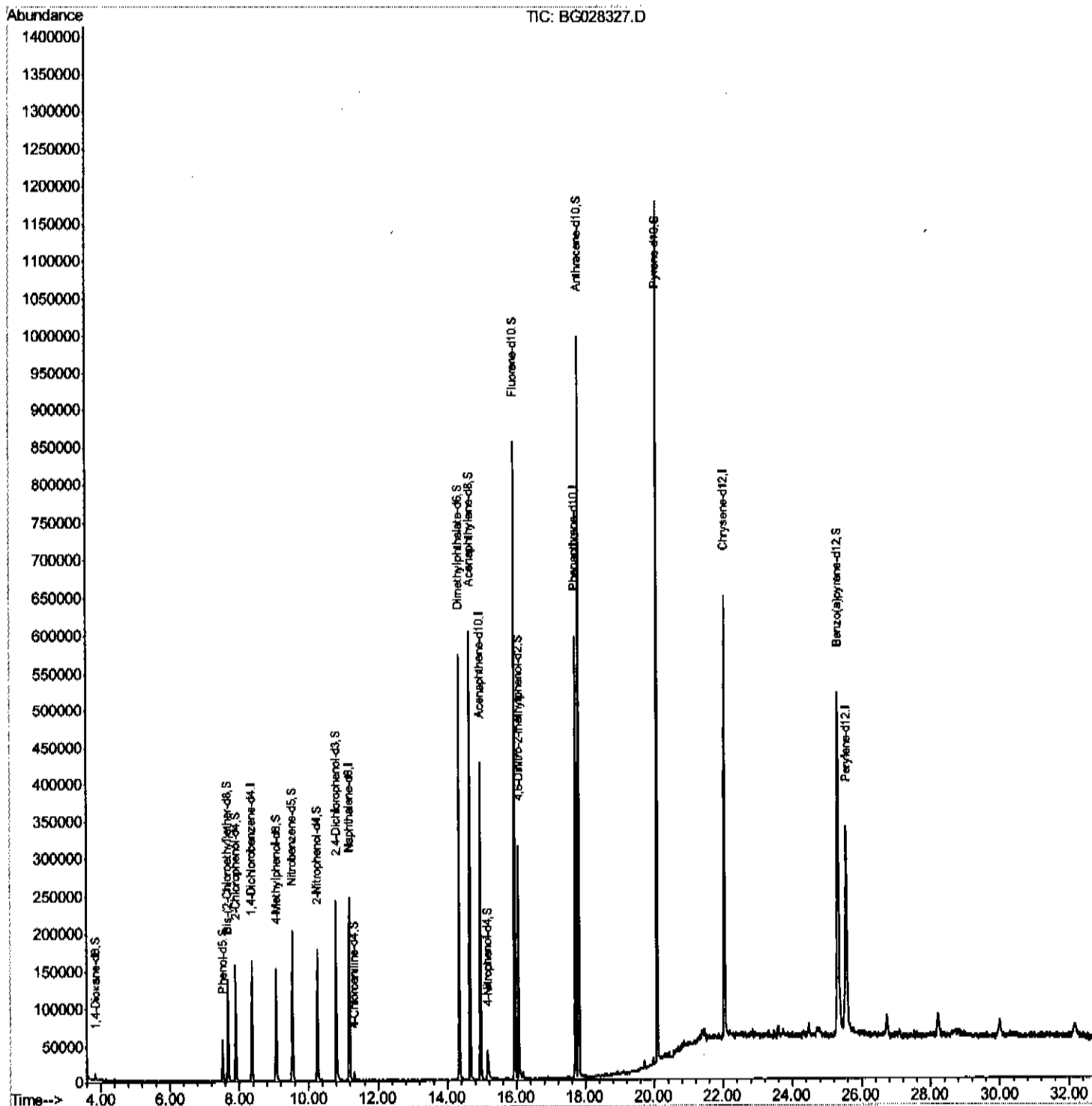
Data Path : Z:\HPCHEM1\BNA_G\DATA\BG081417\
Data File : BG028327.D
Acq On : 15 Aug 2017 00:43
Operator : SJ/JU
Sample : I4673-02
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0AJ5

