

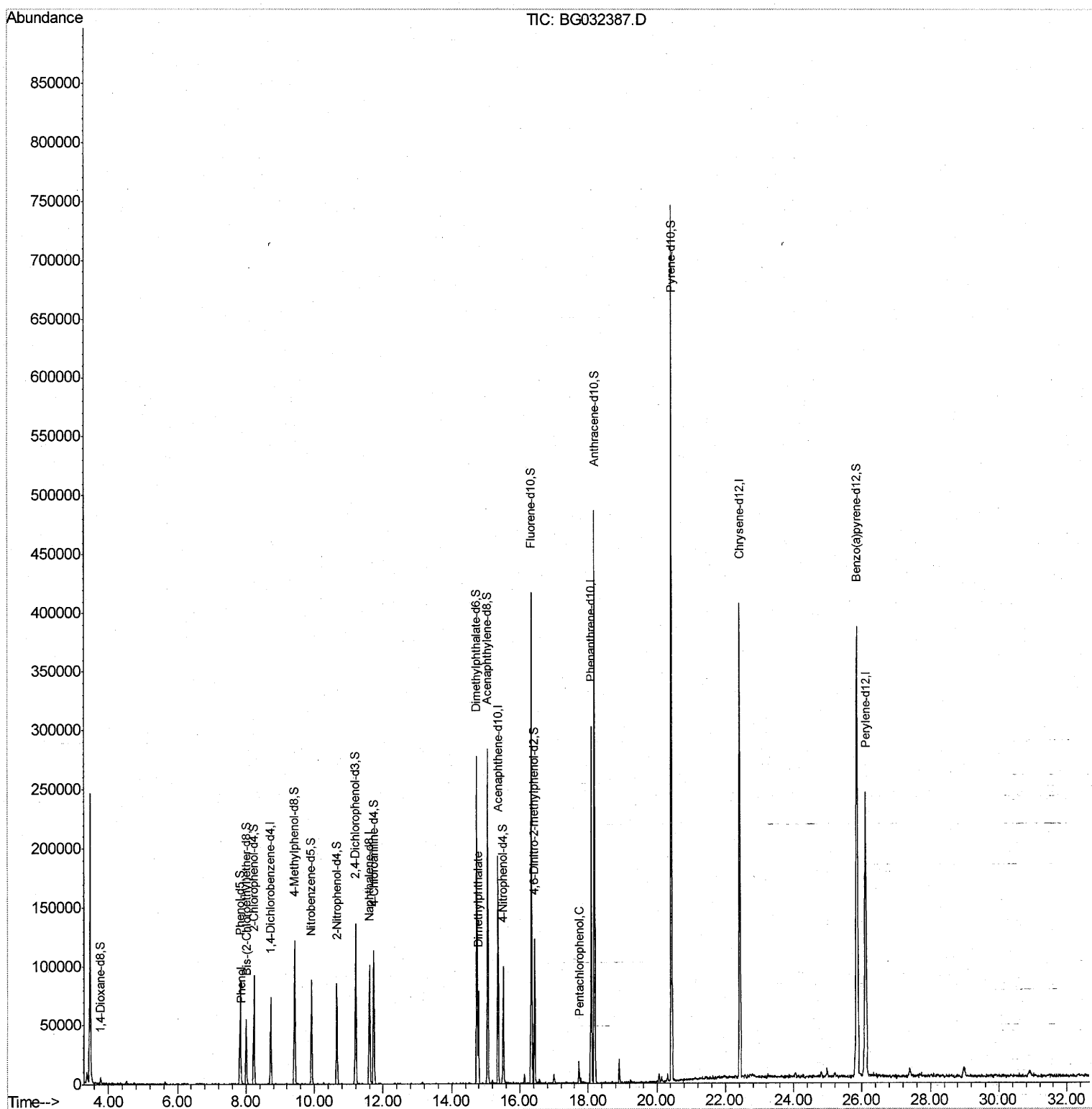
Data Path : Z:\HPCHEM1\BNA G\DATA\BG012818\
Data File : BG032387.D
Acq On : 28 Jan 2018 14:01
Operator : SJ/JU
Sample : J1296-02
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampled :
F6C20

Manual Integrations
APPROVED

Sohil
1/29/2018 5:49:55 PM

Quant Time: Jan 29 05:44:24 2018
Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM-EPA-BG012218.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Jan 29 04:59:53 2018
Response via : Initial Calibration