Manual Integrations
APPROVED

Sohil
8/15/2017 6:01:54 PM

Quant Time: Aug 15 06:58:07 2017
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG080217.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Aug 15 06:41:43 2017
Response via : Initial Calibration



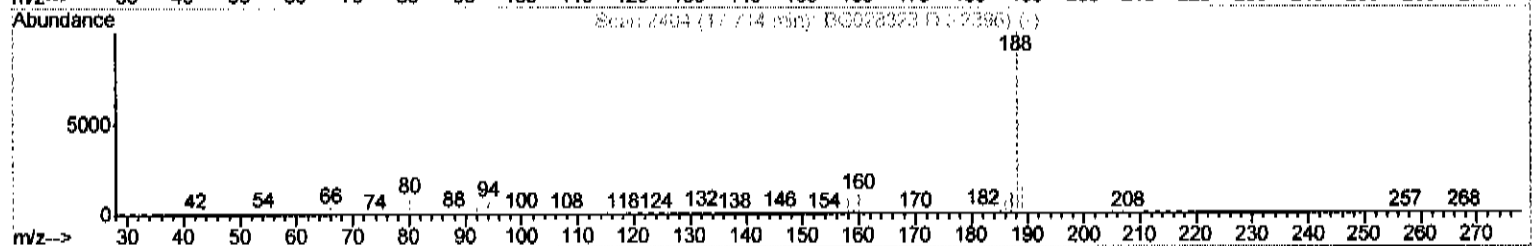
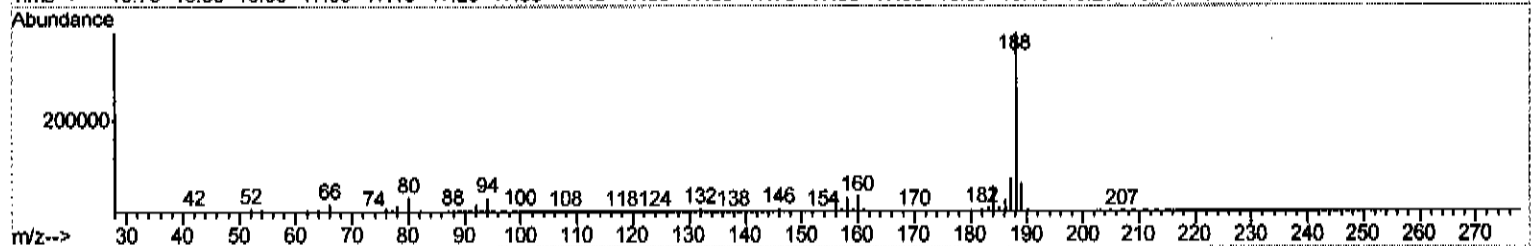
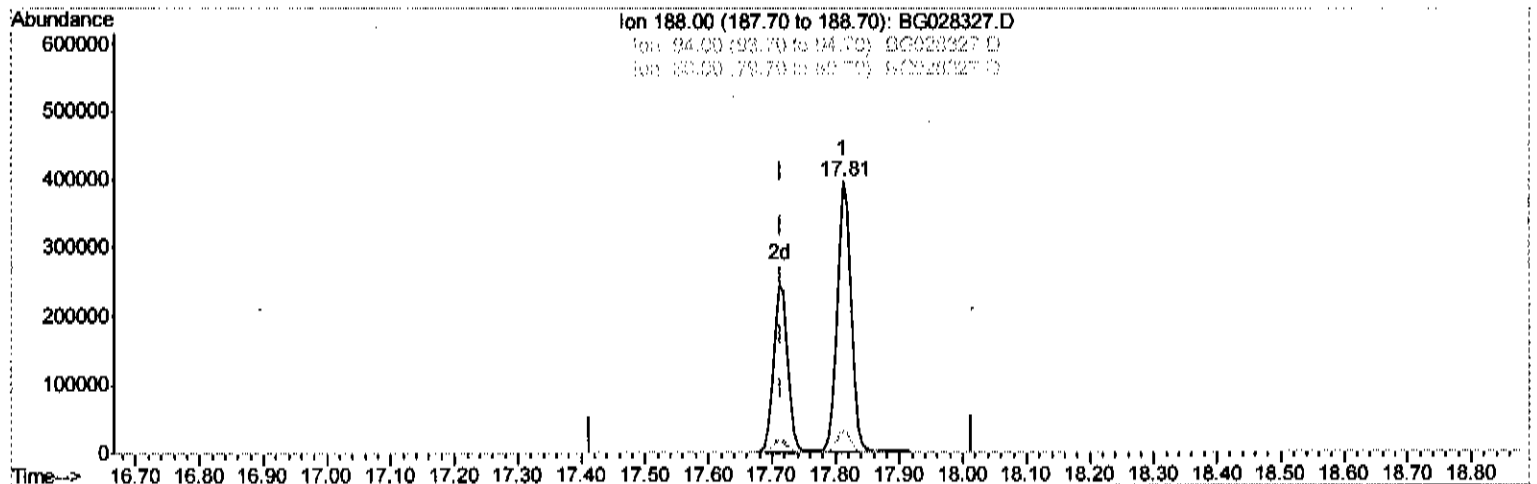
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Quant Time: Aug 15 06:42:45 2017
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TIC: BG028327.D

(61) Phenanthrene-d10 (I)

17.812min (+0.098) 20.00ng/ul

response 597286

Ion	Exp%	Act%
188.00	100	100
94.00	8.80	7.89
80.00	9.50	7.96
0.00	0.00	0.00

Quantitation Report (Qedit)

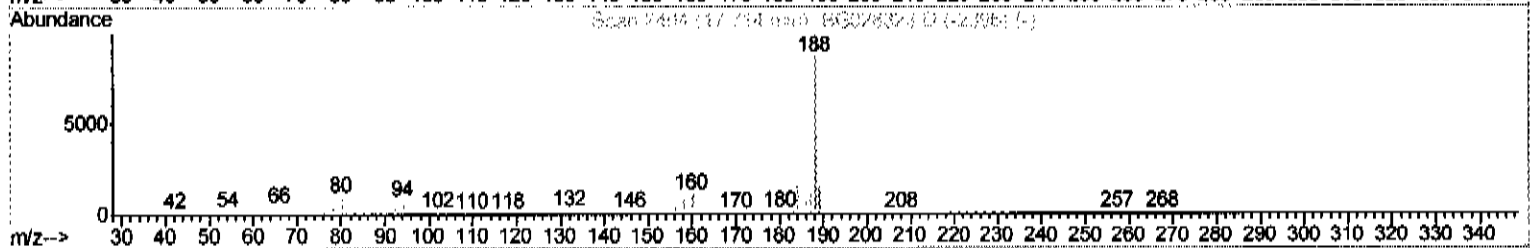
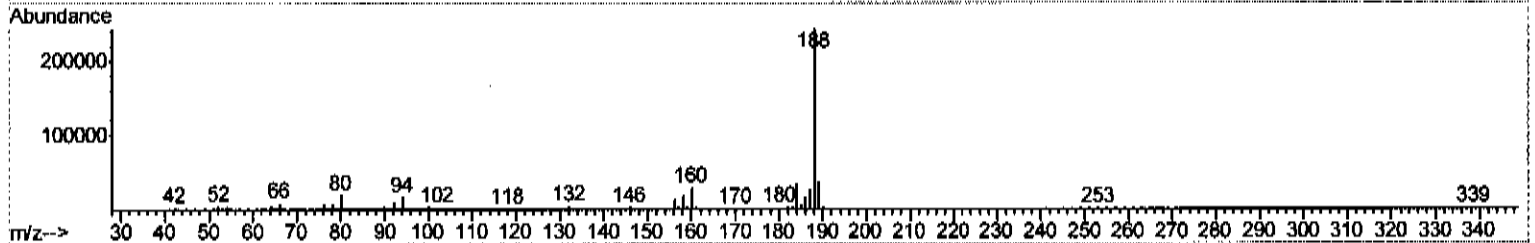
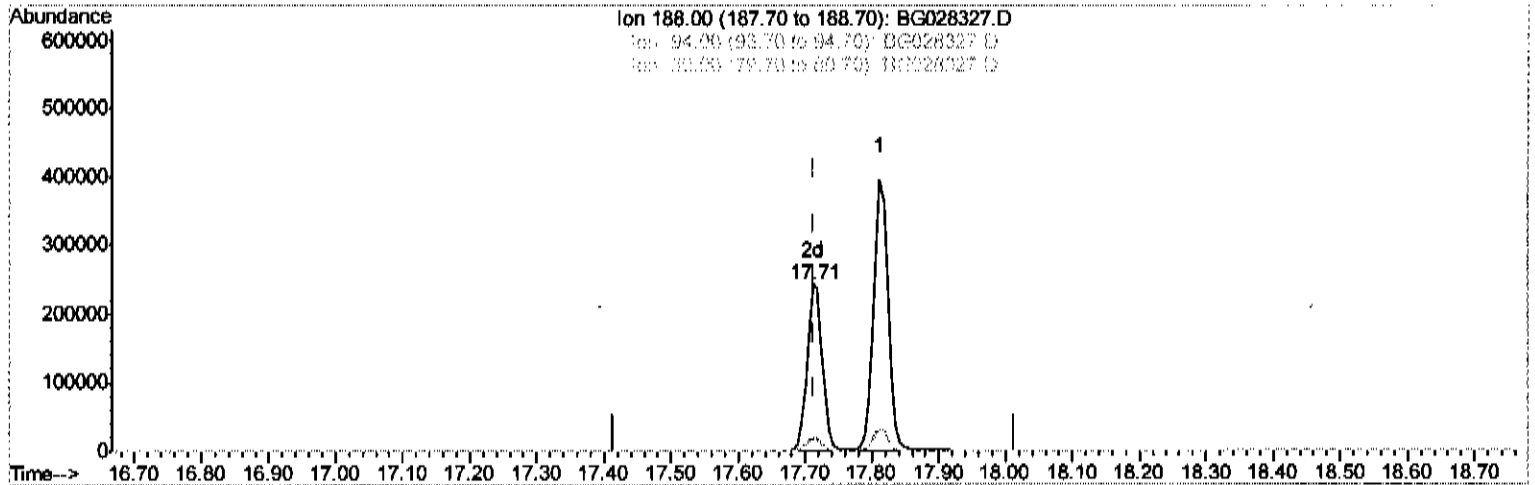
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Manual Integrations
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(61) Phenanthrene-d10 (I)