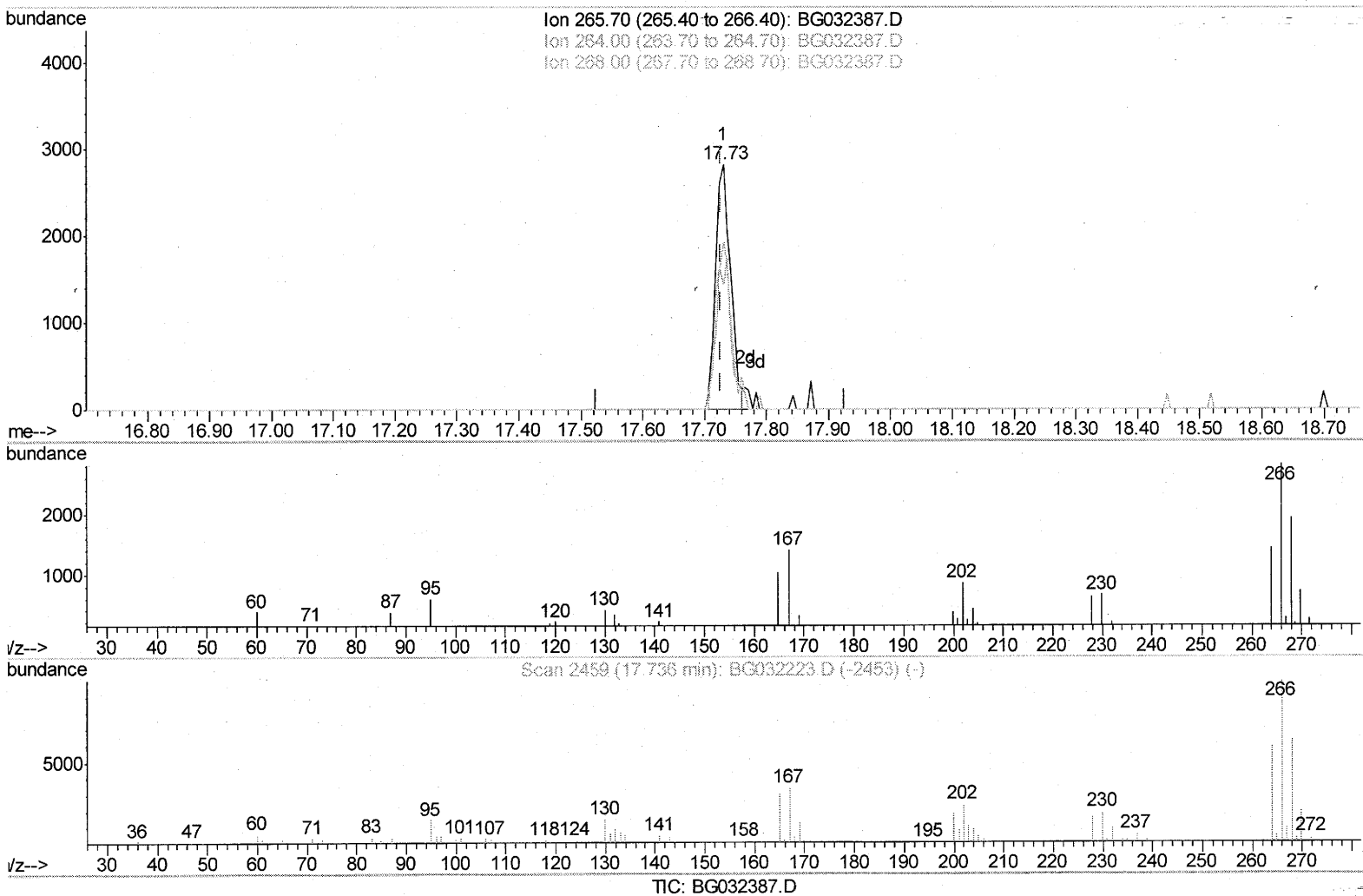
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 Response via : Initial Calibration



(68) Pentachlorophenol (C)