17.712min (-0.002) 20.00ng/ul m

response 383070

Ion	Exp%	Act%
188.00	100	100
94.00	8.80	6.60#
80.00	9.50	8.17
0.00	0.00	0.00

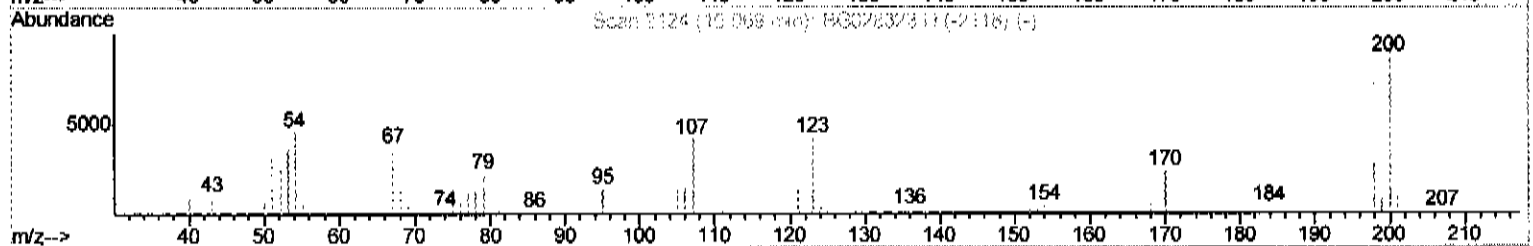
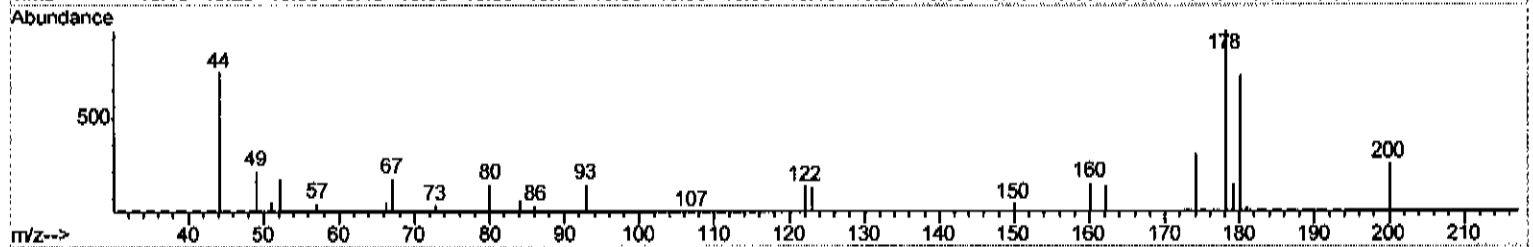
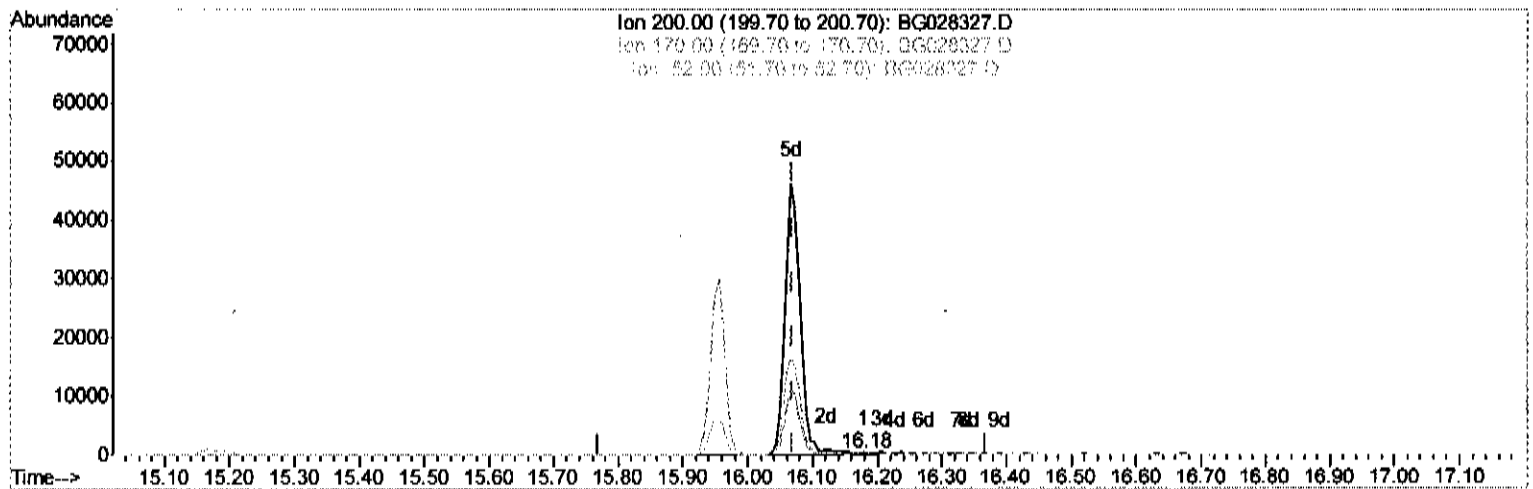
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ALS Vial : 13 Sample Multiplier: 1