17.730min (+0.004) 3.10ng/ul

response 4706

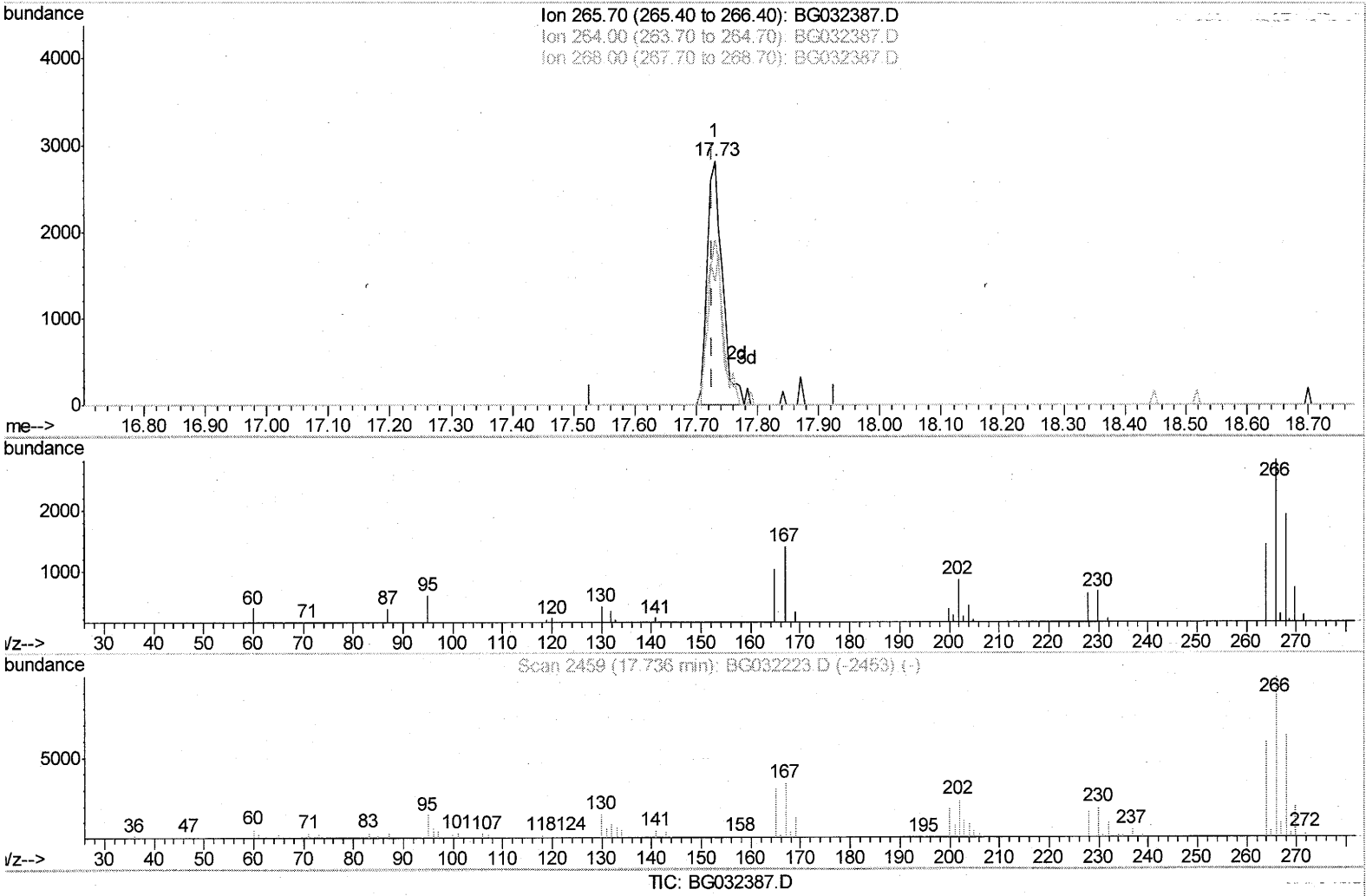
Ion	Exp%	Act%
265.70	100	100
264.00	60.30	51.08
268.00	65.40	75.19
0.00	0.00	0.00

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 Response via : Initial Calibration



(68) Pentachlorophenol (C)

17.730min (+0.004) 3.21ng/ul m

response 4871

Ion	Exp%	Act%
265.70	100	100
264.00	60.30	51.08
268.00	65.40	68.34
0.00	0.00	0.00

SJ/JU

Data Path : Z:\HPCHEM1\BNA G\DATA\BG012818\
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Manual Integrations
APPROVED

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.72	152	25321	20.00	ng/ul	0.00
18) Naphthalene-d8	11.60	136	99048	20.00	ng/ul	0.00
35) Acenaphthene-d10	15.36	164	71425	20.00	ng/ul	0.00
61) Phenanthrene-d10	18.09	188	203587	20.00	ng/ul	0.00
77) Chrysene-d12	22.42	240	307313	20.00	ng/ul	0.00
85) Perylene-d12	26.10	264	321843	20.00	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.78	96	2385	5.70	ng/uL	0.00
5) Phenol-d5	7.83	99	51269	29.47	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	8.01	67	23647	28.81	ng/ul	0.00
9) 2-Chlorophenol-d4	8.24	132	50412	31.05	ng/ul	0.00
13) 4-Methylphenol-d8	9.42	113	49317	31.22	ng/ul	0.00
19) Nitrobenzene-d5	9.92	128	24478	33.91	ng/ul	0.00
22) 2-Nitrophenol-d4	10.66	143	29779	34.42	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	11.20	165	60631	31.92	ng/ul	0.00
29) 4-Chloroaniline-d4	11.73	131	67847	51.24	ng/ul	0.00
43) Dimethylphthalate-d6	14.73	166	204343	34.27	ng/ul	0.00
46) Acenaphthylene-d8	15.06	160	238602	32.64	ng/ul	0.00
51) 4-Nitrophenol-d4	15.51	143	27782	34.96	ng/ul	0.00
57) Fluorene-d10	16.34	176	186663	33.62	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	16.43	200	37481	29.84	ng/ul	0.00
70) Anthracene-d10	18.19	188	315674	32.24	ng/ul	0.00
78) Pyrene-d10	20.43	212	451453	31.65	ng/ul	0.00
89) Benzo(a)pyrene-d12	25.84	264	496569	33.38	ng/ul	0.01
Target Compounds						
6) Phenol	7.85	94	3687	2.162	ng/ul	92
44) Dimethylphthalate	14.78	163	54421	9.508	ng/ul	98
68) Pentachlorophenol	17.73	266	4871m	3.208	ng/ul	

SS/29/18

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