Instrument :
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ClientSampleId :
C0AJ5

Manual Integrations
APPROVED

Sohil
8/15/2017 6:01:54 PM

Quant Time: Aug 15 06:57:43 2017
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TIC: BG028327.D

(62) 4,6-Dinitro-2-methylphenol-d2 (S)

16.179min (+0.110) 0.04ng/ul

response 110

Ion	Exp%	Act%
200.00	100	100
170.00	23.20	0.00#
52.00	47.20	82.48#
0.00	0.00	0.00

Quantitation Report (Qedit)

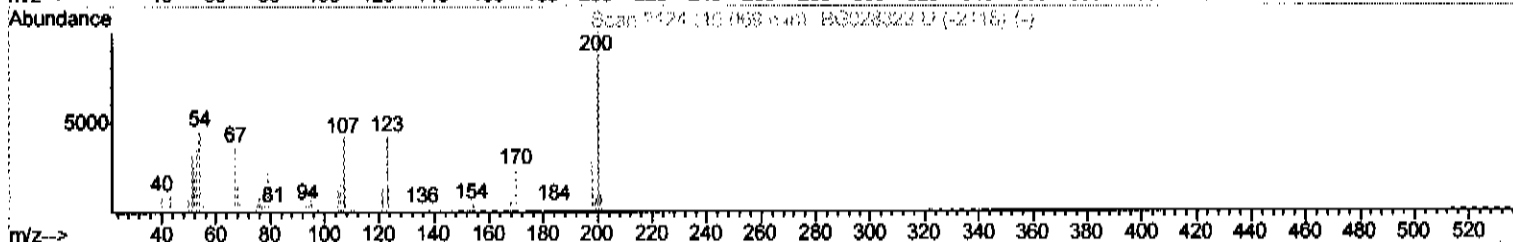
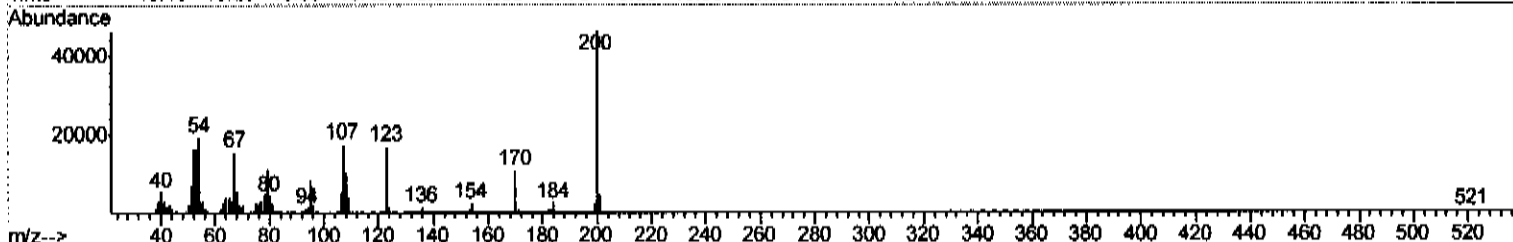
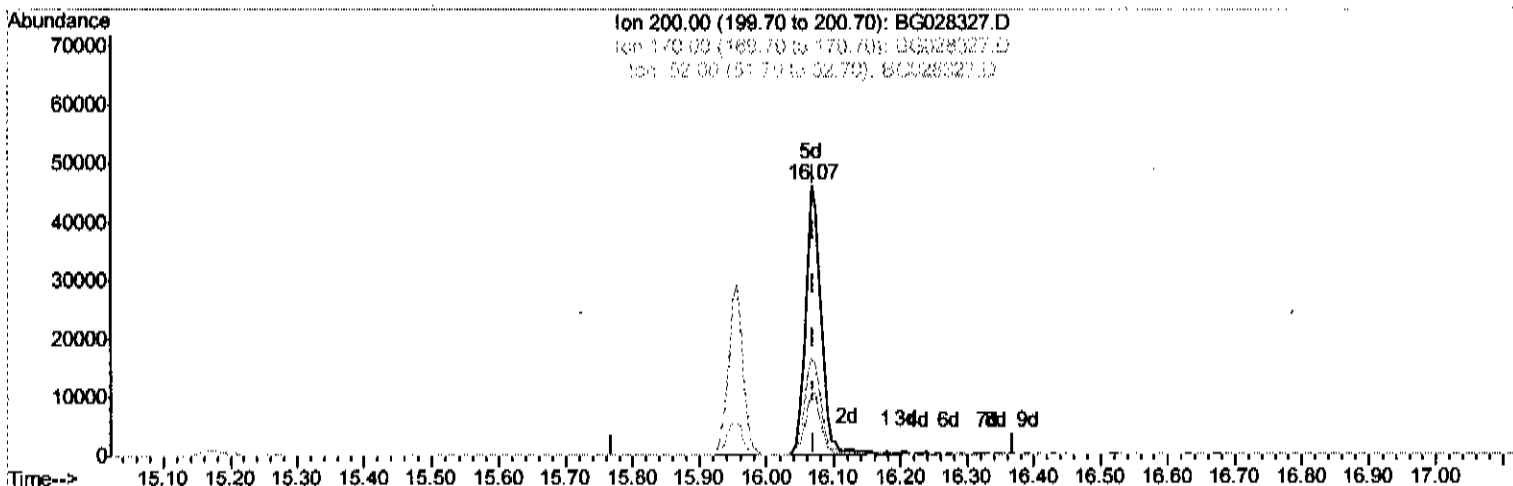
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 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
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Manual Integrations
 APPROVED

Sohil
 8/15/2017 6:01:54 PM

Quant Time: Aug 15 06:57:43 2017
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TIC: BG028327.D

(62) 4,6-Dinitro-2-methylphenol-d2 (S)

16.067min (-0.002) 29.05ng/ul m

response 72288

Ion	Exp%	Act%
200.00	100	100
170.00	23.20	23.68
52.00	47.20	36.08#
0.00	0.00	0.00

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.36	152	45458	20.00	ng/ul	0.00
18) Naphthalene-d8	11.18	136	210089	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.96	164	154564	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.71	188	383070m	20.00	ng/ul	0.00
75) Chrysene-d12	22.05	240	400577	20.00	ng/ul	0.00
83) Perylene-d12	25.56	264	369566	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.83	96	4175	4.28	ng/uL	0.00
5) Phenol-d5	7.51	99	33114	6.63	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.67	67	72975	23.08	ng/ul	0.00
9) 2-Chlorophenol-d4	7.89	132	74238	23.57	ng/ul	0.00
13) 4-Methylphenol-d8	9.06	113	60821	14.58	ng/ul	0.00
19) Nitrobenzene-d5	9.53	128	45303	27.58	ng/ul	0.00
22) 2-Nitrophenol-d4	10.25	143	55599	30.87	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.80	165	100772	27.04	ng/ul	0.00
29) 4-Chloroaniline-d4	11.31	131	6832	1.65	ng/ul	0.00
43) Dimethylphthalate-d6	14.35	166	397290	28.90	ng/ul	0.00
46) Acenaphthylene-d8	14.66	160	444491	28.68	ng/ul	0.00
51) 4-Nitrophenol-d4	15.17	143	15047	6.88	ng/ul	0.00
57) Fluorene-d10	15.96	176	350238	28.40	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	16.07	200	72288m	29.05	ng/ul	0.00
70) Anthracene-d10	17.81	188	597286	32.52	ng/ul	0.00
76) Pyrene-d10	20.09	212	656058	31.94	ng/ul	0.00
87) Benzo(a)pyrene-d12	25.32	264	560379	32.39	ng/ul	0.00

Target Compounds Qvalue

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